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# Information Theory and Network Coding

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To my parents and my family

# Preface

Cover and Thomas wrote a book on information theory [72] ten years ago which covers most of the major topics with considerable depth. Their book has since become the standard textbook in the field, and it was no doubt a remarkable success. Instead of writing another comprehensive textbook on the subject, which has become more difficult as new results keep emerging, my goal is to write a book on the fundamentals of the subject in a unified and coherent manner.

During the last ten years, significant progress has been made in understanding the entropy function and information inequalities of discrete random variables. The results along this direction not only are of core interest in information theory, but also have applications in network coding theory, probability theory, group theory, Kolmogorov complexity, and possibly physics. This book is an up-to-date treatment of information theory for discrete random variables, which forms the foundation of the theory at large. There are eight chapters on classical topics (Chapters 1, 2, 3, 4, 5, 8, 9, and 10), five chapters on fundamental tools (Chapters 6, 7, 12, 13, and 14), and three chapters on selected topics (Chapters 11, 15, and 16). The chapters are arranged according to the logical order instead of the chronological order of the results in the literature.

# What is in this book

Out of the sixteen chapters in this book, the first thirteen chapters are basic topics, while the last three chapters are advanced topics for the more enthusiastic reader. A brief rundown of the chapters will give a better idea of what is in this book.

Chapter 1 is a very high level introduction to the nature of information theory and the main results in Shannon's original paper in 1948 which founded the field. There are also pointers to Shannon's biographies and his works.

Chapter 2 introduces Shannon's information measures and their basic properties. Useful identities and inequalities in information theory are derived and explained. Extra care is taken in handling joint distributions with

## VIII Preface

zero probability masses. The chapter ends with a section on the entropy rate of a stationary information source.

Chapter 3 is a discussion of zero-error data compression by uniquely decodable codes, with prefix codes as a special case. A proof of the entropy bound for prefix codes which involves neither the Kraft inequality nor the fundamental inequality is given. This proof facilitates the discussion of the redundancy of prefix codes.

Chapter 4 is a thorough treatment of weak typicality. The weak asymptotic equipartition property and the source coding theorem are discussed. An explanation of the fact that a good data compression scheme produces almost i.i.d. bits is given. There is also a brief discussion of the Shannon-McMillan-Breiman theorem.

Chapter 5 introduces a new definition of strong typicality which does not involve the cardinalities of the alphabet sets. The treatment of strong typicality here is more detailed than Berger [27] but less abstract than Csiszár and Körner [75]. A new exponential convergence result is proved in Theorem 5.3.

Chapter 6 is an introduction to the theory of *I*-Measure which establishes a one-to-one correspondence between Shannon's information measures and set theory. A number of examples are given to show how the use of information diagrams can simplify the proofs of many results in information theory. Most of these examples are previously unpublished. In particular, Example 6.15 is a generalization of Shannon's perfect secrecy theorem.

Chapter 7 explores the structure of the *I*-Measure for Markov structures. Set-theoretic characterizations of full conditional independence and Markov random field are discussed. The treatment of Markov random field here is perhaps too specialized for the average reader, but the structure of the *I*-Measure and the simplicity of the information diagram for a Markov chain is best explained as a special case of a Markov random field.

Chapter 8 consists of a new treatment of the channel coding theorem. Specifically, a graphical model approach is employed to explain the conditional independence of random variables. Great care is taken in discussing feedback.

Chapter 9 is an introduction to rate-distortion theory. The version of the rate-distortion theorem here, proved by using strong typicality, is a stronger version of the original theorem obtained by Shannon.

In Chapter 10, the Blahut-Arimoto algorithms for computing channel capacity and the rate-distortion function are discussed, and a simplified proof for convergence is given. Great care is taken in handling distributions with zero probability masses.

Chapter 11 is an introduction to network coding theory. The surprising fact that coding at the intermediate nodes can improve the throughput when an information source is multicast in a point-to-point network is explained. The max-flow bound for network coding with a single information source is explained in detail. Multi-source network coding will be discussed in Chapter 15 after the necessary tools are developed in the next three chapters. Information inequalities are sometimes called the laws of information theory because they govern the impossibilities in information theory. In Chapter 12, the geometrical meaning of information inequalities and the relation between information inequalities and conditional independence are explained in depth. The framework for information inequalities discussed here is the basis of the next two chapters.

Chapter 13 explains how the problem of proving information inequalities can be formulated as a linear programming problem. This leads to a complete characterization of all information inequalities which can be proved by conventional techniques. These are called Shannon-type inequalities, which can now be proved by the software ITIP which comes with this book. It is also shown how Shannon-type inequalities can be used to tackle the implication problem of conditional independence in probability theory.

All information inequalities we used to know were Shannon-type inequalities. Recently, a few non-Shannon-type inequalities have been discovered. This means that there exist laws in information theory beyond those laid down by Shannon. These inequalities and their applications are explained in depth in Chapter 14.

Network coding theory is further developed in Chapter 15. The situation when more than one information source are multicast in a point-to-point network is discussed. The surprising fact that a multi-source problem is not equivalent to a few single-source problems even when the information sources are mutually independent is clearly explained. Implicit and explicit bounds on the achievable coding rate region are discussed. These characterizations on the achievable coding rate region involve almost all the tools that have been developed earlier in the book, in particular, the framework for information inequalities.

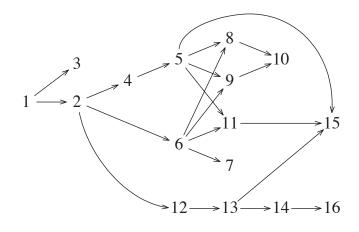
Chapter 16 explains an intriguing relation between information theory and group theory. Specifically, for every information inequality satisfied by any joint distribution, there is a corresponding group inequality satisfied by any finite group and its subgroups, and vice versa. Inequalities of the latter type govern the orders of any finite group and their subgroups. Group-theoretic proofs of Shannon-type information inequalities are given. At the end of this chapter, a group inequality is obtained from a non-Shannon-type inequality discussed in Chapter 14. The meaning and the implication of this inequality are yet to be understood.

#### How to use this book

You are recommended to read the chapters according to the above chart. However, you will not have too much difficulty jumping around in the book because there should be sufficient references to the previous relevant sections.

As a relatively slow thinker, I feel uncomfortable whenever I do not reason in the most explicit way. This probably has helped in writing this book, in which all the derivations are from the first principle. In the book, I try to explain all the subtle mathematical details without sacrificing the big picture.

# X Preface



Interpretations of the results are usually given before the proofs are presented. The book also contains a large number of examples. Unlike the examples in most books which are supplementary, the examples in this book are essential.

This book can be used as a reference book or a textbook. For a twosemester course on information theory, this would be a suitable textbook for the first semester. This would also be a suitable textbook for a one-semester course if only information theory for discrete random variables is covered. If the instructor also wants to include topics on continuous random variables, this book can be used as a textbook or a reference book in conjunction with another suitable textbook. The instructor will find this book a good source for homework problems because many problems here do not appear in any other textbook. A comprehensive instructor's manual is available upon request. Please contact the author at whyeung@ie.cuhk.edu.hk for information and access.

Just like any other lengthy document, this book for sure contains errors and omissions. To alleviate the problem, an errata will be maintained at the book homepage http://www.ie.cuhk.edu.hk/IT\_book/.

Hong Kong, August 2007 Raymond W. Yeung

# Contents

1 The Science of Information	1
------------------------------	---

# Part I Components of Information Theory

<b>2</b>	Info	rmation Measures	7
	2.1	Independence and Markov Chains	7
	2.2	Shannon's Information Measures	12
	2.3	Continuity of Shannon's Information Measures for Fixed	
		Finite Alphabets	18
	2.4	Chain Rules	20
	2.5	Informational Divergence	22
	2.6	The Basic Inequalities	26
	2.7	Some Useful Information Inequalities	28
	2.8	Fano's Inequality	31
	2.9	Maximum Entropy Distributions	35
	2.10	Entropy Rate of Stationary Source	
	App	endix 2.A: Approximation of Random Variables with	
		Countably Infinite Alphabets by Truncation	40
	Prob	blems	42
	Hist	orical Notes	46
3	$Th\epsilon$	I-Measure	49
	3.1	Preliminaries	50
	3.2	The <i>I</i> -Measure for Two Random Variables	
	3.3	Construction of the <i>I</i> -Measure $\mu^*$	
	3.4	$\mu^*$ Can be Negative	
	3.5	Information Diagrams	59
	3.6	Examples of Applications	66
	App	endix 3.A: A Variation of the Inclusion-Exclusion Formula	73
		blems	75

XII	Contents
	Historical Notes
4	Zero-Error Data Compression794.1 The Entropy Bound804.2 Prefix Codes834.2.1 Definition and Existence834.2.2 Huffman Codes864.3 Redundancy of Prefix Codes91Problems95Historical Notes97
5	Weak Typicality995.1 The Weak AEP995.2 The Source Coding Theorem1025.3 Efficient Source Coding1045.4 The Shannon-McMillan-Breiman Theorem105Problems108Historical Notes110
6	Strong Typicality1116.1Strong AEP1116.2Strong Typicality Versus Weak Typicality1196.3Joint Typicality1206.4An Interpretation of the Basic Inequalities129Problems129Historical Notes132
7	Discrete Memoryless Channels1337.1 Definition and Capacity1367.2 The Channel Coding Theorem1457.3 The Converse1477.4 Achievability1537.5 A Discussion1607.6 Feedback Capacity1637.7 Separation of Source and Channel Coding168Problems172Historical Notes176
8	Rate-Distortion Theory1798.1Single-Letter Distortion Measures1808.2The Rate-Distortion Function $R(D)$ 1838.3The Rate-Distortion Theorem1888.4The Converse1968.5Achievability of $R_I(D)$ 198Problems203

	Historical Notes
9	The Blahut-Arimoto Algorithms2079.1Alternating Optimization2089.2The Algorithms2109.2.1Channel Capacity2109.2.2The Rate-Distortion Function2159.3Convergence2189.3.1A Sufficient Condition2189.3.2Convergence to the Channel Capacity222
	Problems
10	Differential Entropy22510.1 Preliminaries22710.2 Definition23010.3 Joint and Conditional Differential Entropy23410.4 The AEP for Continuous Random Variables24210.5 Informational Divergence24510.6 Maximum Differential Entropy Distributions246Problems249Historical Notes250
11	Continuous-Valued Channels25111.1 Discrete-Time Channels25111.2 The Channel Coding Theorem25411.3 Proof of the Channel Coding Theorem256
	11.3.1 The Converse25611.3.2 Achievability25911.4 Memoryless Gaussian Channels26411.5 Parallel Gaussian Channels26611.6 Correlated Gaussian Channels27111.7 The Bandlimited White Gaussian Channel27311.8 The Bandlimited Colored Gaussian Channel28111.9 Zero-Mean Gaussian Noise is the Worst Additive Noise283Problems287Historical Notes289

13	Information Inequalities	
	13.1 The Region $\Gamma_n^*$	315
	13.2 Information Expressions in Canonical Form	316
	13.3 A Geometrical Framework	319
	13.3.1 Unconstrained Inequalities	319
	13.3.2 Constrained Inequalities	320
	13.3.3 Constrained Identities	321
	13.4 Equivalence of Constrained Inequalities	322
	13.5 The Implication Problem of Conditional Independence	326
	Problems	327
	Historical Notes	327
14	Shannon-Type Inequalities	220
14	14.1 The Elemental Inequalities	
	14.1 The Elemental Inequalities	
	14.2.1 Unconstrained Inequalities	
	14.2.1 Constrained Inequalities	
	14.3 A Duality	
	14.4 Machine Proving – ITIP	
	14.5 Tackling the Implication Problem	
	14.6 Minimality of the Elemental Inequalities	
		040
	Appendix 14.A: The Basic Inequalities and the Polymatroidal	
	Appendix 14.A: The Basic Inequalities and the Polymatroidal Axioms	346
	Appendix 14.A: The Basic Inequalities and the Polymatroidal	$\frac{346}{347}$
	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes	346 347 349
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities	346 347 349 351
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$	346 347 349 351 351
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality	346 347 349 351 351 359
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality	346 347 349 351 351 359 364
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications	346 347 349 351 351 359 364 371
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems	346 347 349 351 351 359 364 371 373
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications	346 347 349 351 351 359 364 371 373
15	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems	346 347 349 351 351 359 364 371 373 374
	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems         Historical Notes	346 347 349 351 351 359 364 371 373 374 377
	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         Beyond Shannon-Type Inequalities         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems         Historical Notes         15.4 Applications         Problems         Historical Notes         15.1 Group Preliminaries	346 347 349 351 351 359 364 371 373 374 377 378
	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes <b>Beyond Shannon-Type Inequalities</b> 15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems         Historical Notes <b>Entropy and Groups</b> 16.1 Group Preliminaries         16.2 Group-Characterizable Entropy Functions	346 347 349 351 359 364 371 373 374 377 378 383
	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes <b>Beyond Shannon-Type Inequalities</b> 15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems         Historical Notes <b>Entropy and Groups</b> 16.1 Group Preliminaries         16.2 Group-Characterizable Entropy Functions         16.3 A Group Characterization of $\overline{\Gamma}_n^*$	346 347 349 351 359 364 371 373 374 377 378 383 388
	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes         15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems         Historical Notes         Berropy and Groups         16.1 Group Preliminaries         16.2 Group-Characterizable Entropy Functions         16.3 A Group Characterization of $\overline{\Gamma}_n^*$ 16.4 Information Inequalities and Group Inequalities	346 347 349 351 359 364 371 373 374 377 378 383 388 391
	Appendix 14.A: The Basic Inequalities and the Polymatroidal         Axioms         Problems         Historical Notes <b>Beyond Shannon-Type Inequalities</b> 15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$ 15.2 A Non-Shannon-Type Unconstrained Inequality         15.3 A Non-Shannon-Type Constrained Inequality         15.4 Applications         Problems         Historical Notes <b>Entropy and Groups</b> 16.1 Group Preliminaries         16.2 Group-Characterizable Entropy Functions         16.3 A Group Characterization of $\overline{\Gamma}_n^*$	346 347 349 351 359 364 371 373 374 377 378 383 388 391 395

# Part II Fundamentals of Network Coding

17	Introduction       401         17.1 The Butterfly Network       402         17.2 Wireless and Satellite Communications       405         17.3 Source Separation       407         Problems       408         Historical Notes       410
18	The Max-Flow Bound41318.1 Point-to-Point Communication Networks41318.2 Examples Achieving the Max-Flow Bound41618.3 A Class of Network Codes41918.4 Proof of the Max-Flow Bound421Problems423Historical Notes425
19	Single-Source Linear Network Coding: Acyclic Networks
20	Single-Source Linear Network Coding: Cyclic Networks47520.1 Delay-Free Cyclic Networks47520.2 Convolutional Network Codes47820.3 Decoding of Convolutional Network Codes489Problems493Historical Notes493
21	Multi-Source Network Coding49521.1 The Max-Flow Bounds49521.2 Examples of Application49821.2.1 Multilevel Diversity Coding49821.2.2 Satellite Communication Network50021.3 A Network Code for Acyclic Networks50021.4 The Achievable Information Rate Region50221.5 Explicit Inner and Outer Bounds50521.6 The Converse50621.7 Achievability511

# XVI Contents

21.7.1 Random Code Construction	514
21.7.2 Performance Analysis	517
Problems	
Historical Notes	
References	
Index	

# The Science of Information

In a communication system, we try to convey information from one point to another, very often in a noisy environment. Consider the following scenario. A secretary needs to send facsimiles regularly and she wants to convey as much information as possible on each page. She has a choice of the font size, which means that more characters can be squeezed onto a page if a smaller font size is used. In principle, she can squeeze as many characters as desired on a page by using a small enough font size. However, there are two factors in the system which may cause errors. First, the fax machine has a finite resolution. Second, the characters transmitted may be received incorrectly due to noise in the telephone line. Therefore, if the font size is too small, the characters may not be recognizable on the facsimile. On the other hand, although some characters on the facsimile may not be recognizable, the recipient can still figure out the words from the context provided that the number of such characters is not excessive. In other words, it is not necessary to choose a font size such that all the characters on the facsimile are recognizable almost surely. Then we are motivated to ask: What is the maximum amount of meaningful information which can be conveyed on one page of facsimile?

This question may not have a definite answer because it is not very well posed. In particular, we do not have a precise measure of meaningful information. Nevertheless, this question is an illustration of the kind of fundamental questions we can ask about a communication system.

Information, which is not a physical entity but an abstract concept, is hard to quantify in general. This is especially the case if human factors are involved when the information is utilized. For example, when we play Beethoven's violin concerto from a compact disc, we receive the musical information from the loudspeakers. We enjoy this information because it arouses certain kinds of emotion within ourselves. While we receive the same information every time we play the same piece of music, the kinds of emotions aroused may be different from time to time because they depend on our mood at that particular moment. In other words, we can derive utility from the same information every time in a different way. For this reason, it is extremely difficult to devise 2 1 The Science of Information

a measure which can quantify the amount of information contained in a piece of music.

In 1948, Bell Telephone Laboratories scientist Claude E. Shannon (1916-2001) published a paper entitled "The Mathematical Theory of Communication" [291] which laid the foundation of an important field now known as information theory. In his paper, the model of a point-to-point communication system depicted in Figure 1.1 is considered. In this model, a message is gener-

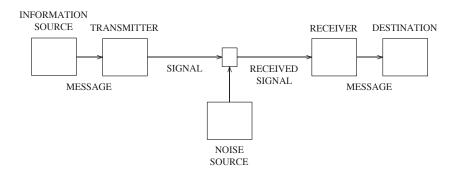


Fig. 1.1. Schematic diagram for a general point-to-point communication system.

ated by the information source. The message is converted by the transmitter into a signal which is suitable for transmission. In the course of transmission, the signal may be contaminated by a noise source, so that the received signal may be different from the transmitted signal. Based on the received signal, the receiver then makes an estimate of the message and deliver it to the destination.

In this abstract model of a point-to-point communication system, one is only concerned about whether the message generated by the source can be delivered correctly to the receiver without worrying about how the message is actually used by the receiver. In a way, Shannon's model does not cover all the aspects of a communication system. However, in order to develop a precise and useful theory of information, the scope of the theory has to be restricted.

In [291], Shannon introduced two fundamental concepts about 'information' from the communication point of view. First, information is *uncertainty*. More specifically, if a piece of information we are interested in is deterministic, then it has no value at all because it is already known with no uncertainty. From this point of view, for example, the continuous transmission of a still picture on a television broadcast channel is superfluous. Consequently, an information source is naturally modeled as a random variable or a random process, and probability is employed to develop the theory of information. Second, information to be transmitted is *digital*. This means that the information source should first be converted into a stream of 0's and 1's called bits, and the remaining task is to deliver these bits to the receiver correctly with no reference to their actual meaning. This is the foundation of all modern digital communication systems. In fact, this work of Shannon appears to contain the first published use of the term *bit*, which stands for <u>binary digit</u>.

In the same work, Shannon also proved two important theorems. The first theorem, called the *source coding theorem*, introduces *entropy* as the fundamental measure of information which characterizes the minimum rate of a source code representing an information source essentially free of error. The source coding theorem is the theoretical basis for *lossless* data compression<sup>1</sup>. The second theorem, called the *channel coding theorem*, concerns communication through a noisy channel. It was shown that associated with every noisy channel is a parameter, called the *capacity*, which is strictly positive except for very special channels, such that information can be communicated reliably through the channel as long as the information rate is less than the capacity. These two theorems, which give fundamental limits in point-to-point communication, are the two most important results in information theory.

In science, we study the laws of Nature which must be obeyed by any physical systems. These laws are used by engineers to design systems to achieve specific goals. Therefore, science is the foundation of engineering. Without science, engineering can only be done by trial and error.

In information theory, we study the fundamental limits in communication regardless of the technologies involved in the actual implementation of the communication systems. These fundamental limits are not only used as guidelines by communication engineers, but they also give insights into what optimal coding schemes are like. Information theory is therefore the science of information.

Since Shannon published his original paper in 1948, information theory has been developed into a major research field in both communication theory and applied probability. After more than half a century's research, it is quite impossible for a book on the subject to cover all the major topics with considerable depth. This book is a modern treatment of information theory for discrete random variables, which is the foundation of the theory at large. The book consists of two parts. The first part, namely Chapter 1 to Chapter 13, is a thorough discussion of the basic topics in information theory, including fundamental results, tools, and algorithms. The second part, namely Chapter 14 to Chapter 16, is a selection of advanced topics which demonstrate the use of the tools developed in the first part of the book. The topics discussed in this part of the book also represent new research directions in the field.

An undergraduate level course on probability is the only prerequisite for this book. For a non-technical introduction to information theory, we refer the reader to *Encyclopedia Britannica* [47]. In fact, we strongly recommend the reader to first read this excellent introduction before starting this book. For biographies of Claude Shannon, a legend of the 20th Century who had made

<sup>&</sup>lt;sup>1</sup> A data compression scheme is lossless if the data can be recovered with an arbitrarily small probability of error.

## 4 1 The Science of Information

fundamental contribution to the Information Age, we refer the readers to [53] and [307]. The latter is also a complete collection of Shannon's papers.

Unlike most branches of applied mathematics in which physical systems are studied, abstract systems of communication are studied in information theory. In reading this book, it is not unusual for a beginner to be able to understand all the steps in a proof but has no idea what the proof is leading to. The best way to learn information theory is to study the materials first and come back at a later time. Many results in information theory are rather subtle, to the extent that an expert in the subject may from time to time realize that his/her understanding of certain basic results has been inadequate or even incorrect. While a novice should expect to raise his/her level of understanding of the subject by reading this book, he/she should not be discouraged to find after finishing the book that there are actually more things yet to be understood. In fact, this is exactly the challenge and the beauty of information theory. **Components of Information Theory** 

Shannon's information measures refer to entropy, conditional entropy, mutual information, and conditional mutual information. They are the most important measures of information in information theory. In this chapter, we introduce these measures and establish some basic properties they possess. The physical meanings of these measures will be discussed in depth in subsequent chapters. We then introduce the informational divergence which measures the "distance" between two probability distributions and prove some useful inequalities in information theory. The chapter ends with a section on the entropy rate of a stationary information source.

# 2.1 Independence and Markov Chains

We begin our discussion in this chapter by reviewing two basic notions in probability: independence of random variables and Markov chain. All the random variables in this book are discrete unless otherwise specified.

Let X be a random variable taking values in an alphabet  $\mathcal{X}$ . The probability distribution for X is denoted as  $\{p_X(x), x \in \mathcal{X}\}$ , with  $p_X(x) = \Pr\{X = x\}$ . When there is no ambiguity,  $p_X(x)$  will be abbreviated as p(x), and  $\{p(x)\}$ will be abbreviated as p(x). The *support* of X, denoted by  $\mathcal{S}_X$ , is the set of all  $x \in \mathcal{X}$  such that p(x) > 0. If  $\mathcal{S}_X = \mathcal{X}$ , we say that p is *strictly positive*. Otherwise, we say that p is not strictly positive, or p contains zero probability masses. All the above notations naturally extend to two or more random variables. As we will see, probability distributions with zero probability masses are very delicate in general, and they need to be handled with great care.

**Definition 2.1.** Two random variables X and Y are independent, denoted by  $X \perp Y$ , if

$$p(x,y) = p(x)p(y)$$
(2.1)

for all x and y (i.e., for all  $(x, y) \in \mathcal{X} \times \mathcal{Y}$ ).

For more than two random variables, we distinguish between two types of independence.

**Definition 2.2 (Mutual Independence).** For  $n \ge 3$ , random variables  $X_1, X_2, \dots, X_n$  are mutually independent if

$$p(x_1, x_2, \cdots, x_n) = p(x_1)p(x_2)\cdots p(x_n)$$
 (2.2)

for all  $x_1, x_2, \cdots, x_n$ .

**Definition 2.3 (Pairwise Independence).** For  $n \ge 3$ , random variables  $X_1, X_2, \dots, X_n$  are pairwise independent if  $X_i$  and  $X_j$  are independent for all  $1 \le i < j \le n$ .

Note that mutual independence implies pairwise independence. We leave it as an exercise for the reader to show that the converse is not true.

**Definition 2.4 (Conditional Independence).** For random variables X, Y, and Z, X is independent of Z conditioning on Y, denoted by  $X \perp Z|Y$ , if

$$p(x, y, z)p(y) = p(x, y)p(y, z)$$
 (2.3)

for all x, y, and z, or equivalently,

$$p(x, y, z) = \begin{cases} \frac{p(x, y)p(y, z)}{p(y)} = p(x, y)p(z|y) & \text{if } p(y) > 0\\ 0 & \text{otherwise.} \end{cases}$$
(2.4)

The first definition of conditional independence above is sometimes more convenient to use because it is not necessary to distinguish between the cases p(y) > 0 and p(y) = 0. However, the physical meaning of conditional independence is more explicit in the second definition.

**Proposition 2.5.** For random variables X, Y, and  $Z, X \perp Z | Y$  if and only if

$$p(x, y, z) = a(x, y)b(y, z)$$
 (2.5)

for all x, y, and z such that p(y) > 0.

*Proof.* The 'only if' part follows immediately from the definition of conditional independence in (2.4), so we will only prove the 'if' part. Assume

$$p(x, y, z) = a(x, y)b(y, z)$$

$$(2.6)$$

for all x, y, and z such that p(y) > 0. Then for such x, y, and z, we have

$$p(x,y) = \sum_{z} p(x,y,z) = \sum_{z} a(x,y)b(y,z) = a(x,y)\sum_{z} b(y,z)$$
(2.7)

and

$$p(y,z) = \sum_{x} p(x,y,z) = \sum_{x} a(x,y)b(y,z) = b(y,z)\sum_{x} a(x,y).$$
(2.8)

Furthermore,

$$p(y) = \sum_{z} p(y, z) = \left(\sum_{x} a(x, y)\right) \left(\sum_{z} b(y, z)\right).$$
(2.9)

Therefore,

$$\frac{p(x,y)p(y,z)}{p(y)} = \frac{\left(a(x,y)\sum_{z}b(y,z)\right)\left(b(y,z)\sum_{x}a(x,y)\right)}{\left(\sum_{x}a(x,y)\right)\left(\sum_{z}b(y,z)\right)}$$
(2.10)

$$= a(x, y)o(y, z)$$
(2.11)  
=  $p(x, y, z).$  (2.12)

For x, y, and z such that p(y) = 0, since

$$0 \le p(x, y, z) \le p(y) = 0, \tag{2.13}$$

we have

$$p(x, y, z) = 0.$$
 (2.14)

Hence,  $X \perp Z | Y$ . The proof is accomplished.  $\Box$ 

**Definition 2.6 (Markov Chain).** For random variables  $X_1, X_2, \dots, X_n$ , where  $n \ge 3$ ,  $X_1 \rightarrow X_2 \rightarrow \dots \rightarrow X_n$  forms a Markov chain if

$$p(x_1, x_2, \cdots, x_n)p(x_2)p(x_3)\cdots p(x_{n-1}) = p(x_1, x_2)p(x_2, x_3)\cdots p(x_{n-1}, x_n)$$
(2.15)

for all  $x_1, x_2, \dots, x_n$ , or equivalently,

$$p(x_1, x_2, \cdots, x_n) = \begin{cases} p(x_1, x_2) p(x_3 | x_2) \cdots p(x_n | x_{n-1}) & \text{if } p(x_2), p(x_3), \cdots, p(x_{n-1}) > 0 \\ 0 & \text{otherwise.} \end{cases}$$
(2.16)

We note that  $X \perp Z | Y$  is equivalent to the Markov chain  $X \to Y \to Z$ .

**Proposition 2.7.**  $X_1 \to X_2 \to \cdots \to X_n$  forms a Markov chain if and only if  $X_n \to X_{n-1} \to \cdots \to X_1$  forms a Markov chain.

9

*Proof.* This follows directly from the symmetry in the definition of a Markov chain in (2.15).  $\Box$ 

In the following, we state two basic properties of a Markov chain. The proofs are left as an exercise.

**Proposition 2.8.**  $X_1 \to X_2 \to \cdots \to X_n$  forms a Markov chain if and only if

$$X_1 \to X_2 \to X_3$$

$$(X_1, X_2) \to X_3 \to X_4$$

$$\vdots$$

$$(X_1, X_2, \cdots, X_{n-2}) \to X_{n-1} \to X_n$$

$$(2.17)$$

form Markov chains.

**Proposition 2.9.**  $X_1 \to X_2 \to \cdots \to X_n$  forms a Markov chain if and only if

$$p(x_1, x_2, \cdots, x_n) = f_1(x_1, x_2) f_2(x_2, x_3) \cdots f_{n-1}(x_{n-1}, x_n)$$
(2.18)

for all  $x_1, x_2, \dots, x_n$  such that  $p(x_2), p(x_3), \dots, p(x_{n-1}) > 0$ .

Note that Proposition 2.9 is a generalization of Proposition 2.5. From Proposition 2.9, one can prove the following important property of a Markov chain. Again, the details are left as an exercise.

**Proposition 2.10 (Markov subchains).** Let  $\mathcal{N}_n = \{1, 2, \dots, n\}$  and let  $X_1 \to X_2 \to \dots \to X_n$  form a Markov chain. For any subset  $\alpha$  of  $\mathcal{N}_n$ , denote  $(X_i, i \in \alpha)$  by  $X_\alpha$ . Then for any disjoint subsets  $\alpha_1, \alpha_2, \dots, \alpha_m$  of  $\mathcal{N}_n$  such that

$$k_1 < k_2 < \dots < k_m \tag{2.19}$$

for all  $k_j \in \alpha_j$ ,  $j = 1, 2, \cdots, m$ ,

$$X_{\alpha_1} \to X_{\alpha_2} \to \dots \to X_{\alpha_m}$$
 (2.20)

forms a Markov chain. That is, a subchain of  $X_1 \to X_2 \to \cdots \to X_n$  is also a Markov chain.

*Example 2.11.* Let  $X_1 \to X_2 \to \cdots \to X_{10}$  form a Markov chain and  $\alpha_1 = \{1, 2\}, \alpha_2 = \{4\}, \alpha_3 = \{6, 8\}$ , and  $\alpha_4 = \{10\}$  be subsets of  $\mathcal{N}_{10}$ . Then Proposition 2.10 says that

$$(X_1, X_2) \to X_4 \to (X_6, X_8) \to X_{10}$$
 (2.21)

also forms a Markov chain.

We have been very careful in handling probability distributions with zero probability masses. In the rest of the section, we show that such distributions are very delicate in general. We first prove the following property of a strictly positive probability distribution involving four random variables<sup>1</sup>.

**Proposition 2.12.** Let  $X_1, X_2, X_3$ , and  $X_4$  be random variables such that  $p(x_1, x_2, x_3, x_4)$  is strictly positive. Then

$$X_1 \perp X_4 | (X_2, X_3) X_1 \perp X_3 | (X_2, X_4)$$
  $\Rightarrow X_1 \perp (X_3, X_4) | X_2.$  (2.22)

*Proof.* If  $X_1 \perp X_4 | (X_2, X_3)$ , we have

$$p(x_1, x_2, x_3, x_4) = \frac{p(x_1, x_2, x_3)p(x_2, x_3, x_4)}{p(x_2, x_3)}.$$
(2.23)

On the other hand, if  $X_1 \perp X_3 | (X_2, X_4)$ , we have

$$p(x_1, x_2, x_3, x_4) = \frac{p(x_1, x_2, x_4)p(x_2, x_3, x_4)}{p(x_2, x_4)}.$$
(2.24)

Equating (2.23) and (2.24), we have

$$p(x_1, x_2, x_3) = \frac{p(x_2, x_3)p(x_1, x_2, x_4)}{p(x_2, x_4)}.$$
(2.25)

Then

$$p(x_1, x_2) = \sum_{x_3} p(x_1, x_2, x_3)$$
(2.26)

$$=\sum_{x_3} \frac{p(x_2, x_3)p(x_1, x_2, x_4)}{p(x_2, x_4)}$$
(2.27)

$$=\frac{p(x_2)p(x_1, x_2, x_4)}{p(x_2, x_4)},$$
(2.28)

or

$$\frac{p(x_1, x_2, x_4)}{p(x_2, x_4)} = \frac{p(x_1, x_2)}{p(x_2)}.$$
(2.29)

Hence from (2.24),

$$p(x_1, x_2, x_3, x_4) = \frac{p(x_1, x_2, x_4)p(x_2, x_3, x_4)}{p(x_2, x_4)} = \frac{p(x_1, x_2)p(x_2, x_3, x_4)}{p(x_2)} \quad (2.30)$$

for all  $x_1, x_2, x_3$ , and  $x_4$ , i.e.,  $X_1 \perp (X_3, X_4) | X_2$ .  $\Box$ 

<sup>&</sup>lt;sup>1</sup> Proposition 2.12 is called the *intersection* axiom in Bayesian networks. See [258].

If  $p(x_1, x_2, x_3, x_4) = 0$  for some  $x_1, x_2, x_3$ , and  $x_4$ , i.e., p is not strictly positive, the arguments in the above proof are not valid. In fact, the proposition may not hold in this case. For instance, let  $X_1 = Y$ ,  $X_2 = Z$ , and  $X_3 = X_4 = (Y, Z)$ , where Y and Z are independent random variables. Then  $X_1 \perp X_4 | (X_2, X_3), X_1 \perp X_3 | (X_2, X_4), \text{ but } X_1 \not\perp (X_3, X_4) | X_2$ . Note that for this construction, p is not strictly positive because  $p(x_1, x_2, x_3, x_4) = 0$  if  $x_3 \neq (x_1, x_2)$  or  $x_4 \neq (x_1, x_2)$ .

The above example is somewhat counter-intuitive because it appears that Proposition 2.12 should hold for all probability distributions via a continuity argument. Specifically, such an argument goes like this. For any distribution p, let  $\{p_k\}$  be a sequence of strictly positive distributions such that  $p_k \to p$ and  $p_k$  satisfies (2.23) and (2.24) for all k, i.e.,

$$p_k(x_1, x_2, x_3, x_4)p_k(x_2, x_3) = p_k(x_1, x_2, x_3)p_k(x_2, x_3, x_4)$$
(2.31)

and

$$p_k(x_1, x_2, x_3, x_4)p_k(x_2, x_4) = p_k(x_1, x_2, x_4)p_k(x_2, x_3, x_4).$$
(2.32)

Then by the proposition,  $p_k$  also satisfies (2.30), i.e.,

$$p_k(x_1, x_2, x_3, x_4)p_k(x_2) = p_k(x_1, x_2)p_k(x_2, x_3, x_4).$$
(2.33)

Letting  $k \to \infty$ , we have

$$p(x_1, x_2, x_3, x_4)p(x_2) = p(x_1, x_2)p(x_2, x_3, x_4)$$
(2.34)

for all  $x_1, x_2, x_3$ , and  $x_4$ , i.e.,  $X_1 \perp (X_3, X_4) | X_2$ . Such an argument would be valid if there always exists a sequence  $\{p_k\}$  as prescribed. However, the existence of the distribution  $p(x_1, x_2, x_3, x_4)$  constructed immediately after Proposition 2.12 simply says that it is not always possible to find such a sequence  $\{p_k\}$ .

Therefore, probability distributions which are not strictly positive can be very delicate. For strictly positive distributions, we see from Proposition 2.5 that their conditional independence structures are closely related to the factorization problem of such distributions, which has been investigated by Chan [57].

# 2.2 Shannon's Information Measures

We begin this section by introducing the *entropy* of a random variable. As we will see shortly, all Shannon's information measures can be expressed as linear combinations of entropies.

**Definition 2.13.** The entropy H(X) of a random variable X is defined as

$$H(X) = -\sum_{x} p(x) \log p(x).$$
 (2.35)

In all definitions of information measures, we adopt the convention that summation is taken over the corresponding support. Such a convention is necessary because  $p(x) \log p(x)$  in (2.35) is undefined if p(x) = 0.

The base of the logarithm in (2.35) can be chosen to be any convenient real number great than 1. We write H(X) as  $H_{\alpha}(X)$  when the base of the logarithm is  $\alpha$ . When the base of the logarithm is 2, the unit for entropy is the *bit*. When the base of the logarithm is *e*, the unit for entropy is the *nat*. When the base of the logarithm is an integer  $D \ge 2$ , the unit for entropy is the *D-it* (*D*-ary digit). In the context of source coding, the base is usually taken to be the size of the code alphabet. This will be discussed in Chapter 4.

In computer science a bit means an entity which can take the value 0 or 1. In information theory the entropy of a random variable is measured in bits. The reader should distinguish these two meanings of a bit from each other carefully.

Let g(X) be any function of a random variable X. We will denote the expectation of g(X) by Eg(X), i.e.,

$$Eg(X) = \sum_{x} p(x)g(x), \qquad (2.36)$$

where the summation is over  $S_X$ . Then the definition of the entropy of a random variable X can be written as

$$H(X) = -E\log p(X). \tag{2.37}$$

Expressions of Shannon's information measures in terms of expectations will be useful in subsequent discussions.

The entropy H(X) of a random variable X is a functional of the probability distribution p(x) which measures the average amount of information contained in X, or equivalently, the average amount of *uncertainty* removed upon revealing the outcome of X. Note that H(X) depends only on p(x), not on the actual values in  $\mathcal{X}$ . Occasionally, we also denote H(X) by H(p).

For  $0 \leq \gamma \leq 1$ , define

$$h_b(\gamma) = -\gamma \log \gamma - (1 - \gamma) \log(1 - \gamma)$$
(2.38)

with the convention  $0 \log 0 = 0$ , so that  $h_b(0) = h_b(1) = 0$ . With this convention,  $h_b(\gamma)$  is continuous at  $\gamma = 0$  and  $\gamma = 1$ .  $h_b$  is called the *binary entropy function*. For a binary random variable X with distribution  $\{\gamma, 1 - \gamma\}$ ,

$$H(X) = h_b(\gamma). \tag{2.39}$$

Figure 2.1 shows the graph  $h_b(\gamma)$  versus  $\gamma$  in the base 2. Note that  $h_b(\gamma)$  achieves the maximum value 1 when  $\gamma = \frac{1}{2}$ .

The definition of the *joint entropy* of two random variables is similar to the definition of the entropy of a single random variable. Extension of this definition to more than two random variables is straightforward.

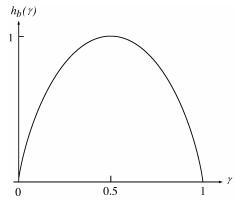


Fig. 2.1.  $h_b(q)$  versus q in the base 2.

**Definition 2.14.** The joint entropy H(X, Y) of a pair of random variables X and Y is defined as

$$H(X,Y) = -\sum_{x,y} p(x,y) \log p(x,y) = -E \log p(X,Y).$$
(2.40)

For two random variables, we define in the following the *conditional entropy* of one random variable when the other random variable is given.

**Definition 2.15.** For random variables X and Y, the conditional entropy of Y given X is defined as

$$H(Y|X) = -\sum_{x,y} p(x,y) \log p(y|x) = -E \log p(Y|X).$$
(2.41)

From (2.41), we can write

$$H(Y|X) = \sum_{x} p(x) \left[ -\sum_{y} p(y|x) \log p(y|x) \right].$$
 (2.42)

The inner sum is the entropy of Y conditioning on a fixed  $x \in S_X$ . Thus we are motivated to express H(Y|X) as

$$H(Y|X) = \sum_{x} p(x)H(Y|X=x),$$
 (2.43)

where

$$H(Y|X = x) = -\sum_{y} p(y|x) \log p(y|x).$$
(2.44)

Observe that the right hand sides of (2.35) and (2.44) have exactly the same form. Similarly, for H(Y|X, Z), we write

2.2 Shannon's Information Measures 15

$$H(Y|X,Z) = \sum_{z} p(z)H(Y|X,Z=z),$$
(2.45)

where

$$H(Y|X, Z = z) = -\sum_{x,y} p(x, y|z) \log p(y|x, z).$$
(2.46)

## Proposition 2.16.

$$H(X,Y) = H(X) + H(Y|X)$$
 (2.47)

and

$$H(X,Y) = H(Y) + H(X|Y).$$
 (2.48)

Proof. Consider

$$H(X,Y) = -E\log p(X,Y) \tag{2.49}$$

$$= -E \log[p(X)p(Y|X)]$$
(2.50)

$$= -E\log p(X) - E\log p(Y|X)$$
(2.51)

$$= H(X) + H(Y|X).$$
 (2.52)

Note that (2.50) is justified because the summation of the expectation is over  $S_{XY}$ , and we have used the linearity of expectation<sup>2</sup> to obtain (2.51). This proves (2.47), and (2.48) follows by symmetry.  $\Box$ 

This proposition has the following interpretation. Consider revealing the outcome of a pair of random variables X and Y in two steps: first the outcome of X and then the outcome of Y. Then the proposition says that the total amount of uncertainty removed upon revealing both X and Y is equal to the sum of the uncertainty removed upon revealing X (uncertainty removed in the first step) and the uncertainty removed upon revealing Y once X is known (uncertainty removed in the second step).

**Definition 2.17.** For random variables X and Y, the mutual information between X and Y is defined as

$$I(X;Y) = \sum_{x,y} p(x,y) \log \frac{p(x,y)}{p(x)p(y)} = E \log \frac{p(X,Y)}{p(X)p(Y)}.$$
 (2.53)

**Remark** I(X;Y) is symmetrical in X and Y.

**Proposition 2.18.** The mutual information between a random variable X and itself is equal to the entropy of X, i.e., I(X; X) = H(X).

 $<sup>^{2}</sup>$  See Problem 6 at the end of the chapter.

*Proof.* This can be seen by considering

$$I(X;X) = E \log \frac{p(X)}{p(X)^2}$$
(2.54)

$$= -E\log p(X) \tag{2.55}$$

$$=H(X). (2.56)$$

The proposition is proved.  $\Box$ 

**Remark** The entropy of X is sometimes called the self-information of X.

#### Proposition 2.19.

$$I(X;Y) = H(X) - H(X|Y),$$
(2.57)

$$I(X;Y) = H(Y) - H(Y|X),$$
(2.58)

and

$$I(X;Y) = H(X) + H(Y) - H(X,Y),$$
(2.59)

provided that all the entropies and conditional entropies are finite (see Example 2.46 in Section 2.8).

The proof of this proposition is left as an exercise.

From (2.57), we can interpret I(X; Y) as the reduction in uncertainty about X when Y is given, or equivalently, the amount of information about X provided by Y. Since I(X; Y) is symmetrical in X and Y, from (2.58), we can as well interpret I(X; Y) as the amount of information about Y provided by X.

The relations between the (joint) entropies, conditional entropies, and mutual information for two random variables X and Y are given in Propositions 2.16 and 2.19. These relations can be summarized by the diagram in Figure 2.2 which is a variation of the Venn diagram<sup>3</sup>. One can check that all the relations between Shannon's information measures for X and Y which are shown in Figure 2.2 are consistent with the relations given in Propositions 2.16 and 2.19. This one-to-one correspondence between Shannon's information measures and set theory is not just a coincidence for two random variables. We will discuss this in depth when we introduce the *I*-Measure in Chapter 3.

Analogous to entropy, there is a conditional version of mutual information called conditional mutual information.

<sup>&</sup>lt;sup>3</sup> The rectangle representing the universal set in a usual Venn diagram is missing in Figure 2.2.

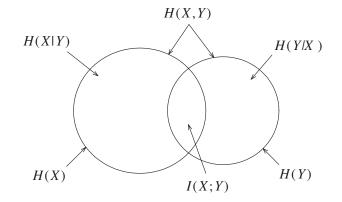


Fig. 2.2. Relationship between entropies and mutual information for two random variables.

**Definition 2.20.** For random variables X, Y and Z, the mutual information between X and Y conditioning on Z is defined as

$$I(X;Y|Z) = \sum_{x,y,z} p(x,y,z) \log \frac{p(x,y|z)}{p(x|z)p(y|z)} = E \log \frac{p(X,Y|Z)}{p(X|Z)p(Y|Z)}.$$
 (2.60)

**Remark** I(X;Y|Z) is symmetrical in X and Y.

Analogous to conditional entropy, we write

$$I(X;Y|Z) = \sum_{z} p(z)I(X;Y|Z=z),$$
(2.61)

where

$$I(X;Y|Z=z) = \sum_{x,y} p(x,y|z) \log \frac{p(x,y|z)}{p(x|z)p(y|z)}.$$
(2.62)

Similarly, when conditioning on two random variables, we write

$$I(X;Y|Z,T) = \sum_{t} p(t)I(X;Y|Z,T=t)$$
(2.63)

where

$$I(X;Y|Z,T=t) = \sum_{x,y,z} p(x,y,z|t) \log \frac{p(x,y|z,t)}{p(x|z,t)p(y|z,t)}.$$
(2.64)

Conditional mutual information satisfies the same set of relations given in Propositions 2.18 and 2.19 for mutual information except that all the terms are now conditioned on a random variable Z. We state these relations in the next two propositions. The proofs are omitted.

**Proposition 2.21.** The mutual information between a random variable X and itself conditioning on a random variable Z is equal to the conditional entropy of X given Z, i.e., I(X; X|Z) = H(X|Z).

# Proposition 2.22.

$$I(X;Y|Z) = H(X|Z) - H(X|Y,Z),$$
(2.65)

$$I(X;Y|Z) = H(Y|Z) - H(Y|X,Z),$$
(2.66)

and

$$I(X;Y|Z) = H(X|Z) + H(Y|Z) - H(X,Y|Z),$$
(2.67)

provided that all the conditional entropies are finite.

To conclude this section, we show that all Shannon's information measures are special cases of conditional mutual information. Let  $\Phi$  be a degenerate random variable, i.e.,  $\Phi$  takes a constant value with probability 1. Consider the mutual information I(X;Y|Z). When X = Y and  $Z = \Phi$ , I(X;Y|Z) becomes the entropy H(X). When X = Y, I(X;Y|Z) becomes the conditional entropy H(X|Z). When  $Z = \Phi$ , I(X;Y|Z) becomes the mutual information I(X;Y). Thus all Shannon's information measures are special cases of conditional mutual information.

# 2.3 Continuity of Shannon's Information Measures for Fixed Finite Alphabets

In this section, we prove that for fixed finite alphabets, all Shannon's information measures are continuous functionals of the joint distribution of the random variables involved. To formulate the notion of continuity, we first introduce the *variational distance*<sup>4</sup> as a distance measure between two probability distributions on a common alphabet.

**Definition 2.23.** Let p and q be two probability distributions on a common alphabet  $\mathcal{X}$ . The variational distance between p and q is defined as

$$V(p,q) = \sum_{x \in \mathcal{X}} |p(x) - q(x)|.$$
 (2.68)

According to (2.35), the entropy of a distribution p on an alphabet  $\mathcal{X}$  is defined as

$$H(p) = -\sum_{x \in \mathcal{S}_p} p(x) \log p(x)$$
(2.69)

where  $S_p$  denotes the support of p and  $S_p \subset \mathcal{X}$ . Let  $\mathcal{P}_{\mathcal{X}}$  be the set of all distributions on  $\mathcal{X}$ . In order for H(p) to be continuous with respect to convergence in variational distance at  $p \in \mathcal{P}_{\mathcal{X}}$ , for any  $\epsilon > 0$ , there exists  $\delta > 0$  such that

<sup>&</sup>lt;sup>4</sup> Also referred to as the  $L_1$  distance in mathematics.

2.3 Continuity of Shannon's Information Measures for Fixed Finite Alphabets 19

$$|H(p) - H(q)| < \epsilon \tag{2.70}$$

for all  $q \in \mathcal{P}_{\mathcal{X}}$  satisfying

$$V(p,q) < \delta, \tag{2.71}$$

or equivalently, for all  $p \in \mathcal{P}_{\mathcal{X}}$ ,

$$\lim_{p' \to p} H(p') = H\left(\lim_{p' \to p} p'\right) = H(p), \qquad (2.72)$$

where the convergence  $p' \rightarrow p$  is in variational distance.

Define a function  $l: [0, \infty) \to \Re$  by

$$l(a) = \begin{cases} a \log a \text{ if } a > 0\\ 0 \quad \text{if } a = 0. \end{cases}$$
(2.73)

Obviously, l(a) is continuous at all a > 0 and is continuous at a = 0 because  $a \log a \to 0$  as  $a \to 0$ . Thus l(a) is a continuous function of a. Then (2.69) can be rewritten as

$$H(p) = -\sum_{x \in \mathcal{X}} l(p(x)), \qquad (2.74)$$

where the summation above is over all x in  $\mathcal{X}$  instead of  $\mathcal{S}_p$ . Upon defining a function  $l_x : \mathcal{P}_{\mathcal{X}} \to \Re$  for all  $x \in \mathcal{X}$  by

$$l_x(p) = l(p(x)),$$
 (2.75)

(2.74) becomes

$$H(p) = -\sum_{x \in \mathcal{X}} l_x(p).$$
(2.76)

Evidently,  $l_x(p)$  is continuous in p (with respect to convergence in variational distance). Since the summation in (2.76) involves a finite number of terms, we conclude that H(p) is a continuous functional of p.

We now proceed to prove the continuity of conditional mutual information which covers all cases of Shannon's information measures. Consider I(X;Y|Z)and let  $p_{XYZ}$  be the joint distribution of X, Y, and Z, where the alphabets  $\mathcal{X}, \mathcal{Y}$ , and  $\mathcal{Z}$  are assumed to be finite. From (2.47) and (2.67), we obtain

$$I(X;Y|Z) = H(X,Z) + H(Y,Z) - H(X,Y,Z) - H(Z).$$
(2.77)

Note that each term on the right hand side above is the unconditional entropy of the corresponding marginal distribution. Then (2.77) can be rewritten as

$$I_{X;Y|Z}(p_{XYZ}) = H(p_{XZ}) + H(p_{YZ}) - H(p_{XYZ}) - H(p_Z), \qquad (2.78)$$

where we have used  $I_{X;Y|Z}(p_{XYZ})$  to denote I(X;Y|Z). It follows that

$$\lim_{\substack{p'_{XYZ} \to p_{XYZ} \\ p'_{XYZ} \to p_{XYZ}}} I_{X;Y|Z}(p'_{XYZ}) = \lim_{\substack{p'_{XYZ} \to p_{XYZ} \\ m_{XYZ} \to p_{XYZ}}} [H(p'_{XZ}) + H(p'_{YZ}) - H(p'_{XYZ}) - H(p'_{Z})]$$
(2.79)  
$$= \lim_{\substack{p'_{XYZ} \to p_{XYZ} \\ m_{XYZ} \to p_{XYZ}}} H(p'_{XZ}) + \lim_{\substack{p'_{XYZ} \to p_{XYZ} \\ m_{XYZ} \to p_{XYZ}}} H(p'_{XYZ}) - \lim_{\substack{p'_{XYZ} \to p_{XYZ} \\ m_{XYZ} \to p_{XYZ}}} H(p'_{XYZ}) - \lim_{\substack{p'_{XYZ} \to p_{XYZ} \\ m_{XYZ} \to p_{XYZ}}} H(p'_{XYZ}).$$
(2.80)

It can readily be proved, for example, that

$$\lim_{p'_{XYZ} \to p_{XYZ}} p'_{XZ} = p_{XZ}, \tag{2.81}$$

so that

$$\lim_{p'_{XYZ} \to p_{XYZ}} H(p'_{XZ}) = H\left(\lim_{p'_{XYZ} \to p_{XYZ}} p'_{XZ}\right) = H(p_{XZ})$$
(2.82)

by the continuity of  $H(\cdot)$  when the alphabets involved are fixed and finite. The details are left as an exercise. Hence, we conclude that

$$\lim_{p'_{XYZ} \to p_{XYZ}} I_{X;Y|Z}(p'_{XYZ}) = H(p_{XZ}) + H(p_{YZ}) - H(p_{XYZ}) - H(p_Z)$$
(2.83)

$$=I_{X;Y|Z}(p_{XYZ}),$$
 (2.84)

i.e.,  $I_{X;Y|Z}(p_{XYZ})$  is a continuous functional of  $p_{XYZ}$ .

We have completed the proof of the continuity of all Shannon's information measures with respect to convergence in variational distance for fixed finite alphabets. However, this result is rather restrictive and need to be applied with caution. It is because fixed finite alphabets are assumed for the random variables involved, and whether continuity holds depends critically on the distance measure used. In fact, Shannon's information measures are everywhere discontinuous with respect to convergence in a number of commonly used distance measures if the alphabets are not fixed. We refer the readers to Problems 27 to 29 for a discussion.

# 2.4 Chain Rules

In this section, we present a collection of information identities known as the chain rules which are often used in information theory.

Proposition 2.24 (Chain Rule for Entropy).

$$H(X_1, X_2, \cdots, X_n) = \sum_{i=1}^n H(X_i | X_1, \cdots, X_{i-1}).$$
 (2.85)

*Proof.* The chain rule for n = 2 has been proved in Proposition 2.16. We prove the chain rule by induction on n. Assume (2.85) is true for n = m, where  $m \ge 2$ . Then

$$H(X_1, \dots, X_m, X_{m+1}) = H(X_1, \dots, X_m) + H(X_{m+1}|X_1, \dots, X_m)$$
(2.86)

$$=\sum_{i=1}^{m} H(X_i|X_1,\cdots,X_{i-1}) + H(X_{m+1}|X_1,\cdots,X_m)$$
(2.87)

$$=\sum_{i=1}^{m+1} H(X_i|X_1,\cdots,X_{i-1}),$$
(2.88)

where in (2.86) we have used (2.47) by letting  $X = (X_1, \dots, X_m)$  and  $Y = X_{m+1}$ , and in (2.87) we have used (2.85) for n = m. This proves the chain rule for entropy.  $\Box$ 

The chain rule for entropy has the following conditional version.

# Proposition 2.25 (Chain Rule for Conditional Entropy).

$$H(X_1, X_2, \cdots, X_n | Y) = \sum_{i=1}^n H(X_i | X_1, \cdots, X_{i-1}, Y).$$
(2.89)

*Proof.* This can be proved by considering

$$H(X_1, X_2, \cdots, X_n | Y) = H(X_1, X_2, \cdots, X_n, Y) - H(Y)$$
(2.90)

$$= H((X_1, Y), X_2, \cdots, X_n) - H(Y)$$
(2.91)

$$= H(X_1, Y) + \sum_{i=2} H(X_i | X_1, \cdots, X_{i-1}, Y) - H(Y)$$
(2.92)

$$= H(X_1|Y) + \sum_{i=2}^{n} H(X_i|X_1, \cdots, X_{i-1}, Y)$$
(2.93)

$$=\sum_{i=1}^{n} H(X_i|X_1,\cdots,X_{i-1},Y),$$
(2.94)

where (2.90) and (2.93) follow from Proposition 2.16, while (2.92) follows from Proposition 2.24.  $\Box$ 

# Proposition 2.26 (Chain Rule for Mutual Information).

$$I(X_1, X_2, \cdots, X_n; Y) = \sum_{i=1}^n I(X_i; Y | X_1, \cdots, X_{i-1}).$$
(2.95)

*Proof.* Consider

$$I(X_1, X_2, \cdots, X_n; Y) = H(X_1, X_2, \cdots, X_n | Y)$$
(2.96)

$$=\sum_{i=1}^{n} [H(X_i|X_1,\cdots,X_{i-1}) - H(X_i|X_1,\cdots,X_{i-1},Y)]$$
(2.97)

$$=\sum_{i=1}^{n} I(X_i; Y|X_1, \cdots, X_{i-1}),$$
(2.98)

where in (2.97), we have invoked both Propositions 2.24 and 2.25. The chain rule for mutual information is proved.  $\Box$ 

**Proposition 2.27 (Chain Rule for Conditional Mutual Information).** For random variables  $X_1, X_2, \dots, X_n$ , Y, and Z,

$$I(X_1, X_2, \cdots, X_n; Y|Z) = \sum_{i=1}^n I(X_i; Y|X_1, \cdots, X_{i-1}, Z).$$
(2.99)

*Proof.* This is the conditional version of the chain rule for mutual information. The proof is similar to that for Proposition 2.25. The details are omitted.  $\Box$ 

# 2.5 Informational Divergence

Let p and q be two probability distributions on a common alphabet  $\mathcal{X}$ . We very often want to measure how much p is different from q, and vice versa. In order to be useful, this measure must satisfy the requirements that it is always nonnegative and it takes the zero value if and only if p = q. We denote the support of p and q by  $\mathcal{S}_p$  and  $\mathcal{S}_q$ , respectively. The *informational divergence* defined below serves this purpose.

**Definition 2.28.** The informational divergence between two probability distributions p and q on a common alphabet  $\mathcal{X}$  is defined as

$$D(p||q) = \sum_{x} p(x) \log \frac{p(x)}{q(x)} = E_p \log \frac{p(X)}{q(X)},$$
(2.100)

where  $E_p$  denotes expectation with respect to p.

In the above definition, in addition to the convention that the summation is taken over  $S_p$ , we further adopt the convention  $c \log \frac{c}{0} = \infty$  for c > 0. With this convention, if  $D(p||q) < \infty$ , then p(x) = 0 whenever q(x) = 0, i.e.,  $S_p \subset S_q$ . In the literature, the informational divergence is also referred to as *relative* entropy or the Kullback-Leibler distance. We note that D(p||q) is not symmetrical in p and q, so it is not a true metric or "distance." Moreover,  $D(\cdot||\cdot)$  does not satisfy the triangular inequality (see Problem 13).

In the rest of the book, the informational divergence will be referred to as *divergence* for brevity. Before we prove that divergence is always nonnegative, we first establish the following simple but important inequality called the *fundamental inequality* in information theory.

Lemma 2.29 (Fundamental Inequality). For any a > 0,

$$\ln a \le a - 1 \tag{2.101}$$

with equality if and only if a = 1.

*Proof.* Let  $f(a) = \ln a - a + 1$ . Then f'(a) = 1/a - 1 and  $f''(a) = -1/a^2$ . Since f(1) = 0, f'(1) = 0, and f''(1) = -1 < 0, we see that f(a) attains its maximum value 0 when a = 1. This proves (2.101). It is also clear that equality holds in (2.101) if and only if a = 1. Figure 2.3 is an illustration of the fundamental inequality.  $\Box$ 

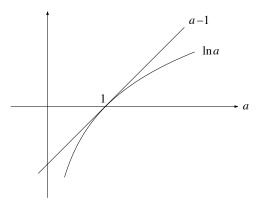


Fig. 2.3. The fundamental inequality  $\ln a \leq a - 1$ .

Corollary 2.30. For any a > 0,

$$\ln a \ge 1 - \frac{1}{a} \tag{2.102}$$

with equality if and only if a = 1.

*Proof.* This can be proved by replacing a by 1/a in (2.101).  $\Box$ 

We can see from Figure 2.3 that the fundamental inequality results from the convexity of the logarithmic function. In fact, many important results in

information theory are also direct or indirect consequences of the convexity of the logarithmic function!

**Theorem 2.31 (Divergence Inequality).** For any two probability distributions p and q on a common alphabet  $\mathcal{X}$ ,

$$D(p\|q) \ge 0 \tag{2.103}$$

with equality if and only if p = q.

*Proof.* If q(x) = 0 for some  $x \in S_p$ , then  $D(p||q) = \infty$  and the theorem is trivially true. Therefore, we assume that q(x) > 0 for all  $x \in S_p$ . Consider

$$D(p||q) = (\log e) \sum_{x \in S_p} p(x) \ln \frac{p(x)}{q(x)}$$
(2.104)

$$\geq (\log e) \sum_{x \in \mathcal{S}_p} p(x) \left( 1 - \frac{q(x)}{p(x)} \right)$$
(2.105)

$$= (\log e) \left[ \sum_{x \in \mathcal{S}_p} p(x) - \sum_{x \in \mathcal{S}_p} q(x) \right]$$
(2.106)

$$\geq 0, \tag{2.107}$$

where (2.105) results from an application of (2.102), and (2.107) follows from

$$\sum_{x \in \mathcal{S}_p} q(x) \le 1 = \sum_{x \in \mathcal{S}_p} p(x).$$
(2.108)

This proves (2.103).

For equality to hold in (2.103), equality must hold in (2.105) for all  $x \in S_p$ and also in (2.107). For the former, we see from Lemma 2.29 that this is the case if and only if

$$p(x) = q(x)$$
 for all  $x \in \mathcal{S}_p$ , (2.109)

which implies

$$\sum_{x \in \mathcal{S}_p} q(x) = \sum_{x \in \mathcal{S}_p} p(x) = 1, \qquad (2.110)$$

i.e., (2.107) holds with equality. Thus (2.109) is a necessary and sufficient condition for equality to hold in (2.103).

It is immediate that p = q implies (2.109), so it remains to prove the converse. Since  $\sum_{x} q(x) = 1$  and  $q(x) \ge 0$  for all x, p(x) = q(x) for all  $x \in S_p$  implies q(x) = 0 for all  $x \notin S_p$ , and therefore p = q. The theorem is proved.  $\Box$ 

We now prove a very useful consequence of the divergence inequality called the *log-sum inequality*.

**Theorem 2.32 (Log-Sum Inequality).** For positive numbers  $a_1, a_2, \cdots$  and nonnegative numbers  $b_1, b_2, \cdots$  such that  $\sum_i a_i < \infty$  and  $0 < \sum_i b_i < \infty$ ,

$$\sum_{i} a_i \log \frac{a_i}{b_i} \ge \left(\sum_{i} a_i\right) \log \frac{\sum_i a_i}{\sum_i b_i}$$
(2.111)

with the convention that  $\log \frac{a_i}{0} = \infty$ . Moreover, equality holds if and only if  $\frac{a_i}{b_i} = \text{constant for all } i$ .

The log-sum inequality can easily be understood by writing it out for the case when there are two terms in each of the summations:

$$a_1 \log \frac{a_1}{b_1} + a_2 \log \frac{a_2}{b_2} \ge (a_1 + a_2) \log \frac{a_1 + a_2}{b_1 + b_2}.$$
 (2.112)

Proof of Theorem 2.32. Let  $a'_i = a_i / \sum_j a_j$  and  $b'_i = b_i / \sum_j b_j$ . Then  $\{a'_i\}$  and  $\{b'_i\}$  are probability distributions. Using the divergence inequality, we have

$$0 \le \sum_{i} a'_i \log \frac{a'_i}{b'_i} \tag{2.113}$$

$$=\sum_{i}\frac{a_{i}}{\sum_{j}a_{j}}\log\frac{a_{i}/\sum_{j}a_{j}}{b_{i}/\sum_{j}b_{j}}$$
(2.114)

$$= \frac{1}{\sum_{j} a_{j}} \left[ \sum_{i} a_{i} \log \frac{a_{i}}{b_{i}} - \left( \sum_{i} a_{i} \right) \log \frac{\sum_{j} a_{j}}{\sum_{j} b_{j}} \right], \qquad (2.115)$$

which implies (2.111). Equality holds if and only if  $a'_i = b'_i$  for all i, or  $\frac{a_i}{b_i} = constant$  for all i. The theorem is proved.  $\Box$ 

One can also prove the divergence inequality by using the log-sum inequality (see Problem 19), so the two inequalities are in fact equivalent. The log-sum inequality also finds application in proving the next theorem which gives a lower bound on the divergence between two probability distributions on a common alphabet in terms of the *variational distance* between them. We will see further applications of the log-sum inequality when we discuss the convergence of some iterative algorithms in Chapter 9.

# Theorem 2.33 (Pinsker's Inequality).

$$D(p||q) \ge \frac{1}{2\ln 2} V^2(p,q).$$
(2.116)

Both divergence and the variational distance can be used as measures of the difference between two probability distributions defined on the same alphabet. Pinsker's inequality has the important implication that for two probability distributions p and q defined on the same alphabet, if D(p||q) is small,

then so is V(p,q). Furthermore, for a sequence of probability distributions  $q_k$ , as  $k \to \infty$ , if  $D(p||q_k) \to 0$ , then  $V(p,q_k) \to 0$ . In other words, convergence in divergence<sup>5</sup> is a stronger notion of convergence than convergence in variational distance.

The proof of Pinsker's inequality as well as its consequence discussed above is left as an exercise (see Problems 22 and 23).

### 2.6 The Basic Inequalities

In this section, we prove that all Shannon's information measures, namely entropy, conditional entropy, mutual information, and conditional mutual information are always nonnegative. By this, we mean that these quantities are nonnegative for all joint distributions for the random variables involved.

**Theorem 2.34.** For random variables X, Y, and Z,

$$I(X;Y|Z) \ge 0,$$
 (2.117)

with equality if and only if X and Y are independent when conditioning on Z.

Proof. Observe that

$$I(X;Y|Z) = \sum_{x,y,z} p(x,y,z) \log \frac{p(x,y|z)}{p(x|z)p(y|z)}$$
(2.118)

$$= \sum_{z} p(z) \sum_{x,y} p(x,y|z) \log \frac{p(x,y|z)}{p(x|z)p(y|z)}$$
(2.119)

$$=\sum_{z} p(z) D(p_{XY|z} || p_{X|z} p_{Y|z}), \qquad (2.120)$$

where we have used  $p_{XY|z}$  to denote  $\{p(x, y|z), (x, y) \in \mathcal{X} \times \mathcal{Y}\}$ , etc. Since for a fixed z, both  $p_{XY|z}$  and  $p_{X|z}p_{Y|z}$  are joint probability distributions on  $\mathcal{X} \times \mathcal{Y}$ , we have

$$D(p_{XY|z} \| p_{X|z} p_{Y|z}) \ge 0.$$
(2.121)

Therefore, we conclude that  $I(X; Y|Z) \ge 0$ . Finally, we see from Theorem 2.31 that I(X; Y|Z) = 0 if and only if for all  $z \in S_z$ ,

$$p(x, y|z) = p(x|z)p(y|z),$$
 (2.122)

<sup>&</sup>lt;sup>5</sup> In Harremoës [141], a sequence of probability distributions  $q_k$  converges to a probability distribution p if  $D(q_k || p) \to 0$  as  $k \to \infty$ , which is different from our discussion here. Nevertheless, as  $k \to 0$ , it holds that if  $D(q_k || p) \to 0$ , then  $V(p, q_k) \to 0$ .

$$p(x, y, z) = p(x, z)p(y|z)$$
 (2.123)

for all x and y. Therefore, X and Y are independent conditioning on Z. The proof is accomplished.  $\Box$ 

As we have seen in Section 2.2 that all Shannon's information measures are special cases of conditional mutual information, we already have proved that all Shannon's information measures are always nonnegative. The nonnegativity of all Shannon's information measures are called the *basic inequalities*.

For entropy and conditional entropy, we offer the following more direct proof for their nonnegativity. Consider the entropy H(X) of a random variable X. For all  $x \in S_X$ , since  $0 < p(x) \le 1$ ,  $\log p(x) \le 0$ . It then follows from the definition in (2.35) that  $H(X) \ge 0$ . For the conditional entropy H(Y|X) of random variable Y given random variable X, since  $H(Y|X = x) \ge 0$  for each  $x \in S_X$ , we see from (2.43) that  $H(Y|X) \ge 0$ .

**Proposition 2.35.** H(X) = 0 if and only if X is deterministic.

*Proof.* If X is deterministic, i.e., there exists  $x^* \in \mathcal{X}$  such that  $p(x^*) = 1$ and p(x) = 0 for all  $x \neq x^*$ , then  $H(X) = -p(x^*)\log p(x^*) = 0$ . On the other hand, if X is not deterministic, i.e., there exists  $x^* \in \mathcal{X}$  such that  $0 < p(x^*) < 1$ , then  $H(X) \ge -p(x^*)\log p(x^*) > 0$ . Therefore, we conclude that H(X) = 0 if and only if X is deterministic.  $\Box$ 

**Proposition 2.36.** H(Y|X) = 0 if and only if Y is a function of X.

*Proof.* From (2.43), we see that H(Y|X) = 0 if and only if H(Y|X = x) = 0 for each  $x \in S_X$ . Then from the last proposition, this happens if and only if Y is deterministic for each given x. In other words, Y is a function of X.  $\Box$ 

**Proposition 2.37.** I(X;Y) = 0 if and only if X and Y are independent.

*Proof.* This is a special case of Theorem 2.34 with Z being a degenerate random variable.  $\Box$ 

One can regard (conditional) mutual information as a measure of (conditional) dependency between two random variables. When the (conditional) mutual information is exactly equal to 0, the two random variables are (conditionally) independent.

We refer to inequalities involving Shannon's information measures only (possibly with constant terms) as *information inequalities*. The basic inequalities are important examples of information inequalities. Likewise, we refer to identities involving Shannon's information measures only as *information identities*. From the information identities (2.47), (2.57), and (2.65), we see that all Shannon's information measures can be expressed as linear combinations of entropies provided that the latter are all finite. Specifically,

$$H(Y|X) = H(X,Y) - H(X),$$
(2.124)

$$I(X;Y) = H(X) + H(Y) - H(X,Y), \qquad (2.125)$$

and

$$I(X;Y|Z) = H(X,Z) + H(Y,Z) - H(X,Y,Z) - H(Z).$$
(2.126)

Therefore, an information inequality is an inequality which involves only entropies.

As we will see later in the book, information inequalities form the most important set of tools for proving converse coding theorems in information theory. Except for a few so-called non-Shannon-type inequalities, all known information inequalities are implied by the basic inequalities. Information inequalities will be studied systematically in Chapters 13, 14, and 15. In the next section, we will prove some consequences of the basic inequalities which are often used in information theory.

# 2.7 Some Useful Information Inequalities

In this section, we prove some useful consequences of the basic inequalities introduced in the last section. Note that the conditional versions of these inequalities can be proved by techniques similar to those used in the proof of Proposition 2.25.

### Theorem 2.38 (Conditioning Does Not Increase Entropy).

$$H(Y|X) \le H(Y) \tag{2.127}$$

with equality if and only if X and Y are independent.

*Proof.* This can be proved by considering

$$H(Y|X) = H(Y) - I(X;Y) \le H(Y), \tag{2.128}$$

where the inequality follows because I(X;Y) is always nonnegative. The inequality is tight if and only if I(X;Y) = 0, which is equivalent by Proposition 2.37 to X and Y being independent.  $\Box$ 

Similarly, it can be shown that  $H(Y|X,Z) \leq H(Y|Z)$ , which is the conditional version of the above proposition. These results have the following interpretation. Suppose Y is a random variable we are interested in, and X and Z are side-information about Y. Then our uncertainty about Y cannot be increased on the average upon receiving side-information X. Once we know X, our uncertainty about Y again cannot be increased on the average upon further receiving side-information Z.

**Remark** Unlike entropy, the mutual information between two random variables can be increased by conditioning on a third random variable. We refer the reader to Section 3.4 for a discussion.

Theorem 2.39 (Independence Bound for Entropy).

$$H(X_1, X_2, \cdots, X_n) \le \sum_{i=1}^n H(X_i)$$
 (2.129)

with equality if and only if  $X_i$ ,  $i = 1, 2, \dots, n$ , are mutually independent.

*Proof.* By the chain rule for entropy,

$$H(X_1, X_2, \cdots, X_n) = \sum_{i=1}^n H(X_i | X_1, \cdots, X_{i-1})$$
(2.130)

$$\leq \sum_{i=1}^{n} H(X_i),$$
 (2.131)

where the inequality follows because we have proved in the last theorem that conditioning does not increase entropy. The inequality is tight if and only if it is tight for each i, i.e.,

$$H(X_i|X_1, \cdots, X_{i-1}) = H(X_i)$$
(2.132)

for  $1 \leq i \leq n$ . From the last theorem, this is equivalent to  $X_i$  being independent of  $X_1, X_2, \dots, X_{i-1}$  for each *i*. Then

$$p(x_1, x_2, \cdots, x_n) = p(x_1, x_2, \cdots, x_{n-1})p(x_n)$$
(2.133)

$$= p(p(x_1, x_2, \cdots, x_{n-2})p(x_{n-1})p(x_n)$$
(2.134)

$$\vdots = p(x_1)p(x_2)\cdots p(x_n)$$
 (2.135)

for all  $x_1, x_2, \dots, x_n$ , i.e.,  $X_1, X_2, \dots, X_n$  are mutually independent.

Alternatively, we can prove the theorem by considering

$$\sum_{i=1}^{n} H(X_i) - H(X_1, X_2, \cdots, X_n)$$
  
=  $-\sum_{i=1}^{n} E \log p(X_i) + E \log p(X_1, X_2, \cdots, X_n)$  (2.136)

$$= -E \log[p(X_1)p(X_2)\cdots p(X_n)] + E \log p(X_1, X_2, \cdots, X_n)$$
(2.137)

$$= E \log \frac{p(X_1, X_2, \cdots, X_n)}{p(X_1)p(X_2)\cdots p(X_n)}$$
(2.138)

$$= D(p_{X_1 X_2 \cdots X_n} \| p_{X_1} p_{X_2} \cdots p_{X_n})$$
(2.139)

$$\geq 0, \tag{2.140}$$

where equality holds if and only if

$$p(x_1, x_2, \cdots, x_n) = p(x_1)p(x_2)\cdots p(x_n)$$
 (2.141)

for all  $x_1, x_2, \dots, x_n$ , i.e.,  $X_1, X_2, \dots, X_n$  are mutually independent.  $\Box$ 

### Theorem 2.40.

$$I(X;Y,Z) \ge I(X;Y),$$
 (2.142)

with equality if and only if  $X \to Y \to Z$  forms a Markov chain.

*Proof.* By the chain rule for mutual information, we have

$$I(X;Y,Z) = I(X;Y) + I(X;Z|Y) \ge I(X;Y).$$
(2.143)

The above inequality is tight if and only if I(X; Z|Y) = 0, or  $X \to Y \to Z$  forms a Markov chain. The theorem is proved.  $\Box$ 

**Lemma 2.41.** If  $X \to Y \to Z$  forms a Markov chain, then

$$I(X;Z) \le I(X;Y) \tag{2.144}$$

and

$$I(X;Z) \le I(Y;Z).$$
 (2.145)

Before proving this inequality we discuss its meaning. Suppose X is a random variable we are interested in, and Y is an observation of X. If we infer X via Y, our uncertainty about X on the average is H(X|Y). Now suppose we process Y (either deterministically or probabilistically) to obtain a random variable Z. If we infer X via Z, our uncertainty about X on the average is H(X|Z). Since  $X \to Y \to Z$  forms a Markov chain, from (2.144), we have

$$H(X|Z) = H(X) - I(X;Z)$$
(2.146)

$$\geq H(X) - I(X;Y) \tag{2.147}$$

$$=H(X|Y), (2.148)$$

i.e., further processing of Y can only increase our uncertainty about X on the average.

Proof of Lemma 2.41. Assume  $X \to Y \to Z$ , i.e.,  $X \perp Z | Y$ . By Theorem 2.34, we have

$$I(X;Z|Y) = 0. (2.149)$$

Then

$$I(X;Z) = I(X;Y,Z) - I(X;Y|Z)$$
(2.150)  
 $\leq I(X;Y,Z)$ 
(2.151)

$$\leq I(X;Y,Z) \tag{2.151}$$

$$= I(X;Y) + I(X;Z|Y)$$
(2.152)

$$= I(X;Y).$$
 (2.153)

In (2.150) and (2.152), we have used the chain rule for mutual information. The inequality in (2.151) follows because I(X;Y|Z) is always nonnegative, and (2.153) follows from (2.149). This proves (2.144).

Since  $X \to Y \to Z$  is equivalent to  $Z \to Y \to X$ , we also have proved (2.145). This completes the proof of the lemma.  $\Box$ 

From Lemma 2.41, we can prove the more general data processing theorem.

**Theorem 2.42 (Data Processing Theorem).** If  $U \to X \to Y \to V$  forms a Markov chain, then

$$I(U;V) \le I(X;Y).$$
 (2.154)

*Proof.* Assume  $U \to X \to Y \to V$ . Then by Proposition 2.10, we have  $U \to X \to Y$  and  $U \to Y \to V$ . From the first Markov chain and Lemma 2.41, we have

$$I(U;Y) \le I(X;Y).$$
 (2.155)

From the second Markov chain and Lemma 2.41, we have

$$I(U;V) \le I(U;Y).$$
 (2.156)

Combining (2.155) and (2.156), we obtain (2.154), proving the theorem.  $\Box$ 

### 2.8 Fano's Inequality

In the last section, we have proved a few information inequalities involving only Shannon's information measures. In this section, we first prove an upper bound on the entropy of a random variable in terms of the size of the alphabet. This inequality is then used in the proof of Fano's inequality, which is extremely useful in proving converse coding theorems in information theory.

**Theorem 2.43.** For any random variable X,

$$H(X) \le \log |\mathcal{X}|,\tag{2.157}$$

where  $|\mathcal{X}|$  denotes the size of the alphabet  $\mathcal{X}$ . This upper bound is tight if and only if X distributes uniformly on  $\mathcal{X}$ .

*Proof.* Let u be the uniform distribution on  $\mathcal{X}$ , i.e.,  $u(x) = |\mathcal{X}|^{-1}$  for all  $x \in \mathcal{X}$ . Then

$$\log |\mathcal{X}| - H(X) = -\sum_{x \in \mathcal{S}_X} p(x) \log |\mathcal{X}|^{-1} + \sum_{x \in \mathcal{S}_X} p(x) \log p(x)$$
(2.158)

$$= -\sum_{x \in \mathcal{S}_X} p(x) \log u(x) + \sum_{x \in \mathcal{S}_X} p(x) \log p(x)$$
(2.159)

$$=\sum_{x\in\mathcal{S}_X} p(x)\log\frac{p(x)}{u(x)}$$
(2.160)

$$= D(p||u) \tag{2.161}$$

$$\geq 0, \tag{2.162}$$

proving (2.157). This upper bound is tight if and if only D(p||u) = 0, which from Theorem 2.31 is equivalent to p(x) = u(x) for all  $x \in \mathcal{X}$ , completing the proof.  $\Box$ 

**Corollary 2.44.** *The entropy of a random variable may take any nonnegative real value.* 

*Proof.* Consider a random variable X and let  $|\mathcal{X}|$  be fixed. We see from the last theorem that  $H(X) = \log |\mathcal{X}|$  is achieved when X distributes uniformly on  $\mathcal{X}$ . On the other hand, H(X) = 0 is achieved when X is deterministic. For any value  $0 < a < \log |\mathcal{X}|$ , by the intermediate value theorem of continuous functions, there exists a distribution for X such that H(X) = a. Then we see that H(X) can take any positive value by letting  $|\mathcal{X}|$  be sufficiently large. This accomplishes the proof.  $\Box$ 

**Remark** Let  $|\mathcal{X}| = D$ , or the random variable X is a D-ary symbol. When the base of the logarithm is D, (2.157) becomes

$$H_D(X) \le 1.$$
 (2.163)

Recall that the unit of entropy is the *D*-it when the logarithm is in the base *D*. This inequality says that a *D*-ary symbol can carry at most 1 *D*-it of information. This maximum is achieved when *X* has a uniform distribution. We already have seen the binary case when we discuss the binary entropy function  $h_b(p)$  in Section 2.2.

We see from Theorem 2.43 that the entropy of a random variable is finite as long as it has a finite alphabet. However, if a random variable has a countable  $alphabet^6$ , its entropy may or may not be finite. This will be shown in the next two examples.

Example 2.45. Let X be a random variable such that

$$\Pr\{X = i\} = 2^{-i},\tag{2.164}$$

 $i = 1, 2, \cdots$ . Then

$$H_2(X) = \sum_{i=1}^{\infty} i2^{-i} = 2, \qquad (2.165)$$

which is finite.

<sup>&</sup>lt;sup>6</sup> An alphabet is countable means that it may be countably infinite.

For a random variable X with a countable alphabet and finite entropy, we show in Appendix 2.10 that the entropy of X can be approximated by the entropy of a truncation of the distribution of X.

*Example 2.46.* Let Y be a random variable which takes values in the subset of pairs of integers

$$\left\{ (i,j) : 1 \le i < \infty \text{ and } 1 \le j \le \frac{2^{2^i}}{2^i} \right\}$$
(2.166)

such that

$$\Pr\{Y = (i, j)\} = 2^{-2^{i}}$$
(2.167)

for all i and j. First, we check that

$$\sum_{i=1}^{\infty} \sum_{j=1}^{2^{2^i}/2^i} \Pr\{Y = (i,j)\} = \sum_{i=1}^{\infty} 2^{-2^i} \left(\frac{2^{2^i}}{2^i}\right) = 1.$$
(2.168)

Then

$$H_2(Y) = -\sum_{i=1}^{\infty} \sum_{j=1}^{2^{2^i}/2^i} 2^{-2^i} \log_2 2^{-2^i} = \sum_{i=1}^{\infty} 1,$$
 (2.169)

which does not converge.

Let X be a random variable and  $\hat{X}$  be an estimate of X which takes value in the same alphabet  $\mathcal{X}$ . Let the probability of error  $P_e$  be

$$P_e = \Pr\{X \neq \hat{X}\}.\tag{2.170}$$

If  $P_e = 0$ , i.e.,  $X = \hat{X}$  with probability 1, then  $H(X|\hat{X}) = 0$  by Proposition 2.36. Intuitively, if  $P_e$  is small, i.e.,  $X = \hat{X}$  with probability close to 1, then  $H(X|\hat{X})$  should be close to 0. Fano's inequality makes this intuition precise.

**Theorem 2.47 (Fano's Inequality).** Let X and  $\hat{X}$  be random variables taking values in the same alphabet  $\mathcal{X}$ . Then

$$H(X|X) \le h_b(P_e) + P_e \log(|\mathcal{X}| - 1),$$
 (2.171)

where  $h_b$  is the binary entropy function.

*Proof.* Define a random variable

$$Y = \begin{cases} 0 \text{ if } X = \hat{X} \\ 1 \text{ if } X \neq \hat{X}. \end{cases}$$
(2.172)

The random variable Y is an indicator of the error event  $\{X \neq \hat{X}\}$ , with  $\Pr\{Y=1\} = P_e$  and  $H(Y) = h_b(P_e)$ . Since Y is a function X and  $\hat{X}$ ,

 $H(X|\hat{X})$ 

$$H(Y|X, \hat{X}) = 0.$$
 (2.173)

Then

$$= H(X|\hat{X}) + H(Y|X,\hat{X})$$
(2.174)

$$=H(X,Y|\hat{X}) \tag{2.175}$$

$$= H(Y|\hat{X}) + H(X|\hat{X},Y)$$
(2.176)

$$\leq H(Y) + H(X|\hat{X}, Y) \tag{2.177}$$

$$= H(Y) + \sum_{\hat{x} \in \mathcal{X}} \left[ \Pr\{\hat{X} = \hat{x}, Y = 0\} H(X | \hat{X} = \hat{x}, Y = 0) + \Pr\{\hat{X} = \hat{x}, Y = 1\} H(X | \hat{X} = \hat{x}, Y = 1) \right].$$
(2.178)

In the above, (2.174) follows from (2.173), (2.177) follows because conditioning does not increase entropy, and (2.178) follows from an application of (2.43). Now X must take the value  $\hat{x}$  if  $\hat{X} = \hat{x}$  and Y = 0. In other words, X is conditionally deterministic given  $\hat{X} = \hat{x}$  and Y = 0. Therefore, by Proposition 2.35,

$$H(X|\hat{X} = \hat{x}, Y = 0) = 0.$$
(2.179)

If  $\hat{X} = \hat{x}$  and Y = 1, then X must take a value in the set  $\{x \in \mathcal{X} : x \neq \hat{x}\}$  which contains  $|\mathcal{X}| - 1$  elements. From the last theorem, we have

$$H(X|\dot{X} = \hat{x}, Y = 1) \le \log(|\mathcal{X}| - 1), \tag{2.180}$$

where this upper bound does not depend on  $\hat{x}$ . Hence,

.

$$H(X|\hat{X}) \leq h_b(P_e) + \left(\sum_{\hat{x}\in\mathcal{X}} \Pr\{\hat{X}=\hat{x}, Y=1\}\right) \log(|\mathcal{X}|-1)$$
(2.181)

$$= h_b(P_e) + \Pr\{Y = 1\} \log(|\mathcal{X}| - 1)$$
(2.182)

$$= h_b(P_e) + P_e \log(|\mathcal{X}| - 1), \qquad (2.183)$$

which completes the proof.  $\hfill\square$ 

Very often, we only need the following simplified version when we apply Fano's inequality. The proof is omitted.

# Corollary 2.48. $H(X|\hat{X}) < 1 + P_e \log |\mathcal{X}|.$

Fano's inequality has the following implication. If the alphabet  $\mathcal{X}$  is finite, as  $P_e \to 0$ , the upper bound in (2.171) tends to 0, which implies  $H(X|\hat{X})$  also tends to 0. However, this is not necessarily the case if  $\mathcal{X}$  is countable, which is shown in the next example.

*Example 2.49.* Let  $\hat{X}$  take the value 0 with probability 1. Let Z be an independent binary random variable taking values in  $\{0, 1\}$ . Define the random variable X by

$$X = \begin{cases} 0 & \text{if } Z = 0 \\ Y & \text{if } Z = 1, \end{cases}$$
(2.184)

where Y is the random variable in Example 2.46 whose entropy is infinity. Let

$$P_e = \Pr\{X \neq X\} = \Pr\{Z = 1\}.$$
 (2.185)

Then

$$H(X|\hat{X}) \tag{2.186}$$

$$=H(X) \tag{2.187}$$

$$\geq H(X|Z)$$

$$= \Pr\{Z = 0\} H(Y|Z = 0) + \Pr\{Z = 1\} H(Y|Z = 1)$$
(2.188)
(2.188)
(2.180)

$$= \Pr\{Z = 0\} H(X|Z = 0) + \Pr\{Z = 1\} H(X|Z = 1)$$
(2.189)  
=  $(1 - P_e) \cdot 0 + P_e \cdot H(Y)$ (2.190)

$$=\infty$$
(2.191)

for any  $P_e > 0$ . Therefore,  $H(X|\hat{X})$  does not tend to 0 as  $P_e \to 0$ .

# 2.9 Maximum Entropy Distributions

In Theorem 2.43, we have proved that for any random variable X,

$$H(X) \le \log |\mathcal{X}|,\tag{2.192}$$

with equality when X is distributed uniformly over its support. In this section, we revisit this result in the context that X is a real random variable.

To simplify our discussion, all the logarithms are in the base e. Consider the following problem: Maximize H(p) over all probability distribution p defined on a countable subset S (possibly infinite) of the set of real numbers, subject to

$$\sum_{x \in \mathcal{S}_p} p(x)r_i(x) = a_i \quad \text{for } 1 \le i \le m.$$
(2.193)

The following theorem provides a solution to this problem.

### Theorem 2.50. Let

$$p^{*}(x) = e^{-\lambda_{0} - \sum_{i=1}^{m} \lambda_{i} r_{i}(x)}$$
(2.194)

for all  $x \in S$ , where  $\lambda_0, \lambda_1, \dots, \lambda_m$  are chosen such that the constraints in (2.193) are satisfied. Then  $p^*$  maximizes H(p) over all probability distribution p on S, subject to the constraints in (2.193).

*Proof.* For any p satisfying the constraints in (2.193), consider

$$H(p^*) - H(p) = -\sum_{x \in S} p^*(x) \ln p^*(x) + \sum_{x \in S_p} p(x) \ln p(x)$$
(2.195)

$$= -\sum_{x \in \mathcal{S}} p^*(x) \left( -\lambda_0 - \sum_i \lambda_i r_i(x) \right) + \sum_{x \in \mathcal{S}_p} p(x) \ln p(x)$$
(2.196)

$$= \lambda_0 \left( \sum_{x \in \mathcal{S}} p^*(x) \right) + \sum_i \lambda_i \left( \sum_{x \in \mathcal{S}} p^*(x) r_i(x) \right) + \sum_{x \in \mathcal{S}_p} p(x) \ln p(x) \quad (2.197)$$

$$= \lambda_0 \cdot 1 + \sum_i \lambda_i a_i + \sum_{x \in \mathcal{S}_p} p(x) \ln p(x)$$
(2.198)

$$=\lambda_0 \left(\sum_{x\in\mathcal{S}_p} p(x)\right) + \sum_i \lambda_i \left(\sum_{x\in\mathcal{S}_p} p(x)r_i(x)\right) + \sum_{x\in\mathcal{S}_p} p(x)\ln p(x) \quad (2.199)$$

$$= -\sum_{x \in \mathcal{S}_p} p(x) \left( -\lambda_0 - \sum_i \lambda_i r_i(x) \right) + \sum_{x \in \mathcal{S}_p} p(x) \ln p(x)$$
(2.200)

$$= -\sum_{x \in S_p} p(x) \ln p^*(x) + \sum_{x \in S_p} p(x) \ln p(x)$$
(2.201)

$$= \sum_{x \in S_p} p(x) \ln \frac{p(x)}{p^*(x)}$$
(2.202)

$$= D(p||p^*)$$
(2.203)

$$\geq 0. \tag{2.204}$$

In the above, (2.200) is obtained from (2.196) by replacing  $p^*(x)$  by p(x) and  $x \in S$  by  $x \in S_p$  in the first summation, while the intermediate steps (2.197) to (2.199) are justified by noting that both  $p^*$  and p satisfy the constraints in (2.193). The last step is an application of the divergence inequality (Theorem 2.31). The proof is accomplished.  $\Box$ 

**Remark** For all  $x \in S$ ,  $p^*(x) > 0$ , so that  $S_p = S$ .

The following corollary of Theorem 2.50 is rather subtle.

**Corollary 2.51.** Let  $p^*$  be a probability distribution defined on S with

$$p^{*}(x) = e^{-\lambda_{0} - \sum_{i=1}^{m} \lambda_{i} r_{i}(x)}$$
(2.205)

for all  $x \in S$ . Then  $p^*$  maximizes H(p) over all probability distribution p defined on S, subject to the constraints

$$\sum_{x \in \mathcal{S}_p} p(x)r_i(x) = \sum_{x \in \mathcal{S}} p^*(x)r_i(x) \quad \text{for } 1 \le i \le m.$$
(2.206)

*Example 2.52.* Let S be finite and let the set of constraints in (2.193) be empty. Then

$$p^*(x) = e^{-\lambda_0},$$
 (2.207)

a constant that does not depend on x. Therefore,  $p^*$  is simply the uniform distribution over S, i.e.,  $p^*(x) = |S|^{-1}$  for all  $x \in S$ .

*Example 2.53.* Let  $S = \{0, 1, 2, \dots\}$ , and let the set of constraints in (2.193) be

$$\sum_{x} p(x)x = a, \qquad (2.208)$$

where  $a \ge 0$ , i.e., the mean of the distribution p is fixed at some nonnegative value a. We now determine  $p^*$  using the prescription in Theorem 2.50. Let

$$q_i = e^{-\lambda_i} \tag{2.209}$$

for i = 0, 1. Then

$$p^*(x) = q_0 q_1^x. (2.210)$$

Evidently,  $p^*$  is a geometric distribution, so that

$$q_0 = 1 - q_1. \tag{2.211}$$

Finally, we invoke the constraint (2.193) on p to obtain  $q_1 = (a+1)^{-1}$ . The details are omitted.

### 2.10 Entropy Rate of Stationary Source

In the previous sections, we have discussed various properties of the entropy of a finite collection of random variables. In this section, we discuss the entropy rate *entropy rate* of a discrete-time information source.

A discrete-time information source  $\{X_k, k \ge 1\}$  is an infinite collection of random variables indexed by the set of positive integers. Since the index set is ordered, it is natural to regard the indices as time indices. We will refer to the random variables  $X_k$  as *letters*.

We assume that  $H(X_k) < \infty$  for all k. Then for any finite subset A of the index set  $\{k : k \ge 1\}$ , we have

$$H(X_k, k \in A) \le \sum_{k \in A} H(X_k) < \infty.$$
(2.212)

However, it is not meaningful to discuss  $H(X_k, k \ge 1)$  because the joint entropy of an infinite collection of letters is infinite except in special cases. On the other hand, since the indices are ordered, we can naturally define the *entropy rate* of an information source, which gives the average entropy per letter of the source.

**Definition 2.54.** The entropy rate of an information source  $\{X_k\}$  is defined by

$$H_X = \lim_{n \to \infty} \frac{1}{n} H(X_1, X_2, \cdots, X_n)$$
(2.213)

when the limit exists.

We show in the next two examples that the entropy rate of a source may or may not exist.

*Example 2.55.* Let  $\{X_k\}$  be an i.i.d. source with generic random variable X. Then

$$\lim_{n \to \infty} \frac{1}{n} H(X_1, X_2, \cdots, X_n) = \lim_{n \to \infty} \frac{n H(X)}{n}$$
(2.214)  
=  $\lim_{n \to \infty} H(X)$ (2.215)

$$=\lim_{n \to \infty} H(X) \tag{2.215}$$

$$=H(X), (2.216)$$

i.e., the entropy rate of an i.i.d. source is the entropy of any of its single letters.

*Example 2.56.* Let  $\{X_k\}$  be a source such that  $X_k$  are mutually independent and  $H(X_k) = k$  for  $k \ge 1$ . Then

$$\frac{1}{n}H(X_1, X_2, \cdots, X_n) = \frac{1}{n}\sum_{k=1}^n k$$
(2.217)

$$=\frac{1}{n}\frac{n(n+1)}{2}$$
(2.218)

$$=\frac{1}{2}(n+1),$$
 (2.219)

which does not converge as  $n \to \infty$  although  $H(X_k) < \infty$  for all k. Therefore, the entropy rate of  $\{X_k\}$  does not exist.

Toward characterizing the asymptotic behavior of  $\{X_k\}$ , it is natural to consider the limit

$$H'_{X} = \lim_{n \to \infty} H(X_{n} | X_{1}, X_{2}, \cdots, X_{n-1})$$
(2.220)

if it exists. The quantity  $H(X_n|X_1, X_2, \dots, X_{n-1})$  is interpreted as the conditional entropy of the next letter given that we know all the past history of the source, and  $H'_X$  is the limit of this quantity after the source has been run for an indefinite amount of time.

**Definition 2.57.** An information source  $\{X_k\}$  is stationary if

$$X_1, X_2, \cdots, X_m \tag{2.221}$$

and

$$X_{1+l}, X_{2+l}, \cdots, X_{m+l} \tag{2.222}$$

have the same joint distribution for any  $m, l \geq 1$ .

In the rest of the section, we will show that stationarity is a sufficient condition for the existence of the entropy rate of an information source.

# **Lemma 2.58.** Let $\{X_k\}$ be a stationary source. Then $H'_X$ exists.

*Proof.* Since  $H(X_n|X_1, X_2, \dots, X_{n-1})$  is lower bounded by zero for all n, it suffices to prove that  $H(X_n|X_1, X_2, \dots, X_{n-1})$  is non-increasing in n to conclude that the limit  $H'_X$  exists. Toward this end, for  $n \ge 2$ , consider

$$H(X_n|X_1, X_2, \cdots, X_{n-1}) \le H(X_n|X_2, X_3, \cdots, X_{n-1})$$
(2.223)

$$= H(X_{n-1}|X_1, X_2, \cdots, X_{n-2}), \qquad (2.224)$$

where the last step is justified by the stationarity of  $\{X_k\}$ . The lemma is proved.  $\Box$ 

**Lemma 2.59 (Cesáro Mean).** Let  $a_k$  and  $b_k$  be real numbers. If  $a_n \to a$  as  $n \to \infty$  and  $b_n = \frac{1}{n} \sum_{k=1}^n a_k$ , then  $b_n \to a$  as  $n \to \infty$ .

*Proof.* The idea of the lemma is the following. If  $a_n \to a$  as  $n \to \infty$ , then the average of the first n terms in  $\{a_k\}$ , namely  $b_n$ , also tends to a as  $n \to \infty$ .

The lemma is formally proved as follows. Since  $a_n \to a$  as  $n \to \infty$ , for every  $\epsilon > 0$ , there exists  $N(\epsilon)$  such that  $|a_n - a| < \epsilon$  for all  $n > N(\epsilon)$ . For  $n > N(\epsilon)$ , consider

$$|b_n - a| = \left| \frac{1}{n} \sum_{i=1}^n a_i - a \right|$$
(2.225)

$$= \left| \frac{1}{n} \sum_{i=1}^{n} (a_i - a) \right|$$
(2.226)

$$\leq \frac{1}{n} \sum_{i=1}^{n} |a_i - a| \tag{2.227}$$

$$= \frac{1}{n} \left( \sum_{i=1}^{N(\epsilon)} |a_i - a| + \sum_{i=N(\epsilon)+1}^n |a_i - a| \right)$$
(2.228)

$$<\frac{1}{n}\sum_{i=1}^{N(\epsilon)}|a_i-a|+\frac{(n-N(\epsilon))\epsilon}{n}$$
(2.229)

$$<\frac{1}{n}\sum_{i=1}^{N(\epsilon)}|a_i-a|+\epsilon.$$
(2.230)

The first term tends to 0 as  $n \to \infty$ . Therefore, for any  $\epsilon > 0$ , by taking n to be sufficiently large, we can make  $|b_n - a| < 2\epsilon$ . Hence  $b_n \to a$  as  $n \to \infty$ , proving the lemma.  $\Box$ 

We now prove that  $H'_X$  is an alternative definition/interpretation of the entropy rate of  $\{X_k\}$  when  $\{X_k\}$  is stationary.

**Theorem 2.60.** For a stationary source  $\{X_k\}$ , the entropy rate  $H_X$  exists, and it is equal to  $H'_X$ .

*Proof.* Since we have proved in Lemma 2.58 that  $H'_X$  always exists for a stationary source  $\{X_k\}$ , in order to prove the theorem, we only have to prove that  $H_X = H'_X$ . By the chain rule for entropy,

$$\frac{1}{n}H(X_1, X_2, \cdots, X_n) = \frac{1}{n}\sum_{k=1}^n H(X_k|X_1, X_2, \cdots, X_{k-1}).$$
 (2.231)

Since

$$\lim_{k \to \infty} H(X_k | X_1, X_2, \cdots, X_{k-1}) = H'_X$$
(2.232)

from (2.220), it follows from Lemma 2.59 that

$$H_X = \lim_{n \to \infty} \frac{1}{n} H(X_1, X_2, \cdots, X_n) = H'_X.$$
 (2.233)

The theorem is proved.  $\Box$ 

In this theorem, we have proved that the entropy rate of a random source  $\{X_k\}$  exists under the fairly general assumption that  $\{X_k\}$  is stationary. However, the entropy rate of a stationary source  $\{X_k\}$  may not carry any physical meaning unless  $\{X_k\}$  is also ergodic. This will be explained when we discuss the Shannon-McMillan-Breiman Theorem in Section 5.4.

# Appendix 2.A: Approximation of Random Variables with Countably Infinite Alphabets by Truncation

Let X be a random variable with a countable alphabet  $\mathcal{X}$  such that  $H(X) < \infty$ . Without loss of generality,  $\mathcal{X}$  is taken to be the set of positive integers. Define a random variable X(m) which takes values in

$$\mathcal{N}_m = \{1, 2, \cdots, m\}$$
(2.234)

such that

$$\Pr\{X(m) = k\} = \frac{\Pr\{X = k\}}{\Pr\{X \in \mathcal{N}_m\}}$$

$$(2.235)$$

for all  $k \in \mathcal{N}_m$ , i.e., the distribution of X(m) is the truncation of the distribution of X up to m.

It is intuitively correct that  $H(X(m)) \to H(X)$  as  $m \to \infty$ , which we formally prove in this appendix. For every  $m \ge 1$ , define the binary random variable

2.10 Entropy Rate of Stationary Source 41

$$B(m) = \begin{cases} 1 \text{ if } X \le m\\ 0 \text{ if } X > m. \end{cases}$$
(2.236)

Consider

$$H(X) = -\sum_{k=1}^{m} \Pr\{X = k\} \log \Pr\{X = k\}$$
$$-\sum_{k=m+1}^{\infty} \Pr\{X = k\} \log \Pr\{X = k\}.$$
 (2.237)

As  $m \to \infty$ ,

$$-\sum_{k=1}^{m} \Pr\{X=k\} \log \Pr\{X=k\} \to H(X).$$
 (2.238)

Since  $H(X) < \infty$ ,

$$-\sum_{k=m+1}^{\infty} \Pr\{X=k\} \log \Pr\{X=k\} \to 0$$
 (2.239)

as  $k \to \infty$ . Now consider

$$H(X) = H(X|B(m)) + I(X;B(m))$$
(2.240)  
=  $H(X|B(m) = 1)\Pr\{B(m) = 1\} + H(X|B(m) = 0)$   
 $\times \Pr\{B(m) = 0\} + I(X;B(m))$ (2.241)

$$= H(X(m)) \Pr\{B(m) = 1\} + H(X|B(m) = 0)$$
  
 
$$\times \Pr\{B(m) = 0\} + I(X; B(m)).$$
(2.242)

As  $m\to\infty,\ H(B(m))\to 0$  since  $\Pr\{B(m)=1\}\to 1.$  This implies  $I(X;B(m))\to 0$  because

$$I(X; B(m)) \le H(B(m)).$$
 (2.243)

In (2.242), we further consider

$$H(X|B(m) = 0)\Pr\{B(m) = 0\}$$
  
=  $-\sum_{k=m+1}^{\infty} \Pr\{X = k\} \log \frac{\Pr\{X = k\}}{\Pr\{B(m) = 0\}}$  (2.244)  
=  $-\sum_{k=m+1}^{\infty} \Pr\{X = k\} (\log \Pr\{X = k\})$   
 $-\log \Pr\{B(m) = 0\})$  (2.245)  
=  $-\sum_{k=m+1}^{\infty} (\Pr\{X = k\} \log \Pr\{X = k\})$ 

$$+\left(\sum_{k=m+1}^{\infty} \Pr\{X=k\}\right) \log \Pr\{B(m)=0\}$$
(2.246)  
$$= -\sum_{k=m+1}^{\infty} \Pr\{X=k\} \log \Pr\{X=k\}$$
  
$$+\Pr\{B(m)=0\} \log \Pr\{B(m)=0\}.$$
(2.247)

As  $m \to \infty$ , the summation above tends to 0 by (2.239). Since  $\Pr\{B(m) = 0\} \to 0$ ,  $\Pr\{B(m) = 0\} \log \Pr\{B(m) = 0\} \to 0$ . Therefore,

$$H(X|B(m) = 0)\Pr\{B(m) = 0\} \to 0,$$
 (2.248)

and we see from (2.242) that  $H(X(m)) \to H(X)$  as  $m \to \infty$ .

# Problems

1. Let X and Y be random variables with alphabets  $\mathcal{X} = \mathcal{Y} = \{1, 2, 3, 4, 5\}$ and joint distribution p(x, y) given by

$$\frac{1}{25} \begin{bmatrix} 1 & 1 & 1 & 1 & 1 \\ 2 & 1 & 2 & 0 & 0 \\ 2 & 0 & 1 & 1 & 1 \\ 0 & 3 & 0 & 2 & 0 \\ 0 & 0 & 1 & 1 & 3 \end{bmatrix}$$

Calculate H(X), H(Y), H(X|Y), H(Y|X), and I(X;Y). 2. Consider any probability distribution p(x, y, z) and let

$$q(x, y, z) = \begin{cases} p(x)p(y)p(z|x, y) \text{ if } p(x, y) > 0\\ 0 & \text{otherwise.} \end{cases}$$

- a) Show that q(x, y, z) is in general not a probability distribution.
- b) By ignoring the fact that q(x, y, z) may not be a probability distribution, application of the divergence inequality  $D(p||q) \ge 0$  would yield the inequality

$$H(X) + H(Y) + H(Z|X,Y) \ge H(X,Y,Z),$$

which indeed is valid for all jointly distributed random variables X, Y, and Z. Explain.

- 3. Prove Propositions 2.8, 2.9, 2.10, 2.19, 2.21, and 2.22.
- 4. Give an example which shows that pairwise independence does not imply mutual independence.
- 5. Verify that p(x, y, z) as defined in Definition 2.4 is a probability distribution. You should exclude all the zero probability masses from the summation carefully.

- 6. Linearity of expectation It is well-known that expectation is linear, i.e., E[f(X) + g(Y)] = Ef(X) + Eg(Y), where the summation in an expectation is taken over the corresponding alphabet. However, we adopt in information theory the convention that the summation in an expectation is taken over the corresponding support. Justify carefully the linearity of expectation under this convention.
- 7. The identity I(X;Y) = H(X) H(X|Y) is invalid if H(X|Y) (and hence H(X)) is equal to infinity. Give an example such that I(X;Y) has a finite value but both H(X) and H(Y|X) are equal to infinity.
- 8. Let  $p'_{XY}$  and  $p_{XY}$  be defined on  $\mathcal{X} \times \mathcal{Y}$ , where  $\mathcal{X}$  and  $\mathcal{Y}$  are fixed finite alphabets. Prove that

$$\lim_{p'_{XY} \to p_{XY}} p'_x = p_X,$$

where the limit is taken with respect to the variational distance.

- 9. Let  $C_{\alpha} = \sum_{n=2}^{\infty} \frac{1}{n(\log n)^{\alpha}}$ .
  - a) Prove that

$$C_{\alpha} \begin{cases} < \infty \text{ if } \alpha > 1 \\ = \infty \text{ if } 0 \le \alpha \le 1 \end{cases}$$

Then

$$p_{\alpha}(n) = [C_{\alpha}n(\log n)^{\alpha}]^{-1}, \quad n = 2, 3, \cdots$$

is a probability distribution for  $\alpha > 1$ .

b) Prove that

$$H(p_{\alpha}) \begin{cases} < \infty \text{ if } \alpha > 2 \\ = \infty \text{ if } 1 < \alpha \le 2 \end{cases}$$

10. Prove that H(p) is concave in p, i.e., for  $0 \le \lambda \le 1$  and  $\overline{\lambda} = 1 - \lambda$ ,

$$\lambda H(p_1) + \overline{\lambda} H(p_2) \le H(\lambda p_1 + \overline{\lambda} p_2).$$

- 11. Let  $(X, Y) \sim p(x, y) = p(x)p(y|x)$ .
  - a) Prove that for fixed p(x), I(X;Y) is a convex functional of p(y|x).
  - b) Prove that for fixed p(y|x), I(X;Y) is a concave functional of p(x).
- 12. Do I(X;Y) = 0 and I(X;Y|Z) = 0 imply each other? If so, give a proof. If not, give a counterexample.
- 13. Give an example for which  $D(\cdot \| \cdot)$  does not satisfy the triangular inequality.
- 14. Let X be a function of Y. Prove that  $H(X) \leq H(Y)$ . Interpret this result.
- 15. Prove that for any  $n \ge 2$ ,

$$H(X_1, X_2, \cdots, X_n) \ge \sum_{i \neq j} H(X_i | X_j).$$

Give the necessary and sufficient condition for equality to hold and construct a nontrivial example that satisfies this condition.

- 44 2 Information Measures
- 16. Prove that for any  $n \ge 2$ ,

$$H(X_1, X_2, \cdots, X_n) \ge \sum_{i=1}^n H(X_i | X_j, j \ne i).$$

17. Prove that

$$\frac{1}{2}[H(X_1, X_2) + H(X_2, X_3) + H(X_1, X_3)] \ge H(X_1, X_2, X_3).$$

Hint: Sum the identities

$$H(X_1, X_2, X_3) = H(X_j, j \neq i) + H(X_i | X_j, j \neq i)$$

for i = 1, 2, 3 and apply the result in Problem 16.

18. Let  $\mathcal{N}_n = \{1, 2, \dots, n\}$  and denote  $H(X_i, i \in \alpha)$  by  $H(X_\alpha)$  for any subset  $\alpha$  of  $\mathcal{N}_n$ . For  $1 \leq k \leq n$ , let

$$H_k = \frac{1}{\binom{n-1}{k-1}} \sum_{\alpha:|\alpha|=k} H(X_\alpha).$$

Prove that

$$H_1 \ge H_2 \ge \cdots \ge H_n$$

This sequence of inequalities, due to Han [134], is a generalization of the independence bound for entropy (Theorem 2.39). See Problem 5 in Chapter 21 for an application of these inequalities.

- 19. Prove the divergence inequality by using the log-sum inequality.
- 20. Prove that D(p||q) is convex in the pair (p,q), i.e., if  $(p_1,q_1)$  and  $(p_2,q_2)$  are two pairs of probability distributions on a common alphabet, then

$$D(\lambda p_1 + \overline{\lambda} p_2 \| \lambda q_1 + \overline{\lambda} q_2) \le \lambda D(p_1 \| q_1) + \overline{\lambda} D(p_2 \| q_2)$$

for all  $0 \leq \lambda \leq 1$ , where  $\overline{\lambda} = 1 - \lambda$ .

- 21. Let  $p_{XY}$  and  $q_{XY}$  be two probability distributions on  $\mathcal{X} \times \mathcal{Y}$ . Prove that  $D(p_{XY} || q_{XY}) \ge D(p_X || q_X)$ .
- 22. Pinsker's inequality Let V(p,q) denotes the variational distance between two probability distributions p and q on a common alphabet  $\mathcal{X}$ . We will determine the largest c which satisfies

$$D(p||q) \ge cd^2(p,q).$$

- a) Let  $A = \{x : p(x) \ge q(x)\}, \hat{p} = \{p(A), 1 p(A)\}, \text{ and } \hat{q} = \{q(A), 1 q(A)\}.$  Show that  $D(p||q) \ge D(\hat{p}||\hat{q})$  and  $V(p,q) = V(\hat{p},\hat{q}).$
- b) Show that toward determining the largest value of c, we only have to consider the case when  $\mathcal{X}$  is binary.

c) By virtue of b), it suffices to determine the largest c such that

$$p\log\frac{p}{q} + (1-p)\log\frac{1-p}{1-q} - 4c(p-q)^2 \ge 0$$

for all  $0 \le p, q \le 1$ , with the convention that  $0 \log \frac{0}{b} = 0$  for  $b \ge 0$ and  $a \log \frac{a}{0} = \infty$  for a > 0. By observing that equality in the above holds if p = q and considering the derivative of the left hand side with respect to q, show that the largest value of c is equal to  $(2 \ln 2)^{-1}$ .

- 23. Let p and  $q_k, k \ge 1$  be probability distributions on a common alphabet. Show that if  $q_k$  converges to p in divergence, then it also converges to p in variational distance.
- 24. Find a necessary and sufficient condition for Fano's inequality to be tight.
- 25. Determine the probability distribution defined on  $\{0, 1, \dots, n\}$  that maximizes the entropy subject to the constraint that the mean is equal to m, where  $0 \le m \le n$ .
- 26. Show that for a stationary source  $\{X_k\}, \frac{1}{n}H(X_1, X_2, \dots, X_n)$  is nonincreasing in n.
- 27. For real numbers  $\alpha > 1$  and  $\beta > 0$  and an integer  $n \ge \alpha$ , define the probability distribution

$$\mathcal{D}_{n}^{(\alpha,\beta)} = \left\{ 1 - \left(\frac{\log \alpha}{\log n}\right)^{\beta}, \underbrace{\frac{1}{n} \left(\frac{\log \alpha}{\log n}\right)^{\beta}, \cdots, \frac{1}{n} \left(\frac{\log \alpha}{\log n}\right)^{\beta}}_{n}, 0, 0, \cdots \right\}.$$

Let  $\nu = \{1, 0, 0, ...\}$  be the deterministic distribution.

- a) Show that  $\lim_{n\to\infty} D\left(\nu||\mathcal{D}_n^{(\alpha,\beta)}\right) = 0.$ b) Determine  $\lim_{n\to\infty} H\left(\mathcal{D}_n^{(\alpha,\beta)}\right).$
- 28. Discontinuity of entropy with respect to convergence in divergence Let  $\mathcal{P}$ be the set of all probability distributions on a countable alphabet. A function  $f: \mathcal{P} \to \Re$  is continuous with respect to convergence in divergence at  $P \in \mathcal{P}$  if for any  $\epsilon > 0$ , there exists  $\delta > 0$  such that  $|f(Q) - f(P)| < \epsilon$ for all  $Q \in \mathcal{P}$  satisfying  $D(P||Q) < \delta$ ; otherwise, f is discontinuous at P.
  - a) Let  $H: \mathcal{P} \to \Re$  be the entropy function. Show that H is discontinuous at the deterministic distribution  $\nu = \{1, 0, 0, \dots, \}$ . Hint: Use the results in Problem 27.
  - b) Show that H is discontinuous at  $P = \{p_0, p_1, p_2, \dots\}$  for all P such that  $H(P) < \infty$ . Hint: Consider the probability distribution

$$Q_n = \left\{ p_0 - \frac{p_0}{\sqrt{\log n}}, p_1 + \frac{p_0}{n\sqrt{\log n}}, p_2 + \frac{p_0}{n\sqrt{\log n}}, \cdots, \\ p_n + \frac{p_0}{n\sqrt{\log n}}, p_{n+1}, p_{n+2}, \cdots \right\}$$

for large n.

- 462 Information Measures
- 29. Discontinuity of entropy with respect to convergence in variational dis*tance* Refer to Problem 28. The continuity of a function  $f : \mathcal{P} \to \Re$  with respect to convergence in variational distance can be defined similarly.
  - a) Show that if a function f is continuous with respect to convergence in variational distance, then it is also continuous with respect to convergence in divergence. Hint: Use Pinsker's inequality.
  - b) Repeat b) in Problem 28 with continuity defined with respect to convergence in variational distance.
- 30. Continuity of the entropy function for a fixed finite alphabet Refer to Problems 28 and 29. Suppose the domain of H is confined to  $\mathcal{P}'$ , the set of all probability distributions on a fixed finite alphabet. Show that H is continuous with respect to convergence in divergence.
- 31. Let  $\mathbf{p} = \{p_1, p_2, \dots, p_n\}$  and  $\mathbf{q} = \{q_1, q_2, \dots, q_n\}$  be two sets of real numbers such that  $p_i \ge p_{i'}$  and  $q_i \ge q_{i'}$  for all i < i'. We say that  $\mathbf{p}$  is *majorized* by  $\mathbf{q}$  if  $\sum_{i=1}^m p_i \le \sum_{j=1}^m q_j$  for all m = 1, 2, ..., n, where equality holds when m = n. A function  $f : \Re^n \to \Re$  is *Schur-concave* if  $f(\mathbf{p}) \ge f(\mathbf{q})$ whenever  $\mathbf{p}$  is majorized by  $\mathbf{q}$ . Now let  $\mathbf{p}$  and  $\mathbf{q}$  be probability distributions. We will show in the following steps that  $H(\cdot)$  is Schur-concave.
  - a) Show that for  $\mathbf{p} \neq \mathbf{q}$ , there exist  $1 \leq j < k \leq n$  which satisfy the following:
    - i) j is the largest index i such that  $p_i < q_i$
    - ii) k is the smallest index i such that i > j and  $p_i > q_i$
    - iii)  $p_i = q_i$  for all j < i < k.
  - b) Consider the distribution  $\mathbf{q}^* = \{q_1^*, q_2^*, \cdots, q_n^*\}$  defined by  $q_i^* = q_i$  for  $i \neq j, k$  and

$$(q_j^*, q_k^*) = \begin{cases} (p_j, q_k + (q_j - p_j)) & \text{if } p_k - q_k \ge q_j - p_j \\ (q_j - (p_k - q_k), p_k) & \text{if } p_k - q_k < q_j - p_j. \end{cases}$$

Note that either  $q_j^* = p_j$  or  $q_k^* = p_k$ . Show that i)  $q_i^* \ge q_{i'}^*$  for all  $i \le i'$ ii)  $\sum_{i=1}^m p_i \le \sum_{i=1}^m q_i^*$  for all  $m = 1, 2, \dots, n$ iii)  $H(a^*) > H(a)$ 

- iii)  $\widetilde{H}(\mathbf{q}^*) \geq H(\mathbf{q})$
- c) Prove that  $H(\mathbf{p}) \geq H(\mathbf{q})$  by induction on the Hamming distance between  $\mathbf{p}$  and  $\mathbf{q}$ , i.e., the number of places where  $\mathbf{p}$  and  $\mathbf{q}$  differ.

In general, if a concave function f is symmetric, i.e.  $f(\mathbf{p}) = f(\mathbf{p}')$  where  $\mathbf{p}'$  is a permutation of  $\mathbf{p}$ , then f is Schur-concave. We refer the reader to [221] for the theory of majorization. (Hardy, Littlewood, and Pólya [140].)

# **Historical Notes**

The concept of entropy has its root in thermodynamics. Shannon [291] was the first to use entropy as a measure of information. Informational divergence was introduced by Kullback and Leibler [194], and it has been studied extensively by Csiszár [73] and Amari [14].

The materials in this chapter can be found in most textbooks in information theory. The main concepts and results are due to Shannon [291]. Pinsker's inequality is due to Pinsker [263]. Fano's inequality has its origin in the converse proof of the channel coding theorem (to be discussed in Chapter 7) by Fano [95]. Generalizations of Fano's inequality which apply to random variables with countable alphabets have been obtained by Han and Verdú [139] and by Ho [148] (see also [151]). Maximum entropy, a concept in statistical mechanics, was expounded in Jaynes [167].

# The *I*-Measure

In Chapter 2, we have shown the relationship between Shannon's information measures for two random variables by the diagram in Figure 2.2. For convenience, Figure 2.2 is reproduced in Figure 3.1 with the random variables X and Y replaced by  $X_1$  and  $X_2$ , respectively. This diagram suggests that

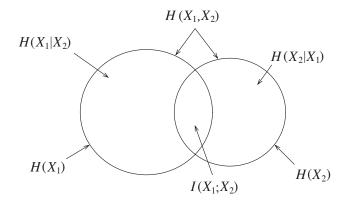


Fig. 3.1. Relationship between entropies and mutual information for two random variables.

Shannon's information measures for any  $n \ge 2$  random variables may have a set-theoretic structure.

In this chapter, we develop a theory which establishes a one-to-one correspondence between Shannon's information measures and set theory in full generality. With this correspondence, manipulations of Shannon's information measures can be viewed as set operations, thus allowing the rich suite of tools in set theory to be used in information theory. Moreover, the structure of Shannon's information measures can easily be visualized by means of an *information diagram* if four or fewer random variables are involved. The use 50 3 The *I*-Measure

of information diagrams simplifies many difficult proofs in information theory problems. More importantly, these results, which may be difficult to discover in the first place, can easily be obtained by inspection of an information diagram.

The main concepts to be used in this chapter are from measure theory. However, it is not necessary for the reader to know measure theory to read this chapter.

### 3.1 Preliminaries

In this section, we introduce a few basic concepts in measure theory which will be used subsequently. These concepts will be illustrated by simple examples.

**Definition 3.1.** The field  $\mathcal{F}_n$  generated by sets  $\tilde{X}_1, \tilde{X}_2, \dots, \tilde{X}_n$  is the collection of sets which can be obtained by any sequence of usual set operations (union, intersection, complement, and difference) on  $\tilde{X}_1, \tilde{X}_2, \dots, \tilde{X}_n$ .

**Definition 3.2.** The atoms of  $\mathcal{F}_n$  are sets of the form  $\cap_{i=1}^n Y_i$ , where  $Y_i$  is either  $\tilde{X}_i$  or  $\tilde{X}_i^c$ , the complement of  $\tilde{X}_i$ .

There are  $2^n$  atoms and  $2^{2^n}$  sets in  $\mathcal{F}_n$ . Evidently, all the atoms in  $\mathcal{F}_n$  are disjoint, and each set in  $\mathcal{F}_n$  can be expressed uniquely as the union of a subset of the atoms of  $\mathcal{F}_n^{-1}$ . We assume that the sets  $\tilde{X}_1, \tilde{X}_2, \dots, \tilde{X}_n$  intersect with each other generically, i.e., all the atoms of  $\mathcal{F}_n$  are nonempty unless otherwise specified.

*Example 3.3.* The sets  $\tilde{X}_1$  and  $\tilde{X}_2$  generate the field  $\mathcal{F}_2$ . The atoms of  $\mathcal{F}_2$  are

$$\tilde{X}_1 \cap \tilde{X}_2, \tilde{X}_1^c \cap \tilde{X}_2, \tilde{X}_1 \cap \tilde{X}_2^c, \tilde{X}_1^c \cap \tilde{X}_2^c,$$
(3.1)

which are represented by the four distinct regions in the Venn diagram in Figure 3.2. The field  $\mathcal{F}_2$  consists of the unions of subsets of the atoms in (3.1). There are a total of 16 sets in  $\mathcal{F}_2$ , which are precisely all the sets which can be obtained from  $\tilde{X}_1$  and  $\tilde{X}_2$  by the usual set operations.

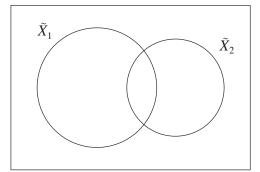
**Definition 3.4.** A real function  $\mu$  defined on  $\mathcal{F}_n$  is called a signed measure if it is set-additive, i.e., for disjoint A and B in  $\mathcal{F}_n$ ,

$$\mu(A \cup B) = \mu(A) + \mu(B).$$
(3.2)

For a signed measure  $\mu$ , we have

$$\mu(\emptyset) = 0, \tag{3.3}$$

<sup>&</sup>lt;sup>1</sup> We adopt the convention that the union of the empty subset of the atoms of  $\mathcal{F}_n$  is the empty set.



**Fig. 3.2.** The Venn diagram for  $\tilde{X}_1$  and  $\tilde{X}_2$ .

which can be seen as follows. For any A in  $\mathcal{F}_n$ ,

$$\mu(A) = \mu(A \cup \emptyset) = \mu(A) + \mu(\emptyset) \tag{3.4}$$

by set-additivity because A and  $\emptyset$  are disjoint, which implies (3.3).

A signed measure  $\mu$  on  $\mathcal{F}_n$  is completely specified by its values on the atoms of  $\mathcal{F}_n$ . The values of  $\mu$  on the other sets in  $\mathcal{F}_n$  can be obtained via set-additivity.

*Example 3.5.* A signed measure  $\mu$  on  $\mathcal{F}_2$  is completely specified by the values

$$\mu(\tilde{X}_1 \cap \tilde{X}_2), \mu(\tilde{X}_1^c \cap \tilde{X}_2), \mu(\tilde{X}_1 \cap \tilde{X}_2^c), \mu(\tilde{X}_1^c \cap \tilde{X}_2^c).$$
(3.5)

The value of  $\mu$  on  $\tilde{X}_1$ , for example, can be obtained as

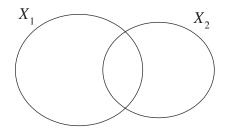
$$\mu(\tilde{X}_1) = \mu((\tilde{X}_1 \cap \tilde{X}_2) \cup (\tilde{X}_1 \cap \tilde{X}_2^c))$$
(3.6)

$$= \mu(\tilde{X}_1 \cap \tilde{X}_2) + \mu(\tilde{X}_1 \cap \tilde{X}_2^c).$$
(3.7)

### 3.2 The *I*-Measure for Two Random Variables

To fix ideas, we first formulate in this section the one-to-one correspondence between Shannon's information measures and set theory for two random variables. For random variables  $X_1$  and  $X_2$ , let  $\tilde{X}_1$  and  $\tilde{X}_2$  be sets corresponding to  $X_1$  and  $X_2$ , respectively. The sets  $\tilde{X}_1$  and  $\tilde{X}_2$  generates the field  $\mathcal{F}_2$  whose atoms are listed in (3.1). In our formulation, we set the universal set  $\Omega$  to  $\tilde{X}_1 \cup \tilde{X}_2$  for reasons which will become clear later. With this choice of  $\Omega$ , the Venn diagram for  $X_1$  and  $X_2$  is represented by the diagram in Figure 3.3. For simplicity, the sets  $X_1$  and  $X_2$  are respectively labeled by  $X_1$  and  $X_2$  in the diagram. We call this the *information diagram* for the random variables  $X_1$ and  $X_2$ . In this diagram, the universal set, which is the union of  $X_1$  and  $X_2$ , is not shown explicitly just as in a usual Venn diagram. Note that with our

51



**Fig. 3.3.** The generic information diagram for  $X_1$  and  $X_2$ .

choice of the universal set, the atom  $\tilde{X}_1^c \cap \tilde{X}_2^c$  degenerates to the empty set, because

$$\tilde{X}_1^c \cap \tilde{X}_2^c = (\tilde{X}_1 \cup \tilde{X}_2)^c = \Omega^c = \emptyset.$$
(3.8)

Thus this atom is not shown in the information diagram in Figure 3.3.

For random variables  $X_1$  and  $X_2$ , the Shannon's information measures are

$$H(X_1), H(X_2), H(X_1|X_2), H(X_2|X_1), H(X_1, X_2), I(X_1; X_2).$$
 (3.9)

Writing  $A \cap B^c$  as A - B, we now define a signed measure<sup>2</sup>  $\mu^*$  by

$$\mu^*(X_1 - X_2) = H(X_1 | X_2) \tag{3.10}$$

$$\mu^*(\tilde{X}_2 - \tilde{X}_1) = H(X_2|X_1), \qquad (3.11)$$

and

$$\mu^*(X_1 \cap X_2) = I(X_1; X_2). \tag{3.12}$$

These are the values of  $\mu^*$  on the nonempty atoms of  $\mathcal{F}_2$  (i.e., atoms of  $\mathcal{F}_2$ other than  $\tilde{X}_1^c \cap \tilde{X}_2^c$ ). The values of  $\mu^*$  on the other sets in  $\mathcal{F}_2$  can be obtained via set-additivity. In particular, the relations

$$\mu^*(\tilde{X}_1 \cup \tilde{X}_2) = H(X_1, X_2) \tag{3.13}$$

$$\mu^*(X_1) = H(X_1), \tag{3.14}$$

and

$$\mu^*(\tilde{X}_2) = H(X_2) \tag{3.15}$$

can readily be verified. For example, (3.13) is seen to be true by considering

$$= \mu^{*}(\tilde{X}_{1} - \tilde{X}_{2}) + \mu^{*}(\tilde{X}_{2} - \tilde{X}_{1}) + \mu^{*}(\tilde{X}_{1} \cap \tilde{X}_{2})$$

$$(3.16)$$

$$U(X + X) + U(X + X) + U(X + X)$$

$$(3.17)$$

$$= H(X_1|X_2) + H(X_2|X_1) + I(X_1;X_2)$$
(3.17)

$$= H(X_1, X_2). (3.18)$$

 $<sup>^2</sup>$  It happens that  $\mu^*$  defined here for n=2 assumes only nonnegative values, but we will see in Section 3.4 that  $\mu^*$  can assume negative values for  $n \geq 3$ .

The right hand sides of (3.10) to (3.15) are the six Shannon's information measures for  $X_1$  and  $X_2$  in (3.9). Now observe that (3.10) to (3.15) are consistent with how the Shannon's information measures on the right hand side are identified in Figure 3.1, with the left circle and the right circle representing the sets  $\tilde{X}_1$  and  $\tilde{X}_2$ , respectively. Specifically, in each of these equations, the left hand side and the right hand side correspond to each other via the following substitution of symbols:

$$H/I \leftrightarrow \mu^*$$

$$, \leftrightarrow \cup$$

$$; \leftrightarrow \cap$$

$$| \leftrightarrow -.$$

$$(3.19)$$

Note that we make no distinction between the symbols H and I in this substitution. Thus, for two random variables  $X_1$  and  $X_2$ , Shannon's information measures can be regarded formally as a signed measure on  $\mathcal{F}_2$ . We will refer to  $\mu^*$  as the *I-Measure* for the random variables  $X_1$  and  $X_2^3$ .

Upon realizing that Shannon's information measures can be viewed as a signed measure, we can apply the rich family of operations in set theory to information theory. This explains why Figure 3.1 or Figure 3.3 represents the relationships among all Shannon's information measures for two random variables correctly. As an example, consider the following set identity which is readily identified in Figure 3.3:

$$\mu^*(\tilde{X}_1 \cup \tilde{X}_2) = \mu^*(\tilde{X}_1) + \mu^*(\tilde{X}_2) - \mu^*(\tilde{X}_1 \cap \tilde{X}_2)$$
(3.20)

This identity is a special case of the inclusion-exclusion formula in set theory. By means of the substitution of symbols in (3.19), we immediately obtain the information identity

$$H(X_1, X_2) = H(X_1) + H(X_2) - I(X_1; X_2).$$
(3.21)

We end this section with a remark. The value of  $\mu^*$  on the atom  $\tilde{X}_1^c \cap \tilde{X}_2^c$ has no apparent information-theoretic meaning. In our formulation, we set the universal set  $\Omega$  to  $\tilde{X}_1 \cup \tilde{X}_2$  so that the atom  $\tilde{X}_1^c \cap \tilde{X}_2^c$  degenerates to the empty set. Then  $\mu^*(\tilde{X}_1^c \cap \tilde{X}_2^c)$  naturally vanishes because  $\mu^*$  is a measure, so that  $\mu^*$ is completely specified by all Shannon's information measures involving the random variables  $X_1$  and  $X_2$ .

# 3.3 Construction of the *I*-Measure $\mu^*$

We have constructed the *I*-Measure for two random variables in the last section. We now construct the *I*-Measure for any  $n \ge 2$  random variables.

<sup>&</sup>lt;sup>3</sup> The reader should not confuse  $\mu^*$  with the probability measure defining the random variables  $X_1$  and  $X_2$ . The former, however, is determined by the latter.

#### 54 3 The *I*-Measure

Consider *n* random variables  $X_1, X_2, \dots, X_n$ . For any random variable *X*, let  $\tilde{X}$  be a set corresponding to *X*. Let

$$\mathcal{N}_n = \{1, 2, \cdots, n\}.$$
 (3.22)

Define the universal set  $\Omega$  to be the union of the sets  $\tilde{X}_1, \tilde{X}_2, \dots, \tilde{X}_n$ , i.e.,

$$\Omega = \bigcup_{i \in \mathcal{N}_n} \tilde{X}_i. \tag{3.23}$$

We use  $\mathcal{F}_n$  to denote the field generated by  $\tilde{X}_1, \tilde{X}_2, \cdots, \tilde{X}_n$ . The set

$$A_0 = \bigcap_{i \in \mathcal{N}_n} \tilde{X}_i^c \tag{3.24}$$

is called the *empty atom* of  $\mathcal{F}_n$  because

$$\bigcap_{i \in \mathcal{N}_n} \tilde{X}_i^c = \left(\bigcup_{i \in \mathcal{N}_n} \tilde{X}_i\right)^c = \Omega^c = \emptyset.$$
(3.25)

All the atoms of  $\mathcal{F}_n$  other than  $A_0$  are called *nonempty atoms*.

Let  $\mathcal{A}$  be the set of all nonempty atoms of  $\mathcal{F}_n$ . Then  $|\mathcal{A}|$ , the cardinality of  $\mathcal{A}$ , is equal to  $2^n - 1$ . A signed measure  $\mu$  on  $\mathcal{F}_n$  is completely specified by the values of  $\mu$  on the nonempty atoms of  $\mathcal{F}_n$ .

To simplify notation, we will use  $X_G$  to denote  $(X_i, i \in G)$  and  $\tilde{X}_G$  to denote  $\bigcup_{i \in G} \tilde{X}_i$  for any nonempty subset G of  $\mathcal{N}_n$ .

### Theorem 3.6. Let

$$\mathcal{B} = \left\{ \tilde{X}_G : G \text{ is a nonempty subset of } \mathcal{N}_n \right\}.$$
(3.26)

Then a signed measure  $\mu$  on  $\mathcal{F}_n$  is completely specified by  $\{\mu(B), B \in \mathcal{B}\}$ , which can be any set of real numbers.

*Proof.* The number of elements in  $\mathcal{B}$  is equal to the number of nonempty subsets of  $\mathcal{N}_n$ , which is  $2^n - 1$ . Thus  $|\mathcal{A}| = |\mathcal{B}| = 2^n - 1$ . Let  $k = 2^n - 1$ . Let **u** be a column k-vector of  $\mu(A), A \in \mathcal{A}$ , and **h** be a column k-vector of  $\mu(B), B \in \mathcal{B}$ . Since all the sets in  $\mathcal{B}$  can expressed uniquely as the union of some nonempty atoms in  $\mathcal{A}$ , by the set-additivity of  $\mu$ , for each  $B \in \mathcal{B}, \mu(B)$ can be expressed uniquely as the sum of some components of **u**. Thus

$$\mathbf{h} = \mathbf{C}_n \mathbf{u},\tag{3.27}$$

where  $\mathbf{C}_n$  is a *unique*  $k \times k$  matrix. On the other hand, it can be shown (see Appendix 3.6) that for each  $A \in \mathcal{A}$ ,  $\mu(A)$  can be expressed as a linear combination of  $\mu(B), B \in \mathcal{B}$  by applications, if necessary, of the following two identities:

3.3 Construction of the *I*-Measure  $\mu^*$  55

$$\mu(A \cap B - C) = \mu(A - C) + \mu(B - C) - \mu(A \cup B - C)$$
(3.28)

$$\mu(A - B) = \mu(A \cup B) - \mu(B). \tag{3.29}$$

However, the existence of the said expression does not imply its uniqueness. Nevertheless, we can write

$$\mathbf{u} = \mathbf{D}_n \mathbf{h} \tag{3.30}$$

for some  $k \times k$  matrix  $\mathbf{D}_n$ . Upon substituting (3.27) into (3.30), we obtain

$$\mathbf{u} = (\mathbf{D}_n \mathbf{C}_n) \mathbf{u},\tag{3.31}$$

which implies that  $\mathbf{D}_n$  is the inverse of  $\mathbf{C}_n$  as (3.31) holds regardless of the choice of  $\mu$ . Since  $\mathbf{C}_n$  is unique, so is  $\mathbf{D}_n$ . Therefore,  $\mu(A), A \in \mathcal{A}$  are uniquely determined once  $\mu(B), B \in \mathcal{B}$  are specified. Hence, a signed measure  $\mu$  on  $\mathcal{F}_n$  is completely specified by  $\{\mu(B), B \in \mathcal{B}\}$ , which can be any set of real numbers. The theorem is proved.  $\Box$ 

We now prove the following two lemmas which are related by the substitution of symbols in (3.19).

### Lemma 3.7.

$$\mu(A \cap B - C) = \mu(A \cup C) + \mu(B \cup C) - \mu(A \cup B \cup C) - \mu(C).$$
(3.32)

*Proof.* From (3.28) and (3.29), we have

$$\mu(A \cap B - C) = \mu(A - C) + \mu(B - C) - \mu(A \cup B - C)$$
(3.33)

$$= (\mu(A \cup C) - \mu(C)) + (\mu(B \cup C) - \mu(C)) -(\mu(A \cup B \cup C) - \mu(C))$$
(3.34)

$$-(\mu(A \cup B \cup C) - \mu(C))$$
(3.34)

$$= \mu(A \cup C) + \mu(B \cup C) - \mu(A \cup B \cup C) - \mu(C).$$
(3.35)

The lemma is proved.  $\hfill\square$ 

#### Lemma 3.8.

$$I(X;Y|Z) = H(X,Z) + H(Y,Z) - H(X,Y,Z) - H(Z).$$
(3.36)

Proof. Consider

$$I(X;Y|Z) = H(X|Z) - H(X|Y,Z)$$
(3.37)  
=  $H(X,Z) - H(Z) - (H(X,Y,Z) - H(Y,Z))$ (3.38)

$$= H(X,Z) + H(Y,Z) - H(X,Y,Z) - H(Z).$$
(3.39)

56 3 The *I*-Measure

The lemma is proved.  $\Box$ 

We now construct the *I*-Measure  $\mu^*$  on  $\mathcal{F}_n$  using Theorem 3.6 by defining

$$\mu^*(\tilde{X}_G) = H(X_G) \tag{3.40}$$

for all nonempty subsets G of  $\mathcal{N}_n$ . In order for  $\mu^*$  to be meaningful, it has to be consistent with all Shannon's information measures (via the substitution of symbols in (3.19)). In that case, the following must hold for all (not necessarily disjoint) subsets G, G', G'' of  $\mathcal{N}_n$  where G and G' are nonempty:

$$\mu^*(\tilde{X}_G \cap \tilde{X}_{G'} - \tilde{X}_{G''}) = I(X_G; X_{G'} | X_{G''}).$$
(3.41)

When  $G'' = \emptyset$ , (3.41) becomes

$$\mu^*(\tilde{X}_G \cap \tilde{X}_{G'}) = I(X_G; X_{G'}).$$
(3.42)

When G = G', (3.41) becomes

$$\mu^*(\tilde{X}_G - \tilde{X}_{G''}) = H(X_G | X_{G''}). \tag{3.43}$$

When G = G' and  $G'' = \emptyset$ , (3.41) becomes

$$\mu^*(X_G) = H(X_G). \tag{3.44}$$

Thus (3.41) covers all the four cases of Shannon's information measures, and it is the necessary and sufficient condition for  $\mu^*$  to be consistent with all Shannon's information measures.

**Theorem 3.9.**  $\mu^*$  is the unique signed measure on  $\mathcal{F}_n$  which is consistent with all Shannon's information measures.

Proof. Consider

$$\mu^{*}(X_{G} \cap X_{G'} - X_{G''}) = \mu^{*}(\tilde{X}_{G \cup G''}) + \mu^{*}(\tilde{X}_{G' \cup G''}) - \mu^{*}(\tilde{X}_{G \cup G' \cup G''}) - \mu^{*}(\tilde{X}_{G''})$$
(3.45)

$$= H(\Lambda_{G\cup G''}) + H(\Lambda_{G'\cup G''}) - H(\Lambda_{G\cup G'\cup G''}) - H(\Lambda_{G''})$$

$$(3.40)$$

$$= I(X_G; X_{G'}|X_{G''}), (3.47)$$

where (3.45) and (3.47) follow from Lemmas 3.7 and 3.8, respectively, and (3.46) follows from (3.40), the definition of  $\mu^*$ . Thus we have proved (3.41), i.e.,  $\mu^*$  is consistent with all Shannon's information measures.

In order that  $\mu^*$  is consistent with all Shannon's information measures, for all nonempty subsets G of  $\mathcal{N}_n$ ,  $\mu^*$  has to satisfy (3.44), which in fact is the definition of  $\mu^*$  in (3.40). Therefore,  $\mu^*$  is the unique signed measure on  $\mathcal{F}_n$ which is consistent with all Shannon's information measures.  $\Box$ 

# 3.4 $\mu^*$ Can be Negative

In the previous sections, we have been cautious in referring to the *I*-Measure  $\mu^*$  as a signed measure instead of a measure<sup>4</sup>. In this section, we show that  $\mu^*$  in fact can take negative values for  $n \geq 3$ .

For n = 2, the three nonempty atoms of  $\mathcal{F}_2$  are

$$\tilde{X}_1 \cap \tilde{X}_2, \tilde{X}_1 - \tilde{X}_2, \tilde{X}_2 - \tilde{X}_1.$$
 (3.48)

The values of  $\mu^*$  on these atoms are respectively

$$I(X_1; X_2), H(X_1|X_2), H(X_2|X_1).$$
(3.49)

These quantities are Shannon's information measures and hence nonnegative by the basic inequalities. Therefore,  $\mu^*$  is always nonnegative for n = 2.

For n = 3, the seven nonempty atoms of  $\mathcal{F}_3$  are

$$\tilde{X}_i - \tilde{X}_{\{j,k\}}, \tilde{X}_i \cap \tilde{X}_j - \tilde{X}_k, \tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3,$$
(3.50)

where  $1 \le i < j < k \le 3$ . The values of  $\mu^*$  on the first two types of atoms are

$$\mu^*(\tilde{X}_i - \tilde{X}_{\{j,k\}}) = H(X_i | X_j, X_k)$$
(3.51)

and

$$\mu^*(\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_k) = I(X_i; X_j | X_k), \qquad (3.52)$$

respectively, which are Shannon's information measures and therefore nonnegative. However,  $\mu^*(\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3)$  does not correspond to a Shannon's information measure. In the next example, we show that  $\mu^*(\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3)$  can actually be negative.

*Example 3.10.* In this example, all entropies are in the base 2. Let  $X_1$  and  $X_2$  be independent binary random variables with

$$\Pr\{X_i = 0\} = \Pr\{X_i = 1\} = 0.5, \tag{3.53}$$

i = 1, 2. Let

$$X_3 = (X_1 + X_2) \mod 2. \tag{3.54}$$

It is easy to check that  $X_3$  has the same marginal distribution as  $X_1$  and  $X_2$ . Thus,

$$H(X_i) = 1 \tag{3.55}$$

for i = 1, 2, 3. Moreover,  $X_1, X_2$ , and  $X_3$  are pairwise independent. Therefore,

$$H(X_i, X_j) = 2$$
 (3.56)

and

<sup>&</sup>lt;sup>4</sup> A measure can assume only nonnegative values.

58 3 The *I*-Measure

$$I(X_i; X_j) = 0 (3.57)$$

for  $1 \leq i < j \leq 3$ . We further see from (3.54) that each random variable is a function of the other two random variables. Then by the chain rule for entropy, we have

$$H(X_1, X_2, X_3) = H(X_1, X_2) + H(X_3 | X_1, X_2)$$
(3.58)

$$= 2 + 0$$
 (3.59)

$$= 2.$$
 (3.60)

Now for  $1 \le i < j < k \le 3$ ,

$$I(X_i; X_j | X_k) = H(X_i, X_k) + H(X_j, X_k) - H(X_1, X_2, X_3) - H(X_k)$$
(3.61)

$$= 2 + 2 - 2 - 1 \tag{3.62}$$

$$= 1,$$
 (3.63)

where we have invoked Lemma 3.8. It then follows that

$$\mu^*(\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3) = \mu^*(\tilde{X}_1 \cap \tilde{X}_2) - \mu^*(\tilde{X}_1 \cap \tilde{X}_2 - \tilde{X}_3)$$
(3.64)

$$= I(X_1; X_2) - I(X_1; X_2 | X_3)$$
(3.65)

$$= 0 - 1$$
 (3.66)

$$= -1.$$
 (3.67)

Thus  $\mu^*$  takes a negative value on the atom  $\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3$ .

Motivated by the substitution of symbols in (3.19) for Shannon's information measures, we will write  $\mu^*(\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3)$  as  $I(X_1; X_2; X_3)$ . In general, we will write

$$\mu^*(\tilde{X}_{G_1} \cap \tilde{X}_{G_2} \cap \dots \cap \tilde{X}_{G_m} - \tilde{X}_F) \tag{3.68}$$

as

$$I(X_{G_1}; X_{G_2}; \cdots; X_{G_m} | X_F)$$
(3.69)

and refer to it as the *mutual information* between  $X_{G_1}, X_{G_2}, \dots, X_{G_m}$  conditioning on  $X_F$ . Then (3.64) in the above example can be written as

$$I(X_1; X_2; X_3) = I(X_1; X_2) - I(X_1; X_2 | X_3).$$
(3.70)

For this example,  $I(X_1; X_2; X_3) < 0$ , which implies

$$I(X_1; X_2 | X_3) > I(X_1; X_2).$$
(3.71)

Therefore, unlike entropy, the mutual information between two random variables can be increased by conditioning on a third random variable. Also, we note in (3.70) that although the expression on the right hand side is not symbolically symmetrical in  $X_1$ ,  $X_2$ , and  $X_3$ , we see from the left hand side that it is in fact symmetrical in  $X_1$ ,  $X_2$ , and  $X_3$ .

### 3.5 Information Diagrams

We have established in Section 3.3 a one-to-one correspondence between Shannon's information measures and set theory. Therefore, it is valid to use an *information diagram*, which is a variation of a Venn diagram, to represent the relationship between Shannon's information measures.

For simplicity, a set  $X_i$  will be labeled by  $X_i$  in an information diagram. We have seen the generic information diagram for n = 2 in Figure 3.3. A generic information diagram for n = 3 is shown in Figure 3.4. The information-

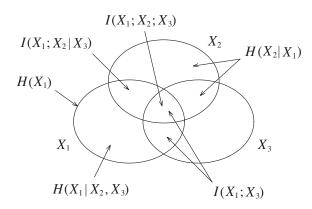


Fig. 3.4. The generic information diagram for  $X_1$ ,  $X_2$ , and  $X_3$ .

theoretic labeling of the values of  $\mu^*$  on some of the sets in  $\mathcal{F}_3$  is shown in the diagram. As an example, the information diagram for the *I*-Measure for random variables  $X_1, X_2$ , and  $X_3$  discussed in Example 3.10 is shown in Figure 3.5.

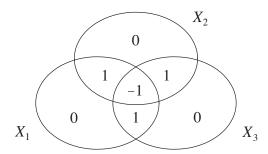


Fig. 3.5. The information diagram for  $X_1$ ,  $X_2$ , and  $X_3$  in Example 3.10.

For  $n \ge 4$ , it is not possible to display an information diagram perfectly in two dimensions. In general, an information diagram for n random variables

needs n-1 dimensions to be displayed perfectly. Nevertheless, for n = 4, an information diagram can be displayed in two dimensions almost perfectly as shown in Figure 3.6. This information diagram is correct in that the region representing the set  $\tilde{X}_4$  splits each atom in Figure 3.4 into two atoms. However, the adjacency of certain atoms are not displayed correctly. For example, the set  $\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_4^c$ , which consists of the atoms  $\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3 \cap \tilde{X}_4^c$  and  $\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3^c \cap \tilde{X}_4^c$ , is not represented by a connected region because the two atoms are not adjacent to each other.

When  $\mu^*$  takes the value zero on an atom A of  $\mathcal{F}_n$ , we do not need to display the atom A in an information diagram because the atom A does not contribute to  $\mu^*(B)$  for any set B containing the atom A. As we will see shortly, this can happen if certain Markov constraints are imposed on the random variables involved, and the information diagram can be simplified accordingly. In a generic information diagram (i.e., when there is no constraint on the random variables), however, all the atoms have to be displayed, as is implied by the next theorem.

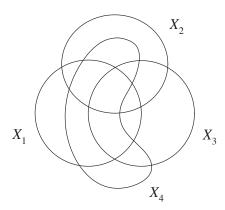
**Theorem 3.11.** If there is no constraint on  $X_1, X_2, \dots, X_n$ , then  $\mu^*$  can take any set of nonnegative values on the nonempty atoms of  $\mathcal{F}_n$ .

*Proof.* We will prove the theorem by constructing a  $\mu^*$  which can take any set of nonnegative values on the nonempty atoms of  $\mathcal{F}_n$ . Recall that  $\mathcal{A}$  is the set of all nonempty atoms of  $\mathcal{F}_n$ . Let  $Y_A, A \in \mathcal{A}$  be mutually independent random variables. Now define the random variables  $X_i, i = 1, 2, \dots, n$  by

 $X_i = (Y_A : A \in \mathcal{A} \text{ and } A \subset \tilde{X}_i).$ (3.72)

We determine the *I*-Measure  $\mu^*$  for  $X_1, X_2, \dots, X_n$  so defined as follows. Since  $Y_A$  are mutually independent, for any nonempty subsets G of  $\mathcal{N}_n$ , we have

$$H(X_G) = \sum_{A \in \mathcal{A}: A \subset \tilde{X}_G} H(Y_A).$$
(3.73)



**Fig. 3.6.** The generic information diagram for  $X_1$ ,  $X_2$ ,  $X_3$ , and  $X_4$ .

On the other hand,

$$H(X_G) = \mu^*(\tilde{X}_G) = \sum_{A \in \mathcal{A}: A \subset \tilde{X}_G} \mu^*(A).$$
(3.74)

Equating the right hand sides of (3.73) and (3.74), we have

A

$$\sum_{\substack{\in \mathcal{A}: A \subset \tilde{X}_G}} H(Y_A) = \sum_{\substack{A \in \mathcal{A}: A \subset \tilde{X}_G}} \mu^*(A).$$
(3.75)

Evidently, we can make the above equality hold for all nonempty subsets G of  $\mathcal{N}_n$  by taking

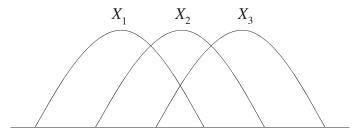
$$\mu^*(A) = H(Y_A) \tag{3.76}$$

for all  $A \in \mathcal{A}$ . By the uniqueness of  $\mu^*$ , this is also the only possibility for  $\mu^*$ . Since  $H(Y_A)$  can take any nonnegative value by Corollary 2.44,  $\mu^*$  can take any set of nonnegative values on the nonempty atoms of  $\mathcal{F}_n$ . The theorem is proved.  $\Box$ 

In the rest of this section, we explore the structure of Shannon's information measures when  $X_1 \to X_2 \to \cdots \to X_n$  forms a Markov chain. To start with, we consider n = 3, i.e.,  $X_1 \to X_2 \to X_3$  forms a Markov chain. Since

$$\mu^*(\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3) = I(X_1; X_3 | X_2) = 0, \qquad (3.77)$$

the atom  $\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3$  does not have to be displayed in an information diagram. As such, in constructing the information diagram, the regions representing the random variables  $X_1$ ,  $X_2$ , and  $X_3$  should overlap with each other such that the region corresponding to the atom  $\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3$  is empty, while the regions corresponding to all other nonempty atoms are nonempty. Figure 3.7 shows such a construction, in which each random variable is represented by a mountain<sup>5</sup>. From Figure 3.7, we see that  $\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3$ , as the only atom



**Fig. 3.7.** The information diagram for the Markov chain  $X_1 \to X_2 \to X_3$ .

<sup>&</sup>lt;sup>5</sup> This form of an information diagram for a Markov chain first appeared in Kawabata [176].

on which  $\mu^*$  may take a negative value, now becomes identical to the atom  $\tilde{X}_1 \cap \tilde{X}_3$ . Therefore, we have

$$I(X_1; X_2; X_3) = \mu^* (\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3)$$
(3.78)

$$=\mu^*(\tilde{X}_1 \cap \tilde{X}_3) \tag{3.79}$$

$$= I(X_1; X_3) (3.80)$$

$$\geq 0. \tag{3.81}$$

Hence, we conclude that when  $X_1 \to X_2 \to X_3$  forms a Markov chain,  $\mu^*$  is always nonnegative.

Next, we consider n = 4, i.e.,  $X_1 \to X_2 \to X_3 \to X_4$  forms a Markov chain. With reference to Figure 3.6, we first show that under this Markov constraint,  $\mu^*$  always vanishes on certain nonempty atoms:

1. The Markov chain  $X_1 \to X_2 \to X_3$  implies

$$I(X_1; X_3; X_4 | X_2) + I(X_1; X_3 | X_2, X_4) = I(X_1; X_3 | X_2) = 0.$$
(3.82)

2. The Markov chain  $X_1 \to X_2 \to X_4$  implies

$$I(X_1; X_3; X_4 | X_2) + I(X_1; X_4 | X_2, X_3) = I(X_1; X_4 | X_2) = 0.$$
(3.83)

3. The Markov chain  $X_1 \to X_3 \to X_4$  implies

$$I(X_1; X_2; X_4 | X_3) + I(X_1; X_4 | X_2, X_3) = I(X_1; X_4 | X_3) = 0.$$
(3.84)

4. The Markov chain  $X_2 \to X_3 \to X_4$  implies

$$I(X_1; X_2; X_4 | X_3) + I(X_2; X_4 | X_1, X_3) = I(X_2; X_4 | X_3) = 0.$$
(3.85)

5. The Markov chain  $(X_1, X_2) \to X_3 \to X_4$  implies

$$\mathbf{K} X_2; X_4 | X_3) + I(X_1; X_4 | X_2, X_3) + I(X_2; X_4 | X_1, X_3) 
= I(X_1, X_2; X_4 | X_3)$$
(3.86)

$$= 0.$$
 (3.87)

Now (3.82) and (3.83) imply

$$I(X_1; X_4 | X_2, X_3) = I(X_1; X_3 | X_2, X_4),$$
(3.88)

(3.84) and (3.88) imply

$$I(X_1; X_2; X_4 | X_3) = -I(X_1; X_3 | X_2, X_4),$$
(3.89)

and (3.85) and (3.89) imply

$$I(X_2; X_4 | X_1, X_3) = I(X_1; X_3 | X_2, X_4).$$
(3.90)

The terms on the left hand sides of (3.88), (3.89), and (3.90) are the three terms on the left hand side of (3.87). Then we substitute (3.88), (3.89), and (3.90) in (3.87) to obtain

$$\mu^* (\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3 \cap \tilde{X}_4^c) = I(X_1; X_3 | X_2, X_4) = 0.$$
(3.91)

From (3.82), (3.88), (3.89), and (3.90), (3.91) implies

$$\mu^*(\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3 \cap \tilde{X}_4) = I(X_1; X_3; X_4 | X_2) = 0$$
(3.92)

$$\mu^*(X_1 \cap X_2^c \cap X_3^c \cap X_4) = I(X_1; X_4 | X_2, X_3) = 0$$
(3.93)

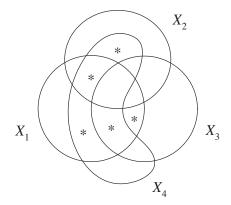
$$\mu^*(X_1 \cap X_2 \cap X_3^c \cap X_4) = I(X_1; X_2; X_4 | X_3) = 0$$
(3.94)

$$\mu^*(X_1^c \cap X_2 \cap X_3^c \cap X_4) = I(X_2; X_4 | X_1, X_3) = 0.$$
(3.95)

From (3.91) to (3.95), we see that  $\mu^*$  always vanishes on the atoms

$$\tilde{X}_{1} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3} \cap \tilde{X}_{4}^{c} 
\tilde{X}_{1} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3} \cap \tilde{X}_{4} 
\tilde{X}_{1} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3}^{c} \cap \tilde{X}_{4} 
\tilde{X}_{1} \cap \tilde{X}_{2} \cap \tilde{X}_{3}^{c} \cap \tilde{X}_{4} 
\tilde{X}_{1}^{c} \cap \tilde{X}_{2} \cap \tilde{X}_{3}^{c} \cap \tilde{X}_{4}$$
(3.96)

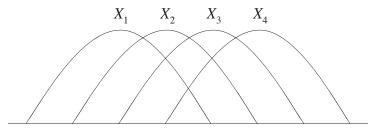
of  $\mathcal{F}_4$ , which we mark by an asterisk in the information diagram in Figure 3.8. In fact, the reader can gain a lot of insight by letting  $I(X_1; X_3 | X_2, X_4) = a \ge 0$ 



**Fig. 3.8.** The atoms of  $\mathcal{F}_4$  on which  $\mu^*$  vanishes when  $X_1 \to X_2 \to X_3 \to X_4$  forms a Markov chain.

in (3.82) and tracing the subsequent steps leading to the above conclusion in the information diagram in Figure 3.6.

It is not necessary to display the five atoms in (3.96) in an information diagram because  $\mu^*$  always vanishes on these atoms. Therefore, in constructing



**Fig. 3.9.** The information diagram for the Markov chain  $X_1 \to X_2 \to X_3 \to X_4$ .

the information diagram, the regions representing the random variables should overlap with each other such that the regions corresponding to these five nonempty atoms are empty, while the regions corresponding to the other ten nonempty atoms, namely

$$\begin{aligned}
\tilde{X}_{1} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3}^{c} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1} \cap \tilde{X}_{2} \cap \tilde{X}_{3}^{c} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1} \cap \tilde{X}_{2} \cap \tilde{X}_{3} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1} \cap \tilde{X}_{2} \cap \tilde{X}_{3} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1}^{c} \cap \tilde{X}_{2} \cap \tilde{X}_{3}^{c} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1}^{c} \cap \tilde{X}_{2} \cap \tilde{X}_{3} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1}^{c} \cap \tilde{X}_{2} \cap \tilde{X}_{3} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1}^{c} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1}^{c} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3} \cap \tilde{X}_{4}^{c} \\
\tilde{X}_{1}^{c} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3} \cap \tilde{X}_{4} \\
\tilde{X}_{1}^{c} \cap \tilde{X}_{2}^{c} \cap \tilde{X}_{3} \cap \tilde{X}_{4},
\end{aligned}$$
(3.97)

are nonempty. Figure 3.9 shows such a construction. The reader should compare the information diagrams in Figures 3.7 and 3.9 and observe that the latter is an extension of the former. From Figure 3.9, we see that the values of  $\mu^*$  on the ten nonempty atoms in (3.97) are equivalent to

$$H(X_{1}|X_{2}, X_{3}, X_{4})$$

$$I(X_{1}; X_{2}|X_{3}, X_{4})$$

$$I(X_{1}; X_{3}|X_{4})$$

$$I(X_{1}; X_{4})$$

$$H(X_{2}|X_{1}, X_{3}, X_{4})$$

$$I(X_{2}; X_{3}|X_{1}; X_{4})$$

$$I(X_{2}; X_{4}|X_{1})$$

$$H(X_{3}|X_{1}, X_{2}, X_{4})$$

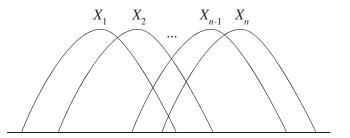
$$I(X_{3}; X_{4}|X_{1}, X_{2})$$

$$H(X_{4}|X_{1}, X_{2}, X_{3}),$$
(3.98)

respectively<sup>6</sup>. Since these are all Shannon's information measures and thus nonnegative, we conclude that  $\mu^*$  is always nonnegative.

When  $X_1 \to X_2 \to \cdots \to X_n$  forms a Markov chain, for n = 3, there is only one nonempty atom, namely  $\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3$ , on which  $\mu^*$  always vanishes. This atom can be determined directly from the Markov constraint  $I(X_1; X_3 | X_2) =$ 0. For n = 4, the five nonempty atoms on which  $\mu^*$  always vanishes are listed in (3.96). The determination of these atoms, as we have seen, is not straightforward. We have also shown that for n = 3 and n = 4,  $\mu^*$  is always nonnegative.

We will extend this theme in Chapter 12 to finite Markov random fields with Markov chains being a special case. For a Markov chain, the information diagram can always be displayed in two dimensions as in Figure 3.10, and  $\mu^*$ is always nonnegative. These will be explained in Chapter 12.



**Fig. 3.10.** The information diagram for the Markov chain  $X_1 \to X_2 \to \cdots \to X_n$ .

 $<sup>^{6}</sup>$  A formal proof will be given in Theorem 12.30.

### 3.6 Examples of Applications

In this section, we give a few examples of applications of information diagrams. These examples show how information diagrams can help solving information theory problems.

The use of an information diagram is highly intuitive. To obtain an information identity from an information diagram is WYSIWYG<sup>7</sup>. However, how to obtain an information inequality from an information diagram needs some explanation.

Very often, we use a Venn diagram to represent a measure  $\mu$  which takes nonnegative values. If we see in the Venn diagram two sets A and B such that A is a subset of B, then we can immediately conclude that  $\mu(A) \leq \mu(B)$ because

$$\mu(B) - \mu(A) = \mu(B - A) \ge 0. \tag{3.99}$$

However, an *I*-Measure  $\mu^*$  can take negative values. Therefore, when we see in an information diagram that *A* is a subset of *B*, we cannot conclude from this fact alone that  $\mu^*(A) \leq \mu^*(B)$  unless we know from the setup of the problem that  $\mu^*$  is nonnegative. (For example,  $\mu^*$  is nonnegative if the random variables involved form a Markov chain.) Instead, information inequalities can be obtained from an information diagram in conjunction with the basic inequalities. The following examples will illustrate how it works.

Example 3.12 (Concavity of Entropy). Let  $X_1 \sim p_1(x)$  and  $X_2 \sim p_2(x)$ . Let

$$X \sim p(x) = \lambda p_1(x) + \bar{\lambda} p_2(x), \qquad (3.100)$$

where  $0 \leq \lambda \leq 1$  and  $\overline{\lambda} = 1 - \lambda$ . We will show that

$$H(X) \ge \lambda H(X_1) + \bar{\lambda} H(X_2). \tag{3.101}$$

Consider the system in Figure 3.11 in which the position of the switch is determined by a random variable Z with

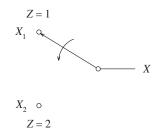


Fig. 3.11. The schematic diagram for Example 3.12.

<sup>&</sup>lt;sup>7</sup> What you see is what you get.

3.6 Examples of Applications 67

$$\Pr\{Z=1\} = \lambda \quad \text{and} \quad \Pr\{Z=2\} = \overline{\lambda}, \tag{3.102}$$

where Z is independent of  $X_1$  and  $X_2$ . The switch takes position *i* if Z = i, i = 1, 2. The random variable Z is called a *mixing random variable* for the probability distributions  $p_1(x)$  and  $p_2(x)$ . Figure 3.12 shows the information diagram for X and Z. From the diagram, we see that  $\tilde{X} - \tilde{Z}$  is a subset of  $\tilde{X}$ . Since  $\mu^*$  is nonnegative for two random variables, we can conclude that

$$\mu^*(\tilde{X}) \ge \mu^*(\tilde{X} - \tilde{Z}),$$
 (3.103)

which is equivalent to

$$H(X) \ge H(X|Z). \tag{3.104}$$

Then

$$H(X) \ge H(X|Z) \tag{3.105}$$

$$= \Pr\{Z=1\}H(X|Z=1) + \Pr\{Z=2\}H(X|Z=2)$$
(3.106)

$$=\lambda H(X_1) + \lambda H(X_2), \qquad (3.107)$$

proving (3.101). This shows that H(X) is a concave functional of p(x).

Example 3.13 (Convexity of Mutual Information). Let

$$(X, Y) \sim p(x, y) = p(x)p(y|x).$$
 (3.108)

We will show that for fixed p(x), I(X;Y) is a convex functional of p(y|x).

Let  $p_1(y|x)$  and  $p_2(y|x)$  be two transition matrices. Consider the system in Figure 3.13 in which the position of the switch is determined by a random variable Z as in the last example, where Z is independent of X, i.e.,

$$I(X;Z) = 0. (3.109)$$

In the information diagram for X, Y, and Z in Figure 3.14, let

$$I(X;Z|Y) = a \ge 0. \tag{3.110}$$

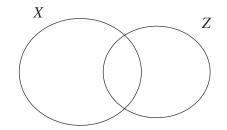


Fig. 3.12. The information diagram for Example 3.12.

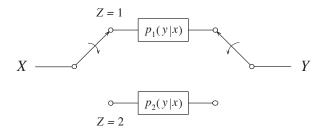


Fig. 3.13. The schematic diagram for Example 3.13.

Since I(X; Z) = 0, we see that

$$I(X;Y;Z) = -a,$$
 (3.111)

because

$$I(X;Z) = I(X;Z|Y) + I(X;Y;Z).$$
(3.112)

Then

$$\begin{split} &I(X;Y) \\ &= I(X;Y|Z) + I(X;Y;Z) \\ &= I(X;Y|Z) - a \\ &\leq I(X;Y|Z) \\ &= \Pr\{Z=1\}I(X;Y|Z=1) + \Pr\{Z=2\}I(X;Y|Z=2) \\ &= \lambda I(p(x),p_1(y|x)) + \bar{\lambda}I(p(x),p_2(y|x)), \end{split}$$
(3.117)

where  $I(p(x), p_i(y|x))$  denotes the mutual information between the input and output of a channel with input distribution p(x) and transition matrix  $p_i(y|x)$ . This shows that for fixed p(x), I(X;Y) is a convex functional of p(y|x).

Example 3.14 (Concavity of Mutual Information). Let

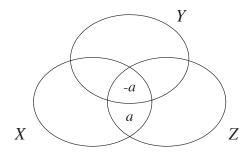


Fig. 3.14. The information diagram for Example 3.13.

3.6 Examples of Applications 69

$$(X, Y) \sim p(x, y) = p(x)p(y|x).$$
 (3.118)

We will show that for fixed p(y|x), I(X;Y) is a concave functional of p(x).

Consider the system in Figure 3.15, where the position of the switch is determined by a random variable Z as in the last example. In this system, when X is given, Y is independent of Z, or  $Z \to X \to Y$  forms a Markov chain. Then  $\mu^*$  is nonnegative, and the information diagram for X, Y, and Z is shown in Figure 3.16.

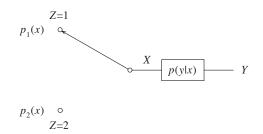


Fig. 3.15. The schematic diagram for Example 3.14.

From Figure 3.16, since  $\tilde{X} \cap \tilde{Y} - \tilde{Z}$  is a subset of  $\tilde{X} \cap \tilde{Y}$  and  $\mu^*$  is nonnegative, we immediately see that

$$I(X;Y) \geq I(X;Y|Z)$$
(3.119)  
=  $\Pr\{Z=1\}I(X;Y|Z=1) + \Pr\{Z=2\}I(X;Y|Z=2)$ (3.120)  
=  $\lambda I(p_1(x), p(y|x)) + \bar{\lambda}I(p_2(x), p(y|x)).$ (3.121)

This shows that for fixed p(y|x), I(X;Y) is a concave functional of p(x).

Example 3.15 (Imperfect Secrecy Theorem)). Let X be the plain text, Y be the cipher text, and Z be the key in a secret key cryptosystem. Since X can be recovered from Y and Z, we have

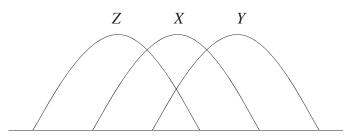


Fig. 3.16. The information diagram for Example 3.14.

$$H(X|Y,Z) = 0. (3.122)$$

We will show that this constraint implies

$$I(X;Y) \ge H(X) - H(Z).$$
 (3.123)

The quantity I(X; Y) is a measure of the security level of the cryptosystem. In general, we want to make I(X; Y) small so that the eavesdropper cannot obtain too much information about the plain text X by observing the cipher text Y. The inequality in (3.123) says that the system can attain a certain level of security only if H(Z) (often called the key length) is sufficiently large. In particular, if perfect secrecy is required, i.e., I(X;Y) = 0, then H(Z) must be at least equal to H(X). This special case is known as Shannon's perfect secrecy theorem [292]<sup>8</sup>.

We now prove (3.123). Let

$$I(X;Y|Z) = a \ge 0$$
 (3.124)

$$I(Y;Z|X) = b \ge 0 \tag{3.125}$$

$$H(Z|X,Y) = c \ge 0,$$
 (3.126)

and

$$I(X;Y;Z) = d. (3.127)$$

(See Figure 3.17.) Since  $I(Y; Z) \ge 0$ ,

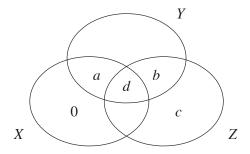


Fig. 3.17. The information diagram for Example 3.15.

$$b+d \ge 0. \tag{3.128}$$

In comparing H(X) with H(Z), we do not have to consider I(X;Z|Y) and I(X;Y;Z) since they belong to both H(X) and H(Z). Then we see from Figure 3.17 that

<sup>&</sup>lt;sup>8</sup> Shannon used a combinatorial argument to prove this theorem. An informationtheoretic proof can be found in Massey [225].

3.6 Examples of Applications 71

$$H(X) - H(Z) = a - b - c.$$
(3.129)

Therefore,

$$I(X;Y) = a + d$$
 (3.130)

$$\geq a - b \tag{3.131}$$

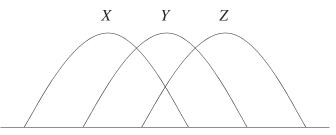
$$\geq a - b - c \tag{3.132}$$

$$=H(X) - H(Z),$$
 (3.133)

where (3.131) and (3.132) follow from (3.128) and (3.126), respectively, proving (3.123).

Note that in deriving our result, the assumptions that H(Y|X, Z) = 0, i.e., the cipher text is a function of the plain text and the key, and I(X; Z) = 0, i.e., the plain text and the key are independent, are not necessary.

*Example 3.16.* Figure 3.18 shows the information diagram for the Markov chain  $X \to Y \to Z$ . From this diagram, we can identify the following two



**Fig. 3.18.** The information diagram for the Markov chain  $X \to Y \to Z$ .

information identities:

$$I(X;Y) = I(X;Y,Z)$$
 (3.134)

$$H(X|Y) = H(X|Y,Z).$$
 (3.135)

Since  $\mu^*$  is nonnegative and  $\tilde{X} \cap \tilde{Z}$  is a subset of  $\tilde{X} \cap \tilde{Y}$ , we have

$$I(X;Z) \le I(X;Y), \tag{3.136}$$

which has already been obtained in Lemma 2.41. Similarly, we can also obtain

$$H(X|Y) \le H(X|Z). \tag{3.137}$$

Example 3.17 (Data Processing Theorem). Figure 3.19 shows the information diagram for the Markov chain  $X \to Y \to Z \to T$ . Since  $\mu^*$  is nonnegative and  $\tilde{X} \cap \tilde{T}$  is a subset of  $\tilde{Y} \cap \tilde{Z}$ , we have

$$I(X;T) \le I(Y;Z),$$
 (3.138)

which is the data processing theorem (Theorem 2.42).

We end this chapter by giving an application of the information diagram for a Markov chain with five random variables.

*Example 3.18.* In this example, we prove with the help of an information diagram that for five random variables X, Y, Z, T, and U such that  $X \to Y \to Z \to T \to U$  forms a Markov chain,

$$H(Y) + H(T) =$$
  
$$I(Z; X, Y, T, U) + I(X, Y; T, U) + H(Y|Z) + H(T|Z).$$
(3.139)

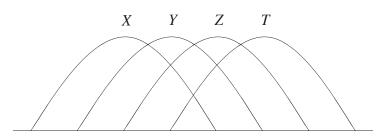
In the information diagram for X, Y, Z, T, and U in Figure 3.20, we first identify the atoms of H(Y) and then the atoms of H(T) by marking each of them by a dot. If an atom belongs to both H(Y) and H(T), it receives two dots. The resulting diagram represents

$$H(Y) + H(T).$$
 (3.140)

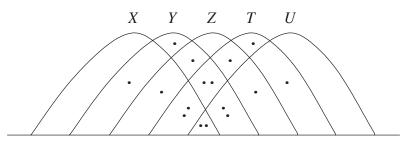
By repeating the same procedure for

$$I(Z; X, Y, T, U) + I(X, Y; T, U) + H(Y|Z) + H(T|Z),$$
(3.141)

we obtain the information diagram in Figure 3.21. Comparing these two information diagrams, we find that they are identical. Hence, the information identity in (3.139) always holds conditioning on the Markov chain

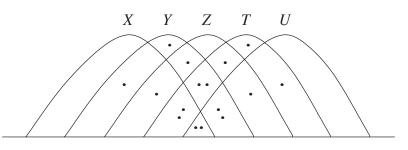


**Fig. 3.19.** The information diagram for the Markov chain  $X \to Y \to Z \to T$ .



**Fig. 3.20.** The atoms of H(Y) + H(T).

 $X \to Y \to Z \to T \to U$ . This identity is critical in proving an outer bound on the achievable coding rate region of the multiple descriptions problem in Fu *et al.* [113]. It is virtually impossible to discover this identity without the help of an information diagram!



**Fig. 3.21.** The atoms of I(Z; X, Y, T, U) + I(X, Y; T, U) + H(Y|Z) + H(T|Z).

## Appendix 3.A: A Variation of the Inclusion-Exclusion Formula

In this appendix, we show that for each  $A \in \mathcal{A}$ ,  $\mu(A)$  can be expressed as a linear combination of  $\mu(B), B \in \mathcal{B}$  via applications of (3.28) and (3.29). We first prove by using (3.28) the following variation of the inclusive-exclusive formula.

**Theorem 3.19.** For a set-additive function  $\mu$ ,

$$\mu\left(\bigcap_{k=1}^{n} A_{k} - B\right) = \sum_{1 \le i \le n} \mu(A_{i} - B) - \sum_{1 \le i < j \le n} \mu(A_{i} \cup A_{j} - B) + \dots + (-1)^{n+1} \mu(A_{1} \cup A_{2} \cup \dots \cup A_{n} - B).$$
(3.142)

*Proof.* The theorem will be proved by induction on n. First, (3.142) is obviously true for n = 1. Assume (3.142) is true for some  $n \ge 1$ . Now consider

$$\mu\left(\bigcap_{k=1}^{n+1} A_k - B\right)$$

$$= \mu\left(\left(\bigcap_{k=1}^n A_k\right) \cap A_{n+1} - B\right)$$

$$= \mu\left(\bigcap_{k=1}^n A_k - B\right) + \mu(A_{n+1} - B) - \mu\left(\left(\bigcap_{k=1}^n A_k\right) \cup A_{n+1} - B\right) \quad (3.144)$$

$$= \left\{ \sum_{1 \le i \le n} \mu(A_i - B) - \sum_{1 \le i < j \le n} \mu(A_i \cup A_j - B) + \cdots + (-1)^{n+1} \mu(A_1 \cup A_2 \cup \cdots \cup A_n - B) \right\} + \mu(A_{n+1} - B)$$

$$-\mu \left( \bigcap_{k=1}^n (A_k \cup A_{n+1}) - B \right)$$

$$= \left\{ \sum_{1 \le i \le n} \mu(A_i - B) - \sum_{1 \le i < j \le n} \mu(A_i \cup A_j - B) + \cdots + (-1)^{n+1} \mu(A_1 \cup A_2 \cup \cdots \cup A_n - B) \right\} + \mu(A_{n+1} - B)$$

$$- \left\{ \sum_{1 \le i \le n} \mu(A_i \cup A_{n+1} - B) - \sum_{1 \le i < j \le n} \mu(A_i \cup A_j \cup A_{n+1} - B) + \cdots + (-1)^{n+1} \mu(A_1 \cup A_2 \cup \cdots \cup A_n \cup A_n \cup A_{n+1} - B) \right\}$$

$$= \sum_{1 \le i \le n+1} \mu(A_i - B) - \sum_{1 \le i < j \le n+1} \mu(A_i \cup A_j - B) + \cdots + (-1)^{n+2} \mu(A_1 \cup A_2 \cup \cdots \cup A_{n+1} - B).$$

$$(3.147)$$

In the above, (3.28) was used in obtaining (3.144), and the induction hypothesis was used in obtaining (3.145) and (3.146). The theorem is proved.  $\Box$ 

Now a nonempty atom of  $\mathcal{F}_n$  has the form

$$\bigcap_{i=1}^{n} Y_i, \tag{3.148}$$

where  $Y_i$  is either  $\tilde{X}_i$  or  $\tilde{X}_i^c$ , and there exists at least one *i* such that  $Y_i = \tilde{X}_i$ . Then we can write the atom in (3.148) as

$$\bigcap_{i:Y_i=\tilde{X}_i} \tilde{X}_i - \left(\bigcup_{j:Y_j=\tilde{X}_j^c} \tilde{X}_j\right).$$
(3.149)

Note that the intersection above is always nonempty. Then using (3.142) and (3.29), we see that for each  $A \in \mathcal{A}$ ,  $\mu(A)$  can be expressed as a linear combination of  $\mu(B), B \in \mathcal{B}$ .

### Problems

1. Show that

$$I(X;Y;Z) = E \log \frac{p(X,Y)p(Y,Z)p(X,Z)}{p(X)p(Y)p(Z)p(X,Y,Z)}$$

and obtain a general formula for  $I(X_1; X_2; \dots; X_n)$ .

- 2. Suppose  $X \perp Y$  and  $X \perp Z$ . Does  $X \perp (Y, Z)$  hold in general?
- 3. Show that I(X;Y;Z) vanishes if at least one of the following conditions hold:
  - a) X, Y, and Z are mutually independent;
  - b)  $X \to Y \to Z$  forms a Markov chain and X and Z are independent.
- 4. a) Verify that I(X;Y;Z) vanishes for the distribution p(x,y,z) given by

$$p(0,0,0) = 0.0625, p(0,0,1) = 0.0772, p(0,1,0) = 0.0625$$
  
 $p(0,1,1) = 0.0625, p(1,0,0) = 0.0625, p(1,0,1) = 0.1103$   
 $p(1,1,0) = 0.1875, p(1,1,1) = 0.375.$ 

- b) Verify that the distribution in part a) does not satisfy the conditions in Problem 3.
- 5. Weak independence X is weakly independent of Y if the rows of the transition matrix [p(x|y)] are linearly dependent.
  - a) Show that if X and Y are independent, then X is weakly independent of Y.
  - b) Show that for random variables X and Y, there exists a random variable Z satisfying
    - i)  $X \to Y \to Z$
    - ii) X and Z are independent
    - iii) Y and Z are not independent
    - if and only if X is weakly independent of Y.

(Berger and Yeung [28].)

- 6. Prove that
  - a)  $I(X;Y;Z) \ge -\min\{I(X;Y|Z), I(Y;Z|X), I(X,Z|Y)\}$
  - b)  $I(X;Y;Z) \le \min\{I(X;Y), I(Y;Z), I(X;Z)\}.$
- 7. a) Prove that if X and Y are independent, then  $I(X,Y;Z) \ge I(X;Y|Z)$ .
  - b) Show that the inequality in part a) is not valid in general by giving a counterexample.
- 8. In Example 3.15, it was shown that  $I(X;Y) \ge H(X) H(Z)$ , where X is the plain text, Y is the cipher text, and Z is the key in a secret key cryptosystem. Give an example of a secret key cryptosystem such that this inequality is tight.
- 9. Secret sharing For a given finite set  $\mathcal{P}$  and a collection  $\mathcal{A}$  of subsets of  $\mathcal{P}$ , a secret sharing scheme is a random variable S and a family of random variables  $\{X_p : p \in \mathcal{P}\}$  such that for all  $A \in \mathcal{A}$ ,

$$H(S|X_A) = 0,$$

and for all  $B \notin \mathcal{A}$ ,

$$H(S|X_B) = H(S)$$

Here, S is the secret and  $\mathcal{P}$  is the set of participants of the scheme. A participant p of the scheme possesses a share  $X_p$  of the secret. The set  $\mathcal{A}$  specifies the access structure of the scheme: For a subset A of  $\mathcal{P}$ , by pooling their shares, if  $A \in \mathcal{A}$ , the participants in A can reconstruct S, otherwise they can know nothing about S.

a) i) Prove that for  $A, B \subset \mathcal{P}$ , if  $B \notin \mathcal{A}$  and  $A \cup B \in \mathcal{A}$ , then

$$H(X_A|X_B) = H(S) + H(X_A|X_B, S).$$

ii) Prove that if  $B \in \mathcal{A}$ , then

$$H(X_A|X_B) = H(X_A|X_B, S).$$

(Capocelli *et al.* [54].)

b) Prove that for  $A, B, C \subset \mathcal{P}$  such that  $A \cup C \in \mathcal{A}$ ,  $B \cup C \in \mathcal{A}$ , and  $C \notin \mathcal{A}$ , then

$$I(X_A; X_B | X_C) \ge H(S).$$

(van Dijk [328].)

- 10. Consider four random variables X, Y, Z, and T which satisfy the following constraints: H(T|X) = H(T), H(T|X, Y) = 0, H(T|Y) = H(T), H(Y|Z) = 0, and H(T|Z) = 0. Prove that
  - a) H(T|X, Y, Z) = I(Z; T|X, Y) = 0.
  - b) I(X;T|Y,Z) = I(X;Y;T|Z) = I(Y;T|X,Z) = 0.
  - c) I(X; Z; T) = I(Y; Z; T) = 0.
  - d) H(Y|X, Z, T) = I(X; Y|Z, T) = 0.
  - e)  $I(X;Y;Z) \ge 0.$
  - f)  $I(X;Z) \ge H(T)$ .

The inequality in f) finds application in a secret sharing problem studied by Blundo *et al.* [41].

In the following, we use  $X \perp Y | Z$  to denote that X and Y are independent given Z.

- 11. a) Prove that under the constraint that  $X \to Y \to Z$  forms a Markov chain,  $X \perp Y | Z$  and  $X \perp Z$  imply  $X \perp Y$ .
  - b) Prove that the implication in a) continues to be valid without the Markov chain constraint.
- 12. a) Show that  $Y \perp Z|T$  does not imply  $Y \perp Z|(X,T)$  by giving a counterexample.
  - b) Prove that  $Y \perp Z | T$  implies  $Y \perp Z | (X, T)$  conditioning on  $X \to Y \to Z \to T$ .

13. Prove that for random variables X, Y, Z, and T,

$$\left. \begin{array}{c} X \perp Z | Y \\ (X,Y) \perp T | Z \\ Y \perp Z | T \\ Y \perp Z | X \\ X \perp T \end{array} \right\} \Rightarrow Y \perp Z.$$

Hint: Observe that  $X \perp Z | Y$  and  $(X, Y) \perp T | Z$  are equivalent to  $X \rightarrow Y \rightarrow Z \rightarrow T$  and use an information diagram.

14. Prove that

$$\begin{cases} X \perp Y \\ X \perp Y | (Z, T) \\ Z \perp T | X \\ Z \perp T | Y \end{cases} \Leftrightarrow \begin{cases} Z \perp T \\ Z \perp T | (X, Y) \\ X \perp Y | Z \\ X \perp Y | T. \end{cases}$$

(Studený [313].)

### **Historical Notes**

The original work on the set-theoretic structure of Shannon's information measures is due to Hu [155]. It was established in this paper that every information identity implies a set identity via a substitution of symbols. This allows the tools for proving information identities to be used in proving set identities. Since the paper was published in Russian, it was largely unknown to the West until it was described in Csiszár and Körner [75]. Throughout the years, the use of Venn diagrams to represent the structure of Shannon's information measures for two or three random variables has been suggested by various authors, for example, Reza [273], Abramson [2], and Papoulis [257], but no formal justification was given until Yeung [357] introduced the *I*-Measure.

McGill [239] proposed a multiple mutual information for any number of random variables which is equivalent to the mutual information between two or more random variables discussed here. Properties of this quantity have been investigated by Kawabata [176] and Yeung [357].

Along a related direction, Han [133] viewed the linear combination of entropies as a vector space and developed a lattice-theoretic description of Shannon's information measures.

In a random experiment, a coin is tossed n times. Let  $X_i$  be the outcome of the *i*th toss, with

$$\Pr\{X_i = \text{HEAD}\} = p \text{ and } \Pr\{X_i = \text{TAIL}\} = 1 - p, \qquad (4.1)$$

where  $0 \le p \le 1$ . It is assumed that  $X_i$  are i.i.d., and the value of p is known. We are asked to describe the outcome of the random experiment without error (with zero error) by using binary symbols. One way to do this is to encode a HEAD by a '0' and a TAIL by a '1.' Then the outcome of the random experiment is encoded into a binary codeword of length n. When the coin is fair, i.e., p = 0.5, this is the best we can do because the probability of every outcome of the experiment is equal to  $2^{-n}$ . In other words, all the outcomes are equally likely.

However, if the coin is biased, i.e.,  $p \neq 0.5$ , the probability of an outcome of the experiment depends on the number of HEADs and the number of TAILs in the outcome. In other words, the probabilities of the outcomes are no longer uniform. It turns out that we can take advantage of this by encoding more likely outcomes into shorter codewords and less likely outcomes into longer codewords. By doing so, it is possible to use fewer than n bits on the average to describe the outcome of the random experiment. In particular, in the extreme case in which p = 0 or 1, we actually do not need to describe the outcome of the experiment because it is deterministic.

At the beginning of Chapter 2, we mentioned that the entropy H(X) measures the amount of information contained in a random variable X. In this chapter, we substantiate this claim by exploring the role of entropy in the context of zero-error data compression.

### 4.1 The Entropy Bound

In this section, we establish that H(X) is a fundamental lower bound on the expected length of the number of symbols needed to describe the outcome of a random variable X with zero error. This is called the *entropy bound*.

**Definition 4.1.** A D-ary source code C for a source random variable X is a mapping from  $\mathcal{X}$  to  $\mathcal{D}^*$ , the set of all finite length sequences of symbols taken from a D-ary code alphabet.

Consider an information source  $\{X_k, k \geq 1\}$ , where  $X_k$  are discrete random variables which take values in the same alphabet. We apply a source code C to each  $X_k$  and concatenate the codewords. Once the codewords are concatenated, the boundaries of the codewords are no longer explicit. In other words, when the code C is applied to a source sequence, a sequence of code symbols are produced, and the codewords may no longer be distinguishable. We are particularly interested in *uniquely decodable codes* which are defined as follows.

**Definition 4.2.** A code C is uniquely decodable if for any finite source sequence, the sequence of code symbols corresponding to this source sequence is different from the sequence of code symbols corresponding to any other (finite) source sequence.

Suppose we use a code C to encode a source file into a coded file. If C is uniquely decodable, then we can always recover the source file from the coded file. An important class of uniquely decodable codes, called *prefix codes*, are discussed in the next section. But we first look at an example of a code which is not uniquely decodable.

*Example 4.3.* Let  $\mathcal{X} = \{A, B, C, D\}$ . Consider the code  $\mathcal{C}$  defined by

x	$\mathcal{C}(x)$
Α	0
В	1
$\mathbf{C}$	01
D	10

Then all the three source sequences AAD, ACA, and AABA produce the code sequence 0010. Thus from the code sequence 0010, we cannot tell which of the three source sequences it comes from. Therefore, C is not uniquely decodable.

In the next theorem, we prove that for any uniquely decodable code, the lengths of the codewords have to satisfy an inequality called the *Kraft inequality*.

#### 4.1 The Entropy Bound 81

**Theorem 4.4 (Kraft Inequality).** Let C be a D-ary source code, and let  $l_1, l_2, \dots, l_m$  be the lengths of the codewords. If C is uniquely decodable, then

$$\sum_{k=1}^{m} D^{-l_k} \le 1. \tag{4.2}$$

*Proof.* Let N be an arbitrary positive integer, and consider the identity

$$\left(\sum_{k=1}^{m} D^{-l_k}\right)^N = \sum_{k_1=1}^{m} \sum_{k_2=1}^{m} \cdots \sum_{k_N=1}^{m} D^{-(l_{k_1}+l_{k_2}+\dots+l_{k_N})}.$$
 (4.3)

By collecting terms on the right-hand side, we write

$$\left(\sum_{k=1}^{m} D^{-l_k}\right)^N = \sum_{i=1}^{Nl_{\max}} A_i D^{-i}$$
(4.4)

where

$$l_{\max} = \max_{1 \le k \le m} l_k \tag{4.5}$$

and  $A_i$  is the coefficient of  $D^{-i}$  in  $\left(\sum_{k=1}^m D^{-l_k}\right)^N$ . Now observe that  $A_i$  gives the total number of sequences of N codewords with a total length of i code symbols. Since the code is uniquely decodable, these code sequences must be distinct, and therefore  $A_i \leq D^i$  because there are  $D^i$  distinct sequences of icode symbols. Substituting this inequality into (4.4), we have

$$\left(\sum_{k=1}^{m} D^{-l_k}\right)^N \le \sum_{i=1}^{Nl_{\max}} 1 = Nl_{\max},\tag{4.6}$$

or

$$\sum_{k=1}^{m} D^{-l_k} \le (N l_{\max})^{1/N}.$$
(4.7)

Since this inequality holds for any N, upon letting  $N \to \infty$ , we obtain (4.2), completing the proof.  $\Box$ 

Let X be a source random variable with probability distribution

$$\{p_1, p_2, \cdots, p_m\},$$
 (4.8)

where  $m \geq 2$ . When we use a uniquely decodable code C to encode the outcome of X, we are naturally interested in the expected length of a codeword, which is given by

$$L = \sum_{i} p_i l_i. \tag{4.9}$$

We will also refer to L as the expected length of the code C. The quantity L gives the average number of symbols we need to describe the outcome of X when the code C is used, and it is a measure of the efficiency of the code C. Specifically, the smaller the expected length L is, the better the code C is.

In the next theorem, we will prove a fundamental lower bound on the expected length of any uniquely decodable *D*-ary code. We first explain why this is the lower bound we should expect. In a uniquely decodable code, we use L *D*-ary symbols on the average to describe the outcome of X. Recall from the remark following Theorem 2.43 that a *D*-ary symbol can carry at most one *D*-it of information. Then the maximum amount of information which can be carried by the codeword on the average is  $L \cdot 1 = L$  *D*-its. Since the code is uniquely decodable, the amount of entropy carried by the codeword on the average is H(X). Therefore, we have

$$H_D(X) \le L. \tag{4.10}$$

In other words, the expected length of a uniquely decodable code is at least the entropy of the source. This argument is rigorized in the proof of the next theorem.

**Theorem 4.5 (Entropy Bound).** Let C be a D-ary uniquely decodable code for a source random variable X with entropy  $H_D(X)$ . Then the expected length of C is lower bounded by  $H_D(X)$ , i.e.,

$$L \ge H_D(X). \tag{4.11}$$

This lower bound is tight if and only if  $l_i = -\log_D p_i$  for all *i*.

*Proof.* Since C is uniquely decodable, the lengths of its codewords satisfy the Kraft inequality. Write

$$L = \sum_{i} p_i \log_D D^{l_i} \tag{4.12}$$

and recall from Definition 2.35 that

$$H_D(X) = -\sum_i p_i \log_D p_i.$$
(4.13)

Then

$$L - H_D(X) = \sum_i p_i \log_D(p_i D^{l_i})$$
(4.14)

$$= (\ln D)^{-1} \sum_{i} p_i \ln(p_i D^{l_i})$$
(4.15)

$$\geq (\ln D)^{-1} \sum_{i} p_i \left( 1 - \frac{1}{p_i D^{l_i}} \right)$$
(4.16)

$$= (\ln D)^{-1} \left[ \sum_{i} p_{i} - \sum_{i} D^{-l_{i}} \right]$$
(4.17)

4.2 Prefix Codes 83

$$\geq (\ln D)^{-1}(1-1) \tag{4.18}$$

$$= 0,$$
 (4.19)

where we have invoked the fundamental inequality in (4.16) and the Kraft inequality in (4.18). This proves (4.11). In order for this lower bound to be tight, both (4.16) and (4.18) have to be tight simultaneously. Now (4.16) is tight if and only if  $p_i D^{l_i} = 1$ , or  $l_i = -\log_D p_i$  for all *i*. If this holds, we have

$$\sum_{i} D^{-l_i} = \sum_{i} p_i = 1, \tag{4.20}$$

i.e., (4.18) is also tight. This completes the proof of the theorem.  $\Box$ 

The entropy bound can be regarded as a generalization of Theorem 2.43, as is seen from the following corollary.

### Corollary 4.6. $H(X) \leq \log |\mathcal{X}|$ .

*Proof.* Considering encoding each outcome of a random variable X by a distinct symbol in  $\{1, 2, \dots, |\mathcal{X}|\}$ . This is obviously a  $|\mathcal{X}|$ -ary uniquely decodable code with expected length 1. Then by the entropy bound, we have

$$H_{|\mathcal{X}|}(X) \le 1, \tag{4.21}$$

which becomes

$$H(X) \le \log |\mathcal{X}| \tag{4.22}$$

when the base of the logarithm is not specified.  $\Box$ 

Motivated by the entropy bound, we now introduce the *redundancy* of a uniquely decodable code.

**Definition 4.7.** The redundancy R of a D-ary uniquely decodable code is the difference between the expected length of the code and the entropy of the source.

We see from the entropy bound that the redundancy of a uniquely decodable code is always nonnegative.

### 4.2 Prefix Codes

### 4.2.1 Definition and Existence

**Definition 4.8.** A code is called a prefix-free code if no codeword is a prefix of any other codeword. For brevity, a prefix-free code will be referred to as a prefix code.

*Example 4.9.* The code C in Example 4.3 is not a prefix code because the codeword 0 is a prefix of the codeword 01, and the codeword 1 is a prefix of the codeword 10. It can easily be checked that the following code C' is a prefix code.

$$\begin{array}{c|c} \mathbf{x} & \mathcal{C}'(x) \\ \hline \mathbf{A} & \mathbf{0} \\ \mathbf{B} & \mathbf{10} \\ \mathbf{C} & \mathbf{110} \\ \mathbf{D} & \mathbf{1111} \end{array}$$

A D-ary tree is a graphical representation of a collection of finite sequences of D-ary symbols. In a D-ary tree, each node has at most D children. If a node has at least one child, it is called an *internal node*, otherwise it is called a *leaf*. The children of an internal node are labeled by the D symbols in the code alphabet.

A *D*-ary prefix code can be represented by a *D*-ary tree with the leaves of the tree being the codewords. Such a tree is called the *code tree* for the prefix code. Figure 4.1 shows the code tree for the prefix code C' in Example 4.9.

As we have mentioned in Section 4.1, once a sequence of codewords are concatenated, the boundaries of the codewords are no longer explicit. Prefix codes have the desirable property that the end of a codeword can be recognized instantaneously so that it is not necessary to make reference to the future codewords during the decoding process. For example, for the source sequence  $BCDAC \cdots$ , the code C' in Example 4.9 produces the code sequence  $1011011110110 \cdots$ . Based on this binary sequence, the decoder can reconstruct the source sequence as follows. The first bit 1 cannot form the first codeword because 1 is not a valid codeword. The first two bits 10 must form the first codeword because it is a valid codeword and it is not the prefix of any other codeword. The same procedure is repeated to locate the end of the next codeword, and the code sequence is parsed as  $10, 110, 1111, 0, 110, \cdots$ . Then the source sequence  $BCDAC \cdots$  can be reconstructed correctly.

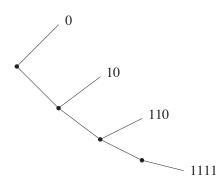


Fig. 4.1. The code tree for the code C'.

Since a prefix code can always be decoded correctly, it is a uniquely decodable code. Therefore, by Theorem 4.4, the codeword lengths of a prefix code also satisfies the Kraft inequality. In the next theorem, we show that the Kraft inequality fully characterizes the existence of a prefix code.

**Theorem 4.10.** There exists a D-ary prefix code with codeword lengths  $l_1$ ,  $l_2, \dots, l_m$  if and only if the Kraft inequality

$$\sum_{k=1}^{m} D^{-l_k} \le 1 \tag{4.23}$$

is satisfied.

*Proof.* We only need to prove the existence of a *D*-ary prefix code with codeword lengths  $l_1, l_2, \dots, l_m$  if these lengths satisfy the Kraft inequality. Without loss of generality, assume that  $l_1 \leq l_2 \leq \dots \leq l_m$ .

Consider all the *D*-ary sequences of lengths less than or equal to  $l_m$  and regard them as the nodes of the full *D*-ary tree of depth  $l_m$ . We will refer to a sequence of length l as a node of order l. Our strategy is to choose nodes as codewords in nondecreasing order of the codeword lengths. Specifically, we choose a node of order  $l_1$  as the first codeword, then a node of order  $l_2$  as the second codeword, so on and so forth, such that each newly chosen codeword is not prefixed by any of the previously chosen codewords. If we can successfully choose all the *m* codewords, then the resultant set of codewords forms a prefix code with the desired set of lengths.

There are  $D^{l_1} > 1$  (since  $l_1 \ge 1$ ) nodes of order  $l_1$  which can be chosen as the first codeword. Thus choosing the first codeword is always possible. Assume that the first *i* codewords have been chosen successfully, where  $1 \le i \le m-1$ , and we want to choose a node of order  $l_{i+1}$  as the (i+1)st codeword such that it is not prefixed by any of the previously chosen codewords. In other words, the (i + 1)st node to be chosen cannot be a descendant of any of the previously chosen codewords. Observe that for  $1 \le j \le i$ , the codeword with length  $l_j$  has  $D^{l_{i+1}-l_j}$  descendents of order  $l_{i+1}$ . Since all the previously chosen codewords are not prefeces of each other, their descendents of order  $l_{i+1}$  do not overlap. Therefore, upon noting that the total number of nodes of order  $l_{i+1}$ is  $D^{l_{i+1}}$ , the number of nodes which can be chosen as the (i + 1)st codeword is

$$D^{l_{i+1}} - D^{l_{i+1}-l_1} - \dots - D^{l_{i+1}-l_i}.$$
(4.24)

If  $l_1, l_2, \dots, l_m$  satisfy the Kraft inequality, we have

$$D^{-l_1} + \dots + D^{-l_i} + D^{-l_{i+1}} \le 1.$$
(4.25)

Multiplying by  $D^{l_{i+1}}$  and rearranging the terms, we have

$$D^{l_{i+1}} - D^{l_{i+1}-l_1} - \dots - D^{l_{i+1}-l_i} \ge 1.$$
(4.26)

The left hand side is the number of nodes which can be chosen as the (i+1)st codeword as given in (4.24). Therefore, it is possible to choose the (i+1)st codeword. Thus we have shown the existence of a prefix code with codeword lengths  $l_1, l_2, \dots, l_m$ , completing the proof.  $\Box$ 

A probability distribution  $\{p_i\}$  such that for all i,  $p_i = D^{-t_i}$ , where  $t_i$  is a positive integer, is called a *D*-adic distribution. When D = 2,  $\{p_i\}$  is called a *dyadic* distribution. From Theorem 4.5 and the above theorem, we can obtain the following result as a corollary.

**Corollary 4.11.** There exists a D-ary prefix code which achieves the entropy bound for a distribution  $\{p_i\}$  if and only if  $\{p_i\}$  is D-adic.

*Proof.* Consider a *D*-ary prefix code which achieves the entropy bound for a distribution  $\{p_i\}$ . Let  $l_i$  be the length of the codeword assigned to the probability  $p_i$ . By Theorem 4.5, for all  $i, l_i = -\log_D p_i$ , or  $p_i = D^{-l_i}$ . Thus  $\{p_i\}$  is *D*-adic.

Conversely, suppose  $\{p_i\}$  is *D*-adic, and let  $p_i = D^{-t_i}$  for all *i*. Let  $l_i = t_i$  for all *i*. Then by the Kraft inequality, there exists a prefix code with codeword lengths  $\{l_i\}$ , because

$$\sum_{i} D^{-l_i} = \sum_{i} D^{-t_i} = \sum_{i} p_i = 1.$$
(4.27)

Assigning the codeword with length  $l_i$  to the probability  $p_i$  for all i, we see from Theorem 4.5 that this code achieves the entropy bound.  $\Box$ 

### 4.2.2 Huffman Codes

As we have mentioned, the efficiency of a uniquely decodable code is measured by its expected length. Thus for a given source X, we are naturally interested in prefix codes which have the minimum expected length. Such codes, called optimal codes, can be constructed by the *Huffman procedure*, and these codes are referred to as *Huffman codes*. In general, there exists more than one optimal code for a source, and some optimal codes cannot be constructed by the Huffman procedure.

For simplicity, we first discuss binary Huffman codes. A binary prefix code for a source X with distribution  $\{p_i\}$  is represented by a binary code tree, with each leaf in the code tree corresponding to a codeword. The Huffman procedure is to form a code tree such that the expected length is minimum. The procedure is described by a very simple rule:

Keep merging the two smallest probability masses until one probability mass (i.e., 1) is left.

The merging of two probability masses corresponds to the formation of an internal node of the code tree. We now illustrate the Huffman procedure by the following example.

*Example 4.12.* Let X be the source with  $\mathcal{X} = \{A, B, C, D, E\}$ , and the probabilities are 0.35, 0.1, 0.15, 0.2, 0.2, respectively. The Huffman procedure is shown in Figure 4.2. In the first step, we merge probability masses 0.1 and 0.15

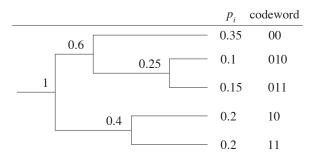


Fig. 4.2. The Huffman procedure.

into a probability mass 0.25. In the second step, we merge probability masses 0.2 and 0.2 into a probability mass 0.4. In the third step, we merge probability masses 0.35 and 0.25 into a probability mass 0.6. Finally, we merge probability masses 0.6 and 0.4 into a probability mass 1. A code tree is then formed. Upon assigning 0 and 1 (in any convenient way) to each pair of branches at an internal node, we obtain the codeword assigned to each source symbol.

In the Huffman procedure, sometimes there is more than one choice of merging the two smallest probability masses. We can take any one of these choices without affecting the optimality of the code eventually obtained.

For an alphabet of size m, it takes m - 1 steps to complete the Huffman procedure for constructing a binary code, because we merge two probability masses in each step. In the resulting code tree, there are m leaves and m - 1internal nodes.

In the Huffman procedure for constructing a D-ary code, the smallest D probability masses are merged in each step. If the resulting code tree is formed in k + 1 steps, where  $k \ge 0$ , then there will be k + 1 internal nodes and D + k(D - 1) leaves, where each leaf corresponds to a source symbol in the alphabet. If the alphabet size m has the form D + k(D - 1), then we can apply the Huffman procedure directly. Otherwise, we need to add a few dummy symbols with probability 0 to the alphabet in order to make the total number of symbols have the form D + k(D - 1).

Example 4.13. If we want to construct a quaternary Huffman code (D = 4) for the source in the last example, we need to add 2 dummy symbols so that the total number of symbols becomes 7 = 4 + (1)3, where k = 1. In general, we need to add at most D - 2 dummy symbols.

In Section 4.1, we have proved the entropy bound for a uniquely decodable code. This bound also applies to a prefix code since a prefix code is uniquely

decodable. In particular, it applies to a Huffman code, which is a prefix code by construction. Thus the expected length of a Huffman code is at least the entropy of the source. In Example 4.12, the entropy H(X) is 2.202 bits, while the expected length of the Huffman code is

$$0.35(2) + 0.1(3) + 0.15(3) + 0.2(2) + 0.2(2) = 2.25.$$

$$(4.28)$$

We now turn to proving the optimality of a Huffman code. For simplicity, we will only prove the optimality of a binary Huffman code. Extension of the proof to the general case is straightforward.

Without loss of generality, assume that

$$p_1 \ge p_2 \ge \dots \ge p_m. \tag{4.29}$$

Denote the codeword assigned to  $p_i$  by  $c_i$ , and denote its length by  $l_i$ . To prove that a Huffman code is actually optimal, we make the following observations.

**Lemma 4.14.** In an optimal code, shorter codewords are assigned to larger probabilities.

*Proof.* Consider  $1 \leq i < j \leq m$  such that  $p_i > p_j$ . Assume that in a code, the codewords  $c_i$  and  $c_j$  are such that  $l_i > l_j$ , i.e., a shorter codeword is assigned to a smaller probability. Then by exchanging  $c_i$  and  $c_j$ , the expected length of the code is changed by

$$(p_i l_j + p_j l_i) - (p_i l_i + p_j l_j) = (p_i - p_j)(l_j - l_i) < 0$$

$$(4.30)$$

since  $p_i > p_j$  and  $l_i > l_j$ . In other words, the code can be improved and therefore is not optimal. The lemma is proved.  $\Box$ 

**Lemma 4.15.** There exists an optimal code in which the codewords assigned to the two smallest probabilities are siblings, i.e., the two codewords have the same length and they differ only in the last symbol.

*Proof.* The reader is encouraged to trace the steps in this proof by drawing a code tree. Consider any optimal code. From the last lemma, the codeword  $c_m$  assigned to  $p_m$  has the longest length. Then the sibling of  $c_m$  cannot be the prefix of another codeword.

We claim that the sibling of  $c_m$  must be a codeword. To see this, assume that it is not a codeword (and it is not the prefix of another codeword). Then we can replace  $c_m$  by its parent to improve the code because the length of the codeword assigned to  $p_m$  is reduced by 1, while all the other codewords remain unchanged. This is a contradiction to the assumption that the code is optimal. Therefore, the sibling of  $c_m$  must be a codeword.

If the sibling of  $c_m$  is assigned to  $p_{m-1}$ , then the code already has the desired property, i.e., the codewords assigned to the two smallest probabilities are siblings. If not, assume that the sibling of  $c_m$  is assigned to  $p_i$ , where  $i < c_m$ 

m-1. Since  $p_i \ge p_{m-1}$ ,  $l_{m-1} \ge l_i = l_m$ . On the other hand, by Lemma 4.14,  $l_{m-1}$  is always less than or equal to  $l_m$ , which implies that  $l_{m-1} = l_m = l_i$ . Then we can exchange the codewords for  $p_i$  and  $p_{m-1}$  without changing the expected length of the code (i.e., the code remains optimal) to obtain the desired code. The lemma is proved.  $\Box$ 

Suppose  $c_i$  and  $c_j$  are siblings in a code tree. Then  $l_i = l_j$ . If we replace  $c_i$  and  $c_j$  by a common codeword at their parent, call it  $c_{ij}$ , then we obtain a reduced code tree, and the probability of  $c_{ij}$  is  $p_i + p_j$ . Accordingly, the probability set becomes a reduced probability set with  $p_i$  and  $p_j$  replaced by a probability  $p_i + p_j$ . Let L and L' be the expected lengths of the original code and the reduced code, respectively. Then

$$L - L' = (p_i l_i + p_j l_j) - (p_i + p_j)(l_i - 1)$$
(4.31)

$$= (p_i l_i + p_j l_i) - (p_i + p_j)(l_i - 1)$$
(4.32)

$$= p_i + p_j, \tag{4.33}$$

which implies

$$L = L' + (p_i + p_j). (4.34)$$

This relation says that the difference between the expected length of the original code and the expected length of the reduced code depends only on the values of the two probabilities merged but not on the structure of the reduced code tree.

### Theorem 4.16. The Huffman procedure produces an optimal prefix code.

Proof. Consider an optimal code in which  $c_m$  and  $c_{m-1}$  are siblings. Such an optimal code exists by Lemma 4.15. Let  $\{p'_i\}$  be the reduced probability set obtained from  $\{p_i\}$  by merging  $p_m$  and  $p_{m-1}$ . From (4.34), we see that L' is the expected length of an optimal code for  $\{p'_i\}$  if and only if L is the expected length of an optimal code for  $\{p'_i\}$ . Therefore, if we can find an optimal code for  $\{p'_i\}$ , we can use it to construct an optimal code for  $\{p_i\}$ . Note that by merging  $p_m$  and  $p_{m-1}$ , the size of the problem, namely the total number of probability masses, is reduced by one. To find an optimal code for  $\{p'_i\}$ , we again merge the two smallest probability in  $\{p'_i\}$ . This is repeated until the size of the problem is eventually reduced to 2, which we know that an optimal code has two codewords of length 1. In the last step of the Huffman procedure, two probability masses are merged, which corresponds to the formation of a code with two codewords of length 1. Thus the Huffman procedure indeed produces an optimal code.  $\Box$ 

We have seen that the expected length of a Huffman code is lower bounded by the entropy of the source. On the other hand, it would be desirable to obtain an upper bound in terms of the entropy of the source. This is given in the next theorem.

**Theorem 4.17.** The expected length of a Huffman code, denoted by  $L_{\text{Huff}}$ , satisfies

$$L_{\text{Huff}} < H_D(X) + 1.$$
 (4.35)

This bound is the tightest among all the upper bounds on  $L_{\text{Huff}}$  which depend only on the source entropy.

*Proof.* We will construct a prefix code with expected length less than H(X) + 1. Then, because a Huffman code is an optimal prefix code, its expected length  $L_{\text{Huff}}$  is upper bounded by H(X) + 1.

Consider constructing a prefix code with codeword lengths  $\{l_i\}$ , where

$$l_i = \lceil -\log_D p_i \rceil. \tag{4.36}$$

Then

$$-\log_D p_i \le l_i < -\log_D p_i + 1, \tag{4.37}$$

or

$$p_i \ge D^{-l_i} > D^{-1} p_i. \tag{4.38}$$

Thus

$$\sum_{i} D^{-l_i} \le \sum_{i} p_i = 1, \tag{4.39}$$

i.e.,  $\{l_i\}$  satisfies the Kraft inequality, which implies that it is possible to construct a prefix code with codeword lengths  $\{l_i\}$ .

It remains to show that L, the expected length of this code, is less than H(X) + 1. Toward this end, consider

$$L = \sum_{i} p_{i} l_{i} \tag{4.40}$$

$$<\sum_{i} p_i(-\log_D p_i + 1)$$
 (4.41)

$$= -\sum_{i} p_i \log_D p_i + \sum_{i} p_i \tag{4.42}$$

$$=H(X)+1,$$
 (4.43)

where (4.41) follows from the upper bound in (4.37). Thus we conclude that

$$L_{\text{Huff}} \le L < H(X) + 1.$$
 (4.44)

To see that this upper bound is the tightest possible, we have to show that there exists a sequence of distributions  $P_k$  such that  $L_{\text{Huff}}$  approaches H(X)+1as  $k \to \infty$ . This can be done by considering the sequence of *D*-ary distributions

$$P_{k} = \left\{ 1 - \frac{D-1}{k}, \frac{1}{k}, \cdots, \frac{1}{k} \right\},$$
(4.45)

where  $k \geq D$ . The Huffman code for each  $P_k$  consists of D codewords of length 1. Thus  $L_{\text{Huff}}$  is equal to 1 for all k. As  $k \to \infty$ ,  $H(X) \to 0$ , and hence  $L_{\text{Huff}}$  approaches H(X) + 1. The theorem is proved.  $\Box$ 

The code constructed in the above proof is known as the Shannon code. The idea is that in order for the code to be near-optimal, we should choose  $l_i$  close to  $-\log p_i$  for all *i*. When  $\{p_i\}$  is *D*-adic,  $l_i$  can be chosen to be exactly  $-\log p_i$  because the latter are integers. In this case, the entropy bound is tight.

From the entropy bound and the above theorem, we have

$$H(X) \le L_{\text{Huff}} < H(X) + 1.$$
 (4.46)

Now suppose we use a Huffman code to encode  $X_1, X_2, \dots, X_n$  which are n i.i.d. copies of X. Let us denote the length of this Huffman code by  $L^n_{\text{Huff}}$ . Then (4.46) becomes

$$nH(X) \le L_{\text{Huff}}^n < nH(X) + 1.$$
 (4.47)

Dividing by n, we obtain

$$H(X) \le \frac{1}{n} L_{\text{Huff}}^n < H(X) + \frac{1}{n}.$$
 (4.48)

As  $n \to \infty$ , the upper bound approaches the lower bound. Therefore,  $n^{-1}L_{\text{Huff}}^n$ , the coding rate of the code, namely the average number of code symbols needed to encode a source symbol, approaches H(X) as  $n \to \infty$ . But of course, as n becomes large, constructing a Huffman code becomes very complicated. Nevertheless, this result indicates that entropy is a fundamental measure of information.

### 4.3 Redundancy of Prefix Codes

The entropy bound for a uniquely decodable code has been proved in Section 4.1. In this section, we present an alternative proof specifically for prefix codes which offers much insight into the redundancy of such codes.

Let X be a source random variable with probability distribution

$$\{p_1, p_2, \cdots, p_m\},$$
 (4.49)

where  $m \geq 2$ . A *D*-ary prefix code for *X* can be represented by a *D*-ary code tree with *m* leaves, where each leaf corresponds to a codeword. We denote the leaf corresponding to  $p_i$  by  $c_i$  and the order of  $c_i$  by  $l_i$ , and assume that the alphabet is

$$\{0, 1, \cdots, D-1\}.$$
 (4.50)

Let  $\mathcal{I}$  be the index set of all the internal nodes (including the root) in the code tree.

Instead of matching codewords by brute force, we can use the code tree of a prefix code for more efficient decoding. To decode a codeword, we trace the path specified by the codeword from the root of the code tree until it terminates at the leaf corresponding to that codeword. Let  $q_k$  be the probability of reaching an internal node  $k \in \mathcal{I}$  during the decoding process. The probability  $q_k$  is called the *reaching probability* of internal node k. Evidently,  $q_k$  is equal to the sum of the probabilities of all the leaves descending from node k.

Let  $\tilde{p}_{k,j}$  be the probability that the *j*th branch of node *k* is taken during the decoding process. The probabilities  $\tilde{p}_{k,j}, 0 \leq j \leq D-1$ , are called the *branching probabilities* of node *k*, and

$$q_k = \sum_j \tilde{p}_{k,j}.$$
(4.51)

Once node k is reached, the *conditional branching distribution* is

$$\left\{\frac{\tilde{p}_{k,0}}{q_k}, \frac{\tilde{p}_{k,1}}{q_k}, \cdots, \frac{\tilde{p}_{k,D-1}}{q_k}\right\}.$$
(4.52)

Then define the *conditional entropy* of node k by

$$h_{k} = H_{D}\left(\left\{\frac{\tilde{p}_{k,0}}{q_{k}}, \frac{\tilde{p}_{k,1}}{q_{k}}, \cdots, \frac{\tilde{p}_{k,D-1}}{q_{k}}\right\}\right),$$
(4.53)

where with a slight abuse of notation, we have used  $H_D(\cdot)$  to denote the entropy in the base D of the conditional branching distribution in the parenthesis. By Theorem 2.43,  $h_k \leq 1$ . The following lemma relates the entropy of X with the structure of the code tree.

### **Lemma 4.18.** $H_D(X) = \sum_{k \in \mathcal{I}} q_k h_k.$

*Proof.* We prove the lemma by induction on the number of internal nodes of the code tree. If there is only one internal node, it must be the root of the tree. Then the lemma is trivially true upon observing that the reaching probability of the root is equal to 1.

Assume the lemma is true for all code trees with n internal nodes. Now consider a code tree with n + 1 internal nodes. Let k be an internal node such that k is the parent of a leaf c with maximum order. Each sibling of c may or may not be a leaf. If it is not a leaf, then it cannot be the ascendent of another leaf because we assume that c is a leaf with maximum order. Now consider revealing the outcome of X in two steps. In the first step, if the outcome of X is not a leaf descending from node k, we identify the outcome exactly, otherwise we identify the outcome to be a child of node k. We call this random variable V. If we do not identify the outcome exactly in the first step, which happens with probability  $q_k$ , we further identify in the second step which of the children (child) of node k the outcome is (there is only one child of node k which can be the outcome if all the siblings of c are not leaves). We call this random variable W. If the second step is not necessary, we assume that W takes a constant value with probability 1. Then X = (V, W).

The outcome of V can be represented by a code tree with n internal nodes which is obtained by pruning the original code tree at node k. Then by the induction hypothesis,

$$H(V) = \sum_{k' \in \mathcal{I} \setminus \{k\}} q_{k'} h_{k'}.$$
(4.54)

By the chain rule for entropy, we have

$$H(X) = H(V) + H(W|V)$$
(4.55)

$$= \sum_{k' \in \mathcal{I} \setminus \{k\}} q_{k'} h_{k'} + (1 - q_k) \cdot 0 + q_k h_k$$
(4.56)

$$=\sum_{k'\in\mathcal{I}}q_{k'}h_{k'}.$$
(4.57)

The lemma is proved.  $\hfill\square$ 

The next lemma expresses the expected length L of a prefix code in terms of the reaching probabilities of the internal nodes of the code tree.

# Lemma 4.19. $L = \sum_{k \in \mathcal{I}} q_k$ .

Proof. Define

$$a_{ki} = \begin{cases} 1 \text{ if leaf } c_i \text{ is a descendent of internal node k} \\ 0 \text{ otherwise.} \end{cases}$$
(4.58)

Then

$$l_i = \sum_{k \in \mathcal{I}} a_{ki}, \tag{4.59}$$

because there are exactly  $l_i$  internal nodes of which  $c_i$  is a descendent if the order of  $c_i$  is  $l_i$ . On the other hand,

$$q_k = \sum_i a_{ki} p_i. \tag{4.60}$$

Then

$$L = \sum_{i} p_i l_i \tag{4.61}$$

$$=\sum_{i} p_i \sum_{k \in \mathcal{I}} a_{ki} \tag{4.62}$$

$$=\sum_{k\in\mathcal{I}}\sum_{i}p_{i}a_{ki}\tag{4.63}$$

$$=\sum_{k\in\mathcal{I}}q_k,\tag{4.64}$$

proving the lemma.  $\Box$ 

Define the *local redundancy* of an internal node k by

$$r_k = q_k (1 - h_k). (4.65)$$

This quantity is local to node k in the sense that it depends only on the branching probabilities of node k, and it vanishes if and only if  $\tilde{p}_{k,j} = q_k/D$  for all j, i.e., if and only if the node is *balanced*. Note that  $r_k \ge 0$  because  $h_k \le 1$ .

The next theorem says that the redundancy R of a prefix code is equal to the sum of the local redundancies of all the internal nodes of the code tree.

**Theorem 4.20 (Local Redundancy Theorem).** Let L be the expected length of a D-ary prefix code for a source random variable X, and R be the redundancy of the code. Then

$$R = \sum_{k \in \mathcal{I}} r_k. \tag{4.66}$$

*Proof.* By Lemmas 4.18 and 4.19, we have

$$R = L - H_D(X) \tag{4.67}$$

$$=\sum_{k\in\mathcal{I}}q_k-\sum_k q_kh_k\tag{4.68}$$

$$=\sum_{k\in\mathcal{I}}q_k(1-h_k)\tag{4.69}$$

$$=\sum_{k\in\mathcal{I}}r_k.$$
(4.70)

The theorem is proved.  $\Box$ 

We now present an slightly different version of the entropy bound.

**Corollary 4.21 (Entropy Bound).** Let R be the redundancy of a prefix code. Then  $R \ge 0$  with equality if and only if all the internal nodes in the code tree are balanced.

*Proof.* Since  $r_k \ge 0$  for all k, it is evident from the local redundancy theorem that  $R \ge 0$ . Moreover R = 0 if and only if  $r_k = 0$  for all k, which means that all the internal nodes in the code tree are balanced.  $\Box$ 

**Remark** Before the entropy bound was stated in Theorem 4.5, we gave the intuitive explanation that the entropy bound results from the fact that a D-ary symbol can carry at most one D-it of information. Therefore, when the entropy bound is tight, each code symbol has to carry exactly one D-it of information. Now consider revealing a random codeword one symbol after another. The above corollary states that in order for the entropy bound to be tight, all the internal nodes in the code tree must be balanced. That is, as long as the codeword is not completed, the next code symbol to be revealed always carries one D-it of information because it distributes uniformly on the alphabet. This is consistent with the intuitive explanation we gave for the entropy bound.

*Example 4.22.* The local redundancy theorem allows us to lower bound the redundancy of a prefix code based on partial knowledge on the structure of the code tree. More specifically,

$$R \ge \sum_{k \in \mathcal{I}'} r_k \tag{4.71}$$

for any subset  $\mathcal{I}'$  of  $\mathcal{I}$ .

Let  $p_{m-1}, p_m$  be the two smallest probabilities in the source distribution. In constructing a binary Huffman code,  $p_{m-1}$  and  $p_m$  are merged. Then the redundancy of a Huffman code is lower bounded by

$$(p_{m-1} + p_m) \left[ 1 - H_2 \left( \left\{ \frac{p_{m-1}}{p_{m-1} + p_m}, \frac{p_m}{p_{m-1} + p_m} \right\} \right) \right], \tag{4.72}$$

the local redundancy of the parent of the two leaves corresponding to  $p_{m-1}$  and  $p_m$ . See Yeung [358] for progressive lower and upper bounds on the redundancy of a Huffman code.

# Problems

- Construct a binary Huffman code for the distribution {0.25, 0.05, 0.1, 0.13, 0.2, 0.12, 0.08, 0.07}.
- 2. Construct a ternary Huffman code for the source distribution in Problem 1.
- 3. Show that a Huffman code is an optimal uniquely decodable code for a given source distribution.
- 4. Construct an optimal binary prefix code for the source distribution in Problem 1 such that all the codewords have even lengths.
- 5. Prove directly that the codeword lengths of a prefix code satisfy the Kraft inequality without using Theorem 4.4.
- 6. Prove that if  $p_1 > 0.4$ , then the shortest codeword of a binary Huffman code has length equal to 1. Then prove that the redundancy of such a Huffman code is lower bounded by  $1 h_b(p_1)$ . (Johnsen [172].)
- 7. *Suffix codes* A code is a suffix code if no codeword is a suffix of any other codeword. Show that a suffix code is uniquely decodable.

### 96 4 Zero-Error Data Compression

8. Fix-free codes A code is a fix-free code if it is both a prefix code and a suffix code. Let  $l_1, l_2, \dots, l_m$  be m positive integers. Prove that if

$$\sum_{k=1}^{m} 2^{-l_k} \le \frac{1}{2},$$

then there exists a binary fix-free code with codeword lengths  $l_1, l_2, \dots, l_m$ . (Ahlswede *et al.* [6].)

- 9. Random coding for prefix codes Construct a binary prefix code with codeword lengths  $l_1 \leq l_2 \leq \cdots \leq l_m$  as follows. For each  $1 \leq k \leq m$ , the codeword with length  $l_k$  is chosen independently from the set of all  $2^{l_k}$ possible binary strings with length  $l_k$  according the uniform distribution. Let  $P_m(good)$  be the probability that the code so constructed is a prefix code.
  - a) Prove that  $P_2(good) = (1 2^{-l_1})^+$ , where

$$(x)^{+} = \begin{cases} x \text{ if } x \ge 0\\ 0 \text{ if } x < 0 \end{cases}$$

b) Prove by induction on m that

$$P_m(good) = \prod_{k=1}^m \left(1 - \sum_{j=1}^{k-1} s^{-l_j}\right)^+$$

c) Observe that there exists a prefix code with codeword lengths  $l_1, l_2, \dots, l_m$  if and only if  $P_m(good) > 0$ . Show that  $P_m(good) > 0$  is equivalent to the Kraft inequality.

By using this random coding method, one can derive the Kraft inequality without knowing the inequality ahead of time. (Ye and Yeung [354].)

10. Let X be a source random variable. Suppose a certain probability mass  $p_k$  in the distribution of X is given. Let

$$l_j = \begin{cases} \lceil -\log p_j \rceil & \text{if } j = k \\ \lceil -\log(p_j + x_j) \rceil & \text{if } j \neq k, \end{cases}$$

where

$$x_j = p_j \left( \frac{p_k - 2^{-\lceil -\log p_k \rceil}}{1 - p_k} \right)$$

- for all  $j \neq k$ .
- a) Show that  $1 \leq l_j \leq \lfloor -\log p_j \rfloor$  for all j.
- b) Show that  $\{l_j\}$  satisfies the Kraft inequality.
- c) Obtain an upper bound on  $L_{\text{Huff}}$  in terms of H(X) and  $p_k$  which is tighter than H(X)+1. This shows that when partial knowledge about the source distribution in addition to the source entropy is available, tighter upper bounds on  $L_{\text{Huff}}$  can be obtained.

(Ye and Yeung [355].)

## **Historical Notes**

The foundation for the material in this chapter can be found in Shannon's original paper [291]. The Kraft inequality for uniquely decodable codes was first proved by McMillan [241]. The proof given here is due to Karush [175]. The Huffman coding procedure was devised and proved to be optimal by Huffman [156]. The same procedure was devised independently by Zimmerman [375]. Linder *et al.* [211] have proved the existence of an optimal prefix code for an infinite source alphabet which can be constructed from Huffman codes for truncations of the source distribution. The local redundancy theorem is due to Yeung [358]. A comprehensive survey of code trees for lossless data compression can be found in Abrahams [1].

# Weak Typicality

In the last chapter, we have discussed the significance of entropy in the context of zero-error data compression. In this chapter and the next, we explore entropy in terms of the asymptotic behavior of i.i.d. sequences. Specifically, two versions of the asymptotic equipartition property (AEP), namely the weak AEP and the strong AEP, are discussed. The role of these AEP's in information theory is analogous to the role of the weak law of large numbers in probability theory. In this chapter, the weak AEP and its relation with the source coding theorem are discussed. All the logarithms are in the base 2 unless otherwise specified.

## 5.1 The Weak AEP

We consider an information source  $\{X_k, k \geq 1\}$  where  $X_k$  are i.i.d. with distribution p(x). We use X to denote the generic random variable and H(X) to denote the common entropy for all  $X_k$ , where  $H(X) < \infty$ . Let  $\mathbf{X} = (X_1, X_2, \dots, X_n)$ . Since  $X_k$  are i.i.d.,

$$p(\mathbf{X}) = p(X_1)p(X_2)\cdots p(X_n).$$
(5.1)

Note that  $p(\mathbf{X})$  is a random variable because it is a function of the random variables  $X_1, X_2, \dots, X_n$ . We now prove an asymptotic property of  $p(\mathbf{X})$  called the *weak asymptotic equipartition property* (weak AEP).

### Theorem 5.1 (Weak AEP I).

$$-\frac{1}{n}\log p(\mathbf{X}) \to H(X) \tag{5.2}$$

in probability as  $n \to \infty$ , i.e., for any  $\epsilon > 0$ , for n sufficiently large,

$$\Pr\left\{\left|-\frac{1}{n}\log p(\mathbf{X}) - H(X)\right| \le \epsilon\right\} > 1 - \epsilon.$$
(5.3)

100 5 Weak Typicality

*Proof.* Since  $X_1, X_2, \dots, X_n$  are i.i.d., by (5.1),

$$-\frac{1}{n}\log p(\mathbf{X}) = -\frac{1}{n}\sum_{k=1}^{n}\log p(X_k).$$
 (5.4)

The random variables  $\log p(X_k)$  are also i.i.d. Then by the weak law of large numbers, the right hand side of (5.4) tends to

$$-E\log p(X) = H(X), \tag{5.5}$$

in probability, proving the theorem.  $\Box$ 

The weak AEP is nothing more than a straightforward application of the weak law of large numbers. However, as we will see shortly, this property has significant implications.

**Definition 5.2.** The weakly typical set  $W_{[X]\epsilon}^n$  with respect to p(x) is the set of sequences  $\mathbf{x} = (x_1, x_2, \dots, x_n) \in \mathcal{X}^n$  such that

$$\left| -\frac{1}{n} \log p(\mathbf{x}) - H(X) \right| \le \epsilon, \tag{5.6}$$

or equivalently,

$$H(X) - \epsilon \le -\frac{1}{n} \log p(\mathbf{x}) \le H(X) + \epsilon, \tag{5.7}$$

where  $\epsilon$  is an arbitrarily small positive real number. The sequences in  $W_{[X]\epsilon}^n$  are called weakly  $\epsilon$ -typical sequences.

The quantity

$$-\frac{1}{n}\log p(\mathbf{x}) = -\frac{1}{n}\sum_{k=1}^{n}\log p(x_k)$$
(5.8)

is called the *empirical entropy* of the sequence **x**. The empirical entropy of a weakly typical sequence is close to the true entropy H(X). The important properties of the set  $W_{[X]_{\epsilon}}^n$  are summarized in the next theorem which we will see is equivalent to the weak AEP.

**Theorem 5.3 (Weak AEP II).** The following hold for any  $\epsilon > 0$ :

1) If 
$$\mathbf{x} \in W_{[X]\epsilon}^n$$
, then

$$2^{-n(H(X)+\epsilon)} \le p(\mathbf{x}) \le 2^{-n(H(X)-\epsilon)}.$$
(5.9)

2) For n sufficiently large,

$$\Pr\{\mathbf{X} \in W^n_{[X]\epsilon}\} > 1 - \epsilon.$$
(5.10)

3) For n sufficiently large,

$$(1-\epsilon)2^{n(H(X)-\epsilon)} \le |W_{[X]\epsilon}^n| \le 2^{n(H(X)+\epsilon)}.$$
(5.11)

*Proof.* Property 1 follows immediately from the definition of  $W_{[X]\epsilon}^n$  in (5.7). Property 2 is equivalent to Theorem 5.1. To prove Property 3, we use the lower bound in (5.9) and consider

$$|W_{[X]\epsilon}^n|^{2^{-n(H(X)+\epsilon)}} \le \Pr\{W_{[X]\epsilon}^n\} \le 1,$$
(5.12)

which implies

$$|W_{[X]\epsilon}^{n}| \le 2^{n(H(X)+\epsilon)}.$$
(5.13)

Note that this upper bound holds for any  $n \ge 1$ . On the other hand, using the upper bound in (5.9) and Theorem 5.1, for n sufficiently large, we have

$$1 - \epsilon \le \Pr\{W_{[X]\epsilon}^n\} \le |W_{[X]\epsilon}^n|^{2^{-n(H(X)-\epsilon)}}.$$
(5.14)

Then

$$|W_{[X]\epsilon}^n| \ge (1-\epsilon)2^{n(H(X)-\epsilon)}.$$
 (5.15)

Combining (5.13) and (5.15) gives Property 3. The theorem is proved.  $\Box$ 

**Remark** Theorem 5.3 is a consequence of Theorem 5.1. However, Property 2 in Theorem 5.3 is equivalent to Theorem 5.1. Therefore, Theorem 5.1 and Theorem 5.3 are equivalent, and they will both be referred to as the weak AEP.

The weak AEP has the following interpretation. Suppose  $\mathbf{X} = (X_1, X_2, \cdots, X_n)$  is drawn i.i.d. according to p(x), where n is large. After the sequence is drawn, we ask what the probability of occurrence of the sequence is. The weak AEP says that the probability of occurrence of the sequence drawn is close to  $2^{-nH(X)}$  with very high probability. Such a sequence is called a weakly typical sequence. Moreover, the total number of weakly typical sequences is approximately equal to  $2^{nH(X)}$ . The weak AEP, however, does not say that most of the sequences in  $\mathcal{X}^n$  are weakly typical. In fact, the number of weakly typical sequences is in general insignificant compared with the total number of sequences, because

$$\frac{|W_{[X]\delta}^n|}{|\mathcal{X}|^n} \approx \frac{2^{nH(X)}}{2^{n\log|\mathcal{X}|}} = 2^{-n(\log|\mathcal{X}| - H(X))} \to 0$$
(5.16)

as  $n \to \infty$  as long as H(X) is strictly less than  $\log |\mathcal{X}|$ . The idea is that, although the size of the weakly typical set may be insignificant compared with the size of the set of all sequences, the former has almost all the probability.

When n is large, one can almost think of the sequence **X** as being obtained by choosing a sequence from the weakly typical set according to the uniform

### 102 5 Weak Typicality

distribution. Very often, we concentrate on the properties of typical sequences because any property which is proved to be true for typical sequences will then be true with high probability. This in turn determines the average behavior of a large sample.

**Remark** The most likely sequence is in general not weakly typical although the probability of the weakly typical set is close to 1 when n is large. For example, for  $X_k$  i.i.d. with p(0) = 0.1 and p(1) = 0.9,  $(1, 1, \dots, 1)$  is the most likely sequence, but it is not weakly typical because its empirical entropy is not close to the true entropy. The idea is that as  $n \to \infty$ , the probability of every sequence, including that of the most likely sequence, tends to 0. Therefore, it is not necessary for a weakly typical set to include the most likely sequence in order to possess a probability close to 1.

## 5.2 The Source Coding Theorem

To encode a random sequence  $\mathbf{X} = (X_1, X_2, \dots, X_n)$  drawn i.i.d. according to p(x) by a *block code*, we construct a one-to-one mapping from a subset  $\mathcal{A}$ of  $\mathcal{X}^n$  to an index set

$$\mathcal{I} = \{1, 2, \cdots, M\},\tag{5.17}$$

where  $|\mathcal{A}| = M \leq |\mathcal{X}|^n$ . We do not have to assume that  $|\mathcal{X}|$  is finite. The indices in  $\mathcal{I}$  are called *codewords*, and the integer n is called the *block length* of the code. If a sequence  $\mathbf{x} \in \mathcal{A}$  occurs, the encoder outputs the corresponding codeword which is specified by approximately  $\log M$  bits. If a sequence  $\mathbf{x} \notin \mathcal{A}$  occurs, the encoder outputs the constant codeword 1. In either case, the codeword output by the encoder is decoded to the sequence in  $\mathcal{A}$  corresponding to that codeword by the decoder. If a sequence  $\mathbf{x} \notin \mathcal{A}$  occurs, then  $\mathbf{x}$  is decoded correctly by the decoder. If a sequence  $\mathbf{x} \notin \mathcal{A}$  occurs, then  $\mathbf{x}$  is not decoded correctly by the decoder. For such a code, its performance is measured by the coding rate defined as  $n^{-1} \log M$  (in bits per source symbol), and the probability of error is given by

$$P_e = \Pr\{\mathbf{X} \notin \mathcal{A}\}.$$
 (5.18)

If the code is not allowed to make any error, i.e.,  $P_e = 0$ , it is clear that M must be taken to be  $|\mathcal{X}|^n$ , or  $\mathcal{A} = \mathcal{X}^n$ . In that case, the coding rate is equal to  $\log |\mathcal{X}|$ . However, if we allow  $P_e$  to be any small quantity, Shannon [291] showed that there exists a block code whose coding rate is arbitrarily close to H(X) when n is sufficiently large. This is the direct part of Shannon's source coding theorem, and in this sense the source sequence **X** is said to be reconstructed almost perfectly.

We now prove the direct part of the source coding theorem by constructing a desired code. First, we fix  $\epsilon > 0$  and take

5.2 The Source Coding Theorem 103

$$\mathcal{A} = W^n_{[X]\epsilon} \tag{5.19}$$

and

$$M = |\mathcal{A}|. \tag{5.20}$$

For sufficiently large n, by the weak AEP,

$$(1-\epsilon)2^{n(H(X)-\epsilon)} \le M = |\mathcal{A}| = |W_{[X]\epsilon}^n| \le 2^{n(H(X)+\epsilon)}.$$
 (5.21)

Therefore, the coding rate  $n^{-1} \log M$  satisfies

$$\frac{1}{n}\log(1-\epsilon) + H(X) - \epsilon \le \frac{1}{n}\log M \le H(X) + \epsilon.$$
(5.22)

Also by the weak AEP,

$$P_e = \Pr\{\mathbf{X} \notin \mathcal{A}\} = \Pr\{\mathbf{X} \notin W_{[X]\epsilon}^n\} < \epsilon.$$
(5.23)

Letting  $\epsilon \to 0$ , the coding rate tends to H(X), while  $P_e$  tends to 0. This proves the direct part of the source coding theorem.

The converse part of the source coding theorem says that if we use a block code with block length n and coding rate less than  $H(X) - \zeta$ , where  $\zeta > 0$ does not change with n, then  $P_e \to 1$  as  $n \to \infty$ . To prove this, consider any code with block length n and coding rate less than  $H(X) - \zeta$ , so that M, the total number of codewords, is at most  $2^{n(H(X)-\zeta)}$ . We can use some of these codewords for the typical sequences  $\mathbf{x} \in W_{[X]\epsilon}^n$ , and some for the non-typical sequences  $\mathbf{x} \notin W_{[X]\epsilon}^n$ . The total probability of the typical sequences covered by the code, by the weak AEP, is upper bounded by

$$2^{n(H(X)-\zeta)}2^{-n(H(X)-\epsilon)} = 2^{-n(\zeta-\epsilon)}.$$
(5.24)

Therefore, the total probability covered by the code is upper bounded by

$$2^{-n(\zeta-\epsilon)} + \Pr\{\mathbf{X} \notin W_{[X]\epsilon}^n\} < 2^{-n(\zeta-\epsilon)} + \epsilon$$
(5.25)

for n sufficiently large, again by the weak AEP. This probability is equal to  $1 - P_e$  because  $P_e$  is the probability that the source sequence **X** is not covered by the code. Thus

$$1 - P_e < 2^{-n(\zeta - \epsilon)} + \epsilon, \tag{5.26}$$

or

$$P_e > 1 - (2^{-n(\zeta - \epsilon)} + \epsilon).$$
 (5.27)

This inequality holds when n is sufficiently large for any  $\epsilon > 0$ , in particular for  $\epsilon < \zeta$ . Then for any  $\epsilon < \zeta$ ,  $P_e > 1 - 2\epsilon$  when n is sufficiently large. Hence,  $P_e \to 1$  as  $n \to \infty$  and then  $\epsilon \to 0$ . This proves the converse part of the source coding theorem.

104 5 Weak Typicality

### 5.3 Efficient Source Coding

**Theorem 5.4.** Let  $\mathbf{Y} = (Y_1, Y_2, \dots, Y_m)$  be a random binary sequence of length m. Then  $H(\mathbf{Y}) \leq m$  with equality if and only if  $Y_i$  are drawn i.i.d. according to the uniform distribution on  $\{0, 1\}$ .

*Proof.* By the independence bound for entropy,

$$H(\mathbf{Y}) \le \sum_{i=1}^{m} H(Y_i) \tag{5.28}$$

with equality if and only if  $Y_i$  are mutually independent. By Theorem 2.43,

$$H(Y_i) \le \log 2 = 1 \tag{5.29}$$

with equality if and only if  $Y_i$  distributes uniformly on  $\{0, 1\}$ . Combining (5.28) and (5.29), we have

$$H(\mathbf{Y}) \le \sum_{i=1}^{m} H(Y_i) \le m,$$
(5.30)

where this upper bound is tight if and only if  $Y_i$  are mutually independent and each of them distributes uniformly on  $\{0, 1\}$ . The theorem is proved.  $\Box$ 

Let  $\mathbf{Y} = (Y_1, Y_2, \dots, Y_n)$  be a sequence of length n such that  $Y_i$  are drawn i.i.d. according to the uniform distribution on  $\{0, 1\}$ , and let Y denote the generic random variable. Then H(Y) = 1. According to the source coding theorem, for almost perfect reconstruction of  $\mathbf{Y}$ , the coding rate of the source code must be at least 1. It turns out that in this case it is possible to use a source code with coding rate exactly equal to 1 while the source sequence  $\mathbf{Y}$ can be reconstructed with zero error. This can be done by simply encoding all the  $2^n$  possible binary sequences of length n, i.e., by taking  $M = 2^n$ . Then the coding rate is given by

$$n^{-1}\log M = n^{-1}\log 2^n = 1.$$
(5.31)

Since each symbol in **Y** is a bit and the rate of the best possible code describing **Y** is 1 bit per symbol,  $Y_1, Y_2, \dots, Y_n$  are called *fair bits*, with the connotation that they are incompressible.

It turns out that the whole idea of efficient source coding by a block code is to describe the information source by a binary sequence consisting of "almost fair" bits. Consider a sequence of block codes which encode  $\mathbf{X} = (X_1, X_2, \dots, X_n)$  into  $\mathbf{Y} = (Y_1, Y_2, \dots, Y_m)$ , where  $X_k$  are i.i.d. with generic random variable X,  $\mathbf{Y}$  is a binary sequence with length

$$m \approx nH(X),$$
 (5.32)

and  $n \to \infty$ . For simplicity, we assume that the common alphabet  $\mathcal{X}$  is finite. Let  $\hat{\mathbf{X}} \in \mathcal{X}^n$  be the reconstruction of  $\mathbf{X}$  by the decoder and  $P_e$  be the probability of error, i.e.,

$$P_e = \Pr\{\mathbf{X} \neq \mathbf{X}\}. \tag{5.33}$$

Further assume  $P_e \to 0$  as  $n \to \infty$ . We will show that **Y** consists of almost fair bits.

By Fano's inequality,

$$H(\mathbf{X}|\mathbf{X}) \le 1 + P_e \log |\mathcal{X}|^n = 1 + nP_e \log |\mathcal{X}|.$$
(5.34)

Since  $\hat{\mathbf{X}}$  is a function of  $\mathbf{Y}$ ,

$$H(\mathbf{Y}) = H(\mathbf{Y}, \hat{\mathbf{X}}) \ge H(\hat{\mathbf{X}}).$$
(5.35)

It follows that

$$H(\mathbf{Y}) \ge H(\hat{\mathbf{X}}) \tag{5.36}$$

$$\geq I(\mathbf{X}; \mathbf{X}) \tag{5.37}$$

$$= H(\mathbf{X}) - H(\mathbf{X}|\mathbf{X}) \tag{5.38}$$

$$\geq nH(X) - (1 + nP_e \log |\mathcal{X}|) \tag{5.39}$$

$$= n(H(X) - P_e \log |\mathcal{X}|) - 1.$$
 (5.40)

On the other hand, by Theorem 5.4,

$$H(\mathbf{Y}) \le m. \tag{5.41}$$

Combining (5.40) and (5.41), we have

$$n(H(X) - P_e \log |\mathcal{X}|) - 1 \le H(\mathbf{Y}) \le m.$$
(5.42)

Since  $P_e \to 0$  as  $n \to \infty$ , the above lower bound on  $H(\mathbf{Y})$  is approximately equal to

$$nH(X) \approx m \tag{5.43}$$

when n is large (cf. (5.32)). Therefore,

$$H(\mathbf{Y}) \approx m. \tag{5.44}$$

In light of Theorem 5.4,  $\mathbf{Y}$  almost attains the maximum possible entropy. In this sense, we say that  $\mathbf{Y}$  consists of almost fair bits.

# 5.4 The Shannon-McMillan-Breiman Theorem

For an i.i.d. information source  $\{X_k\}$  with generic random variable X and generic distribution p(x), the weak AEP states that

106 5 Weak Typicality

$$-\frac{1}{n}\log p(\mathbf{X}) \to H(X) \tag{5.45}$$

in probability as  $n \to \infty$ , where  $\mathbf{X} = (X_1, X_2, \dots, X_n)$ . Here H(X) is the entropy of the generic random variables X as well as the entropy rate of the source  $\{X_k\}$ .

In Section 2.10, we showed that the entropy rate H of a source  $\{X_k\}$  exists if the source is stationary. The Shannon-McMillan-Breiman theorem states that if  $\{X_k\}$  is also ergodic, then

$$\Pr\left\{-\lim_{n\to\infty}\frac{1}{n}\log\Pr\{\mathbf{X}\}=H\right\}=1.$$
(5.46)

This means that if  $\{X_k\}$  is stationary and ergodic, then  $-\frac{1}{n}\log \Pr\{\mathbf{X}\}$  not only almost always converges, but it also almost always converges to H. For this reason, the Shannon-McMillan-Breiman theorem is also referred to as the weak AEP for ergodic stationary sources.

The formal definition of an ergodic source and the statement of the Shannon-McMillan-Breiman theorem require the use of measure theory which is beyond the scope of this book. We point out that the event in (5.46) involves an infinite collection of random variables which cannot be described by a joint distribution except in very special cases. Without measure theory, the probability of this event in general cannot be properly defined. However, this does not prevent us from developing some appreciation of the Shannon-McMillan-Breiman theorem.

Let  $\mathcal{X}$  be the common alphabet for a stationary source  $\{X_k\}$ . Roughly speaking, a stationary source  $\{X_k\}$  is ergodic if the time average exhibited by a single realization of the source is equal to the ensemble average with probability 1. More specifically, for any  $k_1, k_2, \dots, k_m$ ,

$$\Pr\left\{\lim_{n \to \infty} \frac{1}{n} \sum_{l=0}^{n-1} f(X_{k_1+l}, X_{k_2+l}, \cdots, X_{k_m+l}) = Ef(X_{k_1}, X_{k_2}, \cdots, X_{k_m})\right\} = 1,$$
(5.47)

where f is a function defined on  $\mathcal{X}^m$  which satisfies suitable conditions. For the special case that  $\{X_k\}$  satisfies

$$\Pr\left\{\lim_{n \to \infty} \frac{1}{n} \sum_{l=1}^{n} X_l = E X_k\right\} = 1,$$
(5.48)

we say that  $\{X_k\}$  is mean ergodic.

Example 5.5. The i.i.d. source  $\{X_k\}$  is mean ergodic under suitable conditions because the strong law of the large numbers states that (5.48) is satisfied.

*Example 5.6.* Consider the source  $\{X_k\}$  defined as follows. Let Z be a binary random variable uniformly distributed on  $\{0, 1\}$ . For all k, let  $X_k = Z$ . Then

$$\Pr\left\{\lim_{n \to \infty} \frac{1}{n} \sum_{l=1}^{n} X_l = 0\right\} = \frac{1}{2}$$
(5.49)

and

$$\Pr\left\{\lim_{n \to \infty} \frac{1}{n} \sum_{l=1}^{n} X_l = 1\right\} = \frac{1}{2}.$$
(5.50)

Since  $EX_k = \frac{1}{2}$ ,

$$\Pr\left\{\lim_{n \to \infty} \frac{1}{n} \sum_{l=1}^{n} X_l = E X_k\right\} = 0.$$
(5.51)

Therefore,  $\{X_k\}$  is not mean ergodic and hence not ergodic.

If an information source  $\{X_k\}$  is stationary and ergodic, by the Shannon-McMillan-Breiman theorem,

$$-\frac{1}{n}\log\Pr\{\mathbf{X}\}\approx H\tag{5.52}$$

when n is large. That is, with probability close to 1, the probability of the sequence **X** which occurs is approximately equal to  $2^{-nH}$ . Then by means of arguments similar to the proof of Theorem 5.3, we see that there exist approximately  $2^{nH}$  sequences in  $\mathcal{X}^n$  whose probabilities are approximately equal to  $2^{-nH}$ , and the total probability of these sequences is almost 1. Therefore, by encoding these sequences with approximately nH bits, the source sequence **X** can be recovered with an arbitrarily small probability of error when the block length n is sufficiently large. This is a generalization of the direct part of the source coding theorem which gives a physical meaning to the entropy rate of an ergodic, although the entropy rate always exists, it may not carry any physical meaning.

As an example, by regarding printed English as a stationary ergodic process, Shannon [294] estimated by a guessing game that its entropy rate is about 1.3 bits per letter. Cover and King [70] described a gambling estimate of the entropy rate of printed English which gives 1.34 bits per letter. These results show that it is not possible to describe printed English accurately by using less than about 1.3 bits per letter. 108 5 Weak Typicality

## Problems

- 1. Show that for any  $\epsilon > 0$ ,  $W_{[X]\epsilon}^n$  is nonempty for sufficiently large n.
- 2. The source coding theorem with a general block code In proving the converse of the source coding theorem, we assume that each codeword in  $\mathcal{I}$  corresponds to a unique sequence in  $\mathcal{X}^n$ . More generally, a block code with block length n is defined by an encoding function  $f: \mathcal{X}^n \to \mathcal{I}$  and a decoding function  $g: \mathcal{I} \to \mathcal{X}^n$ . Prove that  $P_e \to 1$  as  $n \to \infty$  even if we are allowed to use a general block code.
- 3. Following Problem 2, we further assume that we can use a block code with probabilistic encoding and decoding. For such a code, encoding is defined by a transition matrix F from  $\mathcal{X}^n$  to  $\mathcal{I}$  and decoding is defined by a transition matrix G from  $\mathcal{I}$  to  $\mathcal{X}^n$ . Prove that  $P_e \to 1$  as  $n \to \infty$  even if we are allowed to use such a code.
- 4. In the discussion in Section 5.3, we made the assumption that the common alphabet  $\mathcal{X}$  is finite. Can you draw the same conclusion when  $\mathcal{X}$  is countable but H(X) is finite? Hint: use Problem 2.
- 5. Alternative definition of weak typicality Let  $\mathbf{X} = (X_1, X_2, \dots, X_n)$  be an i.i.d. sequence whose generic random variable X is distributed with p(x). Let  $q_{\mathbf{x}}$  be the empirical distribution of the sequence  $\mathbf{x}$ , i.e.,  $q_{\mathbf{x}}(x) = n^{-1}N(x;\mathbf{x})$  for all  $x \in \mathcal{X}$ , where  $N(x;\mathbf{x})$  is the number of occurrence of x in  $\mathbf{x}$ .
  - a) Show that for any  $\mathbf{x} \in \mathcal{X}^n$ ,

$$-\frac{1}{n}\log p(\mathbf{x}) = D(q_{\mathbf{x}}||p) + H(q_{\mathbf{x}}).$$

b) Show that for any  $\epsilon > 0$ , the weakly typical set  $W_{[X]\epsilon}^n$  with respect to p(x) is the set of sequences  $\mathbf{x} \in \mathcal{X}^n$  such that

$$|D(q_{\mathbf{x}}||p) + H(q_{\mathbf{x}}) - H(p)| \le \epsilon.$$

c) Show that for sufficiently large n,

$$\Pr\{|D(q_{\mathbf{x}}||p) + H(q_{\mathbf{x}}) - H(p)| \le \epsilon\} > 1 - \epsilon.$$

(Ho and Yeung [150].)

6. Let p and q be two probability distributions on the same alphabet  $\mathcal{X}$  such that  $H(p) \neq H(q)$ . Show that there exists an  $\epsilon > 0$  such that

$$p^{n}\left(\left\{x^{n} \in \mathcal{X}^{n}: \left|-\frac{1}{n}\log p^{n}(x^{n}) - H(q)\right| < \epsilon\right\}\right) \to 0$$

as  $n \to \infty$ . Give an example that  $p \neq q$  but the above convergence does not hold.

7. Let p and q be two probability distributions on the same alphabet  $\mathcal{X}$  with the same support.

a) Prove that for any  $\delta > 0$ ,

$$p^{n}\left(\left\{x^{n} \in \mathcal{X}^{n}: \left|-\frac{1}{n}\log q^{n}(x^{n}) - (H(p) + D(p||q))\right| < \delta\right\}\right) \to 1$$

as  $n \to \infty$ . b) Prove that for any  $\delta > 0$ ,

$$\left| \left\{ x^{n} \in \mathcal{X}^{n} : \left| -\frac{1}{n} \log q^{n}(x^{n}) - (H(p) + D(p||q)) \right| < \delta \right\} \right| \le 2^{n(H(p) + D(p||q) + \delta)}$$

8. Universal source coding Let  $\mathcal{F} = \{\{X_k^{(s)}, k \geq 1\} : s \in \mathcal{S}\}$  be a family of i.i.d. information sources indexed by a finite set  $\mathcal{S}$  with a common alphabet  $\mathcal{X}$ . Define

$$\bar{H} = \max_{s \in \mathcal{S}} H(X^{(s)})$$

where  $X^{(s)}$  is the generic random variable for  $\{X_k^{(s)}, k \ge 1\}$ , and

$$A^n_{\epsilon}(\mathcal{S}) = \bigcup_{s \in \mathcal{S}} W^n_{[X^{(s)}]\epsilon},$$

where  $\epsilon > 0$ .

a) Prove that for all  $s \in \mathcal{S}$ ,

$$\Pr\{\mathbf{X}^{(s)} \in A^n_{\epsilon}(\mathcal{S})\} \to 1$$

as  $n \to \infty$ , where  $\mathbf{X}^{(s)} = (X_1^{(s)}, X_2^{(s)}, \cdots, X_n^{(s)})$ . b) Prove that for any  $\epsilon' > \epsilon$ ,

$$|A^n_{\epsilon}(\mathcal{S})| \le 2^{n(\bar{H}+\epsilon')}$$

for sufficiently large n.

- c) Suppose we know that an information source is in the family  $\mathcal{F}$  but we do not know which one it is. Devise a compression scheme for the information source such that it is asymptotically optimal for every possible source in  $\mathcal{F}$ .
- 9. Let  $\{X_k, k \ge 1\}$  be an i.i.d. information source with generic random variable X and alphabet  $\mathcal{X}$ . Assume

$$\sum_{x} p(x) [\log p(x)]^2 < \infty$$

and define

$$Z_n = -\frac{\log p(\mathbf{X})}{\sqrt{n}} - \sqrt{n}H(X)$$

for  $n = 1, 2, \dots$ . Prove that  $Z_n \to Z$  in distribution, where Z is a Gaussian random variable with mean 0 and variance  $\sum_x p(x) [\log p(x)]^2 - H(X)^2$ .

110 5 Weak Typicality

# **Historical Notes**

The weak asymptotic equipartition property (AEP), which is instrumental in proving the source coding theorem, was first proved by Shannon in his original paper [291]. In this paper, he also stated that this property can be extended to an ergodic stationary source. Subsequently, McMillan [240] and Breiman [46] proved this property for an ergodic stationary source with a finite alphabet. Chung [66] extended the theme to a countable alphabet.

Weak typicality requires that the empirical entropy of a sequence is close to the true entropy. In this chapter, we introduce a stronger notion of typicality which requires that the relative frequency of each possible outcome is close to the corresponding probability. As we will see later, strong typicality is more powerful and flexible than weak typicality as a tool for theorem proving for memoryless problems. However, strong typicality can be used only for random variables with finite alphabets. Throughout this chapter, typicality refers to strong typicality and all the logarithms are in the base 2 unless otherwise specified.

### 6.1 Strong AEP

We consider an information source  $\{X_k, k \geq 1\}$  where  $X_k$  are i.i.d. with distribution p(x). We use X to denote the generic random variable and H(X) to denote the common entropy for all  $X_k$ , where  $H(X) < \infty$ . Let  $\mathbf{X} = (X_1, X_2, \dots, X_n)$ .

**Definition 6.1.** The strongly typical set  $T^n_{[X]\delta}$  with respect to p(x) is the set of sequences  $\mathbf{x} = (x_1, x_2, \dots, x_n) \in \mathcal{X}^n$  such that  $N(x; \mathbf{x}) = 0$  for  $x \notin \mathcal{S}_X$ , and

$$\sum_{x} \left| \frac{1}{n} N(x; \mathbf{x}) - p(x) \right| \le \delta, \tag{6.1}$$

where  $N(x; \mathbf{x})$  is the number of occurrences of x in the sequence  $\mathbf{x}$ , and  $\delta$  is an arbitrarily small positive real number. The sequences in  $T^n_{[X]\delta}$  are called strongly  $\delta$ -typical sequences.

Throughout this chapter, we adopt the convention that all the summations, products, unions, etc, are taken over the corresponding supports unless otherwise specified. The strongly typical set  $T^n_{[X]\delta}$  shares similar properties with

its weakly typical counterpart, which is summarized as the *strong asymptotic equipartition property* (strong AEP) below. The interpretation of the strong AEP is similar to that of the weak AEP.

**Theorem 6.2 (Strong AEP).** There exists  $\eta > 0$  such that  $\eta \to 0$  as  $\delta \to 0$ , and the following hold:

1) If  $\mathbf{x} \in T^n_{[X]\delta}$ , then

$$2^{-n(H(X)+\eta)} \le p(\mathbf{x}) \le 2^{-n(H(X)-\eta)}.$$
(6.2)

2) For n sufficiently large,

$$\Pr\{\mathbf{X} \in T^n_{[X]\delta}\} > 1 - \delta.$$
(6.3)

3) For n sufficiently large,

$$(1-\delta)2^{n(H(X)-\eta)} \le |T_{[X]\delta}^n| \le 2^{n(H(X)+\eta)}.$$
(6.4)

**Proof** To prove Property 1, for  $\mathbf{x} \in T_{[X]\delta}^n$ , we write

$$p(\mathbf{x}) = \prod_{x} p(x)^{N(x;\mathbf{x})}.$$
(6.5)

Then

$$\log p(\mathbf{x}) = \sum_{x} N(x; \mathbf{x}) \log p(x)$$
(6.6)

$$=\sum_{x} (N(x; \mathbf{x}) - np(x) + np(x)) \log p(x)$$
(6.7)

$$= n \sum_{x} p(x) \log p(x) - n \sum_{x} \left(\frac{1}{n} N(x; \mathbf{x}) - p(x)\right) (-\log p(x))$$
(6.8)

$$= -n\left[H(X) + \sum_{x} \left(\frac{1}{n}N(x;\mathbf{x}) - p(x)\right)(-\log p(x))\right].$$
(6.9)

Since  $\mathbf{x} \in T^n_{[X]\delta}$ ,

$$\sum_{x} \left| \frac{1}{n} N(x; \mathbf{x}) - p(x) \right| \le \delta, \tag{6.10}$$

which implies

$$\left|\sum_{x} \left(\frac{1}{n} N(x; \mathbf{x}) - p(x)\right) \left(-\log p(x)\right)\right|$$

6.1 Strong AEP 113

$$\leq \sum_{x} \left| \frac{1}{n} N(x; \mathbf{x}) - p(x) \right| \left( -\log p(x) \right) \tag{6.11}$$

$$\leq -\log\left(\min_{x} p(x)\right) \sum_{x} \left|\frac{1}{n} N(x; \mathbf{x}) - p(x)\right|$$
(6.12)

$$\leq -\delta \log \left( \min_{x} p(x) \right) \tag{6.13}$$

$$=\eta, \tag{6.14}$$

where

$$\eta = -\delta \log\left(\min_{x} p(x)\right) > 0. \tag{6.15}$$

Therefore,

$$-\eta \le \sum_{x} \left(\frac{1}{n} N(x; \mathbf{x}) - p(x)\right) \left(-\log p(x)\right) \le \eta.$$
(6.16)

It then follows from (6.9) that

$$-n(H(X) + \eta) \le \log p(\mathbf{x}) \le -n(H(X) - \eta), \tag{6.17}$$

 $\mathbf{or}$ 

$$2^{-n(H(X)+\eta)} \le p(\mathbf{x}) \le 2^{-n(H(X)-\eta)},\tag{6.18}$$

where  $\eta \to 0$  as  $\delta \to 0$ , proving Property 1.

To prove Property 2, we write

$$N(x; \mathbf{X}) = \sum_{k=1}^{n} B_k(x), \tag{6.19}$$

where

$$B_k(x) = \begin{cases} 1 \text{ if } X_k = x\\ 0 \text{ if } X_k \neq x. \end{cases}$$
(6.20)

Then  $B_k(x), k = 1, 2, \cdots, n$  are i.i.d. random variables with

$$\Pr\{B_k(x) = 1\} = p(x) \tag{6.21}$$

and

$$\Pr\{B_k(x) = 0\} = 1 - p(x). \tag{6.22}$$

Note that

$$EB_k(x) = (1 - p(x)) \cdot 0 + p(x) \cdot 1 = p(x).$$
(6.23)

By the weak law of large numbers, for any  $\delta > 0$  and for any  $x \in \mathcal{X}$ ,

$$\Pr\left\{\left|\frac{1}{n}\sum_{k=1}^{n}B_{k}(x)-p(x)\right|>\frac{\delta}{|\mathcal{X}|}\right\}<\frac{\delta}{|\mathcal{X}|}$$
(6.24)

for n sufficiently large. Then

$$\Pr\left\{ \left| \frac{1}{n} N(x; \mathbf{X}) - p(x) \right| > \frac{\delta}{|\mathcal{X}|} \text{ for some } x \right\}$$
$$= \Pr\left\{ \left| \frac{1}{n} \sum_{k=1}^{n} B_k(x) - p(x) \right| > \frac{\delta}{|\mathcal{X}|} \text{ for some } x \right\}$$
(6.25)

$$=\Pr\left\{\bigcup_{x}\left\{\left|\frac{1}{n}\sum_{k=1}^{n}B_{k}(x)-p(x)\right|>\frac{\delta}{|\mathcal{X}|}\right\}\right\}$$
(6.26)

$$\leq \sum_{x} \Pr\left\{ \left| \frac{1}{n} \sum_{k=1}^{n} B_k(x) - p(x) \right| > \frac{\delta}{|\mathcal{X}|} \right\}$$
(6.27)

$$<\sum_{x} \frac{\delta}{|\mathcal{X}|}$$
 (6.28)

$$=\delta, \tag{6.29}$$

where we have used the union bound<sup>1</sup> to obtain (6.27). Since

$$\sum_{x} \left| \frac{1}{n} N(x; \mathbf{x}) - p(x) \right| > \delta$$
(6.30)

implies

$$\left|\frac{1}{n}N(x;\mathbf{x}) - p(x)\right| > \frac{\delta}{|\mathcal{X}|} \quad \text{for some } x \in \mathcal{X},$$
(6.31)

we have

$$\Pr\left\{\mathbf{X} \in T_{[X]\delta}^{n}\right\}$$
$$= \Pr\left\{\sum_{x} \left|\frac{1}{n}N(x;\mathbf{X}) - p(x)\right| \le \delta\right\}$$
(6.32)

$$= 1 - \Pr\left\{\sum_{x} \left|\frac{1}{n}N(x;\mathbf{X}) - p(x)\right| > \delta\right\}$$
(6.33)

$$\geq 1 - \Pr\left\{ \left| \frac{1}{n} N(x; \mathbf{X}) - p(x) \right| > \frac{\delta}{|\mathcal{X}|} \text{ for some } x \in \mathcal{X} \right\}$$
(6.34)

$$> 1 - \delta, \tag{6.35}$$

proving Property 2.

Finally, Property 3 follows from Property 1 and Property 2 in exactly the same way as in Theorem 5.3, so the proof is omitted.  $\Box$ 

**Remark** Analogous to weak typicality, we note that the upper bound on  $|T_{[X]\delta}^n|$  in Property 3 holds for all  $n \ge 1$ , and for any  $\delta > 0$ , there exists at least one strongly typical sequence when n is sufficiently large. See Problem 1 in Chapter 5.

<sup>&</sup>lt;sup>1</sup> The union bound refers to  $\Pr\{A \cup B\} \leq \Pr\{A\} + \Pr\{B\}$ .

In the rest of the section, we prove an enhancement of Property 2 of the strong AEP which gives an exponential bound on the probability of obtaining a non-typical vector<sup>2</sup>. This result, however, will not be used until Chapter 21.

**Theorem 6.3.** For sufficiently large n, there exists  $\varphi(\delta) > 0$  such that

$$\Pr\{\mathbf{X} \notin T^n_{[X]\delta}\} < 2^{-n\varphi(\delta)}.$$
(6.36)

The proof of this theorem is based on the Chernoff bound [63] which we prove in the next lemma.

**Lemma 6.4 (Chernoff Bound).** Let Y be a real random variable and s be any nonnegative real number. Then for any real number a,

$$\log \Pr\{Y \ge a\} \le -sa + \log E\left[2^{sY}\right] \tag{6.37}$$

and

$$\log \Pr\{Y \le a\} \le sa + \log E\left[2^{-sY}\right]. \tag{6.38}$$

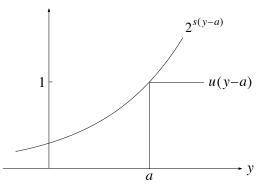
Proof. Let

$$u(y) = \begin{cases} 1 \text{ if } y \ge 0\\ 0 \text{ if } y < 0. \end{cases}$$
(6.39)

Then for any  $s \ge 0$ ,

$$u(y-a) \le 2^{s(y-a)}.$$
(6.40)

This is illustrated in Fig. 6.1. Then



**Fig. 6.1.** An illustration of  $u(y-a) \leq 2^{s(y-a)}$ .

<sup>&</sup>lt;sup>2</sup> This result is due to Ning Cai and Raymond W. Yeung. An alternative proof based on Pinsker's inequality (Theorem 2.33) and the method of types has been given by Prakash Narayan (private communication).

$$E[u(Y-a)] \le E\left[2^{s(Y-a)}\right] = 2^{-sa}E\left[2^{sY}\right].$$
 (6.41)

Since

$$E[u(Y-a)] = \Pr\{Y \ge a\} \cdot 1 + \Pr\{Y < a\} \cdot 0 = \Pr\{Y \ge a\},$$
(6.42)

we see that

$$\Pr\{Y \ge a\} \le 2^{-sa} E\left[2^{sY}\right] = 2^{-sa + \log E\left[2^{sY}\right]}.$$
(6.43)

Then (6.37) is obtained by taking logarithm in the base 2. Upon replacing Y by -Y and a by -a in (6.37), (6.38) is obtained. The lemma is proved.  $\Box$ 

Proof of Theorem 6.3. We will follow the notation in the proof of Theorem 6.2. Consider  $x \in X$  such that p(x) > 0. Applying (6.37), we have

$$\log \Pr\left\{\sum_{k=1}^{n} B_k(x) \ge n\left(p(x) + \delta\right)\right\}$$
  
$$\le -sn\left(p(x) + \delta\right) + \log E\left[2^{s\sum_{k=1}^{n} B_k(x)}\right]$$
(6.44)

$$\stackrel{a)}{=} -sn\left(p(x) + \delta\right) + \log\left(\prod_{k=1}^{n} E\left[2^{sB_k(x)}\right]\right)$$
(6.45)

$$\stackrel{b)}{=} -sn\left(p(x) + \delta\right) + n\log(1 - p(x) + p(x)2^s)$$
(6.46)

$$\leq -sn\left(p(x) + \delta\right) + n(\ln 2)^{-1}(-p(x) + p(x)2^s)$$

$$(6.47)$$

$$= -n \left[ s \left( p(x) + \delta \right) + (\ln 2)^{-1} p(x) (1 - 2^s) \right], \tag{6.48}$$

where

- a) follows because  $B_k(x)$  are mutually independent;
- b) is a direct evaluation of the expectation from the definition of  $B_k(x)$  in (6.20);
- c) follows from the fundamental inequality  $\ln a \le a 1$ .
- In (6.48), upon defining

$$\beta_x(s,\delta) = s \left( p(x) + \delta \right) + (\ln 2)^{-1} p(x) (1 - 2^s), \tag{6.49}$$

we have

$$\log \Pr\left\{\sum_{k=1}^{n} B_k(x) \ge n\left(p(x) + \delta\right)\right\} \le -n\beta_x(s,\delta),\tag{6.50}$$

or

$$\Pr\left\{\sum_{k=1}^{n} B_k(x) \ge n\left(p(x) + \delta\right)\right\} \le 2^{-n\beta_x(s,\delta)}.$$
(6.51)

It is readily seen that

6.1 Strong AEP 117

$$\beta_x(0,\delta) = 0. \tag{6.52}$$

Regarding  $\delta$  as fixed and differentiate with respect to s, we have

$$\beta'_x(s,\delta) = p(x)(1-2^s) + \delta.$$
(6.53)

Then

$$\beta_x'(0,\delta) = \delta > 0 \tag{6.54}$$

and it is readily verified that

$$\beta_x'(s,\delta) \ge 0 \tag{6.55}$$

for

$$0 \le s \le \log\left(1 + \frac{\delta}{p(x)}\right). \tag{6.56}$$

Therefore, we conclude that  $\beta_x(s,\delta)$  is strictly positive for

$$0 < s \le \log\left(1 + \frac{\delta}{p(x)}\right). \tag{6.57}$$

On the other hand, by applying (6.38), we can obtain in the same fashion the bound

$$\log \Pr\left\{\sum_{k=1}^{n} B_k(x) \le n\left(p(x) - \delta\right)\right\} \le -n\sigma_x(s,\delta),\tag{6.58}$$

or

$$\Pr\left\{\sum_{k=1}^{n} B_k(x) \le n\left(p(x) - \delta\right)\right\} \le 2^{-n\sigma_x(s,\delta)},\tag{6.59}$$

where

$$\sigma_x(s,\delta) = -s \left( p(x) - \delta \right) + (\ln 2)^{-1} p(x) (1 - 2^{-s}).$$
(6.60)

Then

$$\sigma_x(0,\delta) = 0,\tag{6.61}$$

 $\quad \text{and} \quad$ 

$$\sigma'_x(s,\delta) = p(x)(2^{-s} - 1) + \delta, \tag{6.62}$$

which is nonnegative for

$$0 \le s \le -\log\left(1 - \frac{\delta}{p(x)}\right). \tag{6.63}$$

In particular,

$$\sigma'_x(0,\delta) = \delta > 0. \tag{6.64}$$

Therefore, we conclude that  $\sigma_x(s, \delta)$  is strictly positive for

$$0 < s \le -\log\left(1 - \frac{\delta}{p(x)}\right). \tag{6.65}$$

By choosing s satisfying

$$0 < s \le \min\left[\log\left(1 + \frac{\delta}{p(x)}\right), -\log\left(1 - \frac{\delta}{p(x)}\right)\right], \tag{6.66}$$

both  $\beta_x(s,\delta)$  and  $\sigma_x(s,\delta)$  are strictly positive. From (6.51) and (6.59), we have

$$\Pr\left\{ \left| \frac{1}{n} \sum_{k=1}^{n} B_k(x) - p(x) \right| \ge \delta \right\}$$
$$= \Pr\left\{ \left| \sum_{k=1}^{n} B_k(x) - np(x) \right| \ge n\delta \right\}$$
(6.67)

$$\leq \Pr\left\{\sum_{k=1}^{n} B_k(x) \geq n\left(p(x) + \delta\right)\right\}$$
$$+\Pr\left\{\sum_{k=1}^{n} B_k(x) \leq n\left(p(x) - \delta\right)\right\}$$
(6.68)

$$\leq 2^{-n\beta_x(s,\delta)} + 2^{-n\sigma_x(s,\delta)} \tag{6.69}$$

$$\leq 2 \cdot 2^{-n \min(\beta_x(s,\delta),\sigma_x(s,\delta))} \tag{6.70}$$

$$=2^{-n\left[\min(\beta_x(s,\delta),\sigma_x(s,\delta))-\frac{1}{n}\right]}$$
(6.71)

$$=2^{-n\varphi_x(\delta)},\tag{6.72}$$

where

$$\varphi_x(\delta) = \min(\beta_x(s,\delta), \sigma_x(s,\delta)) - \frac{1}{n}.$$
(6.73)

Then  $\varphi_x(\delta)$  is strictly positive for sufficiently large n because both  $\beta_x(s,\delta)$ and  $\sigma_x(s,\delta)$  are strictly positive.

Finally, consider

$$\Pr\{\mathbf{X} \in T_{[X]\delta}^n\} = \Pr\left\{\sum_x \left|\frac{1}{n}N(x;\mathbf{X}) - p(x)\right| \le \delta\right\}$$
(6.74)

$$\geq \Pr\left\{ \left| \frac{1}{n} N(x; \mathbf{X}) - p(x) \right| \le \frac{\delta}{|\mathcal{X}|} \text{ for all } x \in \mathcal{X} \right\}$$
(6.75)

$$= 1 - \Pr\left\{ \left| \frac{1}{n} N(x; \mathbf{X}) - p(x) \right| > \frac{\delta}{|\mathcal{X}|} \text{ for some } x \in \mathcal{X} \right\}$$
(6.76)

$$\geq 1 - \sum_{x} \Pr\left\{ \left| \frac{1}{n} N(x; \mathbf{X}) - p(x) \right| > \frac{\delta}{|\mathcal{X}|} \right\}$$
(6.77)

6.2 Strong Typicality Versus Weak Typicality 119

$$=1-\sum_{x}\Pr\left\{\left|\frac{1}{n}\sum_{k=1}^{n}B_{k}(x)-p(x)\right|>\frac{\delta}{|\mathcal{X}|}\right\}$$
(6.78)

$$=1-\sum_{x:p(x)>0}\Pr\left\{\left|\frac{1}{n}\sum_{k=1}^{n}B_{k}(x)-p(x)\right|>\frac{\delta}{|\mathcal{X}|}\right\}$$
(6.79)

$$\geq 1 - \sum_{x:p(x)>0} 2^{-n\varphi_x\left(\frac{\delta}{|\mathcal{X}|}\right)},\tag{6.80}$$

where the last step follows from (6.72). Define

$$\varphi(\delta) = \frac{1}{2} \left[ \min_{x: p(x) > 0} \varphi_x \left( \frac{\delta}{|\mathcal{X}|} \right) \right].$$
(6.81)

Then for sufficiently large n,

=

$$\Pr\{\mathbf{X} \in T^n_{[X]\delta}\} > 1 - 2^{-n\varphi(\delta)},\tag{6.82}$$

or

$$\Pr\{\mathbf{X} \notin T^n_{[X]\delta}\} < 2^{-n\varphi(\delta)},\tag{6.83}$$

where  $\varphi(\delta)$  is strictly positive. The theorem is proved.  $\Box$ 

## 6.2 Strong Typicality Versus Weak Typicality

As we have mentioned at the beginning of the chapter, strong typicality is more powerful and flexible than weak typicality as a tool for theorem proving for memoryless problems, but it can be used only for random variables with finite alphabets. We will prove in the next proposition that strong typicality is stronger than weak typicality in the sense that the former implies the latter.

**Proposition 6.5.** For any  $\mathbf{x} \in \mathcal{X}^n$ , if  $\mathbf{x} \in T^n_{[X]\delta}$ , then  $\mathbf{x} \in W^n_{[X]\eta}$ , where  $\eta \to 0$  as  $\delta \to 0$ .

*Proof.* By Property 1 of strong AEP (Theorem 6.2), if  $\mathbf{x} \in T^n_{[X]\delta}$ , then

$$2^{-n(H(X)+\eta)} \le p(\mathbf{x}) \le 2^{-n(H(X)-\eta)},\tag{6.84}$$

or

$$H(X) - \eta \le -\frac{1}{n}\log p(\mathbf{x}) \le H(X) + \eta, \tag{6.85}$$

where  $\eta \to 0$  as  $\delta \to 0$ . Then  $\mathbf{x} \in W^n_{[X]\eta}$  by Definition 5.2. The proposition is proved.  $\Box$ 

We have proved in this proposition that strong typicality implies weak typicality, but the converse is not true. This idea can be explained without

any detailed analysis. Let X be distributed with p such that p(0) = 0.5, p(1) = 0.25, and p(2) = 0.25. Consider a sequence **x** of length n and let q(i) be the relative frequency of occurrence of symbol i in **x**, i.e.,  $\frac{1}{n}N(i;\mathbf{x})$ , where i = 0, 1, 2. In order for the sequence **x** to be weakly typical, we need

$$-\frac{1}{n}\log p(\mathbf{x}) = -q(0)\log 0.5 - q(1)\log 0.25 - q(2)\log 0.25$$
(6.86)

$$\approx H(X)$$
 (6.87)

$$= -(0.5)\log 0.5 - (0.25)\log 0.25 - (0.25)\log 0.25.$$
(6.88)

Obviously, this can be satisfied by choosing q(i) = p(i) for all *i*. But alternatively, we can choose q(0) = 0.5, q(1) = 0.5, and q(2) = 0. With such a choice of  $\{q(i)\}$ , the sequence **x** is weakly typical with respect to *p* but obviously not strongly typical with respect to *p*, because the relative frequency of occurrence of each symbol *i* is q(i), which is not close to p(i) for i = 1, 2.

Therefore, we conclude that strong typicality is indeed stronger than weak typicality. However, as we have pointed out at the beginning of the chapter, strong typicality can only be used for random variables with finite alphabets.

## 6.3 Joint Typicality

In this section, we discuss strong joint typicality with respect to a bivariate distribution. Generalization to a multivariate distribution is straightforward.

Consider a bivariate information source  $\{(X_k, Y_k), k \ge 1\}$  where  $(X_k, Y_k)$  are i.i.d. with distribution p(x, y). We use (X, Y) to denote the pair of generic random variables.

**Definition 6.6.** The strongly jointly typical set  $T^n_{[XY]\delta}$  with respect to p(x, y) is the set of  $(\mathbf{x}, \mathbf{y}) \in \mathcal{X}^n \times \mathcal{Y}^n$  such that  $N(x, y; \mathbf{x}, \mathbf{y}) = 0$  for  $(x, y) \notin \mathcal{S}_{XY}$ , and

$$\sum_{x} \sum_{y} \left| \frac{1}{n} N(x, y; \mathbf{x}, \mathbf{y}) - p(x, y) \right| \le \delta,$$
(6.89)

where  $N(x, y; \mathbf{x}, \mathbf{y})$  is the number of occurrences of (x, y) in the pair of sequences  $(\mathbf{x}, \mathbf{y})$ , and  $\delta$  is an arbitrarily small positive real number. A pair of sequences  $(\mathbf{x}, \mathbf{y})$  is called strongly jointly  $\delta$ -typical if it is in  $T^n_{[XY]\delta}$ .

Strong typicality satisfies the following *consistency* property.

**Theorem 6.7 (Consistency).** If  $(\mathbf{x}, \mathbf{y}) \in T^n_{[XY]\delta}$ , then  $\mathbf{x} \in T^n_{[X]\delta}$  and  $\mathbf{y} \in T^n_{[Y]\delta}$ .

*Proof.* If  $(\mathbf{x}, \mathbf{y}) \in T^n_{[XY]\delta}$ , then

1216.3 Joint Typicality

$$\sum_{x} \sum_{y} \left| \frac{1}{n} N(x, y; \mathbf{x}, \mathbf{y}) - p(x, y) \right| \le \delta.$$
(6.90)

Upon observing that

$$N(x; \mathbf{x}) = \sum_{y} N(x, y; \mathbf{x}, \mathbf{y}), \qquad (6.91)$$

we have

$$\sum_{x} \left| \frac{1}{n} N(x; \mathbf{x}) - p(x) \right|$$
$$= \sum_{x} \left| \frac{1}{n} \sum_{y} N(x, y; \mathbf{x}, \mathbf{y}) - \sum_{y} p(x, y) \right|$$
(6.92)

$$=\sum_{x}\left|\sum_{y}\left(\frac{1}{n}N(x,y;\mathbf{x},\mathbf{y})-p(x,y)\right)\right|$$
(6.93)

$$\leq \sum_{x} \sum_{y} \left| \frac{1}{n} N(x, y; \mathbf{x}, \mathbf{y}) - p(x, y) \right|$$

$$\leq \delta.$$
(6.94)
(6.95)

$$\leq \delta.$$
 (6.95)

Therefore,  $\mathbf{x} \in T^n_{[X]\delta}$ . Similarly,  $\mathbf{y} \in T^n_{[Y]\delta}$ . The theorem is proved.  $\Box$ 

The following theorem asserts that strong typicality is preserved when a function is applied to a vector.

**Theorem 6.8 (Preservation).** Let Y = f(X). If

$$\mathbf{x} = (x_1, x_2, \cdots, x_n) \in T^n_{[X]\delta},\tag{6.96}$$

then

$$f(\mathbf{x}) = (y_1, y_2, \cdots, y_n) \in T^n_{[Y]\delta},$$
 (6.97)

where  $y_i = f(x_i)$  for  $1 \le i \le n$ .

*Proof.* Consider  $\mathbf{x} \in T^n_{[X]\delta}$ , i.e.,

$$\sum_{x} \left| \frac{1}{n} N(x; \mathbf{x}) - p(x) \right| < \delta.$$
(6.98)

Since Y = f(X),

$$p(y) = \sum_{x \in f^{-1}(y)} p(x)$$
(6.99)

for all  $y \in \mathcal{Y}$ . On the other hand,

$$N(y; f(\mathbf{x})) = \sum_{x \in f^{-1}(y)} N(x; \mathbf{x})$$
(6.100)

for all  $y \in \mathcal{Y}$ . Then

$$\sum_{y} \left| \frac{1}{n} N(y; f(\mathbf{x})) - p(y) \right|$$
$$= \sum_{y} \left| \sum_{x \in f^{-1}(y)} \left( \frac{1}{n} N(x; \mathbf{x}) - p(x) \right) \right|$$
(6.101)

$$\leq \sum_{y} \sum_{x \in f^{-1}(y)} \left| \frac{1}{n} N(x; \mathbf{x}) - p(x) \right|$$
(6.102)

$$=\sum_{x}\left|\frac{1}{n}N(x;\mathbf{x})-p(x)\right|$$
(6.103)

$$<\delta. \tag{6.104}$$

Therefore,  $f(\mathbf{x}) \in T^n_{[Y]\delta}$ , proving the lemma.  $\Box$ 

For a bivariate i.i.d. source  $\{(X_k, Y_k)\}$ , we have the strong joint asymptotic equipartition property (strong JAEP), which can readily be obtained by applying the strong AEP to the source  $\{(X_k, Y_k)\}$ .

### Theorem 6.9 (Strong JAEP). Let

$$(\mathbf{X}, \mathbf{Y}) = ((X_1, Y_1), (X_2, Y_2), \cdots, (X_n, Y_n)), \tag{6.105}$$

where  $(X_i, Y_i)$  are *i.i.d.* with generic pair of random variables (X, Y). Then there exists  $\lambda > 0$  such that  $\lambda \to 0$  as  $\delta \to 0$ , and the following hold:

1) If  $(\mathbf{x}, \mathbf{y}) \in T^n_{[XY]\delta}$ , then

$$2^{-n(H(X,Y)+\lambda)} \le p(\mathbf{x}, \mathbf{y}) \le 2^{-n(H(X,Y)-\lambda)}.$$
(6.106)

2) For n sufficiently large,

$$\Pr\{(\mathbf{X}, \mathbf{Y}) \in T^n_{[XY]\delta}\} > 1 - \delta.$$
(6.107)

3) For n sufficiently large,

$$(1-\delta)2^{n(H(X,Y)-\lambda)} \le |T^n_{[XY]\delta}| \le 2^{n(H(X,Y)+\lambda)}.$$
(6.108)

From the strong JAEP, we can see the following. Since there are approximately  $2^{nH(X,Y)}$  typical  $(\mathbf{x}, \mathbf{y})$  pairs and approximately  $2^{nH(X)}$  typical  $\mathbf{x}$ , for a typical  $\mathbf{x}$ , the number of  $\mathbf{y}$  such that  $(\mathbf{x}, \mathbf{y})$  is jointly typical is approximately

6.3 Joint Typicality 123

$$\frac{2^{nH(X,Y)}}{2^{nH(X)}} = 2^{nH(Y|X)} \tag{6.109}$$

on the average. The next theorem reveals that this is not only true on the average, but it is in fact true for every typical  $\mathbf{x}$  as long as there exists at least one  $\mathbf{y}$  such that  $(\mathbf{x}, \mathbf{y})$  is jointly typical.

**Theorem 6.10.** For any  $\mathbf{x} \in T^n_{[X]\delta}$ , define

$$T^{n}_{[Y|X]\delta}(\mathbf{x}) = \{ \mathbf{y} \in T^{n}_{[Y]\delta} : (\mathbf{x}, \mathbf{y}) \in T^{n}_{[XY]\delta} \}.$$
 (6.110)

If  $|T_{[Y|X]\delta}^n(\mathbf{x})| \geq 1$ , then

$$2^{n(H(Y|X)-\nu)} \le |T^n_{[Y|X]\delta}(\mathbf{x})| \le 2^{n(H(Y|X)+\nu)}, \tag{6.111}$$

where  $\nu \to 0$  as  $n \to \infty$  and  $\delta \to 0$ .

We first prove the following lemma which is along the line of Stirling's approximation [101].

**Lemma 6.11.** For any n > 0,

$$n\ln n - n < \ln n! < (n+1)\ln(n+1) - n.$$
(6.112)

*Proof.* First, we write

$$\ln n! = \ln 1 + \ln 2 + \dots + \ln n. \tag{6.113}$$

Since  $\ln x$  is a monotonically increasing function of x, we have

$$\int_{k-1}^{k} \ln x \, dx < \ln k < \int_{k}^{k+1} \ln x \, dx. \tag{6.114}$$

Summing over  $1 \le k \le n$ , we have

$$\int_{0}^{n} \ln x \, dx < \ln n! < \int_{1}^{n+1} \ln x \, dx, \tag{6.115}$$

or

$$n \ln n - n < \ln n! < (n+1) \ln(n+1) - n.$$
 (6.116)

The lemma is proved.  $\Box$ 

Proof of Theorem 6.10. Let  $\delta$  be a small positive real number and n be a large positive integer to be specified later. Fix an  $\mathbf{x} \in T^n_{[X]\delta}$ , so that

$$\sum_{x} \left| \frac{1}{n} N(X; \mathbf{x}) - p(x) \right| \le \delta.$$
(6.117)

This implies that for all  $x \in \mathcal{X}$ ,

$$\left|\frac{1}{n}N(X;\mathbf{x}) - p(x)\right| \le \delta,\tag{6.118}$$

or

$$p(x) - \delta \le \frac{1}{n} N(x; \mathbf{x}) \le p(x) + \delta.$$
(6.119)

We first prove the upper bound on  $|T_{[Y|X]\delta}^n(\mathbf{x})|$ . For any  $\nu > 0$ , consider

$$2^{-n(H(X)-\nu/2)} \stackrel{a)}{\geq} p(\mathbf{x})$$
 (6.120)

$$=\sum_{\mathbf{y}\in\mathcal{Y}^n} p(\mathbf{x},\mathbf{y}) \tag{6.121}$$

$$\geq \sum_{\mathbf{y}\in T^n_{[Y|X]\delta}(\mathbf{x})} p(\mathbf{x}, \mathbf{y}) \tag{6.122}$$

<sup>b)</sup> 
$$\geq \sum_{\mathbf{y} \in T^n_{[Y|X]\delta}(\mathbf{x})} 2^{-n(H(X)+\nu/2)}$$
 (6.123)

$$\geq |T_{[Y|X]\delta}^{n}(\mathbf{x})|^{2^{-n(H(X)+\nu/2)}}, \qquad (6.124)$$

where a) and b) follow from the strong AEP (Theorem 6.2) and the strong joint AEP (Theorem 6.9), respectively. Then we obtain

$$|T_{[Y|X]\delta}^{n}(\mathbf{x})| \le 2^{-n(H(X)+\nu)},\tag{6.125}$$

which is the upper bound to be proved.

Assume that  $|T_{[Y|X]\delta}^n(\mathbf{x})| \ge 1$ . We now prove the lower bound on  $|T_{[Y|X]\delta}^n(\mathbf{x})|$ . Let

$$\{K(x,y), (x,y) \in \mathcal{X} \times \mathcal{Y}\}$$
(6.126)

be any set of nonnegative integers such that

1.

$$\sum_{y} K(x, y) = N(x; \mathbf{x}) \tag{6.127}$$

for all  $x \in \mathcal{X}$ , and

2. for any  $\mathbf{y} \in \mathcal{Y}^n$ , if

$$N(x, y; \mathbf{x}, \mathbf{y}) = K(x, y) \tag{6.128}$$

for all  $(x, y) \in \mathcal{X} \times \mathcal{Y}$ , then  $(\mathbf{x}, \mathbf{y}) \in T^n_{[XY]\delta}$ .

Then by Definition 6.6,  $\{K(x, y)\}$  satisfy

$$\sum_{x} \sum_{y} \left| \frac{1}{n} K(x, y) - p(x, y) \right| \le \delta, \tag{6.129}$$

6.3 Joint Typicality 125

which implies that for all  $(x, y) \in \mathcal{X} \times \mathcal{Y}$ ,

$$\left|\frac{1}{n}K(x,y) - p(x,y)\right| \le \delta,\tag{6.130}$$

or

$$p(x,y) - \delta \le \frac{1}{n} K(x,y) \le p(x,y) + \delta.$$
(6.131)

Such a set  $\{K(x, y)\}$  exists because  $T^n_{[Y|X]\delta}(\mathbf{x})$  is assumed to be nonempty. Straightforward combinatorics reveals that the number of  $\mathbf{y}$  which satisfy the constraints in (6.128) is equal to

$$M(K) = \prod_{x} \frac{N(x; \mathbf{x})!}{\prod_{y} K(x, y)!},$$
(6.132)

and it is readily seen that

$$|T_{[Y|X]\delta}^n(\mathbf{x})| \ge M(K). \tag{6.133}$$

Using Lemma 6.11, we can lower bound  $\ln M(K)$  as follows.

$$\ln M(K) \geq \sum_{x} \left\{ N(x; \mathbf{x}) \ln N(x; \mathbf{x}) - N(x; \mathbf{x}) - \sum_{y} \left[ (K(x, y) + 1) \ln(K(x, y) + 1) - K(x, y) \right] \right\}$$
(6.134)  

$$\stackrel{a)}{=} \sum_{x} \left[ N(x; \mathbf{x}) \ln N(x; \mathbf{x}) - \sum_{y} (K(x, y) + 1) \ln(K(x, y) + 1) \right]$$
(6.135)  

$$\stackrel{b)}{\geq} \sum_{x} \left\{ N(x; \mathbf{x}) \ln(n (p(x) - \delta)) - \sum_{y} (K(x, y) + 1) \ln\left[ n \left( p(x, y) + \delta + \frac{1}{n} \right) \right] \right\}.$$
(6.136)

In the above, a) follows from (6.127), and b) is obtained by applying the lower bound on  $n^{-1}N(x; \mathbf{x})$  in (6.119) and the upper bound on  $n^{-1}K(x, y)$  in (6.131). Also from (6.127), the coefficient of  $\ln n$  in (6.136) is given by

$$\sum_{x} \left[ N(x; \mathbf{x}) - \sum_{y} (K(x, y) + 1) \right] = -|\mathcal{X}||\mathcal{Y}|.$$
 (6.137)

Let  $\delta$  be sufficiently small and n be sufficiently large so that

$$0 < p(x) - \delta < 1 \tag{6.138}$$

and

$$p(x,y) + \delta + \frac{1}{n} < 1 \tag{6.139}$$

for all x and y. Then in (6.136), both the logarithms

$$\ln(p(x) - \delta) \tag{6.140}$$

and

$$\ln\left(p(x,y) + \delta + \frac{1}{n}\right) \tag{6.141}$$

are negative. Note that the logarithm in (6.140) is well-defined by virtue of (6.138). Rearranging the terms in (6.136), applying the upper bound in (6.119) and the lower bound<sup>3</sup> in (6.131), and dividing by n, we have

$$n^{-1} \ln M(K)$$

$$\geq \sum_{x} (p(x) + \delta) \ln (p(x) - \delta) - \sum_{x} \sum_{y} \left( p(x, y) - \delta + \frac{1}{n} \right)$$

$$\times \ln \left( p(x, y) + \delta + \frac{1}{n} \right) - \frac{|\mathcal{X}| |\mathcal{Y}| \ln n}{n}$$
(6.142)

$$= -H_e(X) + H_e(X,Y) + L_l(n,\delta)$$
(6.143)

$$= H_e(Y|X) + L_l(n,\delta), (6.144)$$

where  $L_l(n, \delta)$  denotes a function of n and  $\delta$  which tends to 0 as  $n \to \infty$  and  $\delta \to 0$ . Changing the base of the logarithm to 2, we have

$$n^{-1}\log M(K) \ge H(Y|X) + L_l(n,\delta).$$
 (6.145)

Then it follows from (6.133) that

$$n^{-1} \log |T_{[Y|X]\delta}^{n}(\mathbf{x})| \ge H(Y|X) + L_{l}(n,\delta).$$
(6.146)

Upon replacing  $L_l(n, \delta)$ , we obtain

$$|T_{[Y|X]\delta}^{n}(\mathbf{x})| \ge 2^{n(H(Y|X)-\nu)},$$
 (6.147)

where  $\nu \to 0$  as  $n \to \infty$  and  $\delta \to 0$  as discussed. The theorem is proved.  $\Box$ 

The above theorem says that for any typical  $\mathbf{x}$ , as long as there is one typical  $\mathbf{y}$  such that  $(\mathbf{x}, \mathbf{y})$  is jointly typical, there are approximately  $2^{nH(Y|X)}$   $\mathbf{y}$  such that  $(\mathbf{x}, \mathbf{y})$  is jointly typical. This theorem has the following corollary that the number of such typical  $\mathbf{x}$  grows with n at almost the same rate as the total number of typical  $\mathbf{x}$ .

<sup>&</sup>lt;sup>3</sup> For the degenerate case when p(x, y) = 1 for some x and y,  $p(x, y) + \delta + \frac{1}{n} > 1$ , and the logarithm in (6.141) is in fact positive. Then the upper bound instead of the lower bound should be applied. The details are omitted.

**Corollary 6.12.** For a joint distribution p(x, y) on  $\mathcal{X} \times \mathcal{Y}$ , let  $S^n_{[X]\delta}$  be the set of all sequences  $\mathbf{x} \in T^n_{[X]\delta}$  such that  $T^n_{[Y|X]\delta}(\mathbf{x})$  is nonempty. Then

$$|S_{[X]\delta}^n| \ge (1-\delta)2^{n(H(X)-\psi)},\tag{6.148}$$

where  $\psi \to 0$  as  $n \to \infty$  and  $\delta \to 0$ .

*Proof.* By the consistency of strong typicality (Theorem 6.7), if  $(\mathbf{x}, \mathbf{y}) \in T^n_{[XY]\delta}$ , then  $\mathbf{x} \in T^n_{[X]\delta}$ . In particular,  $\mathbf{x} \in S^n_{[X]\delta}$ . Then

$$T^n_{[XY]\delta} = \bigcup_{\mathbf{x}\in S^n_{[X]\delta}} \{ (\mathbf{x}, \mathbf{y}) : \mathbf{y} \in T^n_{[Y|X]\delta}(\mathbf{x}) \}.$$
(6.149)

Using the lower bound on  $|T_{[XY]\delta}^n|$  in Theorem 6.9 and the upper bound on  $|T_{[Y|X]\delta}^n(\mathbf{x})|$  in the last theorem, we have

$$(1-\delta)2^{n(H(X,Y)-\lambda)} \le |T^n_{[XY]\delta}| \le |S^n_{[X]\delta}|2^{n(H(Y|X)+\nu)}$$
(6.150)

which implies

$$|S_{[X]\delta}^n| \ge (1-\delta)2^{n(H(X)-(\lambda+\nu))}.$$
(6.151)

The theorem is proved upon letting  $\psi = \lambda + \nu$ .  $\Box$ 

We have established a rich set of structural properties for strong typicality with respect to a bivariate distribution p(x, y), which is summarized in the two-dimensional *strong joint typicality array* in Figure 6.2. In this array, the

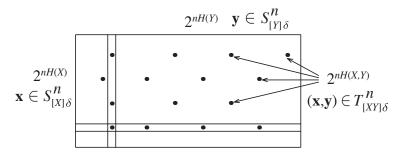


Fig. 6.2. A two-dimensional strong joint typicality array.

rows and the columns are the typical sequences  $\mathbf{x} \in S_{[X]\delta}^n$  and  $\mathbf{y} \in S_{[Y]\delta}^n$ , respectively. The total number of rows and columns are approximately equal to  $2^{nH(X)}$  and  $2^{nH(Y)}$ , respectively. An entry indexed by  $(\mathbf{x}, \mathbf{y})$  receives a dot if  $(\mathbf{x}, \mathbf{y})$  is strongly jointly typical. The total number of dots is approximately equal to  $2^{nH(X,Y)}$ . The number of dots in each row is approximately equal to

 $2^{nH(Y|X)}$ , while the number of dots in each column is approximately equal to  $2^{nH(X|Y)}$ .

For reasons which will become clear in Chapter 16, the strong joint typicality array in Figure 6.2 is said to exhibit an *asymptotic quasi-uniform* structure. By a two-dimensional asymptotic quasi-uniform structure, we mean that in the array all the columns have approximately the same number of dots, and all the rows have approximately the same number of dots. The strong joint typicality array for a multivariate distribution continues to exhibit an asymptotic quasi-uniform structure. The three-dimensional strong joint typicality array with respect to a distribution p(x, y, z) is illustrated in Figure 6.3. As

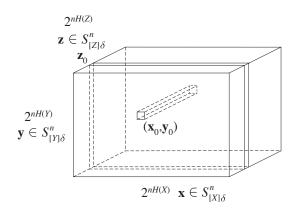


Fig. 6.3. A three-dimensional strong joint typicality array.

before, an entry  $(\mathbf{x}, \mathbf{y}, \mathbf{z})$  receives a dot if  $(\mathbf{x}, \mathbf{y}, \mathbf{z})$  is strongly jointly typical. This is not shown in the figure otherwise it will be very confusing. The total number of dots in the whole array is approximately equal to  $2^{nH(X,Y,Z)}$ . These dots are distributed in the array such that all the planes parallel to each other have approximately the same number of dots, and all the cylinders parallel to each other have approximately the same number of dots. More specifically, the total number of dots on the plane for any fixed  $\mathbf{z}_0 \in S^n_{[Z]\delta}$  (as shown) is approximately equal to  $2^{nH(X,Y|Z)}$ , and the total number of dots in the cylinder for any fixed  $(\mathbf{x}_0, \mathbf{y}_0)$  pair in  $S^n_{[XY]\delta}$  (as shown) is approximately equal to  $2^{nH(Z|X,Y)}$ , so on and so forth.

We see from the strong AEP and Corollary 6.12 that  $S_{[X]\delta}^n$  and  $T_{[X]\delta}^n$  grow with *n* at approximately the same rate. We end this section by stating in the next proposition that  $S_{[X]\delta}^n$  indeed contains almost all the probability when *n* is large. The proof is left as an exercise (see Problem 4).

**Proposition 6.13.** With respect to a joint distribution p(x, y) on  $\mathcal{X} \times \mathcal{Y}$ , for any  $\delta > 0$ ,

$$\Pr\{\mathbf{X} \in S^n_{[X]\delta}\} > 1 - \delta \tag{6.152}$$

for n sufficiently large.

### 6.4 An Interpretation of the Basic Inequalities

The asymptotic quasi-uniform structure exhibited in a strong joint typicality array discussed in the last section is extremely important in information theory. Later in the book, we will see how this structure is involved in proving results such as the channel coding theorem and the rate-distortion theorem. In this section, we show how the basic inequalities can be revealed by examining this structure. It has further been shown by Chan [56] that all unconstrained information inequalities can be obtained from this structure, thus giving a physical meaning to these inequalities.

Consider random variables X, Y, and Z and a fixed  $\mathbf{z} \in S_{[Z]\delta}^n$ , so that  $T_{[XY|Z]\delta}^n(\mathbf{z})$  is nonempty. By the consistency of strong typicality, if  $(\mathbf{x}, \mathbf{y}, \mathbf{z}) \in T_{[XYZ]\delta}^n$ , then  $(\mathbf{x}, \mathbf{z}) \in T_{[XZ]\delta}^n$  and  $(\mathbf{y}, \mathbf{z}) \in T_{[YZ]\delta}^n$ , or  $\mathbf{x} \in T_{[X|Z]\delta}^n(\mathbf{z})$  and  $\mathbf{y} \in T_{[Y|Z]\delta}^n(\mathbf{z})$ , respectively. Thus

$$T^{n}_{[XY|Z]\delta}(\mathbf{z}) \subset T^{n}_{[X|Z]\delta}(\mathbf{z}) \times T^{n}_{[Y|Z]\delta}(\mathbf{z}), \qquad (6.153)$$

which implies

$$|T_{[XY|Z]\delta}^{n}(\mathbf{z})| \leq |T_{[X|Z]\delta}^{n}(\mathbf{z})||T_{[Y|Z]\delta}^{n}(\mathbf{z})|.$$
(6.154)

Applying the lower bound in Theorem 6.10 to  $T^n_{[XY|Z]\delta}(\mathbf{z})$  and the upper bound to  $T^n_{[X|Z]\delta}(\mathbf{z})$  and  $T^n_{[Y|Z]\delta}(\mathbf{z})$ , we have

$$2^{n(H(X,Y|Z)-\zeta)} \le 2^{n(H(X|Z)+\gamma)} 2^{n(H(Y|Z)+\phi)}, \tag{6.155}$$

where  $\zeta, \gamma, \phi \to 0$  as  $n \to \infty$  and  $\delta \to 0$ . Taking logarithm to the base 2 and dividing by n, we obtain

$$H(X, Y|Z) \le H(X|Z) + H(Y|Z)$$
 (6.156)

upon letting  $n \to \infty$  and  $\delta \to 0$ . This inequality is equivalent to

$$I(X;Y|Z) \ge 0.$$
 (6.157)

Thus we have proved the nonnegativity of conditional mutual information. Since all Shannon's information measures are special cases of conditional mutual information, we have proved the nonnegativity of all Shannon's information measures, namely the basic inequalities.

## Problems

1. Show that  $(\mathbf{x}, \mathbf{y}) \in T^n_{[X,Y]\delta}$  and  $(\mathbf{y}, \mathbf{z}) \in T^n_{[Y,Z]\delta}$  do not imply  $(\mathbf{x}, \mathbf{z}) \in T^n_{[X,Z]\delta}$ .

- 130 6 Strong Typicality
- 2. Let  $\mathbf{X} = (X_1, X_2, \dots, X_n)$ , where  $X_k$  are i.i.d. with generic random variable X. Prove that

$$\Pr\{\mathbf{X} \in T^n_{[X]\delta}\} \ge 1 - \frac{|\mathcal{X}|^3}{n\delta^2}$$

for any n and  $\delta > 0$ . This shows that  $\Pr\{\mathbf{X} \in T^n_{[X]\delta}\} \to 1 \text{ as } \delta \to 0 \text{ and } n \to \infty \text{ if } \sqrt{n\delta} \to \infty.$ 

- 3. Prove that for a random variable X with a countable alphabet, Property 2 of the strong AEP holds, while Properties 1 and 3 do not hold.
- 4. Prove Proposition 6.13. Hint: First prove that if  $(\mathbf{X}, \mathbf{Y}) \in T^n_{[XY]\delta}$ , then  $\mathbf{X} \in S^n_{[X]\delta}$ .
- 5. Let  $\mathcal{P}(\mathcal{X})$  be the set of all probability distributions over a finite alphabet  $\mathcal{X}$ . Find a polynomial Q(n) such that for any integer n, there exists a subset  $\mathcal{P}_n(\mathcal{X})$  of  $\mathcal{P}(\mathcal{X})$  such that
  - a)  $|\mathcal{P}_n(\mathcal{X})| \le Q(n);$
  - b) for all  $P \in \mathcal{P}(\mathcal{X})$ , there exists  $P_n \in \mathcal{P}_n(\mathcal{X})$  such that

$$|P_n(x) - P(x)| < \frac{1}{n}$$

for all  $x \in \mathcal{X}$ .

Hint: Let  $\mathcal{P}_n(\mathcal{X})$  be the set of all probability distributions over  $\mathcal{X}$  such that all the probability masses can be expressed as fractions with denominator n.

6. Let p be any probability distribution over a finite set  $\mathcal{X}$  and  $\eta$  be a real number in (0, 1). Prove that for any subset A of  $\mathcal{X}^n$  with  $p^n(A) \ge \eta$ ,

$$|A \cap T^n_{[X]\delta}| \ge 2^{n(H(p)-\delta')},$$

where  $\delta' \to 0$  as  $\delta \to 0$  and  $n \to \infty$ .

In the following problems, for a sequence  $\mathbf{x} \in \mathcal{X}^n$ , let  $q_{\mathbf{x}}$  be the empirical distribution of  $\mathbf{x}$ , i.e.,  $q_{\mathbf{x}}(x) = n^{-1}N(x; \mathbf{x})$  for all  $x \in \mathcal{X}$ . Similarly, for a pair of sequences  $(\mathbf{x}, \mathbf{y}) \in \mathcal{X}^n \times \mathcal{Y}^n$ , let  $q_{\mathbf{x}, \mathbf{y}}$  be the joint empirical distribution of  $(\mathbf{x}, \mathbf{y})$ , i.e.,  $q_{\mathbf{x}, \mathbf{y}}(x, y) = n^{-1}N(x, y; \mathbf{x}, \mathbf{y})$  for all  $(x, y) \in \mathcal{X} \times \mathcal{Y}$ .

7. Alternative definition of strong typicality Show that (6.1) is equivalent to

$$V(q_{\mathbf{x}}, p) \le \delta,$$

where  $V(\cdot, \cdot)$  denotes the variational distance. Thus strong typicality can be regarded as requiring the empirical distribution of a sequence to be close to the probability distribution of the generic random variable in variational distance. Also compare the result here with the alternative definition of weak typicality (Problem 5 in Chapter 5).

- 8. The empirical distribution  $q_{\mathbf{x}}$  of the sequence  $\mathbf{x}$  is also called the *type* of **x**. Assuming that  $\mathcal{X}$  is finite, show that there are a total of  $\binom{n+|\mathcal{X}|-1}{n}$ distinct types  $q_{\mathbf{x}}$ . Hint: There are  $\binom{a+b-1}{a}$  ways to distribute a identical balls in b boxes.
- 9. Unified typicality Let  $\mathbf{X} = (X_1, X_2, \dots, X_n)$  be an i.i.d. sequence whose generic random variable X is distributed with p(x), where the alphabet  $\mathcal{X}$  is countable. For any  $\eta > 0$ , the unified typical set  $U_{[X]\eta}^n$  with respect to p(x) is the set of sequences  $\mathbf{x} \in \mathcal{X}^n$  such that

$$D(q_{\mathbf{x}}||p) + |H(q_{\mathbf{x}}) - H(p)| \le \eta.$$

- a) Show that for any  $\mathbf{x} \in \mathcal{X}^n$ , if  $\mathbf{x} \in U_{[X]\eta}^n$ , then  $\mathbf{x} \in W_{[X]\eta}^n$ . b) Show that for any  $\mathbf{x} \in \mathcal{X}^n$ , if  $\mathbf{x} \in U_{[X]\eta}^n$ , then  $\mathbf{x} \in T_{[X]\delta}^n$ , where  $\delta = \sqrt{\eta \cdot 2 \ln 2}.$

Therefore, unified typicality implies both weak typicality and strong typicality.

- 10. The AEP for unified typicality Unified typicality defined in Problem 9, unlike strong typicality, can be applied to random variables whose alphabets are countable. At the same time, it preserves the essential properties of strong typicality. The following outlines the proof of the AEP which has been discussed in Theorem 5.3 and Theorem 6.2 for weak typicality and strong typicality, respectively.
  - a) Show that

$$2^{-n(H(X)+\eta)} < p(\mathbf{x}) < 2^{-n(H(X)-\eta)}$$

i.e., Property 1 of the AEP.

b) Show that for sufficiently large n,

$$\Pr\{H(q_{\mathbf{x}}) - H(p) > \epsilon\} < \epsilon.$$

Hint: Use the results in Problem 9 above and Problem 5 in Chapter 5. c) It can be proved by means of the result in Problem 9 that

$$\Pr\{H(p) - H(q_{\mathbf{x}}) > \epsilon\} < \epsilon$$

(see Ho and Yeung [150]). By assuming this inequality, prove that

$$\Pr\{|H(q_{\mathbf{x}}) - H(p)| \le \epsilon\} < 1 - 2\epsilon.$$

d) Show that if  $|H(q_{\mathbf{x}}) - H(p)| \le \epsilon$  and  $|D(q_{\mathbf{x}}||p) + H(q_{\mathbf{x}}) - H(p)| \le \epsilon$ , then

$$D(q_{\mathbf{x}}||p) + |H(q_{\mathbf{x}}) - H(p)| \le 3\epsilon.$$

e) Use the results in c) and d) above and the result in c) in Problem 5 in Chapter 5 to show that

$$\Pr\{D(q_{\mathbf{x}} \| p) + |H(q_{\mathbf{x}}) - H(p)| \le \eta\} > 1 - \eta.$$

This proves Property 2 of the AEP. Property 3 of the AEP follows from Property 1 as in the proof of Theorem 5.3.

### 132 6 Strong Typicality

11. Consistency of unified typicality For any  $\eta > 0$ , the unified jointly typical set  $U^n_{[XY]\eta}$  with respect to  $p_{XY}(x, y)$  is the set of sequences  $(\mathbf{x}, \mathbf{y}) \in \mathcal{X}^n \times \mathcal{Y}^n$  such that

$$D(q_{\mathbf{x},\mathbf{y}} || p_{XY}) + |H(q_{\mathbf{x},\mathbf{y}}) - H(p_{XY})| + |H(q_{\mathbf{x}}) - H(p_X)| + |H(q_{\mathbf{y}}) - H(p_Y)| \le \eta.$$

Show that if  $(\mathbf{x}, \mathbf{y}) \in U_{[XY]\eta}^n$ , then  $\mathbf{x} \in U_{[X]\eta}^n$  and  $\mathbf{y} \in U_{[Y]\eta}^n$ .

## **Historical Notes**

Strong typicality was used by Wolfowitz [346] for proving channel coding theorems and by Berger [27] for proving the rate-distortion theorem and various results in multiterminal source coding. The method of types, a refinement of the notion of strong typicality, was systematically developed in the book by Csiszár and Körner [75]. The interpretation of the basic inequalities in Section 6.4 is a preamble to the relation between entropy and groups to be discussed in Chapter 16.

Recently, Ho and Yeung [150] introduced the notion of unified typicality which is stronger than both weak typicality and strong typicality. This notion of typicality can be applied to random variables with countable alphabets, while at the same time preserve the essential properties of strong typicality. See Problems 9, 10, and 11 for a discussion.

In all practical communication systems, when a signal is transmitted from one point to another point, the signal is inevitably contaminated by random noise, i.e., the signal received is correlated with but possibly different from the signal transmitted. We use a *noisy channel* to model such a situation. A noisy channel is a "system" which has one input terminal and one output terminal<sup>1</sup>, with the input connected to the transmission point and the output connected to the receiving point. When the signal is transmitted through the channel, it is distorted in a random way which depends on the channel characteristics. As such, the signal received may be different from the signal transmitted.

In communication engineering, we are interested in conveying messages reliably through a noisy channel at the maximum possible rate. We first look at a simple channel called the *binary symmetric channel* (BSC), which is represented by the transition diagram in Figure 7.1. In this channel both the input X and the output Y take values in the set  $\{0, 1\}$ . There is a certain probability, denoted by  $\epsilon$ , that the output is not equal to the input. That is, if the input is 0, then the output is 0 with probability  $1 - \epsilon$ , and is 1 with probability  $\epsilon$ . Likewise, if the input is 1, then the output is 1 with probability  $1 - \epsilon$ , and is 0 with probability  $\epsilon$ . The parameter  $\epsilon$  is called the *crossover* probability of the BSC.

Let  $\{A, B\}$  be the message set which contains two possible messages to be conveyed through a BSC with  $0 \le \epsilon < 0.5$ . We further assume that the two messages A and B are equally likely. If the message is A, we map it to the codeword 0, and if the message is B, we map it to the codeword 1. This is the simplest example of a *channel code*. The codeword is then transmitted through the channel. Our task is to decode the message based on the output of the channel, and an error is said to occur if the message is decoded incorrectly.

Consider

$$\Pr\{A|Y=0\} = \Pr\{X=0|Y=0\}$$
(7.1)

<sup>&</sup>lt;sup>1</sup> The discussion on noisy channels here is confined to point-to-point channels.

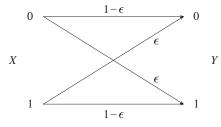


Fig. 7.1. A binary symmetric channel.

$$= \frac{\Pr\{X=0\}\Pr\{Y=0|X=0\}}{\Pr\{Y=0\}}$$
(7.2)

$$=\frac{0.5(1-\epsilon)}{\Pr\{Y=0\}}.$$
(7.3)

Since

$$\Pr\{Y = 0\} = \Pr\{Y = 1\} = 0.5 \tag{7.4}$$

by symmetry<sup>2</sup>, it follows that

$$\Pr\{A|Y=0\} = 1 - \epsilon \tag{7.5}$$

and

$$\Pr\{B|Y=0\} = 1 - \Pr\{A|Y=0\} = \epsilon.$$
(7.6)

Since  $\epsilon < 0.5$ ,

$$\Pr\{B|Y=0\} < \Pr\{A|Y=0\}.$$
(7.7)

Therefore, in order to minimize the probability of error, we decode a received 0 to the message A. By symmetry, we decode a received 1 to the message B.

An error occurs if a 0 is received and the message is B, or if a 1 is received and the message is A. Therefore, the probability of error, denoted by  $P_e$ , is given by

$$P_e = \Pr\{Y = 0\}\Pr\{B|Y = 0\} + \Pr\{Y = 1\}\Pr\{A|Y = 1\}$$
(7.8)

$$= 0.5\epsilon + 0.5\epsilon \tag{7.9}$$

$$=\epsilon,$$
 (7.10)

 $^{2}$  More explicitly,

$$\begin{aligned} \Pr\{Y=0\} &= \Pr\{A\} \Pr\{Y=0|A\} + \Pr\{B\} \Pr\{Y=0|B\} \\ &= 0.5 \, \Pr\{Y=0|X=0\} + 0.5 \, \Pr\{Y=0|X=1\} \\ &= 0.5(1-\epsilon) + 0.5\epsilon \\ &= 0.5. \end{aligned}$$

where (7.9) follows from (7.6) because

$$\Pr\{A|Y=1\} = \Pr\{B|Y=0\} = \epsilon$$
(7.11)

by symmetry.

Let us assume that  $\epsilon \neq 0$ . Then the above scheme obviously does not provide perfectly reliable communication. If we are allowed to use the channel only once, then this is already the best we can do. However, if we are allowed to use the same channel repeatedly, then we can improve the reliability by generalizing the above scheme.

We now consider the following channel code which we refer to as the binary repetition code. Let  $n \ge 1$  be an odd positive integer which is called the block length of the code. In this code, the message A is mapped to the sequence of n 0's, and the message B is mapped to the sequence of n 1's. The codeword, which consists of a sequence of either n 0's or n 1's, is transmitted through the channel in n uses. Upon receiving a sequence of n bits at the output of the channel, we use the majority vote to decode the message, i.e., if there are more 0's than 1's in the sequence, we decode the sequence to the message A, otherwise we decode the sequence to the message B. Note that the block length is chosen to be odd so that there cannot be a tie. When n = 1, this scheme reduces to the previous scheme.

For this more general scheme, we continue to denote the probability of error by  $P_e$ . Let  $N_0$  and  $N_1$  be the number of 0's and 1's in the received sequence, respectively. Clearly,

$$N_0 + N_1 = n. (7.12)$$

For large n, if the message is A, the number of 0's received is approximately equal to

$$E[N_0|A] = n(1-\epsilon) \tag{7.13}$$

and the number of 1's received is approximately equal to

$$E[N_1|A] = n\epsilon \tag{7.14}$$

with high probability by the weak law of large numbers. This implies that the probability of an error, namely the event  $\{N_0 < N_1\}$ , is small because

$$n(1-\epsilon) > n\epsilon \tag{7.15}$$

with the assumption that  $\epsilon < 0.5$ . Specifically,

$$\Pr\{\text{error}|A\} = \Pr\{N_0 < N_1|A\}$$
(7.16)

$$= \Pr\{n - N_1 < N_1 | A\}$$
(7.17)

$$= \Pr\{N_1 > 0.5n|A\} \tag{7.18}$$

$$\leq \Pr\{N_1 > (\epsilon + \phi)n|A\},\tag{7.19}$$

where

$$0 < \phi < 0.5 - \epsilon, \tag{7.20}$$

so that  $\phi$  is positive and

$$\epsilon + \phi < 0.5. \tag{7.21}$$

Note that such a  $\phi$  exists because  $\epsilon < 0.5$ . Then by the weak law of large numbers, the upper bound in (7.19) tends to 0 as  $n \to \infty$ . By symmetry,  $\Pr\{\text{error}|B\}$  also tends to 0 as  $n \to \infty$ . Therefore,

$$P_e = \Pr\{A\}\Pr\{\operatorname{error}|A\} + \Pr\{B\}\Pr\{\operatorname{error}|B\}$$
(7.22)

tends to 0 as  $n \to \infty$ . In other words, by using a long enough repetition code, we can make  $P_e$  arbitrarily small. In this sense, we say that reliable communication is achieved asymptotically.

We point out that for a BSC with  $\epsilon > 0$ , for any given transmitted sequence of length n, the probability of receiving any given sequence of length n is nonzero. It follows that for any two distinct input sequences, there is always a nonzero probability that the same output sequence is produced so that the two input sequences become indistinguishable. Therefore, except for very special channels (e.g., the BSC with  $\epsilon = 0$ ), no matter how the encoding/decoding scheme is devised, a nonzero probability of error is inevitable, and asymptotically reliable communication is the best we can hope for.

Though a rather naive approach, asymptotically reliable communication can be achieved by using the repetition code. The repetition code, however, is not without catch. For a channel code, the *rate* of the code in bit(s) per use, is defined as the ratio of the logarithm of the size of the message set in the base 2 to the block length of the code. Roughly speaking, the rate of a channel code is the average number of bits the channel code attempts to convey through the channel per use of the channel. For a binary repetition code with block length n, the rate is  $\frac{1}{n} \log 2 = \frac{1}{n}$ , which tends to 0 as  $n \to \infty$ . Thus in order to achieve asymptotic reliability by using the repetition code, we cannot communicate through the noisy channel at any positive rate!

In this chapter, we characterize the maximum rate at which information can be communicated through a *discrete memoryless channel* (DMC) with an arbitrarily small probability of error. This maximum rate, which is generally positive, is known as the *channel capacity*. Then we discuss the use of feedback in communicating through a channel, and show that feedback does not increase the capacity. At the end of the chapter, we discuss transmitting an information source through a DMC, and we show that asymptotic optimality can be achieved by separating source coding and channel coding.

### 7.1 Definition and Capacity

**Definition 7.1.** Let  $\mathcal{X}$  and  $\mathcal{Y}$  be discrete alphabets, and p(y|x) be a transition matrix from  $\mathcal{X}$  to  $\mathcal{Y}$ . A discrete channel p(y|x) is a single-input single-output

system with input random variable X taking values in  $\mathcal{X}$  and output random variable Y taking values in  $\mathcal{Y}$  such that

$$\Pr\{X = x, Y = y\} = \Pr\{X = x\}p(y|x)$$
(7.23)

for all  $(x, y) \in \mathcal{X} \times \mathcal{Y}$ .

**Remark** From (7.23), we see that if  $Pr\{X = x\} > 0$ , then

$$\Pr\{Y = y | X = x\} = \frac{\Pr\{X = x, Y = y\}}{\Pr\{X = x\}} = p(y|x).$$
(7.24)

Note that  $\Pr\{Y = y | X = x\}$  is undefined if  $\Pr\{X = x\} = 0$ . Nevertheless, (7.23) is valid for both cases.

We now present an alternative description of a discrete channel. Let  $\mathcal{X}$ and  $\mathcal{Y}$  be discrete alphabets. Let X be a random variable taking values in  $\mathcal{X}$ and p(y|x) be any transition matrix from  $\mathcal{X}$  to  $\mathcal{Y}$ . Define random variables  $Z_x$  with  $\mathcal{Z}_x = \mathcal{Y}$  for  $x \in \mathcal{X}$  such that

$$\Pr\{Z_x = y\} = p(y|x) \tag{7.25}$$

for all  $y \in \mathcal{Y}$ . We assume that  $Z_x, x \in \mathcal{X}$  are mutually independent and also independent of X. Further define the random variable

$$Z = (Z_x : x \in \mathcal{X}), \tag{7.26}$$

called the *noise variable*. Note that Z is independent of X. Now define a random variable taking values in  $\mathcal{Y}$  as

$$Y = Z_x \quad \text{if } X = x. \tag{7.27}$$

Evidently, Y is a function of X and Z. Then for  $x \in \mathcal{X}$  such that  $\Pr\{X = x\} > 0$ , we have

$$\Pr\{X = x, Y = y\} = \Pr\{X = x\}\Pr\{Y = y|X = x\}$$
(7.28)

$$= \Pr\{X = x\} \Pr\{Z_x = y | X = x\}$$
(7.29)

$$= \Pr\{X = x\} \Pr\{Z_x = y\}$$
(7.30)

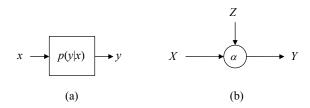
$$= \Pr\{X = x\}p(y|x), \tag{7.31}$$

i.e., (7.23) in Definition 7.1, where (7.30) follows from the assumption that  $Z_x$  is independent of  $Z_x$ . For  $x \in \mathcal{X}$  such that  $\Pr\{X = x\} = 0$ , since  $\Pr\{X = x\} = 0$  implies  $\Pr\{X = x, Y = y\} = 0$ , (7.23) continues to hold. Then by regarding X and Y as the input and output random variables, we have obtained an alternative description of the discrete channel p(y|x).

Since Y is a function of X and Z, we can write

$$Y = \alpha(X, Z). \tag{7.32}$$

Then we have the following equivalent definition for a discrete channel.



**Fig. 7.2.** Illustrations of (a) a discrete channel p(y|x) and (b) a discrete channel  $(\alpha, Z)$ .

**Definition 7.2.** Let  $\mathcal{X}$ ,  $\mathcal{Y}$ , and  $\mathcal{Z}$  be discrete alphabets. Let  $\alpha : \mathcal{X} \times \mathcal{Z} \to \mathcal{Y}$ , and Z be a random variable taking values in  $\mathcal{Z}$ , called the noise variable. A discrete channel  $(\alpha, Z)$  is a single-input single-output system with input alphabet  $\mathcal{X}$  and output alphabet  $\mathcal{Y}$ . For any input random variable X, the noise variable Z is independent of X, and the output random variable Y is given by

$$Y = \alpha(X, Z). \tag{7.33}$$

Figure 7.2 illustrates a discrete channel p(y|x) and a discrete channel  $(\alpha, Z)$ . The next definition gives the condition for the equivalence of the two specifications of a discrete channel according to Definitions 7.1 and 7.2, respectively.

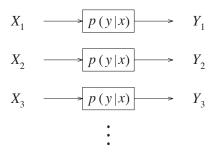
**Definition 7.3.** Two discrete channels p(y|x) and  $(\alpha, Z)$  defined on the same input alphabet  $\mathcal{X}$  and output alphabet  $\mathcal{Y}$  are equivalent if

$$\Pr\{\alpha(x, Z) = y\} = p(y|x) \tag{7.34}$$

for all x and y.

We point out that the adjective "discrete" in a discrete channel refers to the input and output alphabets of the channel being discrete. As part of a discrete-time communication system, a discrete channel can be used repeatedly at every time index  $i = 1, 2, \cdots$ . As the simplest model, we may assume that the noise for the transmission over the channel at different time indices are independent of each other. In the next definition, we will introduce the *discrete memoryless channel* (DMC) as a discrete-time extension of a discrete channel that captures this modeling assumption.

To properly formulate a DMC, we regard it as a subsystem of a discretetime stochastic system which will be referred to as "the system" in the sequel. In such a system, random variables are generated sequentially in discrete-time, and more than one random variable may be generated instantaneously but sequentially at a particular time index.



**Fig. 7.3.** An illustration of a discrete memoryless channel p(y|x).

**Definition 7.4.** A discrete memoryless channel (DMC) p(y|x) is a sequence of replicates of a generic discrete channel p(y|x). These discrete channels are indexed by a discrete-time index i, where  $i \ge 1$ , with the ith channel being available for transmission at time i. Transmission through a channel is assumed to be instantaneous. Let  $X_i$  and  $Y_i$  be respectively the input and the output of the DMC at time i, and let  $T_{i-}$  denote all the random variables that are generated in the system before  $X_i$ . The equality

$$\Pr\{Y_i = y, X_i = x, T_{i-} = t\} = \Pr\{X_i = x, T_{i-} = t\}p(y|x)$$
(7.35)

holds for all  $(x, y, t) \in \mathcal{X} \times \mathcal{Y} \times \mathcal{T}_{i-}$ .

**Remark** Similar to the remark following Definition 7.1, if  $Pr\{X_i = x, T_{i-} = t\} > 0$ , then

$$\Pr\{Y_i = y | X_i = x, T_{i-} = t\} = \frac{\Pr\{Y_i = y, X_i = x, T_{i-} = t\}}{\Pr\{X_i = x, T_{i-} = t\}}$$
(7.36)

$$= p(y|x). \tag{7.37}$$

Note that  $\Pr\{Y_i = y | X_i = x, T_{i-} = t\}$  is undefined if  $\Pr\{X_i = x, T_{i-} = t\} = 0$ . Nevertheless, (7.35) is valid for both cases.

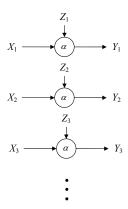
Invoking Proposition 2.5, we see from (7.35) that

$$T_{i-} \to X_i \to Y_i$$
 (7.38)

forms a Markov chain, i.e., the output of the DMC at time i is independent of all the random variables that have already been generated in the system conditioning on the input at time i. This captures the memorylessness of a DMC. Figure 7.3 is an illustration of a DMC p(y|x).

Paralleling Definition 7.2 for a discrete channel, we now present an alternative definition of a DMC.

**Definition 7.5.** A discrete memoryless channel  $(\alpha, Z)$  is a sequence of replicates of a generic discrete channel  $(\alpha, Z)$ . These discrete channels are indexed



**Fig. 7.4.** An illustration of a discrete memoryless channel  $(\alpha, Z)$ .

by a discrete-time index i, where  $i \geq 1$ , with the ith channel being available for transmission at time i. Transmission through a channel is assumed to be instantaneous. Let  $X_i$  and  $Y_i$  be respectively the input and the output of the DMC at time i, and let  $T_{i-}$  denote all the random variables that are generated in the system before  $X_i$ . The noise variable  $Z_i$  for the transmission at time i is a copy of the generic noise variable Z, and it is independent of  $(X_i, T_{i-})$ . The output of the DMC at time i is given by

$$Y_i = \alpha(X_i, Z_i). \tag{7.39}$$

Figure 7.4 is an illustration of a discrete memoryless channel  $(\alpha, Z)$ . We now show the equivalence between Definitions 7.4 and 7.5. Suppose the generic discrete channel p(y|x) in Definition 7.4 is equivalent to the generic discrete channel  $(\alpha, Z)$  in Definition 7.5, i.e., (7.34) holds. Consider the DMC  $(\alpha, Z)$ in Definition 7.5. We will show that it is equivalent to the DMC p(y|x) in Definition 7.3. For the DMC  $(\alpha, Z)$ , consider

$$0 \le I(T_{i-}; Y_i | X_i) \tag{7.40}$$

$$\leq I(T_{i-}; Y_i, X_i, Z_i | X_i) \tag{7.41}$$

$$= I(T_{i-}; X_i, Z_i | X_i)$$
(7.42)

$$= I(T_{i-}; Z_i | X_i) \tag{7.43}$$

$$= 0,$$
 (7.44)

where the first equality follows from (7.39) and the last equality follows from the assumption that  $Z_i$  is independent of  $(X_i, T_{i-})$ . Therefore,

$$I(T_{i-};Y_i|X_i) = 0, (7.45)$$

or  $T_{i-} \to X_i \to Y_i$  forms a Markov chain. It remains to establish (7.35) for all  $(x, y, t) \in \mathcal{X} \times \mathcal{Y} \times \mathcal{T}_{i-}$ . For  $x \in \mathcal{X}$  such that  $\Pr\{X_i = x\} = 0$ , both  $\Pr\{Y_i = y, X_i = x, T_{i-} = t\}$  and  $\Pr\{X_i = x, T_{i-} = t\}$  vanish because they are upper bounded by  $\Pr\{X_i = x\}$ . Therefore (7.35) holds. For  $x \in \mathcal{X}$  such that  $\Pr\{X_i = x\} > 0$ ,

$$\Pr\{Y_i = y, X_i = x, T_{i-} = t\}$$

$$\stackrel{a)}{=} \Pr\{X_i = x, T_{i-} = t\} \Pr\{Y_i = y | X_i = x\}$$
(7.46)

<sup>b)</sup> 
$$= \Pr\{X_i = x, T_{i-} = t\}\Pr\{\alpha(X_i, Z_i) = y | X_i = x\}$$
 (7.47)

$$= \Pr\{X_i = x, T_{i-} = t\} \Pr\{\alpha(x, Z_i) = y | X_i = x\}$$
(7.48)

$$\stackrel{(i)}{=} \Pr\{X_i = x, T_{i-} = t\} \Pr\{\alpha(x, Z_i) = y\}$$
(7.49)

$$\stackrel{_{I}}{=} \Pr\{X_i = x, T_{i-} = t\} \Pr\{\alpha(x, Z) = y\}$$
(7.50)

$$\stackrel{e}{=} \Pr\{X_i = x, T_{i-} = t\} p(y|x), \tag{7.51}$$

where

- a) follows from the Markov chain  $T_{i-} \to X_i \to Y_i$ ;
- b) follows from (7.39);

 $\alpha$ 

- c) follows from Definition 7.5 that  $Z_i$  is independent of  $X_i$ ;
- d) follows from Definition 7.5 that  $Z_i$  and the generic noise variable Z have the same distribution;
- e) follows from (7.34).

Hence, (7.35) holds for all  $(x, y, t) \in \mathcal{X} \times \mathcal{Y} \times \mathcal{T}_{i-}$ , proving that the DMC  $(\alpha, Z)$  in Definition 7.4 is equivalent to the DMC (p(y|x)) in Definition 7.5.

Definition 7.5 renders the following physical conceptualization of a DMC. The DMC can be regarded as a "box" which has only two terminals, the input and the output. The box perfectly shields its contents from the rest of the system. At time *i*, upon the transmission of the input  $X_i$ , the noise variable  $Z_i$  is generated inside the box according to the distribution of the generic noise variable Z. Since the box is perfectly shielded, the generation of the  $Z_i$  is independent of  $X_i$  and any other random variable that has already been generated in the system. Then the function  $\alpha$  is applied to  $(X_i, Z_i)$  to produce the output  $Y_i$ .

In the rest of the section, we will define the capacity of a DMC and discuss some of its basic properties. The capacities of two simple DMCs will also be evaluated explicitly. To keep our discussion simple, we will assume that the alphabets  $\mathcal{X}$  and  $\mathcal{Y}$  are finite.

**Definition 7.6.** The capacity of a discrete memoryless channel p(y|x) is defined as

$$C = \max_{p(x)} I(X;Y),$$
 (7.52)

where X and Y are respectively the input and the output of the generic discrete channel, and the maximum is taken over all input distributions p(x).

From the above definition, we see that

$$C \ge 0 \tag{7.53}$$

because

$$I(X;Y) \ge 0 \tag{7.54}$$

for all input distributions p(x). By Theorem 2.43, we have

$$C = \max_{p(x)} I(X;Y) \le \max_{p(x)} H(X) = \log |\mathcal{X}|.$$
 (7.55)

Likewise, we have

$$C \le \log |\mathcal{Y}|. \tag{7.56}$$

Therefore,

$$C \le \min(\log |\mathcal{X}|, \log |\mathcal{Y}|). \tag{7.57}$$

Since I(X;Y) is a continuous functional of p(x) and the set of all p(x) is a compact set (i.e., closed and bounded) in  $\Re^{|\mathcal{X}|}$ , the maximum value of I(X;Y) can be attained<sup>3</sup>. This justifies taking the maximum rather than the supremum in the definition of channel capacity in (7.52).

We will prove subsequently that C is in fact the maximum rate at which information can be communicated reliably through a DMC. We first give some examples of DMC's for which the capacities can be obtained in closed form. In the following, X and Y denote, respectively, the input and the output of the generic discrete channel, and all logarithms are in the base 2.

Example 7.7 (Binary Symmetric Channel). The binary symmetric channel (BSC) has been shown in Figure 7.1. Alternatively, a BSC can be represented by Figure 7.5. Here, Z is a binary random variable representing the noise of

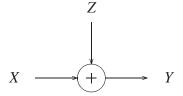


Fig. 7.5. An alternative representation for a binary symmetric channel.

the channel, with

$$\Pr\{Z=0\} = 1 - \epsilon \text{ and } \Pr\{Z=1\} = \epsilon,$$
 (7.58)

and Z is independent of X. Then

<sup>&</sup>lt;sup>3</sup> The assumption that  $\mathcal{X}$  is finite is essential in this argument.

7.1 Definition and Capacity 143

$$Y = X + Z \mod 2. \tag{7.59}$$

This representation for a BSC is in the form prescribed by Definition 7.2.

In order to determine the capacity of the BSC, we first bound I(X;Y) as follows.

$$I(X;Y) = H(Y) - H(Y|X)$$
(7.60)

$$= H(Y) - \sum_{x} p(x)H(Y|X=x)$$
(7.61)

$$=H(Y) - \sum_{x} p(x)h_b(\epsilon)$$
(7.62)

$$=H(Y)-h_b(\epsilon) \tag{7.63}$$

$$\leq 1 - h_b(\epsilon),\tag{7.64}$$

where we have used  $h_b$  to denote the binary entropy function in the base 2. In order to achieve this upper bound, we have to make H(Y) = 1, i.e., the output distribution of the BSC is uniform. This can be done by letting p(x)be the uniform distribution on  $\{0, 1\}$ . Therefore, the upper bound on I(X; Y)can be achieved, and we conclude that

$$C = 1 - h_b(\epsilon)$$
 bit per use. (7.65)

Figure 7.6 is a plot of the capacity C versus the crossover probability  $\epsilon$ . We

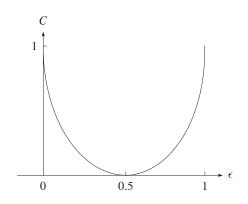


Fig. 7.6. The capacity of a binary symmetric channel.

see from the plot that C attains the maximum value 1 when  $\epsilon = 0$  or  $\epsilon = 1$ , and attains the minimum value 0 when  $\epsilon = 0.5$ . When  $\epsilon = 0$ , it is easy to see that C = 1 is the maximum rate at which information can be communicated through the channel reliably. This can be achieved simply by transmitting unencoded bits through the channel, and no decoding is necessary because all

the bits are received unchanged. When  $\epsilon = 1$ , the same can be achieved with the additional decoding step which complements all the received bits. By doing so, the bits transmitted through the channel can be recovered without error. Thus from the communication point of view, for binary channels, a channel which never makes error and a channel which always makes errors are equally good. When  $\epsilon = 0.5$ , the channel output is independent of the channel input. Therefore, no information can possibly be communicated through the channel.

Example 7.8 (Binary Erasure Channel). The binary erasure channel is shown in Figure 7.7. In this channel, the input alphabet is  $\{0, 1\}$ , while the output

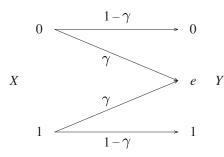


Fig. 7.7. A binary erasure channel.

alphabet is  $\{0, 1, e\}$ . With probability  $\gamma$ , the erasure symbol e is produced at the output, which means that the input bit is lost; otherwise the input bit is reproduced at the output without error. The parameter  $\gamma$  is called the erasure probability.

To determine the capacity of this channel, we first consider

$$C = \max_{r(r)} I(X;Y) \tag{7.66}$$

$$= \max_{p(x)} (H(Y) - H(Y|X))$$
(7.67)

$$= \max_{p(x)} H(Y) - h_b(\gamma). \tag{7.68}$$

Thus we only have to maximize H(Y). To this end, let

$$\Pr\{X = 0\} = a \tag{7.69}$$

and define a binary random variable E by

$$E = \begin{cases} 0 \text{ if } Y \neq e\\ 1 \text{ if } Y = e. \end{cases}$$
(7.70)

The random variable E indicates whether an erasure has occurred, and it is a function of Y. Then

7.2 The Channel Coding Theorem 145

$$H(Y) = H(Y, E) \tag{7.71}$$

$$= H(E) + H(Y|E)$$
 (7.72)

$$= h_b(\gamma) + (1 - \gamma)h_b(a).$$
 (7.73)

Hence,

$$C = \max_{p(x)} H(Y) - h_b(\gamma) \tag{7.74}$$

$$= \max_{a} [h_b(\gamma) + (1 - \gamma)h_b(a)] - h_b(\gamma)$$
(7.75)

$$= (1 - \gamma) \max_{a} h_b(a) \tag{7.76}$$

$$=(1-\gamma),\tag{7.77}$$

where capacity is achieved by letting a = 0.5, i.e., the input distribution is uniform.

It is in general not possible to obtain the capacity of a DMC in closed form, and we have to resort to numerical computation. In Chapter 9 we will discuss the Blahut-Arimoto algorithm for computing channel capacity.

## 7.2 The Channel Coding Theorem

We will justify the definition of the capacity of a DMC by the proving the *channel coding theorem*. This theorem, which consists of two parts, will be formally stated at the end of the section. The direct part of the theorem asserts that information can be communicated through a DMC with an arbitrarily small probability of error at any rate less than the channel capacity. Here it is assumed that the decoder knows the transition matrix of the DMC. The converse part of the theorem asserts that if information is communicated through a DMC at a rate higher than the capacity, then the probability of error is bounded away from zero. For better appreciation of the definition of channel capacity, we will first prove the converse part in Section 7.3 and then prove the direct part in Section 7.4.

**Definition 7.9.** An (n, M) code for a discrete memoryless channel with input alphabet  $\mathcal{X}$  and output alphabet  $\mathcal{Y}$  is defined by an encoding function

$$f: \{1, 2, \cdots, M\} \to \mathcal{X}^n \tag{7.78}$$

and a decoding function

$$g: \mathcal{Y}^n \to \{1, 2, \cdots, M\}. \tag{7.79}$$

The set  $\{1, 2, \dots, M\}$ , denoted by  $\mathcal{W}$ , is called the message set. The sequences  $f(1), f(2), \dots, f(M)$  in  $\mathcal{X}^n$  are called codewords, and the set of codewords is called the codebook.

In order to distinguish a channel code as defined above from a channel code with feedback which will be discussed in Section 7.6, we will refer to the former as a channel code without feedback.

We assume that a message W is randomly chosen from the message set  $\mathcal{W}$  according to the uniform distribution. Therefore,

$$H(W) = \log M. \tag{7.80}$$

With respect to a channel code for a DMC, we let

٦

$$\mathbf{X} = (X_1, X_2, \cdots, X_n) \tag{7.81}$$

and

$$\mathbf{Y} = (Y_1, Y_2, \cdots, Y_n) \tag{7.82}$$

be the input sequence and the output sequence of the channel, respectively. Evidently,

$$\mathbf{X} = f(W). \tag{7.83}$$

We also let

$$\hat{W} = g(\mathbf{Y}) \tag{7.84}$$

be the estimate on the message W by the decoder. Figure 7.8 is the block diagram for a channel code.

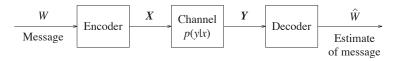


Fig. 7.8. A channel code with block length n.

**Definition 7.10.** For all  $1 \le w \le M$ , let

$$\lambda_w = \Pr\{\hat{W} \neq w | W = w\} = \sum_{\mathbf{y} \in \mathcal{Y}^n : g(\mathbf{y}) \neq w} \Pr\{\mathbf{Y} = \mathbf{y} | \mathbf{X} = f(w)\}$$
(7.85)

be the conditional probability of error given that the message is w.

We now define two performance measures for a channel code.

**Definition 7.11.** The maximal probability of error of an (n, M) code is defined as

$$\lambda_{max} = \max_{w} \lambda_w. \tag{7.86}$$

**Definition 7.12.** The average probability of error of an (n, M) code is defined as

$$P_e = \Pr\{W \neq W\}. \tag{7.87}$$

From the definition of  $P_e$ , we have

$$P_e = \Pr\{\hat{W} \neq W\} \tag{7.88}$$

$$= \sum_{w} \Pr\{W = w\} \Pr\{\hat{W} \neq W | W = w\}$$
(7.89)

$$=\sum_{w}\frac{1}{M}\Pr\{\hat{W}\neq w|W=w\}$$
(7.90)

$$=\frac{1}{M}\sum_{w}\lambda_{w},\tag{7.91}$$

i.e.,  $P_e$  is the arithmetic mean of  $\lambda_w$ ,  $1 \leq w \leq M$ . It then follows that

$$P_e \le \lambda_{max}.\tag{7.92}$$

In fact, it can be readily seen that this inequality remains valid even without the assumption that W distributes uniformly on the message set  $\mathcal{W}$ .

**Definition 7.13.** The rate of an (n, M) channel code is  $n^{-1} \log M$  in bits per use.

**Definition 7.14.** A rate R is asymptotically achievable for a discrete memoryless channel if for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M)code such that

$$\frac{1}{n}\log M > R - \epsilon \tag{7.93}$$

and

$$\lambda_{max} < \epsilon. \tag{7.94}$$

For brevity, an asymptotically achievable rate will be referred to as an achievable rate.

In other words, a rate R is achievable if there exists a sequence of codes whose rates approach R and whose probabilities of error approach zero. We end this section by stating the channel coding theorem, which gives a characterization of all achievable rates. This theorem will be proved in the next two sections.

**Theorem 7.15 (Channel Coding Theorem).** A rate R is achievable for a discrete memoryless channel if and only if  $R \leq C$ , the capacity of the channel.

## 7.3 The Converse

Let us consider a channel code with block length n. The random variables involved in this code are W,  $X_i$  and  $Y_i$  for  $1 \leq i \leq n$ , and  $\hat{W}$ . We see

from the definition of a channel code in Definition 7.9 that all the random variables are generated sequentially according to some deterministic or probabilistic rules. Specifically, the random variables are generated in the order  $W, X_1, Y_1, X_2, Y_2, \dots, X_n, Y_n, \hat{W}$ . The generation of these random variables can be represented by the dependency graph<sup>4</sup> in Figure 7.9. In this graph, a

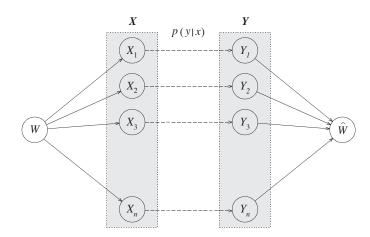


Fig. 7.9. The dependency graph for a channel code without feedback.

node represents a random variable. If there is a (directed) edge from node X to node Y, then node X is called a *parent* of node Y. We further distinguish a *solid* edge and a *dotted* edge: a solid edge represents functional (deterministic) dependency, while a dotted edge represents the probabilistic dependency induced by the transition matrix p(y|x) of the generic discrete channel. For a node X, its parent nodes represent all the random variables on which random variable X depends when it is generated.

We now explain the specific structure of the dependency graph. First,  $X_i$  is a function of W, so each  $X_i$  is connected to W by a solid edge. According to Definition 7.4,

$$T_{i-} = (W, X_1, Y_1, \cdots, X_{i-1}, Y_{i-1}).$$
(7.95)

By (7.35), the Markov chain

$$(W, X_1, Y_1, \cdots, X_{i-1}, Y_{i-1}) \to X_i \to Y_i$$
 (7.96)

prevails. Therefore, the generation of  $Y_i$  depends only on  $X_i$  and not on  $W, X_1, Y_1, \dots, X_{i-1}, Y_{i-1}$ . So,  $Y_i$  is connected to  $X_i$  by a dotted edge representing the discrete channel p(y|x) at time i, and there is no connection between  $Y_i$  and any of the nodes  $W, X_1, Y_1, \dots, X_{i-1}, Y_{i-1}$ . Finally,  $\hat{W}$  is a function of  $Y_1, Y_2, \dots, Y_n$ , so  $\hat{W}$  is connected to each  $Y_i$  by a solid edge.

 $<sup>^{4}</sup>$  A dependency graph can be regarded as a Bayesian network [258].

We will use q to denote the joint distribution of these random variables as well as all the marginals, and let  $x_i$  denote the *i*th component of a sequence  $\mathbf{x}$ . From the dependency graph, we see that for all  $(w, \mathbf{x}, \mathbf{y}, \hat{w}) \in \mathcal{W} \times \mathcal{X}^n \times \mathcal{Y}^n \times \hat{\mathcal{W}}$ such that  $q(\mathbf{x}) > 0$  and  $q(\mathbf{y}) > 0$ ,

$$q(w, \mathbf{x}, \mathbf{y}\,\hat{w}) = q(w) \left(\prod_{i=1}^{n} q(x_i|w)\right) \left(\prod_{i=1}^{n} p(y_i|x_i)\right) q(\hat{w}|\mathbf{y}).$$
(7.97)

Note that q(w) > 0 for all w so that  $q(x_i|w)$  are well-defined, and  $q(x_i|w)$  and  $q(\hat{w}|\mathbf{y})$  are both deterministic. Denote the set of nodes  $X_1, X_2, \dots, X_n$  by  $\mathbf{X}$  and the set of nodes  $Y_1, Y_2, \dots, Y_n$  by  $\mathbf{Y}$ . We notice the following structure in the dependency graph: all the edges from W end in  $\mathbf{X}$ , all the edges from  $\mathbf{X}$  end in  $\mathbf{Y}$ , and all the edges from  $\mathbf{Y}$  end in  $\hat{W}$ . This suggests that the random variables  $W, \mathbf{X}, \mathbf{Y}$ , and  $\hat{W}$  form the Markov chain

$$W \to \mathbf{X} \to \mathbf{Y} \to \hat{W}.$$
 (7.98)

The validity of this Markov chain can be formally justified by applying Proposition 2.9 to (7.97), so that for all  $(w, \mathbf{x}, \mathbf{y}, \hat{w}) \in \mathcal{W} \times \mathcal{X}^n \times \mathcal{Y}^n \times \hat{\mathcal{W}}$  such that  $q(\mathbf{x}) > 0$  and  $q(\mathbf{y}) > 0$ , we can write

$$q(w, \mathbf{x}, \mathbf{y}, \hat{w}) = q(w)q(\mathbf{x}|w)q(\mathbf{y}|\mathbf{x})q(\hat{w}|\mathbf{y}).$$
(7.99)

Now  $q(\mathbf{x}, \mathbf{y})$  is obtained by summing over all w and  $\hat{w}$  in (7.97), and  $q(\mathbf{x})$  is obtained by further summing over all  $\mathbf{y}$ . After some straightforward algebra and using

$$q(\mathbf{y}|\mathbf{x}) = \frac{q(\mathbf{x}, \mathbf{y})}{q(\mathbf{x})}$$
(7.100)

for all **x** such that  $q(\mathbf{x}) > 0$ , we obtain

$$q(\mathbf{y}|\mathbf{x}) = \prod_{i=1}^{n} p(y_i|x_i).$$
 (7.101)

The Markov chain in (7.98) and the relation in (7.101) are apparent from the setup of the problem, and the above justification may seem superfluous. However, the methodology developed here is necessary for handling the more delicate situation which arises when the channel is used with feedback. This will be discussed in Section 7.6.

Consider a channel code whose probability of error is arbitrarily small. Since  $W, \mathbf{X}, \mathbf{Y}$ , and  $\hat{W}$  form the Markov chain in (7.98), the information diagram for these four random variables is as shown in Figure 7.10. Moreover, **X** is a function of W, and  $\hat{W}$  is a function of **Y**. These two relations are equivalent to

$$H(\mathbf{X}|W) = 0, \tag{7.102}$$

and

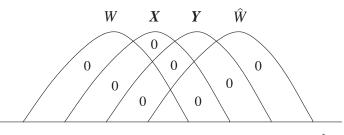


Fig. 7.10. The information diagram for  $W \to \mathbf{X} \to \mathbf{Y} \to \hat{W}$ .

$$H(\hat{W}|\mathbf{Y}) = 0, \tag{7.103}$$

respectively. Since the probability of error is arbitrarily small, W and  $\hat{W}$  are essentially identical. To gain insight into the problem, we assume for the time being that W and  $\hat{W}$  are equivalent, so that

$$H(\hat{W}|W) = H(W|\hat{W}) = 0.$$
(7.104)

Since the *I*-Measure  $\mu^*$  for a Markov chain is nonnegative, the constraints in (7.102) to (7.104) imply that  $\mu^*$  vanishes on all the atoms in Figure 7.10 marked with a '0.' Immediately, we see that

$$H(W) = I(\mathbf{X}; \mathbf{Y}). \tag{7.105}$$

That is, the amount of information conveyed through the channel is essentially the mutual information between the input sequence and the output sequence of the channel.

For a single transmission, we see from the definition of channel capacity that the mutual information between the input and the output cannot exceed the capacity of the channel, i.e., for all  $1 \le i \le n$ ,

$$I(X_i; Y_i) \le C. \tag{7.106}$$

Summing i from 1 to n, we have

$$\sum_{i=1}^{n} I(X_i; Y_i) \le nC.$$
(7.107)

Upon establishing in the next lemma that

$$I(\mathbf{X}; \mathbf{Y}) \le \sum_{i=1}^{n} I(X_i; Y_i),$$
 (7.108)

the converse of the channel coding theorem then follows from

7.3 The Converse 151

$$\frac{1}{n}\log M = \frac{1}{n}H(W) \tag{7.109}$$

$$=\frac{1}{n}I(\mathbf{X};\mathbf{Y}) \tag{7.110}$$

$$\leq \frac{1}{n} \sum_{i=1}^{n} I(X_i; Y_i) \tag{7.111}$$

$$\leq C. \tag{7.112}$$

**Lemma 7.16.** For a discrete memoryless channel used with a channel code without feedback, for any  $n \ge 1$ ,

$$I(\mathbf{X}; \mathbf{Y}) \le \sum_{i=1}^{n} I(X_i; Y_i),$$
 (7.113)

where  $X_i$  and  $Y_i$  are, respectively, the input and the output of the channel at time *i*.

*Proof.* For any  $(\mathbf{x}, \mathbf{y}) \in \mathcal{X}^n \times \mathcal{Y}^n$ , if  $q(\mathbf{x}, \mathbf{y}) > 0$ , then  $q(\mathbf{x}) > 0$  and (7.101) holds. Therefore,

$$q(\mathbf{Y}|\mathbf{X}) = \prod_{i=1}^{n} p(Y_i|X_i)$$
(7.114)

holds for all  $(\mathbf{x}, \mathbf{y})$  in the support of  $q(\mathbf{x}, \mathbf{y})$ . Then

$$-E\log q(\mathbf{Y}|\mathbf{X}) = -E\log\prod_{i=1}^{n} p(Y_i|X_i) = -\sum_{i=1}^{n} E\log p(Y_i|X_i), \quad (7.115)$$

or

$$H(\mathbf{Y}|\mathbf{X}) = \sum_{i=1}^{n} H(Y_i|X_i).$$
 (7.116)

Hence,

$$I(\mathbf{X}; \mathbf{Y}) = H(\mathbf{Y}) - H(\mathbf{Y}|\mathbf{X})$$
(7.117)

$$\leq \sum_{i=1}^{n} H(Y_i) - \sum_{i=1}^{n} H(Y_i|X_i)$$
(7.118)

$$=\sum_{i=1}^{n} I(X_i; Y_i).$$
(7.119)

The lemma is proved.  $\Box$ 

We now formally prove the converse of the channel coding theorem. Let R be an achievable rate, i.e., for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M) code such that

$$\frac{1}{n}\log M > R - \epsilon \tag{7.120}$$

and

$$\lambda_{max} < \epsilon. \tag{7.121}$$

Consider

$$\log M \stackrel{a)}{=} H(W) \tag{7.122}$$

$$= H(W|W) + I(W;W)$$
(7.123)

$$\leq H(W|\tilde{W}) + I(\mathbf{X};\mathbf{Y})$$
(7.124)

$$\stackrel{c)}{\leq} H(W|\hat{W}) + \sum_{i=1}^{n} I(X_i; Y_i)$$
(7.125)

$$\stackrel{d)}{\leq} H(W|\hat{W}) + nC, \tag{7.126}$$

where

- a) follows from (7.80);
- b) follows from the data processing theorem since  $W \to \mathbf{X} \to \mathbf{Y} \to \hat{W}$ ;
- c) follows from Lemma 7.16;
- d) follows from (7.107).

From (7.87) and Fano's inequality (cf. Corollary 2.48), we have

$$H(W|W) < 1 + P_e \log M.$$
 (7.127)

Therefore, from (7.126),

$$\log M < 1 + P_e \log M + nC \tag{7.128}$$

$$\leq 1 + \lambda_{max} \log M + nC \tag{7.129}$$

$$< 1 + \epsilon \log M + nC, \tag{7.130}$$

where we have used (7.92) and (7.121), respectively, to obtain the last two inequalities. Dividing by n and rearranging the terms, we have

$$\frac{1}{n}\log M < \frac{\frac{1}{n} + C}{1 - \epsilon},\tag{7.131}$$

and from (7.120), we obtain

$$R - \epsilon < \frac{\frac{1}{n} + C}{1 - \epsilon}.\tag{7.132}$$

For any  $\epsilon > 0$ , the above inequality holds for all sufficiently large n. Letting  $n \to \infty$  and then  $\epsilon \to 0$ , we conclude that

$$R \le C. \tag{7.133}$$

This completes the proof for the converse of the channel coding theorem.

From the above proof, we can obtain an asymptotic bound on  $P_e$  when the rate of the code  $\frac{1}{n} \log M$  is greater than C. Consider (7.128) and obtain

$$P_e \ge 1 - \frac{1 + nC}{\log M} = 1 - \frac{\frac{1}{n} + C}{\frac{1}{n} \log M}.$$
(7.134)

Then

$$P_e \ge 1 - \frac{\frac{1}{n} + C}{\frac{1}{n} \log M} \approx 1 - \frac{C}{\frac{1}{n} \log M} \tag{7.135}$$

when n is large. This asymptotic bound on  $P_e$ , which is strictly positive if  $\frac{1}{n} \log M > C$ , is illustrated in Figure 7.11.

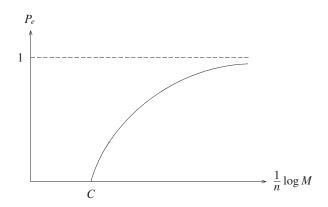


Fig. 7.11. An asymptotic upper bound on  $P_e$ .

In fact, the lower bound in (7.134) implies that  $P_e > 0$  for all n if  $\frac{1}{n} \log M > C$  because if  $P_e^{(n_0)} = 0$  for some  $n_0$ , then for all  $k \ge 1$ , by concatenating k copies of the code, we obtain a code with the same rate and block length equal to  $kn_0$  such that  $P_e^{(kn_0)} = 0$ , which is a contradiction to our conclusion that  $P_e > 0$  when n is large. Therefore, if we use a code whose rate is greater than the channel capacity, the probability of error is non-zero for all block lengths.

The converse of the channel coding theorem we have proved is called the *weak converse*. A stronger version of this result called the *strong converse* can be proved, which says that  $P_e \to 1$  as  $n \to \infty$  if there exists an  $\epsilon > 0$  such that  $\frac{1}{n} \log M \ge C + \epsilon$  for all n.

## 7.4 Achievability

We have shown in the last section that the channel capacity C is an upper bound on all achievable rates for a DMC. In this section, we show that the rate C is achievable, which implies that any rate  $R \leq C$  is achievable.

Consider a DMC p(y|x), and denote the input and the output of the generic discrete channel by X and Y, respectively. For every input distribution p(x), we will prove that the rate I(X;Y) is achievable by showing for large n the existence of a channel code such that

- 1. the rate of the code is arbitrarily close to I(X;Y);
- 2. the maximal probability of error  $\lambda_{max}$  is arbitrarily small.

Then by choosing the input distribution p(x) to be one which achieves the channel capacity, i.e., I(X;Y) = C, we conclude that the rate C is achievable.

Before we prove the achievability of the channel capacity, we first prove the following lemma.

**Lemma 7.17.** Let  $(\mathbf{X}', \mathbf{Y}')$  be *n* i.i.d. copies of a pair of generic random variables (X', Y'), where X' and Y' are independent and have the same marginal distributions as X and Y, respectively. Then

$$\Pr\{(\mathbf{X}', \mathbf{Y}') \in T^n_{[XY]\delta}\} \le 2^{-n(I(X;Y)-\tau)},$$
(7.136)

where  $\tau \to 0$  and  $\delta \to 0$ .

Proof. Consider

$$\Pr\{(\mathbf{X}', \mathbf{Y}') \in T^n_{[XY]\delta}\} = \sum_{(\mathbf{x}, \mathbf{y}) \in T^n_{[XY]\delta}} p(\mathbf{x}) p(\mathbf{y}).$$
(7.137)

By the consistency of strong typicality, for  $(\mathbf{x}, \mathbf{y}) \in T^n_{[XY]\delta}$ ,  $\mathbf{x} \in T^n_{[X]\delta}$  and  $\mathbf{y} \in T^n_{[Y]\delta}$ . By the strong AEP, all the  $p(\mathbf{x})$  and  $p(\mathbf{y})$  in the above summation satisfy

$$p(\mathbf{x}) \le 2^{-n(H(X)-\eta)}$$
 (7.138)

and

$$p(\mathbf{y}) \le 2^{-n(H(Y)-\zeta)},$$
(7.139)

where  $\eta, \zeta \to 0$  as  $\delta \to 0$ . Again by the strong JAEP,

$$|T^n_{[XY]\delta}| \le 2^{n(H(X,Y)+\xi)},\tag{7.140}$$

where  $\xi \to 0$  as  $\delta \to 0$ . Then from (7.137), we have

$$\Pr\{(\mathbf{X}', \mathbf{Y}') \in T^{n}_{[XY]\delta}\} < 2^{n(H(X,Y)+\xi)} \cdot 2^{-n(H(X)-\eta)} \cdot 2^{-n(H(Y)-\zeta)}$$
(7.141)

$$= 2^{-n(H(X)+H(Y)-H(X,Y)-\xi-\eta-\zeta)}$$
(7.142)

$$=2^{-n(I(X;Y)-\xi-\eta-\zeta)}$$
(7.143)

$$=2^{-n(I(X;Y)-\tau)}, (7.144)$$

where

$$\tau = \xi + \eta + \zeta \to 0 \tag{7.145}$$

as  $\delta \to 0$ . The lemma is proved.  $\Box$ 

Fix any  $\epsilon > 0$  and let  $\delta$  be a small positive quantity to be specified later. Toward proving the existence of a desired code, we fix an input distribution p(x) for the generic discrete channel p(y|x), and let M be an *even* integer satisfying

$$I(X;Y) - \frac{\epsilon}{2} < \frac{1}{n} \log M < I(X;Y) - \frac{\epsilon}{4},$$
 (7.146)

where n is sufficiently large. We now describe a *random coding scheme* in the following steps:

- 1. Construct the codebook C of an (n, M) code randomly by generating M codewords in  $\mathcal{X}^n$  independently and identically according to  $p(x)^n$ . Denote these codewords by  $\tilde{\mathbf{X}}(1), \tilde{\mathbf{X}}(2), \dots, \tilde{\mathbf{X}}(M)$ .
- 2. Reveal the codebook  ${\mathcal C}$  to both the encoder and the decoder.
- 3. A message W is chosen from  $\mathcal{W}$  according to the uniform distribution.
- 4. The sequence  $\mathbf{X} = \mathbf{X}(W)$ , namely the Wth codeword in the codebook  $\mathcal{C}$ , is transmitted through the channel.
- 5. The channel outputs a sequence  $\mathbf{Y}$  according to

$$\Pr\{\mathbf{Y} = \mathbf{y} | \tilde{\mathbf{X}}(W) = \mathbf{x}\} = \prod_{i=1}^{n} p(y_i | x_i)$$
(7.147)

(cf. (7.101) and Remark 1 in Section 7.6).

6. The sequence  $\mathbf{Y}$  is decoded to the message w if  $(\tilde{\mathbf{X}}(w), \mathbf{Y}) \in T^n_{[XY]\delta}$  and there does not exists  $w' \neq w$  such that  $(\tilde{\mathbf{X}}(w'), \mathbf{Y}) \in T^n_{[XY]\delta}$ . Otherwise,  $\mathbf{Y}$  is decoded to a constant message in  $\mathcal{W}$ . Denote by  $\hat{W}$  the message to which  $\mathbf{Y}$  is decoded.

**Remark 1** There are a total of  $|\mathcal{X}|^{Mn}$  possible codebooks which can be constructed in Step 1 of the random coding scheme, where we regard two codebooks whose sets of codewords are permutations of each other as two different codebooks.

**Remark 2** Strong typicality is used in defining the decoding function in Step 6. This is made possible by the assumption that the alphabets  $\mathcal{X}$  and  $\mathcal{Y}$  are finite.

We now analyze the performance of this random coding scheme. Let

$$Err = \{\hat{W} \neq W\} \tag{7.148}$$

be the event of a decoding error. In the following, we analyze  $\Pr{Err}$ , the probability of a decoding error for the random code constructed above. For all  $1 \le w \le M$ , define the event

$$E_w = \{ (\tilde{\mathbf{X}}(w), \mathbf{Y}) \in T^n_{[XY]\delta} \}.$$
(7.149)

Now

$$\Pr\{Err\} = \sum_{w=1}^{M} \Pr\{Err|W = w\} \Pr\{W = w\}.$$
 (7.150)

Since  $\Pr\{Err|W = w\}$  are identical for all w by symmetry in the code construction, we have

$$\Pr\{Err\} = \Pr\{Err|W = 1\} \sum_{w=1}^{M} \Pr\{W = w\}$$
(7.151)

$$= \Pr\{Err|W = 1\}, \tag{7.152}$$

i.e., we can assume without loss of generality that the message 1 is chosen. Then decoding is correct if the received sequence  $\mathbf{Y}$  is decoded to the message 1. This is the case if  $E_1$  occurs but  $E_w$  does not occur for all  $2 \le w \le M$ . It follows that<sup>5</sup>

$$\Pr\{Err^{c}|W=1\} \ge \Pr\{E_{1} \cap E_{2}^{c} \cap E_{3}^{c} \cap \dots \cap E_{M}^{c}|W=1\},$$
(7.153)

which implies

$$Pr\{Err|W = 1\} = 1 - Pr\{Err^{c}|W = 1\}$$
(7.154)

$$\leq 1 - \Pr\{E_1 \cap E_2^c \cap E_3^c \cap \dots \cap E_M^c | W = 1\}$$
(7.155)

$$= \Pr\{(E_1 \cap E_2^c \cap E_3^c \cap \dots \cap E_M^c)^c | W = 1\}$$
(7.156)

$$= \Pr\{E_1^c \cup E_2 \cup E_3 \cup \dots \cup E_M | W = 1\}.$$
 (7.157)

By the union bound, we have

$$\Pr\{Err|W=1\} \le \Pr\{E_1^c|W=1\} + \sum_{w=2}^M \Pr\{E_w|W=1\}.$$
 (7.158)

First, conditioning on  $\{W = 1\}$ ,  $(\tilde{\mathbf{X}}(1), \mathbf{Y})$  are *n* i.i.d. copies of the pair of generic random variables (X, Y). By the strong JAEP (Theorem 6.9), for any  $\nu > 0$ ,

$$\Pr\{E_1^c | W = 1\} = \Pr\{(\tilde{\mathbf{X}}(1), \mathbf{Y}) \notin T_{[XY]\delta}^n | W = 1\} < \nu$$
(7.159)

for sufficiently large n. This gives an upper bound on the first term on the right hand side of (7.158).

<sup>&</sup>lt;sup>5</sup> If  $E_1$  does not occur or  $E_w$  occurs for some  $1 \le w \le M$ , the received sequence

 $<sup>\</sup>mathbf{Y}$  is decoded to the constant message, which may happen to be the message 1. Therefore, the inequality in (7.153) is not an equality in general.

Second, conditioning on  $\{W = 1\}$ , for  $2 \leq w \leq M$ ,  $(\hat{\mathbf{X}}(w), \mathbf{Y})$  are *n* i.i.d. copies of the pair of generic random variables (X', Y'), where X' and Y' have the same marginal distributions as X and Y, respectively. Furthermore, from the random coding scheme and the memorylessness of the DMC, it is intuitively correct that X' and Y' are independent because  $\tilde{\mathbf{X}}(1)$  and  $\tilde{\mathbf{X}}(w)$  are independent and the generation of  $\mathbf{Y}$  depends only on  $\tilde{\mathbf{X}}(1)$ .

A formal proof of this claim requires a more detailed analysis. In our random coding scheme, the random variables are generated in the order  $\tilde{\mathbf{X}}(1), \tilde{\mathbf{X}}(2), \dots, \tilde{\mathbf{X}}(M), W, X_1, Y_1, X_2, Y_2, \dots, X_n, Y_n, \hat{W}$ . By considering the joint distribution of these random variables, similar to the discussion in Section 7.3, the Markov chain

$$(\tilde{\mathbf{X}}(1), \tilde{\mathbf{X}}(2), \cdots, \tilde{\mathbf{X}}(M), W) \to \mathbf{X} \to \mathbf{Y} \to \hat{W}$$
 (7.160)

can be established. See Problem 1 for the details. Then for any  $2 \le w \le M$ , from the above Markov chain, we have

$$I(\mathbf{Y}; \tilde{\mathbf{X}}(w), W | \mathbf{X}) = 0.$$
(7.161)

By the chain rule for mutual information, the left hand side can be written as

$$I(\mathbf{Y}; W | \mathbf{X}) + I(\mathbf{Y}; \mathbf{X}(w) | \mathbf{X}, W).$$
(7.162)

By the nonnegativity of conditional mutual information, this implies

$$I(\mathbf{Y}; \mathbf{X}(w) | \mathbf{X}, W) = 0, \qquad (7.163)$$

or

$$\sum_{w=1}^{M} \Pr\{W = w\} I(\mathbf{Y}; \tilde{\mathbf{X}}(w) | \mathbf{X}, W = w) = 0.$$
 (7.164)

Since  $I(\mathbf{Y}; \tilde{\mathbf{X}}(w) | \mathbf{X}, W = w)$  are all nonnegative, we see from the above that they must all vanish. In particular,

$$I(\mathbf{Y}; \mathbf{X}(w) | \mathbf{X}, W = 1) = 0.$$
 (7.165)

Then

$$I(\mathbf{Y}; \tilde{\mathbf{X}}(w) | \tilde{\mathbf{X}}(1), W = 1) = I(\mathbf{Y}; \tilde{\mathbf{X}}(w) | \tilde{\mathbf{X}}(W), W = 1)$$
(7.166)

$$= I(\mathbf{Y}; \tilde{\mathbf{X}}(w) | \mathbf{X}, W = 1)$$
(7.167)

$$= 0.$$
 (7.168)

On the other hand, since  $\tilde{\mathbf{X}}(1), \tilde{\mathbf{X}}(w)$ , and W are mutually independent, we have

$$I(\mathbf{X}(1); \mathbf{X}(w) | W = 1) = 0.$$
 (7.169)

Hence,

$$I(\mathbf{Y}; \mathbf{\tilde{X}}(w)|W = 1)$$
  

$$\leq I(\mathbf{\tilde{X}}(1), \mathbf{Y}; \mathbf{\tilde{X}}(w)|W = 1)$$
(7.170)

$$= I(\tilde{\mathbf{X}}(1); \tilde{\mathbf{X}}(w) | W = 1) + I(\mathbf{Y}; \tilde{\mathbf{X}}(w) | \tilde{\mathbf{X}}(1), W = 1)$$
(7.171)

$$= 0 + 0$$
 (7.172)

$$= 0,$$
 (7.173)

where (7.172) follows from (7.168) and (7.169), proving the claim.

Let us now return to (7.158). For any  $2 \le w \le M$ , it follows from the above claim and Lemma 7.17 that

$$\Pr\{E_w|W=1\}$$
  
= 
$$\Pr\{(\tilde{\mathbf{X}}(w), \mathbf{Y}) \in T^n_{[XY]\delta}|W=1\}$$
 (7.174)

$$\leq 2^{-n(I(X;Y)-\tau)},\tag{7.175}$$

where  $\tau \to 0$  as  $\delta \to 0$ .

From the upper bound in (7.146), we have

$$M < 2^{n(I(X;Y) - \frac{\epsilon}{4})}.$$
(7.176)

Using (7.159), (7.175), and the above upper bound on M, it follows from (7.152) and (7.158) that

$$\Pr\{Err\} < \nu + 2^{n(I(X;Y) - \frac{\epsilon}{4})} \cdot 2^{-n(I(X;Y) - \tau)}$$
(7.177)

$$= \nu + 2^{-n(\frac{\epsilon}{4} - \tau)}.$$
 (7.178)

Since  $\tau \to 0$  as  $\delta \to 0$ , for sufficiently small  $\delta$ , we have

$$\frac{\epsilon}{4} - \tau > 0 \tag{7.179}$$

for any  $\epsilon > 0$ , so that  $2^{-n(\frac{\epsilon}{4}-\tau)} \to 0$  as  $n \to \infty$ . Then by letting  $\nu < \frac{\epsilon}{3}$ , it follows from (7.178) that

$$\Pr\{Err\} < \frac{\epsilon}{2} \tag{7.180}$$

for sufficiently large n.

The main idea of the above analysis of  $\Pr{Err}$  is the following. In constructing the codebook, we randomly generate M codewords in  $\mathcal{X}^n$  according to  $p(x)^n$ , and one of the codewords is sent through the channel p(y|x). When nis large, with high probability, the received sequence is jointly typical with the codeword sent with respect to p(x, y). If the number of codewords M grows with n at a rate less than I(X;Y), then the probability that the received sequence is jointly typical with a codeword other than the one sent through the channel is negligible. Accordingly, the message can be decoded correctly with probability arbitrarily close to 1.

In constructing the codebook in Step 1 of the random coding scheme, we choose a codebook C with a certain probability  $Pr\{C\}$  from the ensemble of all possible codebooks. By conditioning on the codebook chosen, we have

7.4 Achievability 159

$$\Pr\{Err\} = \sum_{\mathcal{C}} \Pr\{\mathcal{C}\} \Pr\{Err|\mathcal{C}\}, \qquad (7.181)$$

i.e.,  $\Pr{Err}$  is a weighted average of  $\Pr{Err|\mathcal{C}}$  over all  $\mathcal{C}$  in the ensemble of all possible codebooks, where  $\Pr{Err|\mathcal{C}}$  is the average probability of error of the code, i.e.,  $P_e$ , when the codebook  $\mathcal{C}$  is chosen (cf. Definition 7.12). The reader should compare the two different expansions of  $\Pr{Err}$  in (7.181) and (7.150).

Therefore, there exists at least one codebook  $\mathcal{C}^*$  such that

$$\Pr\{Err|\mathcal{C}^*\} \le \Pr\{Err\} < \frac{\epsilon}{2}.$$
(7.182)

Thus we have shown that for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M) code such that

$$\frac{1}{n}\log M > I(X;Y) - \frac{\epsilon}{2} \tag{7.183}$$

(cf. (7.146)) and

$$P_e < \frac{\epsilon}{2}.\tag{7.184}$$

We are still one step away from proving that the rate I(X;Y) is achievable because we require that  $\lambda_{max}$  instead of  $P_e$  is arbitrarily small. Toward this end, we write (7.184) as

$$\frac{1}{M}\sum_{w=1}^{M}\lambda_w < \frac{\epsilon}{2},\tag{7.185}$$

or

$$\sum_{w=1}^{M} \lambda_w < \left(\frac{M}{2}\right)\epsilon. \tag{7.186}$$

Upon ordering the codewords according to their conditional probabilities of error, we observe that the conditional probabilities of error of the better half of the M codewords are less than  $\epsilon$ , otherwise the conditional probabilities of error of the worse half of the codewords are at least  $\epsilon$ , and they contribute at least  $(\frac{M}{2})\epsilon$  to the summation in (7.186), which is a contradiction.

Thus by discarding the worse half of the codewords in  $C^*$ , for the resulting codebook, the maximal probability of error  $\lambda_{max}$  is less than  $\epsilon$ . Using (7.183) and considering

$$\frac{1}{n}\log\frac{M}{2} = \frac{1}{n}\log M - \frac{1}{n}$$
(7.187)

$$> \left(I(X;Y) - \frac{\epsilon}{2}\right) - \frac{1}{n} \tag{7.188}$$

$$> I(X;Y) - \epsilon \tag{7.189}$$

when n is sufficiently large, we see that the rate of the resulting code is greater than  $I(X;Y) - \epsilon$ . Hence, we conclude that the rate I(X;Y) is achievable.

Finally, upon letting the input distribution p(x) be one which achieves the channel capacity, i.e., I(X;Y) = C, we have proved that the rate C is achievable. This completes the proof of the direct part of the channel coding theorem.

## 7.5 A Discussion

In the last two sections, we have proved the channel coding theorem which asserts that reliable communication through a DMC at rate R is possible if and only if R < C, the channel capacity. By reliable communication at rate R, we mean that the size of the message set grows exponentially with n at rate R, while the message can be decoded correctly with probability arbitrarily close to 1 as  $n \to \infty$ . Therefore, the capacity C is a fundamental characterization of a DMC.

The capacity of a noisy channel is analogous to the capacity of a water pipe in the following way. For a water pipe, if we pump water through the pipe at a rate higher than its capacity, the pipe would burst and water would be lost. For a communication channel, if we communicate through the channel at a rate higher than the capacity, the probability of error is bounded away from zero, i.e., information is lost.

In proving the direct part of the channel coding theorem, we showed that there exists a channel code whose rate is arbitrarily close to C and whose probability of error is arbitrarily close to zero. Moreover, the existence of such a code is guaranteed only when the block length n is large. However, the proof does not indicate how we can find such a codebook. For this reason, the proof we gave is called an existence proof (as oppose to a constructive proof).

For a fixed block length n, we in principle can search through the ensemble of all possible codebooks for a good one, but this is quite prohibitive even for small n because the number of all possible codebooks grows doubly exponentially with n. Specifically, the total number of all possible (n, M) codebooks is equal to  $|\mathcal{X}|^{Mn}$ . When the rate of the code is close to C, M is approximately equal to  $2^{nC}$ . Therefore, the number of codebooks we need to search through is about  $|\mathcal{X}|^{n2^{nC}}$ .

Nevertheless, the proof of the direct part of the channel coding theorem does indicate that if we generate a codebook randomly as prescribed, the codebook is most likely to be good. More precisely, we now show that the probability of choosing a code C such that  $\Pr\{Err|C\}$  is greater than any prescribed  $\psi > 0$  is arbitrarily small when n is sufficiently large. Consider

$$\Pr{Err} = \sum_{\mathcal{C}} \Pr{\mathcal{C}} \Pr{Err|\mathcal{C}}$$

$$= \sum_{\mathcal{C}: \Pr{Err|\mathcal{C}} \le \psi} \Pr{\mathcal{C}} \Pr{Err|\mathcal{C}}$$
(7.190)

7.5 A Discussion 161

$$+\sum_{\mathcal{C}:\Pr\{Err|\mathcal{C}\}>\psi}\Pr\{\mathcal{C}\}\Pr\{Err|\mathcal{C}\}$$
(7.191)

$$\geq \sum_{\mathcal{C}:\Pr\{Err|\mathcal{C}\}>\psi} \Pr\{\mathcal{C}\}\Pr\{Err|\mathcal{C}\}$$
(7.192)

$$> \psi \sum_{\mathcal{C}: \Pr\{Err | \mathcal{C}\} > \psi} \Pr\{\mathcal{C}\},$$
(7.193)

which implies

$$\sum_{\mathcal{C}:\Pr\{Err|\mathcal{C}\}>\psi} \Pr\{\mathcal{C}\} < \frac{\Pr\{Err\}}{\psi}.$$
(7.194)

From (7.182), we have

$$\Pr\{Err\} < \frac{\epsilon}{2} \tag{7.195}$$

for any  $\epsilon > 0$  when n is sufficiently large. Then

$$\sum_{\mathcal{C}:\Pr\{Err|\mathcal{C}\}>\psi} \Pr\{\mathcal{C}\} < \frac{\epsilon}{2\psi}.$$
(7.196)

Since  $\psi$  is fixed, this upper bound can be made arbitrarily small by choosing a sufficiently small  $\epsilon$ .

Although the proof of the direct part of the channel coding theorem does not provide an explicit construction of a good code, it does give much insight into what a good code is like. Figure 7.12 is an illustration of a channel code which achieves the channel capacity. Here we assume that the input distribu-

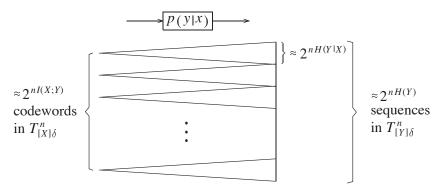


Fig. 7.12. A channel code which achieves capacity.

tion p(x) is one which achieves the channel capacity, i.e., I(X;Y) = C. The idea is that most of the codewords are typical sequences in  $\mathcal{X}^n$  with respect to p(x). (For this reason, the repetition code is not a good code.) When such a codeword is transmitted through the channel, the received sequence is likely

to be one of about  $2^{nH(Y|X)}$  sequences in  $\mathcal{Y}^n$  which are jointly typical with the transmitted codeword with respect to p(x, y). The association between a codeword and the about  $2^{nH(Y|X)}$  corresponding sequences in  $\mathcal{Y}^n$  is shown as a cone in the figure. As we require that the probability of decoding error is small, the cones essentially do not overlap with each other. Since the number of typical sequences with respect to p(y) is about  $2^{nH(Y)}$ , the number of codewords cannot exceed about

$$\frac{2^{nH(Y)}}{2^{nH(Y|X)}} = 2^{nI(X;Y)} = 2^{nC}.$$
(7.197)

This is consistent with the converse of the channel coding theorem. The direct part of the channel coding theorem says that when n is large, as long as the number of codewords generated randomly is not more than about  $2^{n(C-\epsilon)}$ , the overlap among the cones is negligible with high probability.

Therefore, instead of searching through the ensemble of all possible codebooks for a good one, we can generate a codebook randomly, and it is likely to be good. However, such a code is difficult to use due to the following implementation issues.

A codebook with block length n and rate R consists of  $n2^{nR}$  symbols from the input alphabet  $\mathcal{X}$ . This means that the size of the codebook, i.e., the amount of storage required to store the codebook, grows exponentially with n. This also makes the encoding process inefficient.

Another issue is regarding the computation required for decoding. Based on the sequence received at the output of the channel, the decoder needs to decide which of the about  $2^{nR}$  codewords was the one transmitted. This requires an exponential amount of computation.

In practice, we are satisfied with the reliability of communication once it exceeds a certain level. Therefore, the above implementation issues may eventually be resolved with the advancement of microelectronics. But before then, we still have to deal with these issues. For this reason, the entire field of *coding theory* has been developed since the 1950's. Researchers in this field are devoted to searching for good codes and devising efficient decoding algorithms.

In fact, almost all the codes studied in coding theory are *linear codes*. By taking advantage of the linear structures of these codes, efficient encoding and decoding can be achieved. In particular, Berrou *et al.* [32] have proposed a linear code called the *turbo code*<sup>6</sup> in 1993, which is now generally believed to be the practical way to achieve the channel capacity.

Today, channel coding has been widely used in home entertainment systems (e.g., audio CD and DVD), computer storage systems (e.g., CD-ROM, hard disk, floppy disk, and magnetic tape), computer communication, wireless

<sup>&</sup>lt;sup>6</sup> The turbo code is a special case of the class of *Low-density parity-check* (LDPC) codes proposed by Gallager [115] in 1962 (see MacKay [215]). However, the performance of such codes was not known at that time due to lack of high speed computers for simulation.

communication, and deep space communication. The most popular channel codes used in existing systems include the Hamming code, the Reed-Solomon code<sup>7</sup>, the BCH code, and convolutional codes. We refer the interested reader to textbooks on coding theory [37] [209] [341] for discussions of this subject.

### 7.6 Feedback Capacity

Feedback is common in practical communication systems for correcting possible errors which occur during transmission. As an example, during a telephone conversation, we often have to request the speaker to repeat due to poor voice quality of the telephone line. As another example, in data communication the receiver may request a packet to be retransmitted if the *parity check* bits received are incorrect. In general, when feedback from the receiver is available at the transmitter, the transmitter can at any time decide what to transmit next based on the feedback so far, and can potentially transmit information through the channel reliably at a higher rate.

In this section, we study a model in which a DMC is used with complete feedback. The block diagram for the model is shown in Figure 7.13. In this

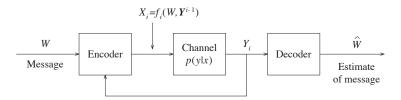


Fig. 7.13. A channel code with feedback.

model, the symbol  $Y_i$  received at the output of the channel at time *i* is available instantaneously at the encoder without error. Then depending on the message W and all the previous feedback  $Y_1, Y_2, \dots, Y_i$ , the encoder decides the value of  $X_{i+1}$ , the next symbol to be transmitted. Such a channel code is formally defined below.

**Definition 7.18.** An (n, M) code with complete feedback for a discrete memoryless channel with input alphabet  $\mathcal{X}$  and output alphabet  $\mathcal{Y}$  is defined by encoding functions

$$f_i: \{1, 2, \cdots, M\} \times \mathcal{Y}^{i-1} \to \mathcal{X}$$
(7.198)

for  $1 \leq i \leq n$  and a decoding function

$$g: \mathcal{Y}^n \to \{1, 2, \cdots, M\}. \tag{7.199}$$

<sup>&</sup>lt;sup>7</sup> The Reed-Solomon code was independently discovered by Arimoto [17].

We will use  $\mathbf{Y}^i$  to denote  $(Y_1, Y_2, \dots, Y_i)$  and  $X_i$  to denote  $f_i(W, \mathbf{Y}^{i-1})$ . We note that a channel code without feedback is a special case of a channel code with complete feedback because for the latter, the encoder can ignore the feedback.

**Definition 7.19.** A rate R is achievable with complete feedback for a discrete memoryless channel p(y|x) if for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M) code with complete feedback such that

$$\frac{1}{n}\log M > R - \epsilon \tag{7.200}$$

and

$$\lambda_{max} < \epsilon. \tag{7.201}$$

**Definition 7.20.** The feedback capacity,  $C_{FB}$ , of a discrete memoryless channel is the supremum of all the rates achievable by codes with complete feedback.

**Proposition 7.21.** The supremum in the definition of  $C_{FB}$  in Definition 7.20 is the maximum.

*Proof.* Consider rates  $\mathbb{R}^{(k)}$  which are achievable with complete feedback such that

$$\lim_{k \to \infty} R^{(k)} = R. \tag{7.202}$$

Then for any  $\epsilon > 0$ , for all k, there exists for sufficiently large n an  $(n, M^{(k)})$  code with complete feedback such that

$$\frac{1}{n}\log M^{(k)} > R^{(k)} - \epsilon \tag{7.203}$$

and

$$\lambda_{max}^{(k)} < \epsilon. \tag{7.204}$$

By virtue of (7.202), let  $k(\epsilon)$  be an integer such that for all  $k > k(\epsilon)$ ,

$$|R - R^{(k)}| < \epsilon, \tag{7.205}$$

which implies

$$R^{(k)} > R - \epsilon. \tag{7.206}$$

Then for all  $k > k(\epsilon)$ ,

$$\frac{1}{n}\log M^{(k)} > R^{(k)} - \epsilon > R - 2\epsilon.$$
(7.207)

Therefore, it follows from (7.207) and (7.204) that R is achievable with complete feedback. This implies that the supremum in Definition 7.20, which can be achieved, is in fact the maximum.  $\Box$ 

Since a channel code without feedback is a special case of a channel code with complete feedback, any rate R achievable by the former is also achievable by the latter. Therefore,

$$C_{FB} \ge C. \tag{7.208}$$

A fundamental question is whether  $C_{FB}$  is greater than C. The answer surprisingly turns out to be negative for a DMC, as we now show. From the description of a channel code with complete feedback, we obtain the dependency graph for the random variables  $W, \mathbf{X}, \mathbf{Y}, \hat{W}$  in Figure 7.14. From this

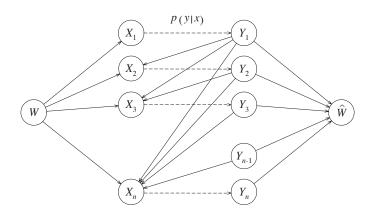


Fig. 7.14. The dependency graph for a channel code with feedback.

dependency graph, we see that

$$q(w, \mathbf{x}, \mathbf{y}, \hat{w}) = q(w) \left(\prod_{i=1}^{n} q(x_i | w, \mathbf{y}^{i-1})\right) \left(\prod_{i=1}^{n} p(y_i | x_i)\right) q(\hat{w} | \mathbf{y})$$
(7.209)

for all  $(w, \mathbf{x}, \mathbf{y}, \hat{w}) \in \mathcal{W} \times \mathcal{X}^n \times \mathcal{Y}^n \times \mathcal{W}$  such that  $q(w, \mathbf{y}^{i-1}), q(x_i) > 0$  for  $1 \leq i \leq n$  and  $q(\mathbf{y}) > 0$ , where  $\mathbf{y}^i = (y_1, y_2, \dots, y_i)$ . Note that  $q(x_i|w, \mathbf{y}^{i-1})$  and  $q(\hat{w}|\mathbf{y})$  are deterministic.

Lemma 7.22. For all  $1 \leq i \leq n$ ,

$$(W, \mathbf{Y}^{i-1}) \to X_i \to Y_i \tag{7.210}$$

forms a Markov chain.

*Proof.* The dependency graph for the random variables  $W, \mathbf{X}^{i}$ , and  $\mathbf{Y}^{i}$  is shown in Figure 7.15. Denote the set of nodes  $W, \mathbf{X}^{i-1}$ , and  $\mathbf{Y}^{i-1}$  by Z. Then we see that all the edges from Z end at  $X_i$ , and the only edge from  $X_i$  ends at  $Y_i$ . This means that  $Y_i$  depends on  $(W, \mathbf{X}^{i-1}, \mathbf{Y}^{i-1})$  only through  $X_i$ , i.e.,

$$(W, \mathbf{X}^{i-1}, \mathbf{Y}^{i-1}) \to X_i \to Y_i$$
 (7.211)

forms a Markov chain, or

$$I(W, \mathbf{X}^{i-1}, \mathbf{Y}^{i-1}; Y_i | X_i) = 0.$$
(7.212)

This can be formally justified by Proposition 2.9, and the details are omitted here. Since

$$0 = I(W, \mathbf{X}^{i-1}, \mathbf{Y}^{i-1}; Y_i | X_i)$$
(7.213)

$$= I(W, \mathbf{Y}^{i-1}; Y_i | X_i) + I(\mathbf{X}^{i-1}; Y_i | W, X_i, \mathbf{Y}^{i-1})$$
(7.214)

and mutual information is nonnegative, we obtain

$$I(W, \mathbf{Y}^{i-1}; Y_i | X_i) = 0, (7.215)$$

or

$$W, \mathbf{Y}^{i-1}) \to X_i \to Y_i$$
 (7.216)

forms a Markov chain. The lemma is proved.  $\Box$ 

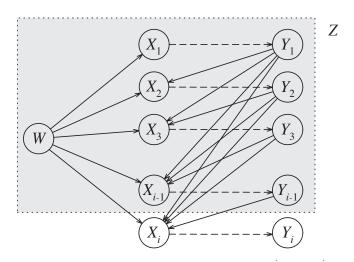


Fig. 7.15. The dependency graph for  $W, \mathbf{X}^i$ , and  $\mathbf{Y}^i$ .

From the definition of  $C_{FB}$  and by virtue of Proposition 7.21, if  $R \leq C_{FB}$ , then R is a rate achievable with complete feedback. We will show that if a rate R is achievable with complete feedback, then  $R \leq C$ . If so, then  $R \leq C_{FB}$ implies  $R \leq C$ , which can be true if and only if  $C_{FB} \leq C$ . Then from (7.208), we can conclude that  $C_{FB} = C$ .

Let R be a rate achievable with complete feedback, i.e., for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M) code with complete feedback such that

#### 7.6 Feedback Capacity 167

$$n^{-1}\log M > R - \epsilon \tag{7.217}$$

and

$$\lambda_{max} < \epsilon. \tag{7.218}$$

Consider

$$\log M = H(W) = I(W; \mathbf{Y}) + H(W|\mathbf{Y})$$
(7.219)

and bound  $I(W; \mathbf{Y})$  and  $H(W|\mathbf{Y})$  as follows. First,

$$I(W; \mathbf{Y}) = H(\mathbf{Y}) - H(\mathbf{Y}|W)$$
(7.220)

$$= H(\mathbf{Y}) - \sum_{i=1}^{n} H(Y_i | \mathbf{Y}^{i-1}, W)$$
 (7.221)

$$\stackrel{a)}{=} H(\mathbf{Y}) - \sum_{i=1}^{n} H(Y_i | \mathbf{Y}^{i-1}, W, X_i)$$
(7.222)

<sup>b)</sup> 
$$= H(\mathbf{Y}) - \sum_{i=1}^{n} H(Y_i|X_i)$$
 (7.223)

$$\leq \sum_{\substack{i=1\\n}}^{n} H(Y_i) - \sum_{i=1}^{n} H(Y_i|X_i)$$
(7.224)

$$=\sum_{i=1}^{n} I(X_i; Y_i)$$
(7.225)

$$\leq nC,$$
 (7.226)

where a) follows because  $X_i$  is a function of W and  $\mathbf{Y}^{i-1}$  and b) follows from Lemma 7.22. Second,

$$H(W|\mathbf{Y}) = H(W|\mathbf{Y}, \hat{W}) \le H(W|\hat{W}).$$
 (7.227)

Thus

$$\log M \le H(W|\hat{W}) + nC, \tag{7.228}$$

which is the same as (7.126). Then by (7.217) and an application of Fano's inequality, we conclude as in the proof for the converse of the channel coding theorem that

$$R \le C. \tag{7.229}$$

Hence, we have proved that  $C_{FB} = C$ .

**Remark 1** The proof for the converse of the channel coding theorem in Section 7.3 depends critically on the Markov chain

$$W \to \mathbf{X} \to \mathbf{Y} \to \hat{W}$$
 (7.230)

and the relation in (7.101) (the latter implies Lemma 7.16). Both of them do not hold in general in the presence of feedback.

#### 168 7 Discrete Memoryless Channels

**Remark 2** The proof for  $C_{FB} = C$  in this section is also a proof for the converse of the channel coding theorem, so we actually do not need the proof in Section 7.3. However, the proof here and the proof in Section 7.3 have different spirits. Without comparing the two proofs, one cannot possibly understand the subtlety of the result that feedback does not increase the capacity of a DMC.

**Remark 3** Although feedback does not increase the capacity of a DMC, the availability of feedback often makes coding much simpler. For some channels, communication through the channel with zero probability of error can be achieved in the presence of feedback by using a *variable-length* channel code. These are discussed in the next example.

Example 7.23. Consider the binary erasure channel in Example 7.8 whose capacity is  $1 - \gamma$ , where  $\gamma$  is the erasure probability. In the presence of complete feedback, for every information bit to be transmitted, the encoder can transmit the same information bit through the channel until an erasure does not occur, i.e., the information bit is received correctly. Then the number of uses of the channel it takes to transmit an information bit through the channel correctly has a truncated geometrical distribution whose mean is  $(1 - \gamma)^{-1}$ . Therefore, the effective rate at which information can be transmitted through the channel is  $1 - \gamma$ . In other words, the channel capacity is achieved by using a very simple variable-length code. Moreover, the channel capacity is achieved with zero probability of error.

In the absence of feedback, the rate  $1 - \gamma$  can also be achieved, but with an arbitrarily small probability of error and a much more complicated code.

To conclude this section, we point out that the memoryless assumption of the channel is essential for drawing the conclusion that feedback does not increase the channel capacity not because the proof presented in this section does not go through without this assumption, but because if the channel has memory, feedback actually can increase the channel capacity. For an illustrating example, see Problem 12.

## 7.7 Separation of Source and Channel Coding

We have so far considered the situation in which we want to convey a message through a DMC, where the message is randomly selected from a finite set according to the uniform distribution. However, in most situations, we want to convey an information source through a DMC. Let  $\{U_k, k > -n\}$  be an ergodic stationary information source with entropy rate H. Denote the common alphabet by  $\mathcal{U}$  and assume that  $\mathcal{U}$  is finite. To convey  $\{U_k\}$  through the channel, we can employ a source code with rate  $R_s$  and a channel code with rate  $R_c$  as shown in Figure 7.16 such that  $R_s < R_c$ .

$$U \longrightarrow \text{source} \xrightarrow{W} \text{channel} \xrightarrow{X} p(y|x) \xrightarrow{Y} \text{channel} \xrightarrow{W} \text{source} \text{decoder} \xrightarrow{W} \hat{U}$$

Fig. 7.16. Separation of source coding and channel coding.

Let  $f^s$  and  $g^s$  be respectively the encoding function and the decoding function of the source code, and  $f^c$  and  $g^c$  be respectively the encoding function and the decoding function of the channel code. The block of n information symbols  $\mathbf{U} = (U_{-(n-1)}, U_{-(n-2)}, \cdots, U_0)$  is first encoded by the source encoder into an index

$$W = f^s(\mathbf{U}),\tag{7.231}$$

called the source codeword. Then W is mapped by the channel encoder to a distinct channel codeword

$$\mathbf{X} = f^c(W), \tag{7.232}$$

where  $\mathbf{X} = (X_1, X_2, \dots, X_n)$ . This is possible because there are about  $2^{nR_s}$  source codewords and about  $2^{nR_c}$  channel codewords, and we assume that  $R_s < R_c$ . Then  $\mathbf{X}$  is transmitted through the DMC p(y|x), and the sequence  $\mathbf{Y} = (Y_1, Y_2, \dots, Y_n)$  is received. Based on  $\mathbf{Y}$ , the channel decoder first estimates W as

$$\hat{W} = g^c(\mathbf{Y}). \tag{7.233}$$

Finally, the source decoder decodes  $\hat{W}$  to

$$\hat{\mathbf{U}} = g^s(\hat{W}). \tag{7.234}$$

For this scheme, an error occurs if  $\mathbf{U} \neq \hat{\mathbf{U}}$ , and we denote the probability of error by  $P_e$ .

We now show that if H < C, the capacity of the DMC p(y|x), then it is possible to convey **U** through the channel with an arbitrarily small probability of error. First, we choose  $R_s$  and  $R_c$  such that

$$H < R_s < R_c < C. \tag{7.235}$$

Observe that if  $\hat{W} = W$  and  $g^s(W) = \mathbf{U}$ , then from (7.234),

$$\hat{\mathbf{U}} = g^s(\hat{W}) = g^s(W) = \mathbf{U}, \qquad (7.236)$$

i.e., an error does not occur. In other words, if an error occurs, either  $\hat{W} \neq W$  or  $g^s(W) \neq \mathbf{U}$ . Then by the union bound, we have

$$P_e \le \Pr\{\hat{W} \neq W\} + \Pr\{g^s(W) \neq \mathbf{U}\}.$$
(7.237)

For any  $\epsilon > 0$  and sufficiently large n, by the Shannon-McMillan-Breiman theorem, there exists a source code such that

$$\Pr\{g^s(W) \neq \mathbf{U}\} \le \epsilon. \tag{7.238}$$

#### 170 7 Discrete Memoryless Channels

By the channel coding theorem, there exists a channel code such that  $\lambda_{max} \leq \epsilon$ , where  $\lambda_{max}$  is the maximal probability of error. This implies

$$\Pr\{\hat{W} \neq W\} = \sum_{w} \Pr\{\hat{W} \neq W | W = w\} \Pr\{W = w\}$$
(7.239)

$$\leq \lambda_{max} \sum_{w} \Pr\{W = w\}$$
(7.240)

$$=\lambda_{max} \tag{7.241}$$

$$\leq \epsilon.$$
 (7.242)

Combining (7.238) and (7.242), we have

=

$$P_e \le 2\epsilon. \tag{7.243}$$

Therefore, we conclude that as long as H < C, it is possible to convey  $\{U_k\}$  through the DMC reliably.

In the scheme we have discussed, source coding and channel coding are separated. In general, source coding and channel coding can be combined. This technique is called *joint source-channel coding*. It is then natural to ask whether it is possible to convey information through the channel reliably at a higher rate by using joint source-channel coding. In the rest of the section, we show that the answer to this question is no to the extent that for asymptotic reliability, we must have  $H \leq C$ . However, whether asymptotical reliability can be achieved for H = C depends on the specific information source and channel.

We base our discussion on the general assumption that complete feedback is available at the encoder as shown in Figure 7.17. Let  $f_i^{sc}$ ,  $1 \le i \le n$ , be

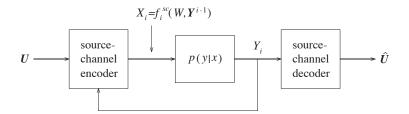


Fig. 7.17. Joint source-channel coding.

the encoding functions and  $g^{sc}$  be the decoding function of the source-channel code. Then

$$X_i = f_i^{sc}(\mathbf{U}, \mathbf{Y}^{i-1}) \tag{7.244}$$

for  $1 \le i \le n$ , where  $\mathbf{Y}^{i-1} = (Y_1, Y_2, \cdots, Y_{i-1})$ , and

$$\hat{\mathbf{U}} = g^{sc}(\mathbf{Y}),\tag{7.245}$$

## 7.7 Separation of Source and Channel Coding 171

where  $\hat{\mathbf{U}} = (\hat{U}_1, \hat{U}_2, \dots, \hat{U}_n)$ . In exactly the same way as we proved (7.226) in the last section, we can prove that

$$I(\mathbf{U}; \mathbf{Y}) \le nC. \tag{7.246}$$

Since  $\hat{\mathbf{U}}$  is a function of  $\mathbf{Y}$ ,

$$I(\mathbf{U}; \hat{\mathbf{U}}) \le I(\mathbf{U}; \hat{\mathbf{U}}, \mathbf{Y}) \tag{7.247}$$

$$=I(\mathbf{U};\mathbf{Y})\tag{7.248}$$

$$\leq nC.$$
 (7.249)

For any  $\epsilon > 0$ ,

$$H(\mathbf{U}) \ge n(H - \epsilon) \tag{7.250}$$

for sufficiently large n. Then

$$n(H-\epsilon) \le H(\mathbf{U}) = H(\mathbf{U}|\hat{\mathbf{U}}) + I(\mathbf{U};\hat{\mathbf{U}}) \le H(\mathbf{U}|\hat{\mathbf{U}}) + nC.$$
(7.251)

Applying Fano's inequality (Corollary 2.48), we obtain

$$n(H-\epsilon) \le 1 + nP_e \log |\mathcal{U}| + nC, \tag{7.252}$$

or

$$H - \epsilon \le \frac{1}{n} + P_e \log |\mathcal{U}| + C. \tag{7.253}$$

For asymptotic reliability,  $P_e \to 0$  as  $n \to \infty$ . Therefore, by letting  $n \to \infty$  and then  $\epsilon \to 0$ , we conclude that

$$H \le C. \tag{7.254}$$

This result, sometimes called the *separation theorem for source and channel coding*, says that asymptotic optimality can be achieved by separating source coding and channel coding. This theorem has significant engineering implication because the source code and the channel code can be designed separately without losing asymptotic optimality. Specifically, we only need to design the best source code for the information source and design the best channel code for the channel. Moreover, separation of source coding and channel coding facilitates the transmission of different information sources on the same channel because we need only change the source code for different information sources. Likewise, separation of source coding and channel coding also facilitates the transmission of an information source on different channels because we need only change the channel code for different channels.

We remark that although asymptotic optimality can be achieved by separating source coding and channel coding, for finite block length, the probability of error generally can be reduced by using joint source-channel coding. 172 7 Discrete Memoryless Channels

## Problems

In the following,  $\mathbf{X} = (X_1, X_2, \dots, X_n)$ ,  $\mathbf{x} = (x_1, x_2, \dots, x_n)$ , and so on.

- 1. Refer to the discussion in Section 7.4.
  - a) Construct the dependency graph for the random variables involved in the random coding scheme.
  - b) By considering the joint distribution of these random variables, prove the Markov chain in (7.160).
- 2. Show that the capacity of a DMC with complete feedback cannot be increased by using probabilistic encoding and/or decoding schemes.
- 3. Memory increases capacity Consider a BSC with crossover probability  $0 < \epsilon < 1$  represented by  $X_i = Y_i + Z_i \mod 2$ , where  $X_i, Y_i$ , and  $Z_i$  are respectively the input, the output, and the noise variable at time *i*. Then

$$\Pr\{Z_i = 0\} = 1 - \epsilon \quad \text{and} \quad \Pr\{Z_i = 1\} = \epsilon$$

for all *i*. We assume that  $\{X_i\}$  and  $\{Z_i\}$  are independent, but we make no assumption that  $Z_i$  are i.i.d. so that the channel may have memory. a) Prove that

$$I(\mathbf{X}; \mathbf{Y}) \le n - h_b(\epsilon).$$

- b) Show that the upper bound in a) can be achieved by letting  $X_i$  be i.i.d. bits taking the values 0 and 1 with equal probability and  $Z_1 = Z_2 = \cdots = Z_n$ .
- c) Show that with the assumptions in b),  $I(\mathbf{X}; \mathbf{Y}) > nC$ , where  $C = 1 h_b(\epsilon)$  is the capacity of the BSC if it is memoryless.
- 4. Consider the channel in Problem 3, Part b).
  - a) Show that the channel capacity is not increased by feedback.
  - b) Devise a coding scheme without feedback that achieves the channel capacity.
- 5. In Remark 1 toward the end of Section 7.6, it was mentioned that in the presence of feedback, both the Markov chain  $W \to \mathbf{X} \to \mathbf{Y} \to \hat{W}$  and Lemma 7.16 do not hold in general. Give examples to substantiate this remark.
- 6. Prove that when a DMC is used with complete feedback,

$$\Pr\{Y_i = y_i | \mathbf{X}^i = \mathbf{x}^i, \mathbf{Y}^{i-1} = \mathbf{y}^{i-1}\} = \Pr\{Y_i = y_i | X_i = x_i\}$$

for all  $i \ge 1$ . This relation, which is a consequence of the causality of the code, says that given the current input, the current output does not depend on all the past inputs and outputs of the DMC.

7. Let

$$P(\epsilon) = \begin{bmatrix} 1 - \epsilon & \epsilon \\ \epsilon & 1 - \epsilon \end{bmatrix}$$

be the transition matrix for a BSC with crossover probability  $\epsilon$ . Define a \* b = (1 - a)b + a(1 - b) for  $0 \le a, b \le 1$ .

- a) Prove that a DMC with transition matrix  $P(\epsilon_1)P(\epsilon_2)$  is equivalent to a BSC with crossover probability  $\epsilon_1 * \epsilon_2$ . Such a channel is the cascade of two BSC's with crossover probabilities  $\epsilon_1$  and  $\epsilon_2$ , respectively.
- b) Repeat a) for a DMC with transition matrix  $P(\epsilon_2)P(\epsilon_1)$ .
- c) Prove that

$$1 - h_b(\epsilon_1 * \epsilon_2) \le \min(1 - h_b(\epsilon_1), 1 - h_b(\epsilon_2)).$$

This means that the capacity of the cascade of two BSC's is upper bounded by the capacity of either of the two BSC's.

- d) Prove that a DMC with transition matrix  $P(\epsilon)^n$  is equivalent to a BSC with crossover probabilities  $\frac{1}{2}(1-(1-2\epsilon)^n)$ .
- 8. Symmetric channel A DMC is symmetric if the rows of the transition matrix p(y|x) are permutations of each other and so are the columns. Determine the capacity of such a channel.

See Section 4.5 in Gallager [117] for a more general discussion.

- 9. Let  $C_1$  and  $C_2$  be the capacities of two DMC's with transition matrices  $P_1$  and  $P_2$ , respectively, and let C be the capacity of the DMC with transition matrix  $P_1P_2$ . Prove that  $C \leq \min(C_1, C_2)$ .
- 10. Two parallel channels Let  $C_1$  and  $C_2$  be the capacities of two DMC's  $p_1(y_1|x_1)$  and  $p_2(y_2|x_2)$ , respectively. Determine the capacity of the DMC

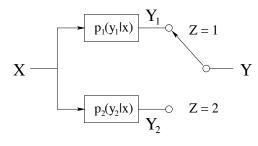
$$p(y_1, y_2|x_1, x_2) = p_1(y_1|x_1)p_2(y_2|x_2).$$

Hint: Prove that

$$I(X_1, X_2; Y_1, Y_2) \le I(X_1; Y_1) + I(X_2; Y_2)$$

if  $p(y_1, y_2|x_1, x_2) = p_1(y_1|x_1)p_2(y_2|x_2)$ .

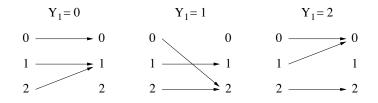
11. In the system below, there are two channels with transition matrices  $p_1(y_1|x)$  and  $p_2(y_2|x)$ . These two channels have a common input alphabet  $\mathcal{X}$  and output alphabets  $\mathcal{Y}_1$  and  $\mathcal{Y}_2$ , repespectively, where  $\mathcal{Y}_1$  and  $\mathcal{Y}_2$  are disjoint. The position of the switch is determined by a random variable Z which is independent of X, where  $\Pr\{Z=1\} = \lambda$ .



- 174 7 Discrete Memoryless Channels
  - a) Show that

$$I(X;Y) = \lambda I(X;Y_1) + (1-\lambda)I(X;Y_2).$$

- b) The capacity of the system is given by  $C = \max_{p(x)} I(X;Y)$ . Show that  $C \leq \lambda C_1 + (1-\lambda)C_2$ , where  $C_i = \max_{p(x)} I(X;Y_i)$  is the capacity of the channel with transition matrix  $p_i(y_i|x)$ , i = 1, 2.
- c) If both  $C_1$  and  $C_2$  can be achieved by a common input distribution, show that  $C = \lambda C_1 + (1 - \lambda)C_2$ .
- 12. Feedback increases capacity Consider a ternary channel with memory with input/output alphabet  $\{0, 1, 2\}$  as follows. At time 1, the output of the channel  $Y_1$  has a uniform distribution on  $\{0, 1, 2\}$  and is independent of the input  $X_1$  (i.e., the channel outputs each of the values 0, 1, and 2 with probability  $\frac{1}{3}$  regardless of the input). At time 2, the transition from  $X_2$  to  $Y_2$  which depends on the value of  $Y_1$  is depicted below:



For every two subsequent transmissions, the channel replicates itself independently. So we only need to consider the first two transmissions. In the sequel, we regard this channel as described by a generic discrete channel (with transmission duration equals 2) with two input symbols  $X_1$  and  $X_2$ and two output symbols  $Y_1$  and  $Y_2$ , and we will refer to this channel as the *block* channel.

- a) Determine the capacity this block channel when it is used without feedback. Hint: Use the results in Problems 8 and 11.
- b) Consider the following coding scheme when the block channel is used with feedback. Let the message W = (W<sub>1</sub>, W<sub>2</sub>) with W<sub>1</sub> = {0,1,2} and W<sub>2</sub> = {0,1}. Let W<sub>1</sub> and W<sub>2</sub> be independent, and each of them distributes uniformly on its alphabet. First, Let X<sub>1</sub> = W<sub>1</sub> and transmit X<sub>1</sub> through the channel to obtain Y<sub>1</sub>, which is independent of X<sub>1</sub>. Then based on the value of Y<sub>1</sub>, we determine X<sub>2</sub> as follows:

  i) If Y<sub>1</sub> = 0, let X<sub>2</sub> = 0 if W<sub>2</sub> = 0, and let X<sub>2</sub> = 1 if W<sub>2</sub> = 1.
  ii) If Y<sub>1</sub> = 1, let X<sub>2</sub> = 1 if W<sub>2</sub> = 0, and let X<sub>2</sub> = 2 if W<sub>2</sub> = 1.

  iii) If Y<sub>1</sub> = 2, let X<sub>2</sub> = 0 if W<sub>2</sub> = 0, and let X<sub>2</sub> = 2 if W<sub>2</sub> = 1.
  iii) If Y<sub>1</sub> = 2, let X<sub>2</sub> = 0 if W<sub>2</sub> = 0, and let X<sub>2</sub> = 2 if W<sub>2</sub> = 1.

coding scheme, show that for the capacity of this block channel can be increased by feedback.

13. Channel with memory and directed information The memorylessness of a DMC is characterized by the Markov chain  $T_{i-} \to X_i \to Y_i$  according to

the discussion following Definition 7.4. In general, a channel with memory satisfies the Markov chain  $T'_{i-} \to (\mathbf{X}^i, \mathbf{Y}^{i-1}) \to Y_i$ , where  $T'_{i-}$  denotes all the random variables generated in the system before  $X_i$  (i.e., the random variables denoted by  $T_{i-}$ ) except for  $\mathbf{X}^{i-1}$  and  $\mathbf{Y}^{i-1}$ . Consider the use of such a channel in the presence of complete feedback.

- a) Give the dependency graph for all the random variables involved in the coding scheme. Note that the memory of the channel is manifested by the dependence of  $Y_i$  on  $\mathbf{X}^{i-1}$  and  $\mathbf{Y}^{i-1}$  (in addition to its dependence on  $X_i$ ) for  $1 \leq i \leq n$ .
- b) Verify the correctness of the following derivation:

$$\begin{split} I(W;\mathbf{Y}) &= H(\mathbf{Y}) - H(\mathbf{Y}|W) \\ &= \sum_{i=1}^{n} [H(Y_{i}|\mathbf{Y}^{i-1}) - H(Y_{i}|W,\mathbf{Y}^{i-1})] \\ &\leq \sum_{i=1}^{n} [H(Y_{i}|\mathbf{Y}^{i-1}) - H(Y_{i}|W,\mathbf{X}^{i},\mathbf{Y}^{i-1})] \\ &= \sum_{i=1}^{n} [H(Y_{i}|\mathbf{Y}^{i-1}) - H(Y_{i}|\mathbf{X}^{i},\mathbf{Y}^{i-1})] \\ &= \sum_{i=1}^{n} I(Y_{i};\mathbf{X}^{i}|\mathbf{Y}^{i-1}). \end{split}$$

The above upper bound on  $I(W; \mathbf{Y})$ , denoted by  $I(\mathbf{X} \to \mathbf{Y})$ , is called the *directed information* from  $\mathbf{X}$  to  $\mathbf{Y}$ .

- c) Show that the inequality in the derivation in b) is in fact an equality. Hint: Use Definition 7.18.
- d) In the spirit of the informal discussion in Section 7.3, we impose the constraint  $H(W|\mathbf{Y}) = 0$ . Show that

$$H(W) = I(\mathbf{X} \to \mathbf{Y}).$$

This is the generalization of (7.105) for a channel with memory in the presence of complete feedback.

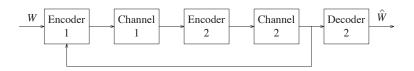
e) Show that  $I(\mathbf{X} \to \mathbf{Y}) = I(\mathbf{X}; \mathbf{Y})$  if the channel code does not make use of the feedback. Hint: First show that

$$H(Y_i|\mathbf{X}^i, \mathbf{Y}^{i-1}) = H(Y_i|W, \mathbf{X}^i, \mathbf{Y}^{i-1}) = H(Y_i|W, \mathbf{X}, \mathbf{Y}^{i-1}).$$

(Marko [220] and Massey [224].)

- 14. Maximum likelihood decoding In maximum likelihood decoding for a given channel and a given codebook, if a received sequence  $\mathbf{y}$  is decoded to a codeword  $\mathbf{x}$ , then  $\mathbf{x}$  maximizes  $\Pr{\{\mathbf{y}|\mathbf{x}'\}}$  among all codewords  $\mathbf{x}'$  in the codebook.
  - a) Prove that maximum likelihood decoding minimizes the average probability of error.

- 176 7 Discrete Memoryless Channels
  - b) Does maximum likelihood decoding also minimize the maximal probability of error? Give an example if your answer is no.
- 15. Minimum distance decoding The Hamming distance between two binary sequences  $\mathbf{x}$  and  $\mathbf{y}$ , denoted by  $d(\mathbf{x}, \mathbf{y})$ , is the number of places where  $\mathbf{x}$  and  $\mathbf{y}$  differ. In minimum distance decoding for a memoryless BSC, if a received sequence  $\mathbf{y}$  is decoded to a codeword  $\mathbf{x}$ , then  $\mathbf{x}$  minimizes  $d(\mathbf{x}', \mathbf{y})$  over all codewords  $\mathbf{x}'$  in the codebook. Prove that minimum distance decoding is equivalent to maximum likelihood decoding if the crossover probability of the BSC is less than 0.5.
- 16. The following figure shows a communication system with two DMC's with complete feedback. The capacities of the two channels are respectively  $C_1$  and  $C_2$ .



- a) Give the dependency graph for all the random variables involved in the coding scheme.
- b) Prove that the capacity of the system is  $\min(C_1, C_2)$ .
- 17. Binary arbitrarily varying channel Consider a memoryless BSC whose crossover probability is time-varying. Specifically, the crossover probability  $\epsilon(i)$  at time *i* is an arbitrary value in  $[\epsilon_1, \epsilon_2]$ , where  $0 \le \epsilon_1 < \epsilon_2 < 0.5$ . Prove that the capacity of this channel is  $1 h_b(\epsilon_2)$ . (Ahlswede and Wolfowitz [12].)
- 18. Consider a BSC with crossover probability  $\epsilon \in [\epsilon_1, \epsilon_2]$ , where  $0 < \epsilon_1 < \epsilon_2 < 0.5$ , but the exact value of  $\epsilon$  is unknown. Prove that the capacity of this channel is  $1 h_b(\epsilon_2)$ .

## **Historical Notes**

The concept of channel capacity was introduced in Shannon's original paper [291], where he stated the channel coding theorem and outlined a proof. The first rigorous proof was due to Feinstein [98]. The random coding error exponent was developed by Gallager [116] in a simplified proof.

The converse of the channel coding theorem was proved by Fano [95], where he used an inequality now bearing his name. The strong converse was first proved by Wolfowitz [345]. An iterative algorithm for calculating the channel capacity developed independently by Arimoto [18] and Blahut [35] will be discussed in Chapter 9. Shannon [295] proved that the capacity of a discrete memoryless channel cannot be increased by feedback.

The definition of a discrete memoryless channel in this chapter is new. With this definition, coding over such a channel with or without feedback can be rigorously formulated.

Let H be the entropy rate of an information source. By the source coding theorem, it is possible to design a source code with rate R which reconstructs the source sequence  $\mathbf{X} = (X_1, X_2, \dots, X_n)$  with an arbitrarily small probability of error provided R > H and the block length n is sufficiently large. However, there are situations in which we want to convey an information source by a source code with rate less than H. Then we are motivated to ask: what is the best we can do when R < H?

A natural approach is to design a source code such that for part of the time the source sequence is reconstructed correctly, while for the other part of the time the source sequence is reconstructed incorrectly, i.e., an error occurs. In designing such a code, we try to minimize the probability of error. However, this approach is not viable asymptotically because the converse of the source coding theorem says that if R < H, then the probability of error inevitably tends to 1 as  $n \to \infty$ .

Therefore, if R < H, no matter how the source code is designed, the source sequence is almost always reconstructed incorrectly when n is large. An alternative approach is to design a source code called a *rate-distortion code* which reproduces the source sequence with distortion. In order to formulate the problem properly, we need a *distortion measure* between each source sequence and each reproduction sequence. Then we try to design a rate-distortion code which with high probability reproduces the source sequence with a distortion within a tolerance level.

Clearly, a smaller distortion can potentially be achieved if we are allowed to use a higher coding rate. *Rate-distortion theory*, the subject matter of this chapter, gives a characterization of the asymptotic optimal tradeoff between the coding rate of a rate-distortion code for a given information source and the allowed distortion in the reproduction sequence with respect to a distortion measure.

## 8.1 Single-Letter Distortion Measures

Let  $\{X_k, k \ge 1\}$  be an i.i.d. information source with generic random variable X. We assume that the source alphabet  $\mathcal{X}$  is finite. Let p(x) be the probability distribution of X, and we assume without loss of generality that the support of X is equal to  $\mathcal{X}$ . Consider a source sequence

$$\mathbf{x} = (x_1, x_2, \cdots, x_n) \tag{8.1}$$

and a reproduction sequence

$$\hat{\mathbf{x}} = (\hat{x}_1, \hat{x}_2, \cdots, \hat{x}_n). \tag{8.2}$$

The components of  $\hat{\mathbf{x}}$  can take values in  $\mathcal{X}$ , but more generally, they can take values in any finite set  $\hat{\mathcal{X}}$  which may be different from  $\mathcal{X}$ . The set  $\hat{\mathcal{X}}$ , which is also assumed to be finite, is called the reproduction alphabet. To measure the distortion between  $\mathbf{x}$  and  $\hat{\mathbf{x}}$ , we introduce the single-letter distortion measure and the average distortion measure.

**Definition 8.1.** A single-letter distortion measure is a mapping

$$d: \mathcal{X} \times \hat{\mathcal{X}} \to \Re^+, \tag{8.3}$$

where  $\Re^+$  is the set of nonnegative real numbers<sup>1</sup>. The value  $d(x, \hat{x})$  denotes the distortion incurred when a source symbol x is reproduced as  $\hat{x}$ .

**Definition 8.2.** The average distortion between a source sequence  $\mathbf{x} \in \mathcal{X}^n$ and a reproduction sequence  $\hat{\mathbf{x}} \in \hat{\mathcal{X}}^n$  induced by a single-letter distortion measure d is defined by

$$d(\mathbf{x}, \hat{\mathbf{x}}) = \frac{1}{n} \sum_{k=1}^{n} d(x_k, \hat{x}_k).$$
(8.4)

In Definition 8.2, we have used d to denote both the single-letter distortion measure and the average distortion measure, but this abuse of notation should cause no ambiguity. Henceforth, we will refer to a single-letter distortion measure simply as a distortion measure.

Very often, the source sequence  $\mathbf{x}$  represents quantized samples of a continuous signal, and the user attempts to recognize certain objects and derive meaning from the reproduction sequence  $\hat{\mathbf{x}}$ . For example,  $\mathbf{x}$  may represent a video signal, an audio signal, or an image. The ultimate purpose of a distortion measure is to reflect the distortion between  $\mathbf{x}$  and  $\hat{\mathbf{x}}$  as *perceived* by the user. This goal is difficult to achieve in general because measurements of the distortion between  $\mathbf{x}$  and  $\hat{\mathbf{x}}$  must be made within context unless the symbols in  $\mathcal{X}$  carry no physical meaning. Specifically, when the user derives meaning

<sup>&</sup>lt;sup>1</sup> Note that  $d(x, \hat{x})$  is finite for all  $(x, \hat{x}) \in \mathcal{X} \times \hat{\mathcal{X}}$ .

from  $\hat{\mathbf{x}}$ , the distortion in  $\hat{\mathbf{x}}$  as perceived by the user depends on the context. For example, the perceived distortion is small for a portrait contaminated by a fairly large noise, while the perceived distortion is large for the image of a book page contaminated by the same noise. Hence, a good distortion measure should be context dependent.

Although the average distortion is not necessarily the best way to measure the distortion between a source sequence and a reproduction sequence, it has the merit of being simple and easy to use. Moreover, rate-distortion theory, which is based on the average distortion measure, provides a framework for data compression when distortion is inevitable.

*Example 8.3.* When the symbols in  $\mathcal{X}$  and  $\hat{\mathcal{X}}$  represent real values, a popular distortion measure is the square-error distortion measure which is defined by

$$d(x, \hat{x}) = (x - \hat{x})^2.$$
(8.5)

The average distortion measure so induced is often referred to as the meansquare error.

*Example 8.4.* When  $\mathcal{X}$  and  $\hat{\mathcal{X}}$  are identical and the symbols in  $\mathcal{X}$  do not carry any particular meaning, a frequently used distortion measure is the Hamming distortion measure, which is defined by

$$d(x, \hat{x}) = \begin{cases} 0 \text{ if } x = \hat{x} \\ 1 \text{ if } x \neq \hat{x}. \end{cases}$$
(8.6)

The Hamming distortion measure indicates the occurrence of an error. In particular, for an estimate  $\hat{X}$  of X, we have

$$Ed(X, \hat{X}) = \Pr\{X = \hat{X}\} \cdot 0 + \Pr\{X \neq \hat{X}\} \cdot 1 = \Pr\{X \neq \hat{X}\}, \quad (8.7)$$

i.e., the expectation of the Hamming distortion measure between X and  $\hat{X}$  is the probability of error.

For  $\mathbf{x} \in \mathcal{X}^n$  and  $\hat{\mathbf{x}} \in \hat{\mathcal{X}}^n$ , the average distortion  $d(\mathbf{x}, \hat{\mathbf{x}})$  induced by the Hamming distortion measure gives the frequency of error in the reproduction sequence  $\hat{\mathbf{x}}$ .

**Definition 8.5.** For a distortion measure d, for each  $x \in \mathcal{X}$ , let  $\hat{x}^*(x) \in \hat{\mathcal{X}}$ minimize  $d(x, \hat{x})$  over all  $\hat{x} \in \hat{\mathcal{X}}$ . A distortion measure d is said to be normal if

$$c_x \stackrel{\text{def}}{=} d(x, \hat{x}^*(x)) = 0 \tag{8.8}$$

for all  $x \in \mathcal{X}$ .

The square-error distortion measure and the Hamming distortion measure are examples of normal distortion measures. Basically, a normal distortion measure is one which allows X to be reproduced with zero distortion. Although

a distortion measure d is not normal in general, a normalization of d can always be obtained by defining the distortion measure

$$\tilde{d}(x,\hat{x}) = d(x,\hat{x}) - c_x \tag{8.9}$$

for all  $(x, \hat{x}) \in \mathcal{X} \times \hat{\mathcal{X}}$ . Evidently,  $\tilde{d}$  is a normal distortion measure, and it is referred to as the normalization of d.

*Example 8.6.* Let d be a distortion measure defined by

Then  $\tilde{d}$ , the normalization of d, is given by

$$\begin{array}{c|c} \tilde{d}(x,\hat{x}) & a & b & c \\ \hline 1 & 0 & 5 & 3 \\ 2 & 1 & 0 & 5 \end{array}$$

Note that for every  $x \in \mathcal{X}$ , there exists an  $\hat{x} \in \hat{\mathcal{X}}$  such that  $\tilde{d}(x, \hat{x}) = 0$ .

Let  $\hat{X}$  be any estimate of X which takes values in  $\hat{\mathcal{X}}$ , and denote the joint distribution for X and  $\hat{X}$  by  $p(x, \hat{x})$ . Then

$$Ed(X, \hat{X}) = \sum_{x} \sum_{\hat{x}} p(x, \hat{x}) d(x, \hat{x})$$
 (8.10)

$$=\sum_{x}\sum_{\hat{x}}p(x,\hat{x})\left[\tilde{d}(x,\hat{x})+c_{x}\right]$$
(8.11)

$$= E\tilde{d}(X, \hat{X}) + \sum_{x} p(x) \sum_{\hat{x}} p(\hat{x}|x) c_x$$
(8.12)

$$= E\tilde{d}(X,\hat{X}) + \sum_{x} p(x)c_x\left(\sum_{\hat{x}} p(\hat{x}|x)\right)$$
(8.13)

$$= E\tilde{d}(X,\hat{X}) + \sum_{x} p(x)c_x \tag{8.14}$$

$$= E\tilde{d}(X,\hat{X}) + \Delta, \tag{8.15}$$

where

$$\Delta = \sum_{x} p(x)c_x \tag{8.16}$$

is a constant which depends only on p(x) and d but not on the conditional distribution  $p(\hat{x}|x)$ . In other words, for a given X and a distortion measure d, the expected distortion between X and an estimate  $\hat{X}$  of X is always reduced by a constant upon using  $\tilde{d}$  instead of d as the distortion measure. For reasons which will be explained in Section 8.3, it is sufficient for us to assume that a distortion measure is normal.

**Definition 8.7.** Let  $\hat{x}^*$  minimizes  $Ed(X, \hat{x})$  over all  $\hat{x} \in \hat{\mathcal{X}}$ , and define

$$D_{max} = Ed(X, \hat{x}^*).$$
 (8.17)

 $\hat{x}^*$  is the best estimate of X if we know nothing about X, and  $D_{max}$  is the minimum expected distortion between X and a constant estimate of X. The significance of  $D_{max}$  can be seen by taking the reproduction sequence  $\hat{\mathbf{X}}$  to be  $(\hat{x}^*, \hat{x}^*, \dots, \hat{x}^*)$ . Since  $d(X_k, \hat{x}^*)$  are i.i.d., by the weak law of large numbers

$$d(\mathbf{X}, \hat{\mathbf{X}}) = \frac{1}{n} \sum_{k=1}^{n} d(X_k, \hat{x}^*) \to Ed(X, \hat{x}^*) = D_{max}$$
(8.18)

in probability, i.e., for any  $\epsilon > 0$ ,

$$\Pr\{d(\mathbf{X}, \mathbf{X}) > D_{max} + \epsilon\} \le \epsilon \tag{8.19}$$

for sufficiently large n. Note that  $\hat{\mathbf{X}}$  is a constant sequence which does not depend on  $\mathbf{X}$ . In other words, even when no description of  $\mathbf{X}$  is available, we can still achieve an average distortion no more than  $D_{max} + \epsilon$  with probability arbitrarily close to 1 when n is sufficiently large.

The notation  $D_{max}$  may seem confusing because the quantity stands for the minimum rather than the maximum expected distortion between X and a constant estimate of X. But we see from the above discussion that this notation is in fact appropriate because  $D_{max}$  is the maximum distortion we have to be concerned about. Specifically, it is not meanful to impose a constraint  $D \ge D_{max}$  on the reproduction sequence because it can be achieved even without receiving any information about the sequence produced by the source.

## 8.2 The Rate-Distortion Function R(D)

Throughout this chapter, all the discussions are with respect to an i.i.d. information source  $\{X_k, k \ge 1\}$  with generic random variable X and a distortion measure d. All logarithms are in the base 2 unless otherwise specified.

**Definition 8.8.** An (n, M) rate-distortion code is defined by an encoding function

$$f: \mathcal{X}^n \to \{1, 2, \cdots, M\} \tag{8.20}$$

and a decoding function

$$g: \{1, 2, \cdots, M\} \to \hat{\mathcal{X}}^n. \tag{8.21}$$

The set  $\{1, 2, \dots, M\}$ , denoted by  $\mathcal{I}$ , is called the index set. The reproduction sequences  $g(f(1)), g(f(2)), \dots, g(f(M))$  in  $\hat{X}^n$  are called codewords, and the set of codewords is called the codebook.

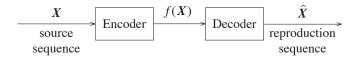


Fig. 8.1. A rate-distortion code with block length n.

Figure 8.1 is an illustration of a rate-distortion code.

**Definition 8.9.** The rate of an (n, M) rate-distortion code is  $n^{-1} \log M$  in bits per symbol.

**Definition 8.10.** A rate-distortion pair (R, D) is asymptotically achievable if for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M) rate-distortion code such that

$$\frac{1}{n}\log M \le R + \epsilon \tag{8.22}$$

and

$$\Pr\{d(\mathbf{X}, \mathbf{X}) > D + \epsilon\} \le \epsilon, \tag{8.23}$$

where  $\hat{\mathbf{X}} = g(f(\mathbf{X}))$ . For brevity, an asymptotically achievable pair will be referred to as an achievable pair.

**Remark** It is clear from the definition that if (R, D) is achievable, then (R', D) and (R, D') are also achievable for all  $R' \ge R$  and  $D' \ge D$ .

**Definition 8.11.** The rate-distortion region is the subset of  $\Re^2$  containing all achievable pairs (R, D).

**Theorem 8.12.** The rate-distortion region is closed and convex.

*Proof.* We first show that the rate-distortion region is closed. Consider achievable rate-distortion pairs  $(R^{(k)}, D^{(k)})$  such that

$$\lim_{k \to \infty} (R^{(k)}, D^{(k)}) = (R, D).$$
(8.24)

Then for any  $\epsilon > 0$ , for all k, there exists for sufficiently large n an  $(n, M^{(k)})$  code such that

$$\frac{1}{n}\log M^{(k)} \le R^{(k)} + \epsilon \tag{8.25}$$

and

$$\Pr\{d(\mathbf{X}^{(k)}, \hat{\mathbf{X}}^{(k)}) > D^{(k)} + \epsilon\} \le \epsilon,$$
(8.26)

where  $f^{(k)}$  and  $g^{(k)}$  are respectively the encoding function and the decoding function of the  $(n, M^{(k)})$  code, and  $\hat{\mathbf{X}}^{(k)} = g^{(k)}(f^{(k)}(\mathbf{X}))$ . By virtue of (8.24), let  $k(\epsilon)$  be an integer such that for all  $k > k(\epsilon)$ ,

$$|R - R^{(k)}| < \epsilon \tag{8.27}$$

8.2 The Rate-Distortion Function R(D) 185

and

$$|D - D^{(k)}| < \epsilon, \tag{8.28}$$

which imply

$$R^{(k)} < R + \epsilon \tag{8.29}$$

and

$$D^{(k)} < D + \epsilon, \tag{8.30}$$

respectively. Then for all  $k > k(\epsilon)$ ,

$$\frac{1}{n}\log M^{(k)} \le R^{(k)} + \epsilon < R + 2\epsilon \tag{8.31}$$

and

$$\Pr\{d(\mathbf{X}^{(k)}, \hat{\mathbf{X}}^{(k)}) > D + 2\epsilon\} \le \Pr\{d(\mathbf{X}^{(k)}, \hat{\mathbf{X}}^{(k)}) > D^{(k)} + \epsilon\}$$
(8.32)

$$\leq \epsilon.$$
 (8.33)

Note that (8.32) follows because

$$D + 2\epsilon > D^{(k)} + \epsilon \tag{8.34}$$

by (8.30). From (8.31) and (8.33), we see that (R, D) is also achievable. Thus we have proved that the rate-distortion region is closed.

We will prove the convexity of the rate-distortion region by a time-sharing argument whose idea is the following. Roughly speaking, if we can use a code  $C_1$  to achieve  $(R^{(1)}, D^{(1)})$  and a code  $C_2$  to achieve  $(R^{(2)}, D^{(2)})$ , then for any rational number  $\lambda$  between 0 and 1, we can use  $C_1$  for a fraction  $\lambda$  of the time and use  $C_2$  for a fraction  $\bar{\lambda}$  of the time to achieve  $(R^{(\lambda)}, D^{(\lambda)})$ , where

$$R^{(\lambda)} = \lambda R^{(1)} + \bar{\lambda} R^{(2)}, \qquad (8.35)$$

$$D^{(\lambda)} = \lambda D^{(1)} + \bar{\lambda} D^{(2)}, \qquad (8.36)$$

and  $\bar{\lambda} = 1 - \lambda$ . Since the rate-distortion region is closed as we have proved,  $\lambda$  can be taken as any real number between 0 and 1, and the convexity of the region follows.

We now give a formal proof for the convexity of the rate-distortion region. Let

$$\lambda = \frac{r}{r+s},\tag{8.37}$$

where r and s are positive integers. Then  $\lambda$  is a rational number between 0 and 1. We now prove that if  $(R^{(1)}, D^{(1)})$  and  $(R^{(2)}, D^{(2)})$  are achievable, then  $(R^{(\lambda)}, D^{(\lambda)})$  is also achievable. Assume  $(R^{(1)}, D^{(1)})$  and  $(R^{(2)}, D^{(2)})$  are achievable. Then for any  $\epsilon > 0$  and sufficiently large n, there exist an  $(n, M^{(1)})$  code and an  $(n, M^{(2)})$  code such that

$$\frac{1}{n}\log M^{(i)} \le R^{(i)} + \epsilon \tag{8.38}$$

and

$$\Pr\{d(\mathbf{X}, \hat{\mathbf{X}}^{(i)}) > D^{(i)} + \epsilon\} \le \epsilon,$$
(8.39)

$$i = 1, 2$$
. Let

$$M(\lambda) = (M^{(1)})^r (M^{(2)})^s \tag{8.40}$$

and

$$n(\lambda) = (r+s)n. \tag{8.41}$$

We now construct an  $(n(\lambda), M(\lambda))$  code by concatenating r copies of the  $(n, M^{(1)})$  code followed by s copies of the  $(n, M^{(2)})$  code. We call these r + s codes subcodes of the  $(n(\lambda), M(\lambda))$  code. For this code, let

$$\mathbf{Y} = (\mathbf{X}(1), \mathbf{X}(2), \cdots, \mathbf{X}(r+s)) \tag{8.42}$$

and

$$\hat{\mathbf{Y}} = (\hat{\mathbf{X}}(1), \hat{\mathbf{X}}(2), \cdots, \hat{\mathbf{X}}(r+s)), \qquad (8.43)$$

where  $\mathbf{X}(j)$  and  $\hat{\mathbf{X}}(j)$  are the source sequence and the reproduction sequence of the *j*th subcode, respectively. Then for this  $(n(\lambda), M(\lambda))$  code,

$$\frac{1}{n(\lambda)}\log M(\lambda) = \frac{1}{(r+s)n}\log[(M^{(1)})^r (M^{(2)})^s]$$
(8.44)

$$= \frac{1}{(r+s)n} (r \log M^{(1)} + s \log M^{(2)})$$
(8.45)

$$= \lambda \left(\frac{1}{n} \log M^{(1)}\right) + \bar{\lambda} \left(\frac{1}{n} \log M^{(2)}\right)$$
(8.46)

$$\leq \lambda(R^{(1)} + \epsilon) + \bar{\lambda}(R^{(2)} + \epsilon) \tag{8.47}$$

$$= (\lambda R^{(1)} + \lambda R^{(2)}) + \epsilon \tag{8.48}$$

$$=R^{(\lambda)}+\epsilon, \tag{8.49}$$

where (8.47) follows from (8.38), and

$$\Pr\{d(\mathbf{Y}, \hat{\mathbf{Y}}) > D^{(\lambda)} + \epsilon\} = \Pr\left\{\frac{1}{r+s} \sum_{j=1}^{r+s} d(\mathbf{X}(j), \hat{\mathbf{X}}(j)) > D^{(\lambda)} + \epsilon\right\}$$
(8.50)

$$\leq \Pr\left\{ d(\mathbf{X}(j), \hat{\mathbf{X}}(j)) > D^{(1)} + \epsilon \text{ for some } 1 \leq j \leq r \text{ or} \\ d(\mathbf{X}(j), \hat{\mathbf{X}}(j)) > D^{(2)} + \epsilon \text{ for some } r + 1 \leq j \leq r + s \right\}$$
(8.51)

$$\leq \sum_{j=1}^{r} \Pr\{d(\mathbf{X}(j), \hat{\mathbf{X}}(j)) > D^{(1)} + \epsilon\} + \sum_{j=r+1}^{r+s} \Pr\{d(\mathbf{X}(j), \hat{\mathbf{X}}(j)) > D^{(2)} + \epsilon\}$$
(8.52)

$$\leq (r+s)\epsilon,\tag{8.53}$$

where (8.52) follows from the union bound and (8.53) follows from (8.39). Hence, we conclude that the rate-distortion pair  $(R^{(\lambda)}, D^{(\lambda)})$  is achievable. This completes the proof of the theorem.  $\Box$ 

**Definition 8.13.** The rate-distortion function R(D) is the minimum of all rates R for a given distortion D such that (R, D) is achievable.

**Definition 8.14.** The distortion-rate function D(R) is the minimum of all distortions D for a given rate R such that (R, D) is achievable.

Both the functions R(D) and D(R) are equivalent descriptions of the boundary of the rate-distortion region. They are sufficient to describe the rate-distortion region because the region is closed. Note that in defining R(D), the minimum instead of the infimum is taken because for a fixed D, the set of all R such that (R, D) is achievable is closed and lower bounded by zero. Similarly, the minimum instead of the infimum is taken in defining D(R). In the subsequent discussions, only R(D) will be used.

**Theorem 8.15.** The following properties hold for the rate-distortion function R(D):

1. R(D) is non-increasing in D. 2. R(D) is convex. 3. R(D) = 0 for  $D \ge D_{max}$ . 4.  $R(0) \le H(X)$ .

*Proof.* From the remark following Definition 8.10, since (R(D), D) is achievable, (R(D), D') is also achievable for all  $D' \ge D$ . Therefore,  $R(D) \ge R(D')$  because R(D') is the minimum of all R such that (R, D') is achievable. This proves Property 1.

Property 2 follows immediately from the convexity of the rate-distortion region which was proved in Theorem 8.12. From the discussion toward the end of the last section, we see for any  $\epsilon > 0$ , it is possible to achieve

$$\Pr\{d(\mathbf{X}, \mathbf{X}) > D_{max} + \epsilon\} \le \epsilon \tag{8.54}$$

for sufficiently large n with no description of **X** available. Therefore, (0, D) is achievable for all  $D \ge D_{max}$ , proving Property 3.

Property 4 is a consequence of the assumption that the distortion measure d is normalized, which can be seen as follows. By the source coding theorem, for any  $\epsilon > 0$ , by using a rate no more than  $H(X) + \epsilon$ , we can describe the source sequence **X** of length n with probability of error less than  $\epsilon$  when n is sufficiently large. Since d is normalized, for each  $k \geq 1$ , let

$$\hat{X}_k = \hat{x}^*(X_k) \tag{8.55}$$

(cf. Definition 8.5), so that whenever an error does not occur,

$$d(X_k, \hat{X}_k) = d(X_k, \hat{x}^*(X_k)) = 0$$
(8.56)

by (8.8) for each k, and

$$d(\mathbf{X}, \hat{\mathbf{X}}) = \frac{1}{n} \sum_{k=1}^{n} d(X_k, \hat{X}_k) = \frac{1}{n} \sum_{k=1}^{n} d(X_k, \hat{x}^*(X_k)) = 0.$$
(8.57)

Therefore,

$$\Pr\{d(\mathbf{X}, \mathbf{X}) > \epsilon\} \le \epsilon, \tag{8.58}$$

which shows that the pair (H(X), 0) is achievable. This in turn implies that  $R(0) \leq H(X)$  because R(0) is the minimum of all R such that (R, 0) is achievable.  $\Box$ 

Figure 8.2 is an illustration of a rate-distortion function R(D). The reader

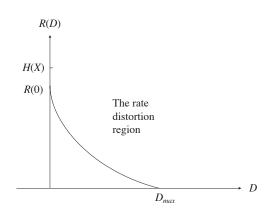


Fig. 8.2. A rate-distortion function R(D).

should note the four properties of R(D) in Theorem 8.15. The rate-distortion theorem, which will be stated in the next section, gives a characterization of R(D).

## 8.3 The Rate-Distortion Theorem

**Definition 8.16.** For  $D \ge 0$ , the information rate-distortion function is defined by

$$R_I(D) = \min_{\hat{X}: Ed(X, \hat{X}) \le D} I(X; \hat{X}).$$
(8.59)

In defining  $R_I(D)$ , the minimization is taken over all random variables  $\hat{X}$  jointly distributed with X such that

$$Ed(X, \hat{X}) \le D. \tag{8.60}$$

Since p(x) is given, the minimization is taken over the set of all  $p(\hat{x}|x)$  such that (8.60) is satisfied, namely the set

$$\left\{ p(\hat{x}|x) : \sum_{x,\hat{x}} p(x)p(\hat{x}|x)d(x,\hat{x}) \le D \right\}.$$
 (8.61)

Since this set is compact in  $\Re^{|\mathcal{X}||\hat{\mathcal{X}}|}$  and  $I(X; \hat{X})$  is a continuous functional of  $p(\hat{x}|x)$ , the minimum value of  $I(X; \hat{X})$  can be attained<sup>2</sup>. This justifies taking the minimum instead of the infimum in the definition of  $R_I(D)$ .

We have seen in Section 8.1 that we can obtain a normalization d for any distortion measure d with

$$E\tilde{d}(X,\hat{X}) = Ed(X,\hat{X}) - \Delta$$
(8.62)

for any  $\hat{X}$ , where  $\Delta$  is a constant which depends only on p(x) and d. Thus if d is not normal, we can always replace d by  $\tilde{d}$  and D by  $D - \Delta$  in the definition of  $R_I(D)$  without changing the minimization problem. Therefore, we do not lose any generality by assuming that a distortion measure d is normal.

## Theorem 8.17 (The Rate-Distortion Theorem). $R(D) = R_I(D)$ .

The rate-distortion theorem, which is the main result in rate-distortion theory, says that the minimum coding rate for achieving a distortion D is  $R_I(D)$ . This theorem will be proved in the next two sections. In the next section, we will prove the converse of this theorem, i.e.,  $R(D) \ge R_I(D)$ , and in Section 8.5, we will prove the achievability of  $R_I(D)$ , i.e.,  $R(D) \le R_I(D)$ .

In order for  $R_I(D)$  to be a characterization of R(D), it has to satisfy the same properties as R(D). In particular, the four properties of R(D) in Theorem 8.15 should also be satisfied by  $R_I(D)$ .

**Theorem 8.18.** The following properties hold for the information rate-distortion function  $R_I(D)$ :

1.  $R_I(D)$  is non-increasing in D. 2.  $R_I(D)$  is convex. 3.  $R_I(D) = 0$  for  $D \ge D_{max}$ . 4.  $R_I(0) \le H(X)$ .

<sup>&</sup>lt;sup>2</sup> The assumption that both  $\mathcal{X}$  and  $\hat{\mathcal{X}}$  are finite is essential in this argument.

*Proof.* Referring to the definition of  $R_I(D)$  in (8.59), for a larger D, the minimization is taken over a larger set. Therefore,  $R_I(D)$  is non-increasing in D, proving Property 1.

To prove Property 2, consider any  $D^{(1)}, D^{(2)} \ge 0$  and let  $\lambda$  be any number between 0 and 1. Let  $\hat{X}^{(i)}$  achieves  $R_I(D^{(i)})$  for i = 1, 2, i.e.,

$$R_I(D^{(i)}) = I(X; \hat{X}^{(i)}), \tag{8.63}$$

where

$$Ed(X, \hat{X}^{(i)}) \le D^{(i)},$$
 (8.64)

and let  $\hat{X}^{(i)}$  be defined by the transition matrix  $p_i(\hat{x}|x)$ . Let  $\hat{X}^{(\lambda)}$  be jointly distributed with X which is defined by

$$p_{\lambda}(\hat{x}|x) = \lambda p_1(\hat{x}|x) + \bar{\lambda} p_2(\hat{x}|x), \qquad (8.65)$$

where  $\bar{\lambda} = 1 - \lambda$ . Then

$$Ed(X, \hat{X}^{(\lambda)}) = \sum_{x, \hat{x}} p(x) p_{\lambda}(\hat{x}|x) d(x, \hat{x})$$
(8.66)

$$= \sum_{x,\hat{x}} p(x)(\lambda p_1(\hat{x}|x) + \bar{\lambda} p_2(\hat{x}|x))d(x,\hat{x})$$
(8.67)

$$=\lambda\left(\sum_{x,\hat{x}}p(x)p_1(\hat{x}|x)d(x,\hat{x})\right)+\bar{\lambda}\left(\sum_{x,\hat{x}}p(x)p_2(\hat{x}|x)d(x,\hat{x})\right)$$
(8.68)

$$= \lambda E d(X, \hat{X}^{(1)}) + \bar{\lambda} E d(X, \hat{X}^{(2)})$$
(8.69)
(8.69)

$$\leq \lambda D^{(2)} + \lambda D^{(2)} \tag{8.70}$$
$$= D^{(\lambda)}, \tag{8.71}$$

where

$$D^{(\lambda)} = \lambda D^{(1)} + \bar{\lambda} D^{(2)}, \qquad (8.72)$$

and (8.70) follows from (8.64). Now consider

$$\lambda R_I(D^{(1)}) + \bar{\lambda} R_I(D^{(2)}) = \lambda I(X; \hat{X}^{(1)}) + \bar{\lambda} I(X; \hat{X}^{(2)})$$
(8.73)

$$\geq I(X; \hat{X}^{(\lambda)}) \tag{8.74}$$

$$\geq R_I(D^{(\lambda)}),\tag{8.75}$$

where the inequality in (8.74) follows from the convexity of mutual information with respect to the transition matrix  $p(\hat{x}|x)$  (see Example 3.13), and the inequality in (8.75) follows from (8.71) and the definition of  $R_I(D)$ . Therefore, we have proved Property 2.

To prove Property 3, let  $\hat{X}$  take the value  $\hat{x}^*$  as defined in Definition 8.7 with probability 1. Then

8.3 The Rate-Distortion Theorem 191

$$I(X; \hat{X}) = 0 (8.76)$$

and

$$Ed(X; \hat{X}) = Ed(X; \hat{x}^*) = D_{max}.$$
 (8.77)

Then for  $D \geq D_{max}$ ,

$$R_I(D) \le I(X;X) = 0.$$
 (8.78)

On the other hand, since  $R_I(D)$  is nonnegative, we conclude that

$$R_I(D) = 0.$$
 (8.79)

This proves Property 3.

Finally, to prove Property 4, we let

$$\hat{X} = \hat{x}^*(X),$$
 (8.80)

where  $\hat{x}^*(x)$  is defined in Definition 8.5. Then

$$Ed(X, \hat{X}) = Ed(X, \hat{x}^*(X))$$
 (8.81)

$$=\sum_{x} p(x)d(x, \hat{x}^{*}(x))$$
(8.82)

$$= 0$$
 (8.83)

by (8.8) since we assume that d is a normal distortion measure. Moreover,

$$R_I(0) \le I(X; \hat{X}) \le H(X).$$
 (8.84)

Then Property 4 and hence the theorem is proved.  $\Box$ 

**Corollary 8.19.** If  $R_I(0) > 0$ , then  $R_I(D)$  is strictly decreasing for  $0 \le D \le D_{max}$ , and the inequality constraint in Definition 8.16 for  $R_I(D)$  can be replaced by an equality constraint.

*Proof.* Assume that  $R_I(0) > 0$ . We first show that  $R_I(D) > 0$  for  $0 \le D < D_{max}$  by contradiction. Suppose  $R_I(D') = 0$  for some  $0 \le D' < D_{max}$ , and let  $R_I(D')$  be achieved by some  $\hat{X}$ . Then

$$R_I(D') = I(X; \hat{X}) = 0 \tag{8.85}$$

implies that X and  $\hat{X}$  are independent, or

$$p(x, \hat{x}) = p(x)p(\hat{x})$$
 (8.86)

for all x and  $\hat{x}$ . It follows that

$$D' \ge Ed(X, \hat{X}) \tag{8.87}$$

$$=\sum_{x}\sum_{\hat{x}} p(x,\hat{x})d(x,\hat{x})$$
(8.88)

$$= \sum_{x} \sum_{\hat{x}} p(x) p(\hat{x}) d(x, \hat{x})$$
(8.89)

$$= \sum_{\hat{x}} p(\hat{x}) \sum_{x} p(x) d(x, \hat{x})$$
(8.90)

$$=\sum_{\hat{x}} p(\hat{x}) E d(X, \hat{x}) \tag{8.91}$$

$$\geq \sum_{\hat{x}} p(\hat{x}) E d(X, \hat{x}^*) \tag{8.92}$$

$$=\sum_{\hat{x}} p(\hat{x}) D_{max} \tag{8.93}$$

$$= D_{max}, \tag{8.94}$$

where  $\hat{x}^*$  and  $D_{max}$  are defined in Definition 8.7. This leads to a contradiction because we have assumed that  $0 \leq D' < D_{max}$ . Therefore, we conclude that  $R_I(D) > 0$  for  $0 \leq D < D_{max}$ .

Since  $R_I(0) > 0$  and  $R_I(D_{max}) = 0$ , and  $R_I(D)$  is non-increasing and convex from the above theorem,  $R_I(D)$  must be strictly decreasing for  $0 \le D \le D_{max}$ . We now prove by contradiction that the inequality constraint in Definition 8.16 for  $R_I(D)$  can be replaced by an equality constraint. Assume that  $R_I(D)$  is achieved by some  $\hat{X}^*$  such that

$$Ed(X, \hat{X}^*) = D'' < D.$$
 (8.95)

Then

$$R_I(D'') = \min_{\hat{X}: Ed(X, \hat{X}) \le D''} I(X; \hat{X}) \le I(X; \hat{X}^*) = R_I(D).$$
(8.96)

This is a contradiction because  $R_I(D)$  is strictly decreasing for  $0 \le D \le D_{max}$ . Hence,

$$Ed(X, \hat{X}^*) = D.$$
 (8.97)

This implies that the inequality constraint in Definition 8.16 for  $R_I(D)$  can be replaced by an equality constraint.  $\Box$ 

**Remark** In all problems of interest,  $R(0) = R_I(0) > 0$ . Otherwise, R(D) = 0 for all  $D \ge 0$  because R(D) is nonnegative and non-increasing.

Example 8.20 (Binary Source). Let X be a binary random variable with

$$\Pr\{X=0\} = 1 - \gamma \text{ and } \Pr\{X=1\} = \gamma.$$
 (8.98)

Let  $\hat{\mathcal{X}} = \{0, 1\}$  be the reproduction alphabet for X, and let d be the Hamming distortion measure. We first consider the case that  $0 \le \gamma \le \frac{1}{2}$ . Then if we make

a guess on the value of X, we should guess 0 in order to minimize the expected distortion. Therefore,  $\hat{x}^* = 0$  and

$$D_{max} = Ed(X,0) \tag{8.99}$$

$$= \Pr\{X = 1\} \tag{8.100}$$

$$=\gamma.$$
 (8.101)

We will show that for  $0 \le \gamma \le \frac{1}{2}$ ,

$$R_I(D) = \begin{cases} h_b(\gamma) - h_b(D) \text{ if } 0 \le D < \gamma \\ 0 & \text{if } D \ge \gamma. \end{cases}$$
(8.102)

Let  $\hat{X}$  be an estimate of X taking values in  $\hat{\mathcal{X}}$ , and let Y be the Hamming distortion measure between X and  $\hat{X}$ , i.e.,

$$Y = d(X, \tilde{X}).$$
 (8.103)

Observe that conditioning on  $\hat{X}$ , X and Y determine each other. Therefore,

$$H(X|\hat{X}) = H(Y|\hat{X}).$$
 (8.104)

Then for  $D < \gamma = D_{max}$  and any  $\hat{X}$  such that

$$Ed(X, \hat{X}) \le D,\tag{8.105}$$

we have

$$I(X; \hat{X}) = H(X) - H(X|\hat{X})$$
(8.106)

$$= h_b(\gamma) - H(Y|X) \tag{8.107}$$

$$\geq h_b(\gamma) - H(Y) \tag{8.108}$$

$$= h_b(\gamma) - h_b(\Pr\{X \neq \hat{X}\})$$
(8.109)
$$= h_b(\gamma) - h_b(\Pr\{X \neq \hat{X}\})$$
(8.110)

$$\geq h_b(\gamma) - h_b(D), \tag{8.110}$$

where the last inequality is justified because

$$\Pr\{X \neq \hat{X}\} = Ed(X, \hat{X}) \le D \tag{8.111}$$

and  $h_b(a)$  is increasing for  $0 \le a \le \frac{1}{2}$ . Minimizing over all  $\hat{X}$  satisfying (8.105) in (8.110), we obtain the lower bound

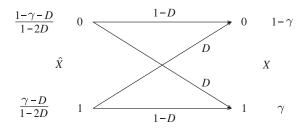
$$R_I(D) \ge h_b(\gamma) - h_b(D). \tag{8.112}$$

To show that this lower bound is achievable, we need to construct an  $\hat{X}$  such that the inequalities in both (8.108) and (8.110) are tight. The tightness of the inequality in (8.110) simply says that

$$\Pr\{X \neq \hat{X}\} = D,\tag{8.113}$$

while the tightness of the inequality in (8.108) says that Y should be independent of  $\hat{X}$ .

It would be more difficult to make Y independent of  $\hat{X}$  if we specify  $\hat{X}$  by  $p(\hat{x}|x)$ . Instead, we specify the joint distribution of X and  $\hat{X}$  by means of a reverse binary symmetric channel (BSC) with crossover probability D as the shown in Figure 8.3. Here, we regard  $\hat{X}$  as the input and X as the



**Fig. 8.3.** Achieving  $R_I(D)$  for a binary source via a reverse binary symmetric channel.

output of the BSC. Then Y is independent of the input  $\hat{X}$  because the error event is independent of the input for a BSC, and (8.113) is satisfied by setting the crossover probability to D. However, we need to ensure that the marginal distribution of X so specified is equal to p(x). Toward this end, we let

$$\Pr\{\hat{X} = 1\} = \alpha, \tag{8.114}$$

and consider

$$Pr\{X = 1\} = Pr\{\hat{X} = 0\}Pr\{X = 1|\hat{X} = 0\}$$
$$+Pr\{\hat{X} = 1\}Pr\{X = 1|\hat{X} = 1\},$$
(8.115)

 $\operatorname{or}$ 

$$\gamma = (1 - \alpha)D + \alpha(1 - D),$$
 (8.116)

which gives

$$\alpha = \frac{\gamma - D}{1 - 2D}.\tag{8.117}$$

Since

$$D < D_{max} = \gamma \le \frac{1}{2},\tag{8.118}$$

we have  $\alpha \geq 0$ . On the other hand,

$$\gamma, D \le \frac{1}{2} \tag{8.119}$$

gives

$$\gamma + D \le 1. \tag{8.120}$$

8.3 The Rate-Distortion Theorem 195

This implies

$$\gamma - D \le 1 - 2D, \tag{8.121}$$

or  $\alpha \leq 1$ . Therefore,

$$0 \le \alpha = \Pr\{\hat{X} = 1\} \le 1$$
 (8.122)

and

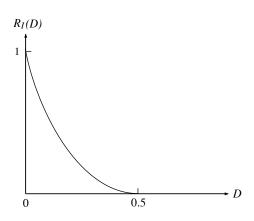
$$0 \le 1 - \alpha = \Pr\{X = 0\} \le 1.$$
(8.123)

Hence, we have shown that the lower bound on  $R_I(D)$  in (8.110) can be achieved, and  $R_I(D)$  is as given in (8.102).

For  $\frac{1}{2} \leq \gamma \leq 1$ , by exchanging the roles of the symbols 0 and 1 in the above argument, we obtain  $R_I(D)$  as in (8.102) except that  $\gamma$  is replaced by  $1 - \gamma$ . Combining the two cases, we have

$$R_I(D) = \begin{cases} h_b(\gamma) - h_b(D) \text{ if } 0 \le D < \min(\gamma, 1 - \gamma) \\ 0 \qquad \text{if } D \ge \min(\gamma, 1 - \gamma). \end{cases}$$
(8.124)

for  $0 \leq \gamma \leq 1$ . The function  $R_I(D)$  for  $\gamma = \frac{1}{2}$  is illustrated in Figure 8.4.



**Fig. 8.4.** The function  $R_I(D)$  for the uniform binary source with the Hamming distortion measure.

**Remark** In the above example, we see that  $R_I(0) = h_b(\gamma) = H(X)$ . Then by the rate-distortion theorem, H(X) is the minimum rate of a rate-distortion code which achieves an arbitrarily small average Hamming distortion. It is tempting to regarding this special case of the rate-distortion theorem as a version of the source coding theorem and conclude that the rate-distortion theorem is a generalization of the source coding theorem. However, this is incorrect because the rate-distortion theorem only guarantees that the *average* Hamming distortion between **X** and  $\hat{\mathbf{X}}$  is small with probability arbitrarily close to 1, but the source coding theorem guarantees that  $\mathbf{X} = \hat{\mathbf{X}}$  with probability arbitrarily close to 1, which is much stronger.

It is in general not possible to obtain the rate-distortion function in closed form, and we have to resort to numerical computation. In Chapter 9, we will discuss the Blahut-Arimoto algorithm for computing the rate-distortion function.

## 8.4 The Converse

In this section, we prove that the rate-distortion function R(D) is lower bounded by the information rate-distortion function  $R_I(D)$ , i.e.,  $R(D) \ge R_I(D)$ . Specifically, we will prove that for any achievable rate-distortion pair  $(R, D), R \ge R_I(D)$ . Then by fixing D and minimizing R over all achievable pairs (R, D), we conclude that  $R(D) \ge R_I(D)$ .

Let (R, D) be any achievable rate-distortion pair. Then for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M) code such that

$$\frac{1}{n}\log M \le R + \epsilon \tag{8.125}$$

and

$$\Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon\} \le \epsilon, \tag{8.126}$$

where  $\hat{\mathbf{X}} = g(f(\mathbf{X}))$ . Then

$$n(R+\epsilon) \stackrel{a)}{\geq} \log M \tag{8.127}$$
  
>  $H(f(\mathbf{X}))$   
(8.128)

$$\geq H(g(f(\mathbf{X})))$$
 (8.129)  
 $\geq H(g(f(\mathbf{X})))$ 

$$=H(\hat{\mathbf{X}}) \tag{8.130}$$

$$= H(\hat{\mathbf{X}}) - H(\hat{\mathbf{X}}|\mathbf{X})$$
(8.131)

$$= I(\hat{\mathbf{X}}; \mathbf{X}) \tag{8.132}$$

$$= H(\mathbf{X}) - H(\mathbf{X}|\hat{\mathbf{X}}) \tag{8.133}$$

$$=\sum_{k=1}^{n} H(X_k) - \sum_{k=1}^{n} H(X_k | \hat{\mathbf{X}}, X_1, X_2, \cdots, X_{k-1})$$
(8.134)

$$\stackrel{b)}{\geq} \sum_{k=1}^{n} H(X_k) - \sum_{k=1}^{n} H(X_k | \hat{X}_k)$$
(8.135)

$$=\sum_{k=1}^{n} [H(X_k) - H(X_k | \hat{X}_k)]$$
(8.136)

$$=\sum_{k=1}^{n} I(X_k; \hat{X}_k)$$
(8.137)

$$\stackrel{c)}{\geq} \sum_{k=1}^{n} R_I(Ed(X_k, \hat{X}_k))$$
(8.138)

8.4 The Converse 197

$$= n \left[ \frac{1}{n} \sum_{k=1}^{n} R_I(Ed(X_k, \hat{X}_k)) \right]$$
(8.139)

$$\stackrel{d)}{\geq} nR_I\left(\frac{1}{n}\sum_{k=1}^n Ed(X_k, \hat{X}_k)\right) \tag{8.140}$$

$$= nR_I(Ed(\mathbf{X}, \hat{\mathbf{X}})). \tag{8.141}$$

In the above,

- a) follows from (8.125);
- b) follows because conditioning does not increase entropy;
- c) follows from the definition of  $R_I(D)$  in Definition 8.16;
- d) follows from the convexity of  $R_I(D)$  proved in Theorem 8.18 and Jensen's inequality.

Now let

$$d_{max} = \max_{x,\hat{x}} d(x,\hat{x}) \tag{8.142}$$

be the maximum value which can be taken by the distortion measure d. The reader should not confuse  $d_{max}$  with  $D_{max}$  in Definition 8.7. Then from (8.126), we have

$$Ed(\mathbf{X}, \hat{\mathbf{X}})$$

$$= E[d(\mathbf{X}, \hat{\mathbf{X}})|d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon] \Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon\}$$

$$+ E[d(\mathbf{X}, \hat{\mathbf{X}})|d(\mathbf{X}, \hat{\mathbf{X}}) \le D + \epsilon] \Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) \le D + \epsilon\}$$
(8.143)

$$\leq d_{max} \cdot \epsilon + (D+\epsilon) \cdot 1 \tag{8.144}$$

$$= D + (d_{max} + 1)\epsilon. \tag{8.145}$$

This shows that if the probability that the average distortion between **X** and  $\hat{\mathbf{X}}$  exceeds  $D + \epsilon$  is small, then the expected average distortion between **X** and  $\hat{\mathbf{X}}$  can exceed D only by a small amount<sup>3</sup>. Following (8.141), we have

$$R + \epsilon \ge R_I(Ed(\mathbf{X}, \mathbf{\hat{X}})) \tag{8.146}$$

$$\geq R_I(D + (d_{max} + 1)\epsilon), \qquad (8.147)$$

where the last inequality follows from (8.145) because  $R_I(D)$  is non-increasing in D. We note that the convexity of  $R_I(D)$  implies that it is a continuous function of D. Then taking the limit as  $\epsilon \to 0$ , we obtain

$$R \ge \lim_{\epsilon \to 0} R_I (D + (d_{max} + 1)\epsilon) \tag{8.148}$$

$$=R_{I}\left(D+\left(d_{max}+1\right)\lim_{\epsilon\to0}\epsilon\right)$$
(8.149)

$$= R_I(D),$$
 (8.150)

=

 $<sup>^{3}</sup>$  The converse is not true.

where we have invoked the continuity of  $R_I(D)$  in obtaining (8.149). Upon minimizing R over all achievable pairs (R, D) for a fixed D in (8.150), we have proved that

$$R(D) \ge R_I(D). \tag{8.151}$$

This completes the proof for the converse of the rate-distortion theorem.

## 8.5 Achievability of $R_I(D)$

In this section, we prove that the rate-distortion function R(D) is upper bounded by the information rate-distortion function  $R_I(D)$ , i.e.,  $R(D) \leq R_I(D)$ . Then by combining with the result that  $R(D) \geq R_I(D)$  from the last section, we conclude that  $R(D) = R_I(D)$ , and the rate-distortion theorem is proved.

For any  $0 \leq D \leq D_{max}$ , we will prove that for every random variable  $\hat{X}$  taking values in  $\hat{\mathcal{X}}$  such that

$$Ed(X, \hat{X}) \le D,\tag{8.152}$$

the rate-distortion pair  $(I(X; \hat{X}), D)$  is achievable. This will be proved by showing for sufficiently large *n* the existence of a rate-distortion code such that

1. the rate of the code is not more than  $I(X; \hat{X}) + \epsilon$ ;

2.  $d(\mathbf{X}, \hat{\mathbf{X}}) \leq D + \epsilon$  with probability almost 1.

Then by minimizing  $I(X; \hat{X})$  over all  $\hat{X}$  satisfying (8.152), we conclude that the rate-distortion pair  $(R_I(D), D)$  is achievable, which implies  $R_I(D) \geq R(D)$  because R(D) is the minimum of all R such that (R, D) is achievable.

Fix any  $0 \leq D \leq D_{max}$  and any  $\epsilon > 0$ , and let  $\delta$  be a small positive quantity to be specified later. Toward proving the existence of a desired code, we fix a random variable  $\hat{X}$  which satisfies (8.152) and let M be an integer satisfying

$$I(X;\hat{X}) + \frac{\epsilon}{2} \le \frac{1}{n} \log M \le I(X;\hat{X}) + \epsilon, \qquad (8.153)$$

where n is sufficiently large.

We now describe a random coding scheme in the following steps:

- 1. Construct a codebook C of an (n, M) code by randomly generating M codewords in  $\hat{\mathcal{X}}^n$  independently and identically according to  $p(\hat{x})^n$ . Denote these codewords by  $\hat{\mathbf{X}}(1), \hat{\mathbf{X}}(2), \dots, \hat{\mathbf{X}}(M)$ .
- 2. Reveal the codebook C to both the encoder and the decoder.
- 3. The source sequence **X** is generated according to  $p(x)^n$ .
- 4. The encoder encodes the source sequence **X** into an index K in the set  $\mathcal{I} = \{1, 2, \dots, M\}$ . The index K takes the value *i* if

a) 
$$(\mathbf{X}, \mathbf{X}(i)) \in T^n_{[X\hat{X}]\delta}$$

b) for all  $i' \in \mathcal{I}$ , if  $(\mathbf{X}, \hat{\mathbf{X}}(i')) \in T^n_{[X\hat{X}]\delta}$ , then  $i' \leq i$ ;

otherwise, K takes the constant value 1.

- 5. The index K is delivered to the decoder.
- 6. The decoder outputs  $\mathbf{X}(K)$  as the reproduction sequence  $\mathbf{X}$ .

**Remark** Strong typicality is used in defining the encoding function in Step 4. This is made possible by the assumption that both the source alphabet  $\mathcal{X}$  and the reproduction alphabet  $\hat{\mathcal{X}}$  are finite.

Let us further explain the encoding scheme described in Step 4. After the source sequence  $\mathbf{X}$  is generated, we search through all the codewords in the codebook C for those which are jointly typical with  $\mathbf{X}$  with respect to  $p(x, \hat{x})$ . If there is at least one such codeword, we let i be the largest index of such codewords and let K = i. If such a codeword does not exist, we let K = 1.

The event  $\{K = 1\}$  occurs in one of the following two scenarios:

- 1.  $\hat{X}(1)$  is the only codeword in  $\mathcal{C}$  which is jointly typical with **X**.
- 2. No codeword in C is jointly typical with **X**.

In either scenario, **X** is not jointly typical with the codewords  $\hat{X}(2), \hat{X}(3), \dots, \hat{X}(M)$ . In other words, if K = 1, then **X** is jointly typical with none of the codewords  $\hat{X}(2), \hat{X}(3), \dots, \hat{X}(M)$ .

Define

$$E_{i} = \left\{ (\mathbf{X}, \hat{\mathbf{X}}(i)) \in T^{n}_{[X\hat{X}]\delta} \right\}$$
(8.154)

to be the event that **X** is jointly typical with the codeword  $\hat{X}(i)$ . We see from the above discussion that

$$\{K=1\} \subset E_2^c \cap E_3^c \cap \dots \cap E_M^c. \tag{8.155}$$

Since the codewords are generated i.i.d., conditioning on  $\{\mathbf{X} = \mathbf{x}\}$  for any  $\mathbf{x} \in \mathcal{X}^n$ , the events  $E_i$  are mutually independent<sup>4</sup>, and they all have the same probability. Then for any  $\mathbf{x} \in \mathcal{X}^n$ ,

$$\Pr\{K=1|\mathbf{X}=\mathbf{x}\} \le \Pr\{E_2^c \cap E_3^c \cap \dots \cap E_M^c | \mathbf{X}=\mathbf{x}\}$$
(8.156)

$$=\prod_{i=2}^{m} \Pr\{E_i^c | \mathbf{X} = \mathbf{x}\}$$
(8.157)

$$= (\Pr\{E_1^c | \mathbf{X} = \mathbf{x}\})^{M-1}$$
(8.158)

$$= (1 - \Pr\{E_1 | \mathbf{X} = \mathbf{x}\})^{M-1}.$$
 (8.159)

We now obtain a lower bound on  $\Pr\{E_1 | \mathbf{X} = \mathbf{x}\}$  for  $\mathbf{x} \in S_{[X]\delta}^n$ , where

$$S_{[X]\delta}^{n} = \{ \mathbf{x} \in T_{[X]\delta}^{n} : |T_{[\hat{X}|X]\delta}^{n}(\mathbf{x})| \ge 1 \}$$
(8.160)

<sup>&</sup>lt;sup>4</sup> Without conditioning on  $\{\mathbf{X} = \mathbf{x}\}$ , the events  $E_i$  are not mutually independent because they depend on each other through  $\mathbf{X}$ .

(cf. Section 6.3). Consider

$$\Pr\{E_1 | \mathbf{X} = \mathbf{x}\} = \Pr\left\{ (\mathbf{x}, \hat{\mathbf{X}}(1)) \in T^n_{[X\hat{X}]\delta} \right\}$$
(8.161)

$$=\sum_{\hat{\mathbf{x}}:(\mathbf{x},\hat{\mathbf{x}})\in T^n_{[X\hat{X}]\delta}} p(\hat{\mathbf{x}}).$$
(8.162)

The summation above is over all  $\hat{\mathbf{x}}$  such that  $(\mathbf{x}, \hat{\mathbf{x}}) \in T^n_{[X\hat{X}]\delta}$ . From the consistency of strong typicality (Theorem 6.7), if  $(\mathbf{x}, \hat{\mathbf{x}}) \in T^n_{[X\hat{X}]\delta}$ , then  $\hat{\mathbf{x}} \in T^n_{[\hat{X}]\delta}$ . By the strong AEP (Theorem 6.2), all  $p(\hat{\mathbf{x}})$  in the above summation satisfy

$$p(\hat{\mathbf{x}}) \ge 2^{-n(H(\hat{X})+\eta)},$$
 (8.163)

where  $\eta \to 0$  as  $\delta \to 0$ . By Theorem 6.10,

$$|T_{[\hat{X}|X]\delta}^{n}(\mathbf{x})| \ge 2^{n(H(\hat{X}|X)-\xi)},$$
(8.164)

where  $\xi \to 0$  as  $\delta \to 0$ . Then from (8.162), we have

$$\Pr\{E_1 | \mathbf{X} = \mathbf{x}\} \ge 2^{n(H(\hat{X}|X) - \xi)} 2^{-n(H(\hat{X}) + \eta)}$$
(8.165)

$$=2^{-n(H(\hat{X})-H(\hat{X}|X)+\xi+\eta)}$$
(8.166)

$$= 2^{-n(I(X;\hat{X})+\zeta)},$$
(8.167)

where

$$\zeta = \xi + \eta \to 0 \tag{8.168}$$

as  $\delta \to 0$ . Following (8.159), we have

$$\Pr\{K = 1 | \mathbf{X} = \mathbf{x}\} \le \left[1 - 2^{-n(I(X;\hat{X}) + \zeta)}\right]^{M-1}.$$
(8.169)

The lower bound in (8.153) implies

$$M \ge 2^{n(I(X;\hat{X}) + \frac{\epsilon}{2})}.$$
(8.170)

Then upon taking natural logarithm in (8.169), we obtain

$$\ln \Pr\{K = 1 | \mathbf{X} = \mathbf{x}\} \le (M - 1) \ln \left[ 1 - 2^{-n(I(X; \hat{X}) + \zeta)} \right]$$
(8.171)

$$\stackrel{a)}{\leq} \left( 2^{n(I(X;\hat{X}) + \frac{\epsilon}{2})} - 1 \right) \ln \left[ 1 - 2^{-n(I(X;\hat{X}) + \zeta)} \right]$$
(8.172)

<sup>b)</sup> 
$$\leq -\left(2^{n(I(X;\hat{X})+\frac{\epsilon}{2})}-1\right)2^{-n(I(X;\hat{X})+\zeta)}$$
 (8.173)

$$= -\left[2^{n(\frac{\epsilon}{2}-\zeta)} - 2^{-n(I(X;\hat{X})+\zeta)}\right].$$
(8.174)

#### 8.5 Achievability of $R_I(D)$ 201

In the above, a) follows from (8.170) by noting that the logarithm in (8.171) is negative, and b) follows from the fundamental inequality  $\ln a \leq a - 1$ . By letting  $\delta$  be sufficiently small so that

$$\frac{\epsilon}{2} - \zeta > 0, \tag{8.175}$$

the above upper bound on  $\ln \Pr\{K = 1 | \mathbf{X} = \mathbf{x}\}$  tends to  $-\infty$  as  $n \to \infty$ , i.e.,  $\Pr\{K = 1 | \mathbf{X} = \mathbf{x}\} \to 0$  as  $n \to \infty$ . This implies

$$\Pr\{K=1|\mathbf{X}=\mathbf{x}\} \le \frac{\epsilon}{2} \tag{8.176}$$

for sufficiently large n. It then follows that

$$Pr\{K = 1\}$$

$$= \sum_{\mathbf{x} \in S_{[X]\delta}^{n}} Pr\{K = 1 | \mathbf{X} = \mathbf{x}\} Pr\{\mathbf{X} = \mathbf{x}\}$$

$$+ \sum_{\mathbf{x} \notin S_{[X]\delta}^{n}} Pr\{K = 1 | \mathbf{X} = \mathbf{x}\} Pr\{\mathbf{X} = \mathbf{x}\}$$

$$(8.177)$$

$$\leq \sum_{\mathbf{x}\in S_{[X]\delta}^n} \frac{\epsilon}{2} \cdot \Pr\{\mathbf{X} = \mathbf{x}\} + \sum_{\mathbf{x}\notin S_{[X]\delta}^n} 1 \cdot \Pr\{\mathbf{X} = \mathbf{x}\}$$
(8.179)

$$= \frac{\epsilon}{2} \cdot \Pr\{\mathbf{X} \in S^n_{[X]\delta}\} + \Pr\{\mathbf{X} \notin S^n_{[X]\delta}\}$$
(8.180)

$$\leq \frac{\epsilon}{2} \cdot 1 + (1 - \Pr\{\mathbf{X} \in S^n_{[X]\delta}\}) \tag{8.181}$$

$$<\frac{\epsilon}{2}+\delta,\tag{8.182}$$

where we have invoked Proposition 6.13 in the last step. By letting  $\delta$  be sufficiently small so that

$$\delta < \frac{\epsilon}{2} \tag{8.183}$$

and (8.175) is satisfied, we obtain

$$\Pr\{K=1\} < \epsilon. \tag{8.184}$$

The main idea of the above upper bound on  $\Pr\{K = 1\}$  for sufficiently large *n* is the following. In constructing the codebook, we randomly generate M codewords in  $\hat{\mathcal{X}}^n$  according to  $p(\hat{x})^n$ . If M grows with *n* at a rate higher than  $I(X; \hat{X})$ , then the probability that there exists at least one codeword which is jointly typical with the source sequence **X** with respect to  $p(x, \hat{x})$ is very high when *n* is large. Further, the average distortion between **X** and such a codeword is close to  $Ed(X, \hat{X})$  because the empirical joint distribution of the symbol pairs in **X** and such a codeword is close to  $p(x, \hat{x})$ . Then by letting the reproduction sequence  $\hat{\mathbf{X}}$  be such a codeword, the average distortion

between **X** and  $\hat{\mathbf{X}}$  is less than  $D + \epsilon$  with probability arbitrarily close to 1 since  $Ed(X, \hat{X}) \leq D$ . These will be formally shown in the rest of the proof.

Now for sufficiently large n, consider

$$\Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon\}$$

$$= \Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon | K = 1\} \Pr\{K = 1\}$$

$$+ \Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon | K \neq 1\} \Pr\{K \neq 1\}$$

$$\leq 1 \cdot \epsilon + \Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon | K \neq 1\} \cdot 1$$
(8.186)
$$\hat{\mathbf{X}} = \{\mathbf{X}, \hat{\mathbf{X}} = 1\}$$

$$(8.186)$$

$$= \epsilon + \Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon | K \neq 1\}.$$
(8.187)

We will show that by choosing the value of  $\delta$  carefully, it is possible to make  $d(\mathbf{X}, \hat{\mathbf{X}})$  always less than or equal to  $D + \epsilon$  provided  $K \neq 1$ . Since  $(\mathbf{X}, \hat{\mathbf{X}}) \in$  $T^n_{[X\hat{X}]\delta}$  conditioning on  $\{K \neq 1\}$ , we have

$$d(\mathbf{X}, \dot{\mathbf{X}}) = \frac{1}{n} \sum_{k=1}^{n} d(X_k, \hat{X}_k)$$
(8.188)

$$= \frac{1}{n} \sum_{x,\hat{x}} d(x,\hat{x}) N(x,\hat{x}|\mathbf{X},\hat{\mathbf{X}})$$
(8.189)

$$= \frac{1}{n} \sum_{x,\hat{x}} d(x,\hat{x}) (np(x,\hat{x}) + N(x,\hat{x}|\mathbf{X},\hat{\mathbf{X}}) - np(x,\hat{x}))$$
(8.190)

$$= \left[\sum_{x,\hat{x}} p(x,\hat{x})d(x,\hat{x})\right] + \left[\sum_{x,\hat{x}} d(x,\hat{x})\left(\frac{1}{n}N(x,\hat{x}|\mathbf{X},\hat{\mathbf{X}}) - p(x,\hat{x})\right)\right]$$
(8.191)

$$= Ed(X,\hat{X}) + \sum_{x,\hat{x}} d(x,\hat{x}) \left(\frac{1}{n}N(x,\hat{x}|\mathbf{X},\hat{\mathbf{X}}) - p(x,\hat{x})\right)$$
(8.192)

$$\leq Ed(X,\hat{X}) + \sum_{x,\hat{x}} d(x,\hat{x}) \left| \frac{1}{n} N(x,\hat{x}|\mathbf{X},\hat{\mathbf{X}}) - p(x,\hat{x}) \right|$$

$$(8.193)$$

$$\stackrel{a)}{\leq} Ed(X,\hat{X}) + d_{max} \sum_{x,\hat{x}} \left| \frac{1}{n} N(x,\hat{x} | \mathbf{X}, \hat{\mathbf{X}}) - p(x,\hat{x}) \right|$$
(8.194)

$$\stackrel{b)}{\leq} Ed(X, \hat{X}) + d_{max}\delta \tag{8.195}$$

$$\stackrel{(5)}{\leq} D + d_{max}\delta,\tag{8.196}$$

where

- a) follows from the definition of  $d_{max}$  in (8.142);
- b) follows because  $(\mathbf{X}, \hat{\mathbf{X}}) \in T^n_{[X\hat{X}]\delta}$ ;

8.5 Achievability of  $R_I(D)$  203

c) follows from (8.152).

Therefore, by taking

$$\delta \le \frac{\epsilon}{d_{max}},\tag{8.197}$$

we obtain

$$d(\mathbf{X}, \hat{\mathbf{X}}) \le D + d_{max}\left(\frac{\epsilon}{d_{max}}\right) = D + \epsilon$$
 (8.198)

if  $K \neq 1$ . Therefore,

$$\Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon | K \neq 1\} = 0,$$
(8.199)

and it follows that from (8.187) that

$$\Pr\{d(\mathbf{X}, \hat{\mathbf{X}}) > D + \epsilon\} \le \epsilon.$$
(8.200)

Thus we have shown that for sufficiently large n, there exists an (n, M) random code which satisfies

$$\frac{1}{n}\log M \le I(X;\hat{X}) + \epsilon \tag{8.201}$$

(this follows from the upper bound in (8.153)) and (8.200). This implies the existence of an (n, M) rate-distortion code which satisfies (8.201) and (8.200). Therefore, the rate-distortion pair  $(I(X; \hat{X}), D)$  is achievable. Then upon minimizing over all  $\hat{X}$  which satisfy (8.152), we conclude that the rate-distortion pair  $(R_I(D), D)$  is achievable, which implies  $R_I(D) \geq R(D)$ . The proof is completed.

## Problems

- 1. Obtain the forward channel description of R(D) for the binary source with the Hamming distortion measure.
- 2. Binary covering radius The Hamming ball with center  $\mathbf{c} = (c_1, c_2, \dots, c_n) \in \{0, 1\}^n$  and radius r is the set

$$S_r(\mathbf{c}) = \left\{ \mathbf{x} \in \{0, 1\}^n : \sum_{i=1}^n |x_i - c_i| \le r \right\}.$$

Let  $M_{r,n}$  be the minimum number M such that there exists Hamming balls  $S_r(\mathbf{c}_j)$ ,  $j = 1, 2, \dots, M$  such that for all  $\mathbf{x} \in \{0, 1\}^n$ ,  $\mathbf{x} \in S_r(\mathbf{c}_j)$  for some j.

a) Show that

$$M_{r,n} \ge \frac{2^n}{\sum_{k=0}^r \binom{n}{k}}.$$

204 8 Rate-Distortion Theory

- b) What is the relation between  $M_{r,n}$  and the rate-distortion function for the binary source with the Hamming distortion measure?
- 3. Consider a source random variable X with the Hamming distortion measure.
  - a) Prove that

$$R(D) \ge H(X) - D\log(|\mathcal{X}| - 1) - h_b(D)$$

for  $0 \leq D \leq D_{max}$ .

b) Show that the above lower bound on R(D) is tight if X distributes uniformly on  $\mathcal{X}$ .

See Jerohin [171] (also see [75], p.133) for the tightness of this lower bound for a general source. This bound is a special case of the Shannon lower bound for the rate-distortion function [296] (also see [72], p.369).

4. Product source Let X and Y be two independent source random variables with reproduction alphabets  $\hat{\mathcal{X}}$  and  $\hat{\mathcal{Y}}$  and distortion measures  $d_x$  and  $d_y$ , and the rate-distortion functions for X and Y are denoted by  $R_x(D_x)$ and  $R_y(D_y)$ , respectively. Now for the product source (X, Y), define a distortion measure  $d: \mathcal{X} \times \mathcal{Y} \to \hat{\mathcal{X}} \times \hat{\mathcal{Y}}$  by

$$d((x, y), (\hat{x}, \hat{y})) = d_x(x, \hat{x}) + d_y(y, \hat{y}).$$

Prove that the rate-distortion function R(D) for (X, Y) with distortion measure d is given by

$$R(D) = \min_{D_x + D_y = D} (R_x(D_x) + R_y(D_y)).$$

Hint: Prove that  $I(X, Y; \hat{X}, \hat{Y}) \ge I(X; \hat{X}) + I(Y; \hat{Y})$  if X and Y are independent. (Shannon [296].)

5. Compound source Let  $\Theta$  be an index set and  $\mathcal{Z}_{\Theta} = \{X_{\theta} : \theta \in \Theta\}$  be a collection of source random variables. The random variables in  $\mathcal{Z}_{\Theta}$  have a common source alphabet  $\mathcal{X}$ , a common reproduction alphabet  $\hat{\mathcal{X}}$ , and a common distortion measure d. A compound source is an i.i.d. information source whose generic random variable is  $X_{\Phi}$ , where  $\Phi$  is equal to some  $\theta \in \Theta$  but we do not know which one it is. The rate-distortion function  $R_{\Phi}(D)$  for  $X_{\Phi}$  has the same definition as the rate-distortion function defined in this chapter except that (8.23) is replaced by

$$\Pr\{d(\mathbf{X}_{\theta}, \mathbf{\hat{X}}) > D + \epsilon\} \le \epsilon \quad \text{for all } \theta \in \Theta.$$

Show that

$$R_{\varPhi}(D) = \sup_{\theta \in \Theta} R_{\theta}(D),$$

where  $R_{\theta}(D)$  is the rate-distortion function for  $X_{\theta}$ .

6. Show that asymptotic optimality can always be achieved by separating rate-distortion coding and channel coding when the information source is i.i.d. (with a single-letter distortion measure) and the channel is memory-less.

- 7. Slepian-Wolf coding Let  $\epsilon, \gamma$ , and  $\delta$  be small positive quantities. For  $1 \leq i \leq 2^{n(H(Y|X)+\epsilon)}$ , randomly and independently select with replacement  $2^{n(I(X;Y)-\gamma)}$  sequences from  $T^n_{[Y]\delta}$  according to the uniform distribution to form a bin  $B_i$ . Let  $(\mathbf{x}, \mathbf{y})$  be a fixed pair of sequences in  $T^n_{[XY]\delta}$ . Prove the following by choosing  $\epsilon, \gamma$ , and  $\delta$  appropriately:
  - a) the probability that **y** is in some  $B_i$  tends to 1 as  $n \to \infty$ ;
  - b) given that  $\mathbf{y} \in B_i$ , the probability that there exists another  $\mathbf{y}' \in B_i$  such that  $(\mathbf{x}, \mathbf{y}') \in T^n_{[XY]\delta}$  tends to 0 as  $n \to \infty$ .

Let  $(\mathbf{X}, \mathbf{Y}) \sim p^n(x, y)$ . The results in a) and b) say that if  $(\mathbf{X}, \mathbf{Y})$  is jointly typical, which happens with probability close to 1 for large n, then it is very likely that  $\mathbf{Y}$  is in some bin  $B_i$ , and that  $\mathbf{Y}$  is the unique vector in  $B_i$  which is jointly typical with  $\mathbf{X}$ . If  $\mathbf{X}$  is available as side-information, then by specifying the index of the bin containing  $\mathbf{Y}$ , which takes about  $2^{nH(Y|X)}$  bits,  $\mathbf{Y}$  can be uniquely specified. Note that no knowledge about  $\mathbf{X}$  is involved in specifying the index of the bin containing  $\mathbf{Y}$ . This is the basis of the Slepian-Wolf coding [306] which launched the whole area of multiterminal source coding (see Berger [27]).

## **Historical Notes**

Transmission of an information source with distortion was first conceived by Shannon in his 1948 paper [291]. He returned to the problem in 1959 and proved the rate-distortion theorem [296]. The normalization of the ratedistortion function is due to Pinkston [261]. The rate-distortion theorem proved here is a stronger version of the original theorem. Extensions of the theorem to more general sources were proved in the book by Berger [26]. An iterative algorithm for computing the rate-distortion function developed by Blahut [35] will be discussed in Chapter 9. Rose [283] has developed an algorithm for the same purpose based on a mapping approach.

For a discrete memoryless channel p(y|x), the capacity

$$C = \max_{r(x)} I(X;Y), \tag{9.1}$$

where X and Y are respectively the input and the output of the generic channel and r(x) is the input distribution, characterizes the maximum asymptotically achievable rate at which information can be transmitted through the channel reliably. The expression for C in (9.1) is called a *single-letter characterization* because it depends only the transition matrix of the generic channel but not on the block length n of a code for the channel. When both the input alphabet  $\mathcal{X}$  and the output alphabet  $\mathcal{Y}$  are finite, the computation of C becomes a finite-dimensional maximization problem.

For an i.i.d. information source  $\{X_k, k \ge 1\}$  with generic random variable X, the rate-distortion function

$$R(D) = \min_{\substack{Q(\hat{x}|x): Ed(X, \hat{X}) \le D}} I(X; \hat{X})$$
(9.2)

characterizes the minimum asymptotically achievable rate of a rate-distortion code which reproduces the information source with an average distortion no more than D with respect to a single-letter distortion measure d. Again, the expression for R(D) in (9.2) is a single-letter characterization because it depends only on the generic random variable X but not on the block length n of a rate-distortion code. When both the source alphabet  $\mathcal{X}$  and the reproduction alphabet  $\hat{\mathcal{X}}$  are finite, the computation of R(D) becomes a finite-dimensional minimization problem.

Unless for very special cases, it is not possible to obtain an expression for C or R(D) in closed form, and we have to resort to numerical computation. However, computing these quantities is not straightforward because the associated optimization problem is nonlinear. In this chapter, we discuss the *Blahut-Arimoto algorithms* (henceforth the BA algorithms), which is an iterative algorithm devised for this purpose.

In order to better understand how and why the BA algorithm works, we will first describe the algorithm in a general setting in the next section. Specializations of the algorithm for the computation of C and R(D) will be discussed in Section 9.2, and convergence of the algorithm will be proved in Section 9.3.

## 9.1 Alternating Optimization

In this section, we describe an alternating optimization algorithm. This algorithm will be specialized in the next section for computing the channel capacity and the rate-distortion function.

Consider the double supremum

$$\sup_{\mathbf{u}_1 \in A_1} \sup_{\mathbf{u}_2 \in A_2} f(\mathbf{u}_1, \mathbf{u}_2), \tag{9.3}$$

where  $A_i$  is a convex subset of  $\Re^{n_i}$  for i = 1, 2, and f is a function defined on  $A_1 \times A_2$ . The function f is bounded from above, and is continuous and has continuous partial derivatives on  $A_1 \times A_2$ . Further assume that for all  $\mathbf{u}_2 \in A_2$ , there exists a unique  $c_1(\mathbf{u}_2) \in A_1$  such that

$$f(c_1(\mathbf{u}_2), \mathbf{u}_2) = \max_{\mathbf{u}_1' \in A_1} f(\mathbf{u}_1', \mathbf{u}_2),$$
(9.4)

and for all  $\mathbf{u}_1 \in A_1$ , there exists a unique  $c_2(\mathbf{u}_1) \in A_2$  such that

$$f(\mathbf{u}_1, c_2(\mathbf{u}_1)) = \max_{\mathbf{u}_2' \in A_2} f(\mathbf{u}_1, \mathbf{u}_2').$$
(9.5)

Let  $\mathbf{u} = (\mathbf{u}_1, \mathbf{u}_2)$  and  $A = A_1 \times A_2$ . Then (9.3) can be written as

$$\sup_{\mathbf{u}\in A} f(\mathbf{u}). \tag{9.6}$$

In other words, the supremum of f is taken over a subset of  $\Re^{n_1+n_2}$  which is equal to the Cartesian product of two convex subsets of  $\Re^{n_1}$  and  $\Re^{n_2}$ , respectively.

We now describe an alternating optimization algorithm for computing  $f^*$ , the value of the double supremum in (9.3). Let  $\mathbf{u}^{(k)} = (\mathbf{u}_1^{(k)}, \mathbf{u}_2^{(k)})$  for  $k \ge 0$ which are defined as follows. Let  $\mathbf{u}_1^{(0)}$  be an arbitrarily chosen vector in  $A_1$ , and let  $\mathbf{u}_2^{(0)} = c_2(\mathbf{u}_1^{(0)})$ . For  $k \ge 1$ ,  $\mathbf{u}^{(k)}$  is defined by

$$\mathbf{u}_1^{(k)} = c_1(\mathbf{u}_2^{(k-1)}) \tag{9.7}$$

and

$$\mathbf{u}_2^{(k)} = c_2(\mathbf{u}_1^{(k)}). \tag{9.8}$$

In other words,  $\mathbf{u}_1^{(k)}$  and  $\mathbf{u}_2^{(k)}$  are generated in the order  $\mathbf{u}_1^{(0)}$ ,  $\mathbf{u}_2^{(0)}$ ,  $\mathbf{u}_1^{(1)}$ ,  $\mathbf{u}_2^{(1)}$ ,  $\mathbf{u}_1^{(2)}$ ,  $\mathbf{u}_1^{(2)}$ ,  $\mathbf{u}_2^{(2)}$ ,  $\cdots$ , where each vector in the sequence is a function of the previous vector except that  $\mathbf{u}_1^{(0)}$  is arbitrarily chosen in  $A_1$ . Let

$$f^{(k)} = f(\mathbf{u}^{(k)}). \tag{9.9}$$

Then from (9.4) and (9.5),

$$f^{(k)} = f(\mathbf{u}_1^{(k)}, \mathbf{u}_2^{(k)}) \tag{9.10}$$

$$\geq f(\mathbf{u}_1^{(k)}, \mathbf{u}_2^{(k-1)}) \tag{9.11}$$

$$\geq f(\mathbf{u}_1^{(k-1)}, \mathbf{u}_2^{(k-1)}) \tag{9.12}$$

$$= f^{(k-1)} (9.13)$$

for  $k \geq 1$ . Since the sequence  $f^{(k)}$  is non-decreasing, it must converge because f is bounded from above. We will show in Section 9.3 that  $f^{(k)} \to f^*$  if f is concave. Figure 9.1 is an illustration of the alternating maximization algorithm, where in this case both  $n_1$  and  $n_2$  are equal to 1, and  $f^{(k)} \to f^*$ .

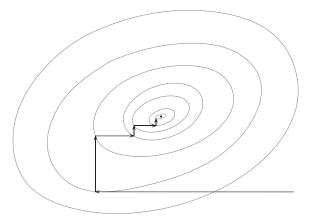


Fig. 9.1. Alternating optimization.

The alternating optimization algorithm can be explained by the following analogy. Suppose a hiker wants to reach the summit of a mountain. Starting from a certain point in the mountain, the hiker moves north-south and eastwest alternately. (In our problem, the north-south and east-west directions can be multi-dimensional.) In each move, the hiker moves to the highest possible point. The question is whether the hiker can eventually approach the summit starting from any point in the mountain.

Replacing f by -f in (9.3), the double supremum becomes the double infimum

$$\inf_{\mathbf{u}_1 \in A_1} \inf_{\mathbf{u}_2 \in A_2} f(\mathbf{u}_1, \mathbf{u}_2). \tag{9.14}$$

All the previous assumptions on  $A_1$ ,  $A_2$ , and f remain valid except that f is now assumed to be bounded from below instead of bounded from above. The double infimum in (9.14) can be computed by the same alternating optimization algorithm. Note that with f replaced by -f, the maximums in (9.4) and (9.5) become minimums, and the inequalities in (9.11) and (9.12) are reversed.

### 9.2 The Algorithms

In this section, we specialize the alternating optimization algorithm described in the last section to compute the channel capacity and the rate-distortion function. The corresponding algorithms are known as the BA algorithms.

#### 9.2.1 Channel Capacity

We will use **r** to denote an input distribution r(x), and we write  $\mathbf{r} > 0$  if **r** is strictly positive, i.e., r(x) > 0 for all  $x \in \mathcal{X}$ . If **r** is not strictly positive, we write  $\mathbf{r} \ge 0$ . Similar notations will be introduced as appropriate.

**Lemma 9.1.** Let r(x)p(y|x) be a given joint distribution on  $\mathcal{X} \times \mathcal{Y}$  such that  $\mathbf{r} > 0$ , and let  $\mathbf{q}$  be a transition matrix from  $\mathcal{Y}$  to  $\mathcal{X}$ . Then

$$\max_{\mathbf{q}} \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)} = \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q^*(x|y)}{r(x)}, \quad (9.15)$$

where the maximization is taken over all q such that

$$q(x|y) = 0 \quad if and only if \quad p(y|x) = 0, \tag{9.16}$$

and

$$q^*(x|y) = \frac{r(x)p(y|x)}{\sum_{x'} r(x')p(y|x')},$$
(9.17)

*i.e.*, the maximizing **q** is the which corresponds to the input distribution **r** and the transition matrix p(y|x).

In (9.15) and the sequel, we adopt the convention that the summation is taken over all x and y such that r(x) > 0 and p(y|x) > 0. Note that the right hand side of (9.15) gives the mutual information I(X;Y) when **r** is the input distribution for the generic channel p(y|x).

Proof. Let

$$w(y) = \sum_{x'} r(x')p(y|x')$$
(9.18)

#### 9.2 The Algorithms 211

in (9.17). We assume with loss of generality that for all  $y \in \mathcal{Y}$ , p(y|x) > 0 for some  $x \in \mathcal{X}$ . Since  $\mathbf{r} > 0$ , w(y) > 0 for all y, and hence  $q^*(x|y)$  is well-defined. Rearranging (9.17), we have

$$r(x)p(y|x) = w(y)q^*(x|y).$$
(9.19)

Consider

$$\sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q^*(x|y)}{r(x)} - \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}$$
$$= \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q^*(x|y)}{q(x|y)}$$
(9.20)

$$= \sum_{y} \sum_{x} w(y) q^{*}(x|y) \log \frac{q^{*}(x|y)}{q(x|y)}$$
(9.21)

$$= \sum_{y} w(y) \sum_{x} q^{*}(x|y) \log \frac{q^{*}(x|y)}{q(x|y)}$$
(9.22)

$$=\sum_{y} w(y) D(q^*(x|y) || q(x|y))$$
(9.23)

$$\geq 0, \tag{9.24}$$

where (9.21) follows from (9.19), and the last step is an application of the divergence inequality. Then the proof is completed by noting in (9.17) that  $\mathbf{q}^*$  satisfies (9.16) because  $\mathbf{r} > 0$ .  $\Box$ 

**Theorem 9.2.** For a discrete memoryless channel p(y|x),

$$C = \sup_{\mathbf{r}>0} \max_{\mathbf{q}} \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}, \tag{9.25}$$

where the maximization is taken over all  $\mathbf{q}$  which satisfies (9.16).

*Proof.* Let  $I(\mathbf{r}, \mathbf{p})$  denote the mutual information I(X; Y) when  $\mathbf{r}$  is the input distribution for the generic channel p(y|x). Then we can write

$$C = \max_{\mathbf{r} \ge 0} I(\mathbf{r}, \mathbf{p}). \tag{9.26}$$

Let  $\mathbf{r}^*$  achieves C. If  $\mathbf{r}^* > 0$ , then

$$C = \max_{\mathbf{r} \ge 0} I(\mathbf{r}, \mathbf{p}) \tag{9.27}$$

$$= \max_{\mathbf{r}>0} I(\mathbf{r}, \mathbf{p}) \tag{9.28}$$

$$= \max_{\mathbf{r}>0} \max_{\mathbf{q}} \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}$$
(9.29)

$$= \sup_{\mathbf{r}>0} \max_{\mathbf{q}} \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}, \qquad (9.30)$$

where (9.29) follows from Lemma 9.1 (and the maximization is over all **q** which satisfies (9.16)).

Next, we consider the case when  $\mathbf{r}^* \ge 0$ . Since  $I(\mathbf{r}, \mathbf{p})$  is continuous in  $\mathbf{r}$ , for any  $\epsilon > 0$ , there exists  $\delta > 0$  such that if

$$\|\mathbf{r} - \mathbf{r}^*\| < \delta, \tag{9.31}$$

then

$$C - I(\mathbf{r}, \mathbf{p}) < \epsilon, \tag{9.32}$$

where  $\|\mathbf{r} - \mathbf{r}^*\|$  denotes the Euclidean distance between  $\mathbf{r}$  and  $\mathbf{r}^*$ . In particular, there exists  $\tilde{\mathbf{r}} > 0$  which satisfies (9.31) and (9.32). Then

$$C = \max_{\mathbf{r} \ge 0} I(\mathbf{r}, \mathbf{p}) \tag{9.33}$$

$$\geq \sup_{\mathbf{r}>0} I(\mathbf{r}, \mathbf{p}) \tag{9.34}$$

$$\geq I(\tilde{\mathbf{r}}, \mathbf{p}) \tag{9.35}$$

$$> C - \epsilon,$$
 (9.36)

where the last step follows because  $\tilde{\mathbf{r}}$  satisfies (9.32). Thus we have

$$C - \epsilon < \sup_{\mathbf{r} > 0} I(\mathbf{r}, \mathbf{p}) \le C.$$
(9.37)

Finally, by letting  $\epsilon \to 0$ , we conclude that

$$C = \sup_{\mathbf{r}>0} I(\mathbf{r}, \mathbf{p}) = \sup_{\mathbf{r}>0} \max_{\mathbf{q}} \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}.$$
(9.38)

This accomplishes the proof.  $\hfill\square$ 

Now for the double supremum in (9.3), let

$$f(\mathbf{r}, \mathbf{q}) = \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}, \qquad (9.39)$$

with  $\mathbf{r}$  and  $\mathbf{q}$  playing the roles of  $\mathbf{u}_1$  and  $\mathbf{u}_2$ , respectively. Let

$$A_1 = \{ (r(x), x \in \mathcal{X}) : r(x) > 0 \text{ and } \sum_x r(x) = 1 \}, \qquad (9.40)$$

and

$$A_{2} = \{ (q(x|y), (x, y) \in \mathcal{X} \times \mathcal{Y}) : q(x|y) > 0 \\ \text{if } p(x|y) > 0, q(x|y) = 0 \text{ if } p(y|x) = 0, \\ \text{and } \sum_{x} q(x|y) = 1 \text{ for all } y \in \mathcal{Y} \}.$$
(9.41)

Then  $A_1$  is a subset of  $\Re^{|\mathcal{X}|}$  and  $A_2$  is a subset of  $\Re^{|\mathcal{X}||\mathcal{Y}|}$ , and it is readily checked that both  $A_1$  and  $A_2$  are convex. For all  $\mathbf{r} \in A_1$  and  $\mathbf{q} \in A_2$ , by Lemma 9.1,

9.2 The Algorithms 213

$$f(\mathbf{r}, \mathbf{q}) = \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}$$
(9.42)

$$\leq \sum_{x} \sum_{y} r(x)p(y|x) \log \frac{q^*(x|y)}{r(x)}$$
(9.43)

$$=I(X;Y) \tag{9.44}$$

$$\leq H(X) \tag{9.45}$$

$$\leq \log |\mathcal{X}|. \tag{9.46}$$

Thus f is bounded from above. Since for all  $\mathbf{q} \in A_2$ , q(x|y) = 0 for all x and y such that p(x|y) = 0, these components of  $\mathbf{q}$  are degenerated. In fact, these components of  $\mathbf{q}$  do not appear in the definition of  $f(\mathbf{r}, \mathbf{q})$  in (9.39), which can be seen as follows. Recall the convention that the double summation in (9.39) is over all x and y such that r(x) > 0 and p(y|x) > 0. If q(x|y) = 0, then p(y|x) = 0, and hence the corresponding term is not included in the double summation. Therefore, it is readily seen that f is continuous and has continuous partial derivatives on A because all the probabilities involved in the double summation in (9.39) are strictly positive. Moreover, for any given  $\mathbf{r} \in A_1$ , by Lemma 9.1, there exists a unique  $\mathbf{q} \in A_2$  which maximizes f. It will be shown shortly that for any given  $\mathbf{q} \in A_2$ , there also exists a unique  $\mathbf{r} \in A_1$  which maximizes f.

The double supremum in (9.3) now becomes

$$\sup_{\mathbf{r}\in A_1} \sup_{\mathbf{q}\in A_2} \sum_x \sum_y r(x) p(y|x) \log \frac{q(x|y)}{r(x)},\tag{9.47}$$

which by Theorem 9.2 is equal to C, where the supremum over all  $\mathbf{q} \in A_2$  is in fact a maximum. We then apply the alternating optimization algorithm in the last section to compute C. First, we arbitrarily choose a *strictly positive* input distribution in  $A_1$  and let it be  $\mathbf{r}^{(0)}$ . Then we define  $\mathbf{q}^{(0)}$  and in general  $\mathbf{q}^{(k)}$  for  $k \geq 0$  by

$$q^{(k)}(x|y) = \frac{r^{(k)}(x)p(y|x)}{\sum_{x'} r^{(k)}(x')p(y|x')}$$
(9.48)

in view of Lemma 9.1. In order to define  $\mathbf{r}^{(1)}$  and in general  $\mathbf{r}^{(k)}$  for  $k \ge 1$ , we need to find the  $\mathbf{r} \in A_1$  which maximizes f for a given  $\mathbf{q} \in A_2$ , where the constraints on  $\mathbf{r}$  are

$$\sum_{x} r(x) = 1 \tag{9.49}$$

and

$$r(x) > 0 \quad \text{for all } x \in \mathcal{X}.$$
 (9.50)

We now use the method of Lagrange multipliers to find the best  $\mathbf{r}$  by ignoring temporarily the positivity constraints in (9.50). Let

$$J = \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)} - \lambda \sum_{x} r(x).$$
(9.51)

For convenience sake, we assume that the logarithm is the natural logarithm. Differentiating with respect to r(x) gives

$$\frac{\partial J}{\partial r(x)} = \sum_{y} p(y|x) \log q(x|y) - \log r(x) - 1 - \lambda.$$
(9.52)

Upon setting  $\frac{\partial J}{\partial r(x)} = 0$ , we have

1

$$\log r(x) = \sum_{y} p(y|x) \log q(x|y) - 1 - \lambda,$$
(9.53)

or

$$\mathbf{r}(x) = e^{-(\lambda+1)} \prod_{y} q(x|y)^{p(y|x)}.$$
(9.54)

By considering the normalization constraint in (9.49), we can eliminate  $\lambda$  and obtain

$$r(x) = \frac{\prod_{y} q(x|y)^{p(y|x)}}{\sum_{x'} \prod_{y} q(x'|y)^{p(y|x')}}.$$
(9.55)

The above product is over all y such that p(y|x) > 0, and q(x|y) > 0 for all such y. This implies that both the numerator and the denominator on the right hand side above are positive, and therefore r(x) > 0. In other words, the  $\mathbf{r}$  thus obtained happen to satisfy the positivity constraints in (9.50) although these constraints were ignored when we set up the Lagrange multipliers. We will show in Section 9.3.2 that f is concave. Then  $\mathbf{r}$  as given in (9.55), which is unique, indeed achieves the maximum of f for a given  $\mathbf{q} \in A_2$  because  $\mathbf{r}$  is in the interior of  $A_1$ . In view of (9.55), we define  $\mathbf{r}^{(k)}$  for  $k \geq 1$  by

$$r^{(k)}(x) = \frac{\prod_{y} q^{(k-1)}(x|y)^{p(y|x)}}{\sum_{x'} \prod_{y} q^{(k-1)}(x'|y)^{p(y|x')}}.$$
(9.56)

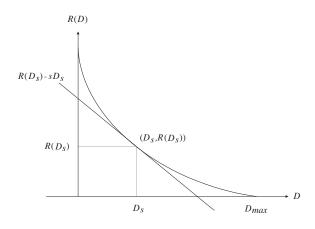
The vectors  $\mathbf{r}^{(k)}$  and  $\mathbf{q}^{(k)}$  are defined in the order  $\mathbf{r}^{(0)}$ ,  $\mathbf{q}^{(0)}$ ,  $\mathbf{r}^{(1)}$ ,  $\mathbf{q}^{(1)}$ ,  $\mathbf{r}^{(2)}$ ,  $\mathbf{q}^{(2)}$ ,..., where each vector in the sequence is a function of the previous vector except that  $\mathbf{r}^{(0)}$  is arbitrarily chosen in  $A_1$ . It remains to show by induction that  $\mathbf{r}^{(k)} \in A_1$  for  $k \ge 1$  and  $\mathbf{q}^{(k)} \in A_2$  for  $k \ge 0$ . If  $\mathbf{r}^{(k)} \in A_1$ , i.e.,  $\mathbf{r}^{(k)} > 0$ , then we see from (9.48) that  $q^{(k)}(x|y) = 0$  if and only if p(x|y) = 0, i.e.,  $\mathbf{q}^{(k)} \in A_2$ . On the other hand, if  $\mathbf{q}^{(k)} \in A_2$ , then we see from (9.56) that  $\mathbf{r}^{(k+1)} > 0$ , i.e.,  $\mathbf{r}^{(k+1)} \in A_2$ . Therefore,  $\mathbf{r}^{(k)} \in A_1$  and  $\mathbf{q}^{(k)} \in A_2$  for all  $k \ge 0$ . Upon determining  $(\mathbf{r}^{(k)}, \mathbf{q}^{(k)})$ , we can compute  $f^{(k)} = f(\mathbf{r}^{(k)}, \mathbf{q}^{(k)})$  for all k. It will be shown in Section 9.3 that  $f^{(k)} \to C$ .

#### 9.2.2 The Rate-Distortion Function

This discussion in this section is analogous to the discussion in Section 9.2.1. Some of the details will be omitted for brevity.

For all problems of interest, R(0) > 0. Otherwise, R(D) = 0 for all  $D \ge 0$  since R(D) is nonnegative and non-increasing. Therefore, we assume without loss of generality that R(0) > 0.

We have shown in Corollary 8.19 that if R(0) > 0, then R(D) is strictly decreasing for  $0 \le D \le D_{max}$ . Since R(D) is convex, for any  $s \le 0$ , there exists a point on the R(D) curve for  $0 \le D \le D_{max}$  such that the slope of a tangent<sup>1</sup> to the R(D) curve at that point is equal to s. Denote such a point on the R(D) curve by  $(D_s, R(D_s))$ , which is not necessarily unique. Then this tangent intersects with the ordinate at  $R(D_s) - sD_s$ . This is illustrated in Figure 9.2.



**Fig. 9.2.** A tangent to the R(D) curve with slope equal to s.

Let  $I(\mathbf{p}, \mathbf{Q})$  denote the mutual information  $I(X, \hat{X})$  and  $D(\mathbf{p}, \mathbf{Q})$  denote the expected distortion  $Ed(X, \hat{X})$  when  $\mathbf{p}$  is the distribution for X and  $\mathbf{Q}$  is the transition matrix from  $\mathcal{X}$  to  $\hat{\mathcal{X}}$  defining  $\hat{X}$ . Then for any  $\mathbf{Q}$ ,  $(I(\mathbf{p}, \mathbf{Q}), D(\mathbf{p}, \mathbf{Q}))$ is a point in the rate-distortion region, and the line with slope s passing through  $(I(\mathbf{p}, \mathbf{Q}), D(\mathbf{p}, \mathbf{Q}))$  intersects the ordinate at  $I(\mathbf{p}, \mathbf{Q}) - sD(\mathbf{p}, \mathbf{Q})$ . Since the R(D) curve defines the boundary of the rate-distortion region and it is above the tangent in Figure 9.2, we see that

$$R(D_s) - sD_s = \min_{\mathbf{Q}} [I(\mathbf{p}, \mathbf{Q}) - sD(\mathbf{p}, \mathbf{Q})].$$
(9.57)

<sup>&</sup>lt;sup>1</sup> We say that a line is a tangent to the R(D) curve if it touches the R(D) curve from below.

For each  $s \leq 0$ , if we can find a  $\mathbf{Q}_s$  which achieves the above minimum, then the line passing through  $(0, I(\mathbf{p}, \mathbf{Q}_s) - sD(\mathbf{p}, \mathbf{Q}_s))$ , i.e., the tangent in Figure 9.2, gives a tight lower bound on the R(D) curve. In particular, if  $(R(D_s), D_s)$  is unique,

$$D_s = D(\mathbf{p}, \mathbf{Q}_s) \tag{9.58}$$

and

$$R(D_s) = I(\mathbf{p}, \mathbf{Q}_s). \tag{9.59}$$

By varying over all  $s \leq 0$ , we can then trace out the whole R(D) curve. In the rest of the section, we will devise an iterative algorithm for the minimization problem in (9.57).

**Lemma 9.3.** Let  $p(x)Q(\hat{x}|x)$  be a given joint distribution on  $\mathcal{X} \times \hat{\mathcal{X}}$  such that  $\mathbf{Q} > 0$ , and let  $\mathbf{t}$  be any distribution on  $\hat{\mathcal{X}}$  such that  $\mathbf{t} > 0$ . Then

$$\min_{\mathbf{t}>0} \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t(\hat{x})} = \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t^*(\hat{x})}, \quad (9.60)$$

where

$$t^*(\hat{x}) = \sum_{x} p(x)Q(\hat{x}|x),$$
(9.61)

*i.e.*, the minimizing  $t(\hat{x})$  is the distribution on  $\hat{\mathcal{X}}$  corresponding to the input distribution  $\mathbf{p}$  and the transition matrix  $\mathbf{Q}$ .

*Proof.* It suffices to prove that

$$\sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t(\hat{x})} \ge \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t^*(\hat{x})}$$
(9.62)

for all  $\mathbf{t} > 0$ . The details are left as an exercise. Note in (9.61) that  $\mathbf{t}^* > 0$  because  $\mathbf{Q} > 0$ .  $\Box$ 

Since  $I(\mathbf{p}, \mathbf{Q})$  and  $D(\mathbf{p}, \mathbf{Q})$  are continuous in  $\mathbf{Q}$ , via an argument similar to the one we used in the proof of Theorem 9.2, we can replace the minimum over all  $\mathbf{Q}$  in (9.57) by the infimum over all  $\mathbf{Q} > 0$ . By noting that the right hand side of (9.60) is equal to  $I(\mathbf{p}, \mathbf{Q})$  and

$$D(\mathbf{p}, \mathbf{Q}) = \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x)d(x, \hat{x}), \qquad (9.63)$$

we can apply Lemma 9.3 to obtain

$$R(D_s) - sD_s = \inf_{Q>0} \left[ \min_{t>0} \sum_{x,\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t(\hat{x})} - s \sum_{x,\hat{x}} p(x)Q(\hat{x}|x)d(x,\hat{x}) \right]$$
(9.64)

$$= \inf_{Q>0} \min_{\mathbf{t}>0} \left[ \sum_{x,\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t(\hat{x})} - s \sum_{x,\hat{x}} p(x)Q(\hat{x}|x)d(x,\hat{x}) \right].$$
(9.65)

#### 9.2 The Algorithms 217

Now in the double infimum in (9.14), let

$$f(\mathbf{Q}, \mathbf{t}) = \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t(\hat{x})}$$
$$-s \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x)d(x, \hat{x}), \qquad (9.66)$$

$$A_{1} = \left\{ (Q(\hat{x}|x), (x, \hat{x}) \in \mathcal{X} \times \hat{\mathcal{X}}) : Q(\hat{x}|x) > 0, \\ \sum_{\hat{x}} Q(\hat{x}|x) = 1 \text{ for all } x \in \mathcal{X} \right\},$$

$$(9.67)$$

and

$$A_2 = \{ (t(\hat{x}), \hat{x} \in \hat{\mathcal{X}}) : t(\hat{x}) > 0 \text{ and } \sum_{\hat{x}} t(\hat{x}) = 1 \},$$
(9.68)

with **Q** and **t** playing the roles of  $\mathbf{u}_1$  and  $\mathbf{u}_2$ , respectively. Then  $A_1$  is a subset of  $\Re^{|\mathcal{X}||\hat{\mathcal{X}}|}$  and  $A_2$  is a subset of  $\Re^{|\hat{\mathcal{X}}|}$ , and it is readily checked that both  $A_1$  and  $A_2$  are convex. Since  $s \leq 0$ ,

$$f(\mathbf{Q}, \mathbf{t}) = \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t(\hat{x})} - s \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x)d(x, \hat{x})$$
(9.69)

$$\geq \sum_{x} \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t^{*}(\hat{x})} + 0$$
(9.70)

$$=I(X;\hat{X}) \tag{9.71}$$

$$\geq 0. \tag{9.72}$$

Therefore, f is bounded from below.

The double infimum in (9.14) now becomes

$$\inf_{\mathbf{Q}\in A_1} \inf_{\mathbf{t}\in A_2} \left[ \sum_x \sum_{\hat{x}} p(x)Q(\hat{x}|x) \log \frac{Q(\hat{x}|x)}{t(\hat{x})} - s \sum_x \sum_{\hat{x}} p(x)Q(\hat{x}|x)d(x,\hat{x}) \right],\tag{9.73}$$

where the infimum over all  $\mathbf{t} \in A_2$  is in fact a minimum. We then apply the alternating optimization algorithm described in Section 9.2 to compute  $f^*$ , the value of (9.73). First, we arbitrarily choose a *strictly positive* transition matrix in  $A_1$  and let it be  $\mathbf{Q}^{(0)}$ . Then we define  $\mathbf{t}^{(0)}$  and in general  $\mathbf{t}^{(k)}$  for  $k \geq 1$  by

$$t^{(k)}(\hat{x}) = \sum_{x} p(x)Q^{(k)}(\hat{x}|x)$$
(9.74)

in view of Lemma 9.3. In order to define  $\mathbf{Q}^{(1)}$  and in general  $\mathbf{Q}^{(k)}$  for  $k \ge 1$ , we need to find the  $\mathbf{Q} \in A_1$  which minimizes f for a given  $\mathbf{t} \in A_2$ , where the constraints on  $\mathbf{Q}$  are

$$Q(\hat{x}|x) > 0 \quad \text{for all } (x, \hat{x}) \in \mathcal{X} \times \hat{\mathcal{X}}, \tag{9.75}$$

and

$$\sum_{\hat{x}} Q(\hat{x}|x) = 1 \quad \text{for all } x \in \mathcal{X}.$$
(9.76)

As we did for the computation of the channel capacity, we first ignore the positivity constraints in (9.75) when setting up the Lagrange multipliers. Then we obtain

$$Q(\hat{x}|x) = \frac{t(\hat{x})e^{sd(x,\hat{x})}}{\sum_{\hat{x}'} t(\hat{x}')e^{sd(x,\hat{x}')}} > 0.$$
(9.77)

The details are left as an exercise. We then define  $\mathbf{Q}^{(k)}$  for  $k \geq 1$  by

$$Q^{(k)}(\hat{x}|x) = \frac{t^{(k-1)}(\hat{x})e^{sd(x,\hat{x})}}{\sum_{\hat{x}'} t^{(k-1)}(\hat{x}')e^{sd(x,\hat{x}')}}.$$
(9.78)

It will be shown in the next section that  $f^{(k)} = f(\mathbf{Q}^{(k)}, \mathbf{t}^{(k)}) \to f^*$  as  $k \to \infty$ . If there exists a unique point  $(R(D_s), D_s)$  on the R(D) curve such that the slope of a tangent at that point is equal to s, then

$$(I(\mathbf{p}, \mathbf{Q}^{(k)}), D(\mathbf{p}, \mathbf{Q}^{(k)})) \to (R(D_s), D_s).$$

$$(9.79)$$

Otherwise,  $(I(\mathbf{p}, \mathbf{Q}^{(k)}), D(\mathbf{p}, \mathbf{Q}^{(k)}))$  is arbitrarily close to the segment of the R(D) curve at which the slope is equal to s when k is sufficiently large. These facts are easily shown to be true.

#### 9.3 Convergence

In this section, we first prove that if f is concave, then  $f^{(k)} \to f^*$ . We then apply this sufficient condition to prove the convergence of the BA algorithm for computing the channel capacity. The convergence of the BA algorithm for computing the rate-distortion function can be proved likewise. The details are omitted.

#### 9.3.1 A Sufficient Condition

In the alternating optimization algorithm in Section 9.1, we see from (9.7) and (9.8) that

$$\mathbf{u}^{(k+1)} = (\mathbf{u}_1^{(k+1)}, \mathbf{u}_2^{(k+1)}) = (c_1(\mathbf{u}_2^{(k)}), c_2(c_1(\mathbf{u}_2^{(k)})))$$
(9.80)

for  $k \geq 0$ . Define

9.3 Convergence 219

$$\Delta f(\mathbf{u}) = f(c_1(\mathbf{u}_2), c_2(c_1(\mathbf{u}_2))) - f(\mathbf{u}_1, \mathbf{u}_2).$$
(9.81)

Then

$$f^{(k+1)} - f^{(k)} = f(\mathbf{u}^{(k+1)}) - f(\mathbf{u}^{(k)})$$
(9.82)

$$= f(c_1(\mathbf{u}_2^{(k)}), c_2(c_1(\mathbf{u}_2^{(k)}))) - f(\mathbf{u}_1^{(k)}, \mathbf{u}_2^{(k)})$$
(9.83)

$$= \Delta f(\mathbf{u}^{(k)}). \tag{9.84}$$

We will prove that f being concave is sufficient for  $f^{(k)} \to f^*$ . To this end, we first prove that if f is concave, then the algorithm cannot be trapped at  $\mathbf{u}$  if  $f(\mathbf{u}) < f^*$ .

**Lemma 9.4.** Let f be concave. If  $f^{(k)} < f^*$ , then  $f^{(k+1)} > f^{(k)}$ .

*Proof.* We will prove that  $\Delta f(\mathbf{u}) > 0$  for any  $\mathbf{u} \in A$  such that  $f(\mathbf{u}) < f^*$ . Then if  $f^{(k)} = f(\mathbf{u}^{(k)}) < f^*$ , we see from (9.84) that

$$f^{(k+1)} - f^{(k)} = \Delta f(\mathbf{u}^{(k)}) > 0, \qquad (9.85)$$

and the lemma is proved.

Consider any  $\mathbf{u} \in A$  such that  $f(\mathbf{u}) < f^*$ . We will prove by contradiction that  $\Delta f(\mathbf{u}) > 0$ . Assume  $\Delta f(\mathbf{u}) = 0$ . Then it follows from (9.81) that

$$f(c_1(\mathbf{u}_2), c_2(c_1(\mathbf{u}_2))) = f(\mathbf{u}_1, \mathbf{u}_2).$$
(9.86)

Now we see from (9.5) that

$$f(c_1(\mathbf{u}_2), c_2(c_1(\mathbf{u}_2))) \ge f(c_1(\mathbf{u}_2), \mathbf{u}_2).$$
 (9.87)

If  $c_1(\mathbf{u}_2) \neq \mathbf{u}_1$ , then

$$f(c_1(\mathbf{u}_2), \mathbf{u}_2) > f(\mathbf{u}_1, \mathbf{u}_2)$$
 (9.88)

because  $c_1(\mathbf{u}_2)$  is unique. Combining (9.87) and (9.88), we have

$$f(c_1(\mathbf{u}_2), c_2(c_1(\mathbf{u}_2))) > f(\mathbf{u}_1, \mathbf{u}_2),$$
 (9.89)

which is a contradiction to (9.86). Therefore,

$$\mathbf{u}_1 = c_1(\mathbf{u}_2). \tag{9.90}$$

Using this, we see from (9.86) that

$$f(\mathbf{u}_1, c_2(\mathbf{u}_1)) = f(\mathbf{u}_1, \mathbf{u}_2),$$
 (9.91)

which implies

$$\mathbf{u}_2 = c_2(\mathbf{u}_1). \tag{9.92}$$

because  $c_2(c_1(\mathbf{u}_2))$  is unique.

Since  $f(\mathbf{u}) < f^*$ , there exists  $\mathbf{v} \in A$  such that

$$f(\mathbf{u}) < f(\mathbf{v}). \tag{9.93}$$

Consider

$$\mathbf{v} - \mathbf{u} = (\mathbf{v}_1 - \mathbf{u}_1, 0) + (0, \mathbf{v}_2 - \mathbf{u}_2).$$
 (9.94)

Let  $\tilde{\mathbf{z}}$  be the unit vector in the direction of  $\mathbf{v} - \mathbf{u}$ ,  $\mathbf{z}_1$  be the unit vector in the direction of  $(\mathbf{v}_1 - \mathbf{u}_1, 0)$ , and  $\mathbf{z}_2$  be the unit vector in the direction of  $(\mathbf{v}_2 - \mathbf{u}_2, 0)$ . Then

$$\|\mathbf{v} - \mathbf{u}\|\tilde{\mathbf{z}} = \|\mathbf{v}_1 - \mathbf{u}_1\|\mathbf{z}_1 + \|\mathbf{v}_2 - \mathbf{u}_2\|\mathbf{z}_2,$$
 (9.95)

or

$$\tilde{\mathbf{z}} = \alpha_1 \mathbf{z}_1 + \alpha_2 \mathbf{z}_2, \tag{9.96}$$

where

$$\alpha_i = \frac{\|\mathbf{v}_i - \mathbf{u}_i\|}{\|\mathbf{v} - \mathbf{u}\|},\tag{9.97}$$

i = 1, 2. Figure 9.3 is an illustration of the vectors  $\mathbf{u}, \mathbf{v}, \tilde{\mathbf{z}}, \mathbf{z}_1$ , and  $\mathbf{z}_2$ .

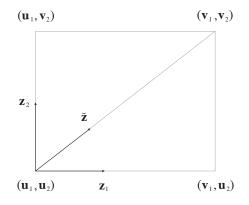


Fig. 9.3. The vectors  $\mathbf{u}$ ,  $\mathbf{v}$ ,  $\tilde{\mathbf{z}}$ ,  $\mathbf{z}_1$ , and  $\mathbf{z}_2$ .

We see from (9.90) that f attains its maximum value at  $\mathbf{u} = (\mathbf{u}_1, \mathbf{u}_2)$ when  $\mathbf{u}_2$  is fixed. In particular, f attains its maximum value at  $\mathbf{u}$  alone the line passing through  $(\mathbf{u}_1, \mathbf{u}_2)$  and  $(\mathbf{v}_1, \mathbf{u}_2)$ . Let  $\nabla f$  denotes the gradient of f. Since f is continuous and has continuous partial derivatives, the directional derivative of f at  $\mathbf{u}$  in the direction of  $\mathbf{z}_1$  exists and is given by  $\nabla f \cdot \mathbf{z}_1$ . It follows from the concavity of f that f is concave along the line passing through  $(\mathbf{u}_1, \mathbf{u}_2)$  and  $(\mathbf{v}_1, \mathbf{u}_2)$ . Since f attains its maximum value at  $\mathbf{u}$ , the derivative of f along the line passing through  $(\mathbf{u}_1, \mathbf{u}_2)$  and  $(\mathbf{v}_1, \mathbf{u}_2)$  vanishes. Then we see that

$$\nabla f \cdot \mathbf{z}_1 = 0. \tag{9.98}$$

Similarly, we see from (9.92) that

9.3 Convergence 221

$$\nabla f \cdot \mathbf{z}_2 = 0. \tag{9.99}$$

Then from (9.96), the directional derivative of f at **u** in the direction of  $\tilde{\mathbf{z}}$  is given by

$$\nabla f \cdot \tilde{\mathbf{z}} = \alpha_1 (\nabla f \cdot \mathbf{z}_1) + \alpha_2 (\nabla f \cdot \mathbf{z}_2) = 0.$$
(9.100)

Since f is concave along the line passing through  $\mathbf{u}$  and  $\mathbf{v}$ , this implies

$$f(\mathbf{u}) \ge f(\mathbf{v}),\tag{9.101}$$

which is a contradiction to (9.93). Hence, we conclude that  $\Delta f(\mathbf{u}) > 0$ .  $\Box$ 

Although we have proved that the algorithm cannot be trapped at **u** if  $f(\mathbf{u}) < f^*$ ,  $f^{(k)}$  does not necessarily converge to  $f^*$  because the increment in  $f^{(k)}$  in each step may be arbitrarily small. In order to prove the desired convergence, we will show in next theorem that this cannot be the case.

**Theorem 9.5.** If f is concave, then  $f^{(k)} \rightarrow f^*$ .

*Proof.* We have already shown in Section 9.1 that  $f^{(k)}$  necessarily converges, say to f'. Hence, for any  $\epsilon > 0$  and all sufficiently large k,

$$f' - \epsilon \le f^{(k)} \le f'. \tag{9.102}$$

Let

$$\gamma = \min_{\mathbf{u} \in A'} \Delta f(\mathbf{u}), \tag{9.103}$$

where

$$A' = \{ \mathbf{u} \in A : f' - \epsilon \le f(\mathbf{u}) \le f' \}.$$

$$(9.104)$$

Since f has continuous partial derivatives,  $\Delta f(\mathbf{u})$  is a continuous function of **u**. Then the minimum in (9.103) exists because A' is compact<sup>2</sup>.

We now show that  $f' < f^*$  will lead to a contradiction if f is concave. If  $f' < f^*$ , then from Lemma 9.4, we see that  $\Delta f(\mathbf{u}) > 0$  for all  $\mathbf{u} \in A'$  and hence  $\gamma > 0$ . Since  $f^{(k)} = f(\mathbf{u}^{(k)})$  satisfies (9.102),  $\mathbf{u}^{(k)} \in A'$ , and

$$f^{(k+1)} - f^{(k)} = \Delta f(\mathbf{u}^{(k)}) \ge \gamma$$
 (9.105)

for all sufficiently large k. Therefore, no matter how smaller  $\gamma$  is,  $f^{(k)}$  will eventually be greater than f', which is a contradiction to  $f^{(k)} \to f'$ . Hence, we conclude that  $f^{(k)} \to f^*$ .  $\Box$ 

 $<sup>^2</sup>$   $A^\prime$  is compact because it is the inverse image of a closed interval under a continuous function and A is bounded.

#### 9.3.2 Convergence to the Channel Capacity

In order to show that the BA algorithm for computing the channel capacity converges as intended, i.e.,  $f^{(k)} \to C$ , we only need to show that the function f defined in (9.39) is concave. Toward this end, for

$$f(\mathbf{r}, \mathbf{q}) = \sum_{x} \sum_{y} r(x) p(y|x) \log \frac{q(x|y)}{r(x)}$$
(9.106)

defined in (9.39), we consider two ordered pairs  $(\mathbf{r}_1, \mathbf{q}_1)$  and  $(\mathbf{r}_2, \mathbf{q}_2)$  in A, where  $A_1$  and  $A_2$  are defined in (9.40) and (9.41), respectively. For any  $0 \leq \lambda \leq 1$  and  $\bar{\lambda} = 1 - \lambda$ , an application of the log-sum inequality (Theorem 2.32) gives

$$(\lambda r_1(x) + \bar{\lambda} r_2(x)) \log \frac{\lambda r_1(x) + \bar{\lambda} r_2(x)}{\lambda q_1(x|y) + \bar{\lambda} q_2(x|y)} \\ \leq \lambda r_1(x) \log \frac{r_1(x)}{q_1(x|y)} + \bar{\lambda} r_2(x) \log \frac{r_2(x)}{q_2(x|y)}.$$
(9.107)

Taking reciprocal in the logarithms yields

$$(\lambda r_1(x) + \bar{\lambda} r_2(x)) \log \frac{\lambda q_1(x|y) + \lambda q_2(x|y)}{\lambda r_1(x) + \bar{\lambda} r_2(x)}$$
  

$$\geq \lambda r_1(x) \log \frac{q_1(x|y)}{r_1(x)} + \bar{\lambda} r_2(x) \log \frac{q_2(x|y)}{r_2(x)}, \qquad (9.108)$$

and upon multiplying by p(y|x) and summing over all x and y, we obtain

$$f(\lambda \mathbf{r}_1 + \bar{\lambda} \mathbf{r}_2, \lambda \mathbf{q}_1 + \bar{\lambda} \mathbf{q}_2) \ge \lambda f(\mathbf{r}_1, \mathbf{q}_1) + \bar{\lambda} f(\mathbf{r}_2, \mathbf{q}_2).$$
(9.109)

Therefore, f is concave. Hence, we have shown that  $f^{(k)} \to C$ .

## Problems

- 1. Implement the BA algorithm for computing channel capacity.
- 2. Implement the BA algorithm for computing the rate-distortion function.
- 3. Explain why in the BA Algorithm for computing channel capacity, we should not choose an initial input distribution which contains zero probability masses.
- 4. Prove Lemma 9.3.
- 5. Consider  $f(\mathbf{Q}, \mathbf{t})$  in the BA algorithm for computing the rate-distortion function.
  - a) Show that for fixed s and t,  $f(\mathbf{Q}, \mathbf{t})$  is minimized by

$$Q(\hat{x}|x) = \frac{t(\hat{x})e^{sd(x,\hat{x})}}{\sum_{\hat{x}'} t(\hat{x}')e^{sd(x,\hat{x}')}}$$

b) Show that  $f(\mathbf{Q}, \mathbf{t})$  is convex.

## **Historical Notes**

An iterative algorithm for computing the channel capacity was developed by Arimoto [18], where the convergence of the algorithm was proved. Blahut [35] independently developed two similar algorithms, the first for computing the channel capacity and the second for computing the rate-distortion function. The convergence of Blahut's second algorithm was proved by Csiszár [74]. These two algorithms are now commonly referred to as the Blahut-Arimoto algorithms. The simplified proof of convergence in this chapter is based on Yeung and Berger [362].

The Blahut-Arimoto algorithms are special cases of a general iterative algorithm due to Csiszár and Tusnády [79] which also include the EM algorithm [83] for fitting models from incomplete data and the algorithm for finding the log-optimal portfolio for a stock market due to Cover [68].

Our discussion in the previous chapters involved only discrete random variables. The actual values taken by these random variables did not play any role in establishing the results. In this chapter and the next chapter, our discussion will involve random variables taking real values. As we will see, the values taken by these random variables *do* play a crucial role in the discussion.

Let X be a real random variable with *cumulative distribution function* (CDF)  $F_X(x) = \Pr\{X \leq x\}$ , which by definition is right-continuous. The random variable X is said to be

- discrete if  $F_X(x)$  increases only at a countable number of values of x;
- continuous if  $F_X(x)$  is continuous, or equivalently,  $\Pr\{X = x\} = 0$  for every value of x;
- *mixed* if it is neither discrete nor continuous.

If  $F_X(x)$  is differentiable with derivative  $f_X(x)$ , then  $f_X(x)$  is called the *probability density function*<sup>1</sup> (pdf) of X. The support of X, denoted by  $S_X$ , is the set of all x such that  $F_X(x) > F_X(x-\epsilon)$  for all  $\epsilon > 0$ . For a function g defined on  $S_X$ , we write

$$Eg(X) = \int_{\mathcal{S}_X} g(x) dF_X(x), \qquad (10.1)$$

where the right hand side is a *Lebesgue-Stieltjes integration* which covers all the cases (i.e., discrete, continuous, and mixed) for the CDF  $F_X(x)$ . It may be regarded as a notation for the expectation of g(X) with respect to  $F_X(x)$ if the reader is not familiar with measure theory.

Let X and Y be two real random variables with joint CDF  $F_{XY}(x,y) = \Pr\{X \leq x, Y \leq y\}$ . The marginal CDF of X is given by  $F_X(x) = F_{XY}(x,\infty)$  (likewise for Y). If  $\frac{\partial^2}{\partial x \partial y} F_{XY}(x,y)$  exists, it is called the joint pdf of X and Y and is denoted by  $f_{XY}(x,y)$ . For  $x \in \mathcal{S}_X$ , the conditional CDF of Y given  $\{X = x\}$  is defined as

<sup>&</sup>lt;sup>1</sup> More generally, the pdf of X can still be defined if  $F_X(x)$  is differentiable at all but a countable number of values of x.

$$F_{Y|X}(y|x) = \int_{-\infty}^{y} f_{Y|X}(v|x)dv,$$
(10.2)

where

$$f_{Y|X}(y|x) = \frac{f_{XY}(x,y)}{f_X(x)}$$
(10.3)

is the conditional pdf of Y given  $\{X = x\}$ . All the above notations naturally extend to two or more real random variables. When there is no ambiguity, the subscripts specifying the random variables will be omitted.

All the random variables in this chapter are assumed to be real<sup>2</sup>. The variance of a random variable X is defined as

$$\operatorname{var} X = E(X - EX)^2 = EX^2 - (EX)^2.$$
(10.4)

The *covariance* between two random variables X and Y is defined as

$$cov(X, Y) = E(X - EX)(Y - EY) = E(XY) - (EX)(EY).$$
 (10.5)

For a random vector  $\mathbf{X} = [X_1 X_2 \cdots X_n]^{\top}$ , the covariance matrix is defined as

$$K_{\mathbf{X}} = E(\mathbf{X} - E\mathbf{X})(\mathbf{X} - E\mathbf{X})^{\top} = [\operatorname{cov}(X_i, X_j)], \qquad (10.6)$$

and the correlation matrix is defined as

$$\tilde{K}_{\mathbf{X}} = E\mathbf{X}\mathbf{X}^{\top} = [EX_i X_j]. \tag{10.7}$$

Then

$$K_{\mathbf{X}} = E(\mathbf{X} - E\mathbf{X})(\mathbf{X} - E\mathbf{X})^{\top}$$
(10.8)  
$$= E[\mathbf{X}\mathbf{X}^{\top} - \mathbf{X}(E\mathbf{X}^{\top}) - (E\mathbf{X})\mathbf{X}^{\top} + (E\mathbf{X})(E\mathbf{X}^{\top})]$$
(10.9)

$$= E[\mathbf{X}\mathbf{X}^{\top} - \mathbf{X}(E\mathbf{X}^{\top}) - (E\mathbf{X})\mathbf{X}^{\top} + (E\mathbf{X})(E\mathbf{X}^{\top})]$$
(10.9)  
$$= E[\mathbf{X}\mathbf{X}^{\top} - (E\mathbf{X})(E\mathbf{X}^{\top}) + (E\mathbf{X})(E\mathbf{X}^{\top})]$$
(10.10)

$$= E\mathbf{X}\mathbf{X}^{\top} - (E\mathbf{X})(E\mathbf{X}^{\top}) - (E\mathbf{X})(E\mathbf{X}^{\top}) + (E\mathbf{X})(E\mathbf{X}^{\top})$$
(10.10)  
$$= E\mathbf{X}\mathbf{X}^{\top} - (E\mathbf{X})(E\mathbf{X}^{\top})$$
(10.11)

$$= E\mathbf{X}\mathbf{X}^{\top} - (E\mathbf{X})(E\mathbf{X}^{\top}) \tag{10.11}$$

$$= K_{\mathbf{X}} - (E\mathbf{X})(E\mathbf{X})^{\top}.$$
(10.12)

This implies that if  $E\mathbf{X} = 0$ , then  $\tilde{K}_{\mathbf{X}} = K_{\mathbf{X}}$ . It can readily be verified that in general,

$$\tilde{K}_{\mathbf{X}} = K_{\mathbf{X}+E\mathbf{X}} \tag{10.13}$$

and

$$K_{\mathbf{X}} = \tilde{K}_{\mathbf{X}-E\mathbf{X}}.\tag{10.14}$$

Therefore, a correlation matrix is a covariance matrix, and vice versa. When there is no ambiguity, the subscripts in  $K_{\mathbf{X}}$  and  $\tilde{K}_{\mathbf{X}}$  will be omitted.

<sup>&</sup>lt;sup>2</sup> If a real random variable X is discrete, i.e., the support  $S_X$  is a countable subset of  $\Re$ , all the information measures involving X are unchanged if the alphabet  $\mathcal{X}$ is replaced by any countable index set.

Let  $\mathcal{N}(\mu, \sigma^2)$  denote the Gaussian distribution with mean  $\mu$  and variance  $\sigma^2$ , i.e., the pdf of the distribution is given by

$$f(x) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$
(10.15)

for  $-\infty < x < \infty$ . Similarly, let  $\mathcal{N}(\boldsymbol{\mu}, K)$  denote the multivariate Gaussian distribution with mean  $\boldsymbol{\mu}$  and covariance matrix K, i.e., the joint pdf is given by

$$f(\mathbf{x}) = \frac{1}{\left(\sqrt{2\pi}\right)^n |K|^{1/2}} e^{-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^\top K^{-1}(\mathbf{x} - \boldsymbol{\mu})}$$
(10.16)

for all  $\mathbf{x} \in \Re^n$ , where K is a symmetric positive definite matrix<sup>3</sup> and |K| is the determinant of K.

In the rest of the chapter, we will define various information measures under suitable conditions. Whenever these information measures are subsequently mentioned, they are assumed to be defined.

## **10.1** Preliminaries

In this section, we present some preliminary results on matrices and linear transformation of random variables. All the matrices and vectors are assumed to be real.

**Definition 10.1.** A square matrix K is symmetric if  $K^{\top} = K$ .

**Definition 10.2.** An  $n \times n$  matrix K is positive definite if

$$\mathbf{x}^{\top} K \mathbf{x} > 0 \tag{10.17}$$

for all nonzero column n-vector  $\mathbf{x}$ , and is positive semidefinite if

$$\mathbf{x}^{\top} K \mathbf{x} \ge 0 \tag{10.18}$$

for all column n-vector  $\mathbf{x}$ .

**Proposition 10.3.** A covariance (correlation) matrix is both symmetric and positive semidefinite.

*Proof.* Omitted.  $\Box$ 

If a matrix K is symmetric, it can be diagonalized as

$$K = Q \Lambda Q^{\top}, \tag{10.19}$$

where  $\Lambda$  is a diagonal matrix and Q (also  $Q^{\top}$ ) is an orthogonal matrix, i.e.,

 $<sup>^{3}</sup>$  See Definitions 10.1 and 10.2.

$$Q^{-1} = Q^{\top}, \tag{10.20}$$

or

$$QQ^{\top} = Q^{\top}Q = I. \tag{10.21}$$

The latter says that the rows (columns) of Q form an orthonormal system. Since

$$|Q|^{2} = |Q||Q^{\top}| = |QQ^{\top}| = |I| = 1,$$
(10.22)

we have

$$Q| = |Q^{\top}| = 1. \tag{10.23}$$

From (10.19) and (10.21), we have

$$KQ = (Q\Lambda Q^{\top})Q = Q\Lambda (Q^{\top}Q) = Q\Lambda.$$
(10.24)

Let  $\lambda_i$  and  $\mathbf{q}_i$  denote the *i*th diagonal element of  $\Lambda$  and the *i*th column of Q, respectively. Then (10.24) can be written as

$$K\mathbf{q}_i = \lambda_i \mathbf{q}_i \tag{10.25}$$

for all *i*, i.e.,  $\mathbf{q}_i$  is an eigenvector of *K* with eigenvalue  $\lambda_i$ . The next proposition further shows that these eigenvalues are nonnegative if *K* is positive semidefinite.

**Proposition 10.4.** The eigenvalues of a positive semidefinite matrix are nonnegative.

*Proof.* Let K be a positive semidefinite matrix, and let  $\mathbf{q}$  be an eigenvector of K with eigenvalue  $\lambda$ , i.e.,

$$K\mathbf{q} = \lambda \mathbf{q}.\tag{10.26}$$

Since K is positive semidefinite,

$$0 \le \mathbf{q}^{\top} K \mathbf{q} = \mathbf{q}^{\top} (\lambda \mathbf{q}) = \lambda(\mathbf{q}^{\top} \mathbf{q}).$$
(10.27)

Then we conclude that  $\lambda \geq 0$  because  $\mathbf{q}^{\top}\mathbf{q} \geq 0$ .  $\Box$ 

The above discussions on diagonalization apply to a covariance matrix because a covariance matrix is both symmetric and positive semidefinite. As we will see, by diagonalizing the covariance matrix, a set of correlated random variables can be decorrelated by an orthogonal transformation. On the other hand, a set of correlated random variables can be regarded as an orthogonal transformation of a set of uncorrelated random variables. This is particularly important in the context of Gaussian random variables because a set of jointly distributed Gaussian random variables are mutually independent if and only if they are uncorrelated. **Proposition 10.5.** Let  $\mathbf{Y} = A\mathbf{X}$ , where  $\mathbf{X}$  and  $\mathbf{Y}$  are column vectors of n random variables and A is an  $n \times n$  matrix. Then

$$K_{\mathbf{Y}} = AK_{\mathbf{X}}A^{\top} \tag{10.28}$$

and

$$\tilde{K}_{\mathbf{Y}} = A \tilde{K}_{\mathbf{X}} A^{\top}.$$
(10.29)

*Proof.* To prove (10.28), consider

$$K_{\mathbf{Y}} = E(\mathbf{Y} - E\mathbf{Y})(\mathbf{Y} - E\mathbf{Y})^{\top}$$
(10.30)

$$= E[A(\mathbf{X} - E\mathbf{X})][A(\mathbf{X} - E\mathbf{X})]^{\top}$$
(10.31)

$$= E[A(\mathbf{X} - E\mathbf{X})(\mathbf{X} - E\mathbf{X})^{\top}A^{\top}]$$
(10.32)

$$= A[E(\mathbf{X} - E\mathbf{X})(\mathbf{X} - E\mathbf{X})^{\top}]A^{\top}$$
(10.33)

$$= AK_{\mathbf{X}}A^{\top}. \tag{10.34}$$

The proof of (10.29) is similar.  $\Box$ 

**Proposition 10.6.** Let  $\mathbf{X}$  and  $\mathbf{Y}$  be column vectors of n random variables such that

$$\mathbf{Y} = Q^{\top} \mathbf{X},\tag{10.35}$$

where  $Q\Lambda Q^{\top}$  is a diagonalization of  $K_{\mathbf{X}}$ . Then  $K_{\mathbf{Y}} = \Lambda$ , i.e., the random variables in  $\mathbf{Y}$  are uncorrelated and  $\operatorname{var} Y_i = \lambda_i$ , the *i*th diagonal element of  $\Lambda$ .

**Remark** The matrix  $K_{\mathbf{X}}$  is positive semidefinite, so that  $\lambda_i$ , being an eigenvalue of  $K_{\mathbf{X}}$ , is nonnegative by Proposition 10.4, as required for being the variance of a random variable.

Proof of Proposition 10.6. By Proposition 10.5,

$$K_{\mathbf{Y}} = Q^{\top} K_{\mathbf{X}} Q \tag{10.36}$$

$$=Q^{+}(QAQ^{+})Q \tag{10.37}$$

$$= (Q^{\top}Q)\Lambda(Q^{\top}Q) \tag{10.38}$$

$$= \Lambda. \tag{10.39}$$

Since  $K_{\mathbf{Y}} = \Lambda$  is a diagonal matrix, the random variables in  $\mathbf{Y}$  are uncorrelated. The variance of  $Y_i$  is given by the *i*th diagonal element of  $K_{\mathbf{Y}} = \Lambda$ , i.e.,  $\lambda_i$ . The proposition is proved.  $\Box$ 

**Corollary 10.7.** Let **X** be a column vector of n random variables such that  $QAQ^{\top}$  is a diagonalization of  $K_{\mathbf{X}}$ . Then

$$\mathbf{X} = Q\mathbf{Y},\tag{10.40}$$

where  $\mathbf{Y}$  is the column vector of n uncorrelated random variables prescribed in Proposition 10.6.

**Proposition 10.8.** Let  $\mathbf{X}$ ,  $\mathbf{Y}$ , and  $\mathbf{Z}$  be vectors of n random variables such that  $\mathbf{X}$  and  $\mathbf{Z}$  are independent and  $\mathbf{Y} = \mathbf{X} + \mathbf{Z}$ . Then

$$K_{\mathbf{Y}} = K_{\mathbf{X}} + K_{\mathbf{Z}}.\tag{10.41}$$

*Proof.* Omitted.  $\Box$ 

In communication engineering, the second moment of a random variable X is very often referred to as the *energy* of X. The total energy of a random vector  $\mathbf{X}$  is then equal to  $E \sum_{i} X_{i}^{2}$ . The following proposition shows that the total energy of a random vector is preserved by an orthogonal transformation.

**Proposition 10.9.** Let  $\mathbf{Y} = Q\mathbf{X}$ , where  $\mathbf{X}$  and  $\mathbf{Y}$  are column vectors of n random variables and Q is an orthogonal matrix. Then

$$E\sum_{i=1}^{n} Y_i^2 = E\sum_{i=1}^{n} X_i^2.$$
 (10.42)

Proof. Consider

$$\sum_{i=1}^{n} Y_i^2 = \mathbf{Y}^{\top} \mathbf{Y} \tag{10.43}$$

$$= (Q\mathbf{X})^{\top}(Q\mathbf{X}) \tag{10.44}$$
$$= \mathbf{X}^{\top}(Q^{\top}Q)\mathbf{X} \tag{10.45}$$

$$= \mathbf{X}^{\top} (Q^{\top} Q) \mathbf{X}$$
(10.45)

$$= \mathbf{X} \mathbf{X} \mathbf{X}$$
(10.46)

$$= \sum_{i=1}^{N} X_i^2. \tag{10.47}$$

The proposition is proved upon taking expectation on both sides.  $\Box$ 

#### 10.2 Definition

We now introduce the *differential entropy* for continuous random variables as the analog of the entropy for discrete random variables.

**Definition 10.10.** The differential entropy h(X) of a continuous random variable X with pdf f(x) is defined as

$$h(X) = -\int_{\mathcal{S}} f(x) \log f(x) dx = -E \log f(X).$$
(10.48)

The entropy of a discrete random variable X is a measure of the average amount of information contained in X, or equivalently, the average amount of uncertainty removed upon revealing the outcome of X. This was justified by the asymptotic achievability of the entropy bound for zero-error data compression discussed in Chapter 4 as well as the source coding theorem discussed in Chapter 5.

However, although entropy and differential entropy have similar mathematical forms, the latter does not serve as a measure of the average amount of information contained in a continuous random variable. In fact, a continuous random variable generally contains an infinite amount of information, as explained in the following example.

Example 10.11. Let X be uniformly distributed on [0, 1). Then we can write

$$X = .X_1 X_2 X_3 \cdots, (10.49)$$

the dyadic expansion of X, where  $X_1, X_2, X_3, \cdots$  is a sequence of fair bits<sup>4</sup>. Then

$$H(X) = H(X_1, X_2, X_3, \cdots)$$
(10.50)

$$=\sum_{i=1}^{\infty}H(X_i) \tag{10.51}$$

$$=\sum_{i=1}^{\infty}1\tag{10.52}$$

$$=\infty.$$
 (10.53)

In the following, we give two examples in which the differential entropy can be evaluated explicitly.

*Example 10.12 (Uniform Distribution).* Let X be uniformly distributed on [0, a). Then

$$h(X) = -\int_0^a \frac{1}{a} \log \frac{1}{a} dx = \log a.$$
(10.54)

From this example, we see immediately that h(X) < 0 if a < 1. This poses no contradiction because as we have mentioned, the differential entropy does not serve as a measure of the average amount of information contained in X. The physical meaning of differential entropy will be understood through the AEP for continuous random variables to be discussed in Section 10.4.

Example 10.13 (Gaussian Distribution). iGaussian Distribution Let  $X \sim \mathcal{N}(0, \sigma^2)$ . Then taking the natural logarithm, we have

 $<sup>^4</sup>$  Fair bits refer to i.i.d. bits, each being distributed uniformly on  $\{0,1\}.$  See Section 5.3.

$$h(X) = -\int f(x)\ln f(x)dx \tag{10.55}$$

$$= -\int f(x) \left( -\frac{x^2}{2\sigma^2} - \ln\sqrt{2\pi\sigma^2} \right) dx \qquad (10.56)$$

$$= \frac{1}{2\sigma^2} \int x^2 f(x) dx + \ln \sqrt{2\pi\sigma^2} \int f(x) dx$$
 (10.57)

$$= \frac{EX^2}{2\sigma^2} + \frac{1}{2}\ln(2\pi\sigma^2)$$
(10.58)

$$= \frac{\operatorname{var} X + (EX)^2}{2\sigma^2} + \frac{1}{2}\ln(2\pi\sigma^2)$$
(10.59)

$$= \frac{\sigma^2 + 0}{2\sigma^2} + \frac{1}{2}\ln(2\pi\sigma^2)$$
(10.60)

$$=\frac{1}{2} + \frac{1}{2}\ln(2\pi\sigma^2) \tag{10.61}$$

$$= \frac{1}{2}\ln e + \frac{1}{2}\ln(2\pi\sigma^2)$$
(10.62)

$$=\frac{1}{2}\ln(2\pi e\sigma^2)\tag{10.63}$$

in nats. Changing the base of the logarithm to any chosen positive value, we obtain

$$h(X) = \frac{1}{2}\log(2\pi e\sigma^2).$$
 (10.64)

The following two basic properties of differential entropy can readily be proved from the definition.

## Theorem 10.14 (Translation).

$$h(X+c) = h(X).$$
 (10.65)

*Proof.* Let Y = X + c, so that  $f_Y(y) = f_X(y - c)$  and  $S_Y = \{x + c : x \in S_X\}$ . Letting x = y - c in (10.48), we have

$$h(X) = -\int_{\mathcal{S}_X} f_X(x) \log f_X(x) dx \tag{10.66}$$

$$= -\int_{\mathcal{S}_Y} f_X(y-c) \log f_X(y-c) dy$$
 (10.67)

$$= -\int_{\mathcal{S}_Y} f_Y(y) \log f_Y(y) dy \tag{10.68}$$

$$=h(Y) \tag{10.69}$$

$$=h(X+c),$$
 (10.70)

accomplishing the proof.  $\Box$ 

Theorem 10.15 (Scaling). For  $a \neq 0$ ,

$$h(aX) = h(X) + \log|a|.$$
(10.71)

*Proof.* Let Y = aX, so that  $f_Y(y) = \frac{1}{|a|} f_X(\frac{y}{a})$  and  $S_Y = \{ax : x \in S_X\}$ . Letting  $x = \frac{y}{a}$  in (10.48), we have

$$h(X) = -\int_{\mathcal{S}_X} f_X(x) \log f_X(x) dx \tag{10.72}$$

$$= -\int_{\mathcal{S}_Y} f_X\left(\frac{y}{a}\right) \log f_X\left(\frac{y}{a}\right) \frac{dy}{|a|}$$
(10.73)

$$= -\int_{\mathcal{S}_Y} \frac{1}{|a|} f_X\left(\frac{y}{a}\right) \left[ \log\left(\frac{1}{|a|} f_X\left(\frac{y}{a}\right)\right) + \log|a| \right] dy \tag{10.74}$$

$$= -\int_{\mathcal{S}_{Y}} f_{Y}(y) \log f_{Y}(y) dy - \log |a| \int_{\mathcal{S}_{Y}} f_{Y}(y) dy$$
(10.75)

$$=h(Y) - \log|a| \tag{10.76}$$

$$= h(aX) - \log|a|.$$
 (10.77)

Hence,

$$h(aX) = h(X) + \log|a|,$$
(10.78)

accomplishing the proof.  $\hfill\square$ 

*Example 10.16.* We illustrate Theorem 10.14 and Theorem 10.15 by means of the Gaussian distribution. Let  $X \sim \mathcal{N}(\mu_X, \sigma_X^2)$ . By Theorem 10.14 (and Example 10.13),

$$h(X) = \frac{1}{2}\log(2\pi e\sigma_X^2).$$
 (10.79)

Let Y = aX. Then  $Y \sim \mathcal{N}(\mu_Y, \sigma_Y^2)$ , where  $\mu_Y = a\mu_X$  and  $\sigma_Y^2 = a^2\sigma_X^2$ . By (10.79),

$$h(Y) = \frac{1}{2}\log(2\pi e\sigma_Y^2) = \frac{1}{2}\log(2\pi ea^2\sigma_X^2) = \frac{1}{2}\log(2\pi e\sigma_X^2) + \log|a|, \quad (10.80)$$

which is consistent with Theorem 10.15.

Theorem 10.14 says that the differential entropy of a random variable is unchanged by translation. Theorem 10.15 says that the differential entropy of a random variable is generally changed by scaling. Specifically, if |a| > 1, the differential entropy is increased by  $\log |a|$ . If |a| < 1, the differential entropy is decreased by  $-\log |a|$  (note that  $-\log |a| > 0$ ). If a = -1, the differential entropy is unchanged.

These properties suggest that the differential entropy of a random variable depends only on the "spread" of the pdf. More specifically, the differential entropy increases with the "spread" of the pdf. This point will be further elaborated in Section 10.6.

## 10.3 Joint Differential Entropy, Conditional (Differential) Entropy, and Mutual Information

The definition for differential entropy is readily extended to multiple continuous random variables. In the rest of the chapter, we let  $\mathbf{X} = [X_1 X_2 \cdots X_n]$ .

**Definition 10.17.** The joint differential entropy  $h(\mathbf{X})$  of a random vector  $\mathbf{X}$  with joint pdf  $f(\mathbf{x})$  is defined as

$$h(\mathbf{X}) = -\int_{\mathcal{S}} f(\mathbf{x}) \log f(\mathbf{x}) d\mathbf{x} = -E \log f(\mathbf{X}).$$
(10.81)

It follows immediately from the above definition that if  $X_1, X_2, \dots, X_n$  are mutually independent, then

$$h(\mathbf{X}) = \sum_{i=1}^{n} h(X_i).$$
 (10.82)

The following two theorems are straightforward generalizations of Theorems 10.14 and 10.15, respectively. The proofs are omitted.

Theorem 10.18 (Translation).

$$h(\mathbf{X} + \mathbf{c}) = h(\mathbf{X}). \tag{10.83}$$

**Theorem 10.19 (Scaling).** Let A be a nonsingular  $n \times n$  matrix. Then

$$h(A\mathbf{X}) = h(\mathbf{X}) + \log|\det(A)|. \tag{10.84}$$

Theorem 10.20 (Multivariate Gaussian Distribution). Let  $\mathbf{X} \sim \mathcal{N}(\boldsymbol{\mu}, K)$ . Then

$$h(\mathbf{X}) = \frac{1}{2} \log \left[ (2\pi e)^n |K| \right].$$
(10.85)

*Proof.* Let K be diagonalizable as  $Q\Lambda Q^{\top}$ . Write  $\mathbf{X} = Q\mathbf{Y}$  as in Corollary 10.7, where the random variables in  $\mathbf{Y}$  are uncorrelated with var  $Y_i = \lambda_i$ , the *i*th diagonal element of  $\Lambda$ . Since  $\mathbf{X}$  is Gaussian, so is  $\mathbf{Y}$ . Then the random variables in  $\mathbf{Y}$  are mutually independent because they are uncorrelated. Now consider

10.3 Joint and Conditional Differential Entropy 235

$$h(\mathbf{X}) = h(Q\mathbf{Y}) \tag{10.86}$$

$$\stackrel{a)}{=} h(\mathbf{Y}) + \log |\det(Q)| \tag{10.87}$$

$$\stackrel{b)}{=} h(\mathbf{Y}) + 0 \tag{10.88}$$

$$\stackrel{c)}{=} \sum_{i=1}^{n} h(Y_i) \tag{10.89}$$

$$\stackrel{d)}{=} \sum_{i=1}^{n} \frac{1}{2} \log(2\pi e\lambda_i) \tag{10.90}$$

$$= \frac{1}{2} \log \left[ (2\pi e)^n \prod_{i=1}^n \lambda_i \right]$$
(10.91)

$$\stackrel{e)}{=} \frac{1}{2} \log[(2\pi e)^n |\Lambda|]$$
(10.92)

$$\stackrel{f}{=} \frac{1}{2} \log[(2\pi e)^n |K|] \tag{10.93}$$

In the above

- a) follows from Theorem 10.19;
- b) follows from (10.23);
- c) follows from (10.82) since  $Y_1, Y_2, \dots, Y_n$  are mutually independent;
- d) follows from Example 10.16;
- e) follows because  $\Lambda$  is a diagonal matrix;
- f) follows because

$$|\Lambda| = |Q||\Lambda||Q^{\top}| = |Q\Lambda Q^{\top}| = |K|.$$
(10.94)

The theorem is proved.  $\Box$ 

In describing a communication system, we very often specify the relation between two random variables X and Y through a conditional distribution p(y|x) (if Y is discrete) or a conditional pdf f(y|x) (if Y is continuous) defined for all x, even though certain x may not be in  $\mathcal{S}_X$ . This is made precise by the following two definitions.

**Definition 10.21.** Let X and Y be two jointly distributed random variables with Y being discrete. The random variable Y is related to the random variable X through a conditional distribution p(y|x) defined for all x means that for all x and y,

$$\Pr\{X \le x, Y = y\} = \int_{-\infty}^{x} p_{Y|X}(y|u) dF_X(u).$$
(10.95)

**Definition 10.22.** Let X and Y be two jointly distributed random variables with Y being continuous. The random variable Y is related to the random variable X through a conditional pdf f(y|x) defined for all x means that for all x and y,

$$F_{XY}(x,y) = \int_{-\infty}^{x} F_{Y|X}(y|u) dF_X(u), \qquad (10.96)$$

where

$$F_{Y|X}(y|x) = \int_{-\infty}^{y} f_{Y|X}(v|x)dv.$$
 (10.97)

**Definition 10.23.** Let X and Y be jointly distributed random variables where Y is continuous and is related to X through a conditional pdf f(y|x) defined for all x. The conditional differential entropy of Y given  $\{X = x\}$  is defined as

$$h(Y|X = x) = -\int_{\mathcal{S}_Y(x)} f(y|x) \log f(y|x) dy$$
(10.98)

where  $S_Y(x) = \{y : f(y|x) > 0\}$ , and the conditional differential entropy of Y given X is defined as

$$h(Y|X) = -\int_{\mathcal{S}_X} h(Y|X=x) dF(x) = -E \log f(Y|X).$$
(10.99)

**Definition 10.24.** A pdf f(y) is bounded if there exists a positive real number M such that f(y) < M for all y.

**Definition 10.25.** A conditional pdf f(y|x) defined for all x is bounded if there exists a positive real number M such that f(y|x) < M for all x and y.

**Proposition 10.26.** Let X and Y be jointly distributed random variables where Y is continuous and is related to X through a conditional pdf f(y|x)defined for all x. If f(y|x) is bounded, then f(y) exists and is bounded, and is given by

$$f(y) = \int f(y|x)dF(x).$$
 (10.100)

*Proof.* We first give an informal proof of the proposition to illustrate the idea. From (10.96), we have

$$F(y) = F_{XY}(\infty, y) = \int F(y|x)dF(x).$$
 (10.101)

Assuming that F(y) is differentiable, the proposition can be proved by considering

10.3 Joint and Conditional Differential Entropy 237

$$f(y) = \frac{dF(y)}{dy} \tag{10.102}$$

$$= \frac{d}{dy} \int F(y|x) dF(x)$$
(10.103)

$$= \int \left[\frac{d}{dy}F(y|x)\right]dF(x) \tag{10.104}$$

$$= \int f(y|x)dF(x), \qquad (10.105)$$

provided that (10.104) is justified.

We now present a formal proof. Let M be a positive real number such that

$$f(y|x) < M \tag{10.106}$$

for all x and y. Since

$$f(y|x) = \frac{d}{dy}F(y|x) \tag{10.107}$$

$$= \lim_{\Delta_y \to 0} \frac{1}{\Delta_y} [F(y + \Delta_y | x) - F(y | x)], \qquad (10.108)$$

for all  $\epsilon > 0$ , there exists  $\delta > 0$  such that for all  $\Delta_y \in (-\delta, \delta)$ ,

$$\left|\frac{1}{\Delta_y}[F(y+\Delta_y|x) - F(y|x)]\right| < f(y|x) + \epsilon < M + \epsilon.$$
(10.109)

It follows that

$$\int \left| \frac{1}{\Delta_y} [F(y + \Delta_y | x) - F(y | x)] \right| dF(x) < \int (M + \epsilon) dF(x) \qquad (10.110)$$
  
$$< M + \epsilon. \qquad (10.111)$$

$$A + \epsilon. \tag{10.111}$$

Now consider

$$\lim_{\Delta_y \to 0} \frac{1}{\Delta_y} [F(y + \Delta_y) - F(y)]$$
  
= 
$$\lim_{\Delta_y \to 0} \frac{1}{\Delta_y} \left[ \int F(y + \Delta_y | x) dF(x) - \int F(y | x) dF(x) \right]$$
(10.112)

$$= \lim_{\Delta_y \to 0} \int \frac{1}{\Delta_y} [F(y + \Delta_y | x) - F(y | x)] dF(x)$$
(10.113)

$$= \int \lim_{\Delta_y \to 0} \frac{1}{\Delta_y} [F(y + \Delta_y | x) - F(y | x)] dF(x)$$
(10.114)

$$= \int f(y|x)dF(x) \tag{10.115}$$

$$< M, \tag{10.116}$$

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where we have invoked the dominated convergence theorem<sup>5</sup> in light of (10.111) to obtain (10.114), and (10.116) follows from (10.106). Therefore,

$$f(y) = \frac{dF(y)}{dy} \tag{10.117}$$

$$= \lim_{\Delta_y \to 0} \frac{1}{\Delta_y} [F(y + \Delta_y) - F(y)]$$
(10.118)

$$= \int f(y|x)dF(x) \tag{10.119}$$

$$< M.$$
 (10.120)

This completes the proof.  $\Box$ 

The above proposition says that if Y is related to X through a bounded conditional pdf f(y|x), then the pdf of Y exists regardless of the distribution of X. The next proposition is a generalization to random vectors, and the proof is left as an exercise. As we will see, the theory in the rest of this chapter and in the next chapter will be developed around this important fact.

**Proposition 10.27.** Let  $\mathbf{X}$  and  $\mathbf{Y}$  be jointly distributed random vectors where  $\mathbf{Y}$  is continuous and is related to  $\mathbf{X}$  through a conditional pdf  $f(\mathbf{y}|\mathbf{x})$  defined for all  $\mathbf{x}$ . If  $f(\mathbf{y}|\mathbf{x})$  is bounded, then  $f(\mathbf{y})$  exists and is bounded, and is given by

$$f(\mathbf{y}) = \int f(\mathbf{y}|\mathbf{x}) dF(\mathbf{x}). \tag{10.121}$$

**Definition 10.28.** Let X and Y be jointly distributed random variables where Y is continuous and is related to X through a bounded conditional pdf f(y|x) defined for all x. The mutual information between X and Y is defined as

$$I(X;Y) = \int_{\mathcal{S}_X} \int_{\mathcal{S}_Y(x)} f(y|x) \log \frac{f(y|x)}{f(y)} dy \, dF(x)$$
(10.122)

$$= E \log \frac{f(Y|X)}{f(Y)}, \qquad (10.123)$$

where f(y) exists and is given in (10.100) by Proposition 10.26. When both X and Y are continuous,

<sup>5</sup> The dominated convergence theorem states that

$$\lim_{k \to \infty} \int_S f_k d\mu = \int_S f d\mu,$$

where  $f_k$  and f are real functions defined on S such that  $f_k \to f$  pointwise, provided that  $|f_k| < g$  pointwise on S and  $\int_S |g| d\mu < \infty$ . See [284] for details.

10.3 Joint and Conditional Differential Entropy 239

$$I(X;Y) = E \log \frac{f(Y|X)}{f(Y)} = E \log \frac{f(X,Y)}{f(X)f(Y)}.$$
 (10.124)

Together with our discussion on discrete random variables in Chapter 2, the mutual information I(X; Y) is defined when each of the random variables involved can be either discrete or continuous. In the same way, we can define the conditional mutual information I(X; Y|T).

**Definition 10.29.** Let X, Y, and T be jointly distributed random variables where Y is continuous and is related to (X,T) through a bounded conditional pdf f(y|x,t) defined for all x and t. The mutual information between X and Y given T is defined as

$$I(X;Y|T) = \int_{\mathcal{S}_T} I(X;Y|T=t) dF(t) = E \log \frac{f(Y|X,T)}{f(Y|T)},$$
 (10.125)

where

$$I(X;Y|T=t) = \int_{\mathcal{S}_X(t)} \int_{\mathcal{S}_Y(x,t)} f(y|x,t) \log \frac{f(y|x,t)}{f(y|t)} dy \, dF(x|t). \quad (10.126)$$

We now give a physical interpretation of I(X;Y) when the joint pdf f(x,y) exists. For simplicity, we assume that f(x,y) > 0 for all x and y. Let  $\Delta$  be a small positive quantity. For all integer *i*, define the interval

$$A_x^i = [i\Delta, (i+1)\Delta) \tag{10.127}$$

in  $\Re$ , and for all integer j, define the interval

$$A_{y}^{j} = [j\Delta, (j+1)\Delta). \tag{10.128}$$

For all integers i and j, define the set

$$A_{xy}^{i,j} = A_x^i \times A_y^j, \tag{10.129}$$

which corresponds to a "rectangle" in  $\Re^2$ .

We now introduce two discrete random variables  $\hat{X}_{\Delta}$  and  $\hat{Y}_{\Delta}$  defined by

$$\begin{cases} \dot{X}_{\Delta} = i \text{ if } X \in A_x^i \\ \dot{Y}_{\Delta} = j \text{ if } Y \in A_y^j. \end{cases}$$
(10.130)

The random variables  $\hat{X}_{\Delta}$  and  $\hat{Y}_{\Delta}$  are quantizations of the continuous random variables X and Y, respectively. For all i and j, let  $(x_i, y_j) \in A_{xy}^{i,j}$ . Then

240 10 Differential Entropy

$$I(\hat{X}_{\Delta}; \hat{Y}_{\Delta}) = \sum_{i} \sum_{j} \Pr\{(\hat{X}_{\Delta}, \hat{Y}_{\Delta}) = (i, j)\} \log \frac{\Pr\{(\hat{X}_{\Delta}, \hat{Y}_{\Delta}) = (i, j)\}}{\Pr\{\hat{X}_{\Delta} = i\} \Pr\{\hat{Y}_{\Delta} = j\}}$$
(10.131)

$$\approx \sum_{i} \sum_{j} f(x_i, y_j) \Delta^2 \log \frac{f(x_i, y_j) \Delta^2}{(f(x_i)\Delta)(f(y_j)\Delta)}$$
(10.132)

$$= \sum_{i} \sum_{j} f(x_i, y_j) \Delta^2 \log \frac{f(x_i, y_j)}{f(x_i) f(y_j)}$$
(10.133)

$$\approx \int \int f(x,y) \log \frac{f(x,y)}{f(x)f(y)} dxdy$$
(10.134)

$$=I(X;Y).$$
 (10.135)

Therefore, I(X;Y) it can be interpreted as the limit of  $I(\hat{X}_{\Delta};\hat{Y}_{\Delta})$  as  $\Delta \to 0$ . This interpretation carries over to the case when X and Y have a general joint distribution<sup>6</sup> (see Dobrushin [85]). As  $I(\hat{X}_{\Delta};\hat{Y}_{\Delta})$  is always nonnegative, this suggests that I(X;Y) is also always nonnegative, which will be established in Theorem 10.33.

**Definition 10.30.** Let X be a continuous random variable and Y be a discrete random variable, where X is related to Y through a bounded conditional pdf f(x|y). The conditional entropy of Y given X is defined as

$$H(Y|X) = H(Y) - I(X;Y), (10.136)$$

where I(X;Y) is defined as in Definition 10.28.

**Proposition 10.31.** For two random variables X and Y,

$$h(Y) = h(Y|X) + I(X;Y)$$
(10.137)

if Y is continuous, and

$$H(Y) = H(Y|X) + I(X;Y)$$
(10.138)

if Y is discrete.

## Proposition 10.32 (Chain Rule for Differential Entropy).

$$h(X_1, X_2, \cdots, X_n) = \sum_{i=1}^n h(X_i | X_1, \cdots, X_{i-1}).$$
(10.139)

 $^{6}$  In the general setting, the mutual information between X and Y is defined as

$$I(X;Y) = \int_{\mathcal{S}_{XY}} \left( \log \frac{dP_{XY}}{d(P_X \times P_Y)} \right) dP_{XY},$$

where  $P_{XY}$ ,  $P_X$ , and  $P_Y$  are the probability measures of (X, Y), X, and Y, respectively, and  $\frac{dP_{XY}}{d(P_X \times P_Y)}$  denotes the Radon-Nikodym derivative of  $P_{XY}$  with respect to the product measure  $P_X \times P_Y$ .

The proofs of these propositions are left as an exercise.

#### Theorem 10.33.

$$I(X;Y) \ge 0,$$
 (10.140)

with equality if and only if X is independent of Y.

Proof. Consider

$$I(X;Y) = \int_{\mathcal{S}_X} \int_{\mathcal{S}_Y(x)} f(y|x) \log \frac{f(y|x)}{f(y)} dy \, dF_X(x)$$
(10.141)

$$\geq (\log e) \int_{\mathcal{S}_X} \int_{\mathcal{S}_Y(x)} f(y|x) \left(1 - \frac{f(y)}{f(y|x)}\right) dy \, dF_X(x) \tag{10.142}$$

$$= (\log e) \int_{\mathcal{S}_X} \left[ \int_{\mathcal{S}_Y(x)} f(y|x) dy - \int_{\mathcal{S}_Y(x)} f(y) dy \right] dF_X(x)$$
(10.143)

$$\geq (\log e) \int_{\mathcal{S}_X} (1-1)dF_X(x) \tag{10.144}$$

$$= 0,$$
 (10.145)

where (10.142) results from an application of the fundamental inequality (Corollary 2.102), and (10.144) follows from

$$\int_{\mathcal{S}_Y(x)} f(y) dy \le 1. \tag{10.146}$$

This proves (10.140).

For equality to hold in (10.140), equality must hold in (10.142) for all  $x \in S_X$  and all  $y \in S_Y(x)$ , and equality must hold in (10.144) for all  $x \in S_X$ . For the former, this is the case if and only if

$$f(y|x) = f(y)$$
 for all  $x \in \mathcal{S}_X$  and  $y \in \mathcal{S}_Y(x)$ , (10.147)

which implies

$$\int_{\mathcal{S}_{Y}(x)} f(y) dy = \int_{\mathcal{S}_{Y}(x)} f(y|x) dy = 1,$$
(10.148)

i.e., equality holds in (10.144). Thus (10.147) is a necessary and sufficient condition for equality to hold in (10.140).

It is immediate that if X and Y are independent, then (10.147) holds. It remains to prove the converse. To this end, observe that (10.148), implied by (10.147), is equivalent to that f(y) = 0 on  $S_Y \setminus S_Y(x)$  a.e. (almost everywhere). By the definition of  $S_Y$ , this means that  $S_Y \setminus S_Y(x) \subset S_Y^c$ , or  $S_Y = S_Y(x)$ . Since this holds for all  $x \in S_X$ , we conclude that f(y|x) = f(y) for all  $(x, y) \in S_X \times S_Y$ , i.e., X and Y are independent. The theorem is proved.  $\Box$  242 10 Differential Entropy

Corollary 10.34.

$$I(X;Y|T) \ge 0,$$
 (10.149)

with equality if and only if X is independent of Y conditioning on T.

*Proof.* This follows directly from (10.125).  $\Box$ 

Corollary 10.35 (Conditioning Does Not Increase Differential Entropy).

$$h(X|Y) \le h(X) \tag{10.150}$$

with equality if and only if X and Y are independent.

Corollary 10.36 (Independence Bound for Differential Entropy).

$$h(X_1, X_2, \cdots, X_n) \le \sum_{i=1}^n h(X_i)$$
 (10.151)

with equality if and only if  $i = 1, 2, \dots, n$  are mutually independent.

# 10.4 The AEP for Continuous Random Variables

The Weak AEP for discrete random variables discussed in Chapter 5 states that for n i.i.d. random variables  $X_1, X_2, \dots, X_n$  with generic discrete random variable  $X, p(X_1, X_2, \dots, X_n)$  is close to  $2^{-nH(X)}$  with high probability when n is large (Theorem 5.1, Weak AEP I). This fundamental property of entropy leads to the definition of weak typicality, and as a consequence, the total number of weakly typical sequences is approximately equal to  $2^{nH(X)}$ (Theorem 5.3, Weak AEP II).

In the following, we develop the AEP for continuous random variables in the same way we developed the Weak AEP for discrete random variables. Some of the proofs are exactly the same as their discrete analogs, and they are omitted. We note that for continuous random variables, the notion of strong typicality does not apply because the probability that a continuous random variable takes a particular value is equal to zero.

## Theorem 10.37 (AEP I for Continuous Random Variables).

$$-\frac{1}{n}\log f(\mathbf{X}) \to h(X) \tag{10.152}$$

in probability as  $n \to \infty$ , i.e., for any  $\epsilon > 0$ , for n sufficiently large,

$$\Pr\left\{\left|-\frac{1}{n}\log f(\mathbf{X}) - h(X)\right| < \epsilon\right\} > 1 - \epsilon.$$
(10.153)

**Definition 10.38.** The typical set  $W_{[X]\epsilon}^n$  with respect to f(x) is the set of sequences  $\mathbf{x} = (x_1, x_2, \dots, x_n) \in \mathcal{X}^n$  such that

$$\left| -\frac{1}{n} \log f(\mathbf{x}) - h(X) \right| < \epsilon, \tag{10.154}$$

or equivalently,

$$h(X) - \epsilon < -\frac{1}{n}\log f(\mathbf{x}) < h(X) + \epsilon, \qquad (10.155)$$

where  $\epsilon$  is an arbitrarily small positive real number. The sequences in  $W_{[X]\epsilon}^n$  are called  $\epsilon$ -typical sequences.

The quantity

$$-\frac{1}{n}\log f(\mathbf{x}) = -\frac{1}{n}\sum_{k=1}^{n}\log f(x_k)$$
(10.156)

is called the *empirical differential entropy* of the sequence  $\mathbf{x}$ . The empirical differential entropy of a typical sequence is close to the true differential entropy h(X).

If the pdf f(x) is continuous, we see from (10.156) that the empirical differential entropy is continuous in  $\mathbf{x}$  with respect to the Euclidean distance in  $\Re^n$ , so that if  $\mathbf{x}$  is  $\epsilon$ -typical, then all the sequences in the neighborhood of  $\mathbf{x}$  are also  $\epsilon$ -typical. As such, the number of  $\epsilon$ -typical sequences is uncountably. Therefore, it is not meaningful to discuss the cardinality of a typical set as in the discrete case. Instead, the "size" of a typical set is measured by its volume.

**Definition 10.39.** The volume of a set A in  $\Re^n$  is defined as

$$\operatorname{Vol}(A) = \int_{A} d\mathbf{x}.$$
 (10.157)

**Theorem 10.40 (AEP II for Continuous Random Variables).** *The following hold for any*  $\epsilon > 0$ *:* 

1) If  $\mathbf{x} \in W_{[X]\epsilon}^n$ , then

$$2^{-n(h(X)+\epsilon)} < f(\mathbf{x}) < 2^{-n(h(X)-\epsilon)}.$$
(10.158)

2) For n sufficiently large,

$$\Pr\{\mathbf{X} \in W_{[X]\epsilon}^n\} > 1 - \epsilon.$$
(10.159)

3) For n sufficiently large,

$$(1-\epsilon)2^{n(h(X)-\epsilon)} < \operatorname{Vol}(W_{[X]\epsilon}^n) < 2^{n(h(X)+\epsilon)}.$$
(10.160)

#### 244 10 Differential Entropy

*Proof.* Property 1 follows immediately from the definition of  $W_{[X]\epsilon}^n$  in (10.155). Property 2 is equivalent to Theorem 10.37. To prove Property 3, we use the lower bound in (10.158) and consider

$$1 \ge \Pr\{W_{[X]\epsilon}^n\} \tag{10.161}$$

$$= \int_{W_{[X]\epsilon}^n} f(\mathbf{x}) \, d\mathbf{x} \tag{10.162}$$

$$> \int_{W_{[X]\epsilon}^n} 2^{-n(h(X)+\epsilon)} \, d\mathbf{x} \tag{10.163}$$

$$> 2^{-n(h(X)+\epsilon)} \int_{W_{[X]\epsilon}^n} d\mathbf{x}$$
(10.164)

$$=2^{-n(h(X)+\epsilon)}\operatorname{Vol}(W^n_{[X]\epsilon}),\qquad(10.165)$$

which implies

$$\operatorname{Vol}(W_{[X]\epsilon}^n) < 2^{n(h(X)+\epsilon)}.$$
(10.166)

Note that this upper bound holds for any  $n \ge 1$ . On the other hand, using the upper bound in (10.158) and Theorem 10.37, for n sufficiently large, we have

$$1 - \epsilon < \Pr\{W_{[X]\epsilon}^n\} \tag{10.167}$$

$$= \int_{W_{[\mathbf{X}]\epsilon}^{n}} f(\mathbf{x}) \, d\mathbf{x} \tag{10.168}$$

$$<\int_{W_{[X]\epsilon}^n} 2^{-n(h(X)-\epsilon)} \, d\mathbf{x} \tag{10.169}$$

$$= 2^{-n(h(X)-\epsilon)} \operatorname{Vol}(W^n_{[X]\epsilon}).$$
(10.170)

Then

$$\operatorname{Vol}(W_{[X]\epsilon}^n) > (1-\epsilon)2^{n(h(X)-\epsilon)}.$$
(10.171)

Combining (10.166) and (10.171) gives Property 3. The theorem is proved.  $\Box$ 

From the AEP for continuous random variables, we see that the volume of the typical set is approximately equal to  $2^{nh(X)}$  when *n* is large. This gives the following physical interpretations of differential entropy. First, the fact that h(X) can be negative does not incur any difficulty because  $2^{nh(X)}$  is always positive. Second, if the differential entropy is large, then the volume of the typical set is large; if the differential entropy is small (not in magnitude but in value), then the volume of the typical set is small.

## **10.5** Informational Divergence

We first extend the definition of informational divergence introduced in Section 2.5 to pdf's.

**Definition 10.41.** Let f and g be two pdf's defined on  $\Re^n$  with supports  $S_f$ and  $S_q$ , respectively. The informational divergence between f and g is defined as

$$D(f||g) = \int_{\mathcal{S}_f} f(x) \log \frac{f(x)}{g(x)} dx = E_f \log \frac{f(X)}{g(X)},$$
(10.172)

where  $E_f$  denotes expectation with respect to f.

**Remark** In the above definition, we adopt the convention  $c \log \frac{c}{0} = \infty$  for c > 0. Furthermore, if the set

$$\mathcal{S}_f \setminus \mathcal{S}_g = \{ x : f(x) > 0 \text{ and } g(x) = 0 \}$$

$$(10.173)$$

has zero Lebesgue measure, the integral is evaluated by excluding this set from  $\mathcal{S}_f$ . Therefore, if  $D(f||g) < \infty$ , then  $\mathcal{S}_f \setminus \mathcal{S}_g$  has zero Lebesgue measure, i.e.,  $\mathcal{S}_f$  is essentially a subset of  $\mathcal{S}_g$ .

Theorem 10.42 (Divergence Inequality). Let f and g be two pdf's defined on  $\Re^n$ . Then

$$D(f||g) \ge 0, \tag{10.174}$$

with equality if and only if f = g a.e.

Proof. Consider

$$D(f||g) = \int_{\mathcal{S}_f} f(x) \log \frac{f(x)}{g(x)} dx \tag{10.175}$$

$$= (\log e) \int_{\mathcal{S}_f} f(x) \ln \frac{f(x)}{g(x)} dx \qquad (10.176)$$

$$\geq (\log e) \int_{\mathcal{S}_f} f(x) \left( 1 - \frac{g(x)}{f(x)} \right) dx \tag{10.177}$$

$$= (\log e) \left[ \int_{\mathcal{S}_f} f(x) dx - \int_{\mathcal{S}_f} g(x) dx \right]$$
(10.178)  
> 0, (10.179)

$$\geq 0, \tag{10.179}$$

where (10.177) follows from the fundamental inequality (Corollary 2.102) and (10.179) follows from

$$\int_{\mathcal{S}_f} g(x) dx \le 1 = \int_{\mathcal{S}_f} f(x) dx.$$
(10.180)

Equality holds in (10.177) if and only if f(x) = g(x) on  $S_f$  a.e., which implies

246 10 Differential Entropy

$$\int_{\mathcal{S}_f} g(x)dx = \int_{\mathcal{S}_f} f(x)dx = 1,$$
(10.181)

i.e., equality holds in (10.179). Then we see from (10.181) that g(x) = 0 on  $S_f^c$  a.e. Hence, we conclude that equality holds in (10.174) if and only if f = g a.e. The theorem is proved.  $\Box$ 

## **10.6** Maximum Differential Entropy Distributions

In Section 2.9, we have discussed maximum entropy distributions for a discrete random variable. We now extend this theme to multiple continuous random variables. Specifically, we are interested in the following problem: Maximize h(f) over all pdf f defined on a subset S of  $\Re^n$ , subject to

$$\int_{\mathcal{S}_f} r_i(\mathbf{x}) f(\mathbf{x}) d\mathbf{x} = a_i \quad \text{for } 1 \le i \le m,$$
(10.182)

where  $S_f \subset S$  and  $r_i(\mathbf{x})$  is defined for all  $\mathbf{x} \in S$ .

Theorem 10.43. Let

$$f^*(\mathbf{x}) = e^{-\lambda_0 - \sum_{i=1}^m \lambda_i r_i(\mathbf{x})}$$
(10.183)

for all  $\mathbf{x} \in S$ , where  $\lambda_0, \lambda_1, \dots, \lambda_m$  are chosen such that the constraints in (10.182) are satisfied. Then  $f^*$  maximizes h(f) over all pdf f defined on S, subject to the constraints in (10.182).

*Proof.* The proof is analogous to that of Theorem 2.50. The details are omitted.  $\Box$ 

**Corollary 10.44.** Let  $f^*$  be a pdf defined on S with

$$f^*(\mathbf{x}) = e^{-\lambda_0 - \sum_{i=1}^m \lambda_i r_i(\mathbf{x})}$$
(10.184)

for all  $\mathbf{x} \in S$ . Then  $f^*$  maximizes h(f) over all pdf f defined on S, subject to the constraints

$$\int_{\mathcal{S}_f} r_i(\mathbf{x}) f(\mathbf{x}) d\mathbf{x} = \int_{\mathcal{S}} r_i(\mathbf{x}) f^*(\mathbf{x}) d\mathbf{x} \quad \text{for } 1 \le i \le m.$$
(10.185)

**Theorem 10.45.** Let X be a continuous random variable with  $EX^2 = \kappa$ . Then

$$h(X) \le \frac{1}{2}\log(2\pi e\kappa),$$
 (10.186)

with equality if and only if  $X \sim \mathcal{N}(0, \kappa)$ .

*Proof.* The problem here is to maximize h(f) subject to the constraint

$$\int x^2 f(x) dx = \kappa. \tag{10.187}$$

An application of Theorem 10.43 yields

$$f^*(x) = ae^{-bx^2} \tag{10.188}$$

which is identified as a Gaussian distribution with zero mean. In order that the constraint (10.187) is satisfied, we must have

$$a = \frac{1}{\sqrt{2\pi\kappa}}$$
 and  $b = \frac{1}{2\kappa}$ . (10.189)

Hence, in light of (10.64) in Example 10.13, we have proved (10.186) with equality if and only if  $X \sim \mathcal{N}(0, \kappa)$ .  $\Box$ 

**Theorem 10.46.** Let X be a continuous random variable with mean  $\mu$  and variance  $\sigma^2$ . Then

$$h(X) \le \frac{1}{2}\log(2\pi e\sigma^2),$$
 (10.190)

with equality if and only if  $X \sim \mathcal{N}(\mu, \sigma^2)$ .

*Proof.* Let  $X' = X - \mu$ . Then

$$EX' = E(X - \mu) = EX - \mu = 0$$
(10.191)

and

$$E(X')^2 = E(X - \mu)^2 = \operatorname{var} X = \sigma^2.$$
 (10.192)

Applying Theorem 10.14 and Theorem 10.45, we have

$$h(X) = h(X') \le \frac{1}{2}\log(2\pi e\sigma^2),$$
 (10.193)

and equality holds if and only if  $X' \sim \mathcal{N}(0, \sigma^2)$ , or  $X \sim \mathcal{N}(\mu, \sigma^2)$ . The theorem is proved.  $\Box$ 

**Remark** Theorem 10.45 says that with the constraint  $EX^2 = \kappa$ , the differential entropy is maximized by the distribution  $\mathcal{N}(0,\kappa)$ . If we impose the additional constraint that EX = 0, then  $\operatorname{var} X = EX^2 = \kappa$ . By Theorem 10.46, the differential entropy is still maximized by  $\mathcal{N}(0,\kappa)$ .

We have mentioned at the end of Section 10.2 that the differential entropy of a random variable increases with the "spread" of the pdf. Though a simple consequence of Theorem 10.45, the above theorem makes this important interpretation precise. By rewriting the upper bound in (10.193), we obtain 248 10 Differential Entropy

$$h(X) \le \log \sigma + \frac{1}{2} \log(2\pi e).$$
 (10.194)

That is, the differential entropy is at most equal to the logarithm of the standard deviation plus a constant. In particular, the differential entropy tends to  $-\infty$  as the standard deviation tends to 0.

The next two theorems are the vector generalizations of Theorems 10.45 and 10.46.

**Theorem 10.47.** Let X be a vector of n continuous random variables with correlation matrix  $\tilde{K}$ . Then

$$h(\mathbf{X}) \le \frac{1}{2} \log\left[ (2\pi e)^n |\tilde{K}| \right],$$
 (10.195)

with equality if and only if  $\mathbf{X} \sim \mathcal{N}(0, \tilde{K})$ .

*Proof.* By Theorem 10.43, the joint pdf that maximizes  $h(\mathbf{X})$  has the form

$$f^*(\mathbf{x}) = e^{-\lambda_0 - \sum_{i,j} \lambda_{ij} x_i x_j} = e^{-\lambda_0 - \mathbf{x}^\top L \mathbf{x}},$$
(10.196)

where  $L = [\lambda_{ij}]$ . Thus  $f^*$  is a multivariate Gaussian distribution with zero mean. Therefore,

$$cov(X_i, X_j) = EX_i X_j - (EX_i)(EX_j) = EX_i X_j$$
 (10.197)

for all *i* and *j*. Since  $f^*$  is constrained by  $\tilde{K}$ ,  $\lambda_0$  and *L* have the unique solution given by

$$e^{-\lambda_0} = \frac{1}{\left(\sqrt{2\pi}\right)^n |\tilde{K}|^{1/2}}$$
(10.198)

and

$$L = \frac{1}{2}\tilde{K}^{-1},\tag{10.199}$$

so that

$$f^*(\mathbf{x}) = \frac{1}{\left(\sqrt{2\pi}\right)^n |\tilde{K}|^{1/2}} e^{-\frac{1}{2}\mathbf{x}^\top \tilde{K}^{-1}\mathbf{x}},$$
 (10.200)

the joint pdf of  $\mathbf{X} \sim \mathcal{N}(0, \tilde{K})$ . Hence, by Theorem 10.20, we have proved (10.195) with equality if and only if  $\mathbf{X} \sim \mathcal{N}(0, \tilde{K})$ .  $\Box$ 

**Theorem 10.48.** Let **X** be a vector of n continuous random variables with mean  $\mu$  and covariance matrix K. Then

$$h(\mathbf{X}) \le \frac{1}{2} \log\left[(2\pi e)^n |K|\right],$$
 (10.201)

with equality if and only if  $\mathbf{X} \sim \mathcal{N}(\boldsymbol{\mu}, K)$ .

*Proof.* Similar to the proof of Theorem 10.46.  $\Box$ 

# Problems

- 1. Prove Propositions 10.3 and 10.8.
- 2. Show that the joint pdf of a multivariate Gaussian distribution integrates to 1.
- 3. Show that a symmetric positive definite matrix is a covariance matrix.
- 4. Let

$$K = \begin{bmatrix} 7/4 & \sqrt{2}/4 & -3/4 \\ \sqrt{2}/4 & 5/2 & -\sqrt{2}/4 \\ -3/4 & -\sqrt{2}/4 & 7/4 \end{bmatrix}.$$

- a) Find the eigenvalues and eigenvectors of K.
- b) Show that K is positive definite.
- c) Suppose K is the covariance matrix of a random vector X = [X<sub>1</sub> X<sub>2</sub> X<sub>3</sub>]<sup>⊤</sup>.
  i) Find the coefficient of correlation between X<sub>i</sub> and X<sub>j</sub> for 1 ≤ i < j ≤ 3.</li>
  - ii) Find an uncorrelated random vector  $\mathbf{Y} = [Y_1 \ Y_2 \ Y_3]$  such that  $\mathbf{X}$  is a linear transformation of  $\mathbf{Y}$ .
  - iii) Determine the covariance matrix of Y.
- 5. Prove Theorem 10.19.
- 6. For continuous random variables X and Y, discuss why I(X; X) is not equal to h(X).
- 7. Prove Proposition 10.27.
- 8. Verify Corollary 10.44.
- 9. Each of the following continuous distributions can be obtained as the distribution that maximizes the differential entropy subject to a suitable set of constraints:
  - a) the exponential distribution,

$$f(x) = \lambda e^{-\lambda x}$$

for  $x \ge 0$ , where  $\lambda > 0$ ;

b) the Laplace distribution,

$$f(x) = \frac{1}{2}\lambda e^{-\lambda|x|}$$

for  $-\infty < x < \infty$ , where  $\lambda > 0$ ;

c) the gamma distribution,

$$f(x) = \frac{\lambda}{\Gamma(\alpha)} (\lambda x)^{\alpha - 1} e^{-\lambda x}$$

for  $x \ge 0$ , where  $\lambda, \alpha > 0$  and  $\Gamma(z) = \int_0^\infty t^{z-1} e^{-t} dt$ ; d) the beta distribution,

$$f(x) = \frac{\Gamma(p+q)}{\Gamma(p)\Gamma(q)} x^{p-1} (1-x)^{q-1}$$

for  $0 \le x \le 1$ , where p, q > 0;

- 250 10 Differential Entropy
  - e) the Cauchy distribution,

$$f(x) = \frac{1}{\pi(1+x^2)}$$

for  $-\infty < x < \infty$ .

Identify the corresponding set of constraints for each of these distributions.
10. Let μ be the mean of a continuous random variable X defined on R<sup>+</sup>. Obtain an upper bound on h(X) in terms of μ.

- 11. The inequality in (10.194) gives an upper bound on the differential entropy in terms of the variance. Can you give an upper bound on the variance in terms of the differential entropy?
- 12. For i = 1, 2, suppose  $f_i$  maximizes h(f) over all the pdf's defined on  $S_i \subset \Re^n$  subject to the constraints in (10.182), where  $S_1 \subset S_2$ . Show that  $h(f_1) \leq h(f_2)$ .
- 13. Hadamard's inequality Show that for a positive semidefinite matrix K,  $|K| \leq \prod_{i=1}^{n} K_{ii}$ , with equality if and only if K is diagonal. Hint: Consider the differential entropy of a multivariate Gaussian distribution.
- 14. Let  $K_{\mathbf{X}}$  and  $\tilde{K}_{\mathbf{X}}$  be the covariance matrix and the correlation matrix of a random vector  $\mathbf{X}$ , respectively. Show that  $|K_{\mathbf{X}}| \leq |\tilde{K}_{\mathbf{X}}|$ . This is a generalization of var $X \leq EX^2$  for a random variable X. Hint: Consider a multivariate Gaussian distribution with another multivariate Gaussian distribution with zero mean and the same correlation matrix.

# **Historical Notes**

The concept of differential entropy was introduced by Shannon [291]. Informational divergence and mutual information were subsequently defined in Kolmogorov [184] and Pinsker [263] in the general setting of measure theory. A measure theoretic treatment of information theory for continuous systems can be found in the book by Ihara [161].

The treatment in this chapter and the next chapter aims to keep the generality of the results without resorting to heavy use of measure theory. The bounds in Section 10.6 for differential entropy subject to constraints are developed in the spirit of maximum entropy expounded in Jayes [167].

In Chapter 7, we have studied the discrete memoryless channel. For such a channel, transmission is in discrete time, and the input and output are discrete. In a physical communication system, the input and output of a channel often take continuous real values. If transmission is in continuous time, the channel is called a *waveform channel*.

In this chapter, we first discuss discrete-time channels with real input and output. We will then extend our discussion to waveform channels. All the logarithms in this chapter are in the base 2.

## **11.1 Discrete-Time Channels**

**Definition 11.1.** Let f(y|x) be a conditional pdf defined for all x, where f(y|x) is bounded (cf. Definition 10.25) and

$$-\int_{\mathcal{S}_Y(x)} f(y|x) \log f(y|x) dy \tag{11.1}$$

is uniformly bounded for all x. A discrete-time continuous channel f(y|x) is a system with input random variable X and output random variable Y such that Y is related to X through f(y|x) (cf. Definition 10.22).

**Remark** The integral in (11.1) is precisely the conditional differential entropy h(Y|X = x) defined in (10.98), which is required to be uniformly bounded in this definition of a discrete-time continuous channel.

**Definition 11.2.** Let  $\alpha : \Re \times \Re \to \Re$ , and Z be a real random variable, called the noise variable. A discrete-time continuous channel  $(\alpha, Z)$  is a system with a real input and a real output. For any input random variable X, the noise random variable Z is independent of X, and the output random variable Y is given by

$$Y = \alpha(X, Z). \tag{11.2}$$

For brevity, a discrete-time continuous channel will be referred to as a continuous channel.

**Definition 11.3.** Two continuous channels f(y|x) and  $(\alpha, Z)$  are equivalent if for every input distribution F(x),

$$\Pr\{\alpha(X,Z) \le y, X \le x\} = \int_{-\infty}^{x} \int_{-\infty}^{y} f_{Y|X}(v|u) dv \, dF_X(u)$$
(11.3)

for all x and y.

**Remark** In the above definitions, the input random variable X is not necessarily continuous.

Definitions 11.1 and 11.2 are two definitions for a continuous channel which are analogous to Definitions 7.1 and 7.2 for a discrete channel. While Definitions 7.1 and 7.2 are equivalent, Definition 11.2 is more general than Definition 11.1. For a continuous channel defined in Definition 11.2, the noise random variable Z may not have a pdf, and the function  $\alpha(x, \cdot)$  may be manyto-one. As a result, the corresponding conditional pdf f(y|x) as required in Definition 11.1 may not exist. In this chapter, we confine our discussion to continuous channels that can be defined by Definition 11.1 (and hence also by Definition 11.2).

**Definition 11.4.** A continuous memoryless channel (CMC) f(y|x) is a sequence of replicates of a generic continuous channel f(y|x). These continuous channels are indexed by a discrete-time index i, where  $i \ge 1$ , with the ith channel being available for transmission at time i. Transmission through a channel is assumed to be instantaneous. Let  $X_i$  and  $Y_i$  be respectively the input and the output of the CMC at time i, and let  $T_{i-}$  denote all the random variables that are generated in the system before  $X_i$ . The Markov chain  $T_{i-} \to X_i \to Y_i$ holds, and

$$\Pr\{Y_i \le y, X_i \le x\} = \int_{-\infty}^x \int_{-\infty}^y f_{Y|X}(v|u) dv \, dF_X(u).$$
(11.4)

**Definition 11.5.** A continuous memoryless channel  $(\alpha, Z)$  is a sequence of replicates of a generic continuous channel  $(\alpha, Z)$ . These continuous channels are indexed by a discrete-time index i, where  $i \ge 1$ , with the *i*th channel being available for transmission at time i. Transmission through a channel is assumed to be instantaneous. Let  $X_i$  and  $Y_i$  be respectively the input and the output of the CMC at time i, and let  $T_{i-}$  denote all the random variables that are generated in the system before  $X_i$ . The noise variable  $Z_i$  for the transmission at time i is a copy of the generic noise variable Z, and it is independent of  $(X_i, T_{i-})$ . The output of the CMC at time i is given by

$$Y_i = \alpha(X_i, Z_i). \tag{11.5}$$

**Definition 11.6.** Let  $\kappa$  be a real function. An average input constraint  $(\kappa, P)$  for a CMC is the requirement that for any codeword  $(x_1, x_2, \dots, x_n)$  transmitted over the channel,

$$\frac{1}{n}\sum_{i=1}^{n}\kappa(x_i) \le P. \tag{11.6}$$

For brevity, an average input constraint is referred to as an input constraint.

**Definition 11.7.** The capacity of a continuous memoryless channel f(y|x) with input constraint  $(\kappa, P)$  is defined as

$$C(P) = \sup_{F(x): E\kappa(X) \le P} I(X;Y), \tag{11.7}$$

where X and Y are respectively the input and output of the generic continuous channel, and F(x) is the distribution of X.

**Theorem 11.8.** C(P) is non-decreasing, concave, and left-continuous.

*Proof.* In the definition of C(P), the supremum is taken over a larger set for a larger P. Therefore, C(P) is non-decreasing in P.

We now show that C(P) is concave. Let i = 1, 2. For an input distribution  $F_j(x)$ , denote the corresponding input and output random variables by  $X_j$  and  $Y_j$ , respectively. Then for any  $P_j$ , for all  $\epsilon > 0$ , there exists  $F_j(x)$  such that

$$E\kappa(X_j) \le P_j \tag{11.8}$$

and

$$I(X_j; Y_j) \ge C(P_j) - \epsilon. \tag{11.9}$$

For  $0 \leq \lambda \leq 1$ , let  $\bar{\lambda} = 1 - \lambda$  and define the random variable

$$X^{(\lambda)} \sim \lambda F_1(x) + \bar{\lambda} F_2(x). \tag{11.10}$$

Then

$$E\kappa(X^{(\lambda)}) = \lambda E\kappa(X_1) + \bar{\lambda} E\kappa(X_2) \le \lambda P_1 + \bar{\lambda} P_2, \qquad (11.11)$$

By the concavity of mutual information with respect to the input distribution<sup>1</sup>, we have

$$I(X^{(\lambda)}; Y^{(\lambda)}) \ge \lambda I(X_1; Y_1) + \bar{\lambda} I(X_2; Y_2)$$
(11.12)

$$\geq \lambda(C(P_1) - \epsilon) + \bar{\lambda}(C(P_2) - \epsilon) \tag{11.13}$$

$$=\lambda C(P_1) + \bar{\lambda}C(P_2) - \epsilon. \tag{11.14}$$

Then

<sup>&</sup>lt;sup>1</sup> Specifically, we refer to the inequality (3.121) in Example 3.14 with X and Y being real random variables related by a conditional pdf f(y|x). The proof of this inequality is left as an exercise.

$$C(\lambda P_1 + \bar{\lambda} P_2) \ge I(X^{(\lambda)}; Y^{(\lambda)}) \ge \lambda C(P_1) + \bar{\lambda} C(P_2) - \epsilon.$$
(11.15)

Letting  $\epsilon \to 0$ , we have

$$C(\lambda P_1 + \bar{\lambda} P_2) \ge \lambda C(P_1) + \bar{\lambda} C(P_2), \qquad (11.16)$$

proving that C(P) is concave.

Finally, we prove that C(P) is left-continuous. Let  $P_1 < P_2$  in (11.16). Since C(P) is non-decreasing, we have

$$C(P_2) \ge C(\lambda P_1 + \bar{\lambda} P_2) \ge \lambda C(P_1) + \bar{\lambda} C(P_2).$$
(11.17)

Letting  $\lambda \to 0$ , we have

$$C(P_2) \ge \lim_{\lambda \to 0} C(\lambda P_1 + \bar{\lambda} P_2) \ge C(P_2), \tag{11.18}$$

which implies

$$\lim_{\lambda \to 0} C(\lambda P_1 + \bar{\lambda} P_2) = C(P_2).$$
(11.19)

Hence, we conclude that

$$\lim_{P \uparrow P_2} C(P) = C(P_2), \tag{11.20}$$

i.e., C(P) is left-continuous. The theorem is proved.  $\Box$ 

# 11.2 The Channel Coding Theorem

**Definition 11.9.** An (n, M) code for a continuous memoryless channel with input constraint  $(\kappa, P)$  is defined by an encoding function

$$e: \{1, 2, \cdots, M\} \to \Re^n \tag{11.21}$$

 $and \ a \ decoding \ function$ 

$$g: \Re^n \to \{1, 2, \cdots, M\}.$$
 (11.22)

The set  $\{1, 2, \dots, M\}$ , denoted by  $\mathcal{W}$ , is called the message set. The sequences  $e(1), e(2), \dots, e(M)$  in  $\mathcal{X}^n$  are called codewords, and the set of codewords is called the codebook. Moreover,

$$\frac{1}{n}\sum_{i=1}^{n}\kappa(x_i(w)) \le P \quad for \ 1 \le w \le M,$$
(11.23)

where  $e(w) = (x_1(w), x_2(w), \dots, x_n(w)).$ 

We assume that a message W is randomly chosen from the message set  $\mathcal{W}$  according to the uniform distribution. Therefore,

$$H(W) = \log M. \tag{11.24}$$

With respect to a channel code for a given CMC, we let

٦

$$\mathbf{X} = (X_1, X_2, \cdots, X_n) \tag{11.25}$$

and

$$\mathbf{Y} = (Y_1, Y_2, \cdots, Y_n)$$
 (11.26)

be the input sequence and the output sequence of the channel, respectively. Evidently,

$$\mathbf{X} = e(W). \tag{11.27}$$

We also let

$$\hat{W} = g(\mathbf{Y}) \tag{11.28}$$

be the estimate on the message W by the decoder.

**Definition 11.10.** For all  $1 \le w \le M$ , let

$$\lambda_w = \Pr\{\hat{W} \neq w | W = w\} = \sum_{\mathbf{y} \in \mathcal{Y}^n: g(\mathbf{y}) \neq w} \Pr\{\mathbf{Y} = \mathbf{y} | \mathbf{X} = e(w)\}$$
(11.29)

be the conditional probability of error given that the message is w.

We now define two performance measures for a channel code.

**Definition 11.11.** The maximal probability of error of an (n, M) code is defined as

$$\lambda_{max} = \max_{w} \lambda_{w}. \tag{11.30}$$

**Definition 11.12.** The average probability of error of an (n, M) code is defined as

$$P_e = \Pr\{\tilde{W} \neq W\}. \tag{11.31}$$

Evidently,  $P_e \leq \lambda_{max}$ .

**Definition 11.13.** A rate R is asymptotically achievable for a continuous memoryless channel if for any  $\epsilon > 0$ , there exists for sufficiently large n an (n, M) code such that

$$\frac{1}{n}\log M > R - \epsilon \tag{11.32}$$

and

$$\lambda_{max} < \epsilon. \tag{11.33}$$

For brevity, an asymptotically achievable rate will be referred to as an achievable rate.

**Theorem 11.14 (Channel Coding Theorem).** A rate R is achievable for a continuous memoryless channel if and only if  $R \leq C$ , the capacity of the channel.

# 11.3 Proof of the Channel Coding Theorem

## 11.3.1 The Converse

We can establish the Markov chain

$$W \to \mathbf{X} \to \mathbf{Y} \to \hat{W}$$
 (11.34)

very much like the discrete case as discussed in Section 7.3. Here, although  $\mathbf{X}$  is a real random vector, it takes only discrete values as it is a function of the message W which is discrete. The only continuous random variable in the above Markov chain is the random vector  $\mathbf{Y}$ , which needs to be handled with caution. The following lemma is essentially the data processing theorem we proved in Theorem 2.42 except that  $\mathbf{Y}$  is continuous. The reader may skip the proof at the first reading.

#### Lemma 11.15.

$$I(W; \hat{W}) \le I(\mathbf{X}; \mathbf{Y}). \tag{11.35}$$

*Proof.* We first consider

$$I(W; \hat{W}) \le I(W, \mathbf{X}; \hat{W}) \tag{11.36}$$

$$= I(\mathbf{X}; \hat{W}) + I(W; \hat{W} | \mathbf{X})$$
(11.37)

$$=I(\mathbf{X};W). \tag{11.38}$$

Note that all the random variables above are discrete. Continuing from the above, we have

$$I(W; \hat{W}) \le I(\mathbf{X}; \hat{W}) \tag{11.39}$$

$$\leq I(\mathbf{X}; \hat{W}) + I(\mathbf{X}; \mathbf{Y} | \hat{W}) \tag{11.40}$$

$$= E \log \frac{p(\mathbf{X}, \hat{W})}{p(\mathbf{X})p(\hat{W})} + E \log \frac{f(\mathbf{Y}|\mathbf{X}, \hat{W})}{f(\mathbf{Y}|\hat{W})}$$
(11.41)

$$= E \log \frac{p(\mathbf{X}, \hat{W}) f(\mathbf{Y} | \mathbf{X}, \hat{W})}{p(\mathbf{X}) [p(\hat{W}) f(\mathbf{Y} | \hat{W})]}$$
(11.42)

$$= E \log \frac{f(\mathbf{Y})p(\mathbf{X}, \hat{W}|\mathbf{Y})}{p(\mathbf{X})[f(\mathbf{Y})p(\hat{W}|\mathbf{Y})]}$$
(11.43)

$$= E \log \frac{p(\mathbf{X}, \hat{W} | \mathbf{Y})}{p(\mathbf{X})p(\hat{W} | \mathbf{Y})}$$
(11.44)

$$= E \log \frac{p(\mathbf{X}|\mathbf{Y})p(W|\mathbf{X},\mathbf{Y})}{p(\mathbf{X})p(\hat{W}|\mathbf{Y})}$$
(11.45)

$$= E \log \frac{p(\mathbf{X}|\mathbf{Y})}{p(\mathbf{X})} + E \log \frac{p(\hat{W}|\mathbf{X},\mathbf{Y})}{p(\hat{W}|\mathbf{Y})}$$
(11.46)

11.3 Proof of the Channel Coding Theorem 257

$$= E \log \frac{f(\mathbf{Y}|\mathbf{X})}{f(\mathbf{Y})} + E \log \frac{p(\mathbf{X}|\mathbf{Y},\hat{W})}{p(\mathbf{X}|\mathbf{Y})}$$
(11.47)

$$= I(\mathbf{X}; \mathbf{Y}) + E \log \frac{p(\mathbf{X}|\mathbf{Y})}{p(\mathbf{X}|\mathbf{Y})}$$
(11.48)

$$= I(\mathbf{X}; \mathbf{Y}) + E \log 1 \tag{11.49}$$

$$=I(\mathbf{X};\mathbf{Y})+0\tag{11.50}$$

$$=I(\mathbf{X};\mathbf{Y}).$$
(11.51)

The above steps are justified as follows:

• The relation

$$f(\mathbf{y}|\mathbf{x}) = \prod_{i=1}^{n} f(y_i|x_i)$$
(11.52)

can be established in exactly the same way as we established (7.101) for the discrete case (when the channel is used without feedback). By our assumption that f(y|x) is bounded,  $f(\mathbf{y}|\mathbf{x})$  is also bounded. Moreover,

$$f(\mathbf{y}|\mathbf{x}, \hat{w}) = \frac{p(\mathbf{x})f(\mathbf{y}|\mathbf{x})p(\hat{w}|\mathbf{y})}{p(\mathbf{x}, \hat{w})}$$
(11.53)

is bounded because  $f(\mathbf{y}|\mathbf{x})$  is bounded. By Proposition 10.26,  $f(\mathbf{y}|\hat{w})$  exists and is bounded. Therefore,  $I(\mathbf{X}; \mathbf{Y}|\hat{W})$  can be defined as in Definition 10.29. The inequality (11.40) follows from Corollary 10.34.

- In (11.43), since  $f(\mathbf{y}|\mathbf{x})$  is bounded, it follows from Proposition 10.26 that  $f(\mathbf{y})$  exists and is bounded.
- (11.47) follows from

$$p(\mathbf{x})f(\mathbf{y}|\mathbf{x}) = f(\mathbf{y})p(\mathbf{x}|\mathbf{y})$$
(11.54)

and

$$p(\mathbf{x}|\mathbf{y})p(\hat{w}|\mathbf{x},\mathbf{y}) = p(\hat{w}|\mathbf{y})p(\mathbf{x}|\mathbf{y},\hat{w}).$$
(11.55)

• (11.48) follows from the Markov chain  $\mathbf{X} \to \mathbf{Y} \to \hat{W}$ .

The proof is accomplished.  $\hfill\square$ 

We now proceed to prove the converse. Let R be an achievable rate, i.e., for any  $\epsilon > 0$ , there exists for sufficiently large n and (n, M) code such that

$$\frac{1}{n}\log M > R - \epsilon \tag{11.56}$$

and

$$\lambda_{max} < \epsilon. \tag{11.57}$$

Consider

$$\log M = H(W) \tag{11.58}$$

$$= H(W|\hat{W}) + I(W;\hat{W})$$
(11.59)

$$\leq H(W|\hat{W}) + I(\mathbf{X}; \mathbf{Y}) \tag{11.60}$$

$$\leq H(W|\hat{W}) + h(\mathbf{Y}) - h(\mathbf{Y}|\mathbf{X}) \tag{11.61}$$

$$\leq H(W|\hat{W}) + \sum_{i=1}^{n} h(Y_i) - h(\mathbf{Y}|\mathbf{X})$$
(11.62)

$$= H(W|\hat{W}) + \sum_{i=1}^{n} h(Y_i) - \sum_{i=1}^{n} h(Y_i|X_i)$$
(11.63)

$$= H(W|\hat{W}) + \sum_{i=1}^{n} I(X_i; Y_i).$$
(11.64)

The above steps are justified as follows:

- (11.60) follows from Lemma 11.15.
- It follows from (11.52) that

$$h(\mathbf{Y}|\mathbf{X}) = \sum_{i=1}^{n} h(Y_i|X_i).$$
 (11.65)

Then (11.1) in Definition 11.1 implies that  $h(Y_i|X_i)$  is finite for all *i*, and hence  $h(\mathbf{Y}|\mathbf{X})$  is also finite.

- From the foregoing,  $f(\mathbf{y})$  exists. Therefore,  $h(\mathbf{Y})$  can be defined according to Definition 10.10 (but  $h(\mathbf{Y})$  may be infinite), and (11.61) follows from Proposition 10.31 because  $h(\mathbf{Y}|\mathbf{X})$  is finite. Note that it is necessary to require  $h(\mathbf{Y}|\mathbf{X})$  to be finite because otherwise  $h(\mathbf{Y})$  is also infinite and Proposition 10.31 cannot be applied.
- (11.62) follows from Corollary 10.36, the independence bound for differential entropy.
- (11.63) from (11.65) above.
- (11.64) follows from Proposition 10.31.

Let V be a mixing random variable distributed uniformly on  $\{1, 2, \dots, n\}$ which is independent of  $X_i$ ,  $1 \leq i \leq n$ . Let  $X = X_V$  and Y be the output of the channel with X being the input. Then

$$E\kappa(X) = EE[\kappa(X)|V]$$
(11.66)

$$=\sum_{i=1}^{n} \Pr\{V=i\} E[\kappa(X)|V=i]$$
(11.67)

$$= \sum_{i=1}^{n} \Pr\{V = i\} E[\kappa(X_i)|V = i]$$
(11.68)

$$=\sum_{i=1}^{n}\frac{1}{n}E\kappa(X_i) \tag{11.69}$$

11.3 Proof of the Channel Coding Theorem 259

$$= E\left[\frac{1}{n}\sum_{i=1}^{n}\kappa(X_i)\right]$$
(11.70)

$$\leq P$$
, (11.71)

where the above inequality follows from (11.23) in the definition of the code. By the concavity of mutual information with respect to the input distribution, we have

$$\frac{1}{n}\sum_{i=1}^{n}I(X_i;Y_i) \le I(X;Y) \le C,$$
(11.72)

where the last inequality holds in light of the definition of C and (11.71). Then it follows from (11.64) that

$$\log M \le H(W|\hat{W}) + nC,\tag{11.73}$$

which is precisely (7.126) in the proof of the converse of the channel coding theorem for the DMC. Following exactly the same steps therein, we conclude that

$$R \le C. \tag{11.74}$$

#### 11.3.2 Achievability

The proof of the achievability of the channel capacity, which involves the construction of a random code, is somewhat different from the construction for the discrete case in Section 7.4. On the one hand, we need to take into account the input constraint. On the other hand, since the input distribution F(x) we use for constructing the random code may not have a pdf, it is difficult to formulate the notion of joint typicality as in the discrete case. Instead, we will introduce a different notion of typicality based on mutual information.

Consider a bivariate information source  $\{(X_k, Y_k), k \ge 1\}$ , where  $(X_k, Y_k)$  are i.i.d. with (X, Y) being the pair of generic real random variables. The conditional pdf f(y|x) exists in the sense prescribed in Definition 10.22. By Proposition 10.26, with the assumption that f(y|x) is bounded, f(y) exists and is bounded, so that the mutual information I(X;Y) can be defined as in Definition 10.28.

**Definition 11.16.** The mutually typical set  $U^n_{[XY]\delta}$  with respect to F(x, y) is the set of  $(\mathbf{x}, \mathbf{y}) \in \mathcal{X}^n \times \mathcal{Y}^n$  such that

$$\left|\frac{1}{n}\log\frac{f(\mathbf{y}|\mathbf{x})}{f(\mathbf{y})} - I(X;Y)\right| \le \delta,\tag{11.75}$$

where

$$f(\mathbf{y}|\mathbf{x}) = \prod_{i=1}^{n} f(y_i|x_i)$$
(11.76)

and

$$f(\mathbf{y}) = \prod_{i=1}^{n} f(y_i), \qquad (11.77)$$

and  $\delta$  is an arbitrarily small positive number. A pair of sequences  $(\mathbf{x}, \mathbf{y})$  is called mutually  $\delta$ -typical if it is in  $U_{[XY]\delta}^n$ .

**Lemma 11.17.** For any  $\delta > 0$ , for sufficiently large n,

$$\Pr\{(\mathbf{X}, \mathbf{Y}) \in U^n_{[XY]\delta}\} \ge 1 - \delta.$$
(11.78)

*Proof.* By (11.76) and (11.77), we write

$$\frac{1}{n}\log\frac{f(\mathbf{Y}|\mathbf{X})}{f(\mathbf{Y})} = \frac{1}{n}\log\prod_{i=1}^{n}\frac{f(Y_i|X_i)}{f(Y_i)} = \frac{1}{n}\sum_{i=1}^{n}\log\frac{f(Y_i|X_i)}{f(Y_i)}.$$
(11.79)

Since  $(X_i, Y_i)$  are i.i.d., so are the random variables  $\log \frac{f(Y_i|X_i)}{f(Y_i)}$ . Thus we conclude by the weak law of large numbers that

$$\frac{1}{n} \sum_{i=1}^{n} \log \frac{f(Y_i|X_i)}{f(Y_i)} \to E \log \frac{f(Y|X)}{f(Y)} = I(X;Y)$$
(11.80)

in probability, i.e., (11.78) holds for sufficiently large n, proving the lemma.

The following lemma is analogous to Lemma 7.17 for the discrete case.

**Lemma 11.18.** Let  $(\mathbf{X}', \mathbf{Y}')$  be *n* i.i.d. copies of a pair of generic random variables (X', Y'), where X' and Y' are independent and have the same marginal distributions as X and Y, respectively. Then

$$\Pr\{(\mathbf{X}', \mathbf{Y}') \in U^n_{[XY]\delta}\} \le 2^{-n(I(X;Y)-\delta)}.$$
(11.81)

*Proof.* For  $(\mathbf{x}, \mathbf{y}) \in U^n_{[XY]\delta}$ , from (11.75), we obtain

$$\frac{1}{n}\log\frac{f(\mathbf{y}|\mathbf{x})}{f(\mathbf{y})} \ge I(X;Y) - \delta,$$
(11.82)

or

$$f(\mathbf{y}|\mathbf{x}) \ge f(\mathbf{y})2^{n(I(X;Y)-\delta)}$$
(11.83)

Then

$$1 \ge \Pr\{(\mathbf{X}, \mathbf{Y}) \in U^n_{[XY]\delta})\}$$
(11.84)

$$= \int \int_{U_{[XY]\delta}^n} f(\mathbf{y}|\mathbf{x}) dF(\mathbf{x})$$
(11.85)

$$\geq 2^{n(I(X;Y)-\delta)} \int \int_{U_{[XY]\delta}^n} f(\mathbf{y}) dF(\mathbf{x})$$
(11.86)

$$=2^{n(I(X;Y)-\delta)}\Pr\{(\mathbf{X}',\mathbf{Y}')\in U^{n}_{[XY]\delta}\},$$
(11.87)

where the last inequality follows from (11.83). Therefore,

$$\Pr\{(\mathbf{X}', \mathbf{Y}') \in U^n_{[XY]\delta}\} \le 2^{-n(I(X;Y)-\delta)},$$
(11.88)

proving the lemma.  $\Box$ 

Fix any  $\epsilon > 0$  and let  $\delta$  be a small quantity to be specified later. Since C(P) is left-continuous, there exists a sufficiently small  $\gamma > 0$  such that

$$C(P-\gamma) > C(P) - \frac{\epsilon}{6}.$$
(11.89)

By the definition of  $C(P - \gamma)$ , there exists an input random variable X such that

$$E\kappa(X) \le P - \gamma \tag{11.90}$$

and

$$I(X;Y) \ge C(P-\gamma) - \frac{\epsilon}{6}.$$
(11.91)

Then choose for a sufficiently large n an even integer M satisfying

$$I(X;Y) - \frac{\epsilon}{6} < \frac{1}{n} \log M < I(X;Y) - \frac{\epsilon}{8},$$
 (11.92)

from which we obtain

$$\frac{1}{n}\log M > I(X;Y) - \frac{\epsilon}{6} \tag{11.93}$$

$$\geq C(P-\gamma) - \frac{\epsilon}{3} \tag{11.94}$$

$$> C(P) - \frac{\epsilon}{2}.\tag{11.95}$$

We now describe a random coding scheme:

- 1. Construct the codebook C of an (n, M) code randomly by generating M codewords in  $\Re^n$  independently and identically according to  $f(x)^n$ . Denote these codewords by  $\tilde{\mathbf{X}}(1), \tilde{\mathbf{X}}(2), \dots, \tilde{\mathbf{X}}(M)$ .
- 2. Reveal the codebook  $\mathcal{C}$  to both the encoder and the decoder.
- 3. A message W is chosen from W according to the uniform distribution.
- 4. The sequence  $\mathbf{X} = \mathbf{\tilde{X}}(W)$ , namely the Wth codeword in the codebook C, is transmitted through the channel.
- 5. The channel outputs a sequence **Y** according to

$$\Pr\{Y_i \le y_i, 1 \le i \le n | \mathbf{X}(W) = \mathbf{x}\} = \prod_{i=1}^n \int_{-\infty}^{y_i} f(y|x_i) dy.$$
(11.96)

This is the continuous analog of (7.101) and can be established similarly.

6. The sequence  $\mathbf{Y}$  is decoded to the message w if  $(\mathbf{X}(w), \mathbf{Y}) \in U^n_{[XY]\delta}$  and there does not exist  $w' \neq w$  such that  $(\mathbf{X}(w'), \mathbf{Y}) \in U^n_{[XY]\delta}$ . Otherwise,  $\mathbf{Y}$  is decoded to a constant message in  $\mathcal{W}$ . Denote by  $\hat{W}$  the message to which  $\mathbf{Y}$  is decoded.

We now analyze the performance of this random coding scheme. Let

$$\tilde{\mathbf{X}}(w) = (\tilde{X}_1(w), \tilde{X}_2(w), \cdots, \tilde{X}_n(w))$$
(11.97)

and define the error event

$$Err = E_e \cup E_d, \tag{11.98}$$

where

$$E_e = \left\{ \frac{1}{n} \sum_{i=1}^n \kappa(\tilde{X}_i(W)) > P \right\}$$
(11.99)

is the event that the input constraint is violated, and

$$E_d = \{ \hat{W} \neq W \}, \tag{11.100}$$

is the event that a decoding error occurs. By symmetry in the code construction,

$$\Pr\{Err\} = \Pr\{Err|W = 1\}$$
(11.101)

$$\leq \Pr\{E_e|W=1\} + \Pr\{E_d|W=1\}.$$
(11.102)

With Lemma 11.18 in place of Lemma 7.17, the analysis of  $Pr\{E_d|W=1\}$  is exactly the same as the analysis of the decoding error in the discrete case. The details are omitted, and we conclude that by choosing  $\delta$  to be a sufficiently small positive quantity,

$$\Pr\{E_d|W=1\} \le \frac{\epsilon}{4} \tag{11.103}$$

for sufficiently large n.

We now analyze  $\Pr\{E_e|W=1\}$ . By the weak law of large numbers,

$$\Pr\{E_e | W = 1\} = \Pr\left\{ \left. \frac{1}{n} \sum_{i=1}^n \kappa(\tilde{X}_i(1)) > P \right| W = 1 \right\}$$
(11.104)

$$= \Pr\left\{\frac{1}{n}\sum_{i=1}^{n}\kappa(\tilde{X}_{i}(1)) > P\right\}$$
(11.105)

$$= \Pr\left\{\frac{1}{n}\sum_{i=1}^{n}\kappa(\tilde{X}_{i}(1)) > (P-\nu) + \nu\right\}$$
(11.106)

$$\leq \Pr\left\{\frac{1}{n}\sum_{i=1}^{n}\kappa(\tilde{X}_{i}(1)) > E\kappa(X) + \nu\right\}$$
(11.107)

$$\leq \frac{\epsilon}{4} \tag{11.108}$$

for sufficiently large n. It then follows from (11.102) and (11.103) that

$$\Pr\{Err\} \le \frac{\epsilon}{2} \tag{11.109}$$

for sufficiently large n.

It remains to show the existence of a codebook such that  $\lambda_{max} < \epsilon$  and the input constraint (11.23) is satisfied by every codeword. Now consider

$$\Pr\{Err\} = \sum_{\mathcal{C}} \Pr\{\mathcal{C}\} \Pr\{Err|\mathcal{C}\}, \qquad (11.110)$$

where  $\Pr{\mathcal{C}}$  is the probability of choosing a codebook  $\mathcal{C}$  from the ensemble of all possible codebooks in Step 1 of the random coding scheme. In light of (11.109), there exists at least one codebook  $\mathcal{C}^*$  such that

$$\Pr\{Err|\mathcal{C}^*\} \le \frac{\epsilon}{2}.$$
(11.111)

Furthermore,

$$\Pr\{Err|\mathcal{C}^*\} = \sum_{w=1}^{M} \Pr\{W = w|\mathcal{C}^*\} \Pr\{Err|\mathcal{C}^*, W = w\}$$
(11.112)

$$= \sum_{w=1}^{M} \Pr\{W = w\} \Pr\{Err | \mathcal{C}^*, W = w\}$$
(11.113)

$$= \frac{1}{M} \sum_{w=1}^{M} \Pr\{Err | \mathcal{C}^*, W = w\}.$$
 (11.114)

By discarding the worst half of the codewords in  $\mathcal{C}^*$ , if a codeword  $\tilde{\mathbf{X}}(w)$  remains in  $\mathcal{C}^*$ , then

$$\Pr\{Err|\mathcal{C}^*, W = w\} \le \epsilon. \tag{11.115}$$

Since  $Err = E_e \cup E_d$ , this implies

$$\Pr\{E_e | \mathcal{C}^*, W = w\} \le \epsilon \tag{11.116}$$

and

$$\Pr\{E_d | \mathcal{C}^*, W = w\} \le \epsilon, \tag{11.117}$$

where the latter implies  $\lambda_{max} \leq \epsilon$  for the codebook  $\mathcal{C}^*$ . Finally, observe that conditioning on  $\{\mathcal{C}^*, W = w\}$ , the codeword  $\tilde{\mathbf{X}}(w)$  is deterministic. Therefore,  $\Pr\{E_e|\mathcal{C}^*, W = w\}$  is equal to 1 if the codeword  $\tilde{\mathbf{X}}(w)$  violates the input constraint (11.23), otherwise it is equal to 0. Then (11.116) implies that for every codeword  $\tilde{\mathbf{X}}(w)$  that remains in  $\mathcal{C}^*$ ,  $\Pr\{E_e|\mathcal{C}^*, W = w\} = 0$ , i.e., the input constraint is satisfied. This completes the proof.

## **11.4 Memoryless Gaussian Channels**

In communication engineering, the Gaussian channel is the most commonly used model for a noisy channel with real input and output. The reasons are two-fold. First, the Gaussian channel is highly analytically tractable. Second, the Gaussian noise can be regarded as the worst kind of additive noise subject to a constraint on the noise power. This will be discussed in Section 11.9.

We first give two equivalent definitions of a Gaussian channel.

**Definition 11.19 (Gaussian Channel).** A Gaussian channel with noise energy N is a continuous channel with the following two equivalent specifications:

1. 
$$f(y|x) = \frac{1}{\sqrt{2\pi N}} e^{-\frac{(y-x)^2}{2N}}$$
.  
2.  $Z \sim \mathcal{N}(0, N)$  and  $\alpha(X, Z) = X + Z$ 

**Definition 11.20 (Memoryless Gaussian Channel).** A memoryless Gaussian channel with noise power N and input power constraint P is a memoryless continuous channel with the generic continuous channel being the Gaussian channel with noise energy N. The input power constraint P refers to the input constraint ( $\kappa$ , P) with  $\kappa(x) = x^2$ .

Using the formula in Definition 11.7 for the capacity of a CMC, the capacity of a Gaussian channel can be evaluated.

**Theorem 11.21 (Capacity of a Memoryless Gaussian Channel).** The capacity of a memoryless Gaussian channel with noise power N and input power constraint P is

$$\frac{1}{2}\log(1+\frac{P}{N}).$$
(11.118)

The capacity is achieved by the input distribution  $\mathcal{N}(0, P)$ .

We first prove the following lemma.

**Lemma 11.22.** Let Y = X + Z. Then h(Y|X) = h(Z|X) provided that  $f_{Z|X}(z|x)$  exists for all  $x \in S_X$ .

*Proof.* For all  $x \in S_X$ , since  $f_{Z|X}(z|x)$  exists,  $f_{Y|X}(y|x)$  also exists and is given by

$$f_{Y|X}(y|x) = f_{Z|X}(y - x|x).$$
(11.119)

Then

$$h(Y|X) = \int h(Y|X=x)dF_X(x)$$
(11.120)

$$= \int h(X+Z|X=x)dF_X(x)$$
 (11.121)

$$= \int h(x+Z|X=x)dF_X(x)$$
 (11.122)

$$= \int h(Z|X=x)dF_X(x) \tag{11.123}$$

$$=h(Z|X) \tag{11.124}$$

In the above, (11.120) and (11.124) follow from (10.99), while (11.123) follows from the translation property of differential entropy (Theorem 10.14).  $\Box$ 

**Remark** Since Y and Z uniquely determine each other given X, it is tempting to write h(Y|X) = h(Z|X) immediately. However, this interpretation is incorrect because differential entropy is not the same as entropy.

Proof of Theorem 11.21. Let F(x) be the CDF of the input random variable X such that  $EX^2 \leq P$ , where X is not necessarily continuous. Since  $Z \sim \mathcal{N}(0, N), f(y|x)$ , as given by (11.119), exists and is bounded. By Proposition 10.26, f(y) exists and hence h(Y) is defined. Since Z is independent of X, by Lemma 11.22,

$$h(Y|X) = h(Z|X) = h(Z).$$
 (11.125)

Then

$$I(X;Y) = h(Y) - h(Y|X)$$
(11.126)

$$= h(Y) - h(Z), \tag{11.127}$$

where (11.126) follows from Proposition 10.31 and (11.127) follows from (11.125).

Since Y = X + Z and Z is independent of X, we have

$$EY^2 = E(X+Z)^2 (11.128)$$

$$= EX^{2} + 2(EX)(EZ) + EZ^{2}$$
(11.129)

$$= EX^{2} + 2(EX)(0) + EZ^{2}$$
(11.130)  
=  $EX^{2} + EZ^{2}$ (11.131)

$$= EX^2 + EZ^2 \tag{11.131}$$

$$\leq P + N. \tag{11.132}$$

Given the above energy constraint on Y, by Theorem 10.45, we have

$$h(Y) \le \frac{1}{2} \log[2\pi e(P+N)],$$
 (11.133)

with equality if  $Y \sim \mathcal{N}(0, P + N)$ .

Recall from Example 10.13 that

$$h(Z) = \frac{1}{2}\log(2\pi eN).$$
(11.134)

It then follows from (11.127), (11.133), and (11.134) that

$$I(X;Y) = h(Y) - h(Z)$$
(11.135)

$$\leq \frac{1}{2}\log[2\pi e(P+N)] - \frac{1}{2}\log(2\pi eN)$$
(11.136)

$$=\frac{1}{2}\log\left(1+\frac{P}{N}\right).\tag{11.137}$$

Evidently, this upper bound is tight if  $X \sim \mathcal{N}(0, P)$ , because then

$$Y = X + Z \sim \mathcal{N}(0, P + N).$$
(11.138)

Therefore,

$$C = \sup_{F(x): EX^2 \le P} I(X;Y)$$
(11.139)

$$= \max_{F(x): EX^2 \le P} I(X;Y)$$
(11.140)

$$=\frac{1}{2}\log\left(1+\frac{P}{N}\right).\tag{11.141}$$

The theorem is proved.  $\Box$ 

Theorem 11.21 says that the capacity of a memoryless Gaussian channel depends only on the ratio of the input power constraint P to the noise power N. This important quantity is called the *signal-to-noise ratio*. Note that no matter how small the signal-to-noise ratio is, the capacity is still strictly positive. That is, reliable communication can be achieved, though at a low rate, even when the noise power is much larger than the signal power. We also see that the capacity is infinite if there is no constraint on the input power.

# **11.5** Parallel Gaussian Channels

In Section 11.4, we have discussed the capacity of a memoryless Gaussian channel. Now suppose k such channels are available for communication, where  $k \geq 1$ . This is illustrated in Figure 11.1, with  $X_i$ ,  $Y_i$ , and  $Z_i$  being the input, the output, and the noise variable of the *i*th channel, respectively, where  $Z_i \sim \mathcal{N}(0, N_i)$  and  $Z_i$ ,  $1 \leq i \leq k$  are independent.

We are interested in the capacity of such a system of parallel Gaussian channels, with the total input power constraint

$$E\sum_{i=1}^{k} X_i^2 \le P.$$
(11.142)

Let  $\mathbf{X} = [X_1 X_2 \cdots X_k], \mathbf{Y} = [Y_1 Y_2 \cdots Y_k], \text{ and } \mathbf{Z} = [Z_1 Z_2 \cdots Z_k].$  Then

$$f_{\mathbf{Y}|\mathbf{X}}(\mathbf{y}|\mathbf{x}) = \prod_{i=1}^{k} f_{Y_i|X_i}(y_i|x_i)$$
(11.143)

$$=\prod_{i=1}^{k} f_{Z_i|X_i}(y_i - x_i|x_i)$$
(11.144)

$$=\prod_{i=1}^{k} f_{Z_i}(y_i - x_i).$$
(11.145)

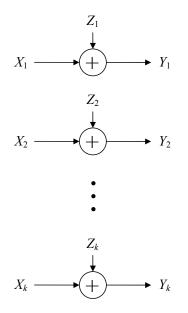


Fig. 11.1. A system of parallel Gaussian channels.

With the existence of  $f(\mathbf{y}|\mathbf{x})$ , by extending Definition 10.23, we have

$$h(\mathbf{Y}|\mathbf{X}) = -\int_{\mathcal{S}_{\mathbf{X}}} \int_{\mathcal{S}_{\mathbf{Y}}(\mathbf{x})} f(\mathbf{y}|\mathbf{x}) \log f(\mathbf{y}|\mathbf{x}) d\mathbf{y} dF(\mathbf{x}).$$
(11.146)

In light of (11.145), since  $f_{Z_i}(z_i)$  is bounded for each *i*, so is  $f_{\mathbf{Y}|\mathbf{X}}(\mathbf{y}|\mathbf{x})$ . Then by Proposition 10.27, we see that  $f(\mathbf{y})$  exists and therefore  $h(\mathbf{Y})$  is defined. By extending Definition 10.28, we have

$$I(\mathbf{X};\mathbf{Y}) = \int_{\mathcal{S}_{\mathbf{X}}} \int_{\mathcal{S}_{\mathbf{Y}}(\mathbf{x})} f(\mathbf{y}|\mathbf{x}) \log \frac{f(\mathbf{y}|\mathbf{x})}{f(\mathbf{y})} d\mathbf{y} dF(\mathbf{x}).$$
(11.147)

It then follows from Definition 11.7 that the capacity of the system is given by

$$C(P) = \sup_{F(\mathbf{x}):E\sum_{i}X_{i}^{2} \leq P} I(\mathbf{X};\mathbf{Y}), \qquad (11.148)$$

where  $F(\mathbf{x})$  is the joint CDF of the input vector  $\mathbf{X}$ . As we will see, the supremum above is indeed a maximum.

When we calculated the capacity of the memoryless Gaussian channel in Theorem 11.21, we obtained in (11.131) that

$$EY^2 = EX^2 + EZ^2, (11.149)$$

i.e., the output energy is equal to the sum of the input energy and the noise energy, provided that the noise has zero mean. By exactly the same argument, we see that

$$EY_i^2 = EX_i^2 + EZ_i^2 (11.150)$$

for all i.

Toward calculating C(P), consider

$$I(\mathbf{X}; \mathbf{Y}) = h(\mathbf{Y}) - h(\mathbf{Y}|\mathbf{X})$$
(11.151)

$$= h(\mathbf{Y}) - h(\mathbf{Z}|\mathbf{X}) \tag{11.152}$$

$$=h(\mathbf{Y})-h(\mathbf{Z}) \tag{11.153}$$

$$= h(\mathbf{Y}) - \sum_{i=1}^{\kappa} h(Z_i)$$
(11.154)

$$= h(\mathbf{Y}) - \frac{1}{2} \sum_{i=1}^{k} \log(2\pi e N_i)$$
(11.155)

$$\leq \sum_{i=1}^{k} h(Y_i) - \frac{1}{2} \sum_{i=1}^{k} \log(2\pi e N_i)$$
(11.156)

$$\leq \frac{1}{2} \sum_{i=1}^{k} \log[2\pi e(EY_i^2)] - \frac{1}{2} \sum_{i=1}^{k} \log(2\pi eN_i)$$
(11.157)

$$= \frac{1}{2} \sum_{i=1}^{k} \log(EY_i^2) - \frac{1}{2} \sum_{i=1}^{k} \log N_i$$
(11.158)

$$= \frac{1}{2} \sum_{i=1}^{k} \log(EX_i^2 + EZ_i^2) - \frac{1}{2} \sum_{i=1}^{k} \log N_i$$
(11.159)

$$= \frac{1}{2} \sum_{i=1}^{k} \log(P_i + N_i) - \frac{1}{2} \sum_{i=1}^{k} \log N_i$$
(11.160)

$$= \frac{1}{2} \sum_{i=1}^{k} \log\left(1 + \frac{P_i}{N_i}\right), \qquad (11.161)$$

where  $P_i = EX_i^2$  is the input power of the *i*th channel. In the above, (11.152) is the vector generalization of Lemma 11.22, (11.154) follows because  $Z_i$  are independent, and (11.159) follows from (11.150).

Equality holds in (11.156) and (11.157) if  $Y_i$ ,  $1 \leq i \leq k$  are independent and  $Y_i \sim \mathcal{N}(0, P_i + N_i)$ . This happens when  $X_i$  are independent of each other and  $X_i \sim \mathcal{N}(0, P_i)$ . Therefore, maximizing  $I(\mathbf{X}; \mathbf{Y})$  becomes maximizing  $\sum_i \log(P_i + N_i)$  in (11.160) with the constraint  $\sum_i P_i \leq P$  and  $P_i \geq 0$  for all *i*. In other words, we are to find the optimal input power allocation among the channels. Comparing (11.161) with (11.141), we see that the capacity of the system of parallel Gaussian channels is equal to the sum of the capacities of the individual Gaussian channels when the input power is optimally allocated.

Toward this end, we first apply the method of Lagrange multipliers by temporarily ignoring the nonnegativity constraints on  $P_i$ . Observe that in

#### 11.5 Parallel Gaussian Channels 269

order for the summation  $\sum_i \log(P_i + N_i)$  in (11.160) to be maximized,  $\sum_i P_i = P$  must hold because  $\log(P_i + N_i)$  is increasing in  $P_i$ . Therefore, the inequality constraint  $\sum_i P_i \leq P$  can be replaced by the equality constraint  $\sum_i P_i = P$ . Let

$$J = \sum_{i=1}^{k} \log \left( P_i + N_i \right) - \mu \sum_{i=1}^{k} P_i.$$
 (11.162)

Differentiating with respect to  $P_i$  gives

$$0 = \frac{\partial J}{\partial P_i} = \frac{1}{P_i + N_i} - \mu, \qquad (11.163)$$

which implies

$$P_i = \frac{1}{\mu} - N_i. \tag{11.164}$$

Upon letting  $\nu = \frac{1}{\mu}$ , we have

$$P_i = \nu - N_i, \tag{11.165}$$

where  $\nu$  is chosen such that

$$\sum_{i=1}^{k} P_i = \sum_{i=1}^{k} (\nu - N_i) = P.$$
(11.166)

However,  $P_i$  as given in (11.165) is not guaranteed to be nonnegative, so it may not be a valid solution. Nevertheless, (11.165) suggests the general solution to be proved in the next proposition.

## Proposition 11.23. The problem

For given 
$$\lambda_i \geq 0$$
, maximize  $\sum_{i=1}^k \log(a_i + \lambda_i)$  subject to  $\sum_i a_i \leq P$   
and  $a_i \geq 0$ 

has the solution

$$a_i^* = (\nu - \lambda_i)^+, \quad 1 \le i \le k,$$
 (11.167)

where

$$(x)^{+} = \begin{cases} x & \text{if } x \ge 0\\ 0 & \text{if } x = 0 \end{cases}$$
(11.168)

and  $\nu$  satisfies

$$\sum_{i=1}^{k} (\nu - \lambda_i)^+ = P.$$
(11.169)

*Proof.* Rewrite the maximization problem as

For given  $\lambda_i \geq 0$ , maximize  $\sum_i \log (a_i + \lambda_i)$  subject to

$$\sum_{i=1}^{k} a_i \le P \tag{11.170}$$

$$-a_i \le 0, \quad 1 \le i \le k.$$
 (11.171)

We will prove the proposition by verifying that the proposed solution in (11.167) satisfies the Karush-Kuhn-Tucker (KKT) condition. This is done by finding nonnegative  $\mu$  and  $\mu_i$  satisfying the equations

$$\frac{1}{2(a_i^* + \lambda_i)} - \mu + \mu_i = 0 \tag{11.172}$$

$$\mu\left(P - \sum_{i=1}^{k} a_i^*\right) = 0 \tag{11.173}$$

$$\mu_i a_i^* = 0, \quad 1 \le i \le k, \tag{11.174}$$

where  $\mu$  and  $\mu_i$  are the multipliers associated with the constraints in (11.170) and (11.171), respectively.

To avoid triviality, assume P > 0 so that  $\nu > 0$ , and observe that there exists at least one *i* such that  $a_i^* > 0$ . For those *i*, (11.174) implies

$$\mu_i = 0. \tag{11.175}$$

On the other hand,

$$a_i^* = (\nu - \lambda_i)^+ = \nu - \lambda_i.$$
 (11.176)

Substituting (11.175) and (11.176) into (11.172), we obtain

$$\mu = \frac{1}{2\nu} > 0. \tag{11.177}$$

For those *i* such that  $a_i^* = 0$ , it follows from (11.167) that  $\nu \leq \lambda_i$ . From (11.177) and (11.172), we obtain

$$\mu_i = \frac{1}{2\nu} - \frac{1}{2\lambda_i} \ge 0. \tag{11.178}$$

Thus we have found nonnegative  $\mu$  and  $\mu_i$  that satisfy (11.172) to (11.174), verifying the KKT condition. The proposition is proved.  $\Box$ 

Hence, following (11.161) and applying the above proposition with  $a_i = P_i$ and  $\lambda_i = N_i$ , we see that the capacity of the system of parallel Gaussian channels is equal to

$$\frac{1}{2} \sum_{i=1}^{k} \log\left(1 + \frac{P_i^*}{N_i}\right), \qquad (11.179)$$

where  $\{P_i^*, 1 \leq i \leq k\}$  is the optimal input power allocation among the channels given by

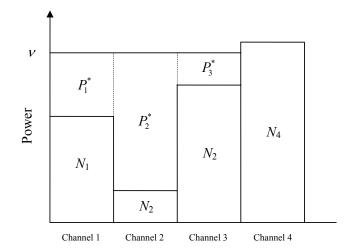


Fig. 11.2. Water-filling for parallel Gaussian channels.

$$P_i^* = (\nu - N_i)^+, \quad 1 \le i \le k \tag{11.180}$$

with  $\nu$  satisfying

$$\sum_{i=1}^{k} (\nu - N_i)^+ = P.$$
(11.181)

The process of obtaining  $\{P_i^*\}$ , called *water-filling*, is illustrated in Figure 11.2. One can image that an amount P of water is poured into a reservoir with an uneven bottom, and  $\nu$  is the level the water rises to. Under this scheme, high input power is allocated to a channel with low noise power. For a channel with noise power higher than  $\nu$ , no input power is allocated, i.e., the channel is not used.

## 11.6 Correlated Gaussian Channels

In this section, we generalize the results in the last section to the case when the noise variables  $Z_i$ ,  $1 \le i \le k$  are correlated with covariance matrix  $K_{\mathbf{Z}}$ . We continue to assume that  $Z_i$  has zero mean for all *i*, i.e.,  $\mathbf{Z} \sim \mathcal{N}(0, K_{\mathbf{Z}})$ , and the total input power constraint

$$E\sum_{i=1}^{k} X_i^2 \le P \tag{11.182}$$

prevails.

We will derive the capacity of such a system of correlated Gaussian channels by decorrelating the noise vector  $\mathbf{Z}$ . Let  $K_{\mathbf{Z}}$  be diagonalizable as  $QAQ^{\top}$ and consider

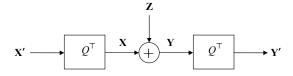


Fig. 11.3. An equivalent system of parallel Gaussian channels.

$$\mathbf{Y} = \mathbf{X} + \mathbf{Z}.\tag{11.183}$$

Then

$$Q^{\top}\mathbf{Y} = Q^{\top}\mathbf{X} + Q^{\top}\mathbf{Z}.$$
 (11.184)

Upon letting

$$\mathbf{X}' = Q^{\top} \mathbf{X}$$
(11.185)  
 
$$\mathbf{Y}' = Q^{\top} \mathbf{Y}$$
(11.186)

and

$$\mathbf{Z}' = Q^{\top} \mathbf{Z},\tag{11.187}$$

we obtain

$$\mathbf{Y}' = \mathbf{X}' + \mathbf{Z}'. \tag{11.188}$$

Note that

$$E\mathbf{Z}' = E(Q\mathbf{Z}) = Q(E\mathbf{Z}) = Q \cdot 0 = 0,$$
 (11.189)

and  $\mathbf{Z}'$  is jointly Gaussian because it is a linear transformation of  $\mathbf{Z}$ . By Proposition 10.5,

$$K_{\mathbf{Z}'} = Q^{\top} K_{\mathbf{Z}} Q = Q^{\top} (Q \Lambda Q^{\top}) Q = \Lambda.$$
(11.190)

Thus the random variables in  $\mathbf{Z}'$  are uncorrelated. Since  $\mathbf{Z}'$  is jointly Gaussian, this implies that the random variables in  $\mathbf{Z}'$  are mutually independent. Therefore, we conclude that  $Z'_i \sim \mathcal{N}(0, \lambda_i)$ , where  $\lambda_i$  is the *i*th diagonal element of  $\Lambda$ .

We are then motivated to convert the given system of correlated Gaussian channels into the system shown in Figure 11.3, with  $\mathbf{X}'$  and  $\mathbf{Y}'$  being the input and output, respectively. Note that in this system,  $\mathbf{X}'$  and  $\mathbf{Y}'$  are related to  $\mathbf{X}$  and  $\mathbf{Y}$  as prescribed in (11.185) and (11.186), respectively. We then see from (11.188) that  $\mathbf{Z}'$  is the equivalent noise vector of the system with  $Z'_i$ being the noise variable of the *i*th channel. Hence, the system in Figure 11.3 is a system of parallel Gaussian channels. By Proposition 10.9, the total input power constraint in (11.182) for the original system translates to the total input power constraint

$$E\sum_{i=1}^{\kappa} (X_i')^2 \le P \tag{11.191}$$

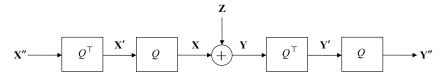


Fig. 11.4. A system identical to the system of correlated Gaussian channels.

for the system in Figure 11.3.

The question is whether the capacity of the system in Figure 11.3 is the same as the capacity of the original system. Let us called these two capacities C' and C, respectively. Intuitively, C' and C should be the same because the matrix Q is invertible. A formal proof goes as follows. We remind the reader that the capacity of a channel is the highest possible asymptotic rate at which information can be transmitted reliably through the channel by means of any encoding/decoding process. In Figure 11.3, by regarding the transformation Q on  $\mathbf{X}'$  as part of the encoding process and the transformation  $Q^{\top}$  on  $\mathbf{Y}$  as part of the decoding process, we see that  $C' \leq C$ . Now further convert the system in Figure 11.3 into the system in Figure 11.4 with input  $\mathbf{X}''$  and output  $\mathbf{Y}''$ , and call the capacity of this system C''. By repeating the same argument, we see that  $C' \leq C'$ . Thus  $C'' \leq C' \leq C$ . However, the system in Figure 11.4 is equivalent to the original system because  $Q^{\top}Q = I$ . Therefore, C'' = C, which implies C' = C.

Upon converting the given system of correlated Gaussian channels into an equivalent system of parallel Gaussian channels, we see that the capacity of the system is equal to

$$\frac{1}{2}\sum_{i=1}^{k}\log\left(1+\frac{a_i^*}{\lambda_i}\right) \tag{11.192}$$

where  $a_i^*$  is the optimal power allocated to the *i*th channel in the equivalent system, and its value can be obtained by water-filling as prescribed in Proposition 11.23. The reader should compare (11.192) with the formula in (11.179) for the capacity of parallel Gaussian channels.

Let  $A^*$  be the  $k \times k$  diagonal matrix with  $a_i^*$  being the *i*th diagonal element. From the discussion in the last section, the optimal distribution for the input  $\mathbf{X}'$  in the equivalent system of parallel channels is  $\mathcal{N}(0, A^*)$ . Accordingly, the distribution of  $\mathbf{X}$  is  $\mathcal{N}(0, QA^*Q^{\top})$ . We leave it as an exercise for the reader to verify that this indeed gives the optimal input distribution for the original system of correlated Gaussian channels.

## 11.7 The Bandlimited White Gaussian Channel

In this section, we discuss a bandlimited waveform channel with zero-mean *additive white Gaussian noise* (AWGN). In the rest of this chapter, the letters

j and f are reserved for  $\sqrt{-1}$  and "frequency," respectively. We begin with a few definitions from signal analysis. All the signals are assumed to be real.

**Definition 11.24.** The Fourier transform of a signal g(t) is defined as

$$G(f) = \int_{-\infty}^{\infty} g(t)e^{-j2\pi ft}dt.$$
 (11.193)

The signal g(t) can be recovered from G(f) as

$$g(t) = \int_{-\infty}^{\infty} G(f) e^{j2\pi ft} df,$$
 (11.194)

and g(t) is called the inverse Fourier transform of G(f). The functions g(t)and G(f) are said to form a transform pair, denoted by

$$g(t) \rightleftharpoons G(f). \tag{11.195}$$

The variables t and f are referred to as time and frequency, respectively.

In general, the Fourier transform of a signal g(t) may not exist. A sufficient condition for the Fourier transform of g(t) to exist is that g(t) has finite energy, i.e.,

$$\int_{-\infty}^{\infty} |g(t)|^2 dt < \infty.$$
(11.196)

A signal with finite energy is called an *energy signal*.

**Definition 11.25.** Let  $g_1(t)$  and  $g_2(t)$  be a pair of energy signals. The crosscorrelation function for  $g_1(t)$  and  $g_2(t)$  is defined as

$$R_{12}(\tau) = \int_{-\infty}^{\infty} g_1(t)g_2(t-\tau)dt.$$
 (11.197)

**Proposition 11.26.** For a pair of energy signals  $g_1(t)$  and  $g_2(t)$ ,

$$R_{12}(\tau) \rightleftharpoons G_1(f)G_2^*(f), \qquad (11.198)$$

where  $G_2^*(f)$  denotes the complex conjugate of  $G_2(f)$ .

**Definition 11.27.** For a wide-sense stationary process  $\{X(t), -\infty < t < \infty\}$ , the autocorrelation function is defined as

$$R_X(\tau) = E[X(t+\tau)X(t)],$$
(11.199)

which does not depend on t, and the power spectral density is defined as

$$S_X(f) = \int_{-\infty}^{\infty} R_X(\tau) e^{-j2\pi f\tau} d\tau, \qquad (11.200)$$

*i.e.*,

$$R_X(\tau) \rightleftharpoons S_X(f). \tag{11.201}$$

**Definition 11.28.** Let  $\{(X(t), Y(t)), -\infty < t < \infty\}$  be a bivariate wide-sense stationary process. Their cross-correlation functions are defined as

$$R_{XY}(\tau) = E[X(t+\tau)Y(t)]$$
(11.202)

and

$$R_{YX}(\tau) = E[Y(t+\tau)X(t)], \qquad (11.203)$$

which do not depend on t. The cross-spectral densities are defined as

$$S_{XY}(f) = \int_{-\infty}^{\infty} R_{XY}(\tau) e^{-j2\pi f\tau} d\tau \qquad (11.204)$$

and

$$S_{YX}(f) = \int_{-\infty}^{\infty} R_{YX}(\tau) e^{-j2\pi f\tau} d\tau,$$
 (11.205)

*i.e.*,

$$R_{XY}(\tau) \rightleftharpoons S_{XY}(f)$$
 (11.206)

and

$$R_{YX}(\tau) \rightleftharpoons S_{YX}(f).$$
 (11.207)

We now describe the simplest nontrivial model for a waveform channel. In wired-line and wireless communication, the frequency spectrum of the medium is often partitioned into a number of communication channels, where each channel occupies a certain frequency band. Consider such a channel that occupies the frequency band  $[f_l, f_h]$  with  $0 \le f_l < f_h$ , where

$$W = f_h - f_l \tag{11.208}$$

is called the *bandwidth*. The input process X(t) is contaminated by a zeromean additive white Gaussian noise process with power  $\frac{N_0}{2}$ , i.e.,

$$S_Z(f) = \frac{N_0}{2}, \quad -\infty < f < \infty.$$
 (11.209)

In reality, such a noise process cannot exist because its total power is infinite. For practical purposes, one can regard the power spectral density to be constant within the range of interest of the problem.

Let h(t) be the impulse response of an ideal bandpass filter for the frequency band  $[f_l, f_h]$ , i.e.,

$$H(f) = \begin{cases} 1 \text{ if } f_l \le |f| \le f_h \\ 0 \text{ otherwise.} \end{cases}$$
(11.210)

At the receiver for this channel, the ideal bandpass filter h(t) is applied to the received signal in order to filter out the frequency components due to other channels. Effectively, we can regard this filtered version of the received signal given by

276 11 Continuous-Valued Channels

$$Y(t) = [X(t) + Z(t)] * h(t) = X(t) * h(t) + Z(t) * h(t)$$
(11.211)

as the output of the channel, where \* denotes convolution in the time domain. Letting

$$X'(t) = X(t) * h(t)$$
(11.212)

and

$$Z'(t) = Z(t) * h(t), \qquad (11.213)$$

(11.211) can be written as

$$Y(t) = X'(t) + Z'(t).$$
(11.214)

The only difference between X(t) and X'(t) is that all the frequency components of X'(t) are in  $[f_l, f_h]$ , while X(t) can have frequency components outside this range. However, even if such frequency components exist in X(t), they are filtered out by the ideal bandpass filter h(t) and do not appear in the output process Y(t). Therefore, we can regard X'(t) instead of X(t) as the input process of the channel. By the same argument, we regard Z'(t) instead of Z(t) as the noise process of the channel. As for the memoryless Gaussian channel discussed in the last section, we impose an average power constraint P on the input process X'(t).

For simplicity, we consider in this section the case that the channel we have described occupies the frequency band [0, W]. This channel, called the *bandlimited white Gaussian channel*, is the special case of the general model with  $f_l = 0$ .

While a rigorous formulation of the bandlimited white Gaussian channel involves mathematical tools beyond the scope of this book, we will nevertheless give a heuristic argument that suggests the formula for the channel capacity. The sampling theorem in signal analysis will allow us to "convert" this waveform channel into a memoryless Gaussian channel discussed in the last section.

**Theorem 11.29 (Sampling Theorem).** Let g(t) be a signal with Fourier transform G(f) that vanishes for  $f \notin [-W, W]$ . Then

$$g(t) = \sum_{i=-\infty}^{\infty} g\left(\frac{i}{2W}\right) \operatorname{sinc}(2Wt - i)$$
(11.215)

for  $-\infty < t < \infty$ , where

$$\operatorname{sinc}(t) = \frac{\sin(\pi t)}{\pi t},\tag{11.216}$$

called the sinc function, is defined to be 1 at t = 0 by continuity.

Letting

$$g_i = \frac{1}{\sqrt{2W}} g\left(\frac{i}{2W}\right) \tag{11.217}$$

11.7 The Bandlimited White Gaussian Channel 277

and

$$\psi_i(t) = \sqrt{2W}\operatorname{sinc}(2Wt - i), \qquad (11.218)$$

the formula in (11.215) can be rewritten as

$$g(t) = \sum_{i=-\infty}^{\infty} g_i \psi_i(t).$$
(11.219)

**Proposition 11.30.**  $\psi_i(t)$ ,  $-\infty < i < \infty$  form an orthonormal basis for signals which are bandlimited to [0, W].

*Proof.* Since

$$\psi_i(t) = \psi_0\left(t - \frac{i}{2W}\right),\tag{11.220}$$

 $\psi_i(t)$  and  $\psi_0(t)$  have the same energy. We first show that

$$\int_{-\infty}^{\infty} \operatorname{sinc}^2(2Wt) dt = \frac{1}{2W}.$$
 (11.221)

This integral is difficult to evaluate directly. Instead we consider

$$\operatorname{sinc}(2Wt) \rightleftharpoons \frac{1}{2W} \operatorname{rect}\left(\frac{f}{2W}\right),$$
 (11.222)

where

$$\operatorname{rect}(f) = \begin{cases} 1 - \frac{1}{2} \le f \le \frac{1}{2} \\ 0 \text{ otherwise.} \end{cases}$$
(11.223)

Then by Rayleigh's energy theorem, we have

$$\int_{-\infty}^{\infty} \operatorname{sinc}^2(2Wt) dt = \int_{-\infty}^{\infty} \left(\frac{1}{2W}\right)^2 \operatorname{rect}^2\left(\frac{f}{2W}\right) df = \frac{1}{2W}.$$
 (11.224)

It then follows that

$$\int_{-\infty}^{\infty} \psi_i^2(t) dt = \int_{-\infty}^{\infty} \psi_0^2(t) dt \tag{11.225}$$

$$= (\sqrt{2W})^2 \int_{-\infty}^{\infty} \operatorname{sinc}^2(2Wt) dt \qquad (11.226)$$

$$=2W\left(\frac{1}{2W}\right) \tag{11.227}$$

$$= 1.$$
 (11.228)

Next, we show that

$$\int_{-\infty}^{\infty} \operatorname{sinc}(2Wt - i) \operatorname{sinc}(2Wt - i') dt \qquad (11.229)$$

#### 278 11 Continuous-Valued Channels

vanishes whenever  $i \neq i'$ . Again, this integral is difficult to evaluate directly. Since (11.224) implies that both  $\operatorname{sinc}(2Wt - i)$  and  $\operatorname{sinc}(2Wt - i')$  have finite energy, we can consider their cross-correlation function, denoted by  $R_{ii'}(\tau)$ . Now

$$\operatorname{sinc}(2Wt - i) \rightleftharpoons e^{-j2\pi f\left(\frac{i}{2W}\right)} \left(\frac{1}{2W}\right) \operatorname{rect}\left(\frac{f}{2W}\right) := G_i(f) \qquad (11.230)$$

and

$$\operatorname{sinc}(2Wt - i') \rightleftharpoons e^{-j2\pi f\left(\frac{i'}{2W}\right)} \left(\frac{1}{2W}\right) \operatorname{rect}\left(\frac{f}{2W}\right) := G_{i'}(f). \quad (11.231)$$

Then we have

$$R_{ii'}(\tau) \rightleftharpoons G_i(f)G_{i'}^*(f), \qquad (11.232)$$

and the integral in (11.229) is given by

$$R_{ii'}(0) = \int_{-\infty}^{\infty} G_i(f) G_{i'}^*(f) df, \qquad (11.233)$$

which vanishes whenever  $i \neq i'$ . Therefore,

$$\int_{-\infty}^{\infty} \psi_i(t)\psi_{i'}(t)dt = 2W \int_{-\infty}^{\infty} \operatorname{sinc}(2Wt - i)\operatorname{sinc}(2Wt - i')dt \quad (11.234)$$
  
= 0. (11.235)

Together with (11.228), this shows that  $\psi_i(t), -\infty < i < \infty$  form an orthonormal set. Finally, since g(t) in (11.219) is any signal bandlimited to [0, W], we conclude that  $\psi_i(t), -\infty < i < \infty$  form an orthonormal basis for such signals. The theorem is proved.  $\Box$ 

Let us return to our discussion of the waveform channel. The sampling theorem implies that the input process X'(t), assuming the existence of the Fourier transform, can be written as

$$X'(t) = \sum_{i=-\infty}^{\infty} X'_{i}\psi_{i}(t),$$
(11.236)

where

$$X_i' = \frac{1}{\sqrt{2W}} X'\left(\frac{i}{2W}\right),\tag{11.237}$$

and there is a one-to-one correspondence between X'(t) and  $\{X'_i, -\infty < i < \infty\}$ . The same applies to (a realization of) the output process Y(t), which we assume can be written as

$$Y(t) = \sum_{i=-\infty}^{\infty} Y'_{i} \psi_{i}(t), \qquad (11.238)$$

11.7 The Bandlimited White Gaussian Channel 279

where

$$Y_i = \frac{1}{\sqrt{2W}} Y\left(\frac{i}{2W}\right). \tag{11.239}$$

With these assumptions on X'(t) and Y(t), the waveform channel is equivalent to a discrete-time channel defined at  $t = \frac{i}{2W}$ , with the *i*th input and output of the channel being  $X'_i$  and  $Y_i$ , respectively.

Toward determining the capacity of this equivalent discrete-time channel, we prove in the following a characterization of the effect of the noise process Z'(t) at the sampling times.

**Proposition 11.31.**  $Z'\left(\frac{i}{2W}\right), -\infty < i < \infty$  are *i.i.d.* Gaussian random variables with zero mean and variance  $N_0W$ .

*Proof.* First of all, Z(t) is a zero-mean Gaussian process and Z'(t) is a filtered version of Z(t), so Z'(t) is also a zero-mean Gaussian process. Consequently,  $Z'\left(\frac{i}{2W}\right)$  are zero-mean Gaussian random variables. The power spectral density of Z'(t) is given by

$$S_{Z'}(f) = \begin{cases} \frac{N_0}{2} & -W \le f \le W\\ 0 & \text{otherwise.} \end{cases}$$
(11.240)

Then the autocorrelation function of Z'(t), which is the inverse Fourier transform of  $S_{Z'}(f)$ , is given by

$$R_{Z'}(\tau) = N_0 W \operatorname{sinc}(2W\tau). \tag{11.241}$$

It is seen that the value of  $R_{Z'}(\tau)$  is equal to 0 when  $\tau = \frac{i}{2W}$  for all  $i \neq 0$ , because the sinc function in (11.216) vanishes at all nonzero integer values of t. This shows that  $Z'\left(\frac{i}{2W}\right), -\infty < i < \infty$  are uncorrelated and hence independent because they are jointly Gaussian. Finally, since  $Z'\left(\frac{i}{2W}\right)$  has zero mean, in light of (11.199), its variance is given by  $R_{Z'}(0) = N_0 W$ .  $\Box$ 

Recall from (11.214) that

$$Y(t) = X'(t) + Z'(t).$$
(11.242)

Then

$$Y\left(\frac{i}{2W}\right) = X'\left(\frac{i}{2W}\right) + Z'\left(\frac{i}{2W}\right).$$
(11.243)

Upon dividing by  $\sqrt{2W}$  and letting

$$Z'_{i} = \frac{1}{\sqrt{2W}} Z'\left(\frac{i}{2W}\right), \qquad (11.244)$$

it follows from (11.237) and (11.239) that

280 11 Continuous-Valued Channels

$$Y_i = X'_i + Z'_i. (11.245)$$

Since  $Z'\left(\frac{i}{2W}\right)$ ,  $-\infty < i < \infty$  are i.i.d. with distribution  $\mathcal{N}(0, N_0 W)$ ,  $Z'_i$ ,  $-\infty < i < \infty$  are i.i.d. with distribution  $\mathcal{N}(0, \frac{N_0}{2})$ .

Thus we have shown that the bandlimited white Gaussian channel is equivalent to a memoryless Gaussian channel with noise power equal to  $\frac{N_0}{2}$ . As we are converting the waveform channel into a discrete-time channel, we need to relate the input power constraint of the waveform channel to the input power constraint of the discrete-time channel. Let P' be the average energy (i.e., the second moment) of the  $X_i$ 's. We now calculate the average power of X'(t)in terms of P'. Since  $\psi_i(t)$  has unit energy, the average contribution to the energy of X'(t) by each sample is P'. As there are 2W samples per unit time and  $\psi_i(t), -\infty < i < \infty$  are orthonormal, X'(t) accumulates energy from the samples at a rate equal to 2WP'. Upon considering

$$2WP' \le P,\tag{11.246}$$

where P is the average power constraint on the input process X'(t), we obtain

$$P' \le \frac{P}{2W},\tag{11.247}$$

i.e., an input power constraint P for the bandlimited Gaussian channel translates to an input power constraint  $\frac{P}{2W}$  for the discrete-time channel. By Theorem 11.21, the capacity of the memoryless Gaussian channel is

$$\frac{1}{2}\log\left(1+\frac{P/2W}{N_0/2}\right) = \frac{1}{2}\log\left(1+\frac{P}{N_0W}\right) \quad \text{bits per sample.} \tag{11.248}$$

Since there are 2W samples per unit time, we conclude that the capacity of the bandlimited Gaussian channel is

$$W \log\left(1 + \frac{P}{N_0 W}\right)$$
 bits per unit time. (11.249)

The argument we have given above is evidently not rigorous because if there is no additional constraint on the  $X_i$ 's other than their average energy not exceeding  $\frac{P}{2W}$ , then X'(t) may not have finite energy. This induces a gap in the argument because the Fourier transform of X'(t) may not exist and hence the sampling theorem cannot be applied.

A rigorous formulation of the bandlimited white Gaussian channel involves the consideration of an input signal of finite duration, which is analogous to a code for the DMC with a finite block length. Since a signal with finite duration cannot be bandlimited, this immediate leads to a contradiction. Overcoming this technical difficulty requires the use of *prolate spheroidal wave functions* [305, 197, 198] which are bandlimited functions with most of the energy on a finite interval. The main idea is that there are approximately 2WT orthonormal basis functions for the set of signals which are bandlimited to W and have most of the energy on [0, T) in time. We refer the reader to Gallager[117] for a rigorous treatment of the bandlimited white Gaussian channel.

#### 11.8 The Bandlimited Colored Gaussian Channel

In the last section, we have discussed the bandlimited white Gaussian channel occupying the frequency band [0, W]. We presented a heuristic argument that led to the formula in (11.249) for the channel capacity. Suppose the channel instead occupies the frequency band  $[f_l, f_h]$ , with  $f_l$  being a multiple of  $W = f_h - f_l$ . Then the noise process Z'(t) has power spectral density

$$S_{Z'}(f) = \begin{cases} \frac{N_0}{2} & \text{if } f_l \le |f| \le f_h \\ 0 & \text{otherwise.} \end{cases}$$
(11.250)

We refer to such a channel as the *bandpass white Gaussian channel*. By an extension of the heuristic argument for the bandlimited white Gaussian channel, which would involve the bandpass version of the sampling theorem, the same formula for the channel capacity can be obtained. The details are omitted here.

We now consider a waveform channel occupying the frequency band [0, W]with input power constraint P and zero-mean additive *colored* Gaussian noise Z(t). We refer to this channel as the *bandlimited colored Gaussian channel*. To analyze the capacity of this channel, divide the interval [0, W] into subintervals  $[f_l^i, f_h^i]$  for  $1 \le i \le k$ , where

$$f_l^i = (i-1)\Delta_k \tag{11.251}$$

$$f_h^i = i\Delta_k, \tag{11.252}$$

and

$$\Delta_k = \frac{W}{k} \tag{11.253}$$

is the width of each subinterval. As an approximation, assume  $S_Z(f)$  is equal to a constant  $S_{Z,i}$  over the subinterval  $[f_l^i, f_h^i]$ . Then the channel consists of k sub-channels, with the *i*th sub-channel being a bandpass (bandlimited if i = 1) white Gaussian channel occupying the frequency band  $[f_l^i, f_h^i]$ . Thus by letting  $N_0 = 2S_{Z,i}$  in (11.250), we obtain from (11.249) that the capacity of the *i*th sub-channel is equal to

$$\Delta_k \log\left(1 + \frac{P_i}{2S_{Z,i}\Delta_k}\right) \tag{11.254}$$

if  $P_i$  is the input power allocated to that sub-channel.

The noise process of the *i*th sub-channel, denoted by  $Z'_i(t)$ , is obtained by passing Z(t) through the ideal bandpass filter with frequency response

$$H_i(f) = \begin{cases} 1 \text{ if } f_l^i \le |f| \le f_h^i \\ 0 \text{ otherwise.} \end{cases}$$
(11.255)

It can be shown (see Problem 10) that the noise processes  $Z_i(t)$ ,  $1 \le i \le k$  are independent. By converting each sub-channel into an equivalent memoryless

#### 282 11 Continuous-Valued Channels

Gaussian channel as discussed in the last section, we see that the k subchannels can be regarded as a system of parallel Gaussian channels. Thus the channel capacity is equal to the sum of the capacities of the individual sub-channels when the power allocation among the k sub-channels is optimal.

Let  $P_i^*$  be the optimal power allocation for the *i*th sub-channel. Then it follows from (11.254) that the channel capacity is equal to

$$\sum_{i=1}^{k} \Delta_k \log\left(1 + \frac{P_i^*}{2S_{Z,i}\Delta_k}\right) = \sum_{i=1}^{k} \Delta_k \log\left(1 + \frac{\frac{P_i^*}{2\Delta_k}}{S_{Z,i}}\right),$$
(11.256)

where by Lemma 11.23,

$$\frac{P_i^*}{2\Delta_k} = (\nu - S_{Z,i})^+, \qquad (11.257)$$

or

$$P_i^* = 2\Delta_k (\nu - S_{Z,i})^+, \qquad (11.258)$$

with

$$\sum_{i=1}^{k} P_i^* = P. \tag{11.259}$$

Then from (11.258) and (11.259), we obtain

$$\sum_{i=1}^{k} (\nu - S_{Z,i})^{+} \Delta_{k} = \frac{P}{2}.$$
(11.260)

As  $k \to \infty$ , following (11.256) and (11.257),

$$\sum_{i=1}^{k} \Delta_k \log \left( 1 + \frac{\frac{P_i^*}{2\Delta_k}}{S_{Z,i}} \right) = \sum_{i=1}^{k} \Delta_k \log \left( 1 + \frac{(\nu - S_{Z,i})^+}{S_{Z,i}} \right)$$
(11.261)

$$+ \int_{0}^{W} \log\left(1 + \frac{(\nu - S_Z(f))^+}{S_Z(f)}\right) df$$
 (11.262)

$$= \frac{1}{2} \int_{-W}^{W} \log\left(1 + \frac{(\nu - S_Z(f))^+}{S_Z(f)}\right) df, \quad (11.263)$$

and following (11.260),

$$\sum_{i=1}^{k} (\nu - S_{Z,i})^{+} \Delta_{k} \to \int_{0}^{W} (\nu - S_{Z}(f))^{+} df$$
 (11.264)

$$= \frac{1}{2} \int_{-W}^{W} (\nu - S_Z(f))^+ df, \qquad (11.265)$$

where (11.263) and (11.265) are obtained by noting that

$$S_{Z'}(-f) = S_{Z'}(f) \tag{11.266}$$

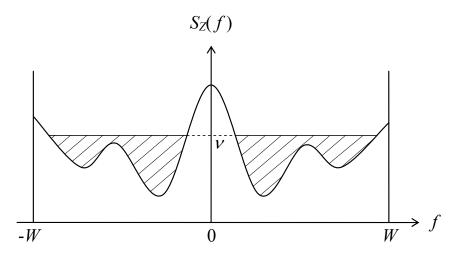


Fig. 11.5. Water-filling for the bandlimited colored Gaussian channel.

for  $-\infty < f < \infty$  (see Problem 9). Hence, we conclude that the capacity of the bandlimited colored Gaussian channel is equal to

$$\frac{1}{2} \int_{-W}^{W} \log\left(1 + \frac{(\nu - S_Z(f))^+}{S_Z(f)}\right) df \quad \text{bits per unit time,}$$
(11.267)

where  $\nu$  satisfies

$$\int_{-W}^{W} (\nu - S_Z(f))^+ df = P \tag{11.268}$$

in view of (11.260). Figure 11.5 is an illustration of the water-filling process for determining  $\nu$ , where the amount of water to be poured into the reservoir is equal to P.

# 11.9 Zero-Mean Gaussian Noise is the Worst Additive Noise

In the last section, we derived the capacity for a system of correlated Gaussian channels, where the noise vector is a zero-mean Gaussian random vector. In this section, we show that in terms of the capacity of the system, the zero-mean Gaussian noise is the worst additive noise given that the noise vector has a bounded joint pdf and a fixed correlation matrix. Note that the diagonal elements of this correlation matrix specify the power of the individual noise variables, while the other elements in the matrix give a characterization of the correlation between the noise variables.

**Theorem 11.32.** For a fixed zero-mean Gaussian random vector  $\mathbf{X}^*$ , let

28411 Continuous-Valued Channels

$$\mathbf{Y} = \mathbf{X}^* + \mathbf{Z},\tag{11.269}$$

where the joint pdf of  $\mathbf{Z}$  exists and is bounded, and  $\mathbf{Z}$  is independent of  $\mathbf{X}^*$ . Under the constraint that the correlation matrix of  $\mathbf{Z}$  is equal to K, where K is any symmetric positive definite matrix,  $I(\mathbf{X}^*; \mathbf{Y})$  is minimized if and only if  $\mathbf{Z} \sim \mathcal{N}(0, K)$ .

Lemma 11.33. Let  $\mathbf{X}$  be a zero-mean random vector and

$$\mathbf{Y} = \mathbf{X} + \mathbf{Z},\tag{11.270}$$

where  $\mathbf{Z}$  is independent of  $\mathbf{X}$ . Then

$$\tilde{K}_{\mathbf{Y}} = \tilde{K}_{\mathbf{X}} + \tilde{K}_{\mathbf{Z}}.$$
(11.271)

*Proof.* For any i and j, consider

$$EY_i Y_j = E(X_i + Z_i)(X_j + Z_j)$$
(11.272)

$$= E(X_i X_j + X_i Z_j + Z_i X_j + Z_i Z_j)$$
(11.273)  
$$= EX_i X_j + EX_i Z_j + EZ_i X_j + EZ_i Z_j$$
(11.274)

$$= EX_iX_j + EX_iZ_j + EZ_iX_j + EZ_iZ_j$$
(11.274)

$$= EX_iX_j + (EX_i)(EZ_j) + (EZ_i)(EX_j) + EZ_iZ_j$$
(11.275)

$$= EX_iX_j + (0)(EZ_j) + (EZ_i)(0) + EZ_iZ_j$$
(11.276)

$$= EX_i X_j + EZ_i Z_j, (11.277)$$

where (11.276) follows from the assumption that  $X_i$  has zero mean for all *i*. The proposition is proved.  $\Box$ 

**Lemma 11.34.** Let  $\mathbf{Y}^* \sim \mathcal{N}(0, K)$  and  $\mathbf{Y}$  be any random vector with correlation matrix K. Then

$$\int f_{\mathbf{Y}^*}(\mathbf{y}) \log f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y} = \int_{\mathcal{S}_{\mathbf{Y}}} f_{\mathbf{Y}}(\mathbf{y}) \log f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y}.$$
 (11.278)

*Proof.* The random vector  $\mathbf{Y}^*$  has joint pdf

$$f_{\mathbf{Y}^*}(\mathbf{y}) = \frac{1}{\left(\sqrt{2\pi}\right)^k |K|^{1/2}} e^{-\frac{1}{2}(\mathbf{y}^\top K^{-1} \mathbf{y})}$$
(11.279)

for all  $\mathbf{y} \in \Re^k$ . Since  $E\mathbf{Y}^* = 0$ ,  $\tilde{K}_{\mathbf{Y}^*} = K_{\mathbf{Y}^*} = K$ . Therefore,  $\mathbf{Y}^*$  and  $\mathbf{Y}$  have the same correlation matrix. Consider

11.9 Zero-Mean Gaussian Noise is the Worst Additive Noise 285

$$\int \left[\ln f_{\mathbf{Y}^*}(\mathbf{y})\right] f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y}$$
$$= \int \left[-\frac{1}{2} (\mathbf{y}^\top K^{-1} \mathbf{y}) - \ln \left[(\sqrt{2\pi})^k |K|^{1/2}\right]\right] f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y}$$
(11.280)

$$= -\frac{1}{2} \int (\mathbf{y}^{\top} K^{-1} \mathbf{y}) f_{\mathbf{Y}^{*}}(\mathbf{y}) d\mathbf{y} - \ln\left[ (\sqrt{2\pi})^{k} |K|^{1/2} \right]$$
(11.281)

$$= -\frac{1}{2} \int \left[ \sum_{i,j} (K^{-1})_{ij} y_i y_j \right] f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y} - \ln\left[ (\sqrt{2\pi})^k |K|^{1/2} \right]$$
(11.282)

$$= -\frac{1}{2} \sum_{i,j} (K^{-1})_{ij} \int (y_i y_j) f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y} - \ln\left[ (\sqrt{2\pi})^k |K|^{1/2} \right]$$
(11.283)

$$= -\frac{1}{2} \sum_{i,j} (K^{-1})_{ij} \int_{\mathcal{S}_{\mathbf{Y}}} (y_i y_j) f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y} - \ln\left[ (\sqrt{2\pi})^k |K|^{1/2} \right]$$
(11.284)

$$= \int_{\mathcal{S}_{\mathbf{Y}}} \left[ -\frac{1}{2} \mathbf{y}^{\top} K^{-1} \mathbf{y} - \ln \left[ (\sqrt{2\pi})^k |K|^{1/2} \right] \right] f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y}$$
(11.285)

$$= \int_{\mathcal{S}_{\mathbf{Y}}} \left[ \ln f_{\mathbf{Y}^*}(\mathbf{y}) \right] f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y}.$$
(11.286)

In the above, (11.284) is justified because  $\mathbf{Y}$  and  $\mathbf{Y}^*$  have the same correlation matrix, and (11.285) is obtained by backtracking the manipulations from (11.280) to (11.283) with  $f_{\mathbf{Y}}(\mathbf{y})$  in place of  $f_{\mathbf{Y}^*}(\mathbf{y})$ . The lemma is proved upon changing the base of the logarithm.  $\Box$ 

Proof of Theorem 11.32. Let  $\mathbf{Z}^* \sim \mathcal{N}(0, K)$  such that  $\mathbf{Z}^*$  is independent of  $\mathbf{X}^*$ , and let

$$Y^* = X^* + Z^*. (11.287)$$

Obviously, the support of  $\mathbf{Y}^*$  is  $\Re^k$  because  $\mathbf{Y}^*$  has a multivariate Gaussian distribution. Note that the support of  $\mathbf{Y}$  is also  $\Re^k$  regardless of the distribution of Z because the support of  $\mathbf{X}^*$  is  $\Re^k$ . We need to prove that for any random vector  $\mathbf{Z}$  with correlation matrix K, where  $\mathbf{Z}$  is independent of  $\mathbf{X}^*$  and the joint pdf of  $\mathbf{Z}$  exists and is bounded,

$$I(\mathbf{X}^*; \mathbf{Y}^*) \le I(\mathbf{X}^*; \mathbf{Y}). \tag{11.288}$$

Since  $E\mathbf{Z}^* = 0$ ,  $\tilde{K}_{\mathbf{Z}^*} = K_{\mathbf{Z}^*} = K$ . Therefore,  $\mathbf{Z}^*$  and  $\mathbf{Z}$  have the same correlation matrix. By noting that  $\mathbf{X}^*$  has zero mean, we apply Lemma 11.33 to see that  $\mathbf{Y}^*$  and  $\mathbf{Y}$  have the same correlation matrix.

The inequality in (11.288) can be proved by considering

$$I(\mathbf{X}^*; \mathbf{Y}^*) - I(\mathbf{X}^*; \mathbf{Y})$$

$$\stackrel{a)}{=} h(\mathbf{Y}^*) - h(\mathbf{Z}^*) - h(\mathbf{Y}) + h(\mathbf{Z})$$

$$= -\int f_{\mathbf{Y}^*}(\mathbf{y}) \log f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y} + \int f_{\mathbf{Z}^*}(\mathbf{z}) \log f_{\mathbf{Z}^*}(\mathbf{z}) d\mathbf{z}$$
(11.289)

28611 Continuous-Valued Channels

$$+ \int f_{\mathbf{Y}}(\mathbf{y}) \log f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y} - \int_{\mathcal{S}_{\mathbf{Z}}} f_{\mathbf{Z}}(\mathbf{z}) \log f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z}$$
(11.290)  
$$\stackrel{b)}{=} - \int f_{\mathbf{Y}}(\mathbf{y}) \log f_{\mathbf{Y}^{*}}(\mathbf{y}) d\mathbf{y} + \int f_{\mathbf{Z}}(\mathbf{z}) \log f_{\mathbf{Z}^{*}}(\mathbf{z}) d\mathbf{z}$$

$$+ \int f_{\mathbf{Y}}(\mathbf{y}) \log f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y} + \int_{\mathcal{S}_{\mathbf{Z}}} f_{\mathbf{Z}}(\mathbf{z}) \log f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z} + \int f_{\mathbf{Y}}(\mathbf{y}) \log f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y} - \int_{\mathcal{S}_{\mathbf{Z}}} f_{\mathbf{Z}}(\mathbf{z}) \log f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z}$$
(11.291)

$$= \int \log\left(\frac{f_{\mathbf{Y}}(\mathbf{y})}{f_{\mathbf{Y}^*}(\mathbf{y})}\right) f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y} + \int_{\mathcal{S}_{\mathbf{Z}}} \log\left(\frac{f_{\mathbf{Z}^*}(\mathbf{z})}{f_{\mathbf{Z}}(\mathbf{z})}\right) f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z}$$
(11.292)

$$\stackrel{(z)}{=} \int_{\mathcal{S}_{\mathbf{Z}}} \int \log\left(\frac{f_{\mathbf{Y}}(\mathbf{y})f_{\mathbf{Z}^*}(\mathbf{z})}{f_{\mathbf{Y}^*}(\mathbf{y})f_{\mathbf{Z}}(\mathbf{z})}\right) f_{\mathbf{Y}\mathbf{Z}}(\mathbf{y}, \mathbf{z}) d\mathbf{y} d\mathbf{z}$$
(11.293)

$$\stackrel{d}{\leq} \log\left(\int_{\mathcal{S}_{\mathbf{Z}}} \int \frac{f_{\mathbf{Y}}(\mathbf{y}) f_{\mathbf{Z}^*}(\mathbf{z})}{f_{\mathbf{Y}^*}(\mathbf{y}) f_{\mathbf{Z}}(\mathbf{z})} f_{\mathbf{Y}\mathbf{Z}}(\mathbf{y}, \mathbf{z}) d\mathbf{y} d\mathbf{z}\right)$$
(11.294)

$$\stackrel{e}{=} \log\left(\int \left[\frac{1}{f_{\mathbf{Y}^*}(\mathbf{y})} \int_{\mathcal{S}_{\mathbf{Z}}} f_{\mathbf{X}^*}(\mathbf{y} - \mathbf{z}) f_{\mathbf{Z}^*}(\mathbf{z}) d\mathbf{z}\right] f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y}\right)$$
(11.295)

$$\stackrel{J}{\leq} \log\left(\int \frac{f_{\mathbf{Y}^*}(\mathbf{y})}{f_{\mathbf{Y}^*}(\mathbf{y})} f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y}\right)$$
(11.296)  
= 0. (11.297)

= 0.

The above steps are explained as follows:

We assume that the pdf of **Z** exists and is bounded. Therefore,  $h(\mathbf{Z})$  is ٠ defined. Moreover,

$$f_{\mathbf{Y}|\mathbf{X}^*}(\mathbf{y}|\mathbf{x}) = f_{\mathbf{Z}}(\mathbf{y} - \mathbf{x})$$
(11.298)

is bounded. By Proposition 10.26,  $f_{\mathbf{Y}}(\mathbf{y})$  exists and hence  $h(\mathbf{Y})$  is defined.

- In b), we have replaced  $f_{\mathbf{Y}^*}(\mathbf{y})$  by  $f_{\mathbf{Y}}(\mathbf{y})$  in the first integral and replaced ٠  $f_{\mathbf{Z}^*}(\mathbf{z})$  by  $f_{\mathbf{Z}}(\mathbf{z})$  in the second integral. The former is justified by an application of Lemma 11.34 to  $\mathbf{Y}^*$  and  $\mathbf{Y}$  by noting that  $\mathbf{Y}^*$  is a zero-mean Gaussian random vector and  $\mathbf{Y}^*$  and  $\mathbf{Y}$  have the same correlation matrix. The latter is justified similarly.
- To justify c), we need  $S_{\mathbf{YZ}} = \Re^k \times S_{\mathbf{Z}}$ , which can be seen by noting that •

$$f_{\mathbf{Y}\mathbf{Z}}(\mathbf{y}, \mathbf{z}) = f_{\mathbf{Y}|\mathbf{Z}}(\mathbf{y}|\mathbf{z}) f_{\mathbf{Z}}(\mathbf{z}) = f_{\mathbf{X}^*}(\mathbf{y} - \mathbf{z}) f_{\mathbf{Z}}(\mathbf{z}) > 0$$
(11.299)

for all  $\mathbf{y} \in \Re^k$  and all  $\mathbf{z} \in \mathcal{S}_{\mathbf{Z}}$ .

- d) follows from Jensen's inequality and the concavity of the logarithmic • function.
- e) follows from (11.299).
- f) follows because ٠

$$\int_{\mathcal{S}_{\mathbf{z}}} f_{\mathbf{X}^*}(\mathbf{y} - \mathbf{z}) f_{\mathbf{Z}^*}(\mathbf{z}) d\mathbf{z} = \int_{\mathcal{S}_{\mathbf{z}}} f_{\mathbf{Y}^* | \mathbf{Z}^*}(\mathbf{y} | \mathbf{z}) f_{\mathbf{Z}^*}(\mathbf{z}) d\mathbf{z}$$
(11.300)

$$\leq \int f_{\mathbf{Y}^*|\mathbf{Z}^*}(\mathbf{y}|\mathbf{z}) f_{\mathbf{Z}^*}(\mathbf{z}) d\mathbf{z}$$
(11.301)

$$= f_{\mathbf{Y}^*}(\mathbf{y}). \tag{11.302}$$

Equality holds in (11.294) if and only if

$$f_{\mathbf{Y}}(\mathbf{y})f_{\mathbf{Z}^*}(\mathbf{z}) = f_{\mathbf{Y}^*}(\mathbf{y})f_{\mathbf{Z}}(\mathbf{z}) \quad \text{for all } \mathbf{y} \in \Re^k, \, \mathbf{z} \in \mathcal{S}_{\mathbf{Z}}.$$
(11.303)

If  $f_{\mathbf{Z}}(\mathbf{z}) = f_{\mathbf{Z}^*}(\mathbf{z})$  for all  $\mathbf{z} \in S_{\mathbf{Z}}$ , then  $S_{\mathbf{Z}} = \Re^k$  and  $\mathbf{Z} \sim \mathcal{N}(0, K)$ . This implies  $f_{\mathbf{Y}}(\mathbf{y}) = f_{\mathbf{Y}^*}(\mathbf{y})$  for all  $\mathbf{y} \in \Re^k$  in view of (11.269) and (11.287), so that (11.303) holds. Thus  $\mathbf{Z} \sim \mathcal{N}(0, K)$  is a sufficient condition for equality to hold in (11.294). Conversely, if equality holds in (11.294), we obtain from (11.303) that

$$f_{\mathbf{Z}^*}(\mathbf{z}) \int f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y} = f_{\mathbf{Z}}(\mathbf{z}) \int f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y}.$$
 (11.304)

Since

$$\int f_{\mathbf{Y}}(\mathbf{y}) d\mathbf{y} = \int f_{\mathbf{Y}^*}(\mathbf{y}) d\mathbf{y} = 1, \qquad (11.305)$$

we see that  $f_{\mathbf{Z}}(\mathbf{z}) = f_{\mathbf{Z}^*}(\mathbf{z})$  for all  $\mathbf{z} \in \mathcal{S}_{\mathbf{Z}}$ , so that  $\mathcal{S}_{\mathbf{Z}} = \Re^k$  and  $\mathbf{Z} \sim \mathcal{N}(0, K)$ . Hence, we conclude that  $\mathbf{Z} \sim \mathcal{N}(0, K)$  is a necessary and sufficient condition for  $I(\mathbf{X}^*; \mathbf{Y})$  to be minimized. The theorem is proved.  $\Box$ 

Consider the system of correlated Gaussian channels discussed in the last section. Denote the noise vector by  $\mathbf{Z}^*$  and its correlation matrix by K. Note that K is also the covariance matrix of  $\mathbf{Z}^*$  because  $\mathbf{Z}^*$  has zero mean. In other words,  $\mathbf{Z}^* \sim \mathcal{N}(0, K)$ . Refer to this system as the zero-mean Gaussian system and let  $C^*$  be its capacity. Then consider another system with exactly the same specification except that the noise vector, denoted by  $\mathbf{Z}$ , may neither be zero-mean nor Gaussian. We, however, require that the joint pdf of  $\mathbf{Z}$  exists and is bounded. Refer to this system as the alternative system and let C be its capacity.

We now apply Theorem 11.32 to show that  $C \ge C^*$ . Let  $\mathbf{X}^*$  be the input random vector that achieves the capacity of the zero-mean Gaussian system. We have mentioned at the end of Section 11.6 that  $\mathbf{X}^*$  is a zero-mean Gaussian random vector. Let  $\mathbf{Y}^*$  and  $\mathbf{Y}$  be defined in (11.287) and (11.269), which correspond to the outputs of the zero-mean Gaussian system and the alternative system, respectively when  $\mathbf{X}^*$  is the input of both systems. Then

$$C \ge I(\mathbf{X}^*; \mathbf{Y}) \ge I(\mathbf{X}^*; \mathbf{Y}^*) = C^*,$$
 (11.306)

where the second inequality follows from (11.288) in the proof of Theorem 11.32. Hence, we conclude that the zero-mean Gaussian noise is indeed the worst additive noise subject to a constraint on the correlation matrix.

# Problems

In the following, X, Y, Z, etc denote vectors of random variables.

- 288 11 Continuous-Valued Channels
- 1. Verify the two properties in Theorem 11.8 for the capacity of the memoryless Gaussian channel.
- 2. Let X and Y be two jointly distributed random variables with Y being continuous. The random variable Y is related to the random variable X through a conditional pdf f(y|x) defined for all x (cf. Definition 10.22). Prove that I(X;Y) is concave in F(x).
- 3. Refer to Lemma 11.18 and prove that

$$\Pr\{(\mathbf{X}', \mathbf{Y}') \in U^n_{[XY]\delta}\} \ge (1-\delta)2^{-n(I(X;Y)-\delta)}$$

for n sufficiently large.

4. Show that the capacity of a continuous memoryless channel is not changed if (11.23) is replaced by

$$E\left[\frac{1}{n}\sum_{i=1}^{n}\kappa(x_i(W))\right] \le P_i$$

i.e., the average input constraint is satisfied on the average by a randomly selected codeword instead of by every codeword in the codebook.

- 5. Show that  $R_{ii'}(0)$  in (11.233) vanishes if and only if  $i \neq i'$ .
- 6. Let  $\mathbf{Y} = \mathbf{X} + \mathbf{Z}$ , where  $\mathbf{Z}$  is independent of  $\mathbf{X}$ . Show that  $K_{\mathbf{Y}} = K_{\mathbf{X}} + K_{\mathbf{Z}}$ . Note that unlike Lemma 11.33, it is not necessary to assume that either  $\mathbf{X}$  or  $\mathbf{Z}$  has zero mean.
- 7. Consider a system of Gaussian channels with noise vector  $\mathbf{Z} \sim (0, K_{\mathbf{Z}})$  and input power constraint equal to 3. Determine the capacity of the system for the following two cases: a)

$$K_{\mathbf{Z}} = \begin{bmatrix} 4 \ 0 \ 0 \\ 0 \ 5 \ 0 \\ 0 \ 0 \ 2 \end{bmatrix};$$

b)

$$K_{\mathbf{Z}} = \begin{bmatrix} 7/4 & \sqrt{2}/4 & -3/4 \\ \sqrt{2}/4 & 5/2 & -\sqrt{2}/4 \\ -3/4 & -\sqrt{2}/4 & 7/4 \end{bmatrix}.$$

For b), you may use the results in Problem 4 in Chapter 10.

- 8. In the system of correlated Gaussian channels, let  $K_{\mathbf{Z}}$  be diagonalizable as  $QAQ^{\top}$ . Let  $A^*$  be the  $k \times k$  diagonal matrix with  $a_i^*$  being the *i*th diagonal element, where  $a_i^*$  is prescribed in (11.192). Show that  $\mathcal{N}(0, QA^*Q^{\top})$  is the optimal input distribution.
- 9. Show that for a wide-sense stationary process X(t),  $S_X(f) = S_X(-f)$  for all f.

- 10. Consider a zero-mean white Gaussian noise process Z(t). Let  $h_1(t)$  and  $h_2(t)$  be two impulse responses such that the supports of  $H_1(f)$  and  $H_2(f)$  do not overlap.
  - a) Show that for any  $t \neq t'$ , the two random variables  $Z(t) * h_1(t)$  and  $Z(t) * h_2(t)$  are independent.
  - b) Show that the two processes  $Z(t) * h_1(t)$  and  $Z(t) * h_2(t)$  are independent.
  - c) Repeat a) and b) if Z(t) is a zero-mean colored Gaussian noise process. Hint: Regard Z(t) as obtained by passing a zero-mean white Gaussian noise process through a coloring filter.
- 11. Interpret the bandpass white Gaussian channel as a special case of the bandlimited colored Gaussian channel in terms of the channel capacity.
- 12. Independent Gaussian noise is the worst Let C be the capacity of a system of k Gaussian channels with  $Z_i \sim \mathcal{N}(0, N_i)$ . By ignoring the possible correlation among the noise variables, we can use the channels in the system independently as parallel Gaussian channels. Thus C is lower bounded by the expression in (11.179). In this sense, a Gaussian noise vector is the worst if its components are uncorrelated. Justify this claim analytically. Hint: Show that  $I(\mathbf{X}; \mathbf{Y}) \geq \sum_i I(X_i; Y_i)$  if  $X_i$  are independent.

# **Historical Notes**

Channels with additive Gaussian noise were first analyzed by Shannon in [291], where the formula for the capacity of the bandlimited white Gaussian channel was given. A rigorous proof of this capacity formula was obtained by Wyner [348]. The water-filling solution to the capacity of the bandlimited colored Gaussian channel was developed by Shannon in [293] and was proved rigorously by Pinsker [263]. The discussion in this chapter on the continuous memoryless channel with an average input constraint is adapted from the discussions in the book by Gallager [117] and the book by Ihara [161], where in the former a comprehensive treatment of waveform channels can also be found. The Gaussian noise being the worst additive noise was proved by Ihara [160]. The proof presented here is based on Diggavi and Cover [84].

We have proved in Section 3.5 that if  $X_1 \to X_2 \to X_3 \to X_4$  forms a Markov chain, the *I*-Measure  $\mu^*$  always vanishes on the five atoms

$$\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3 \cap \tilde{X}_4^c$$

$$\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3 \cap \tilde{X}_4$$

$$\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3^c \cap \tilde{X}_4$$

$$\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3^c \cap \tilde{X}_4$$

$$\tilde{X}_1^c \cap \tilde{X}_2 \cap \tilde{X}_3^c \cap \tilde{X}_4.$$
(12.1)

Consequently, the *I*-Measure  $\mu^*$  is completely specified by the values of  $\mu^*$  on the other ten nonempty atoms of  $\mathcal{F}_4$ , and the information diagram for four random variables forming a Markov chain can be displayed in two dimensions as in Figure 3.11.

Figure 12.1 is a graph which represents the Markov chain  $X_1 \to X_2 \to X_3 \to X_4$ . The observant reader would notice that  $\mu^*$  always vanishes on a



**Fig. 12.1.** The graph representing the Markov chain  $X_1 \to X_2 \to X_3 \to X_4$ .

nonempty atom A of  $\mathcal{F}_4$  if and only if the graph in Figure 12.1 becomes disconnected upon removing all the vertices corresponding to the complemented set variables in A. For example,  $\mu^*$  always vanishes on the atom  $\tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3 \cap \tilde{X}_4^c$ , and the graph in Figure 12.1 becomes disconnected upon removing vertices 2 and 4. On the other hand,  $\mu^*$  does not necessarily vanish on the atom  $\tilde{X}_1^c \cap \tilde{X}_2 \cap \tilde{X}_3 \cap \tilde{X}_4^c$ , and the graph in Figure 12.1 remains connected upon removing vertices 1 and 4. This observation will be explained in a more general setting in the subsequent sections.

The theory of *I*-Measure establishes a one-to-one correspondence between Shannon's information measures and set theory. Based on this theory, we develop in this chapter a set-theoretic characterization of a Markov structure called *full conditional mutual independence*. A Markov chain, and more generally a Markov random field, is a collection of full conditional mutual independencies. We will show that if a collection of random variables forms a Markov random field, then the structure of  $\mu^*$  can be simplified. In particular, when the random variables form a Markov chain,  $\mu^*$  exhibits a very simple structure so that the information diagram can be displayed in two dimensions regardless of the length of the Markov chain, and  $\mu^*$  is always nonnegative. (See also Sections 3.5 and 3.6.)

The topics to be covered in this chapter are fundamental. Unfortunately, the proofs of the results are very heavy. The reader may skip this chapter without affecting further reading of the book.

### 12.1 Conditional Mutual Independence

In this section, we explore the effect of conditional mutual independence on the structure of the *I*-Measure  $\mu^*$ . We begin with a simple example.

*Example 12.1.* Let X, Y, and Z be mutually independent random variables. Then

$$I(X;Y) = I(X;Y;Z) + I(X;Y|Z) = 0.$$
(12.2)

Since  $I(X; Y|Z) \ge 0$ , we let

$$I(X;Y|Z) = a \ge 0,$$
 (12.3)

so that

$$I(X;Y;Z) = -a.$$
 (12.4)

Similarly,

$$I(Y;Z) = I(X;Y;Z) + I(Y;Z|X) = 0$$
(12.5)

and

$$I(X;Z) = I(X;Y;Z) + I(X;Z|Y) = 0.$$
(12.6)

Then from (12.4), we obtain

$$I(Y;Z|X) = I(X;Z|Y) = a.$$
 (12.7)

The relations (12.3), (12.4), and (12.7) are shown in the information diagram in Figure 12.2, which indicates that X, Y, and Z are pairwise independent.

We have proved in Theorem 2.39 that X, Y, and Z are mutually independent if and only if

$$H(X, Y, Z) = H(X) + H(Y) + H(Z).$$
(12.8)

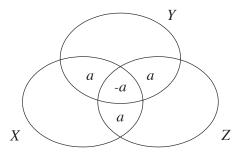


Fig. 12.2. X, Y, and Z are pairwise independent.

By counting atoms in the information diagram, we see that

$$0 = H(X) + H(Y) + H(Z) - H(X, Y, Z)$$
(12.9)

$$= I(X;Y|Z) + I(Y;Z|X) + I(X;Z|Y) + 2I(X;Y;Z)$$
(12.10)

$$= a.$$
 (12.11)

Thus a = 0, which implies

$$I(X;Y|Z), I(Y;Z|X), I(X;Z|Y), I(X;Y;Z)$$
(12.12)

are all equal to 0. Equivalently,  $\mu^*$  vanishes on

$$\tilde{X} \cap \tilde{Y} - \tilde{Z}, \tilde{Y} \cap \tilde{Z} - \tilde{X}, \tilde{X} \cap \tilde{Z} - \tilde{Y}, \tilde{X} \cap \tilde{Y} \cap \tilde{Z},$$
(12.13)

which are precisely the atoms in the intersection of any two of the set variables  $\tilde{X}, \tilde{Y}$ , and  $\tilde{Z}$ .

Conversely, if  $\mu^*$  vanishes on the sets in (12.13), then we see from (12.10) that (12.8) holds, i.e., X, Y, and Z are mutually independent. Therefore, X, Y, and Z are mutually independent if and only if  $\mu^*$  vanishes on the sets in (12.13). This is shown in the information diagram in Figure 12.3.

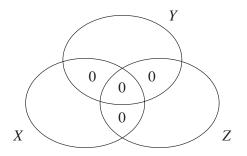


Fig. 12.3. X, Y, and Z are mutually independent.

The theme of this example will be extended to conditional mutual independence among collections of random variables in Theorem 12.9, which is the main result in this section. In the rest of the section, we will develop the necessary tools for proving this theorem. At first reading, the reader should try to understand the results by studying the examples without getting into the details of the proofs.

In Theorem 2.39, we have proved that  $X_1, X_2, \dots, X_n$  are mutually independent if and only if

$$H(X_1, X_2, \cdots, X_n) = \sum_{i=1}^n H(X_i).$$
 (12.14)

By conditioning on a random variable Y, one can readily prove the following.

**Theorem 12.2.**  $X_1, X_2, \dots, X_n$  are mutually independent conditioning on Y if and only if

$$H(X_1, X_2, \cdots, X_n | Y) = \sum_{i=1}^n H(X_i | Y).$$
(12.15)

We now prove two alternative characterizations of conditional mutual independence.

**Theorem 12.3.**  $X_1, X_2, \dots, X_n$  are mutually independent conditioning on Y if and only if for all  $1 \le i \le n$ ,

$$I(X_i; X_j, j \neq i | Y) = 0, \qquad (12.16)$$

*i.e.*,  $X_i$  and  $(X_j, j \neq i)$  are independent conditioning on Y.

**Remark** A conditional independency is a special case of a conditional mutual independency. However, this theorem says that a conditional mutual independency is equivalent to a set of conditional independencies.

Proof of Theorem 12.3. It suffices to prove that (12.15) and (12.16) are equivalent. Assume (12.15) is true, so that  $X_1, X_2, \dots, X_n$  are mutually independent conditioning on Y. Then for all  $i, X_i$  is independent of  $(X_j, j \neq i)$  conditioning on Y. This proves (12.16).

Now assume that (12.16) is true for all  $1 \le i \le n$ . Consider

$$0 = I(X_i; X_j, j \neq i | Y)$$

$$= I(X_i; X_1, X_2, \cdots, X_{i-1} | Y)$$
(12.17)

$$+I(X_i; X_{i+1}, \cdots, X_n | Y, X_1, X_2, \cdots, X_{i-1}).$$
(12.18)

Since mutual information is always nonnegative, this implies

$$I(X_i; X_1, \cdots, X_{i-1} | Y) = 0, (12.19)$$

or  $X_i$  and  $(X_1, X_2, \dots, X_{i-1})$  are independent conditioning on Y. Therefore,  $X_1, X_2, \dots, X_n$  are mutually independent conditioning on Y (see the proof of Theorem 2.39), proving (12.15). Hence, the theorem is proved.  $\Box$  **Theorem 12.4.**  $X_1, X_2, \dots, X_n$  are mutually independent conditioning on Y if and only if

$$H(X_1, X_2, \cdots, X_n | Y) = \sum_{i=1}^n H(X_i | Y, X_j, j \neq i).$$
(12.20)

*Proof.* It suffices to prove that (12.15) and (12.20) are equivalent. Assume (12.15) is true, so that  $X_1, X_2, \dots, X_n$  are mutually independent conditioning on Y. Since for all  $i, X_i$  is independent of  $X_j, j \neq i$  conditioning on Y,

$$H(X_i|Y) = H(X_i|Y, X_j, j \neq i)$$
(12.21)

Therefore, (12.15) implies (12.20).

Now assume that (12.20) is true. Consider

$$H(X_{1}, X_{2}, \dots, X_{n}|Y) = \sum_{i=1}^{n} H(X_{i}|Y, X_{1}, \dots, X_{i-1})$$
(12.22)  
$$= \sum_{i=1}^{n} [H(X_{i}|Y, X_{j}, j \neq i) + I(X_{i}; X_{i+1}, \dots, X_{n}|Y, X_{1}, \dots, X_{i-1})]$$
(12.23)  
$$= \sum_{i=1}^{n} H(X_{i}|Y, X_{j}, j \neq i) + \sum_{i=1}^{n} I(X_{i}; X_{i+1}, \dots, X_{n}|Y, X_{1}, \dots, X_{i-1}).$$
(12.24)

Then (12.20) implies

$$\sum_{i=1}^{n} I(X_i; X_{i+1}, \cdots, X_n | Y, X_1, \cdots, X_{i-1}) = 0.$$
 (12.25)

Since all the terms in the above summation are nonnegative, they must all be equal to 0. In particular, for i = 1, we have

$$I(X_1; X_2, \cdots, X_n | Y) = 0.$$
(12.26)

By symmetry, it can be shown that

$$I(X_i; X_j, j \neq i | Y) = 0$$
(12.27)

for all  $1 \le i \le n$ . Then this implies (12.15) by the last theorem, completing the proof.  $\Box$ 

**Theorem 12.5.** Let C and  $Q_i$  be disjoint index sets and  $W_i$  be a subset of  $Q_i$  for  $1 \le i \le k$ , where  $k \ge 2$ . Assume that there exist at least two i such that  $W_i \ne \emptyset$ . Let  $X_{Q_i} = (X_l, l \in Q_i), 1 \le i \le k$ , and  $X_C = (X_l, l \in C)$  be collections of random variables. If  $X_{Q_i}, 1 \le i \le k$ , are mutually independent conditioning on  $X_C$ , then  $X_{W_i}$  such that  $W_i \ne \emptyset$  are mutually independent conditioning on  $(X_C, X_{Q_i-W_i}, 1 \le i \le k)$ .

We first give an example before we prove the theorem.

*Example 12.6.* Suppose  $X_1$ ,  $(X_2, X_3, X_4)$ , and  $(X_5, X_6)$  are mutually independent conditioning on  $X_7$ . By Theorem 12.5,  $X_1$ ,  $X_2$ , and  $(X_5, X_6)$  are mutually independent conditioning on  $(X_3, X_4, X_7)$ .

Proof of Theorem 12.5. Assume  $X_{Q_i}, 1 \leq i \leq k$ , are mutually independent conditioning on  $X_C$ , i.e.,

$$H(X_{Q_i}, 1 \le i \le k | X_C) = \sum_{i=1}^k H(X_{Q_i} | X_C).$$
(12.28)

Consider

$$H(X_{W_i}, 1 \le i \le k | X_C, X_{Q_i - W_i}, 1 \le i \le k) = H(X_{Q_i}, 1 \le i \le k | X_C) - H(X_{Q_i - W_i}, 1 \le i \le k | X_C)$$
(12.29)

$$= \sum_{i=1}^{k} H(X_{Q_i}|X_C) - \sum_{i=1}^{k} H(X_{Q_i-W_i}|X_C, X_{Q_j-W_j}, 1 \le j \le i-1)$$
(12.30)

$$\geq \sum_{i=1}^{k} H(X_{Q_i}|X_C, X_{Q_j - W_j}, 1 \leq j \leq i - 1) \\ - \sum_{i=1}^{k} H(X_{Q_i - W_i}|X_C, X_{Q_j - W_j}, 1 \leq j \leq i - 1)$$
(12.31)

$$=\sum_{i=1}^{k} H(X_{W_i}|X_C, X_{Q_j - W_j}, 1 \le j \le i)$$
(12.32)

$$\geq \sum_{i=1}^{k} H(X_{W_i}|X_C, X_{Q_j - W_j}, 1 \leq j \leq k).$$
(12.33)

In the second step we have used (12.28), and the two inequalities follow because conditioning does not increase entropy. On the other hand, by the chain rule for entropy, we have

$$H(X_{W_i}, 1 \le i \le k | X_C, X_{Q_i - W_i}, 1 \le i \le k)$$
  
=  $\sum_{i=1}^k H(X_{W_i} | X_C, (X_{Q_j - W_j}, 1 \le j \le k), (X_{W_l}, 1 \le l \le i - 1)).$   
(12.34)

Therefore, it follows from (12.33) that

$$\sum_{i=1}^{k} H(X_{W_i}|X_C, X_{Q_j - W_j}, 1 \le j \le k)$$
(12.35)

$$\leq H(X_{W_i}, 1 \leq i \leq k | X_C, X_{Q_i - W_i}, 1 \leq i \leq k)$$
(12.36)

$$=\sum_{i=1}^{n} H(X_{W_i}|X_C, (X_{Q_j-W_j}, 1 \le j \le k), (X_{W_l}, 1 \le l \le i-1)).$$
(12.37)

However, since conditioning does not increase entropy, the *i*th term in the summation in (12.35) is lower bounded by the *i*th term in the summation in (12.37). Thus we conclude that the inequality in (12.36) is an equality. Hence, the conditional entropy in (12.36) is equal to the summation in (12.35), i.e.,

$$H(X_{W_i}, 1 \le i \le k | X_C, X_{Q_i - W_i}, 1 \le i \le k)$$
(12.38)

$$=\sum_{i=1}^{n} H(X_{W_i}|X_C, X_{Q_j-W_j}, 1 \le j \le k).$$
(12.39)

The theorem is proved.  $\Box$ 

Theorem 12.5 specifies a set of conditional mutual independencies (CMI's) which is implied by a CMI. This theorem is crucial for understanding the effect of a CMI on the structure of the *I*-Measure  $\mu^*$ , which we discuss next.

**Lemma 12.7.** Let  $(Z_{i1}, \dots, Z_{it_i}), 1 \leq i \leq r$ , be r collections of random variables, where  $r \geq 2$ , and let Y be a random variable, such that  $(Z_{i1}, \dots, Z_{it_i}), 1 \leq i \leq r$ , are mutually independent conditioning on Y. Then

$$\mu^* \left( \bigcap_{i=1}^r \bigcap_{j=1}^{t_i} \tilde{Z}_{ij} - \tilde{Y} \right) = 0.$$
 (12.40)

We first prove the following set identity which will be used in proving this lemma.

**Lemma 12.8.** Let S and T be disjoint index sets, and  $A_i$  and B be sets. Let  $\mu$  be a set-additive function. Then

$$\mu\left(\left(\bigcap_{i\in S} A_i\right) \cap \left(\bigcap_{j\in T} A_j\right) - B\right)$$
$$= \sum_{S'\subset S} \sum_{T'\subset T} (-1)^{|S'|+|T'|} \left(\mu(A_{S'}-B) + \mu(A_{T'}-B) - \mu(A_{S'\cup T'}-B)\right),$$
(12.41)

where  $A_{S'}$  denotes  $\cup_{i \in S'} A_i$ .

*Proof.* The right hand side of (12.41) is equal to

$$\sum_{S' \subset S} \sum_{T' \subset T} (-1)^{|S'| + |T'|} \mu(A_{S'} - B) + \sum_{S' \subset S} \sum_{T' \subset T} (-1)^{|S'| + |T'|} \mu(A_{T'} - B)$$
$$- \sum_{S' \subset S} \sum_{T' \subset T} (-1)^{|S'| + |T'|} \mu(A_{S' \cup T'} - B).$$
(12.42)

Now

$$\sum_{S' \subset S} \sum_{T' \subset T} (-1)^{|S'| + |T'|} \mu(A_{S'} - B) = \sum_{S' \subset S} (-1)^{|S'|} \mu(A_{S'} - B) \sum_{T' \subset T} (-1)^{|T'|}.$$
(12.43)

Since

$$\sum_{T' \subset T} (-1)^{|T'|} = \sum_{k=0}^{|T|} {|T| \choose k} (-1)^k = 0$$
(12.44)

by the binomial formula<sup>1</sup>, we conclude that

$$\sum_{S' \subset S} \sum_{T' \subset T} (-1)^{|S'| + |T'|} \mu(A_{S'} - B) = 0.$$
 (12.45)

Similarly,

$$\sum_{S' \subseteq S} \sum_{T' \subseteq T} (-1)^{|S'| + |T'|} \mu(A_{T'} - B) = 0.$$
(12.46)

Therefore, (12.41) is equivalent to

$$\mu\left(\left(\bigcap_{i\in S} A_i\right) \cap \left(\bigcap_{j\in T} A_j\right) - B\right) = \sum_{S'\subset S} \sum_{T'\subset T} (-1)^{|S'| + |T'| + 1} \mu(A_{S'\cup T'} - B)$$
(12.47)

<sup>1</sup> This can be obtained by letting a = 1 and b = -1 in the binomial formula

$$(a+b)^{|T|} = \sum_{k=0}^{|T|} {|T| \choose k} a^k b^{|T|-k}.$$

which can readily be obtained from Theorem 3.142. Hence, the lemma is proved.  $\ \square$ 

*Proof of Lemma 12.7.* We first prove the lemma for r = 2. By Lemma 12.8,

$$\mu^* \left( \bigcap_{i=1}^2 \bigcap_{j=1}^{t_i} \tilde{Z}_{ij} - \tilde{Y} \right) =$$

$$\sum_{S' \subset \{1, \cdots, t_1\}} \sum_{T' \subset \{1, \cdots, t_2\}} (-1)^{|S'| + |T'|} \left[ \mu^* \left( \bigcup_{j \in S'} \tilde{Z}_{1j} - \tilde{Y} \right) + \mu^* \left( \bigcup_{k \in T'} \tilde{Z}_{2k} - \tilde{Y} \right) - \mu^* \left( \left( \bigcup_{j \in S'} \tilde{Z}_{1j} \right) \cup \left( \bigcup_{k \in T'} \tilde{Z}_{2k} \right) - \tilde{Y} \right) \right]. \quad (12.48)$$

The expression in the square bracket is equal to

$$H(Z_{1j}, j \in S'|Y) + H(Z_{2k}, k \in T'|Y)$$
  
-H((Z<sub>1j</sub>, j \in S'), (Z<sub>2k</sub>, k \in T')|Y), (12.49)

which is equal to 0 because  $(Z_{1j}, j \in S')$  and  $(Z_{2k}, k \in T')$  are independent conditioning on Y. Therefore the lemma is proved for r = 2.

For r > 2, we write

$$\mu^* \left( \bigcap_{i=1}^r \bigcap_{j=1}^{t_i} \tilde{Z}_{ij} - \tilde{Y} \right) = \mu^* \left( \left( \bigcap_{i=1}^{r-1} \bigcap_{j=1}^{t_i} \tilde{Z}_{ij} \right) \cap \left( \bigcap_{j=1}^{t_r} \tilde{Z}_{rj} \right) - \tilde{Y} \right). \quad (12.50)$$

Since  $((Z_{i1}, \dots, Z_{it_i}), 1 \leq i \leq r-1)$  and  $(Z_{r1}, \dots, Z_{rt_r})$  are independent conditioning on Y, upon applying the lemma for r = 2, we see that

$$\mu^* \left( \bigcap_{i=1}^r \bigcap_{j=1}^{t_i} \tilde{Z}_{ij} - \tilde{Y} \right) = 0.$$
 (12.51)

The lemma is proved.  $\Box$ 

**Theorem 12.9.** Let T and  $Q_i, 1 \le i \le k$ , be disjoint index sets, where  $k \ge 2$ , and let  $X_{Q_i} = (X_l, l \in Q_i), 1 \le i \le k$ , and  $X_T = (X_l, l \in T)$  be collections of random variables. Then  $X_{Q_i}, 1 \le i \le k$ , are mutually independent conditioning on  $X_T$  if and only if for any  $W_1, W_2, \dots, W_k$ , where  $W_i \subset Q_i, 1 \le i \le k$ , if there exist at least two i such that  $W_i \ne \emptyset$ , then

$$\mu^*\left(\left(\bigcap_{i=1}^k\bigcap_{j\in W_i}\tilde{X}_j\right) - \tilde{X}_{T\cup(\bigcup_{i=1}^k(Q_i - W_i))}\right) = 0.$$
(12.52)

We first give an example before proving this fundamental result. The reader should compare this example with Example 12.6.

*Example 12.10.* Suppose  $X_1, (X_2, X_3, X_4)$ , and  $(X_5, X_6)$  are mutually independent conditioning on  $X_7$ . By Theorem 12.9,

$$\mu^* (\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_5 \cap \tilde{X}_6 - (\tilde{X}_3 \cup \tilde{X}_4 \cup \tilde{X}_7)) = 0.$$
 (12.53)

However, the theorem does not say, for instance, that

$$\mu^* (\tilde{X}_2 \cap \tilde{X}_4 - (\tilde{X}_1 \cup \tilde{X}_3 \cup \tilde{X}_5 \cup \tilde{X}_6 \cup \tilde{X}_7))$$
(12.54)

is equal to 0.

Proof of Theorem 12.9. We first prove the 'if' part. Assume that for any  $W_1, W_2, \dots, W_k$ , where  $W_i \subset Q_i, 1 \leq i \leq k$ , if there exist at least two *i* such that  $W_i \neq \emptyset$ , then (12.52) holds. Consider

$$H(X_{Q_i}, 1 \le i \le k | X_T) = \mu^* \left( \tilde{X}_{\bigcup_{i=1}^k Q_i} - \tilde{X}_T \right)$$
(12.55)

$$=\sum_{B\in S}\mu^{*}(B),$$
 (12.56)

where S consists of sets of the form

$$\left(\bigcap_{i=1}^{k}\bigcap_{j\in W_{i}}\tilde{X}_{j}\right) - \tilde{X}_{T\cup\left(\bigcup_{i=1}^{k}(Q_{i}-W_{i})\right)}$$
(12.57)

with  $W_i \subset Q_i$  for  $1 \leq i \leq k$  and there exists at least one *i* such that  $W_i \neq \emptyset$ . By our assumption, if  $B \in S$  is such that there exist at least two *i* for which  $W_i \neq \emptyset$ , then  $\mu^*(B) = 0$ . Therefore, if  $\mu^*(B)$  is possibly nonzero, then *B* must be such that there exists a unique *i* for which  $W_i \neq \emptyset$ . Now for  $1 \leq i \leq k$ , let  $S_l$  be the set consisting of sets of the form in (12.57) with  $W_i \subset Q_i, W_i \neq \emptyset$ , and  $W_l = \emptyset$  for  $l \neq i$ . In other words,  $S_i$  consists of atoms of the form

$$\left(\bigcap_{j\in W_i} \tilde{X}_j\right) - \tilde{X}_{T\cup(\cup_{l\neq i}Q_l)\cup(Q_i - W_i)}$$
(12.58)

with  $W_i \subset Q_i$  and  $W_i \neq \emptyset$ . Then

$$\sum_{B \in S} \mu^*(B) = \sum_{i=1}^k \sum_{B \in S_i} \mu^*(B).$$
(12.59)

Now

$$X_{Q_i} - X_{T \cup (\bigcup_{l \neq i} Q_l)} = \bigcup_{\substack{W_i \subset Q_i \\ W_i \neq \emptyset}} \left[ \left( \bigcap_{j \in W_i} \tilde{X}_j \right) - \tilde{X}_{T \cup (\bigcup_{l \neq i} Q_l) \cup (Q_i - W_i)} \right]$$
(12.60)

$$=\bigcup_{B\in S_i} B.$$
(12.61)

Since  $\mu^*$  is set-additive, we have

$$\mu^* \left( \tilde{X}_{Q_i} - \tilde{X}_{T \cup (\cup_{l \neq i} Q_l)} \right) = \sum_{B \in S_i} \mu^*(B).$$
 (12.62)

Hence, from (12.56) and (12.59), we have

$$H(X_{Q_i}, 1 \le i \le k | X_T)$$
  
=  $\sum_{i=1}^k \sum_{B \in S_i} \mu^*(B)$  (12.63)

$$=\sum_{i=1}^{k} \mu^{*} \left( \tilde{X}_{Q_{i}} - \tilde{X}_{T \cup (\cup_{l \neq i} Q_{l})} \right)$$
(12.64)

$$=\sum_{i=1}^{k} H(X_{Q_i}|X_T, X_{Q_l}, l \neq i), \qquad (12.65)$$

where (12.64) follows from (12.62). By Theorem 12.4,  $X_{Q_i}, 1 \leq i \leq k$ , are mutually independent conditioning on  $X_T$ .

We now prove the 'only if' part. Assume  $X_{Q_i}, 1 \leq i \leq k$ , are mutually independent conditioning on  $X_T$ . For any collection of sets  $W_1, W_2, \dots, W_k$ , where  $W_i \subset Q_i, 1 \leq i \leq k$ , if there exist at least two *i* such that  $W_i \neq \emptyset$ , by Theorem 12.5,  $X_{W_i}, 1 \leq i \leq k$ , are mutually independent conditioning on  $(X_T, X_{Q_i - W_i}, 1 \leq i \leq k)$ . By Lemma 12.7, we obtain (12.52). The theorem is proved.  $\Box$ 

# 12.2 Full Conditional Mutual Independence

**Definition 12.11.** A conditional mutual independency on  $X_1, X_2, \dots, X_n$  is full if all  $X_1, X_2, \dots, X_n$  are involved. Such a conditional mutual independency is called a full conditional mutual independency (FCMI).

*Example 12.12.* For n = 5,

 $X_1, X_2, X_4$ , and  $X_5$  are mutually independent conditioning on  $X_3$ 

is an FCMI. However,

 $X_1, X_2$ , and  $X_5$  are mutually independent conditioning on  $X_3$  is not an FCMI because  $X_4$  is not involved.

As in the previous chapters, we let

$$\mathcal{N}_n = \{1, 2, \cdots, n\}.$$
 (12.66)

In Theorem 12.9, if

$$T \cup \left(\bigcup_{i=1}^{k} Q_i\right) = \mathcal{N}_n, \qquad (12.67)$$

then the tuple  $(T, Q_i, 1 \le i \le k)$  defines the following FCMI on  $X_1, X_2, \cdots, X_n$ :

 $K: X_{Q_1}, X_{Q_2}, \cdots, X_{Q_k}$  are mutually independent conditioning on  $X_T$ .

We will denote K by  $(T, Q_i, 1 \le i \le k)$ .

**Definition 12.13.** Let  $K = (T, Q_i, 1 \le i \le k)$  be an FCMI on  $X_1, X_2, \dots, X_n$ . The image of K, denoted by Im(K), is the set of all atoms of  $\mathcal{F}_n$  which has the form of the set in (12.57), where  $W_i \subset Q_i, 1 \le i \le k$ , and there exist at least two i such that  $W_i \ne \emptyset$ .

**Proposition 12.14.** Let  $K = (T, Q_1, Q_2)$  be an FCI (full conditional independency) on  $X_1, X_2, \dots, X_n$ . Then

$$Im(K) = \{ A \in \mathcal{A} : A \subset (\tilde{X}_{Q_1} \cap \tilde{X}_{Q_2} - \tilde{X}_T) \}.$$
 (12.68)

**Proposition 12.15.** Let  $K = (T, Q_i, 1 \le i \le k)$  be an FCMI on  $X_1, X_2, \cdots, X_n$ . Then

$$Im(K) = \left\{ A \in \mathcal{A} : A \subset \bigcup_{1 \le i < j \le k} (\tilde{X}_{Q_i} \cap \tilde{X}_{Q_j} - \tilde{X}_T) \right\}.$$
 (12.69)

These two propositions greatly simplify the description of Im(K). Their proofs are elementary and they are left as exercises. We first illustrate these two propositions in the following example.

*Example 12.16.* Consider n = 4 and FCMI's  $K_1 = (\{3\}, \{1\}, \{2, 4\})$  and  $K_2 = (\emptyset, \{1\}, \{2, 3\}, \{4\})$ . Then

$$Im(K_1) = \{ A \in \mathcal{A} : A \subset (\tilde{X}_1 \cap \tilde{X}_{\{2,4\}} - \tilde{X}_3) \}$$
(12.70)

and

$$Im(K_2) = \{A \in \mathcal{A} : A \subset (\tilde{X}_1 \cap \tilde{X}_{\{2,3\}}) \cup (\tilde{X}_{\{2,3\}} \cap \tilde{X}_4) \cup (\tilde{X}_1 \cap \tilde{X}_4)\}.$$
(12.71)

**Theorem 12.17.** Let K be an FCMI on  $X_1, X_2, \dots, X_n$ . Then K holds if and only if  $\mu^*(A) = 0$  for all  $A \in Im(K)$ .

*Proof.* First, (12.67) is true if K is an FCMI. Then the set in (12.57) can be written as

$$\left(\bigcap_{j\in\cup_{i=1}^{k}W_{i}}\tilde{X}_{j}\right)-\tilde{X}_{\mathcal{N}_{n}-\cup_{i=1}^{k}W_{i}},$$
(12.72)

which is seen to be an atom of  $\mathcal{F}_n$ . The theorem can then be proved by a direct application of Theorem 12.9 to the FCMI K.  $\Box$ 

Let  $A = \bigcap_{i=1}^{n} \tilde{Y}_i$  be a nonempty atom of  $\mathcal{F}_n$ . Define the set

$$U_A = \{i \in \mathcal{N}_n : \tilde{Y}_i = \tilde{X}_i^c\}.$$
(12.73)

Note that A is uniquely specified by  $U_A$  because

$$A = \left(\bigcap_{i \in \mathcal{N}_n - U_A} \tilde{X}_i\right) \cap \left(\bigcap_{i \in U_A} \tilde{X}_i^c\right) = \left(\bigcap_{i \in \mathcal{N}_n - U_A} \tilde{X}_i\right) - \tilde{X}_{U_A}.$$
 (12.74)

Define  $w(A) = n - |U_A|$  as the *weight* of the atom A, the number of  $\tilde{X}_i$  in A which are not complemented. We now show that an FCMI  $K = (T, Q_i, 1 \le i \le k)$  is uniquely specified by Im(K). First, by letting  $W_i = Q_i$  for  $1 \le i \le k$  in Definition 12.13, we see that the atom

$$\left(\bigcap_{j\in\cup_{i=1}^{k}Q_{i}}\tilde{X}_{j}\right)-\tilde{X}_{T}$$
(12.75)

is in Im(K), and it is the unique atom in Im(K) with the largest weight. From this atom, T can be determined. To determine  $Q_i, 1 \le i \le k$ , we define a relation q on  $T^c = \mathcal{N}_n \setminus T$  as follows. For  $l, l' \in T^c$ , (l, l') is in q if and only if

- i) l = l'; or
- ii) there exists an atom of the form

$$\tilde{X}_l \cap \tilde{X}_{l'} \cap \bigcap_{\substack{1 \le j \le n \\ j \ne l, l'}} \tilde{Y}_j \tag{12.76}$$

in  $\mathcal{A} - Im(K)$ , where  $\tilde{Y}_j = \tilde{X}_j$  or  $\tilde{X}_j^c$ .

Recall that  $\mathcal{A}$  is the set of nonempty atoms of  $\mathcal{F}_n$ . The idea of ii) is that (l, l') is in q if and only if  $l, l' \in Q_i$  for some  $1 \leq i \leq k$ . Then q is reflexive and symmetric by construction, and is transitive by virtue of the structure of Im(K). In other words, q is an *equivalence relation* which partitions  $T^c$  into  $\{Q_i, 1 \leq i \leq k\}$ . Therefore, K and Im(K) uniquely specify each other.

The image of an FCMI K completely characterizes the effect of K on the *I*-Measure for  $X_1, X_2, \dots, X_n$ . The joint effect of more than one FCMI can easily be described in terms of the images of the individual FCMI's. Let

$$\Pi = \{K_l, 1 \le l \le m\}$$
(12.77)

be a set of FCMI's. By Theorem 12.9,  $K_l$  holds if and only if  $\mu^*$  vanishes on the atoms in  $Im(K_l)$ . Then  $K_l, 1 \leq l \leq m$  hold simultaneously if and only if  $\mu^*$  vanishes on the atoms in  $\cup_{l=1}^k Im(K_l)$ . This is summarized as follows.

**Definition 12.18.** The image of a set of FCMI's  $\Pi = \{K_l, 1 \leq l \leq m\}$  is defined as

$$Im(\Pi) = \bigcup_{l=1}^{k} Im(K_l).$$
 (12.78)

**Theorem 12.19.** Let  $\Pi$  be a set of FCMI's for  $X_1, X_2, \dots, X_n$ . Then  $\Pi$  holds if and only if  $\mu^*(A) = 0$  for all  $A \in Im(\Pi)$ .

In probability problems, we are often given a set of conditional independencies and we need to see whether another given conditional independency is logically implied. This is called the *implication problem* which will be discussed in detail in Section 13.5. The next theorem gives a solution to this problem if only FCMI's are involved.

**Theorem 12.20.** Let  $\Pi_1$  and  $\Pi_2$  be two sets of FCMI's. Then  $\Pi_1$  implies  $\Pi_2$  if and only if  $Im(\Pi_2) \subset Im(\Pi_1)$ .

Proof. We first prove that if  $Im(\Pi_2) \subset Im(\Pi_1)$ , then  $\Pi_1$  implies  $\Pi_2$ . Assume  $Im(\Pi_2) \subset Im(\Pi_1)$  and  $\Pi_1$  holds. Then by Theorem 12.19,  $\mu^*(A) = 0$  for all  $A \in Im(\Pi_1)$ . Since  $Im(\Pi_2) \subset Im(\Pi_1)$ , this implies that  $\mu^*(A) = 0$  for all  $A \in Im(\Pi_2)$ . Again by Theorem 12.19, this implies  $\Pi_2$  also holds. Therefore, if  $Im(\Pi_2) \subset Im(\Pi_1)$ , then  $\Pi_1$  implies  $\Pi_2$ .

We now prove that if  $\Pi_1$  implies  $\Pi_2$ , then  $Im(\Pi_2) \subset Im(\Pi_1)$ . To prove this, we assume that  $\Pi_1$  implies  $\Pi_2$  but  $Im(\Pi_2) \not\subset Im(\Pi_1)$ , and we will show that this leads to a contradiction. Fix a nonempty atom  $A \in Im(\Pi_2) - Im(\Pi_1)$ . By Theorem 3.11, we can construct random variables  $X_1, X_2, \dots, X_n$  such that  $\mu^*$  vanishes on all the atoms of  $\mathcal{F}_n$  except for A. Then  $\mu^*$  vanishes on all the atoms in  $Im(\Pi_1)$  but not on all the atoms in  $Im(\Pi_2)$ . By Theorem 12.19, this implies that for  $X_1, X_2, \dots, X_n$  so constructed,  $\Pi_1$  holds but  $\Pi_2$  does not hold. Therefore,  $\Pi_1$  does not imply  $\Pi_2$ , which is a contradiction. The theorem is proved.  $\Box$ 

**Remark** In the course of proving this theorem and all its preliminaries, we have used nothing more than the basic inequalities. Therefore, we have shown that the basic inequalities are a sufficient set of tools to solve the implication problem if only FCMI's are involved.

**Corollary 12.21.** Two sets of FCMI's are equivalent if and only if their images are identical.

*Proof.* Two set of FCMI's  $\Pi_1$  and  $\Pi_2$  are equivalent if and only if

$$\Pi_1 \Rightarrow \Pi_2 \quad \text{and} \quad \Pi_2 \Rightarrow \Pi_1.$$
(12.79)

Then by the last theorem, this is equivalent to  $Im(\Pi_2) \subset Im(\Pi_1)$  and  $Im(\Pi_1) \subset Im(\Pi_2)$ , i.e.,  $Im(\Pi_2) = Im(\Pi_1)$ . The corollary is proved.  $\Box$ 

Thus a set of FCMI's is completely characterized by its image. A set of FCMI's is a set of probabilistic constraints, but the characterization by its image is purely set-theoretic! This characterization offers an intuitive set-theoretic interpretation of the joint effect of FCMI's on the *I*-Measure for  $X_1, X_2, \dots, X_n$ . For example,  $Im(K_1) \cap Im(K_2)$  is interpreted as the effect commonly due to  $K_1$  and  $K_2$ ,  $Im(K_1) - Im(K_2)$  is interpreted as the effect due to  $K_1$  but not  $K_2$ , etc. We end this section with an example.

Example 12.22. Consider n = 4. Let

$$K_1 = (\emptyset, \{1, 2, 3\}, \{4\}), \quad K_2 = (\emptyset, \{1, 2, 4\}, \{3\})$$
(12.80)

$$K_3 = (\emptyset, \{1, 2\}, \{3, 4\}), \quad K_4 = (\emptyset, \{1, 3\}, \{2, 4\})$$
 (12.81)

and let  $\Pi_1 = \{K_1, K_2\}$  and  $\Pi_2 = \{K_3, K_4\}$ . Then

$$Im(\Pi_1) = Im(K_1) \cup Im(K_2)$$
 (12.82)

and

$$Im(\Pi_2) = Im(K_3) \cup Im(K_4),$$
 (12.83)

where

$$Im(K_1) = \{ A \in \mathcal{A} : A \subset (X_{\{1,2,3\}} \cap X_4) \}$$
(12.84)

$$Im(K_2) = \{A \in \mathcal{A} : A \subset (X_{\{1,2,4\}} \cap X_3)\}$$
(12.85)

$$Im(K_3) = \{A \in \mathcal{A} : A \subset (X_{\{1,2\}} \cap X_{\{3,4\}})\}$$
(12.86)

$$Im(K_4) = \{ A \in \mathcal{A} : A \subset (X_{\{1,3\}} \cap X_{\{2,4\}}) \}.$$
 (12.87)

It can readily be seen by using an information diagram that  $Im(\Pi_1) \subset Im(\Pi_2)$ . Therefore,  $\Pi_2$  implies  $\Pi_1$ . Note that no probabilistic argument is involved in this proof.

# 12.3 Markov Random Field

A Markov random field is a generalization of a discrete time Markov chain in the sense that the time index for the latter, regarded as a chain, is replaced by a general graph for the former. Historically, the study of Markov random field stems from statistical physics. The classical Ising model, which is defined on a rectangular lattice, was used to explain certain empirically observed facts about ferromagnetic materials. In this section, we explore the structure of the *I*-Measure for a Markov random field.

We refer the reader to textbooks on graph theory (e.g. [43]) for formal definitions of the graph-theoretic terminologies to be used in the rest of the chapter. Let G = (V, E) be an undirected graph, where V is the set of vertices and E is the set of edges. We assume that there is no *loop* in G, i.e., there is no edge in G which connects a vertex to itself. For any (possibly empty) subset U of V, denote by  $G \setminus U$  the graph obtained from G by eliminating all the vertices in U and all the edges joining a vertex in U.

The connectivity of a graph partitions the graph into subgraphs called *components*, i.e., two vertices are in the same component if and only if they are connected. Let s(U) be the number of distinct components in  $G \setminus U$ . Denote the sets of vertices of these components by  $V_1(U), V_2(U), \dots, V_{s(U)}(U)$ . If s(U) > 1, we say that U is a cutset in G.

**Definition 12.23 (Markov random field).** Let G = (V, E) be an undirected graph with  $V = \mathcal{N}_n = \{1, 2, \dots, n\}$ , and let  $X_i$  be a random variable corresponding to vertex i. Then  $X_1, X_2, \dots, X_n$  form a Markov random field represented by G if for all cutsets U in G, the sets of random variables  $X_{V_1(U)}, X_{V_2(U)}, \dots, X_{V_{s(U)}(U)}$  are mutually independent conditioning on  $X_U$ .

This definition of a Markov random field is referred to as the global Markov property in the literature. If  $X_1, X_2, \dots, X_n$  form a Markov random field represented by a graph G, we also say that  $X_1, X_2, \dots, X_n$  form a Markov graph G. When G is a chain, we say that  $X_1, X_2, \dots, X_n$  form a Markov chain.

In the definition of a Markov random field, each cutset U in G specifies an FCMI on  $X_1, X_2, \dots, X_n$ , denoted by [U]. Formally,

 $[U]: X_{V_1(U)}, \dots, X_{V_{s(U)}(U)}$  are mutually independent conditioning on  $X_U$ .

For a collection of cutsets  $U_1, U_2, \dots, U_k$  in G, we introduce the notation

$$[U_1, U_2, \cdots, U_k] = [U_1] \land [U_2] \land \cdots \land [U_k]$$
(12.88)

where ' $\wedge$ ' denotes 'logical AND.' Using this notation,  $X_1, X_2, \dots, X_n$  form a Markov graph G if and only if

$$[U \subset V : U \neq V \text{ and } s(U) > 1]$$

$$(12.89)$$

holds. Therefore, a Markov random field is simply a collection of FCMI's induced by a graph.

We now define two types of nonempty atoms of  $\mathcal{F}_n$  with respect to a graph G. Recall the definition of the set  $U_A$  for a nonempty atom A of  $\mathcal{F}_n$  in (12.73).

**Definition 12.24.** For a nonempty atom A of  $\mathcal{F}_n$ , if  $s(U_A) = 1$ , i.e.,  $G \setminus U_A$  is connected, then A is a Type I atom, otherwise A is a Type II atom. The sets of all Type I and Type II atoms of  $\mathcal{F}_n$  are denoted by  $T_1$  and  $T_2$ , respectively.

**Theorem 12.25.**  $X_1, X_2, \dots, X_n$  form a Markov graph G if and only if  $\mu^*$  vanishes on all Type II atoms.

Before we prove this theorem, we first state the following proposition which is the graph-theoretic analog of Theorem 12.5. The proof is trivial and is omitted. This proposition and Theorem 12.5 together establish an analogy between the structure of conditional mutual independence and the connectivity of a graph. This analogy will play a key role in proving Theorem 12.25.

**Proposition 12.26.** Let C and  $Q_i$  be disjoint subsets of the vertex set V of a graph G and  $W_i$  be a subset of  $Q_i$  for  $1 \le i \le k$ , where  $k \ge 2$ . Assume that there exist at least two i such that  $W_i \ne \emptyset$ . If  $Q_i, 1 \le i \le k$ , are disconnected in  $G \setminus C$ , then those  $W_i$  which are nonempty are disconnected in  $G \setminus (C \cup \bigcup_{i=1}^k (Q_i - W_i))$ .

*Example 12.27.* In the graph G in Figure 12.4,  $\{1\}$ ,  $\{2, 3, 4\}$ , and  $\{5, 6\}$  are disjoint in  $G \setminus \{7\}$ . Then Proposition 12.26 says that  $\{1\}$ ,  $\{2\}$ , and  $\{5, 6\}$  are disjoint in  $G \setminus \{3, 4, 7\}$ .

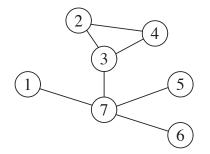


Fig. 12.4. The graph G in Example 12.27.

Proof of Theorem 12.25. Recall the definition of the set  $U_A$  for a nonempty atom A in (12.73). We note that  $\{U_A, A \in \mathcal{A}\}$  contains precisely all the proper subsets of  $\mathcal{N}_n$ . Thus the set of FCMI's specified by the graph G can be written as

$$[U_A : A \in \mathcal{A} \text{ and } s(U_A) > 1] \tag{12.90}$$

(cf. (12.89)). By Theorem 12.19, it suffices to prove that

$$Im([U_A : A \in \mathcal{A} \text{ and } s(U_A) > 1]) = T_2,$$
 (12.91)

where  $T_2$  was defined in Definition 12.24.

We first prove that

$$T_2 \subset Im([U_A : A \in \mathcal{A} \text{ and } s(U_A) > 1]).$$
(12.92)

Consider an atom  $A \in T_2$  so that  $s(U_A) > 1$ . In Definition 12.13, let  $T = U_A$ ,  $k = s(U_A)$ , and  $Q_i = V_i(U_A)$  for  $1 \le i \le s(U_A)$ . By considering  $W_i = V_i(U_A)$  for  $1 \le i \le s(U_A)$ , we see that  $A \in Im([U_A])$ . Therefore,

$$T_2 = \{A \in \mathcal{A} : s(U_A) > 1\}$$
(12.93)

$$\subset \bigcup_{A \in \mathcal{A}: s(U_A) > 1} Im([U_A])$$
(12.94)

$$= Im([U_A : A \in \mathcal{A} \text{ and } s(U_A) > 1]).$$
(12.95)

We now prove that

$$Im([U_A : A \in \mathcal{A} \text{ and } s(U_A) > 1]) \subset T_2.$$
(12.96)

Consider  $A \in Im([U_A : A \in \mathcal{A} \text{ and } s(U_A) > 1])$ . Then there exists  $A^* \in \mathcal{A}$  with  $s(U_{A^*}) > 1$  such that  $A \in Im([U_{A^*}])$ . From Definition 12.13,

$$A = \left(\bigcap_{j \in \bigcup_{i=1}^{s(U_{A^*})} W_i} \tilde{X}_j\right) - \tilde{X}_{U_{A^*} \cup \left(\bigcup_{i=1}^{s(U_{A^*})} (V_i(U_{A^*}) - W_i)\right)},$$
(12.97)

where  $W_i \subset V_i(U_{A^*})$ ,  $1 \leq i \leq s(U_{A^*})$ , and there exist at least two *i* such that  $W_i \neq \emptyset$ . It follows from (12.97) and the definition of  $U_A$  that

$$U_A = U_{A^*} \cup \bigcup_{i=1}^{s(U_{A^*})} (V_i(U_{A^*}) - W_i).$$
(12.98)

With  $U_{A^*}$  playing the role of C and  $V_i(U_{A^*})$  playing the role of  $Q_i$  in Proposition 12.26, we see by applying the proposition that those (at least two)  $W_i$  which are nonempty are disjoint in

$$G \setminus \left( U_{A^*} \cup \left( \bigcup_{i=1}^{s(U_{A^*})} (V_i(U_{A^*}) - W_i) \right) \right) = G \setminus U_A.$$
(12.99)

This implies  $s(U_A) > 1$ , i.e.,  $A \in T_2$ . Therefore, we have proved (12.96), and hence the theorem is proved.  $\Box$ 

*Example 12.28.* With respect to the graph G in Figure 12.5, the Type II atoms are

$$\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3^c \cap \tilde{X}_4, \ \tilde{X}_1^c \cap \tilde{X}_2 \cap \tilde{X}_3^c \cap \tilde{X}_4, \ \tilde{X}_1 \cap \tilde{X}_2^c \cap \tilde{X}_3^c \cap \tilde{X}_4, \ (12.100)$$

while the other twelve nonempty atoms of  $\mathcal{F}_4$  are Type I atoms. The random variables  $X_1, X_2, X_3$ , and  $X_4$  form a Markov graph G if and only if  $\mu^*(A) = 0$  for all Type II atoms A.

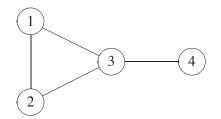


Fig. 12.5. The graph G in Example 12.28.

# 12.4 Markov Chain

When the graph G representing a Markov random field is a chain, the Markov random field becomes a Markov chain. In this section, we will show that the information diagram for a Markov chain can be displayed in two dimensions. We will also show that the *I*-Measure  $\mu^*$  for a Markov chain is always nonnegative. This characteristic of  $\mu^*$  facilitates the use of the information diagram because if B is seen to be a subset of B' in the information diagram, then

$$\mu^*(B') = \mu^*(B) + \mu^*(B' - B) \ge \mu^*(B).$$
(12.101)

These two properties are not possessed by a general Markov random field.

Without loss of generality, we assume that the Markov chain is represented by the graph G in Figure 12.6. This corresponds to the Markov chain  $X_1 \rightarrow$ 



**Fig. 12.6.** The graph G representing the Markov chain  $X_1 \to X_2 \to \cdots \to X_n$ .

 $X_2 \to \cdots \to X_n$ . We first prove the following characterization of a Type I atom for a Markov chain.

**Lemma 12.29.** For the Markov chain represented by the graph G in Figure 12.6, a nonempty atom A of  $\mathcal{F}_n$  is a Type I atom if and only if

$$\mathcal{N}_n \backslash U_A = \{l, l+1, \cdots, u\},\tag{12.102}$$

where  $1 \leq l \leq u \leq n$ , i.e., the indices of the set variables in A which are not complemented are consecutive.

*Proof.* It is easy to see that for a nonempty atom A, if (12.102) is satisfied, then  $G \setminus U_A$  is connected, i.e.,  $s(U_A) = 1$ . Therefore, A is a Type I atom of  $\mathcal{F}_n$ . On the other hand, if (12.102) is not satisfied, then  $G \setminus U_A$  is not connected, i.e.,  $s(U_A) > 1$ , or A is a Type II atom of  $\mathcal{F}_n$ . The lemma is proved.  $\Box$ 

We now show how the information diagram for a Markov chain with any length  $n \geq 3$  can be constructed in two dimensions. Since  $\mu^*$  vanishes on all the Type II atoms of  $\mathcal{F}_n$ , it is not necessary to display these atoms in the information diagram. In constructing the information diagram, the regions representing the random variables  $X_1, X_2, \dots, X_n$  should overlap with each other such that the regions corresponding to all the Type II atoms are empty, while the regions corresponding to all the Type I atoms are nonempty. Figure 12.7 shows such a construction. Note that this information diagram includes Figures 3.7 and 3.9 as special cases, which are information diagrams for Markov chains with lengths 3 and 4, respectively.

We have already shown that  $\mu^*$  is nonnegative for a Markov chain with length 3 or 4. Toward proving that this is true for any length  $n \ge 3$ , it suffices to show that  $\mu^*(A) \ge 0$  for all Type I atoms A of  $\mathcal{F}_n$  because  $\mu^*(A) = 0$  for all Type II atoms A of  $\mathcal{F}_n$ . We have seen in Lemma 12.29 that for a Type I atom A of  $\mathcal{F}_n$ ,  $U_A$  has the form as prescribed in (12.102). Consider any such atom A. Then an inspection of the information diagram in Figure 12.7 reveals that

$$\mu^{*}(A) = \mu^{*}(\tilde{X}_{l} \cap \tilde{X}_{l+1} \cap \dots \cap \tilde{X}_{u} - \tilde{X}_{U_{A}})$$
(12.103)

$$= I(X_l; X_u | X_{U_A}) (12.104)$$

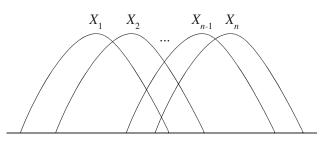
$$\geq 0. \tag{12.105}$$

This shows that  $\mu^*$  is always nonnegative. However, since Figure 12.7 involves an indefinite number of random variables, we give a formal proof of this result in the following theorem.

**Theorem 12.30.** For a Markov chain  $X_1 \to X_2 \to \cdots \to X_n$ ,  $\mu^*$  is nonnegative.

*Proof.* Since  $\mu^*(A) = 0$  for all Type II atoms A of  $\mathcal{F}_n$ , it suffices to show that  $\mu^*(A) \ge 0$  for all Type I atoms A of  $\mathcal{F}_n$ . We have seen in Lemma 12.29 that for a Type I atom A of  $\mathcal{F}_n$ ,  $U_A$  has the form as prescribed in (12.102). Consider any such atom A and define the set

$$W = \{l+1, \cdots, u-1\}.$$
 (12.106)



**Fig. 12.7.** The information diagram for the Markov chain  $X_1 \to X_2 \to \cdots \to X_n$ .

Then

$$I(X_l; X_u | X_{U_A}) = \mu^* (\tilde{X}_l \cap \tilde{X}_u - \tilde{X}_{U_A})$$
(12.107)

$$= \mu^* \left( \bigcup_{S \subset W} \left( \tilde{X}_l \cap \left( \bigcap_{t \in S} \tilde{X}_t \right) \cap \tilde{X}_u - \tilde{X}_{U_{A \cup (W \setminus S)}} \right) \right)$$
(12.108)

$$= \sum_{S \subset W} \mu^* \left( \tilde{X}_l \cap \left( \bigcap_{t \in S} \tilde{X}_t \right) \cap \tilde{X}_u - \tilde{X}_{U_{A \cup (W \setminus S)}} \right).$$
(12.109)

In the above summation, except for the atom corresponding to S = W, namely  $(\tilde{X}_l \cap \tilde{X}_{l+1} \cap \cdots \cap \tilde{X}_u - \tilde{X}_{U_A})$ , all the atoms are Type II atoms. Therefore,

$$I(X_{l}; X_{u} | X_{U_{A}}) = \mu^{*}(\tilde{X}_{l} \cap \tilde{X}_{l+1} \cap \dots \cap \tilde{X}_{u} - \tilde{X}_{U_{A}}).$$
(12.110)

Hence,

$$\mu^*(A) = \mu^*(\tilde{X}_l \cap \tilde{X}_{l+1} \cap \dots \cap \tilde{X}_u - \tilde{X}_{U_A})$$
(12.111)

$$= I(X_l; X_u | X_{U_A}) \tag{12.112}$$

$$\geq 0. \tag{12.113}$$

The theorem is proved.  $\Box$ 

### Problems

- 1. Prove Proposition 12.14 and Proposition 12.15.
- 2. In Example 12.22, it was shown that  $\Pi_2$  implies  $\Pi_1$ . Show that  $\Pi_1$  does not imply  $\Pi_2$ . Hint: Use an information diagram to determine  $Im(\Pi_2) \setminus Im(\Pi_1)$ .
- 3. Alternative definition of the global Markov property: For any partition  $\{U, V_1, V_2\}$  of V such that the sets of vertices  $V_1$  and  $V_2$  are disconnected in  $G \setminus U$ , the sets of random variables  $X_{V_1}$  and  $X_{V_2}$  are independent conditioning on  $X_U$ .

Show that this definition is equivalent to the global Markov property in Definition 12.23.

- 4. The local Markov property: For  $1 \leq i \leq n$ ,  $X_i$  and  $X_{V-N_i-i}$  are independent conditioning on  $X_{N_i}$ , where  $N_i$  is the set of neighbors<sup>2</sup> of vertex *i* in *G*.
  - a) Show that the global Markov property implies the local Markov property.

<sup>&</sup>lt;sup>2</sup> Vertices i and j in an undirected graph are neighbors if i and j are connected by an edge.

- 312 12 Markov Structures
  - b) Show that the local Markov property does not imply the global Markov property by giving a counterexample. Hint: Consider a joint distribution which is not strictly positive.
- 5. Construct a Markov random field whose *I*-Measure  $\mu^*$  can take negative values. Hint: Consider a Markov "star."
- 6. a) Show that  $X_1, X_2, X_3$ , and  $X_4$  are mutually independent if and only if

$$X_1 \perp (X_2, X_3, X_4), \ X_2 \perp (X_3, X_4) | X_1, \ X_3 \perp X_4 | (X_1, X_2).$$

Hint: Use an information diagram.

b) Generalize the result in a) to n random variables.

7. Determine the Markov random field with four random variables  $X_1, X_2, X_3$ , and  $X_4$  which is characterized by the following conditional independencies:

$$\begin{aligned} & (X_1, X_2, X_5) \perp X_4 | X_3 \\ & X_2 \perp (X_4, X_5) | (X_1, X_3) \\ & X_1 \perp (X_3, X_4) | (X_2, X_5). \end{aligned}$$

What are the other conditional independencies pertaining to this Markov random field?

### **Historical Notes**

A Markov random field can be regarded as a generalization of a discrete-time Markov chain. Historically, the study of Markov random field stems from statistical physics. The classical Ising model, which is defined on a rectangular lattice, was used to explain certain empirically observed facts about ferromagnetic materials. The foundation of the theory of Markov random fields can be found in Preston [267] or Spitzer [310].

The structure of the *I*-Measure for a Markov chain was first investigated in the unpublished work of Kawabata [176]. Essentially the same result was independently obtained by R. W. Yeung eleven years later in the context of the *I*-Measure, and the result was eventually published in Kawabata and Yeung [177]. Full conditional independencies were shown to be axiomatizable by Malvestuto [218]. The results in this chapter are due to Yeung *et al.* [364], where they obtained a set-theoretic characterization of full conditional independencies and investigated the structure of the *I*-Measure for a Markov random field. In this paper, they also obtained a hypergraph characterization of a Markov random field based on the *I*-Measure characterization in Theorem 12.25. Ge and Ye [119] have applied these results to characterize a class of graphical models for conditional independence of random variables.

# Information Inequalities

An information expression f refers to a linear combination<sup>1</sup> of Shannon's information measures involving a finite number of random variables. For example,

$$H(X,Y) + 2I(X;Z)$$
 (13.1)

and

$$I(X;Y) - I(X;Y|Z)$$
 (13.2)

are information expressions. An information inequality has the form

$$f \ge c, \tag{13.3}$$

where the constant c is usually equal to zero. We consider non-strict inequalities only because these are usually the form of inequalities in information theory. Likewise, an information identity has the form

$$f = c. \tag{13.4}$$

We point out that an information identity f = c is equivalent to the pair of information inequalities  $f \ge c$  and  $f \le c$ .

An information inequality or identity is said to *always hold* if it holds for any joint distribution for the random variables involved. For example, we say that the information inequality

$$I(X;Y) \ge 0 \tag{13.5}$$

always holds because it holds for any joint distribution p(x, y). On the other hand, we say that an information inequality does not always hold if there exists a joint distribution for which the inequality does not hold. Consider the information inequality

<sup>&</sup>lt;sup>1</sup> More generally, an information expression can be nonlinear, but they do not appear to be useful in information theory.

314 13 Information Inequalities

$$I(X;Y) \le 0. \tag{13.6}$$

Since

$$I(X;Y) \ge 0 \tag{13.7}$$

always holds, (13.6) is equivalent to

$$I(X;Y) = 0, (13.8)$$

which holds if and only if X and Y are independent. In other words, (13.6) does not hold if X and Y are not independent. Therefore, we say that (13.6) does not always hold.

As we have seen in the previous chapters, information inequalities are the major tools for proving converse coding theorems. These inequalities govern the impossibilities in information theory. More precisely, information inequalities imply that certain things cannot happen. As such, they are referred to as the *laws of information theory*.

The basic inequalities form the most important set of information inequalities. In fact, almost all the information inequalities known to date are implied by the basic inequalities. These are called *Shannon-type inequalities*. On the other hand, if an information inequality always holds but is not implied by the basic inequalities, then it is called a *non-Shannon-type inequality*. We have not yet explained what it means by that an inequality is or is not implied by the basic inequalities, but this will become clear later in the chapter.

Let us now rederive the inequality obtained in Example 3.15 (Imperfect secrecy theorem) without using an information diagram. In this example, three random variables X, Y, and Z are involved, and the setup of the problem is equivalent to the constraint

$$H(X|Y,Z) = 0. (13.9)$$

Then

$$= H(X) + H(Y) - H(X,Y)$$
(13.10)

$$= H(X) + H(Y) - [H(X, Y, Z) - H(Z|X, Y)]$$
(13.11)

$$\geq H(X) + H(Y) - H(X, Y, Z)$$
(13.12)

$$= H(X) + H(Y) - [H(Z) + H(Y|Z) + H(X|Y,Z)]$$
(13.13)

$$= H(X) - H(Z) + I(Y;Z) - H(X|Y,Z)$$
(13.14)

$$\geq H(X) - H(Z), \tag{13.15}$$

where we have used

$$H(Z|X,Y) \ge 0 \tag{13.16}$$

in obtaining (13.12), and

I(X:Y)

$$I(Y;Z) \ge 0 \tag{13.17}$$

and (13.9) in obtaining (13.15). This derivation is less transparent than the one we presented in Example 3.15, but the point here is that the final inequality we obtain in (13.15) can be proved by invoking the basic inequalities (13.16) and (13.17). In other words, (13.15) is implied by the basic inequalities. Therefore, it is a (constrained) Shannon-type inequality.

We are motivated to ask the following two questions:

- 1. How can Shannon-type inequalities be characterized? That is, given an information inequality, how can we tell whether it is implied by the basic inequalities?
- 2. Are there any non-Shannon-type information inequalities?

These are two very fundamental questions in information theory. We point out that the first question naturally comes before the second question because if we cannot characterize all Shannon-type inequalities, even if we are given a non-Shannon-type inequality, we cannot tell that it actually is one.

In this chapter, we develop a geometric framework for information inequalities which allows them to be studied systematically. This framework naturally leads to an answer to the first question, which makes machine-proving of all Shannon-type inequalities possible. This will be discussed in the next chapter. The second question will be answered positively in Chapter 15. In other words, there *do* exist laws in information theory beyond those laid down by Shannon.

## 13.1 The Region $\Gamma_n^*$

Let

$$\mathcal{N}_n = \{1, 2, \cdots, n\},\tag{13.18}$$

where  $n \geq 2$ , and let

$$\Theta = \{X_i, i \in \mathcal{N}_n\} \tag{13.19}$$

be any collection of n random variables. Associated with  $\Theta$  are

$$k = 2^n - 1 \tag{13.20}$$

joint entropies. For example, for n = 3, the 7 joint entropies associated with random variables  $X_1, X_2$ , and  $X_3$  are

$$H(X_1), H(X_2), H(X_3), H(X_1, X_2),$$
  
$$H(X_2, X_3), H(X_1, X_3), H(X_1, X_2, X_3).$$
 (13.21)

Let  $\Re$  denote the set of real numbers. For any nonempty subset  $\alpha$  of  $\mathcal{N}_n$ , let

$$X_{\alpha} = (X_i, i \in \alpha) \tag{13.22}$$

and

31613 Information Inequalities

$$H_{\Theta}(\alpha) = H(X_{\alpha}). \tag{13.23}$$

For a fixed  $\Theta$ , we can then view  $H_{\Theta}$  as a set function from  $2^{\mathcal{N}_n}$  to  $\Re$  with  $H_{\Theta}(\emptyset) = 0$ , i.e., we adopt the convention that the entropy of an empty set of random variable is equal to zero. For this reason, we call  $H_{\Theta}$  the *entropy* function of  $\Theta$ .

Let  $\mathcal{H}_n$  be the k-dimensional Euclidean space with the coordinates labeled by  $h_{\alpha}, \alpha \in 2^{\mathcal{N}_n} \setminus \{\emptyset\}$ , where  $h_{\alpha}$  corresponds to the value of  $H_{\Theta}(\alpha)$  for any collection  $\Theta$  of *n* random variables. We will refer to  $\mathcal{H}_n$  as the *entropy space* for n random variables. Then an entropy function  $H_{\Theta}$  can be represented by a column vector in  $\mathcal{H}_n$ . On the other hand, a column vector  $\mathbf{h} \in \mathcal{H}_n$  is called *entropic* if **h** is equal to the entropy function  $H_{\Theta}$  of some collection  $\Theta$  of n random variables. We are motivated to define the following region in  $\mathcal{H}_n$ :

$$\Gamma_n^* = \{ \mathbf{h} \in \mathcal{H}_n : \mathbf{h} \text{ is entropic} \}.$$
(13.24)

For convenience, the vectors in  $\Gamma_n^*$  will also be referred to as entropy functions. As an example, for n = 3, the coordinates of  $\mathcal{H}_3$  are labeled by

$$h_1, h_2, h_3, h_{12}, h_{13}, h_{23}, h_{123},$$
 (13.25)

where  $h_{123}$  denotes  $h_{\{1,2,3\}}$ , etc, and  $\Gamma_3^*$  is the region in  $\mathcal{H}_3$  of all entropy functions for 3 random variables.

While further characterizations of  $\Gamma_n^*$  will be given later, we first point out a few basic properties of  $\Gamma_n^*$ :

- Γ<sub>n</sub><sup>\*</sup> contains the origin.
   Γ<sub>n</sub><sup>\*</sup>, the closure of Γ<sub>n</sub><sup>\*</sup>, is convex.
   Γ<sub>n</sub><sup>\*</sup> is in the nonnegative orthant of the entropy space H<sub>n</sub><sup>2</sup>.

The origin of the entropy space corresponds to the entropy function of n degenerate random variables taking constant values. Hence, Property 1 follows. Property 2 will be proved in Chapter 15. Properties 1 and 2 imply that  $\overline{\Gamma}_n^*$ is a convex cone. Property 3 is true because the coordinates in the entropy space  $\mathcal{H}_n$  correspond to joint entropies, which are always nonnegative.

### **13.2** Information Expressions in Canonical Form

Any Shannon's information measure other than a joint entropy can be expressed as a linear combination of joint entropies by application of one of the following information identities:

$$H(X|Y) = H(X,Y) - H(Y)$$
(13.26)

$$I(X;Y) = H(X) + H(Y) - H(X,Y)$$
(13.27)

$$I(X;Y|Z) = H(X,Z) + H(Y,Z) - H(X,Y,Z) - H(Z).$$
(13.28)

<sup>&</sup>lt;sup>2</sup> The nonnegative orthant of  $\mathcal{H}^n$  is the region  $\{\mathbf{h} \in \mathcal{H}_n : h_\alpha \geq 0 \text{ for all } \alpha \in$  $2^{\mathcal{N}_n} \setminus \{\emptyset\}\}.$ 

The first and the second identity are special cases of the third identity, which has already been proved in Lemma 3.8. Thus any information expression which involves n random variables can be expressed as a linear combination of the kassociated joint entropies. We call this the *canonical form* of an information expression. When we write an information expression f as  $f(\mathbf{h})$ , it means that f is in canonical form. Since an information expression in canonical form is a linear combination of the joint entropies, it has the form

$$\mathbf{b}^{\mathsf{T}}\mathbf{h} \tag{13.29}$$

where  $\mathbf{b}^{\top}$  denotes the transpose of a constant column vector  $\mathbf{b}$  in  $\Re^k$ .

The identities in (13.26) to (13.28) provide a way to express every information expression in canonical form. However, it is not clear whether such a canonical form is unique. To illustrate the point, we consider obtaining the canonical form of H(X|Y) in two ways. First,

$$H(X|Y) = H(X,Y) - H(Y).$$
 (13.30)

Second,

$$H(X|Y) = H(X) - I(X;Y)$$
(13.31)

$$= H(X) - (H(Y) - H(Y|X))$$
(13.32)  
= H(X) - (H(Y) - H(X,Y) + H(X)) (13.33)

$$= H(X) - (H(Y) - H(X, Y) + H(X))$$
(13.33)  
$$H(X, Y) - H(Y)$$
(13.41)

$$= H(X,Y) - H(Y).$$
(13.34)

Thus it turns out that we can obtain the same canonical form for H(X|Y) via two different expansions. This is not accidental, as it is implied by the uniqueness of the canonical form which we will prove shortly.

Recall from the proof of Theorem 3.6 that the vector  $\mathbf{h}$  represents the values of the *I*-Measure  $\mu^*$  on the unions in  $\mathcal{F}_n$ . Moreover,  $\mathbf{h}$  is related to the values of  $\mu^*$  on the atoms of  $\mathcal{F}_n$ , represented as  $\mathbf{u}$ , by

$$\mathbf{h} = \mathbf{C}_n \mathbf{u} \tag{13.35}$$

where  $\mathbf{C}_n$  is a unique  $k \times k$  matrix (cf. (3.27)). We now state the following lemma which is a rephrase of Theorem 3.11. This lemma is essential for proving the next theorem which implies the uniqueness of the canonical form.

Lemma 13.1. Let

$$\Psi_n^* = \{ \mathbf{u} \in \Re^k : \mathbf{C}_n \mathbf{u} \in \Gamma_n^* \}.$$
(13.36)

Then the nonnegative orthant of  $\Re^k$  is a subset of  $\Psi_n^*$ .

**Theorem 13.2.** Let f be an information expression. Then the unconstrained information identity f = 0 always holds if and only if f is the zero function.

#### 318 13 Information Inequalities

*Proof.* Without loss of generality, assume f is in canonical form and let

$$f(\mathbf{h}) = \mathbf{b}^{\top} \mathbf{h}. \tag{13.37}$$

Assume f = 0 always holds and f is not the zero function, i.e.,  $\mathbf{b} \neq 0$ . We will show that this leads to a contradiction. Now f = 0, or more precisely the set

$$\{\mathbf{h}: \mathbf{b}^{\top}\mathbf{h} = 0\},\tag{13.38}$$

is a hyperplane<sup>3</sup> in the entropy space which has zero Lebesgue measure<sup>4</sup>. If f = 0 always holds, i.e., it holds for all joint distributions, then  $\Gamma_n^*$  must be contained in the hyperplane f = 0, otherwise there exists an  $\mathbf{h}_0 \in \Gamma_n^*$  which is not on f = 0, i.e.,  $f(\mathbf{h}_0) \neq 0$ . Since  $\mathbf{h}_0 \in \Gamma_n^*$ , it corresponds to the entropy function of some joint distribution. This means that there exists a joint distribution such that  $f(\mathbf{h}) = 0$  does not hold, which cannot be true because f = 0 always holds.

If  $\Gamma_n^*$  has positive Lebesgue measure, it cannot be contained in the hyperplane f = 0 which has zero Lebesgue measure. Therefore, it suffices to show that  $\Gamma_n^*$  has positive Lebesgue measure. To this end, we see from Lemma 13.1 that the nonnegative orthant of  $\mathcal{H}_n$ , which has positive Lebesgue measure, is a subset of  $\Psi_n^*$ . Thus  $\Psi_n^*$  has positive Lebesgue measure. Since  $\Gamma_n^*$  is an invertible transformation of  $\Psi_n^*$ , its Lebesgue measure is also positive.

Therefore,  $\Gamma_n^*$  is not contained in the hyperplane f = 0, which implies that there exists a joint distribution for which f = 0 does not hold. This leads to a contradiction because we have assumed that f = 0 always holds. Hence, we have proved that if f = 0 always holds, then f must be the zero function.

Conversely, if f is the zero function, then it is trivial that f = 0 always holds. The theorem is proved.  $\Box$ 

Corollary 13.3. The canonical form of an information expression is unique.

*Proof.* Let  $f_1$  and  $f_2$  be canonical forms of an information expression g. Since

$$g = f_1 \tag{13.39}$$

and

$$g = f_2 \tag{13.40}$$

always hold,

$$f_1 - f_2 = 0 \tag{13.41}$$

always holds. By the above theorem,  $f_1 - f_2$  is the zero function, which implies that  $f_1$  and  $f_2$  are identical. The corollary is proved.  $\Box$ 

<sup>&</sup>lt;sup>3</sup> If  $\mathbf{b} = 0$ , then  $\{\mathbf{h} : \mathbf{b}^{\top} \mathbf{h} = 0\}$  is equal to  $\mathcal{H}_n$ .

<sup>&</sup>lt;sup>4</sup> The Lebesque measure can be thought of as "volume" in the Euclidean space if the reader is not familiar with measure theory.

Due to the uniqueness of the canonical form of an information expression, it is an easy matter to check whether for two information expressions  $f_1$  and  $f_2$  the unconstrained information identity

$$f_1 = f_2$$
 (13.42)

always holds. All we need to do is to express  $f_1 - f_2$  in canonical form. If all the coefficients are zero, then (13.42) always holds, otherwise it does not.

### 13.3 A Geometrical Framework

In the last section, we have seen the role of the region  $\Gamma_n^*$  in proving unconstrained information identities. In this section, we explain the geometrical meanings of unconstrained information inequalities, constrained information inequalities, and constrained information identities in terms of  $\Gamma_n^*$ . Without loss of generality, we assume that all information expressions are in canonical form.

#### 13.3.1 Unconstrained Inequalities

Consider an unconstrained information inequality  $f \ge 0$ , where  $f(\mathbf{h}) = \mathbf{b}^{\top} \mathbf{h}$ . Then  $f \ge 0$  corresponds to the set

$$\{\mathbf{h} \in \mathcal{H}_n : \mathbf{b}^\top \mathbf{h} \ge 0\} \tag{13.43}$$

which is a half-space in the entropy space  $\mathcal{H}_n$  containing the origin. Specifically, for any  $\mathbf{h} \in \mathcal{H}_n$ ,  $f(\mathbf{h}) \geq 0$  if and only if  $\mathbf{h}$  belongs to this set. For simplicity, we will refer to this set as the half-space  $f \geq 0$ . As an example, for n = 2, the information inequality

$$I(X_1; X_2) = H(X_1) + H(X_2) - H(X_1, X_2) \ge 0,$$
(13.44)

written as

$$h_1 + h_2 - h_{12} \ge 0, \tag{13.45}$$

corresponds to the half-space

$$\{\mathbf{h} \in \mathcal{H}_n : h_1 + h_2 - h_{12} \ge 0\}.$$
(13.46)

in the entropy space  $\mathcal{H}_2$ .

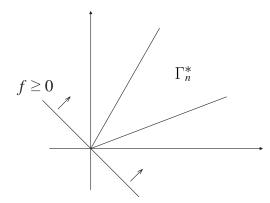
Since an information inequality always holds if and only if it is satisfied by the entropy function of any joint distribution for the random variables involved, we have the following geometrical interpretation of an information inequality:

 $f \ge 0$  always holds if and only if  $\Gamma_n^* \subset {\mathbf{h} \in \mathcal{H}_n : f(\mathbf{h}) \ge 0}.$ 

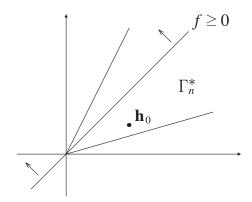
#### 320 13 Information Inequalities

This gives a complete characterization of all unconstrained inequalities in terms of  $\Gamma_n^*$ . If  $\Gamma_n^*$  is known, we in principle can determine whether any information inequality involving n random variables always holds.

The two possible cases for  $f \ge 0$  are illustrated in Figure 13.1 and Figure 13.2. In Figure 13.1,  $\Gamma_n^*$  is completely included in the half-space  $f \ge 0$ ,



**Fig. 13.1.** An illustration for  $f \ge 0$  always holds.



**Fig. 13.2.** An illustration for  $f \ge 0$  not always holds.

so  $f \ge 0$  always holds. In Figure 13.2, there exists a vector  $\mathbf{h}_0 \in \Gamma_n^*$  such that  $f(\mathbf{h}_0) < 0$ . Thus the inequality  $f \ge 0$  does not always hold.

### 13.3.2 Constrained Inequalities

In information theory, we very often deal with information inequalities (identities) with certain constraints on the joint distribution for the random variables involved. These are called constrained information inequalities (identities), and the constraints on the joint distribution can usually be expressed as linear constraints on the entropies. The following are such examples:

- 1.  $X_1, X_2$ , and  $X_3$  are mutually independent if and only if  $H(X_1, X_2, X_3) = H(X_1) + H(X_2) + H(X_3)$ .
- 2.  $X_1, X_2$ , and  $X_3$  are pairwise independent if and only if  $I(X_1; X_2) = I(X_2; X_3) = I(X_1; X_3) = 0$ .
- 3.  $X_1$  is a function of  $X_2$  if and only if  $H(X_1|X_2) = 0$ .
- 4.  $X_1 \rightarrow X_2 \rightarrow X_3 \rightarrow X_4$  forms a Markov chain if and only if  $I(X_1; X_3 | X_2) = 0$  and  $I(X_1, X_2; X_4 | X_3) = 0$ .

Suppose there are q linear constraints on the entropies given by

$$\mathbf{Qh} = 0, \tag{13.47}$$

where  $\mathbf{Q}$  is a  $q \times k$  matrix. Here we do not assume that the q constraints are linearly independent, so  $\mathbf{Q}$  is not necessarily full rank. Let

$$\Phi = \{ \mathbf{h} \in \mathcal{H}_n : \mathbf{Q}\mathbf{h} = 0 \}.$$
(13.48)

In other words, the q constraints confine **h** to a linear subspace  $\Phi$  in the entropy space. Parallel to our discussion on unconstrained inequalities, we have the following geometrical interpretation of a constrained inequality:

Under the constraint 
$$\Phi$$
,  $f \ge 0$  always holds if and only if  $(\Gamma_n^* \cap \Phi) \subset {\mathbf{h} : f(\mathbf{h}) \ge 0}.$ 

This gives a complete characterization of all constrained inequalities in terms of  $\Gamma_n^*$ . Note that  $\Phi = \mathcal{H}_n$  when there is no constraint on the entropies. In this sense, an unconstrained inequality is a special case of a constrained inequality.

The two cases of  $f \ge 0$  under the constraint  $\Phi$  are illustrated in Figure 13.3 and Figure 13.4. Figure 13.3 shows the case when  $f \ge 0$  always holds under the constraint  $\Phi$ . Note that  $f \ge 0$  may or may not always hold when there is no constraint. Figure 13.4 shows the case when  $f \ge 0$  does not always hold under the constraint  $\Phi$ . In this case,  $f \ge 0$  does not always hold when there is no constraint, because

$$(\Gamma_n^* \cap \Phi) \not\subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\}$$
(13.49)

implies

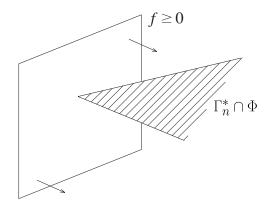
$$\Gamma_n^* \not\subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\}. \tag{13.50}$$

#### 13.3.3 Constrained Identities

As we have pointed out at the beginning of the chapter, an identity

$$f = 0$$
 (13.51)

always holds if and only if both the inequalities  $f \ge 0$  and  $f \le 0$  always hold. Then following our discussion on constrained inequalities, we have



**Fig. 13.3.** An illustration for  $f \ge 0$  always holds under the constraint  $\Phi$ .

Under the constraint  $\Phi$ , f = 0 always holds if and only if  $(\Gamma_n^* \cap \Phi) \subset {\mathbf{h} : f(\mathbf{h}) \ge 0} \cap {\mathbf{h} : f(\mathbf{h}) \le 0}$ ,

or

Under the constraint  $\Phi$ , f = 0 always holds if and only if  $(\Gamma_n^* \cap \Phi) \subset {\mathbf{h} : f(\mathbf{h}) = 0}.$ 

This condition says that the intersection of  $\Gamma_n^*$  and  $\Phi$  is contained in the hyperplane f = 0.

# 13.4 Equivalence of Constrained Inequalities

When there is no constraint on the entropies, two information inequalities

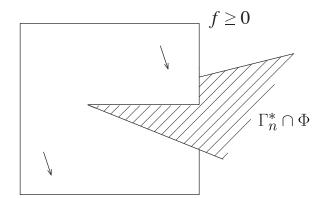


Fig. 13.4. An illustration for  $f \ge 0$  not always holds under the constraint  $\Phi$ .

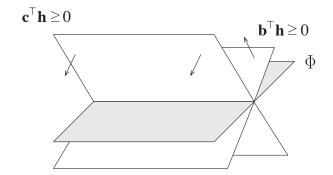
13.4 Equivalence of Constrained Inequalities 323

$$\mathbf{b}^{\mathsf{T}}\mathbf{h} \ge 0 \tag{13.52}$$

and

$$\mathbf{c}^{\top}\mathbf{h} \ge 0 \tag{13.53}$$

are equivalent if and only if  $\mathbf{c} = a\mathbf{b}$ , where *a* is a positive constant. However, this is not the case under a non-trivial constraint  $\Phi \neq \mathcal{H}_n$ . This situation is illustrated in Figure 13.5. In this figure, although the inequalities in (13.52)



**Fig. 13.5.** Equivalence of  $\mathbf{b}^{\top}\mathbf{h} \ge 0$  and  $\mathbf{c}^{\top}\mathbf{h} \ge 0$  under the constraint  $\boldsymbol{\Phi}$ .

and (13.53) correspond to different half-spaces in the entropy space, they actually impose the same constraint on  $\mathbf{h}$  when  $\mathbf{h}$  is confined to  $\Phi$ .

In this section, we present a characterization of (13.52) and (13.53) being equivalent under a set of linear constraint  $\Phi$ . The reader may skip this section at first reading.

Let r be the rank of **Q** in (13.47). Since **h** is in the null space of **Q**, we can write

$$\mathbf{h} = \mathbf{\hat{Q}}\mathbf{h}',\tag{13.54}$$

where  $\tilde{\mathbf{Q}}$  is a  $k \times (k-r)$  matrix such that the rows of  $\tilde{\mathbf{Q}}^{\top}$  form a basis of the orthogonal complement of the row space of  $\mathbf{Q}$ , and  $\mathbf{h}'$  is a column (k-r)-vector. Then using (13.54), (13.52) and (13.53) can be written as

$$\mathbf{b}^{\top}\mathbf{Q}\mathbf{h}' \ge 0 \tag{13.55}$$

and

$$\mathbf{c}^{\top}\tilde{\mathbf{Q}}\mathbf{h}' \ge 0, \tag{13.56}$$

respectively in terms of the set of basis given by the columns of  $\hat{\mathbf{Q}}$ . Then (13.55) and (13.56) are equivalent if and only if

$$\mathbf{c}^{\top}\tilde{\mathbf{Q}} = a\mathbf{b}^{\top}\tilde{\mathbf{Q}},\tag{13.57}$$

### 324 13 Information Inequalities

where a is a positive constant, or

$$(\mathbf{c} - a\mathbf{b})^{\top} \tilde{\mathbf{Q}} = 0. \tag{13.58}$$

In other words,  $(\mathbf{c} - a\mathbf{b})^{\top}$  is in the orthogonal complement of the row space of  $\tilde{\mathbf{Q}}^{\top}$ , i.e.,  $(\mathbf{c} - a\mathbf{b})^{\top}$  is in the row space of  $\mathbf{Q}$ . Let  $\mathbf{Q}'$  be an  $r \times k$  matrix whose row space is the same as that of  $\mathbf{Q}$ . ( $\mathbf{Q}$  can be taken as  $\mathbf{Q}'$  if  $\mathbf{Q}$  is full rank.) Since the rank of  $\mathbf{Q}$  is r and  $\mathbf{Q}'$  has r rows, the rows of  $\mathbf{Q}'$  form a basis for the row space of  $\mathbf{Q}$ , and  $\mathbf{Q}'$  is full rank. Then from (13.58), (13.55) and (13.56) are equivalent under the constraint  $\Phi$  if and only if

$$\mathbf{c} = a\mathbf{b} + (\mathbf{Q}')^{\top}\mathbf{e} \tag{13.59}$$

for some positive constant a and some column r-vector  $\mathbf{e}$ .

Suppose for given **b** and **c**, we want to see whether (13.55) and (13.56) are equivalent under the constraint  $\Phi$ . We first consider the case when either  $\mathbf{b}^{\top}$  or  $\mathbf{c}^{\top}$  is in the row space of **Q**. This is actually not an interesting case because if  $\mathbf{b}^{\top}$ , for example, is in the row space of **Q**, then

$$\mathbf{b}^{\top} \mathbf{\hat{Q}} = 0 \tag{13.60}$$

in (13.55), which means that (13.55) imposes no additional constraint under the constraint  $\Phi$ .

**Theorem 13.4.** If either  $\mathbf{b}^{\top}$  or  $\mathbf{c}^{\top}$  is in the row space of  $\mathbf{Q}$ , then  $\mathbf{b}^{\top}\mathbf{h} \geq 0$ and  $\mathbf{c}^{\top}\mathbf{h} \geq 0$  are equivalent under the constraint  $\Phi$  if and only if both  $\mathbf{b}^{\top}$  and  $\mathbf{c}^{\top}$  are in the row space of  $\mathbf{Q}$ .

The proof of this theorem is left as an exercise. We now turn to the more interesting case when neither  $\mathbf{b}^{\top}$  nor  $\mathbf{c}^{\top}$  is in the row space of  $\mathbf{Q}$ . The following theorem gives an explicit condition for (13.55) and (13.56) to be equivalent under the constraint  $\boldsymbol{\Phi}$ .

**Theorem 13.5.** If neither  $\mathbf{b}^{\top}$  nor  $\mathbf{c}^{\top}$  is in the row space of  $\mathbf{Q}$ , then  $\mathbf{b}^{\top}\mathbf{h} \ge 0$ and  $\mathbf{c}^{\top}\mathbf{h} \ge 0$  are equivalent under the constraint  $\Phi$  if and only if

$$\begin{bmatrix} (\mathbf{Q}')^\top \mathbf{b} \end{bmatrix} \begin{bmatrix} \mathbf{e} \\ a \end{bmatrix} = \mathbf{c}.$$
 (13.61)

has a unique solution with a > 0, where  $\mathbf{Q}'$  is any matrix whose row space is the same as that of  $\mathbf{Q}$ .

*Proof.* For  $\mathbf{b}^{\top}$  and  $\mathbf{c}^{\top}$  not in the row space of  $\mathbf{Q}$ , we want to see when we can find unknowns a and  $\mathbf{e}$  satisfying (13.59) with a > 0. To this end, we write (13.59) in matrix form as (13.61). Since  $\mathbf{b}$  is not in the column space of  $(\mathbf{Q}')^{\top}$  and  $(\mathbf{Q}')^{\top}$  is full rank,  $[(\mathbf{Q}')^{\top} \mathbf{b}]$  is also full rank. Then (13.61) has either a unique solution or no solution. Therefore, the necessary and sufficient condition for (13.55) and (13.56) to be equivalent is that (13.61) has a unique solution and a > 0. The theorem is proved.  $\Box$ 

*Example 13.6.* Consider three random variables  $X_1, X_2$ , and  $X_3$  with the Markov constraint

$$I(X_1; X_3 | X_2) = 0, (13.62)$$

which is equivalent to

$$H(X_1, X_2) + H(X_2, X_3) - H(X_1, X_2, X_3) - H(X_2) = 0.$$
(13.63)

In terms of the coordinates in the entropy space  $\mathcal{H}_3$ , this constraint is written as

$$\mathbf{Qh} = 0, \tag{13.64}$$

where

$$\mathbf{Q} = \begin{bmatrix} 0 \ -1 \ 0 \ 1 \ 1 \ 0 \ -1 \end{bmatrix} \tag{13.65}$$

and

$$\mathbf{h} = [h_1 \ h_2 \ h_3 \ h_{12} \ h_{23} \ h_{13} \ h_{123}]^{\top}.$$
(13.66)

We now show that under the constraint in (13.64), the inequalities

$$H(X_1|X_3) - H(X_1|X_2) \ge 0 \tag{13.67}$$

and

$$I(X_1; X_2 | X_3) \ge 0 \tag{13.68}$$

are in fact equivalent. Toward this end, we write (13.67) and (13.68) as  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  and  $\mathbf{c}^{\top}\mathbf{h} \geq 0$ , respectively, where

$$\mathbf{b} = \begin{bmatrix} 0 \ 1 \ -1 \ -1 \ 0 \ 1 \ 0 \end{bmatrix}^{\top} \tag{13.69}$$

and

$$\mathbf{c} = \begin{bmatrix} 0 \ 0 \ -1 \ 0 \ 1 \ 1 \ -1 \end{bmatrix}^{\top}. \tag{13.70}$$

Since  $\mathbf{Q}$  is full rank, we may take  $\mathbf{Q}' = \mathbf{Q}$ . Upon solving

$$\begin{bmatrix} \mathbf{Q}^\top \mathbf{b} \end{bmatrix} \begin{bmatrix} \mathbf{e} \\ a \end{bmatrix} = \mathbf{c}, \tag{13.71}$$

we obtain the unique solution a = 1 > 0 and  $\mathbf{e} = 1$  ( $\mathbf{e}$  is a  $1 \times 1$  matrix). Therefore, (13.67) and (13.68) are equivalent under the constraint in (13.64).

Under the constraint  $\Phi$ , if neither  $\mathbf{b}^{\top}$  nor  $\mathbf{c}^{\top}$  is in the row space of  $\mathbf{Q}$ , it can be shown that the identities

$$\mathbf{b}^{\mathsf{T}}\mathbf{h} = 0 \tag{13.72}$$

and

$$\mathbf{c}^{\top}\mathbf{h} = 0 \tag{13.73}$$

are equivalent if and only if (13.61) has a unique solution. We leave the proof as an exercise.

## 13.5 The Implication Problem of Conditional Independence

We use  $X_{\alpha} \perp X_{\beta} | X_{\gamma}$  to denote the conditional independency (CI)

 $X_{\alpha}$  and  $X_{\beta}$  are conditionally independent given  $X_{\gamma}$ .

We have proved in Theorem 2.34 that  $X_{\alpha} \perp X_{\beta} | X_{\gamma}$  is equivalent to

$$I(X_{\alpha}; X_{\beta} | X_{\gamma}) = 0. \tag{13.74}$$

When  $\gamma = \emptyset$ ,  $X_{\alpha} \perp X_{\beta} | X_{\gamma}$  becomes an unconditional independency which we regard as a special case of a conditional independency. When  $\alpha = \beta$ , (13.74) becomes

$$H(X_{\alpha}|X_{\gamma}) = 0, \tag{13.75}$$

which we see from Proposition 2.36 that  $X_{\alpha}$  is a function of  $X_{\gamma}$ . For this reason, we also regard functional dependency as a special case of conditional independency.

In probability problems, we are often given a set of CI's and we need to determine whether another given CI is logically implied. This is called the *implication problem*, which is perhaps the most basic problem in probability theory. We have seen in Section 12.2 that the implication problem has a solution if only full conditional mutual independencies are involved. However, the general problem is extremely difficult, and it has recently been solved only up to four random variables by Matúš [230].

We end this section by explaining the relation between the implication problem and the region  $\Gamma_n^*$ . A CI involving random variables  $X_1, X_2, \dots, X_n$ has the form

$$X_{\alpha} \perp X_{\beta} | X_{\gamma}, \tag{13.76}$$

where  $\alpha, \beta, \gamma \in \mathcal{N}_n$ . Since  $I(X_{\alpha}; X_{\beta} | X_{\gamma}) = 0$  is equivalent to

$$H(X_{\alpha\cup\gamma}) + H(X_{\beta\cup\gamma}) - H(X_{\alpha\cup\beta\cup\gamma}) - H(X_{\gamma}) = 0, \qquad (13.77)$$

 $X_{\alpha} \perp X_{\beta} | X_{\gamma}$  corresponds to the hyperplane

$$\{\mathbf{h}: h_{\alpha\cup\gamma} + h_{\beta\cup\gamma} - h_{\alpha\cup\beta\cup\gamma} - h_{\gamma} = 0\}.$$
(13.78)

For a CI K, we denote the hyperplane in  $\mathcal{H}_n$  corresponding to K by  $\mathcal{E}(K)$ .

Let  $\Pi = \{K_l\}$  be a collection of CI's, and we want to determine whether  $\Pi$  implies a given CI K. This would be the case if and only if the following is true:

For all 
$$\mathbf{h} \in \Gamma_n^*$$
, if  $\mathbf{h} \in \bigcap_l \mathcal{E}(K_l)$ , then  $\mathbf{h} \in \mathcal{E}(K)$ .

Equivalently,

$$\Pi$$
 implies K if and only if  $\left(\bigcap_{l} \mathcal{E}(K_{l})\right) \cap \Gamma_{n}^{*} \subset \mathcal{E}(K).$ 

Therefore, the implication problem can be solved if  $\Gamma_n^*$  can be characterized. Hence, the region  $\Gamma_n^*$  is not only of fundamental importance in information theory, but is also of fundamental importance in probability theory.

### Problems

- 1. Symmetrical information expressions An information expression is said to be symmetrical if it is identical under every permutation of the random variables involved. However, sometimes a symmetrical information expression cannot be readily recognized symbolically. For example,  $I(X_1; X_2) - I(X_1; X_2 | X_3)$  is symmetrical in  $X_1, X_2$ , and  $X_3$  but it is not symmetrical symbolically. Devise a general method for recognizing symmetrical information expressions.
- 2. The canonical form of an information expression is unique when there is no constraint on the random variables involved. Show by an example that this does not hold when certain constraints are imposed on the random variables involved.
- 3. Alternative canonical form Denote  $\cap_{i \in G} \tilde{X}_i$  by  $\check{X}_G$  and let

 $\mathcal{C} = \{\check{X}_G : G \text{ is a nonempty subset of } \mathcal{N}_n\}.$ 

- a) Prove that a signed measure  $\mu$  on  $\mathcal{F}_n$  is completely specified by  $\{\mu(C), \mu(C), \mu$  $C \in \mathcal{C}$ , which can be any set of real numbers.
- b) Prove that an information expression involving  $X_1, X_2, \dots, X_n$  can be expressed uniquely as a linear combination of  $\mu^*(X_G)$ , where G are nonempty subsets of  $\mathcal{N}_n$ .
- 4. Uniqueness of the canonical form for nonlinear information expressions Consider a function  $f: \Re^k \to \Re$ , where  $k = 2^n - 1$  such that  $\{\mathbf{h} \in \Re^k :$  $f(\mathbf{h}) = 0$  has zero Lebesgue measure.
  - a) Prove that f cannot be identically zero on  $\Gamma_n^*$ .
  - b) Use the result in a) to show the uniqueness of the canonical form for the class of information expressions of the form  $q(\mathbf{h})$  where q is a polynomial.

(Yeung [360].)

5. Prove that under the constraint  $\mathbf{Qh} = 0$ , if neither  $\mathbf{b}^{\top}$  nor  $\mathbf{c}^{\top}$  is in the row space of **Q**, the identities  $\mathbf{b}^{\top}\mathbf{h} = 0$  and  $\mathbf{c}^{\top}\mathbf{h} = 0$  are equivalent if and only if (13.61) has a unique solution.

### **Historical Notes**

The uniqueness of the canonical form for linear information expressions was first proved by Han [133]. The same result was independently obtained in

327

### 328 13 Information Inequalities

the book by Csiszár and Körner [75]. The geometrical framework for information inequalities is due to Yeung [360]. The characterization of equivalent constrained inequalities in Section 13.4 first appeared in the book by Yeung [361].

# Shannon-Type Inequalities

The basic inequalities form the most important set of information inequalities. In fact, almost all the information inequalities known to date are implied by the basic inequalities. These are called *Shannon-type inequalities*. In this chapter, we show that verification of Shannon-type inequalities can be formulated as a linear programming problem, thus enabling machine-proving of all such inequalities.

### 14.1 The Elemental Inequalities

Consider the conditional mutual information

$$I(X, Y; X, Z, U|Z, T),$$
 (14.1)

in which the random variables X and Z appear more than once. It is readily seen that I(X,Y;X,Z,U|Z,T) can be written as

$$H(X|Z,T) + I(Y;U|X,Z,T),$$
(14.2)

where in both H(X|Z,T) and I(Y;U|X,Z,T), each random variable appears only once.

A Shannon's information measure is said to be *reducible* if there exists a random variable which appears more than once in the information measure, otherwise the information measure is said to be *irreducible*. Without loss of generality, we will consider irreducible Shannon's information measures only, because a reducible Shannon's information measure can always be written as the sum of irreducible Shannon's information measures.

The nonnegativity of all Shannon's information measures form a set of inequalities called the basic inequalities. The set of basic inequalities, however, is not minimal in the sense that some basic inequalities are implied by the others. For example,

$$H(X|Y) \ge 0 \tag{14.3}$$

330 14 Shannon-Type Inequalities

and

$$I(X;Y) \ge 0, \tag{14.4}$$

which are both basic inequalities involving random variables X and Y, imply

$$H(X) = H(X|Y) + I(X;Y) \ge 0, \tag{14.5}$$

which again is a basic inequality involving X and Y.

Let  $\mathcal{N}_n = \{1, 2, \dots, n\}$ , where  $n \geq 2$ . Unless otherwise specified, all information expressions in this chapter involve some or all of the random variables  $X_1, X_2, \dots, X_n$ . The value of n will be specified when necessary. Through application of the identities

$$H(X) = H(X|Y) + I(X;Y)$$
 (14.6)

$$H(X) = H(X) + I(X, T)$$

$$H(X,Y) = H(X) + H(Y|X)$$

$$I(X;Y,Z) = I(X;Y) + I(X;Z|Y)$$

$$H(X|Z) = H(X|Y,Z) + I(X;Y|Z)$$
(14.9)

$$I(X;Y,Z) = I(X;Y) + I(X;Z|Y)$$
(14.8)

$$H(X|Z) = H(X|Y,Z) + I(X;Y|Z)$$
(14.9)

$$H(X, Y|Z) = H(X|Z) + H(Y|X, Z)$$
(14.10)

$$I(X;Y,Z|T) = I(X;Y|T) + I(X;Z|Y,T),$$
(14.11)

any Shannon's information measure can be expressed as the sum of Shannon's information measures of the following two *elemental forms*:

 $\begin{array}{ll} \mathrm{i}) & H(X_i|X_{\mathcal{N}_n-\{i\}}), i\in\mathcal{N}_n\\ \mathrm{ii}) & I(X_i;X_j|X_K), \, \mathrm{where} \,\, i\neq j \,\, \mathrm{and} \,\, K\subset\mathcal{N}_n-\{i,j\}. \end{array}$ 

This will be illustrated in the next example. It is not difficult to check that the total number of the two elemental forms of Shannon's information measures for n random variables is equal to

$$m = n + \binom{n}{2} 2^{n-2}.$$
 (14.12)

The proof of (14.12) is left as an exercise.

Example 14.1. We can expand  $H(X_1, X_2)$  into a sum of elemental forms of Shannon's information measures for n = 3 by applying the identities in (14.6) to (14.11) as follows:

$$H(X_1, X_2) = H(X_1) + H(X_2|X_1)$$
(14.13)

$$= H(X_1|X_2, X_3) + I(X_1; X_2, X_3) + H(X_2|X_1, X_3)$$

$$+ I(X_1, X_2|X_1, X_3) + I(X_1, X$$

$$+I(X_2; X_3|X_1)$$

$$= H(X_1|X_2, X_3) + I(X_1; X_2) + I(X_1; X_3|X_2)$$
(14.14)

$$+H(X_2|X_1, X_3) + I(X_2; X_3|X_1).$$
(14.15)

The nonnegativity of the two elemental forms of Shannon's information measures form a proper subset of the set of basic inequalities. We call the m inequalities in this smaller set the *elemental inequalities*. They are equivalent to the basic inequalities because each basic inequality which is not an elemental inequality can be obtained as the sum of a set of elemental inequalities in view of (14.6) to (14.11). This will be illustrated in the next example. The proof for the minimality of the set of elemental inequalities is deferred to Section 14.6.

*Example 14.2.* In the last example, we expressed  $H(X_1, X_2)$  as

$$H(X_1|X_2, X_3) + I(X_1; X_2) + I(X_1; X_3|X_2) + H(X_2|X_1, X_3) + I(X_2; X_3|X_1).$$
(14.16)

All the five Shannon's information measures in the above expression are in elemental form for n = 3. Then the basic inequality

$$H(X_1, X_2) \ge 0 \tag{14.17}$$

can be obtained as the sum of the following elemental inequalities:

$$H(X_1|X_2, X_3) \ge 0 \tag{14.18}$$

$$I(X_1; X_2) \ge 0 \tag{14.19}$$

$$I(X_1; X_3 | X_2) \ge 0 \tag{14.20}$$

$$H(X_2|X_1, X_3) \ge 0 \tag{14.21}$$

$$I(X_2; X_3 | X_1) \ge 0. \tag{14.22}$$

### 14.2 A Linear Programming Approach

Recall from Section 13.2 that any information expression can be expressed uniquely in canonical form, i.e., a linear combination of the  $k = 2^n - 1$  joint entropies involving some or all of the random variables  $X_1, X_2, \dots, X_n$ . If the elemental inequalities are expressed in canonical form, they become linear inequalities in the entropy space  $\mathcal{H}_n$ . Denote this set of inequalities by  $\mathbf{Gh} \geq 0$ , where  $\mathbf{G}$  is an  $m \times k$  matrix, and define

$$\Gamma_n = \{ \mathbf{h} : \mathbf{G}\mathbf{h} \ge 0 \}. \tag{14.23}$$

We first show that  $\Gamma_n$  is a pyramid in the nonnegative orthant of the entropy space  $\mathcal{H}_n$ . Evidently,  $\Gamma_n$  contains the origin. Let  $\mathbf{e}_j, 1 \leq j \leq k$ , be the column k-vector whose *j*th component is equal to 1 and all the other components are equal to 0. Then the inequality

$$\mathbf{e}_{i}^{\top}\mathbf{h} \ge 0 \tag{14.24}$$

#### 332 14 Shannon-Type Inequalities

corresponds to the nonnegativity of a joint entropy, which is a basic inequality. Since the set of elemental inequalities is equivalent to the set of basic inequalities, if  $\mathbf{h} \in \Gamma_n$ , i.e.,  $\mathbf{h}$  satisfies all the elemental inequalities, then  $\mathbf{h}$  also satisfies the basic inequality in (14.24). In other words,

$$\Gamma_n \subset \{\mathbf{h} : \mathbf{e}_i^{\top} \mathbf{h} \ge 0\} \tag{14.25}$$

for all  $1 \leq j \leq k$ . This implies that  $\Gamma_n$  is in the nonnegative orthant of the entropy space. Since  $\Gamma_n$  contains the origin and the constraints  $\mathbf{Gh} \geq 0$  are linear, we conclude that  $\Gamma_n$  is a pyramid in the nonnegative orthant of  $\mathcal{H}_n$ .

Since the elemental inequalities are satisfied by the entropy function of any *n* random variables  $X_1, X_2, \dots, X_n$ , for any **h** in  $\Gamma_n^*$ , **h** is also in  $\Gamma_n$ , i.e.,

$$\Gamma_n^* \subset \Gamma_n. \tag{14.26}$$

Therefore, for any unconstrained inequality  $f \ge 0$ , if

$$\Gamma_n \subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\},\tag{14.27}$$

then

$$\Gamma_n^* \subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\},\tag{14.28}$$

i.e.,  $f \ge 0$  always holds. In other words, (14.27) is a sufficient condition for  $f \ge 0$  to always hold. Moreover, an inequality  $f \ge 0$  such that (14.27) is satisfied is implied by the basic inequalities, because if **h** satisfies the basic inequalities, i.e.,  $\mathbf{h} \in \Gamma_n$ , then **h** satisfies  $f(\mathbf{h}) \ge 0$ .

For constrained inequalities, following our discussion in Section 13.3, we impose the constraint

$$\mathbf{Qh} = 0 \tag{14.29}$$

and let

$$\boldsymbol{\Phi} = \{ \mathbf{h} : \mathbf{Q}\mathbf{h} = 0 \}. \tag{14.30}$$

For an inequality  $f \ge 0$ , if

$$(\Gamma_n \cap \Phi) \subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\},\tag{14.31}$$

then by (14.26),

$$(\Gamma_n^* \cap \Phi) \subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\},\tag{14.32}$$

i.e.,  $f \geq 0$  always holds under the constraint  $\Phi$ . In other words, (14.31) is a sufficient condition for  $f \geq 0$  to always hold under the constraint  $\Phi$ . Moreover, an inequality  $f \geq 0$  under the constraint  $\Phi$  such that (14.31) is satisfied is implied by the basic inequalities and the constraint  $\Phi$ , because if  $\mathbf{h} \in \Phi$  and  $\mathbf{h}$  satisfies the basic inequalities, i.e.,  $\mathbf{h} \in \Gamma_n \cap \Phi$ , then  $\mathbf{h}$  satisfies  $f(\mathbf{h}) \geq 0$ .

### 14.2.1 Unconstrained Inequalities

To check whether an unconstrained inequality  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is a Shannon-type inequality, we need to check whether  $\Gamma_n$  is a subset of  $\{\mathbf{h} : \mathbf{b}^{\top}\mathbf{h} \geq 0\}$ . The following theorem induces a computational procedure for this purpose.

**Theorem 14.3.**  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is a Shannon-type inequality if and only if the minimum of the problem

Minimize 
$$\mathbf{b}^{\top}\mathbf{h}$$
, subject to  $\mathbf{G}\mathbf{h} \ge 0$  (14.33)

is zero. In this case, the minimum occurs at the origin.

**Remark** The idea of this theorem is illustrated in Figure 14.1 and Figure 14.2. In Figure 14.1,  $\Gamma_n$  is contained in  $\{\mathbf{h} : \mathbf{b}^\top \mathbf{h} \ge 0\}$ . The minimum

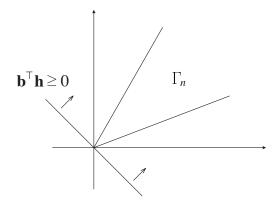


Fig. 14.1.  $\Gamma_n$  is contained in  $\{\mathbf{h} : \mathbf{b}^\top \mathbf{h} \ge 0\}$ .

of  $\mathbf{b}^{\top}\mathbf{h}$  subject to  $\Gamma_n$  occurs at the origin with the minimum equal to 0. In Figure 14.2,  $\Gamma_n$  is not contained in  $\{\mathbf{h} : \mathbf{b}^{\top}\mathbf{h} \ge 0\}$ . The minimum of  $\mathbf{b}^{\top}\mathbf{h}$ subject to  $\Gamma_n$  is  $-\infty$ . A formal proof of the theorem is given next.

Proof of Theorem 14.3. We have to prove that  $\Gamma_n$  is a subset of  $\{\mathbf{h} : \mathbf{b}^{\top} \mathbf{h} \ge 0\}$  if and only if the minimum of the problem in (14.33) is zero. First of all, since  $0 \in \Gamma_n$  and  $\mathbf{b}^{\top} 0 = 0$  for any  $\mathbf{b}$ , the minimum of the problem in (14.33) is at most 0. Assume  $\Gamma_n$  is a subset of  $\{\mathbf{h} : \mathbf{b}^{\top} \mathbf{h} \ge 0\}$  and the minimum of the problem in (14.33) is negative. Then there exists an  $\mathbf{h} \in \Gamma_n$  such that

$$\mathbf{b}^{\top}\mathbf{h} < 0, \tag{14.34}$$

which implies

$$\Gamma_n \not\subset \{\mathbf{h} : \mathbf{b}^\top \mathbf{h} \ge 0\},\tag{14.35}$$

334 14 Shannon-Type Inequalities

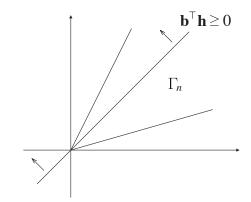


Fig. 14.2.  $\Gamma_n$  is not contained in  $\{\mathbf{h} : \mathbf{b}^\top \mathbf{h} \ge 0\}$ .

which is a contradiction. Therefore, if  $\Gamma_n$  is a subset of  $\{\mathbf{h} : \mathbf{b}^\top \mathbf{h} \ge 0\}$ , then the minimum of the problem in (14.33) is zero.

To prove the converse, assume  $\Gamma_n$  is not a subset of  $\{\mathbf{h} : \mathbf{b}^\top \mathbf{h} \ge 0\}$ , i.e. (14.35) is true. Then there exists an  $\mathbf{h} \in \Gamma_n$  such that

$$\mathbf{b}^{\top}\mathbf{h} < 0. \tag{14.36}$$

This implies that the minimum of the problem in (14.33) is negative, i.e., it is not equal to zero.

Finally, if the minimum of the problem in (14.33) is zero, since the  $\Gamma_n$  contains the origin and  $\mathbf{b}^{\top} 0 = 0$ , the minimum occurs at the origin.  $\Box$ 

By virtue of this theorem, to check whether  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is an unconstrained Shannon-type inequality, all we need to do is to apply the *optimality test* of the *simplex method* [80] to check whether the point  $\mathbf{h} = 0$  is optimal for the minimization problem in (14.33). If  $\mathbf{h} = 0$  is optimal, then  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is an unconstrained Shannon-type inequality, otherwise it is not.

#### 14.2.2 Constrained Inequalities and Identities

To check whether an inequality  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  under the constraint  $\boldsymbol{\Phi}$  is a Shannontype inequality, we need to check whether  $\Gamma_n \cap \boldsymbol{\Phi}$  is a subset of  $\{\mathbf{h} : \mathbf{b}^{\top}\mathbf{h} \geq 0\}$ .

**Theorem 14.4.**  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is a Shannon-type inequality under the constraint  $\Phi$  if and only if the minimum of the problem

Minimize 
$$\mathbf{b}^{\top}\mathbf{h}$$
, subject to  $\mathbf{G}\mathbf{h} \ge 0$  and  $\mathbf{Q}\mathbf{h} = 0$  (14.37)

is zero. In this case, the minimum occurs at the origin.

The proof of this theorem is similar to that for Theorem 14.3, so it is omitted. By taking advantage of the linear structure of the constraint  $\Phi$ , we can reformulate the minimization problem in (14.37) as follows. Let r be the rank of **Q**. Since **h** is in the null space of **Q**, we can write

$$\mathbf{h} = \mathbf{Q}\mathbf{h}',\tag{14.38}$$

where  $\tilde{\mathbf{Q}}$  is a  $k \times (k-r)$  matrix such that the rows of  $\tilde{\mathbf{Q}}^{\top}$  form a basis of the orthogonal complement of the row space of  $\mathbf{Q}$ , and  $\mathbf{h}'$  is a column (k-r)-vector. Then the elemental inequalities can be expressed as

$$\mathbf{G}\mathbf{Q}\mathbf{h}' \ge 0, \tag{14.39}$$

and in terms of  $\mathbf{h}'$ ,  $\Gamma_n$  becomes

$$\Gamma'_{n} = \{ \mathbf{h}' : \mathbf{G}\tilde{\mathbf{Q}}\mathbf{h}' \ge 0 \}, \tag{14.40}$$

which is a pyramid in  $\Re^{k-r}$  (but not necessarily in the nonnegative orthant). Likewise,  $\mathbf{b}^{\top}\mathbf{h}$  can be expressed as  $\mathbf{b}^{\top}\tilde{\mathbf{Q}}\mathbf{h}'$ .

With all the information expressions in terms of  $\mathbf{h}'$ , the problem in (14.37) becomes

Minimize 
$$\mathbf{b}^{\top}\mathbf{Qh}'$$
, subject to  $\mathbf{GQh}' \ge 0.$  (14.41)

Therefore, to check whether  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is a Shannon-type inequality under the constraint  $\Phi$ , all we need to do is to apply the optimality test of the simplex method to check whether the point  $\mathbf{h}' = 0$  is optimal for the problem in (14.41). If  $\mathbf{h}' = 0$  is optimal, then  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is a Shannon-type inequality under the constraint  $\Phi$ , otherwise it is not.

By imposing the constraint  $\Phi$ , the number of elemental inequalities remains the same, while the dimension of the problem decreases from k to k - r.

Finally, to verify that  $\mathbf{b}^{\top}\mathbf{h} = 0$  is a Shannon-type identity under the constraint  $\boldsymbol{\Phi}$ , i.e.,  $\mathbf{b}^{\top}\mathbf{h} = 0$  is implied by the basic inequalities, all we need to do is to verify that both  $\mathbf{b}^{\top}\mathbf{h} \ge 0$  and  $\mathbf{b}^{\top}\mathbf{h} \le 0$  are Shannon-type inequalities under the constraint  $\boldsymbol{\Phi}$ .

### 14.3 A Duality

A nonnegative linear combination is a linear combination whose coefficients are all nonnegative. It is clear that a nonnegative linear combination of basic inequalities is a Shannon-type inequality. However, it is not clear that all Shannon-type inequalities are of this form. By applying the *duality theorem* in linear programming [303], we will see that this is in fact the case.

The *dual* of the *primal* linear programming problem in (14.33) is

Maximize 
$$\mathbf{y}^{\top} \cdot \mathbf{0}$$
 subject to  $\mathbf{y} \ge 0$  and  $\mathbf{y}^{\top} \mathbf{G} \le \mathbf{b}^{\top}$ , (14.42)

336 14 Shannon-Type Inequalities

where

$$\mathbf{y} = [y_1 \cdots y_m]^\top. \tag{14.43}$$

By the duality theorem, if the minimum of the primal problem is zero, which happens when  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is a Shannon-type inequality, the maximum of the dual problem is also zero. Since the cost function in the dual problem is zero, the maximum of the dual problem is zero if and only if the feasible region

$$\Psi = \{ \mathbf{y} : \mathbf{y} \ge 0 \text{ and } \mathbf{y}^{\top} \mathbf{G} \le \mathbf{b}^{\top} \}$$
(14.44)

is nonempty.

**Theorem 14.5.**  $\mathbf{b}^{\top}\mathbf{h} \geq 0$  is a Shannon-type inequality if and only if  $\mathbf{b}^{\top} = \mathbf{x}^{\top}\mathbf{G}$  for some  $\mathbf{x} \geq 0$ , where  $\mathbf{x}$  is a column *m*-vector, i.e.,  $\mathbf{b}^{\top}$  is a nonnegative linear combination of the rows of  $\mathbf{G}$ .

*Proof.* We have to prove that  $\Psi$  is nonempty if and only if  $\mathbf{b}^{\top} = \mathbf{x}^{\top} \mathbf{G}$  for some  $\mathbf{x} \ge 0$ . The feasible region  $\Psi$  is nonempty if and only if

$$\mathbf{b}^{\top} \ge \mathbf{z}^{\top} \mathbf{G} \tag{14.45}$$

for some  $\mathbf{z} \ge 0$ , where  $\mathbf{z}$  is a column *m*-vector. Consider any  $\mathbf{z}$  which satisfies (14.45), and let

$$\mathbf{s}^{\top} = \mathbf{b}^{\top} - \mathbf{z}^{\top} \mathbf{G} \ge 0. \tag{14.46}$$

Denote by  $\mathbf{e}_j$  the column k-vector whose *j*th component is equal to 1 and all the other components are equal to 0,  $1 \leq j \leq k$ . Then  $\mathbf{e}_j^\top \mathbf{h}$  is a joint entropy. Since every joint entropy can be expressed as the sum of elemental forms of Shannon's information measures,  $\mathbf{e}_j^\top$  can be expressed as a nonnegative linear combination of the rows of **G**. Write

$$\mathbf{s} = \left[ s_1 \ s_2 \ \cdots \ s_k \right]^{\top}, \tag{14.47}$$

where  $s_j \ge 0$  for all  $1 \le j \le k$ . Then

$$\mathbf{s}^{\top} = \sum_{j=1}^{k} s_j \mathbf{e}_j^{\top} \tag{14.48}$$

can also be expressed as a nonnegative linear combinations of the rows of **G**, i.e.,

$$\mathbf{s}^{\top} = \mathbf{w}^{\top} \mathbf{G} \tag{14.49}$$

for some  $\mathbf{w} \geq 0$ . From (14.46), we see that

$$\mathbf{b}^{\top} = (\mathbf{w}^{\top} + \mathbf{z}^{\top})\mathbf{G} = \mathbf{x}^{\top}\mathbf{G}, \qquad (14.50)$$

where  $\mathbf{x} \geq 0$ . The proof is accomplished.  $\Box$ 

From this theorem, we see that all Shannon-type inequalities are actually trivially implied by the basic inequalities! However, the verification of a Shannontype inequality requires a computational procedure as described in the last section.

### 14.4 Machine Proving – ITIP

Theorems 14.3 and 14.4 transform the problem of verifying a Shannon-type inequality into a linear programming problem. This enables machine-proving of all Shannon-type inequalities. A software package called ITIP<sup>1</sup>, which runs on MATLAB, has been developed for this purpose. The most updated versions of ITIP can be downloaded from the World Wide Web [366].

Using ITIP is very simple and intuitive. The following examples illustrate the use of ITIP:

1. >> ITIP('H(XYZ) <= H(X) + H(Y) + H(Z)')
True
2. >> ITIP('I(X;Z) = 0','I(X;Z|Y) = 0','I(X;Y) = 0')
True
3. >> ITIP('I(Z;U) - I(Z;U|X) - I(Z;U|Y) <=
0.5 I(X;Y) + 0.25 I(X;ZU) + 0.25 I(Y;ZU)')
Not provable by ITIP</pre>

In the first example, we prove an unconstrained inequality. In the second example, we prove that X and Z are independent if  $X \to Y \to Z$  forms a Markov chain and X and Y are independent. The first identity is what we want to prove, while the second and the third expressions specify the Markov chain  $X \to Y \to Z$  and the independency of X and Y, respectively. In the third example, ITIP returns the clause "Not provable by ITIP," which means that the inequality is not a Shannon-type inequality. This, however, does not mean that the inequality to be proved cannot always hold. In fact, this inequality is one of the known non-Shannon-type inequalities which will be discussed in Chapter 15.

We note that most of the results we have previously obtained by using information diagrams can also be proved by ITIP. However, the advantage of using information diagrams is that one can visualize the structure of the problem. Therefore, the use of information diagrams and ITIP very often complement each other. In the rest of the section, we give a few examples which demonstrate the use of ITIP. The features of ITIP are described in details in the **readme** file.

*Example 14.6.* By Proposition 2.10, the long Markov chain  $X \to Y \to Z \to T$  implies the two short Markov chains  $X \to Y \to Z$  and  $Y \to Z \to T$ . We want to see whether the two short Markov chains also imply the long Markov chain. If so, they are equivalent to each other.

Using ITIP, we have

```
>> ITIP('X/Y/Z/T', 'X/Y/Z', 'Y/Z/T')
Not provable by ITIP
```

<sup>&</sup>lt;sup>1</sup> ITIP stands for Information-Theoretic Inequality Prover.

#### 338 14 Shannon-Type Inequalities

In the above, we have used a macro in ITIP to specify the three Markov chains. The above result from ITIP says that the long Markov chain cannot be proved from the two short Markov chains by means of the basic inequalities. This strongly suggests that the two short Markov chains is weaker than the long Markov chain. However, in order to prove that this is in fact the case, we need an explicit construction of a joint distribution for X, Y, Z, and T which satisfies the two short Markov chains but not the long Markov chain. Toward this end, we resort to the information diagram in Figure 14.3. The Markov

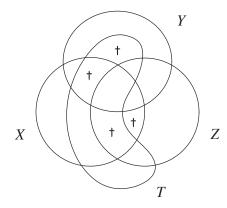


Fig. 14.3. The information diagram for X, Y, Z, and T in Example 14.6.

chain  $X \to Y \to Z$  is equivalent to I(X; Z|Y) = 0, i.e.,

$$\mu^*(\tilde{X} \cap \tilde{Y}^c \cap \tilde{Z} \cap \tilde{T}) + \mu^*(\tilde{X} \cap \tilde{Y}^c \cap \tilde{Z} \cap \tilde{T}^c) = 0.$$
(14.51)

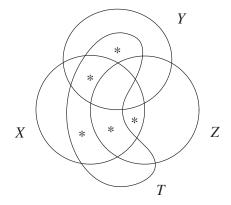
Similarly, the Markov chain  $Y \to Z \to T$  is equivalent to

$$\mu^*(\tilde{X} \cap \tilde{Y} \cap \tilde{Z}^c \cap \tilde{T}) + \mu^*(\tilde{X}^c \cap \tilde{Y} \cap \tilde{Z}^c \cap \tilde{T}) = 0.$$
(14.52)

The four atoms involved in the constraints (14.51) and (14.52) are marked by a dagger in Figure 14.3. In Section 3.5, we have seen that the Markov chain  $X \to Y \to Z \to T$  holds if and only if  $\mu^*$  takes zero value on the set of atoms in Figure 14.4 which are marked with an asterisk<sup>2</sup>. Comparing Figure 14.3 and Figure 14.4, we see that the only atom marked in Figure 14.4 but not in Figure 14.3 is  $\tilde{X} \cap \tilde{Y}^c \cap \tilde{Z}^c \cap \tilde{T}$ . Thus if we can construct a  $\mu^*$  such that it takes zero value on all atoms except for  $\tilde{X} \cap \tilde{Y}^c \cap \tilde{Z}^c \cap \tilde{T}$ , then the corresponding joint distribution satisfies the two short Markov chains but not the long Markov chain. This would show that the two short Markov chains are in fact weaker than the long Markov chain. Following Theorem 3.11, such a  $\mu^*$  can be constructed.

 $<sup>^{2}</sup>$  This information diagram is essentially a reproduction of Figure 3.8.

#### 14.4 Machine Proving – ITIP 339



**Fig. 14.4.** The atoms of  $\mathcal{F}_4$  on which  $\mu^*$  vanishes when  $X \to Y \to Z \to T$  forms a Markov chain.

In fact, the required joint distribution can be obtained by simply letting X = T = U, where U is any random variable such that H(U) > 0, and letting Y and Z be degenerate random variables taking constant values. Then it is easy to see that  $X \to Y \to Z$  and  $Y \to Z \to T$  hold, while  $X \to Y \to Z \to T$  does not hold.

*Example 14.7.* The data processing theorem says that if  $X \to Y \to Z \to T$  forms a Markov chain, then

$$I(Y;Z) \ge I(X;T).$$
 (14.53)

We want to see whether this inequality holds under the weaker condition that  $X \to Y \to Z$  and  $Y \to Z \to T$  form two short Markov chains. By using ITIP, we can show that (14.53) is not a Shannon-type inequality under the Markov conditions

$$I(X;Z|Y) = 0 (14.54)$$

and

$$I(Y;T|Z) = 0. (14.55)$$

This strongly suggests that (14.53) does not always hold under the constraint of the two short Markov chains. However, this has to be proved by an explicit construction of a joint distribution for X, Y, Z, and T which satisfies (14.54)and (14.55) but not (14.53). The construction at the end of the last example serves this purpose.

Example 14.8 (Secret Sharing [40][290]). Consider the following secret sharing problem. Let S be a secret to be encoded into three pieces, X, Y, and Z. The scheme has to satisfy the following two secret sharing requirements:

1. S can be recovered from any two of the three encoded pieces.

- 340 14 Shannon-Type Inequalities
- 2. No information about S can be obtained from any one of the three encoded pieces.

The first requirement is equivalent to the constraints

$$H(S|X,Y) = H(S|Y,Z) = H(S|X,Z) = 0,$$
(14.56)

while the second requirement is equivalent to the constraints

$$I(S;X) = I(S;Y) = I(S;Z) = 0.$$
(14.57)

Since the secret S can be recovered if all X, Y, and Z are known,

$$H(X) + H(Y) + H(Z) \ge H(S).$$
 (14.58)

We are naturally interested in the maximum constant c which satisfies

$$H(X) + H(Y) + H(Z) \ge cH(S).$$
 (14.59)

We can explore the possible values of c by ITIP. After a few trials, we find that ITIP returns a "True" for all  $c \leq 3$ , and returns the clause "Not provable by ITIP" for any c slightly larger than 3, say 3.0001. This means that the maximum value of c is lower bounded by 3. This lower bound is in fact tight, as we can see from the following construction. Let S and N be mutually independent ternary random variables uniformly distributed on  $\{0, 1, 2\}$ , and define

$$X = N \tag{14.60}$$

$$Y = S + N \mod 3,\tag{14.61}$$

and

$$Z = S + 2N \mod 3.$$
(14.62)

Then it is easy to verify that

$$S = Y - X \mod 3 \tag{14.63}$$

$$= 2Y - Z \mod 3 \tag{14.64}$$

$$= Z - 2X \mod 3.$$
 (14.65)

Thus the requirements in (14.56) are satisfied. It is also readily verified that the requirements in (14.57) are satisfied. Finally, all S, X, Y, and Z distribute uniformly on  $\{0, 1, 2\}$ . Therefore,

$$H(X) + H(Y) + H(Z) = 3H(S).$$
(14.66)

This proves that the maximum constant c which satisfies (14.59) is 3.

Using the approach in this example, almost all information-theoretic bounds reported in the literature for this class of problems can be obtained when a definite number of random variables are involved.

### 14.5 Tackling the Implication Problem

We have already mentioned in Section 13.5 that the implication problem of conditional independence is extremely difficult except for the special case that only full conditional mutual independencies are involved. In this section, we employ the tools we have developed in this chapter to tackle this problem.

In Bayesian network (see [258]), the following four axioms are often used for proving implications of conditional independencies:

• Symmetry:

$$X \perp Y | Z \Leftrightarrow Y \perp X | Z \tag{14.67}$$

• Decomposition:

$$X \perp (Y,T)|Z \Rightarrow (X \perp Y|Z) \& (X \perp T|Z)$$
(14.68)

• Weak Union:

$$X \perp (Y,T)|Z \Rightarrow X \perp Y|Z,T \tag{14.69}$$

• Contraction:

$$(X \perp Y|Z) \& (X \perp T|Y,Z) \Rightarrow X \perp (Y,T)|Z.$$
(14.70)

These axioms form a system called *semi-graphoid* and were first proposed by Dawid [82] as heuristic properties of conditional independence.

The axiom of symmetry is trivial in the context of probability<sup>3</sup>. The other three axioms can be summarized by

$$X \perp (Y,T)|Z \Leftrightarrow (X \perp Y|Z) \& (X \perp T|Y,Z).$$
(14.71)

This can easily be proved as follows. Consider the identity

$$I(X;Y,T|Z) = I(X;Y|Z) + I(X;T|Y,Z).$$
(14.72)

Since conditional mutual informations are always nonnegative by the basic inequalities, if I(X; Y, T|Z) vanishes, I(X; Y|Z) and I(X; T|Y, Z) also vanish, and vice versa. This proves (14.71). In other words, (14.71) is the result of a specific application of the basic inequalities. Therefore, any implication which can be proved by invoking these four axioms can also be proved by ITIP.

In fact, ITIP is considerably more powerful than the above four axioms. This will be shown in the next example in which we give an implication which can be proved by ITIP but not by these four axioms<sup>4</sup>. We will see some implications which cannot be proved by ITIP when we discuss non-Shannon-type inequalities in the next chapter.

<sup>&</sup>lt;sup>3</sup> These four axioms can be used beyond the context of probability.

 $<sup>^4</sup>$  This example is due to Zhen Zhang, private communication.

#### 342 14 Shannon-Type Inequalities

Example 14.9. We will show that

$$\begin{array}{l}
I(X;Y|Z) = 0 \\
I(X;T|Z) = 0 \\
I(X;T|Y) = 0 \\
I(X;Z|Y) = 0 \\
I(X;Z|T) = 0
\end{array} \Rightarrow I(X;Y|T) = 0 \tag{14.73}$$

can be proved by invoking the basic inequalities. First, we write

$$I(X;Y|Z) = I(X;Y|Z,T) + I(X;Y;T|Z).$$
(14.74)

Since I(X; Y|Z) = 0 and  $I(X; Y|Z, T) \ge 0$ , we let

$$I(X;Y|Z,T) = a \tag{14.75}$$

for some nonnegative real number a, so that

$$I(X;Y;T|Z) = -a (14.76)$$

from (14.74). In the information diagram in Figure 14.5, we mark the atom I(X;Y|Z,T) by a "+" and the atom I(X;Y;T|Z) by a "-." Then we write

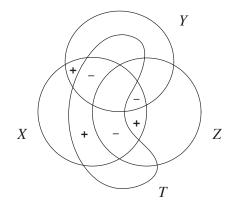


Fig. 14.5. The information diagram for X, Y, Z, and T.

$$I(X;T|Z) = I(X;Y;T|Z) + I(X;T|Y,Z).$$
(14.77)

Since I(X;T|Z) = 0 and I(X;Y;T|Z) = -a, we get

$$I(X;T|Y,Z) = a.$$
 (14.78)

In the information diagram, we mark the atom I(X;T|Y,Z) with a "+." Continue in this fashion, the five CI's on the left hand side of (14.73) imply that all the atoms marked with a "+" in the information diagram take the value a, while all the atoms marked with a "-" take the value -a. From the information diagram, we see that

$$I(X;Y|T) = I(X;Y;Z|T) + I(X;Y|Z,T) = (-a) + a = 0,$$
(14.79)

which proves our claim. Since we base our proof on the basic inequalities, this implication can also be proved by ITIP.

Due to the form of the five given CI's in (14.73), none of the axioms in (14.68) to (14.70) can be applied. Thus we conclude that the implication in (14.73) cannot be proved by the four axioms in (14.67) to (14.70).

### 14.6 Minimality of the Elemental Inequalities

We have already seen in Section 14.1 that the set of basic inequalities is not minimal in the sense that in the set, some inequalities are implied by the others. We then showed that the set of basic inequalities is equivalent to the smaller set of elemental inequalities. Again, we can ask whether the set of elemental inequalities is minimal.

In this section, we prove that the set of elemental inequalities is minimal. This result is important for efficient implementation of ITIP because it says that we cannot consider a smaller set of inequalities. The proof, however, is rather technical. The reader may skip this proof without missing the essence of this chapter.

The elemental inequalities in set-theoretic notations have one of the following two forms:

1. 
$$\mu(\tilde{X}_i - \tilde{X}_{\mathcal{N}_n - \{i\}}) \ge 0,$$
  
2.  $\mu(\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K) \ge 0, i \ne j \text{ and } K \subset \mathcal{N}_n - \{i, j\},$ 

where  $\mu$  denotes a set-additive function defined on  $\mathcal{F}_n$ . They will be referred to as  $\alpha$ -inequalities and  $\beta$ -inequalities, respectively.

We are to show that all the elemental inequalities are nonredundant, i.e., none of them is implied by the others. For an  $\alpha$ -inequality

$$\mu(\ddot{X}_i - \ddot{X}_{\mathcal{N}_n - \{i\}}) \ge 0, \tag{14.80}$$

since it is the only elemental inequality which involves the atom  $\tilde{X}_i - \tilde{X}_{\mathcal{N}_n - \{i\}}$ , it is clearly not implied by the other elemental inequalities. Therefore we only need to show that all  $\beta$ -inequalities are nonredundant. To show that a  $\beta$ inequality is nonredundant, it suffices to show that there exists a measure  $\hat{\mu}$  on  $\mathcal{F}_n$  which satisfies all other elemental inequalities except for that  $\beta$ -inequality.

We will show that the  $\beta$ -inequality

$$\mu(\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K) \ge 0 \tag{14.81}$$

#### 344 14 Shannon-Type Inequalities

is nonredundant. To facilitate our discussion, we denote  $\mathcal{N}_n - K - \{i, j\}$  by L(i, j, K), and we let  $C_{ij|K}(S), S \subset L(i, j, K)$  be the atoms in  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ , where

$$C_{ij|K}(S) = \tilde{X}_i \cap \tilde{X}_j \cap \tilde{X}_S \cap \tilde{X}_K^c \cap \tilde{X}_{L(i,j,K)-S}^c.$$
(14.82)

We first consider the case when  $L(i, j, K) = \emptyset$ , i.e.,  $K = \mathcal{N}_n - \{i, j\}$ . We construct a measure  $\hat{\mu}$  by

$$\hat{\mu}(A) = \begin{cases} -1 \text{ if } A = \tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K \\ 1 \text{ otherwise,} \end{cases}$$
(14.83)

where  $A \in \mathcal{A}$ . In other words,  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$  is the only atom with measure -1; all other atoms have measure 1. Then  $\hat{\mu}(\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K) < 0$  is trivially true. It is also trivial to check that for any  $i' \in \mathcal{N}_n$ ,

$$\hat{\mu}(\tilde{X}_{i'} - \tilde{X}_{\mathcal{N}_n - \{i'\}}) = 1 \ge 0, \qquad (14.84)$$

and for any  $(i', j', K') \neq (i, j, K)$  such that  $i' \neq j'$  and  $K' \subset \mathcal{N}_n - \{i', j'\}$ ,

$$\hat{\mu}(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) = 1 \ge 0 \tag{14.85}$$

if  $K' = \mathcal{N}_n - \{i', j'\}$ . On the other hand, if K' is a proper subset of  $\mathcal{N}_n - \{i', j'\}$ , then  $\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}$  contains at least two atoms, and therefore

$$\hat{\mu}(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) \ge 0.$$
(14.86)

This completes the proof for the  $\beta$ -inequality in (14.81) to be nonredundant when  $L(i, j, K) = \phi$ .

We now consider the case when  $L(i, j, K) \neq \phi$ , or  $|L(i, j, K)| \geq 1$ . We construct a measure  $\hat{\mu}$  as follows. For the atoms in  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ , let

$$\hat{\mu}(C_{ij|K}(S)) = \begin{cases} (-1)^{|S|} - 1 \ S = L(i, j, K) \\ (-1)^{|S|} \ S \neq L(i, j, K). \end{cases}$$
(14.87)

For  $C_{ij|K}(S)$ , if |S| is odd, it is referred to as an *odd atom* of  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ , and if |S| is even, it is referred to as an *even atom* of  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ . For any atom  $A \notin \tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ , we let

$$\hat{\mu}(A) = 1.$$
 (14.88)

This completes the construction of  $\hat{\mu}$ .

We first prove that

$$\hat{\mu}(\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K) < 0. \tag{14.89}$$

Consider

14.6 Minimality of the Elemental Inequalities 345

$$\hat{\mu}(\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K) = \sum_{S \subset L(i,j,K)} \hat{\mu}(C_{ij|K}(S))$$
$$= \left(\sum_{r=0}^{|L(i,j,K)|} \binom{|L(i,j,K)|}{r} (-1)^r\right) - 1$$
$$= -1,$$

where the last equality follows from the binomial formula

$$\sum_{r=0}^{n} \binom{n}{r} (-1)^{r} = 0$$
 (14.90)

for  $n \ge 1$ . This proves (14.89).

Next we prove that  $\hat{\mu}$  satisfies all  $\alpha$ -inequalities. We note that for any  $i' \in \mathcal{N}_n$ , the atom  $\tilde{X}_{i'} - \tilde{X}_{\mathcal{N}_n - \{i'\}}$  is not in  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ . Thus

$$\hat{\mu}(\tilde{X}_{i'} - \tilde{X}_{\mathcal{N}_n - \{i\}}) = 1 \ge 0.$$
(14.91)

It remains to prove that  $\hat{\mu}$  satisfies all  $\beta$ -inequalities except for (14.81), i.e., for any  $(i', j', K') \neq (i, j, K)$  such that  $i' \neq j'$  and  $K' \subset \mathcal{N}_n - \{i', j'\}$ ,

$$\hat{\mu}(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) \ge 0.$$
(14.92)

Consider

$$\hat{\mu}(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) 
= \hat{\mu}((\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) \cap (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)) 
+ \hat{\mu}((\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) - (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)).$$
(14.93)

The nonnegativity of the second term above follows from (14.88). For the first term,

$$(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) \cap (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)$$
(14.94)

is nonempty if and only if

$$\{i', j'\} \cap K = \phi \text{ and } \{i, j\} \cap K' = \phi.$$
 (14.95)

If this condition is not satisfied, then the first term in (14.93) becomes  $\hat{\mu}(\phi) = 0$ , and (14.92) follows immediately.

Let us assume that the condition in (14.95) is satisfied. Then by simple counting, we see that the number atoms in

$$(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) \cap (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)$$
(14.96)

is equal to  $2^{\varphi}$ , where

$$\varphi = n - |\{i, j\} \cup \{i', j'\} \cup K \cup K'|.$$
(14.97)

### 346 14 Shannon-Type Inequalities

For example, for n = 6, there are  $4 = 2^2$  atoms in

$$(\tilde{X}_1 \cap \tilde{X}_2) \cap (\tilde{X}_1 \cap \tilde{X}_3 - \tilde{X}_4), \tag{14.98}$$

namely  $\tilde{X}_1 \cap \tilde{X}_2 \cap \tilde{X}_3 \cap \tilde{X}_4^c \cap Y_5 \cap Y_6$ , where  $Y_i = \tilde{X}_i$  or  $\tilde{X}_i^c$  for i = 5, 6. We check that

$$\varphi = 6 - |\{1, 2\} \cup \{1, 3\} \cup \phi \cup \{4\}| = 2.$$
(14.99)

We first consider the case when  $\varphi = 0$ , i.e.,

$$\mathcal{N}_n = \{i, j\} \cup \{i', j'\} \cup K \cup K'.$$
(14.100)

Then

$$(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) \cap (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)$$
(14.101)

contains exactly one atom. If this atom is an even atom of  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ , then the first term in (14.93) is either 0 or 1 (cf., (14.87)), and (14.92) follows immediately. If this atom is an odd atom of  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$ , then the first term in (14.93) is equal to -1. This happens if and only if  $\{i, j\}$  and  $\{i', j'\}$  have one common element, which implies that  $(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) - (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)$ is nonempty. Therefore the second term in (14.93) is at least 1, and hence (14.92) follows.

Finally, we consider the case when  $\varphi \geq 1$ . Using the binomial formula in (14.90), we see that the number of odd atoms and even atoms of  $\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K$  in

$$(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) \cap (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)$$
(14.102)

are the same. Therefore the first term in (14.93) is equal to -1 if

$$C_{ij|K}(L(i,j,K)) \in \tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'},$$
 (14.103)

and is equal to 0 otherwise. The former is true if and only if  $K' \subset K$ , which implies that  $(\tilde{X}_{i'} \cap \tilde{X}_{j'} - \tilde{X}_{K'}) - (\tilde{X}_i \cap \tilde{X}_j - \tilde{X}_K)$  is nonempty, or that the second term is at least 1. Thus in either case (14.92) is true. This completes the proof that (14.81) is nonredundant.

# Appendix 14.A: The Basic Inequalities and the Polymatroidal Axioms

In this appendix, we show that the basic inequalities for a collection of n random variables  $\Theta = \{X_i, i \in \mathcal{N}_n\}$  is equivalent to the following polymatroidal axioms: For all  $\alpha, \beta \in \mathcal{N}_n$ ,

- P1.  $H_{\Theta}(\emptyset) = 0.$
- P2.  $H_{\Theta}(\alpha) \leq H_{\Theta}(\beta)$  if  $\alpha \subset \beta$ .
- P3.  $H_{\Theta}(\alpha) + H_{\Theta}(\beta) \ge H_{\Theta}(\alpha \cap \beta) + H_{\Theta}(\alpha \cup \beta).$

We first show that the polymatroidal axioms imply the basic inequalities. From P1 and P2, since  $\emptyset \subset \alpha$  for any  $\alpha \subset \mathcal{N}_n$ , we have

$$H_{\Theta}(\alpha) \ge H_{\Theta}(\emptyset) = 0, \tag{14.104}$$

or

$$H(X_{\alpha}) \ge 0. \tag{14.105}$$

This shows that entropy is nonnegative.

In P2, letting  $\gamma = \beta \backslash \alpha$ , we have

$$H_{\Theta}(\alpha) \le H_{\Theta}(\alpha \cup \gamma), \tag{14.106}$$

or

$$H(X_{\gamma}|X_{\alpha}) \ge 0. \tag{14.107}$$

Here,  $\gamma$  and  $\alpha$  are disjoint subsets of  $\mathcal{N}_n$ .

In P3, letting  $\gamma = \beta \setminus \alpha$ ,  $\delta = \alpha \cap \beta$ , and  $\sigma = \alpha \setminus \beta$ , we have

$$H_{\Theta}(\sigma \cup \delta) + H_{\Theta}(\gamma \cup \delta) \ge H_{\Theta}(\delta) + H_{\Theta}(\sigma \cup \delta \cup \gamma), \qquad (14.108)$$

or

$$I(X_{\sigma}; X_{\gamma} | X_{\delta}) \ge 0. \tag{14.109}$$

Again,  $\sigma, \delta$ , and  $\gamma$  are disjoint subsets of  $\mathcal{N}_n$ . When  $\delta = \emptyset$ , from P3, we have

$$I(X_{\sigma}; X_{\gamma}) \ge 0. \tag{14.110}$$

Thus P1 to P3 imply that entropy is nonnegative, and that conditional entropy, mutual information, and conditional mutual information are nonnegative provided that they are irreducible. However, it has been shown in Section 14.1 that a reducible Shannon's information measure can always be written as the sum of irreducible Shannon's information measures. Therefore, we have shown that the polymatroidal axioms P1 to P3 imply the basic inequalities.

The converse is trivial and the proof is omitted.

# Problems

- 1. Prove (14.12) for the total number of elemental forms of Shannon's information measures for n random variables.
- 2. Shannon-type inequalities for n random variables  $X_1, X_2, \dots, X_n$  refer to all information inequalities implied by the basic inequalities for these n random variables. Show that no new information inequality can be generated by considering the basic inequalities for more than n random variables.

- 348 14 Shannon-Type Inequalities
- 3. Show by an example that the decomposition of an information expression into a sum of elemental forms of Shannon's information measures is not unique.
- 4. Elemental forms of conditional independencies Consider random variables  $X_1, X_2, \dots, X_n$ . A conditional independency is said to be elemental if it corresponds to setting an elemental form of Shannon's information measure to zero. Show that any conditional independency involving  $X_1, X_2, \dots, X_n$  is equivalent to a collection of elemental conditional independencies.
- 5. Symmetrical information inequalities
  - a) Show that every symmetrical information expression (cf. Problem 1 in Chapter 13) involving random variable  $X_1, X_2, \dots, X_n$  can be written in the form

$$E = \sum_{k=0}^{n-1} a_k c_k^{(n)}$$

where

$$c_0^{(n)} = \sum_{i=1}^n H(X_i | X_{N-i})$$

and for  $1 \leq k \leq n-1$ ,

$$c_k^{(n)} = \sum_{\substack{1 \le i < j \le n \\ K \subset N - \{i,j\}, |K| = k-1}} I(X_i; X_j | X_K).$$

Note that  $c_0^{(n)}$  is the sum of all Shannon's information measures of the first elemental form, and for  $1 \le k \le n-1$ ,  $c_k^{(n)}$  is the sum of all Shannon's information measures of the second elemental form conditioning on k-1 random variables.

- b) Show that  $E \ge 0$  always holds if  $a_k \ge 0$  for all k.
- c) Show that if  $E \ge 0$  always holds, then  $a_k \ge 0$  for all k. Hint: Construct random variables  $X_1, X_2, \dots, X_n$  for each  $0 \le k \le n-1$  such that  $c_k^{(n)} > 0$  and  $c_{k'}^{(n)} = 0$  for all  $0 \le k' \le n-1$  and  $k' \ne k$ . (Han [134].)
- 6. Strictly positive probability distributions It was shown in Proposition 2.12 that

$$X_1 \perp X_4 | (X_2, X_3) X_1 \perp X_3 | (X_2, X_4)$$
  $\Rightarrow$   $X_1 \perp (X_3, X_4) | X_2$ 

if  $p(x_1, x_2, x_3, x_4) > 0$  for all  $x_1, x_2, x_3$ , and  $x_4$ . Show by using ITIP that this implication is not implied by the basic inequalities. This strongly suggests that this implication does not hold in general, which was shown to be the case by the construction following Proposition 2.12.

7. a) Verify by ITIP that

$$I(X_1, X_2; Y_1, Y_2) \le I(X_1; Y_1) + I(X_2; Y_2)$$

under the constraint  $H(Y_1, Y_2|X_1, X_2) = H(Y_1|X_1) + H(Y_2|X_2)$ . This constrained inequality was used in Problem 10 in Chapter 7 to obtain the capacity of two parallel channels.

b) Verify by ITIP that

$$I(X_1, X_2; Y_1, Y_2) \ge I(X_1; Y_1) + I(X_2; Y_2)$$

under the constraint  $I(X_1; X_2) = 0$ . This constrained inequality was used in Problem 4 in Chapter 8 to obtain the rate-distortion function for a product source.

- 8. Verify by ITIP the information identity in Example 3.18.
- 9. Repeat Problem 10 in Chapter 3 with the help of ITIP.
- Prove the implications in Problem 14 in Chapter 3 by ITIP and show that they cannot be deduced from the semi-graphoidal axioms. (Studený [313].)

## **Historical Notes**

For almost half a century, all information inequalities known in the literature are consequences of the basic inequalities due to Shannon [291]. Fujishige [114] showed that the entropy function is a polymatroid (see Appendix 14.6). Yeung [360] showed that verification of all such inequalities, referred to Shannon-type inequalities, can be formulated as a linear programming problem if the number of random variables involved is fixed. ITIP, a software package for this purpose, was developed by Yeung and Yan [366]. Non-Shannon-type inequalities, which have been discovered only recently, will be discussed in the next chapter.

The recent interest in the implication problem of conditional independence has been fueled by Bayesian networks. For a number of years, researchers in Bayesian networks generally believed that the semi-graphoidal axioms form a complete set of axioms for conditional independence until it was refuted by Studený [313].

In Chapter 13, we introduced the regions  $\Gamma_n^*$  and  $\Gamma_n$  in the entropy space  $\mathcal{H}_n$  for *n* random variables. From  $\Gamma_n^*$ , one in principle can determine whether any information inequality always holds. The region  $\Gamma_n$ , defined by the set of all basic inequalities (equivalently all elemental inequalities) involving *n* random variables, is an outer bound on  $\Gamma_n^*$ . From  $\Gamma_n$ , one can determine whether any information inequality is implied by the basic inequalities. If so, it is called a Shannon-type inequality. Since the basic inequalities always hold, so do all Shannon-type inequalities. In the last chapter, we have shown how machine-proving of all Shannon-type inequalities can be made possible by taking advantage of the linear structure of  $\Gamma_n$ .

If the two regions  $\Gamma_n^*$  and  $\Gamma_n$  are identical, then all information inequalities which always hold are Shannon-type inequalities, and hence all information inequalities can be completely characterized. However, if  $\Gamma_n^*$  is a proper subset of  $\Gamma_n$ , then there exist constraints on an entropy function which are not implied by the basic inequalities. Such a constraint, if in the form of an inequality, is referred to a non-Shannon-type inequality.

There is a point here which needs further explanation. The fact that  $\Gamma_n^* \neq \Gamma_n$  does not necessarily imply the existence of a non-Shannon-type inequality. As an example, suppose  $\Gamma_n$  contains all but an isolated point in  $\Gamma_n^*$ . Then this does not lead to the existence of a non-Shannon-type inequality for n random variables.

In this chapter, we present characterizations of  $\Gamma_n^*$  which are more refined than  $\Gamma_n$ . These characterizations lead to the existence of non-Shannon-type inequalities for  $n \geq 4$ .

# 15.1 Characterizations of $\Gamma_2^*$ , $\Gamma_3^*$ , and $\overline{\Gamma}_n^*$

Recall from the proof of Theorem 3.6 that the vector  $\mathbf{h}$  represents the values of the *I*-Measure  $\mu^*$  on the unions in  $\mathcal{F}_n$ . Moreover,  $\mathbf{h}$  is related to the values of  $\mu^*$  on the atoms of  $\mathcal{F}_n$ , represented as  $\mathbf{u}$ , by

$$\mathbf{h} = \mathbf{C}_n \mathbf{u} \tag{15.1}$$

where  $\mathbf{C}_n$  is a unique  $k \times k$  matrix with  $k = 2^n - 1$  (cf. (3.27)).

Let  $\mathcal{I}_n$  be the k-dimensional Euclidean space with the coordinates labeled by the components of **u**. Note that each coordinate in  $\mathcal{I}_n$  corresponds to the value of  $\mu^*$  on a nonempty atom of  $\mathcal{F}_n$ . Recall from Lemma 13.1 the definition of the region

$$\Psi_n^* = \{ \mathbf{u} \in \mathcal{I}_n : \mathbf{C}_n \mathbf{u} \in \Gamma_n^* \}, \tag{15.2}$$

which is obtained from the region  $\Gamma_n^*$  via the linear transformation induced by  $\mathbf{C}_n^{-1}$ . Analogously, we define the region

$$\Psi_n = \{ \mathbf{u} \in \mathcal{I}_n : \mathbf{C}_n \mathbf{u} \in \mathcal{\Gamma}_n \}.$$
(15.3)

The region  $\Gamma_n^*$ , as we will see, is extremely difficult to characterize for a general n. Therefore, we start our discussion with the simplest case, namely n = 2.

## Theorem 15.1. $\Gamma_2^* = \Gamma_2$ .

*Proof.* For n = 2, the elemental inequalities are

$$H(X_1|X_2) = \mu^*(X_1 - X_2) \ge 0$$
(15.4)

$$H(X_2|X_1) = \mu^* (\tilde{X}_2 - \tilde{X}_1) \ge 0 \tag{15.5}$$

$$I(X_1; X_2) = \mu^*(X_1 \cap X_2) \ge 0.$$
(15.6)

Note that the quantities on the left hand sides above are precisely the values of  $\mu^*$  on the atoms of  $\mathcal{F}_2$ . Therefore,

$$\Psi_2 = \{ \mathbf{u} \in \mathcal{I}_2 : \mathbf{u} \ge 0 \},\tag{15.7}$$

i.e.,  $\Psi_2$  is the nonnegative orthant of  $\mathcal{I}_2$ . Since  $\Gamma_2^* \subset \Gamma_2$ ,  $\Psi_2^* \subset \Psi_2$ . On the other hand,  $\Psi_2 \subset \Psi_2^*$  by Lemma 13.1. Thus  $\Psi_2^* = \Psi_2$ , which implies  $\Gamma_2^* = \Gamma_2$ . The proof is accomplished.  $\Box$ 

Next, we prove that Theorem 15.1 cannot even be generalized to n = 3.

## Theorem 15.2. $\Gamma_3^* \neq \Gamma_3$ .

*Proof.* For n = 3, the elemental inequalities are

$$H(X_i|X_j, X_k) = \mu^* (\tilde{X}_i - \tilde{X}_j - \tilde{X}_k) \ge 0$$
(15.8)

$$I(X_i; X_j | X_k) = \mu^* (X_i \cap X_j - X_k) \ge 0,$$
(15.9)

and

$$I(X_i; X_j) = \mu^* (\tilde{X}_i \cap \tilde{X}_j)$$
(15.10)

$$= \mu^* (X_i \cap X_j \cap X_k) + \mu^* (X_i \cap X_j - X_k)$$
(15.11)

$$\geq 0 \tag{15.12}$$

15.1 Characterizations of  $\Gamma_2^*$ ,  $\Gamma_3^*$ , and  $\overline{\Gamma}_n^*$  353

for  $1 \leq i < j < k \leq 3$ . For  $\mathbf{u} \in \mathcal{I}_3$ , let

$$\mathbf{u} = (u_1, u_2, u_3, u_4, u_5, u_6, u_7), \tag{15.13}$$

where  $u_i, 1 \leq i \leq 7$  correspond to the values

$$\mu^{*}(\tilde{X}_{1} - \tilde{X}_{2} - \tilde{X}_{3}), \ \mu^{*}(\tilde{X}_{2} - \tilde{X}_{1} - \tilde{X}_{3}), \ \mu^{*}(\tilde{X}_{3} - \tilde{X}_{1} - \tilde{X}_{2}), \\
\mu^{*}(\tilde{X}_{1} \cap \tilde{X}_{2} - \tilde{X}_{3}), \ \mu^{*}(\tilde{X}_{1} \cap \tilde{X}_{3} - \tilde{X}_{2}), \ \mu^{*}(\tilde{X}_{2} \cap \tilde{X}_{3} - \tilde{X}_{1}), \\
\mu^{*}(\tilde{X}_{1} \cap \tilde{X}_{2} \cap \tilde{X}_{3}),$$
(15.14)

respectively. These are the values of  $\mu^*$  on the nonempty atoms of  $\mathcal{F}_3$ . Then from (15.8), (15.9), and (15.12), we see that

$$\Psi_3 = \{ \mathbf{u} \in \mathcal{I}_3 : u_i \ge 0, \ 1 \le i \le 6; \ u_j + u_7 \ge 0, \ 4 \le j \le 6 \}.$$
(15.15)

It is easy to check that the point (0, 0, 0, a, a, a, -a) for any  $a \ge 0$  is in  $\Psi_3$ . This is illustrated in Figure 15.1, and it is readily seen that the relations

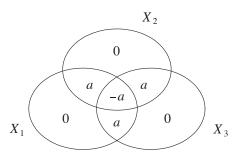


Fig. 15.1. The set-theoretic structure of the point (0, 0, 0, a, a, a, -a) in  $\Psi_3$ .

$$H(X_i|X_j, X_k) = 0 (15.16)$$

and

$$I(X_i; X_j) = 0 (15.17)$$

for  $1 \le i < j < k \le 3$  are satisfied, i.e., each random variable is a function of the other two, and the three random variables are pairwise independent.

Let  $S_{X_i}$  be the support of  $X_i$ , i = 1, 2, 3. For any  $x_1 \in S_{X_1}$  and  $x_2 \in S_{X_2}$ , since  $X_1$  and  $X_2$  are independent, we have

$$p(x_1, x_2) = p(x_1)p(x_2) > 0.$$
(15.18)

Since  $X_3$  is a function of  $X_1$  and  $X_2$ , there is a unique  $x_3 \in \mathcal{S}_{X_3}$  such that

$$p(x_1, x_2, x_3) = p(x_1, x_2) = p(x_1)p(x_2) > 0.$$
 (15.19)

Now since  $X_2$  is a function of  $X_1$  and  $X_3$ , and  $X_1$  and  $X_3$  are independent, we can write

$$p(x_1, x_2, x_3) = p(x_1, x_3) = p(x_1)p(x_3).$$
 (15.20)

Equating (15.19) and (15.20), we have

$$p(x_2) = p(x_3). \tag{15.21}$$

Now consider any  $x'_2 \in \mathcal{S}_{X_2}$  such that  $x'_2 \neq x_2$ . Since  $X_2$  and  $X_3$  are independent, we have

$$p(x'_2, x_3) = p(x'_2)p(x_3) > 0.$$
(15.22)

Since  $X_1$  is a function of  $X_2$  and  $X_3$ , there is a unique  $x'_1 \in \mathcal{S}_{X_1}$  such that

$$p(x'_1, x'_2, x_3) = p(x'_2, x_3) = p(x'_2)p(x_3) > 0.$$
(15.23)

Now since  $X_2$  is a function of  $X_1$  and  $X_3$ , and  $X_1$  and  $X_3$  are independent, we can write

$$p(x'_1, x'_2, x_3) = p(x'_1, x_3) = p(x'_1)p(x_3).$$
(15.24)

Similarly, since  $X_3$  is a function of  $X_1$  and  $X_2$ , and  $X_1$  and  $X_2$  are independent, we can write

$$p(x'_1, x'_2, x_3) = p(x'_1, x'_2) = p(x'_1)p(x'_2).$$
(15.25)

Equating (15.24) and (15.25), we have

$$p(x_2') = p(x_3), \tag{15.26}$$

and from (15.21), we have

$$p(x_2') = p(x_2). \tag{15.27}$$

Therefore  $X_2$  must have a uniform distribution on its support. The same can be proved for  $X_1$  and  $X_3$ . Now from Figure 15.1,

$$H(X_1) = H(X_1|X_2, X_3) + I(X_1; X_2|X_3) + I(X_1; X_3|X_2) +I(X_1; X_2; X_3)$$
(15.28)  
$$= 0 + a + a + (-a)$$
(15.29)

$$=a,$$
(15.30)

and similarly

$$H(X_2) = H(X_3) = a. (15.31)$$

Then the only values that a can take are  $\log M$ , where M (a positive integer) is the cardinality of the supports of  $X_1, X_2$ , and  $X_3$ . In other words, if a is not equal to  $\log M$  for some positive integer M, then the point (0, 0, 0, a, a, a, -a) is not in  $\Psi_3^*$ . This proves that  $\Psi_3^* \neq \Psi_3$ , which implies  $\Gamma_3^* \neq \Gamma_3$ . The theorem is proved.  $\Box$ 

The proof above has the following interpretation. For  $\mathbf{h} \in \mathcal{H}_3$ , let

15.1 Characterizations of 
$$\Gamma_2^*$$
,  $\Gamma_3^*$ , and  $\overline{\Gamma}_n^*$  355

$$\mathbf{h} = (h_1, h_2, h_3, h_{12}, h_{13}, h_{23}, h_{123}). \tag{15.32}$$

From Figure 15.1, we see that the point (0, 0, 0, a, a, a, -a) in  $\Psi_3$  corresponds to the point (a, a, a, 2a, 2a, 2a, 2a) in  $\Gamma_3$ . Evidently, the point (a, a, a, 2a, 2a, 2a, 2a) in  $\Gamma_3$  satisfies the 6 elemental inequalities given in (15.8) and (15.12) for  $1 \leq i < j < k \leq 3$  with equality. Since  $\Gamma_3$  is defined by all the elemental inequalities, the set

$$\{(a, a, a, 2a, 2a, 2a, 2a) \in \Gamma_3 : a \ge 0\}$$
(15.33)

is in the intersection of 6 hyperplanes in  $\mathcal{H}_3$  (i.e.,  $\Re^7$ ) defining the boundary of  $\Gamma_3$ , and hence it defines an extreme direction of  $\Gamma_3$ . Then the proof says that along this extreme direction of  $\Gamma_3$ , only certain discrete points, namely those points with *a* equals log *M* for some positive integer *M*, are entropic. This is illustrated in Figure 15.2. As a consequence, the region  $\Gamma_3^*$  is not convex.



Having proved that  $\Gamma_3^* \neq \Gamma_3$ , it is natural to conjecture that the gap between  $\Gamma_3^*$  and  $\Gamma_3$  has zero Lebesgue measure. In other words,  $\overline{\Gamma}_3^* = \Gamma_3$ , where  $\overline{\Gamma}_3^*$  is the closure of  $\Gamma_3$ . This conjecture is indeed true and will be proved at the end of the section.

More generally, we are interested in characterizing  $\overline{\Gamma}_n^*$ , the closure of  $\Gamma_n^*$ . Although the region  $\overline{\Gamma}_n^*$  is not sufficient for characterizing all information inequalities, it is actually sufficient for characterizing all unconstrained information inequalities. This can be seen as follows. Following the discussion in Section 13.3.1, an unconstrained information inequality  $f \ge 0$  involving nrandom variables always hold if and only if

$$\Gamma_n^* \subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\}. \tag{15.34}$$

Since  $\{\mathbf{h}: f(\mathbf{h}) \ge 0\}$  is closed, upon taking closure on both sides, we have

$$\overline{\Gamma}_n^* \subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\}. \tag{15.35}$$

On the other hand, if  $f \ge 0$  satisfies (15.35), then

$$\Gamma_n^* \subset \overline{\Gamma}_n^* \subset \{\mathbf{h} : f(\mathbf{h}) \ge 0\}.$$
(15.36)

Therefore, (15.34) and (15.35) are equivalent, and hence  $\overline{\Gamma}_n^*$  is sufficient for characterizing all unconstrained information inequalities.

We will prove in the next theorem an important property of the region  $\overline{\Gamma}_n^*$  for all  $n \geq 2$ . This result will be used in the proof for  $\overline{\Gamma}_3^* = \Gamma_3$ . Further,

this result will be used in Chapter 21 when we use  $\Gamma_n^*$  to characterize the achievable information rate region for multi-source networking coding. It will also be used in Chapter 16 when we establish a fundamental relation between information theory and group theory.

We first prove a simple lemma. In the following, we use  $\mathcal{N}_n$  to denote the set  $\{1, 2, \dots, n\}$ .

**Lemma 15.3.** If **h** and **h'** are in  $\Gamma_n^*$ , then  $\mathbf{h} + \mathbf{h'}$  is in  $\Gamma_n^*$ .

*Proof.* Consider **h** and **h'** in  $\Gamma_n^*$ . Let **h** represents the entropy function for random variables  $X_1, X_2, \dots, X_n$ , and let **h'** represents the entropy function for random variables  $X'_1, X'_2, \dots, X'_n$ . Let  $(X_1, X_2, \dots, X_n)$  and  $(X'_1, X'_2, \dots, X'_n)$  be independent, and define random variables  $Y_1, Y_2, \dots, Y_n$  by

$$Y_i = (X_i, X'_i) \tag{15.37}$$

for all  $i \in \mathcal{N}_n$ . Then for any subset  $\alpha$  of  $\mathcal{N}_n$ ,

$$H(Y_{\alpha}) = H(X_{\alpha}) + H(X'_{\alpha}) = h_{\alpha} + h'_{\alpha}.$$
 (15.38)

Therefore,  $\mathbf{h} + \mathbf{h}'$ , which represents the entropy function for  $Y_1, Y_2, \dots, Y_n$ , is in  $\Gamma_n^*$ . The lemma is proved.  $\Box$ 

**Corollary 15.4.** If  $\mathbf{h} \in \Gamma_n^*$ , then  $k\mathbf{h} \in \Gamma_n^*$  for any positive integer k.

*Proof.* It suffices to write

$$k\mathbf{h} = \underbrace{\mathbf{h} + \mathbf{h} + \dots + \mathbf{h}}_{k} \tag{15.39}$$

and apply Lemma 15.3.  $\Box$ 

**Theorem 15.5.**  $\overline{\Gamma}_n^*$  is a convex cone.

*Proof.* Consider the entropy function for random variables  $X_1, X_2, \dots, X_n$  all taking constant values with probability 1. Then for all subset  $\alpha$  of  $\mathcal{N}_n$ ,

$$H(X_{\alpha}) = 0.$$
 (15.40)

Therefore,  $\Gamma_n^*$  contains the origin in  $\mathcal{H}_n$ .

Let **h** and **h'** in  $\Gamma_n^*$  be the entropy functions for any two sets of random variables  $Y_1, Y_2, \dots, Y_n$  and  $Z_1, Z_2, \dots, Z_n$ , respectively. In view of Corollary 15.4, in order to prove that  $\overline{\Gamma}_n^*$  is a convex cone, we only need to show that if **h** and **h'** are in  $\Gamma_n^*$ , then  $b\mathbf{h} + \overline{b}\mathbf{h'}$  is in  $\overline{\Gamma}_n^*$  for all 0 < b < 1, where  $\overline{b} = 1 - b$ .

Let  $(\mathbf{Y}_1, \mathbf{Y}_2, \dots, \mathbf{Y}_n)$  be k independent copies of  $(Y_1, Y_2, \dots, Y_n)$  and  $(\mathbf{Z}_1, \mathbf{Z}_2, \dots, \mathbf{Z}_n)$  be k independent copies of  $(Z_1, Z_2, \dots, Z_n)$ . Let U be a ternary random variable independent of all other random variables such that

$$\Pr\{U=0\} = 1 - \delta - \mu, \ \Pr\{U=1\} = \delta, \ \Pr\{U=2\} = \mu.$$

Now construct random variables  $X_1, X_2, \dots, X_n$  by letting

$$X_i = \begin{cases} 0 & \text{if } U = 0\\ \mathbf{Y}_i & \text{if } U = 1\\ \mathbf{Z}_i & \text{if } U = 2. \end{cases}$$

Note that  $H(U) \to 0$  as  $\delta, \mu \to 0$ . Then for any nonempty subset  $\alpha$  of  $\mathcal{N}_n$ ,

$$H(X_{\alpha}) \le H(X_{\alpha}, U) \tag{15.41}$$

$$= H(U) + H(X_{\alpha}|U)$$
 (15.42)

$$= H(U) + \delta k H(Y_{\alpha}) + \mu k H(Z_{\alpha}).$$
(15.43)

On the other hand,

$$H(X_{\alpha}) \ge H(X_{\alpha}|U) = \delta k H(Y_{\alpha}) + \mu k H(Z_{\alpha}).$$
(15.44)

Combining the above, we have

$$0 \le H(X_{\alpha}) - (\delta k H(Y_{\alpha}) + \mu k H(Z_{\alpha})) \le H(U).$$
(15.45)

Now take

$$\delta = \frac{b}{k} \tag{15.46}$$

and

$$\mu = \frac{\bar{b}}{k} \tag{15.47}$$

to obtain

$$0 \le H(X_{\alpha}) - (bH(Y_{\alpha}) + \bar{b}H(Z_{\alpha})) \le H(U).$$
(15.48)

By letting k be sufficiently large, the upper bound can be made arbitrarily small. This shows that  $b\mathbf{h} + \bar{b}\mathbf{h}' \in \overline{\Gamma}_n^*$ . The theorem is proved.  $\Box$ 

In the next theorem, we prove that  $\Gamma_3^*$  and  $\Gamma_3$  are almost identical. Analogous to  $\overline{\Gamma}_n^*$ , we will use  $\overline{\Psi}_n^*$  to denote the closure of  $\Psi_n^*$ .

# **Theorem 15.6.** $\overline{\Gamma}_3^* = \Gamma_3$ .

*Proof.* We first note that  $\overline{\Gamma}_3^* = \Gamma_3$  if and only if

$$\overline{\Psi}_3^* = \Psi_3. \tag{15.49}$$

Since

$$\Gamma_3^* \subset \Gamma_3 \tag{15.50}$$

and  $\Gamma_3$  is closed, by taking closure on both sides in the above, we obtain  $\overline{\Gamma}_3^* \subset \Gamma_3$ . This implies that  $\overline{\Psi}_3^* \subset \Psi_3$ . Therefore, in order to prove the theorem, it suffices to show that  $\Psi_3 \subset \overline{\Psi}_3^*$ .

We first show that the point (0, 0, 0, a, a, a, -a) is in  $\overline{\Psi}_3^*$  for all a > 0. Let random variables  $X_1, X_2$ , and  $X_3$  be defined as in Example 3.10, i.e.,  $X_1$  and  $X_2$  are two independent binary random variables taking values in  $\{0, 1\}$  according to the uniform distribution, and

$$X_3 = X_1 + X_2 \mod 2. \tag{15.51}$$

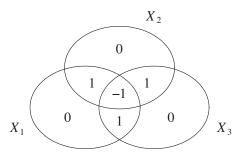
Let  $\mathbf{h} \in \Gamma_3^*$  represents the entropy function for  $X_1, X_2$ , and  $X_3$ , and let

$$\mathbf{u} = \mathbf{C}_3^{-1} \mathbf{h}.\tag{15.52}$$

As in the proof of Theorem 15.2, we let  $u_i$ ,  $1 \le i \le 7$ , be the coordinates of  $\mathcal{I}_3$  which correspond to the values of the quantities in (15.14), respectively. From Example 3.10, we have

$$u_i = \begin{cases} 0 & \text{for } i = 1, 2, 3\\ 1 & \text{for } i = 4, 5, 6\\ -1 & \text{for } i = 7. \end{cases}$$
(15.53)

Thus the point (0, 0, 0, 1, 1, 1, -1) is in  $\Psi_3^*$ , and the *I*-Measure  $\mu^*$  for  $X_1, X_2$ , and  $X_3$  is shown in Figure 15.3. Then by Corollary 15.4, (0, 0, 0, k, k, k, -k) is



**Fig. 15.3.** The *I*-Measure  $\mu^*$  for  $X_1, X_2$ , and  $X_3$  in the proof of Theorem 15.6.

in  $\Psi_3^*$  and hence in  $\overline{\Psi}_3^*$  for all positive integer k. Since  $\overline{\Gamma}_3^*$  contains the origin,  $\overline{\Psi}_3^*$  also contains the origin. By Theorem 15.5,  $\overline{\Gamma}_3^*$  is convex. This implies  $\overline{\Psi}_3^*$  is also convex. Therefore, (0, 0, 0, a, a, a, -a) is in  $\overline{\Psi}_3^*$  for all a > 0.

Consider any  $\mathbf{u} \in \Psi_3$ . Referring to (15.15), we have

$$u_i \ge 0 \tag{15.54}$$

for  $1 \leq i \leq 6$ . Thus  $u_7$  is the only component of **u** which can possibly be negative. We first consider the case when  $u_7 \geq 0$ . Then **u** is in the nonnegative orthant of  $\mathcal{I}_3$ , and by Lemma 13.1, **u** is in  $\Psi_3^*$ . Next, consider the case when  $u_7 < 0$ . Let

$$\mathbf{t} = (0, 0, 0, -u_7, -u_7, -u_7, u_7).$$
(15.55)

15.2 A Non-Shannon-Type Unconstrained Inequality 359

Then

$$\mathbf{u} = \mathbf{w} + \mathbf{t},\tag{15.56}$$

where

$$\mathbf{w} = (u_1, u_2, u_3, u_4 + u_7, u_5 + u_7, u_6 + u_7, 0).$$
(15.57)

Since  $-u_7 > 0$ , we see from the foregoing that  $\mathbf{t} \in \overline{\Psi}_3^*$ . From (15.15), we have

$$u_i + u_7 \ge 0$$
 (15.58)

for i = 4, 5, 6. Thus **w** is in the nonnegative orthant in  $\mathcal{I}_3$  and hence in  $\Psi_3^*$  by Lemma 13.1. Now for any  $\epsilon > 0$ , let  $\mathbf{t}' \in \Psi_3^*$  such that

$$\|\mathbf{t} - \mathbf{t}'\| < \epsilon, \tag{15.59}$$

where  $\|\mathbf{t} - \mathbf{t}'\|$  denotes the Euclidean distance between  $\mathbf{t}$  and  $\mathbf{t}'$ , and let

$$\mathbf{u}' = \mathbf{w} + \mathbf{t}'.\tag{15.60}$$

Since both w and t' are in  $\Psi_3^*$ , by Lemma 15.3, u' is also in  $\Psi_3^*$ , and

$$\|\mathbf{u} - \mathbf{u}'\| = \|\mathbf{t} - \mathbf{t}'\| < \epsilon.$$
(15.61)

Therefore,  $\mathbf{u} \in \overline{\Psi}_3^*$ . Hence,  $\Psi_3 \subset \overline{\Psi}_3^*$ , and the theorem is proved.  $\Box$ 

**Remark 1** Han [135] has found that  $\Gamma_3$  is the smallest cone that contains  $\Gamma_3^*$ . This result together with Theorem 15.5 implies Theorem 15.6. Theorem 15.6 was also obtained by Golić [124], and it is a consequence of the theorem in Matúš [228].

**Remark 2** We have shown that the region  $\overline{\Gamma}_n^*$  completely characterizes all unconstrained information inequalities involving n random variables. Since  $\overline{\Gamma}_3^* = \Gamma_3$ , it follows that there exists no unconstrained information inequalities involving three random variables other than the Shannon-type inequalities. However, whether there exist constrained non-Shannon-type inequalities involving three random variables is still unknown.

## 15.2 A Non-Shannon-Type Unconstrained Inequality

We have proved in Theorem 15.6 at the end of the last section that  $\overline{\Gamma}_3^* = \Gamma_3$ . It is natural to conjecture that this theorem can be generalized to  $n \ge 4$ . If this conjecture is true, then it follows that all unconstrained information inequalities involving a finite number of random variables are Shannon-type inequalities, and they can all be proved by ITIP running on a sufficiently powerful computer. However, it turns out that this is not the case even for n = 4.

We will prove in the next theorem an unconstrained information inequality involving four random variables. Then we will show that this inequality is a non-Shannon-type inequality, and that  $\overline{\Gamma}_4^* \neq \Gamma_4$ .

**Theorem 15.7.** For any four random variables  $X_1, X_2, X_3$ , and  $X_4$ ,

$$2I(X_3; X_4) \le I(X_1; X_2) + I(X_1; X_3, X_4)$$
  
+3I(X\_3; X\_4 | X\_1) + I(X\_3; X\_4 | X\_2). (15.62)

Toward proving this theorem, we introduce two auxiliary random variables  $\tilde{X}_1$  and  $\tilde{X}_2$  jointly distributed with  $X_1, X_2, X_3$ , and  $X_4$  such that  $\tilde{\mathcal{X}}_1 = \mathcal{X}_1$  and  $\tilde{\mathcal{X}}_2 = \mathcal{X}_2$ . To simplify notation, we will use  $p_{1234\tilde{1}\tilde{2}}(x_1, x_2, x_3, x_4, \tilde{x}_1, \tilde{x}_2)$  to denote  $p_{X_1X_2X_3X_4\tilde{X}_1\tilde{X}_2}(x_1, x_2, x_3, x_4, \tilde{x}_1, \tilde{x}_2)$ , etc. The joint distribution for the six random variables  $X_1, X_2, X_3, X_4, \tilde{X}_1$ , and  $\tilde{\mathcal{X}}_2$  is defined by

$$p_{1234\tilde{1}\tilde{2}}(x_1, x_2, x_3, x_4, \tilde{x}_1, \tilde{x}_2) = \begin{cases} \frac{p_{1234}(x_1, x_2, x_3, x_4) p_{1234}(\tilde{x}_1, \tilde{x}_2, x_3, x_4)}{p_{34}(x_3, x_4)} & \text{if } p_{34}(x_3, x_4) > 0\\ 0 & \text{if } p_{34}(x_3, x_4) = 0. \end{cases}$$
(15.63)

Lemma 15.8.

$$(X_1, X_2) \to (X_3, X_4) \to (\tilde{X}_1, \tilde{X}_2)$$
 (15.64)

forms a Markov chain. Moreover,  $(X_1, X_2, X_3, X_4)$  and  $(\tilde{X}_1, \tilde{X}_2, X_3, X_4)$  have the same marginal distribution.

*Proof.* The Markov chain in (15.64) is readily seen by invoking Proposition 2.5. The second part of the lemma is readily seen to be true by noting in (15.63) that  $p_{1234\tilde{1}\tilde{2}}$  is symmetrical in  $X_1$  and  $\tilde{X}_1$  and in  $X_2$  and  $\tilde{X}_2$ .  $\Box$ 

From the above lemma, we see that the pair of auxiliary random variables  $(\tilde{X}_1, \tilde{X}_2)$  corresponds to the pair of random variables  $(X_1, X_2)$  in the sense that  $(\tilde{X}_1, \tilde{X}_2, X_3, X_4)$  have the same marginal distribution as  $(X_1, X_2, X_3, X_4)$ . We need to prove two inequalities regarding these six random variables before we prove Theorem 15.7.

**Lemma 15.9.** For any four random variables  $X_1, X_2, X_3$ , and  $X_4$  and auxiliary random variables  $\tilde{X}_1$  and  $\tilde{X}_2$  as defined in (15.63),

$$I(X_3; X_4) - I(X_3; X_4 | X_1) - I(X_3; X_4 | X_2) \le I(X_1; X_2).$$
(15.65)

Proof. Consider

$$I(X_3; X_4) - I(X_3; X_4 | X_1) - I(X_3; X_4 | X_2)$$

$$\stackrel{a)}{=} [I(X_3; X_4) - I(X_3; X_4 | X_1)] - I(X_3; X_4 | \tilde{X}_2)$$

$$= I(X_1; X_3; X_4) - I(X_3; X_4 | \tilde{X}_2)$$
(15.66)
(15.67)

$$= [I(X_1; X_3; X_4; \tilde{X}_2) + I(X_1; X_3; X_4 | \tilde{X}_2)] - I(X_3; X_4 | \tilde{X}_2)$$
(15.68)

15.2 A Non-Shannon-Type Unconstrained Inequality 361

$$= I(X_1; X_3; X_4; \tilde{X}_2) - [I(X_3; X_4 | \tilde{X}_2) - I(X_1; X_3; X_4 | \tilde{X}_2)]$$
(15.69)

$$= I(X_1; X_3; X_4; X_2) - I(X_3; X_4 | X_1, X_2)$$

$$[I(X_1; X_3; X_4; X_2) - I(X_2; X_4 | X_1, X_2)] = I(X_1; X_1 | X_2, X_1 | X_1, X_2)$$
(15.70)
(15.71)

$$= [I(X_1; X_4; X_2) - I(X_1; X_4; X_2 | X_3)] - I(X_3; X_4 | X_1, X_2)$$
(15.71)  
$$= [I(X_1; \tilde{X}_2) - I(X_1; \tilde{X}_2 | X_4)] - [I(X_1; \tilde{X}_2 | X_3)]$$

$$-I(X_1; \tilde{X}_2 | X_3, X_4)] - I(X_3; X_4 | X_1, \tilde{X}_2)$$
(15.72)

$$\stackrel{b)}{=} I(X_1; \tilde{X}_2) - I(X_1; \tilde{X}_2 | X_4) - I(X_1; \tilde{X}_2 | X_3) -I(X_3; X_4 | X_1, \tilde{X}_2)$$
(15.73)

$$\leq I(X_1; \tilde{X}_2),\tag{15.74}$$

where a) follows because we see from Lemma 15.8 that  $(X_2, X_3, X_4)$  and  $(\tilde{X}_2, X_3, X_4)$  have the same marginal distribution, and b) follows because

$$I(X_1; X_2 | X_3, X_4) = 0 (15.75)$$

from the Markov chain in (15.64). The lemma is proved.  $\Box$ 

**Lemma 15.10.** For any four random variables  $X_1, X_2, X_3$ , and  $X_4$  and auxiliary random variables  $\tilde{X}_1$  and  $\tilde{X}_2$  as defined in (15.63),

$$I(X_3; X_4) - 2I(X_3; X_4 | X_1) \le I(X_1; X_1).$$
(15.76)

*Proof.* Notice that (15.76) can be obtained from (15.65) by replacing  $X_2$  by  $X_1$  and  $\tilde{X}_2$  by  $\tilde{X}_1$  in (15.65). The inequality (15.76) can be proved by replacing  $X_2$  by  $X_1$  and  $\tilde{X}_2$  by  $\tilde{X}_1$  in (15.66) through (15.74) in the proof of the last lemma. The details are omitted.  $\Box$ 

Proof of Theorem 15.7. By adding (15.65) and (15.76), we have

$$\begin{aligned} 2I(X_3; X_4) &- 3I(X_3; X_4 | X_1) - I(X_3; X_4 | X_2) \\ &\leq I(X_1; \tilde{X}_2) + I(X_1; \tilde{X}_1) & (15.77) \\ &= I(X_1; \tilde{X}_2) + [I(X_1; \tilde{X}_1 | \tilde{X}_2) + I(X_1; \tilde{X}_1; \tilde{X}_2)] & (15.78) \\ &= [I(X_1; \tilde{X}_2) + I(X_1; \tilde{X}_1 | \tilde{X}_2)] + I(X_1; \tilde{X}_1; \tilde{X}_2) & (15.79) \\ &= I(X_1; \tilde{X}_1, \tilde{X}_2) + I(X_1; \tilde{X}_1; \tilde{X}_2) & (15.80) \\ &= I(X_1; \tilde{X}_1, \tilde{X}_2) + [I(\tilde{X}_1; \tilde{X}_2) - I(\tilde{X}_1; \tilde{X}_2 | X_1)] & (15.81) \end{aligned}$$

$$\leq I(X_1; \tilde{X}_1, \tilde{X}_2) + I(\tilde{X}_1; \tilde{X}_2)$$
(15.82)

$$\stackrel{a)}{\leq} I(X_1; X_3, X_4) + I(\tilde{X}_1; \tilde{X}_2) \tag{15.83}$$

<sup>b)</sup> = 
$$I(X_1; X_3, X_4) + I(X_1; X_2),$$
 (15.84)

where a) follows from the Markov chain in (15.64), and b) follows because we see from Lemma 15.8 that  $(\tilde{X}_1, \tilde{X}_2)$  and  $(X_1, X_2)$  have the same marginal distribution. Note that the auxiliary random variables  $\tilde{X}_1$  and  $\tilde{X}_2$  disappear in (15.84) after the sequence of manipulations. The theorem is proved.  $\Box$ 

**Theorem 15.11.** The inequality (15.62) is a non-Shannon-type inequality, and  $\overline{\Gamma}_4^* \neq \Gamma_4$ .

*Proof.* Consider for any a > 0 the point  $\tilde{\mathbf{h}}(a) \in \mathcal{H}_4$ , where

$$\tilde{h}_{1}(a) = \tilde{h}_{2}(a) = \tilde{h}_{3}(a) = \tilde{h}_{4}(a) = 2a, 
\tilde{h}_{12}(a) = 4a, \quad \tilde{h}_{13}(a) = \tilde{h}_{14}(a) = 3a, 
\tilde{h}_{23}(a) = \tilde{h}_{24}(a) = \tilde{h}_{34}(a) = 3a, 
\tilde{h}_{123}(a) = \tilde{h}_{124}(a) = \tilde{h}_{134}(a) = \tilde{h}_{234}(a) = \tilde{h}_{1234}(a) = 4a.$$
(15.85)

The set-theoretic structure of  $\tilde{\mathbf{h}}(a)$  is illustrated by the information diagram in Figure 15.4. The reader should check that this information diagram correctly

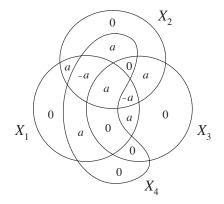


Fig. 15.4. The set-theoretic structure of  $\tilde{\mathbf{h}}(a)$ .

represents  $\tilde{\mathbf{h}}(a)$  as defined. It is also easy to check from this diagram that  $\tilde{\mathbf{h}}(a)$  satisfies all the elemental inequalities for four random variables, and therefore  $\tilde{\mathbf{h}}(a) \in \Gamma_4$ . However, upon substituting the corresponding values in (15.62) for  $\tilde{\mathbf{h}}(a)$  with the help of Figure 15.4, we have

$$2a \le 0 + a + 0 + 0 = a, \tag{15.86}$$

which is a contradiction because a > 0. In other words,  $\tilde{\mathbf{h}}(a)$  does not satisfy (15.62). Equivalently,

$$\mathbf{h}(a) \notin \{\mathbf{h} \in \mathcal{H}_4 : \mathbf{h} \text{ satisfies } (15.62)\}.$$
(15.87)

Since  $\tilde{\mathbf{h}}(a) \in \Gamma_4$ , we conclude that

$$\Gamma_4 \not\subset \{\mathbf{h} \in \mathcal{H}_4 : \mathbf{h} \text{ satisfies } (15.62)\},\tag{15.88}$$

i.e., (15.62) is not implied by the basic inequalities for four random variables. Hence, (15.62) is a non-Shannon-type inequality. Since (15.62) is satisfied by all entropy functions for four random variables, we have

$$\Gamma_4^* \subset \{ \mathbf{h} \in \mathcal{H}_4 : \mathbf{h} \text{ satisfies } (15.62) \}, \tag{15.89}$$

and upon taking closure on both sides, we have

$$\overline{\Gamma}_4^* \subset \{ \mathbf{h} \in \mathcal{H}_4 : \mathbf{h} \text{ satisfies } (15.62) \}.$$
(15.90)

Then (15.87) implies  $\tilde{\mathbf{h}}(a) \notin \overline{\Gamma}_4^*$ . Since  $\tilde{\mathbf{h}}(a) \in \Gamma_4$  and  $\tilde{\mathbf{h}}(a) \notin \overline{\Gamma}_4^*$ , we conclude that  $\overline{\Gamma}_4^* \neq \Gamma_4$ . The theorem is proved.  $\Box$ 

**Remark** We have shown in the proof of Theorem 15.11 that the inequality (15.62) cannot be proved by invoking the basic inequalities for four random variables. However, (15.62) can be proved by invoking the basic inequalities for the six random variables  $X_1, X_2, X_3, X_4, \tilde{X}_1$ , and  $\tilde{X}_2$  with the joint probability distribution  $p_{1234\tilde{12}}$  as constructed in (15.63).

The inequality (15.62) remains valid when the indices 1, 2, 3, and 4 are permuted. Since (15.62) is symmetrical in  $X_3$  and  $X_4$ , 4!/2! = 12 distinct versions of (15.62) can be obtained by permuting the indices, and all these twelve inequalities are simultaneously satisfied by the entropy function of any set of random variables  $X_1, X_2, X_3$ , and  $X_4$ . We will denote these twelve inequalities collectively by  $\langle 15.62 \rangle$ . Now define the region

$$\tilde{\Gamma}_4 = \{ \mathbf{h} \in \Gamma_4 : \mathbf{h} \text{ satisfies } \langle 15.62 \rangle \}.$$
(15.91)

Evidently,

$$\Gamma_4^* \subset \Gamma_4 \subset \Gamma_4. \tag{15.92}$$

Since both  $\tilde{\Gamma}_4$  and  $\Gamma_4$  are closed, upon taking closure, we also have

$$\overline{\Gamma}_4^* \subset \widetilde{\Gamma}_4 \subset \Gamma_4. \tag{15.93}$$

Since  $\langle 15.62 \rangle$  are non-Shannon-type inequalities as we have proved in the last theorem,  $\tilde{\Gamma}_4$  is a proper subset of  $\Gamma_4$  and hence a tighter outer bound on  $\Gamma_4^*$  and  $\overline{\Gamma}_4^*$  than  $\Gamma_4$ .

In the course of proving that (15.62) is of non-Shannon-type, it was shown in the proof of Theorem 15.11 that there exists  $\tilde{\mathbf{h}}(a) \in \Gamma_4$  as defined in (15.85) which does not satisfy (15.62). By investigating the geometrical relation between  $\tilde{\mathbf{h}}(a)$  and  $\Gamma_4$ , we prove in the next theorem that (15.62) in fact induces a class of  $2^{14} - 1$  non-Shannon-type constrained inequalities. Applications of some of these inequalities will be discussed in Section 15.4.

**Theorem 15.12.** The inequality (15.62) is a non-Shannon-type inequality conditioning on setting any nonempty subset of the following 14 Shannon's

information measures to zero:

$$I(X_{1}; X_{2}), I(X_{1}; X_{2}|X_{3}), I(X_{1}; X_{2}|X_{4}), I(X_{1}; X_{3}|X_{4}),$$

$$I(X_{1}; X_{4}|X_{3}), I(X_{2}; X_{3}|X_{4}), I(X_{2}; X_{4}|X_{3}), I(X_{3}; X_{4}|X_{1}),$$

$$I(X_{3}; X_{4}|X_{2}), I(X_{3}; X_{4}|X_{1}, X_{2}), H(X_{1}|X_{2}, X_{3}, X_{4}),$$

$$H(X_{2}|X_{1}, X_{3}, X_{4}), H(X_{3}|X_{1}, X_{2}, X_{4}), H(X_{4}|X_{1}, X_{2}, X_{3}).$$
(15.94)

*Proof.* It is easy to verify from Figure 15.4 that  $\hat{\mathbf{h}}(a)$  lies in exactly 14 hyperplanes in  $\mathcal{H}_4$  (i.e.,  $\Re^{15}$ ) defining the boundary of  $\Gamma_4$  which correspond to setting the 14 Shannon's measures in (15.94) to zero. Therefore,  $\tilde{\mathbf{h}}(a)$  for  $a \ge 0$  define an extreme direction of  $\Gamma_4$ .

Now for any linear subspace  $\Phi$  of  $\mathcal{H}_4$  containing  $\mathbf{h}(a)$ , where a > 0, we have

$$\hat{\mathbf{h}}(a) \in \Gamma_4 \cap \Phi \tag{15.95}$$

and  $\mathbf{h}(a)$  does not satisfy (15.62). Therefore,

$$(\Gamma_4 \cap \Phi) \not\subset \{\mathbf{h} \in \mathcal{H}_4 : \mathbf{h} \text{ satisfies } (15.62)\}.$$
(15.96)

This means that (15.62) is a non-Shannon-type inequality under the constraint  $\Phi$ . From the above, we see that  $\Phi$  can be taken to be the intersection of any nonempty subset of the 14 hyperplanes containing  $\tilde{\mathbf{h}}(a)$ . Thus (15.62) is a non-Shannon-type inequality conditioning on any nonempty subset of the 14 Shannon's measures in (15.94) being equal to zero. Hence, (15.62) induces a class of  $2^{14} - 1$  non-Shannon-type constrained inequalities. The theorem is proved.  $\Box$ 

**Remark** It is not true that the inequality (15.62) is of non-Shannon-type under any constraint. Suppose we impose the constraint

$$I(X_3; X_4) = 0. (15.97)$$

Then the left hand side of (15.62) becomes zero, and the inequality is trivially implied by the basic inequalities because only mutual informations with positive coefficients appear on the right hand side. Then (15.62) becomes a Shannon-type inequality under the constraint in (15.97).

# 15.3 A Non-Shannon-Type Constrained Inequality

In the last section, we proved a non-Shannon-type unconstrained inequality for four random variables which implies  $\overline{\Gamma}_4^* \neq \Gamma_4$ . This inequality induces a region  $\tilde{\Gamma}_4$  which is a tighter outer bound on  $\Gamma_4^*$  and  $\overline{\Gamma}_4^*$  then  $\Gamma_4$ . We further showed that this inequality induces a class of  $2^{14} - 1$  non-Shannon-type constrained inequalities for four random variables. In this section, we prove a non-Shannon-type constrained inequality for four random variables. Unlike the non-Shannon-type unconstrained inequality we proved in the last section, this constrained inequality is not strong enough to imply that  $\Gamma_4^* \neq \Gamma_4$ . However, the latter is not implied by the former.

**Lemma 15.13.** Let  $p(x_1, x_2, x_3, x_4)$  be any probability distribution. Then

$$\tilde{p}(x_1, x_2, x_3, x_4) = \begin{cases} \frac{p(x_1, x_3, x_4)p(x_2, x_3, x_4)}{p(x_3, x_4)} & \text{if } p(x_3, x_4) > 0\\ 0 & \text{if } p(x_3, x_4) = 0 \end{cases}$$
(15.98)

is also a probability distribution. Moreover,

$$\tilde{p}(x_1, x_3, x_4) = p(x_1, x_3, x_4) \tag{15.99}$$

and

$$\tilde{p}(x_2, x_3, x_4) = p(x_2, x_3, x_4)$$
 (15.100)

for all  $x_1, x_2, x_3$ , and  $x_4$ .

*Proof.* The proof for the first part of the lemma is straightforward (see Problem 5 in Chapter 2). The details are omitted here.

To prove the second part of the lemma, it suffices to prove (15.99) for all  $x_1, x_3$ , and  $x_4$  because  $\tilde{p}(x_1, x_2, x_3, x_4)$  is symmetrical in  $x_1$  and  $x_2$ . We first consider  $x_1, x_3$ , and  $x_4$  such that  $p(x_3, x_4) > 0$ . From (15.98), we have

$$\tilde{p}(x_1, x_3, x_4) = \sum_{x_2} \tilde{p}(x_1, x_2, x_3, x_4)$$
(15.101)

$$=\sum_{x_2} \frac{p(x_1, x_3, x_4)p(x_2, x_3, x_4)}{p(x_3, x_4)}$$
(15.102)

$$= \frac{p(x_1, x_3, x_4)}{p(x_3, x_4)} \sum_{x_2} p(x_2, x_3, x_4)$$
(15.103)

$$= \left[\frac{p(x_1, x_3, x_4)}{p(x_3, x_4)}\right] p(x_3, x_4)$$
(15.104)

$$= p(x_1, x_3, x_4). \tag{15.105}$$

For  $x_1, x_3$ , and  $x_4$  such that  $p(x_3, x_4) = 0$ , we have

$$0 \le p(x_1, x_3, x_4) \le p(x_3, x_4) = 0, \tag{15.106}$$

which implies

$$p(x_1, x_3, x_4) = 0. (15.107)$$

Therefore, from (15.98), we have

$$\tilde{p}(x_1, x_3, x_4) = \sum_{x_2} \tilde{p}(x_1, x_2, x_3, x_4)$$
(15.108)

$$=\sum_{x_2} 0$$
 (15.109)

$$= 0$$
 (15.110)

$$= p(x_1, x_3, x_4). \tag{15.111}$$

Thus we have proved (15.99) for all  $x_1, x_3$ , and  $x_4$ , and the lemma is proved.  $\Box$ 

**Theorem 15.14.** For any four random variables  $X_1, X_2, X_3$ , and  $X_4$ , if

$$I(X_1; X_2) = I(X_1; X_2 | X_3) = 0, (15.112)$$

then

$$I(X_3; X_4) \le I(X_3; X_4 | X_1) + I(X_3; X_4 | X_2).$$
 (15.113)

Proof. Consider

$$\begin{split} I(X_3; X_4) &- I(X_3; X_4 | X_1) - I(X_3; X_4 | X_2) \\ &= \sum_{\substack{x_1, x_2, x_3, x_4: \\ p(x_1, x_2, x_3, x_4) > 0}} p(x_1, x_2, x_3, x_4) \log \frac{p(x_3, x_4)p(x_1, x_3)p(x_1, x_4)p(x_2, x_3)p(x_2, x_4)}{p(x_3)p(x_4)p(x_1)p(x_2)p(x_1, x_3, x_4)p(x_2, x_3, x_4)} \\ &= E_p \log \frac{p(X_3, X_4)p(X_1, X_3)p(X_1, X_4)p(X_2, X_3)p(X_2, X_4)}{p(X_3)p(X_4)p(X_1)p(X_2)p(X_1, X_3, X_4)p(X_2, X_3, X_4)}, \end{split}$$

$$(15.114)$$

where we have used  $E_p$  to denote expectation with respect to  $p(x_1, x_2, x_3, x_4)$ . We claim that the above expectation is equal to

$$E_{\tilde{p}} \log \frac{p(X_3, X_4)p(X_1, X_3)p(X_1, X_4)p(X_2, X_3)p(X_2, X_4)}{p(X_3)p(X_4)p(X_1)p(X_2)p(X_1, X_3, X_4)p(X_2, X_3, X_4)},$$
(15.115)

where  $\tilde{p}(x_1, x_2, x_3, x_4)$  is defined in (15.98).

Toward proving that the claim is correct, we note that (15.115) is the sum of a number of expectations with respect to  $\tilde{p}$ . Let us consider one of these expectations, say

$$E_{\tilde{p}} \log p(X_1, X_3) = \sum_{\substack{x_1, x_2, x_3, x_4:\\ \tilde{p}(x_1, x_2, x_3, x_4) > 0}} \tilde{p}(x_1, x_2, x_3, x_4) \log p(x_1, x_3).$$
(15.116)

Note that in the above summation, if  $\tilde{p}(x_1, x_2, x_3, x_4) > 0$ , then from (15.98), we see that

$$p(x_1, x_3, x_4) > 0, (15.117)$$

and hence

$$p(x_1, x_3) > 0. (15.118)$$

Therefore, the summation in (15.116) is always well-defined. Further, it can be written as

$$\sum_{x_1, x_3, x_4} \log p(x_1, x_3) \sum_{x_2: \tilde{p}(x_1, x_2, x_3, x_4) > 0} \tilde{p}(x_1, x_2, x_3, x_4)$$
$$= \sum_{x_1, x_3, x_4} \tilde{p}(x_1, x_3, x_4) \log p(x_1, x_3).$$
(15.119)

Thus  $E_{\tilde{p}} \log p(X_1, X_3)$  depends on  $\tilde{p}(x_1, x_2, x_3, x_4)$  only through  $\tilde{p}(x_1, x_3, x_4)$ , which by Lemma 15.13 is equal to  $p(x_1, x_3, x_4)$ . It then follows that

$$E_{\tilde{p}} \log p(X_1, X_3)$$
  
=  $\sum_{x_1, x_3, x_4} \tilde{p}(x_1, x_3, x_4) \log p(x_1, x_3)$  (15.120)

$$= \sum_{x_1, x_3, x_4} p(x_1, x_3, x_4) \log p(x_1, x_3)$$
(15.121)

$$= E_p \log p(X_1, X_3). \tag{15.122}$$

In other words, the expectation on  $\log p(X_1, X_3)$  can be taken with respect to either  $\tilde{p}(x_1, x_2, x_3, x_4)$  or  $p(x_1, x_2, x_3, x_4)$  without affecting its value. By observing that all the marginals of p in the logarithm in (15.115) involve only subsets of either  $\{X_1, X_3, X_4\}$  or  $\{X_2, X_3, X_4\}$ , we see that similar conclusions can be drawn for all the other expectations in (15.115), and hence the claim is proved.

Thus the claim implies that

$$\begin{split} I(X_3; X_4) &- I(X_3; X_4 | X_1) - I(X_3; X_4 | X_2) \\ &= E_{\tilde{p}} \log \frac{p(X_3, X_4) p(X_1, X_3) p(X_1, X_4) p(X_2, X_3) p(X_2, X_4)}{p(X_3) p(X_4) p(X_1) p(X_2) p(X_1, X_3, X_4) p(X_2, X_3, X_4)} \\ &= \sum_{\substack{x_1, x_2, x_3, x_4:\\ \tilde{p}(x_1, x_2, x_3, x_4) > 0}} \tilde{p}(x_1, x_2, x_3, x_4) \log \frac{p(x_3, x_4) p(x_1, x_3) p(x_1, x_4) p(x_2, x_3) p(x_2, x_4)}{p(x_3) p(x_4) p(x_1) p(x_2) p(x_1, x_3, x_4) p(x_2, x_3, x_4)} \\ &= -\sum_{\substack{x_1, x_2, x_3, x_4:\\ \tilde{p}(x_1, x_2, x_3, x_4) > 0}} \tilde{p}(x_1, x_2, x_3, x_4) \log \frac{\tilde{p}(x_1, x_2, x_3, x_4)}{\hat{p}(x_1, x_2, x_3, x_4)}, \end{split}$$
(15.123)

where

$$\hat{p}(x_1, x_2, x_3, x_4) = \begin{cases}
\frac{p(x_1, x_3)p(x_1, x_4)p(x_2, x_3)p(x_2, x_4)}{p(x_1)p(x_2)p(x_3)p(x_4)} & \text{if } p(x_1), p(x_2), p(x_3), p(x_4) > 0 \\
0 & \text{otherwise.} 
\end{cases}$$
(17.104)

(15.124)

The equality in (15.123) is justified by observing that if  $x_1, x_2, x_3$ , and  $x_4$  are such that  $\tilde{p}(x_1, x_2, x_3, x_4) > 0$ , then

$$p(x_1, x_3), p(x_1, x_4), p(x_2, x_3), p(x_2, x_4), p(x_1), p(x_2), p(x_3), p(x_4)$$
(15.125)

are all strictly positive, and we see from (15.124) that  $\hat{p}(x_1, x_2, x_3, x_4) > 0$ .

To complete the proof, we only need to show that  $\hat{p}(x_1, x_2, x_3, x_4)$  is a probability distribution. Once this is proven, the conclusion of the theorem follows immediately because the summation in (15.123), which is identified as the divergence between  $\tilde{p}(x_1, x_2, x_3, x_4)$  and  $\hat{p}(x_1, x_2, x_3, x_4)$ , is always nonnegative by the divergence inequality (Theorem 2.31). Toward this end, we notice that for  $x_1, x_2$ , and  $x_3$  such that  $p(x_3) > 0$ ,

$$p(x_1, x_2, x_3) = \frac{p(x_1, x_3)p(x_2, x_3)}{p(x_3)}$$
(15.126)

by the assumption

$$I(X_1; X_2 | X_3) = 0, (15.127)$$

and for all  $x_1$  and  $x_2$ ,

$$p(x_1, x_2) = p(x_1)p(x_2)$$
(15.128)

by the assumption

$$I(X_1; X_2) = 0. (15.129)$$

Then

$$\sum_{\substack{x_1, x_2, x_3, x_4 \\ p(x_1, x_2, x_3, x_4) > 0}} \hat{p}(x_1, x_2, x_3, x_4)$$

$$= \sum_{\substack{x_1, x_2, x_3, x_4:\\ \hat{p}(x_1, x_2, x_3, x_4) > 0}} \hat{p}(x_1, x_2, x_3, x_4)$$
(15.130)

$$=\sum_{\substack{x_1,x_2,x_3,x_4:\\p(x_1),p(x_2),p(x_3),p(x_4)>0}}\frac{p(x_1,x_3)p(x_1,x_4)p(x_2,x_3)p(x_2,x_4)}{p(x_1)p(x_2)p(x_3)p(x_4)}$$
(15.131)

$$\stackrel{a)}{=} \sum_{\substack{x_1, x_2, x_3, x_4:\\p(x_1), p(x_2), p(x_3), p(x_4) > 0}} \frac{p(x_1, x_2, x_3)p(x_1, x_4)p(x_2, x_4)}{p(x_1)p(x_2)p(x_4)}$$
(15.132)

$$\stackrel{b)}{=} \sum_{\substack{x_1, x_2, x_3, x_4:\\p(x_1), p(x_2), p(x_3), p(x_4) > 0}} \frac{p(x_1, x_2, x_3)p(x_1, x_4)p(x_2, x_4)}{p(x_1, x_2)p(x_4)}$$
(15.133)

$$=\sum_{\substack{x_1,x_2,x_4:\\p(x_1),p(x_2),p(x_4)>0}}\frac{p(x_1,x_4)p(x_2,x_4)}{p(x_4)}\sum_{x_3:p(x_3)>0}p(x_3|x_1,x_2)$$
(15.134)

$$=\sum_{\substack{x_1,x_2,x_4:\\p(x_1),p(x_2),p(x_4)>0}}\frac{p(x_1,x_4)p(x_2,x_4)}{p(x_4)}$$
(15.135)

15.3 A Non-Shannon-Type Constrained Inequality 369

$$=\sum_{\substack{x_2,x_4:\\p(x_2),p(x_4)>0}} p(x_2,x_4) \sum_{x_1:p(x_1)>0} p(x_1|x_4)$$
(15.136)

$$\stackrel{c)}{=} \sum_{\substack{x_2, x_4:\\p(x_2), p(x_4) > 0}} p(x_2, x_4) \tag{15.137}$$

$$\stackrel{d)}{=} 1, \tag{15.138}$$

where a) and b) follows from (15.126) and (15.128), respectively. The equality in c) is justified as follows. For  $x_1$  such that  $p(x_1) = 0$ ,

$$p(x_1|x_4) = \frac{p(x_1)p(x_4|x_1)}{p(x_4)} = 0.$$
(15.139)

Therefore

$$\sum_{x_1:p(x_1)>0} p(x_1|x_4) = \sum_{x_1} p(x_1|x_4) = 1.$$
(15.140)

Finally, the equality in d) is justified as follows. For  $x_2$  and  $x_4$  such that  $p(x_2)$  or  $p(x_4)$  vanishes,  $p(x_2, x_4)$  must vanish because

$$0 \le p(x_2, x_4) \le p(x_2) \tag{15.141}$$

and

$$0 \le p(x_2, x_4) \le p(x_4). \tag{15.142}$$

Therefore,

$$\sum_{\substack{x_2, x_4:\\(x_2), p(x_4) > 0}} p(x_2, x_4) = \sum_{x_2, x_4} p(x_2, x_4) = 1.$$
(15.143)

The theorem is proved.  $\Box$ 

p

**Theorem 15.15.** The constrained inequality in Theorem 15.14 is a non-Shannon-type inequality.

*Proof.* The theorem can be proved by considering the point  $\mathbf{\tilde{h}}(a) \in \mathcal{H}_4$  for a > 0 as in the proof of Theorem 15.11. The details are left as an exercise.  $\Box$ 

The constrained inequality in Theorem 15.14 has the following geometrical interpretation. The constraints in (15.112) correspond to the intersection of two hyperplanes in  $\mathcal{H}_4$  which define the boundary of  $\Gamma_4$ . Then the inequality (15.62) says that a certain region on the boundary of  $\Gamma_4$  is not in  $\Gamma_4^*$ . It can further be proved by computation<sup>1</sup> that the constrained inequality in Theorem 15.14 is not implied by the twelve distinct versions of the unconstrained inequality in Theorem 15.7 (i.e.,  $\langle 15.62 \rangle$ ) together with the basic inequalities.

We have proved in the last section that the non-Shannon-type inequality (15.62) implies a class of  $2^{14} - 1$  constrained non-Shannon-type inequalities. We end this section by proving a similar result for the non-Shannon-type constrained inequality in Theorem 15.14.

<sup>&</sup>lt;sup>1</sup> Ying-On Yan, private communication.

**Theorem 15.16.** *The inequality* 

$$I(X_3; X_4) \le I(X_3; X_4 | X_1) + I(X_3; X_4 | X_2)$$
(15.144)

is a non-Shannon-type inequality conditioning on setting both  $I(X_1; X_2)$  and  $I(X_1; X_2|X_3)$  and any subset of the following 12 Shannon's information measures to zero:

$$I(X_{1}; X_{2}|X_{4}), I(X_{1}; X_{3}|X_{4}), I(X_{1}; X_{4}|X_{3}),$$

$$I(X_{2}; X_{3}|X_{4}), I(X_{2}; X_{4}|X_{3}), I(X_{3}; X_{4}|X_{1}),$$

$$I(X_{3}; X_{4}|X_{2}), I(X_{3}; X_{4}|X_{1}, X_{2}), H(X_{1}|X_{2}, X_{3}, X_{4}),$$

$$H(X_{2}|X_{1}, X_{3}, X_{4}), H(X_{3}|X_{1}, X_{2}, X_{4}), H(X_{4}|X_{1}, X_{2}, X_{3}).$$
(15.145)

Proof. The proof of this theorem is very similar to the proof of Theorem 15.12. We first note that  $I(X_1; X_2)$  and  $I(X_1; X_2|X_3)$  together with the 12 Shannon's information measures in (15.145) are exactly the 14 Shannon's information measures in (15.94). We have already shown in the proof of Theorem 15.12 that  $\tilde{\mathbf{h}}(a)$  (cf. Figure 15.4) lies in exactly 14 hyperplanes defining the boundary of  $\Gamma_4$  which correspond to setting these 14 Shannon's information measures to zero. We also have shown that  $\tilde{\mathbf{h}}(a)$  for  $a \ge 0$  define an extreme direction of  $\Gamma_4$ .

Denote by  $\Phi_0$  the intersection of the two hyperplanes in  $\mathcal{H}_4$  which correspond to setting  $I(X_1; X_2)$  and  $I(X_1; X_2|X_3)$  to zero. Since  $\tilde{\mathbf{h}}(a)$  for any a > 0 satisfies

$$I(X_1; X_2) = I(X_1; X_2 | X_3) = 0, (15.146)$$

 $\mathbf{h}(a)$  is in  $\Phi_0$ . Now for any linear subspace  $\Phi$  of  $\mathcal{H}_4$  containing  $\mathbf{h}(a)$  such that  $\Phi \subset \Phi_0$ , we have

$$\mathbf{\hat{h}}(a) \in \Gamma_4 \cap \Phi. \tag{15.147}$$

Upon substituting the corresponding values in (15.113) for  $\mathbf{h}(a)$  with the help of Figure 15.4, we have

$$a \le 0 + 0 = 0, \tag{15.148}$$

which is a contradiction because a > 0. Therefore,  $\mathbf{h}(a)$  does not satisfy (15.113). Therefore,

$$(\Gamma_4 \cap \Phi) \not\subset \{\mathbf{h} \in \mathcal{H}_4 : \mathbf{h} \text{ satisfies } (15.113)\}.$$
(15.149)

This means that (15.113) is a non-Shannon-type inequality under the constraint  $\Phi$ . From the above, we see that  $\Phi$  can be taken to be the intersection of  $\Phi_0$  and any subset of the 12 hyperplanes which correspond to setting the 12 Shannon's information measures in (15.145) to zero. Hence, (15.113) is a non-Shannon-type inequality conditioning on  $I(X_1; X_2)$ ,  $I(X_1; X_2|X_3)$ , and any subset of the 12 Shannon's information measures in (15.145) being equal to zero. In other words, the constrained inequality in Theorem 15.14 in fact induces a class of  $2^{12}$  constrained non-Shannon-type inequalities. The theorem is proved.  $\Box$ 

## **15.4 Applications**

As we have mentioned in Chapter 13, information inequalities are the laws of information theory. In this section, we give several applications of the non-Shannon-type inequalities we have proved in this chapter in probability theory and information theory. An application of the unconstrained inequality proved in Section 15.2 in group theory will be discussed in Chapter 16.

*Example 15.17.* For the constrained inequality in Theorem 15.14, if we further impose the constraints

$$I(X_3; X_4 | X_1) = I(X_3; X_4 | X_2) = 0, (15.150)$$

then the right hand side of (15.113) becomes zero. This implies

$$I(X_3; X_4) = 0 \tag{15.151}$$

because  $I(X_3; X_4)$  is nonnegative. This means that

$$\left. \begin{array}{c} X_{1} \perp X_{2} \\ X_{1} \perp X_{2} | X_{3} \\ X_{3} \perp X_{4} | X_{1} \\ X_{3} \perp X_{4} | X_{2} \end{array} \right\} \Rightarrow X_{3} \perp X_{4}.$$
(15.152)

We leave it as an exercise for the reader to show that this implication cannot be deduced from the basic inequalities.

Example 15.18. If we impose the constraints

$$I(X_1; X_2) = I(X_1; X_3, X_4) = I(X_3; X_4 | X_1) = I(X_3; X_4 | X_2) = 0, \quad (15.153)$$

then the right hand side of (15.62) becomes zero, which implies

$$I(X_3; X_4) = 0. (15.154)$$

This means that

$$\left. \begin{array}{c} X_1 \perp X_2 \\ X_1 \perp (X_3, X_4) \\ X_3 \perp X_4 | X_1 \\ X_3 \perp X_4 | X_2 \end{array} \right\} \Rightarrow X_3 \perp X_4. \quad (15.155)$$

Note that (15.152) and (15.155) differ only in the second constraint. Again, we leave it as an exercise for the reader to show that this implication cannot be deduced from the basic inequalities.

Example 15.19. Consider a fault-tolerant data storage system consisting of random variables  $X_1, X_2, X_3, X_4$  such that any three random variables can recover the remaining one, i.e.,

$$H(X_i|X_j, j \neq i) = 0, \quad 1 \le i, j \le 4.$$
(15.156)

We are interested in the set of all entropy functions subject to these constraints, denoted by  $\Upsilon$ , which characterizes the amount of joint information which can possibly be stored in such a data storage system. Let

$$\boldsymbol{\Phi} = \{ \mathbf{h} \in \mathcal{H}_4 : \mathbf{h} \text{ satisfies } (15.156) \}.$$
(15.157)

Then the set  $\Upsilon$  is equal to the intersection between  $\Gamma_4^*$  and  $\Phi$ , i.e.,  $\Gamma_4^* \cap \Phi$ .

Since each constraint in (15.156) is one of the 14 constraints specified in Theorem 15.12, we see that (15.62) is a non-Shannon-type inequality under the constraints in (15.156). Then  $\tilde{\Gamma}_4 \cap \Phi$  (cf. (15.91)) is a tighter outer bound on  $\Upsilon$  than  $\Gamma_4 \cap \Phi$ .

Example 15.20. Consider four random variables  $X_1, X_2, X_3$ , and  $X_4$  such that  $X_3 \to (X_1, X_2) \to X_4$  forms a Markov chain. This Markov condition is equivalent to

$$I(X_3; X_4 | X_1, X_2) = 0. (15.158)$$

It can be proved by invoking the basic inequalities (using ITIP) that

$$I(X_3; X_4) \le I(X_3; X_4 | X_1) + I(X_3; X_4 | X_2) + 0.5I(X_1; X_2) + cI(X_1; X_3, X_4) + (1 - c)I(X_2; X_3, X_4),$$
(15.159)

where  $0.25 \le c \le 0.75$ , and this is the best possible.

Now observe that the Markov condition (15.158) is one of the 14 constraints specified in Theorem 15.12. Therefore, (15.62) is a non-Shannon-type inequality under this Markov condition. By replacing  $X_1$  and  $X_2$  by each other in (15.62), we obtain

$$2I(X_3; X_4) \le I(X_1; X_2) + I(X_2; X_3, X_4) +3I(X_3; X_4 | X_2) + I(X_3; X_4 | X_1).$$
(15.160)

Upon adding (15.62) and (15.160) and dividing by 4, we obtain

$$I(X_3; X_4) \le I(X_3; X_4 | X_1) + I(X_3; X_4 | X_2) + 0.5I(X_1; X_2) + 0.25I(X_1; X_3, X_4) + 0.25I(X_2; X_3, X_4).$$
(15.161)

Comparing the last two terms in (15.159) and the last two terms in (15.161), we see that (15.161) is a sharper upper bound than (15.159).

The Markov chain  $X_3 \to (X_1, X_2) \to X_4$  arises in many communication situations. As an example, consider a person listening to an audio source. Then the situation can be modeled by this Markov chain with  $X_3$  being the sound wave generated at the source,  $X_1$  and  $X_2$  being the sound waves received at the two ear drums, and  $X_4$  being the nerve impulses which eventually arrive at the brain. The inequality (15.161) gives an upper bound on  $I(X_3; X_4)$  which is tighter than what can be implied by the basic inequalities.

There is some resemblance between the constrained inequality (15.161) and the data processing theorem, but there does not seem to be any direct relation between them.

## Problems

- 1. Verify by ITIP that the unconstrained information inequality in Theorem 15.7 is of non-Shannon-type.
- 2. Verify by ITIP and prove analytically that the constrained information inequality in Theorem 15.14 is of non-Shannon-type.
- 3. Use ITIP to verify the unconstrained information inequality in Theorem 15.7. Hint: Create two auxiliary random variables as in the proof of Theorem 15.7 and impose appropriate constraints on the random variables.
- 4. Verify by ITIP that the implications in Examples 15.17 and 15.18 cannot be deduced from the basic inequalities.
- 5. Can you show that the sets of constraints in Examples 15.17 and 15.18 are in fact different?
- 6. Let  $\mathcal{N}_n = \{1, 2, \dots, n\}$ . Consider an information inequality involving random variables  $X_1, X_2, \dots, X_n$ , which can be written as

$$\sum_{\alpha \in 2^{\mathcal{N}_n} \setminus \{\emptyset\}} c_{\alpha} H(X_{\alpha}) \ge 0$$

For  $i \in \mathcal{N}_n$ , let

$$r_i = \sum_{\alpha \in 2^{\mathcal{N}_n} \setminus \{\emptyset\}} c_\alpha \mathbf{1}_\alpha(i),$$

where  $1_{\alpha}(i)$  is equal to 1 if  $i \in \alpha$  and is equal to 0 otherwise.

- a) Show that  $r_i$  is the coefficient associated with  $H(X_i|X_{\mathcal{N}_n-\{i\}})$  in the information inequality.
- b) Show that if the information inequality always holds, then  $r_i \ge 0$  for all  $i \in \mathcal{N}_n$ .

(Chan [58].)

7. Let  $X_i$ ,  $i = 1, 2, \dots, n, Z$ , and T be discrete random variables.

a) Prove that

$$nI(Z;T) - \sum_{j=1}^{n} I(Z;T|X_j) - nI(Z;T|X_i)$$
  
$$\leq I(X_i;Z,T) + \sum_{j=1}^{n} H(X_j) - H(X_1,X_2,\cdots,X_n).$$

Hint: When n = 2, this inequality reduces to the unconstrained non-Shannon-type inequality in Theorem 15.7.

b) Prove that

$$nI(Z;T) - 2\sum_{j=1}^{n} I(Z;T|X_j)$$
  
$$\leq \frac{1}{n}\sum_{i=1}^{n} I(X_i;Z,T) + \sum_{j=1}^{n} H(X_j) - H(X_1,X_2,\cdots,X_n).$$

(Zhang and Yeung [373].)

8. Let  $p(x_1, x_2, x_3, x_4)$  be the joint distribution for random variables  $X_1, X_2, X_3$ , and  $X_4$  such that  $I(X_1; X_2|X_3) = I(X_2; X_4|X_3) = 0$ , and let  $\tilde{p}$  be defined in (15.98).

a) Show that

$$\check{p}(x_1, x_2, x_3, x_4) = \begin{cases} c \cdot \frac{p(x_1, x_2, x_3)p(x_1, x_4)p(x_2, x_4)}{p(x_1, x_2)p(x_4)} & \text{if } p(x_1, x_2), p(x_4) > 0\\ 0 & \text{otherwise} \end{cases}$$

defines a probability distribution for an appropriate  $c \geq 1$ .

- b) Prove that  $\tilde{p}(x_1, x_2, x_3) = p(x_1, x_2, x_3)$  for all  $x_1, x_2$ , and  $x_3$ .
- c) By considering  $D(\tilde{p} \| \check{p}) \ge 0$ , prove that

$$H(X_{13}) + H(X_{14}) + H(X_{23}) + H(X_{24}) + H(X_{34})$$
  

$$\geq H(X_3) + H(X_4) + H(X_{12}) + H(X_{134}) + H(X_{234}),$$

where  $H(X_{134})$  denotes  $H(X_1, X_3, X_4)$ , etc.

d) Prove that under the constraints in (15.112), the inequality in (15.113) is equivalent to the inequality in c).

The inequality in c) is referred to as the *Ingleton inequality* for entropy in the literature. For the origin of the Ingleton inequality, see Problem 9 in Chapter 16. (Matúš [230].)

## **Historical Notes**

In 1986, Pippenger [265] asked whether there exist constraints on the entropy function other than the polymatroidal axioms, which are equivalent to the basic inequalities. He called the constraints on the entropy function the *laws* of information theory. The problem had been open since then until Zhang and Yeung discovered for four random variables first a constrained non-Shannontype inequality [372] and then an unconstrained non-Shannon-type inequality [373]. The inequality reported in [373] has been further generalized by Makarychev *et al.* [217] and Zhang [370]. The existence of these inequalities implies that there are laws in information theory beyond those laid down by Shannon [291].

The non-Shannon-type inequalities that have been discovered induce outer bounds on the region  $\Gamma_4^*$  which are tighter than  $\Gamma_4$ . Matúš and Studený [235] showed that an entropy function in  $\Gamma_4$  is entropic if it satisfies the Ingleton inequality (see Problem 9 in Chapter 16). This gives an inner bound on  $\overline{\Gamma}_4^*$ . A more explicit proof of this inner bound can be found in [373], where the bound was shown not to be tight. Matúš [233] has obtained asymptotically tight inner bounds on  $\overline{\Gamma}_n^*$  by constructing entropy functions from matroids. Dougherty *et al.* [87] discovered a host of unconstrained non-Shannon-type inequalities by means of a computer search based on ITIP and the Markov chain construction in [373] (see Problem 3). Recently, Matúš [234] proved an infinite class of unconstrained non-Shannon-type inequalities, implying that  $\overline{\Gamma}_n^*$  is not polyhedral.

Chan [58] proved a characterization for an inequality for differential entropy in terms of its discrete version. Lněnička [212] proved that the tightness of the continuous version of the inequality reported in [373] can be achieved by a multivariate Gaussian distribution.

In the 1990's, Matúš and Studený [228][235][229] studied the structure of conditional independence (which subsumes the implication problem) of random variables. Matúš [230] finally settled the problem for four random variables by means of a constrained non-Shannon-type inequality which is a variation of the inequality reported in [372].

The von Neumann entropy is an extension of classical entropy (as discussed in this book) to the field of quantum mechanics. The strong subadditivity of the von Neumann entropy proved by Lieb and Ruskai [208] plays the same role as the basic inequalities for classical entropy. Pippenger [266] proved that for a three-party system, there exists no inequality for the von Neumann entropy beyond strong subadditivity. Subsequently, Linden and Winter [210] discovered for a four-party system a constrained inequality for the von Neumann entropy which is independent of strong subadditivity. We refer the reader to the book by Nielsen and Chuang [247] for quantum information theory.

Along a related direction, Hammer *et al.* [131] have shown that all linear inequalities which always hold for Kolmogorov complexity also always hold for entropy, and vice versa.

# **Entropy and Groups**

The group is the first major mathematical structure in abstract algebra, while entropy is the most basic measure of information. Group theory and information theory are two seemingly unrelated subjects which turn out to be intimately related to each other. This chapter explains this intriguing relation between these two fundamental subjects. Those readers who have no knowledge in group theory may skip this introduction and go directly to the next section.

Let  $X_1$  and  $X_2$  be any two random variables. Then

$$H(X_1) + H(X_2) \ge H(X_1, X_2), \tag{16.1}$$

which is equivalent to the basic inequality

$$I(X_1; X_2) \ge 0. \tag{16.2}$$

Let G be any finite group and  $G_1$  and  $G_2$  be subgroups of G. We will show in Section 16.4 that

$$|G||G_1 \cap G_2| \ge |G_1||G_2|, \tag{16.3}$$

where |G| denotes the order of G and  $G_1 \cap G_2$  denotes the intersection of  $G_1$  and  $G_2$  ( $G_1 \cap G_2$  is also a subgroup of G, see Proposition 16.13). By rearranging the terms, the above inequality can be written as

$$\log \frac{|G|}{|G_1|} + \log \frac{|G|}{|G_2|} \ge \log \frac{|G|}{|G_1 \cap G_2|}.$$
(16.4)

By comparing (16.1) and (16.4), one can easily identify the one-to-one correspondence between these two inequalities, namely that  $X_i$  corresponds to  $G_i$ , i = 1, 2, and  $(X_1, X_2)$  corresponds to  $G_1 \cap G_2$ . While (16.1) is true for any pair of random variables  $X_1$  and  $X_2$ , (16.4) is true for any finite group G and subgroups  $G_1$  and  $G_2$ .

Recall from Chapter 13 that the region  $\Gamma_n^*$  characterizes all information inequalities (involving *n* random variables). In particular, we have shown in

#### 378 16 Entropy and Groups

Section 15.1 that the region  $\overline{\Gamma}_n^*$  is sufficient for characterizing all unconstrained information inequalities, i.e., by knowing  $\overline{\Gamma}_n^*$ , one can determine whether any unconstrained information inequality always holds. The main purpose of this chapter is to obtain a characterization of  $\overline{\Gamma}_n^*$  in terms of finite groups. An important consequence of this result is a one-to-one correspondence between unconstrained information inequalities and group inequalities. Specifically, for every unconstrained information inequality, there is a corresponding group inequality, and vice versa. A special case of this correspondence has been given in (16.1) and (16.4).

By means of this result, unconstrained information inequalities can be proved by techniques in group theory, and a certain form of inequalities in group theory can be proved by techniques in information theory. In particular, the unconstrained non-Shannon-type inequality in Theorem 15.7 corresponds to the group inequality

$$|G_1 \cap G_3|^3 |G_1 \cap G_4|^3 |G_3 \cap G_4|^3 |G_2 \cap G_3| |G_2 \cap G_4|$$
  
$$\leq |G_1| |G_1 \cap G_2| |G_3|^2 |G_4|^2 |G_1 \cap G_3 \cap G_4|^4 |G_2 \cap G_3 \cap G_4|, \qquad (16.5)$$

where  $G_i$  are subgroups of a finite group G, i = 1, 2, 3, 4. The meaning of this inequality and its implications in group theory are yet to be understood.

## **16.1 Group Preliminaries**

In this section, we present the definition and some basic properties of a group which are essential for subsequent discussions.

**Definition 16.1.** A group is a set of objects G together with a binary operation on the elements of G, denoted by " $\circ$ " unless otherwise specified, which satisfy the following four axioms:

- 1. Closure For every a, b in  $G, a \circ b$  is also in G.
- 2. Associativity For every a, b, c in  $G, a \circ (b \circ c) = (a \circ b) \circ c$ .
- 3. Existence of Identity There exists an element e in G such that  $a \circ e = e \circ a$ = a for every a in G.
- 4. Existence of Inverse For every a in G, there exists an element b in G such that  $a \circ b = b \circ a = e$ .

**Proposition 16.2.** For any group G, the identity element is unique.

*Proof.* Let e and e' be both identity elements in a group G. Since e is an identity element,

$$e' \circ e = e, \tag{16.6}$$

and since e' is also an identity element,

$$e' \circ e = e'. \tag{16.7}$$

It follows by equating the right hand sides of (16.6) and (16.7) that e = e', which implies the uniqueness of the identity element of a group.  $\Box$ 

(10 11)

**Proposition 16.3.** For every element a in a group G, its inverse is unique. Proof. Let b and b' be inverses of an element a, so that

$$a \circ b = b \circ a = e \tag{16.8}$$

and

$$a \circ b' = b' \circ a = e. \tag{16.9}$$

Then

$$b = b \circ e \tag{16.10}$$

$$= b \circ (a \circ b') \tag{16.11}$$

$$= (b \circ a) \circ b' \tag{16.12}$$
$$= a \circ b' \tag{16.12}$$

$$= e \circ b \tag{10.13}$$

$$= 0$$
, (10.14)

where (16.11) and (16.13) follow from (16.9) and (16.8), respectively, and (16.12) is by associativity. Therefore, the inverse of a is unique.  $\Box$ 

11

Thus the inverse of a group element a is a function of a, and it will be denoted by  $a^{-1}$ .

**Definition 16.4.** The number of elements of a group G is called the order of G, denoted by |G|. If  $|G| < \infty$ , G is called a finite group, otherwise it is called an infinite group.

There is an unlimited supply of examples of groups. Some familiar examples are: the integers under addition, the rationals excluding zero under multiplication, and the set of real-valued  $2 \times 2$  matrices under addition, where addition and multiplication refer to the usual addition and multiplication for real numbers and matrices. In each of these examples, the operation (addition or multiplication) plays the role of the binary operation " $\circ$ " in Definition 16.2.

All the above are examples of infinite groups. In this chapter, however, we are concerned with finite groups. In the following, we discuss two examples of finite groups in details.

Example 16.5 (Modulo 2 Addition). The trivial group consists of only the identity element. The simplest nontrivial group is the group of modulo 2 addition. The order of this group is 2, and the elements are  $\{0, 1\}$ . The binary operation, denoted by "+", is defined by following table:

+	0	1
0	0	1
1	1	0

### 380 16 Entropy and Groups

The four axioms of a group simply say that certain constraints must hold in the above table. We now check that all these axioms are satisfied. First, the closure axiom requires that all the entries in the table are elements in the group, which is easily seen to be the case. Second, it is required that associativity holds. To this end, it can be checked in the above table that for all a, b, and c,

$$a + (b + c) = (a + b) + c.$$
 (16.15)

For example,

$$0 + (1+1) = 0 + 0 = 0, (16.16)$$

while

$$(0+1) + 1 = 1 + 1 = 0, (16.17)$$

which is the same as 0 + (1 + 1). Third, the element 0 is readily identified as the unique identity. Fourth, it is readily seen that an inverse exists for each element in the group. For example, the inverse of 1 is 1, because

$$1 + 1 = 0. \tag{16.18}$$

Thus the above table defines a group of order 2. It happens in this example that the inverse of each element is the element itself, which is not true for a group in general.

We remark that in the context of a group, the elements in the group should be regarded strictly as *symbols* only. In particular, one should not associate group elements with *magnitudes* as we do for real numbers. For instance, in the above example, one should not think of 0 as being less than 1. The element 0, however, is a special symbol which plays the role of the identity of the group.

We also notice that for the group in the above example, a + b is equal to b + a for all group elements a and b. A group with this property is called a *commutative* group or an *Abelian* group<sup>1</sup>.

*Example 16.6 (Symmetric Group).* Consider a permutation of the components of a vector

$$\mathbf{x} = (x_1, x_2, \cdots, x_r) \tag{16.19}$$

given by

$$\sigma[\mathbf{x}] = (x_{\sigma(1)}, x_{\sigma(2)}, \cdots, x_{\sigma(r)}), \qquad (16.20)$$

where

$$\sigma: \{1, 2, \cdots, r\} \to \{1, 2, \cdots, r\}$$
(16.21)

is a one-to-one mapping. The one-to-one mapping  $\sigma$  is called a permutation on  $\{1, 2, \dots, r\}$ , which is represented by

$$\sigma = (\sigma(1), \sigma(2), \cdots, \sigma(r)). \tag{16.22}$$

<sup>&</sup>lt;sup>1</sup> The Abelian group is named after the Norwegian mathematician Niels Henrik Abel (1802-1829).

For two permutations  $\sigma_1$  and  $\sigma_2$ , define  $\sigma_1 \circ \sigma_2$  as the composite function of  $\sigma_1$  and  $\sigma_2$ . For example, for r = 4, suppose

$$\sigma_1 = (2, 1, 4, 3) \tag{16.23}$$

and

$$\sigma_2 = (1, 4, 2, 3). \tag{16.24}$$

Then  $\sigma_1 \circ \sigma_2$  is given by

$$\begin{aligned}
\sigma_1 \circ \sigma_2(1) &= \sigma_1(\sigma_2(1)) = \sigma_1(1) = 2 \\
\sigma_1 \circ \sigma_2(2) &= \sigma_1(\sigma_2(2)) = \sigma_1(4) = 3 \\
\sigma_1 \circ \sigma_2(3) &= \sigma_1(\sigma_2(3)) = \sigma_1(2) = 1 \\
\sigma_1 \circ \sigma_2(4) &= \sigma_1(\sigma_2(4)) = \sigma_1(3) = 4,
\end{aligned}$$
(16.25)

or

$$\sigma_1 \circ \sigma_2 = (2, 3, 1, 4). \tag{16.26}$$

The reader can easily check that

$$\sigma_2 \circ \sigma_1 = (4, 1, 2, 3), \tag{16.27}$$

which is different from  $\sigma_1 \circ \sigma_2$ . Therefore, the operation " $\circ$ " is not commutative.

We now show that the set of all permutations on  $\{1, 2, \dots, r\}$  and the operation " $\circ$ " form a group, called the symmetric group on  $\{1, 2, \dots, r\}$ . First, for two permutations  $\sigma_1$  and  $\sigma_2$ , since both  $\sigma_1$  and  $\sigma_2$  are one-to-one mappings, so is  $\sigma_1 \circ \sigma_2$ . Therefore, the closure axiom is satisfied. Second, for permutations  $\sigma_1$ ,  $\sigma_2$ , and  $\sigma_3$ ,

$$\sigma_1 \circ (\sigma_2 \circ \sigma_3)(i) = \sigma_1(\sigma_2 \circ \sigma_3(i)) \tag{16.28}$$

$$=\sigma_1(\sigma_2(\sigma_3(i)))$$
 (16.29)

$$=\sigma_1 \circ \sigma_2(\sigma_3(i)) \tag{16.30}$$

$$= (\sigma_1 \circ \sigma_2) \circ \sigma_3(i) \tag{16.31}$$

for  $1 \leq i \leq r$ . Therefore, associativity is satisfied. Third, it is clear that the identity map is the identity element. Fourth, for a permutation  $\sigma$ , it is clear that its inverse is  $\sigma^{-1}$ , the inverse mapping of  $\sigma$  which is defined because  $\sigma$  is one-to-one. Therefore, the set of all permutations on  $\{1, 2, \dots, r\}$  and the operation " $\circ$ " form a group. The order of this group is evidently equal to (r!).

**Definition 16.7.** Let G be a group with operation " $\circ$ ", and S be a subset of G. If S is a group with respect to the operation " $\circ$ ", then S is called a subgroup of G.

**Definition 16.8.** Let S be a subgroup of a group G and a be an element of G. The left coset of S with respect to a is the set  $a \circ S = \{a \circ s : s \in S\}$ . Similarly, the right coset of S with respect to a is the set  $S \circ a = \{s \circ a : s \in S\}$ . 382 16 Entropy and Groups

In the sequel, only the left coset will be used. However, any result which applies to the left coset also applies to the right coset, and vice versa. For simplicity,  $a \circ S$  will be denoted by aS.

**Proposition 16.9.** For  $a_1$  and  $a_2$  in G,  $a_1S$  and  $a_2S$  are either identical or disjoint. Further,  $a_1S$  and  $a_2S$  are identical if and only if  $a_1$  and  $a_2$  belong to the same left coset of S.

*Proof.* Suppose  $a_1S$  and  $a_2S$  are not disjoint. Then there exists an element b in  $a_1S \cap a_2S$  such that

$$b = a_1 \circ s_1 = a_2 \circ s_2, \tag{16.32}$$

for some  $s_i$  in S, i = 1, 2. Then

$$a_1 = (a_2 \circ s_2) \circ s_1^{-1} = a_2 \circ (s_2 \circ s_1^{-1}) = a_2 \circ t,$$
(16.33)

where  $t = s_2 \circ s_1^{-1}$  is in S. We now show that  $a_1 S \subset a_2 S$ . For an element  $a_1 \circ s$  in  $a_1 S$ , where  $s \in S$ ,

$$a_1 \circ s = (a_2 \circ t) \circ s = a_2 \circ (t \circ s) = a_2 \circ u,$$
 (16.34)

where  $u = t \circ s$  is in S. This implies that  $a_1 \circ s$  is in  $a_2S$ . Thus,  $a_1S \subset a_2S$ . By symmetry,  $a_2S \subset a_1S$ . Therefore,  $a_1S = a_2S$ . Hence, if  $a_1S$  and  $a_2S$  are not disjoint, then they are identical. Equivalently,  $a_1S$  and  $a_2S$  are either identical or disjoint. This proves the first part of the proposition.

We now prove the second part of the proposition. Since S is a group, it contains e, the identity element. Then for any group element  $a, a = a \circ e$  is in aS because e is in S. If  $a_1S$  and  $a_2S$  are identical, then  $a_1 \in a_1S$  and  $a_2 \in a_2S = a_1S$ . Therefore,  $a_1$  and  $a_2$  belong to the same left coset of S.

To prove the converse, assume  $a_1$  and  $a_2$  belong to the same left coset of S. From the first part of the proposition, we see that a group element belongs to one and only one left coset of S. Since  $a_1$  is in  $a_1S$  and  $a_2$  is in  $a_2S$ , and  $a_1$  and  $a_2$  belong to the same left coset of S, we see that  $a_1S$  and  $a_2S$  are identical. The proposition is proved.  $\Box$ 

**Proposition 16.10.** Let S be a subgroup of a group G and a be an element of G. Then |aS| = |S|, i.e., the numbers of elements in all the left cosets of S are the same, and they are equal to the order of S.

*Proof.* Consider two elements  $a \circ s_1$  and  $a \circ s_2$  in  $a \circ S$ , where  $s_1$  and  $s_2$  are in S such that

$$a \circ s_1 = a \circ s_2. \tag{16.35}$$

Then

$$a^{-1} \circ (a \circ s_1) = a^{-1} \circ (a \circ s_2) \tag{16.36}$$

$$(a^{-1} \circ a) \circ s_1 = (a^{-1} \circ a) \circ s_2 \tag{16.37}$$

$$e \circ s_1 = e \circ s_2 \tag{16.38}$$

$$s_1 = s_2.$$
 (16.39)

Thus each element in S corresponds to a unique element in aS. Therefore, |aS| = |S| for all  $a \in G$ .  $\Box$ 

We are just one step away from obtaining the celebrated Lagrange's theorem stated below.

**Theorem 16.11 (Lagrange's Theorem).** If S is a subgroup of G, then |S| divides |G|.

*Proof.* Since  $a \in aS$  for every  $a \in G$ , every element of G belongs to a left coset of S. Then from Proposition 16.9, we see that the distinct left cosets of S partition G. Therefore |G|, the total number of elements in G, is equal to the number of distinct cosets of S multiplied by the number of elements in each left coset, which is equal to |S| by Proposition 16.10. This implies that |S| divides |G|, proving the theorem.  $\Box$ 

The following corollary is immediate from the proof of Lagrange's Theorem.

**Corollary 16.12.** Let S be a subgroup of a group G. The number of distinct left cosets of S is equal to  $\frac{|G|}{|S|}$ .

### 16.2 Group-Characterizable Entropy Functions

Recall from Chapter 13 that the region  $\Gamma_n^*$  consists of all the entropy functions in the entropy space  $\mathcal{H}_n$  for n random variables. As a first step toward establishing the relation between entropy and groups, we discuss in this section entropy functions in  $\Gamma_n^*$  which can be described by a finite group Gand subgroups  $G_1, G_2, \dots, G_n$ . Such entropy functions are said to be groupcharacterizable. The significance of this class of entropy functions will become clear in the next section.

In the sequel, we will make use of the intersections of subgroups extensively. We first prove that the intersection of two subgroups is also a subgroup.

**Proposition 16.13.** Let  $G_1$  and  $G_2$  be subgroups of a group G. Then  $G_1 \cap G_2$  is also a subgroup of G.

*Proof.* It suffices to show that  $G_1 \cap G_2$  together with the operation "◦" satisfy all the axioms of a group. First, consider two elements a and b of G in  $G_1 \cap G_2$ . Since both a and b are in  $G_1$ ,  $(a \circ b)$  is in  $G_1$ . Likewise,  $(a \circ b)$  is in  $G_2$ . Therefore,  $a \circ b$  is in  $G_1 \cap G_2$ . Thus the closure axiom holds for  $G_1 \cap G_2$ . Second, associativity for  $G_1 \cap G_2$  inherits from G. Third,  $G_1$  and  $G_2$  both contain the identity element because they are groups. Therefore, the identity element is in  $G_1 \cap G_2$ . Fourth, for an element  $a \in G_i$ , since  $G_i$  is a group,  $a^{-1}$ is in  $G_i$ , i = 1, 2. Thus for an element  $a \in G_1 \cap G_2$ ,  $a^{-1}$  is also in  $G_1 \cap G_2$ . Therefore,  $G_1 \cap G_2$  is a group and hence a subgroup of G. □ 384 16 Entropy and Groups

**Corollary 16.14.** Let  $G_1, G_2, \dots, G_n$  be subgroups of a group G. Then  $\bigcap_{i=1}^n G_i$  is also a subgroup of G.

In the rest of the chapter, we let  $\mathcal{N}_n = \{1, 2, \dots, n\}$  and denote  $\bigcap_{i \in \alpha} G_i$  by  $G_{\alpha}$ , where  $\alpha$  is a nonempty subset of  $\mathcal{N}_n$ .

**Lemma 16.15.** Let  $G_i$  be subgroups of a group G and  $a_i$  be elements of G,  $i \in \alpha$ . Then

$$\left|\bigcap_{i\in\alpha}a_{i}G_{i}\right| = \begin{cases} |G_{\alpha}| & \text{if } \bigcap_{i\in\alpha}a_{i}G_{i}\neq\emptyset\\ 0 & \text{otherwise.} \end{cases}$$
(16.40)

*Proof.* For the special case when  $\alpha$  is a singleton, i.e.,  $\alpha = \{i\}$  for some  $i \in \mathcal{N}_n$ , (16.40) reduces to

$$|a_i G_i| = |G_i|, (16.41)$$

which has already been proved in Proposition 16.10.

Let  $\alpha$  be any nonempty subset of  $\mathcal{N}_n$ . If  $\bigcap_{i \in \alpha} a_i G_i = \emptyset$ , then (16.40) is obviously true. If  $\bigcap_{i \in \alpha} a_i G_i \neq \emptyset$ , then there exists  $x \in \bigcap_{i \in \alpha} a_i G_i$  such that for all  $i \in \alpha$ ,

$$x = a_i \circ s_i, \tag{16.42}$$

where  $s_i \in G_i$ . For any  $i \in \alpha$  and for any  $y \in G_\alpha$ , consider

$$x \circ y = (a_i \circ s_i) \circ y = a_i \circ (s_i \circ y). \tag{16.43}$$

Since both  $s_i$  and y are in  $G_i$ ,  $s_i \circ y$  is in  $G_i$ . Thus  $x \circ y$  is in  $a_i G_i$  for all  $i \in \alpha$ , or  $x \circ y$  is in  $\bigcap_{i \in \alpha} a_i G_i$ . Moreover, for  $y, y' \in G_\alpha$ , if  $x \circ y = x \circ y'$ , then y = y'. Therefore, each element in  $G_\alpha$  corresponds to a unique element in  $\bigcap_{i \in \alpha} a_i G_i$ . Hence,

$$|\cap_{i\in\alpha}a_iG_i| = |G_\alpha|,\tag{16.44}$$

proving the lemma.  $\Box$ 

The relation between a finite group G and subgroups  $G_1$  and  $G_2$  is illustrated by the *membership table* in Figure 16.1. In this table, an element of G is represented by a dot. The first column represents the subgroup  $G_1$ , with the dots in the first column being the elements in  $G_1$ . The other columns represent the left cosets of  $G_1$ . By Proposition 16.10, all the columns have the same number of dots. Similarly, the first row represents the subgroup  $G_2$  and the other rows represent the left cosets of  $G_2$ . Again, all the rows have the same number of dots.

The upper left entry in the table represents the subgroup  $G_1 \cap G_2$ . There are  $|G_1 \cap G_2|$  dots in this entry, with one of them representing the identity element. Any other entry represents the intersection between a left coset of  $G_1$  and a left coset of  $G_2$ , and by Lemma 16.15, the number of dots in each of these entries is either equal to  $|G_1 \cap G_2|$  or zero.

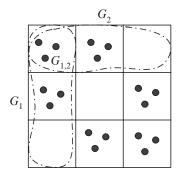


Fig. 16.1. The membership table for a finite group G and subgroups  $G_1$  and  $G_2$ .

Since all the column have the same numbers of dots and all the rows have the same number of dots, we say that the table in Figure 16.1 exhibits a *quasi-uniform* structure. We have already seen a similar structure in Figure 6.1 for the two-dimensional strong joint typicality array, which we reproduce in Figure 16.2. In this array, when n is large, all the columns have approximately the same number of dots and all the rows have approximately the same number of dots. For this reason, we say that the two-dimensional strong typicality array exhibits an *asymptotic* quasi-uniform structure. In a strong typicality array, however, each entry can contain only one dot, while in a membership table, each entry can contain multiple dots.

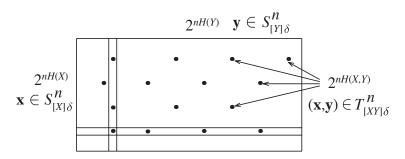


Fig. 16.2. A two-dimensional strong typicality array.

One can make a similar comparison between a strong joint typicality array for any  $n \ge 2$  random variables and the membership table for a finite group with n subgroups. The details are omitted here.

**Theorem 16.16.** Let  $G_i, i \in \mathcal{N}_n$  be subgroups of a group G. Then  $\mathbf{h} \in \mathcal{H}_n$  defined by

$$h_{\alpha} = \log \frac{|G|}{|G_{\alpha}|} \tag{16.45}$$

### 386 16 Entropy and Groups

for all nonempty subset  $\alpha$  of  $\mathcal{N}_n$  is entropic, i.e.,  $\mathbf{h} \in \Gamma_n^*$ .

*Proof.* It suffices to show that there exists a collection of random variables  $X_1, X_2, \dots, X_n$  such that

$$H(X_{\alpha}) = \log \frac{|G|}{|G_{\alpha}|} \tag{16.46}$$

for all nonempty subset  $\alpha$  of  $\mathcal{N}_n$ . We first introduce a uniform random variable  $\Lambda$  defined on the sample space G with probability mass function

$$\Pr\{\Lambda = a\} = \frac{1}{|G|}$$
(16.47)

for all  $a \in G$ . For any  $i \in \mathcal{N}_n$ , let random variable  $X_i$  be a function of  $\Lambda$  such that  $X_i = aG_i$  if  $\Lambda = a$ .

Let  $\alpha$  be a nonempty subset of  $\mathcal{N}_n$ . Since  $X_i = a_i G_i$  for all  $i \in \alpha$  if and only if  $\Lambda$  is equal to some  $b \in \bigcap_{i \in \alpha} a_i G_i$ ,

$$\Pr\{X_i = a_i G_i : i \in \alpha\} = \frac{\left|\bigcap_{i \in \alpha} a_i G_i\right|}{|G|}$$
(16.48)

$$= \begin{cases} \frac{|G_{\alpha}|}{|G|} & \text{if } \bigcap_{i \in \alpha} a_i G_i \neq \emptyset \\ 0 & \text{otherwise} \end{cases}$$
(16.49)

by Lemma 16.15. In other words,  $(X_i, i \in \alpha)$  distributes uniformly on its support whose cardinality is  $\frac{|G|}{|G_{\alpha}|}$ . Then (16.46) follows and the theorem is proved.  $\Box$ 

**Definition 16.17.** Let G be a finite group and  $G_1, G_2, \dots, G_n$  be subgroups of G. Let **h** be a vector in  $\mathcal{H}_n$ . If  $h_\alpha = \log \frac{|G|}{|G_\alpha|}$  for all nonempty subsets  $\alpha$  of  $\mathcal{N}_n$ , then  $(G, G_1, \dots, G_n)$  is a group characterization of **h**.

Theorem 16.16 asserts that certain entropy functions in  $\Gamma_n^*$  have a group characterization. These are called group-characterizable entropy functions, which will be used in the next section to obtain a group characterization of the region  $\overline{\Gamma}_n^*$ . We end this section by giving a few examples of such entropy functions.

*Example 16.18.* Fix any subset  $\beta$  of  $\mathcal{N}_3 = \{1, 2, 3\}$  and define a vector  $\mathbf{h} \in \mathcal{H}_3$  by

$$h_{\alpha} = \begin{cases} \log 2 & \text{if } \alpha \cap \beta \neq \emptyset \\ 0 & \text{otherwise.} \end{cases}$$
(16.50)

We now show that **h** has a group characterization. Let  $G = \{0, 1\}$  be the group of modulo 2 addition in Example 16.5, and for i = 1, 2, 3, let

16.2 Group-Characterizable Entropy Functions 387

$$G_i = \begin{cases} \{0\} \text{ if } i \in \beta \\ G \text{ otherwise.} \end{cases}$$
(16.51)

Then for a nonempty subset  $\alpha$  of  $\mathcal{N}_3$ , if  $\alpha \cap \beta \neq \emptyset$ , there exists an *i* in  $\alpha$  such that *i* is also in  $\beta$ , and hence by definition  $G_i = \{0\}$ . Thus,

$$G_{\alpha} = \bigcap_{i \in \alpha} G_i = \{0\}.$$
(16.52)

Therefore,

$$\log \frac{|G|}{|G_{\alpha}|} = \log \frac{|G|}{|\{0\}|} = \log \frac{2}{1} = \log 2.$$
(16.53)

If  $\alpha \cap \beta = \emptyset$ , then  $G_i = G$  for all  $i \in \alpha$ , and

$$G_{\alpha} = \bigcap_{i \in \alpha} G_i = G. \tag{16.54}$$

Therefore,

$$\log \frac{|G|}{|G_{\alpha}|} = \log \frac{|G|}{|G|} = \log 1 = 0.$$
(16.55)

Then we see from (16.50), (16.53), and (16.55) that

$$h_{\alpha} = \log \frac{|G|}{|G_{\alpha}|} \tag{16.56}$$

for all nonempty subset  $\alpha$  of  $\mathcal{N}_3$ . Hence,  $(G, G_1, G_2, G_3)$  is a group characterization of **h**.

*Example 16.19.* This is a generalization of the last example. Fix any nonempty subset  $\beta$  of  $\mathcal{N}_n$  and define a vector  $\mathbf{h} \in \mathcal{H}_n$  by

$$h_{\alpha} = \begin{cases} \log 2 & \text{if } \alpha \cap \beta \neq \emptyset \\ 0 & \text{otherwise.} \end{cases}$$
(16.57)

Then  $(G, G_1, G_2, \dots, G_n)$  is a group characterization of **h**, where G is the group of modulo 2 addition, and

$$G_i = \begin{cases} \{0\} \text{ if } i \in \beta \\ G \text{ otherwise.} \end{cases}$$
(16.58)

By letting  $\beta = \emptyset$ ,  $\mathbf{h} = 0$ . Thus we see that  $(G, G_1, G_2, \dots, G_n)$  is a group characterization of the origin of  $\mathcal{H}_n$ , with  $G = G_1 = G_2 = \dots = G_n$ .

*Example 16.20.* Define a vector  $\mathbf{h} \in \mathcal{H}_3$  as follows:

$$h_{\alpha} = \min(|\alpha|, 2). \tag{16.59}$$

Let F be the group of modulo 2 addition,  $G = F \times F$ , and

$$G_1 = \{(0,0), (1,0)\}$$
(16.60)  
$$G_1 = \{(0,0), (0,1)\}$$
(16.61)

$$G_2 = \{(0,0), (0,1)\}$$
(16.61)

$$G_3 = \{(0,0), (1,1)\}.$$
 (16.62)

Then  $(G, G_1, G_2, G_3)$  is a group characterization of **h**.

388 16 Entropy and Groups

# 16.3 A Group Characterization of $\overline{\Gamma}_n^*$

We have introduced in the last section the class of entropy functions in  $\Gamma_n^*$ which have a group characterization. However, an entropy function  $\mathbf{h} \in \Gamma_n^*$ may not have a group characterization due to the following observation. Suppose  $\mathbf{h} \in \Gamma_n^*$ . Then there exists a collection of random variables  $X_1, X_2, \cdots, X_n$  such that

$$h_{\alpha} = H(X_{\alpha}) \tag{16.63}$$

for all nonempty subset  $\alpha$  of  $\mathcal{N}_n$ . If  $(G, G_1, \dots, G_n)$  is a group characterization of **h**, then

$$H(X_{\alpha}) = \log \frac{|G|}{|G_{\alpha}|} \tag{16.64}$$

for all nonempty subset of  $\mathcal{N}_n$ . Since both |G| and  $|G_\alpha|$  are integers,  $H(X_\alpha)$  must be the logarithm of a rational number. However, the joint entropy of a set of random variables in general is not necessarily the logarithm of a rational number (see Corollary 2.44). Therefore, it is possible to construct an entropy function  $\mathbf{h} \in \Gamma_n^*$  which has no group characterization.

Although  $\mathbf{h} \in \Gamma_n^*$  does not imply  $\mathbf{h}$  has a group characterization, it turns out that the set of all  $\mathbf{h} \in \Gamma_n^*$  which have a group characterization is almost good enough to characterize the region  $\Gamma_n^*$ , as we will see next.

**Definition 16.21.** Define the following region in  $\mathcal{H}_n$ :

$$\Upsilon_n = \{ \mathbf{h} \in \mathcal{H}_n : \mathbf{h} \text{ has a group characterization} \}.$$
(16.65)

By Theorem 16.16, if  $\mathbf{h} \in \mathcal{H}_n$  has a group characterization, then  $\mathbf{h} \in \Gamma_n^*$ . Therefore,  $\Upsilon_n \subset \Gamma_n^*$ . We will prove as a corollary of the next theorem that  $\overline{con}(\Upsilon_n)$ , the convex closure of  $\Upsilon_n$ , is in fact equal to  $\overline{\Gamma}_n^*$ , the closure of  $\Gamma_n^*$ .

**Theorem 16.22.** For any  $\mathbf{h} \in \Gamma_n^*$ , there exists a sequence  $\{\mathbf{f}^{(r)}\}\$  in  $\Upsilon_n$  such that  $\lim_{r\to\infty} \frac{1}{r} \mathbf{f}^{(r)} = \mathbf{h}$ .

We need the following lemma to prove this theorem. The proof of this lemma resembles the proof of Theorem 6.10. Nevertheless, we give a sketch of the proof for the sake of completeness.

**Lemma 16.23.** Let X be a random variable such that  $|\mathcal{X}| < \infty$  and the distribution  $\{p(x)\}$  is rational, i.e., p(x) is a rational number for all  $x \in \mathcal{X}$ . Without loss of generality, assume p(x) is a rational number with denominator q for all  $x \in \mathcal{X}$ . Then for  $r = q, 2q, 3q, \cdots$ ,

$$\lim_{r \to \infty} \frac{1}{r} \log \frac{r!}{\prod_{x} (rp(x))!} = H(X).$$
(16.66)

*Proof.* Applying Lemma 6.11, we can obtain

$$\frac{1}{r} \ln \frac{r!}{\prod_{x} (rp(x))!} \le -\sum_{x} p(x) \ln p(x) + \frac{r+1}{r} \ln(r+1) - \ln r$$
(16.67)

$$= H_e(X) + \frac{1}{r}\ln r + \left(1 + \frac{1}{r}\right)\ln\left(1 + \frac{1}{r}\right).$$
(16.68)

This upper bound tends to  $H_e(X)$  as  $r \to \infty$ . On the other hand, we can obtain

$$\frac{1}{r}\ln\frac{r!}{\prod_{x}(rp(x))!}$$

$$\geq -\sum_{x}\left(p(x) + \frac{1}{r}\right)\ln\left(p(x) + \frac{1}{r}\right) - \frac{\ln r}{r}.$$
(16.69)

This lower bound also tends to  $H_e(X)$  as  $r \to \infty$ . Then the proof is completed by changing the base of the logarithm if necessary.  $\Box$ 

Proof of Theorem 16.22. For any  $\mathbf{h} \in \Gamma_n^*$ , there exists a collection of random variables  $X_1, X_2, \dots, X_n$  such that

$$h_{\alpha} = H(X_{\alpha}) \tag{16.70}$$

for all nonempty subset  $\alpha$  of  $\mathcal{N}_n$ . We first consider the special case that  $|\mathcal{X}_i| < \infty$  for all  $i \in \mathcal{N}_n$  and the joint distribution of  $X_1, X_2, \dots, X_n$  is rational. We want to show that there exists a sequence  $\{\mathbf{f}^{(r)}\}$  in  $\mathcal{T}_n$  such that  $\lim_{r\to\infty} \frac{1}{r} \mathbf{f}_{-}^{(r)} = \mathbf{h}$ .

Denote  $\prod_{i \in \alpha} \mathcal{X}_i$  by  $\mathcal{X}_\alpha$ . For any nonempty subset  $\alpha$  of  $\mathcal{N}_n$ , let  $Q_\alpha$  be the marginal distribution of  $X_\alpha$ . Assume without loss of generality that for any nonempty subset  $\alpha$  of  $\mathcal{N}_n$  and for all  $a \in \mathcal{X}_\alpha$ ,  $Q_\alpha(a)$  is a rational number with denominator q.

For each  $r = q, 2q, 3q, \cdots$ , fix a sequence

$$\mathbf{x}_{\mathcal{N}_n} = (x_{\mathcal{N}_n,1}, x_{\mathcal{N}_n,2}, \cdots x_{\mathcal{N}_n,r})$$

where for all  $j = 1, 2, \dots, r$ ,  $x_{\mathcal{N}_n, j} = (x_{i,j} : i \in \mathcal{N}_n) \in \mathcal{X}_{\mathcal{N}_n}$ , such that  $N(a|\mathbf{x}_{\mathcal{N}_n})$ , the number of occurrences of a in sequence  $\mathbf{x}_{\mathcal{N}_n}$ , is equal to  $rQ_{\mathcal{N}_n}(a)$  for all  $a \in \mathcal{X}_{\mathcal{N}_n}$ . The existence of such a sequence is guaranteed by that all the values of the joint distribution of  $X_{\mathcal{N}_n}$  are rational numbers with denominator q. Also, we denote the sequence of r elements of  $\mathcal{X}_{\alpha}$ ,  $(x_{\alpha,1}, x_{\alpha,2}, \dots, x_{\alpha,r})$ , where  $x_{\alpha,j} = (x_{i,j} : i \in \alpha)$ , by  $\mathbf{x}_{\alpha}$ . Let  $a \in \mathcal{X}_{\alpha}$ . It is easy to check that  $N(a|\mathbf{x}_{\alpha})$ , the number of occurrences of a in the sequence  $\mathbf{x}_{\alpha}$ , is equal to  $rQ_{\alpha}(a)$  for all  $a \in \mathcal{X}_{\alpha}$ .

#### 390 16 Entropy and Groups

Let G be the group of all permutations on  $\{1, 2, \dots, r\}$ , i.e., the symmetric group on  $\{1, 2, \dots, r\}$  (cf. Example 16.6). The group G depends on r, but for simplicity, we do not state this dependency explicitly. For any  $i \in \mathcal{N}_n$ , define

$$G_i = \{ \sigma \in G : \sigma[\mathbf{x}_i] = \mathbf{x}_i \},\$$

where

$$\sigma[\mathbf{x}_i] = (x_{i,\sigma(1)}, x_{i,\sigma(2)}, \cdots, x_{i,\sigma(r)}).$$
(16.71)

It is easy to check that  $G_i$  is a subgroup of G.

Let  $\alpha$  be a nonempty subset of  $\mathcal{N}_n$ . Then

$$G_{\alpha} = \bigcap_{i \in \alpha} G_i \tag{16.72}$$

$$= \bigcap_{i \in \alpha} \{ \sigma \in G : \sigma[\mathbf{x}_i] = \mathbf{x}_i \}$$
(16.73)

$$= \{ \sigma \in G : \sigma[\mathbf{x}_i] = \mathbf{x}_i \text{ for all } i \in \alpha \}$$
(16.74)

$$= \{ \sigma \in G : \sigma[\mathbf{x}_{\alpha}] = \mathbf{x}_{\alpha} \}, \tag{16.75}$$

where

$$\sigma[\mathbf{x}_{\alpha}] = (x_{\alpha,\sigma(1)}, x_{\alpha,\sigma(2)}, \cdots, x_{\alpha,\sigma(r)}).$$
(16.76)

For any  $a \in \mathcal{X}_{\alpha}$ , define the set

$$L_{\mathbf{x}_{\alpha}}(a) = \{ j \in \{1, 2, \cdots, r\} : x_{\alpha, j} = a \}.$$
 (16.77)

 $L_{\mathbf{x}_{\alpha}}(a)$  contains the "locations" of a in  $\mathbf{x}_{\alpha}$ . Then  $\sigma[\mathbf{x}_{\alpha}] = \mathbf{x}_{\alpha}$  if and only if for all  $a \in \mathcal{X}_{\alpha}, j \in L_{\mathbf{x}_{\alpha}}(a)$  implies  $\sigma(j) \in L_{\mathbf{x}_{\alpha}}(a)$ . Since

$$|L_{\mathbf{x}_{\alpha}}(a)| = N(a|\mathbf{x}_{\alpha}) = rQ_{\alpha}(a), \qquad (16.78)$$

$$|G_{\alpha}| = \prod_{a \in \mathcal{X}_{\alpha}} (rQ_{\alpha}(a))!$$
(16.79)

and therefore

$$\frac{|G|}{|G_{\alpha}|} = \frac{r!}{\prod_{a \in \mathcal{X}_{\alpha}} (rQ_{\alpha}(a))!}.$$
(16.80)

By Lemma 16.23,

$$\lim_{r \to \infty} \frac{1}{r} \log \frac{|G|}{|G_{\alpha}|} = H(X_{\alpha}) = h_{\alpha}.$$
(16.81)

Recall that G and hence all its subgroups depend on r. Define  $\mathbf{f}^{(r)}$  by

$$f_{\alpha}^{(r)} = \log \frac{|G|}{|G_{\alpha}|} \tag{16.82}$$

for all nonempty subset  $\alpha$  of  $\mathcal{N}_n$ . Then  $\mathbf{f}^{(r)} \in \Upsilon_n$  and

16.4 Information Inequalities and Group Inequalities 391

$$\lim_{r \to \infty} \frac{1}{r} \mathbf{f}^{(r)} = \mathbf{h}.$$
 (16.83)

We have already proved the theorem for the special case that  $\mathbf{h}$  is the entropy function of a collection of random variables  $X_1, X_2, \dots, X_n$  with finite alphabets and a rational joint distribution. To complete the proof, we only have to note that for any  $\mathbf{h} \in \Gamma_n^*$ , it is always possible to construct a sequence  $\{\mathbf{h}^{(k)}\}$  in  $\Gamma_n^*$  such that  $\lim_{k\to\infty} \mathbf{h}^{(k)} = \mathbf{h}$ , where  $\mathbf{h}^{(k)}$  is the entropy function of a collection of random variables  $X_1^{(k)}, X_2^{(k)}, \dots, X_n^{(k)}$  with finite alphabets and a rational joint distribution. This can be proved by techniques similar to those used in Appendix 2.10 together with the continuity of the entropy function with respect to the variational distance for a fixed finite support (Section 2.3). The details are omitted here.  $\Box$ 

Corollary 16.24.  $\overline{con}(\Upsilon_n) = \overline{\Gamma}_n^*$ .

Proof. First of all,  $\Upsilon_n \subset \Gamma_n^*$ . By taking convex closure, we have  $\overline{con}(\Upsilon_n) \subset \overline{con}(\Gamma_n^*)$ . By Theorem 15.5,  $\overline{\Gamma}_n^*$  is convex. Therefore,  $\overline{con}(\Gamma_n^*) = \overline{\Gamma}_n^*$ , and we have  $\overline{con}(\Upsilon_n) \subset \overline{\Gamma}_n^*$ . On the other hand, we have shown in Example 16.19 that the origin of  $\mathcal{H}_n$  has a group characterization and therefore is in  $\Upsilon_n$ . It then follows from Theorem 16.22 that  $\overline{\Gamma}_n^* \subset \overline{con}(\Upsilon_n)$ . Hence, we conclude that  $\overline{\Gamma}_n^* = \overline{con}(\Upsilon_n)$ , completing the proof.  $\Box$ 

# 16.4 Information Inequalities and Group Inequalities

We have proved in Section 15.1 that an unconstrained information inequality

$$\mathbf{b}^{\top}\mathbf{h} \ge 0 \tag{16.84}$$

always holds if and only if

$$\overline{\Gamma}_n^* \subset \{ \mathbf{h} \in \mathcal{H}_n : \mathbf{b}^\top \mathbf{h} \ge 0 \}.$$
(16.85)

In other words, all unconstrained information inequalities are fully characterized by  $\overline{\Gamma}_n^*$ . We also have proved at the end of the last section that  $\overline{con}(\Upsilon_n) = \overline{\Gamma}_n^*$ . Since  $\Upsilon_n \subset \Gamma_n^* \subset \overline{\Gamma}_n^*$ , if (16.85) holds, then

$$\Upsilon_n \subset \{ \mathbf{h} \in \mathcal{H}_n : \mathbf{b}^\top \mathbf{h} \ge 0 \}.$$
(16.86)

On the other hand, if (16.86) holds, since  $\{\mathbf{h} \in \mathcal{H}_n : \mathbf{b}^\top \mathbf{h} \ge 0\}$  is closed and convex, by taking convex closure in (16.86), we obtain

$$\overline{\Gamma}_n^* = \overline{con}(\Upsilon_n) \subset \{ \mathbf{h} \in \mathcal{H}_n : \mathbf{b}^\top \mathbf{h} \ge 0 \}.$$
(16.87)

Therefore, (16.85) and (16.86) are equivalent.

Now (16.86) is equivalent to

392 16 Entropy and Groups

$$\mathbf{b}^{\top}\mathbf{h} \ge 0 \text{ for all } \mathbf{h} \in \Upsilon_n. \tag{16.88}$$

Since  $\mathbf{h} \in \Upsilon_n$  if and only if

$$h_{\alpha} = \log \frac{|G|}{|G_{\alpha}|} \tag{16.89}$$

for all nonempty subset  $\alpha$  of  $\mathcal{N}_n$  for some finite group G and subgroups  $G_1$ ,  $G_2, \dots, G_n$ , we see that the inequality (16.84) holds for all random variables  $X_1, X_2, \dots, X_n$  if and only if the inequality obtained from (16.84) by replacing  $h_\alpha$  by  $\log \frac{|G|}{|G_\alpha|}$  for all nonempty subset  $\alpha$  of  $\mathcal{N}_n$  holds for all finite group G and subgroups  $G_1, G_2, \dots, G_n$ . In other words, for every unconstrained information inequality, there is a corresponding group inequality, and vice versa. Therefore, inequalities in information theory can be proved by methods in group theory, and inequalities in group theory can be proved by methods in information theory.

In the rest of the section, we explore this one-to-one correspondence between information theory and group theory. We first give a group-theoretic proof of the basic inequalities in information theory. At the end of the section, we will give an information-theoretic proof for the group inequality in (16.5).

**Definition 16.25.** Let  $G_1$  and  $G_2$  be subgroups of a finite group G. Define

$$G_1 \circ G_2 = \{ a \circ b : a \in G_1 \text{ and } b \in G_2 \}.$$
(16.90)

 $G_1 \circ G_2$  is in general not a subgroup of G. However, it can be shown that  $G_1 \circ G_2$  is a subgroup of G if G is Abelian (see Problem 1).

**Proposition 16.26.** Let  $G_1$  and  $G_2$  be subgroups of a finite group G. Then

$$|G_1 \circ G_2| = \frac{|G_1||G_2|}{|G_1 \cap G_2|}.$$
(16.91)

*Proof.* Fix  $(a_1, a_2) \in G_1 \times G_2$ , Then  $a_1 \circ a_2$  is in  $G_1 \circ G_2$ . Consider any  $(b_1, b_2) \in G_1 \times G_2$  such that

$$b_1 \circ b_2 = a_1 \circ a_2. \tag{16.92}$$

We will determine the number of  $(b_1, b_2)$  in  $G_1 \times G_2$  which satisfies this relation. From (16.92), we have

$$b_1^{-1} \circ (b_1 \circ b_2) = b_1^{-1} \circ (a_1 \circ a_2) \tag{16.93}$$

$$(b_1^{-1} \circ b_1) \circ b_2 = b_1^{-1} \circ a_1 \circ a_2 \tag{16.94}$$

$$b_2 = b_1^{-1} \circ a_1 \circ a_2. \tag{16.95}$$

Then

$$b_2 \circ a_2^{-1} = b_1^{-1} \circ a_1 \circ (a_2 \circ a_2^{-1}) = b_1^{-1} \circ a_1.$$
(16.96)

Let k be this common element in G, i.e.,

$$k = b_2 \circ a_2^{-1} = b_1^{-1} \circ a_1. \tag{16.97}$$

Since  $b_1^{-1} \circ a_1 \in G_1$  and  $b_2 \circ a_2^{-1} \in G_2$ , k is in  $G_1 \cap G_2$ . In other words, for given  $(a_1, a_2) \in G_1 \times G_2$ , if  $(b_1, b_2) \in G_1 \times G_2$  satisfies (16.92), then  $(b_1, b_2)$ satisfies (16.97) for some  $k \in G_1 \cap G_2$ . On the other hand, if  $(b_1, b_2) \in G_1 \times G_2$ satisfies (16.97) for some  $k \in G_1 \cap G_2$ , then (16.96) is satisfied, which implies (16.92). Therefore, for given  $(a_1, a_2) \in G_1 \times G_2$ ,  $(b_1, b_2) \in G_1 \times G_2$  satisfies (16.92) if and only if  $(b_1, b_2)$  satisfies (16.97) for some  $k \in G_1 \cap G_2$ .

Now from (16.97), we obtain

$$b_1(k) = (k \circ a_1^{-1})^{-1} \tag{16.98}$$

and

$$b_2(k) = k \circ a_2, \tag{16.99}$$

where we have written  $b_1$  and  $b_2$  as  $b_1(k)$  and  $b_2(k)$  to emphasize their dependence on k. Now consider  $k, k' \in G_1 \cap G_2$  such that

$$(b_1(k), b_2(k)) = (b_1(k'), b_2(k')).$$
 (16.100)

Since  $b_1(k) = b_1(k')$ , from (16.98), we have

$$(k \circ a_1^{-1})^{-1} = (k' \circ a_1^{-1})^{-1}, (16.101)$$

which implies

$$k = k'.$$
 (16.102)

Therefore, each  $k \in G_1 \cap G_2$  corresponds to a unique pair  $(b_1, b_2) \in G_1 \times G_2$ which satisfies (16.92). Therefore, we see that the number of distinct elements in  $G_1 \circ G_2$  is given by

$$|G_1 \circ G_2| = \frac{|G_1 \times G_2|}{|G_1 \cap G_2|} = \frac{|G_1||G_2|}{|G_1 \cap G_2|},$$
(16.103)

completing the proof.  $\Box$ 

**Theorem 16.27.** Let  $G_1, G_2$ , and  $G_3$  be subgroups of a finite group G. Then

$$|G_3||G_{123}| \ge |G_{13}||G_{23}|. \tag{16.104}$$

Proof. First of all,

$$G_{13} \cap G_{23} = (G_1 \cap G_3) \cap (G_2 \cap G_3) = G_1 \cap G_2 \cap G_3 = G_{123}.$$
(16.105)

By Proposition 16.26, we have

394 16 Entropy and Groups

$$|G_{13} \circ G_{23}| = \frac{|G_{13}||G_{23}|}{|G_{123}|}.$$
(16.106)

It is readily seen that  $G_{13} \circ G_{23}$  is a subset of  $G_3$ , Therefore,

$$|G_{13} \circ G_{23}| = \frac{|G_{13}||G_{23}|}{|G_{123}|} \le |G_3|.$$
(16.107)

The theorem is proved.  $\hfill\square$ 

**Corollary 16.28.** For random variables  $X_1, X_2$ , and  $X_3$ ,

$$I(X_1; X_2 | X_3) \ge 0. \tag{16.108}$$

*Proof.* Let  $G_1, G_2$ , and  $G_3$  be subgroups of a finite group G. Then

$$|G_3||G_{123}| \ge |G_{13}||G_{23}| \tag{16.109}$$

by Theorem 16.27, or

$$\frac{|G|^2}{|G_{13}||G_{23}|} \ge \frac{|G|^2}{|G_3||G_{123}|}.$$
(16.110)

This is equivalent to

$$\log \frac{|G|}{|G_{13}|} + \log \frac{|G|}{|G_{23}|} \ge \log \frac{|G|}{|G_3|} + \log \frac{|G|}{|G_{123}|}.$$
 (16.111)

This group inequality corresponds to the information inequality

$$H(X_1, X_3) + H(X_2, X_3) \ge H(X_3) + H(X_1, X_2, X_3),$$
 (16.112)

which is equivalent to

$$I(X_1; X_2 | X_3) \ge 0. \tag{16.113}$$

The above corollary shows that all the basic inequalities in information theory has a group-theoretic proof. Of course, Theorem 16.27 is also implied by the basic inequalities. As a remark, the inequality in (16.3) is seen to be a special case of Theorem 16.27 by letting  $G_3 = G$ .

We are now ready to prove the group inequality in (16.5). The non-Shannon-type inequality we have proved in Theorem 15.7 can be expressed in canonical form as

$$H(X_1) + H(X_1, X_2) + 2H(X_3) + 2H(X_4) +4H(X_1, X_3, X_4) + H(X_2, X_3, X_4) \leq 3H(X_1, X_3) + 3H(X_1, X_4) + 3H(X_3, X_4) +H(X_2, X_3) + H(X_2, X_4),$$
(16.114)

which corresponds to the group inequality

$$\log \frac{|G|}{|G_1|} + \log \frac{|G|}{|G_{12}|} + 2\log \frac{|G|}{|G_3|} + 2\log \frac{|G|}{|G_4|} + 4\log \frac{|G|}{|G_{134}|} + \log \frac{|G|}{|G_{234}|} \\ \leq 3\log \frac{|G|}{|G_{13}|} + 3\log \frac{|G|}{|G_{14}|} + 3\log \frac{|G|}{|G_{34}|} + \log \frac{|G|}{|G_{23}|} \\ + \log \frac{|G|}{|G_{24}|}.$$
(16.115)

Upon rearranging the terms, we obtain

$$|G_1 \cap G_3|^3 |G_1 \cap G_4|^3 |G_3 \cap G_4|^3 |G_2 \cap G_3| |G_2 \cap G_4|$$
  
$$\leq |G_1| |G_1 \cap G_2| |G_3|^2 |G_4|^2 |G_1 \cap G_3 \cap G_4|^4 |G_2 \cap G_3 \cap G_4|, \qquad (16.116)$$

which is the group inequality in (16.5). The meaning of this inequality and its implications in group theory are yet to be understood.

# Problems

- 1. Let  $G_1$  and  $G_2$  be subgroups of a finite group G. Show that  $G_1 \circ G_2$  is a subgroup if G is Abelian.
- 2. Let  $\mathbf{g}_1$  and  $\mathbf{g}_2$  be group characterizable entropy functions.
  - a) Prove that  $m_1 \mathbf{g}_1 + m_2 \mathbf{g}_2$  is group characterizable, where  $m_1$  and  $m_2$  are any positive integers.
  - b) For any positive real numbers  $a_1$  and  $a_2$ , construct a sequence of group characterizable entropy functions  $\mathbf{f}^{(k)}$  for  $k = 1, 2, \cdots$ , such that

$$\lim_{k \to \infty} \frac{\mathbf{f}^{(k)}}{||\mathbf{f}^{(k)}||} = \frac{\mathbf{h}}{||\mathbf{h}||},$$

where  $\mathbf{h} = a_1 \mathbf{g}_1 + a_2 \mathbf{g}_2$ .

3. Let  $(G, G_1, G_2, \dots, G_n)$  be a group characterization of  $\mathbf{g} \in \Gamma_n^*$ , where **g** is the entropy function for random variables  $X_1, X_2, \dots, X_n$ . Fix any nonempty subset  $\alpha$  of  $\mathcal{N}_n$ , and define **h** by

$$h_{\beta} = g_{\alpha \cup \beta} - g_{\alpha}$$

for all nonempty subsets  $\beta$  of  $\mathcal{N}_n$ . It can easily be checked that  $h_\beta = H(X_\beta | X_\alpha)$ . Show that  $(K, K_1, K_2, \dots, K_n)$  is a group characterization of **h**, where  $K = G_\alpha$  and  $K_i = G_i \cap G_\alpha$ .

4. Let  $(G, G_1, G_2, \dots, G_n)$  be a group characterization of  $\mathbf{g} \in \Gamma_n^*$ , where  $\mathbf{g}$  is the entropy function for random variables  $X_1, X_2, \dots, X_n$ . Show that if  $X_i$  is a function of  $(X_j : j \in \alpha)$ , then  $G_\alpha$  is a subgroup of  $G_i$ .

- 396 16 Entropy and Groups
- 5. Let  $G_1, G_2, G_3$  be subgroups of a finite group G. Prove that

 $|G||G_1 \cap G_2 \cap G_3|^2 \ge |G_1 \cap G_2||G_2 \cap G_3||G_1 \cap G_3|.$ 

Hint: Use the information-theoretic approach.

6. Let  $\mathbf{h} \in \Gamma_2^*$  be the entropy function for random variables  $X_1$  and  $X_2$  such that  $h_1 + h_2 = h_{12}$ , i.e.  $X_1$  and  $X_2$  are independent. Let  $(G, G_1, G_2)$  be a group characterization of **h**, and define a mapping  $L: G_1 \times G_2 \to G$  by

$$L(a,b) = a \circ b.$$

- a) Prove that the mapping L is onto, i.e., for any element  $c \in G$ , there exists  $(a, b) \in G_1 \times G_2$  such that  $a \circ b = c$ .
- b) Prove that  $G_1 \circ G_2$  is a group.
- 7. Denote an entropy function  $\mathbf{h} \in \Gamma_2^*$  by  $(h_1, h_2, h_{12})$ . Construct a group characterization for each of the following entropy functions:
  - a)  $\mathbf{h}_1 = (\log 2, 0, \log 2)$
  - b)  $\mathbf{h}_2 = (0, \log 2, \log 2)$
  - c)  $\mathbf{h}_3 = (\log 2, \log 2, \log 2).$

Verify that  $\Gamma_2$  is the minimal convex set containing the above three entropy functions.

- 8. Denote an entropy function  $\mathbf{h} \in \Gamma_3^*$  by  $(h_1, h_2, h_3, h_{12}, h_{23}, h_{13}, h_{123})$ . Construct a group characterization for each of the following entropy functions:
  - a)  $\mathbf{h}_1 = (\log 2, 0, 0, \log 2, 0, \log 2, \log 2)$
  - b)  $\mathbf{h}_2 = (\log 2, \log 2, 0, \log 2, \log 2, \log 2, \log 2)$

  - d)  $\mathbf{h}_4 = (\log 2, \log 2, \log 2, \log 4, \log 4, \log 4, \log 4, \log 4).$
- 9. Ingleton inequality Let G be a finite Abelian group and  $G_1, G_2, G_3$ , and  $G_4$  be subgroups of G. Let  $(G, G_1, G_2, G_3, G_4)$  be a group characterization of  $\mathbf{g}$ , where  $\mathbf{g}$  is the entropy function for random variables  $X_1, X_2, X_3$ , and  $X_4$ . Prove the following statements: a)

$$(G_1 \cap G_3) \circ (G_1 \cap G_4) | \le |G_1 \cap (G_3 \circ G_4)|$$

Hint: Show that  $(G_1 \cap G_3) \circ (G_1 \cap G_4) \subset G_1 \cap (G_3 \circ G_4)$ . b)

$$|G_1 \circ G_3 \circ G_4| \le \frac{|G_1||G_3 \circ G_4||G_1 \cap G_3 \cap G_4|}{|G_1 \cap G_3||G_1 \cap G_4|}.$$

c)

$$|G_1 \circ G_2 \circ G_3 \circ G_4| \le \frac{|G_1 \circ G_3 \circ G_4| |G_2 \circ G_3 \circ G_4|}{|G_3 \circ G_4|}.$$

d)

$$\leq \frac{|G_1 \circ G_2 \circ G_3 \circ G_4|}{|G_1||G_2||G_3||G_4||G_1 \cap G_3 \cap G_4||G_2 \cap G_3 \cap G_4|} \leq \frac{|G_1||G_2||G_3||G_4||G_1 \cap G_3 \cap G_4||G_2 \cap G_3 \cap G_4||G_3 \cap G_4||G_4 \cap G_4||G_4 \cap G_4 \cap G_4||G_4 \cap G_4 \cap G_4||G_4 \cap G_4 \cap G_4||G_4 \cap G_4 \cap G_4||G_4 \cap G_4 \cap$$

e)

$$|G_1 \cap G_3||G_1 \cap G_4||G_2 \cap G_3||G_2 \cap G_4||G_3 \cap G_4| \leq |G_3||G_4||G_1 \cap G_2||G_1 \cap G_3 \cap G_4||G_2 \cap G_3 \cap G_4|.$$

f)

$$H(X_{13}) + H(X_{14}) + H(X_{23}) + H(X_{24}) + H(X_{34})$$
  

$$\geq H(X_3) + H(X_4) + H(X_{12}) + H(X_{134}) + H(X_{234}),$$

where  $H(X_{134})$  denotes  $H(X_1, X_3, X_4)$ , etc.

g) Is the inequality in f) implied by the basic inequalities? And does it always hold? Explain.

The Ingleton inequality [162] (see also [255]) was originally obtained as a constraint on the rank functions of vector spaces. The inequality in e) was obtained in the same spirit by Chan [55] for subgroups of a finite group. The inequality in f) is referred to as the Ingleton inequality for entropy in the literature. (See also Problem 8 in Chapter 15.)

## **Historical Notes**

The results in this chapter are due to Chan and Yeung [60], whose work was inspired by a one-to-one correspondence between entropy and quasi-uniform arrays previously established by Chan [55] (also Chan [56]). Romashchenko *et al.* [282] have developed an interpretation of Kolmogorov complexity similar to the combinatorial interpretation of entropy in Chan [55].

Fundamentals of Network Coding

# Introduction

For a point-to-point communication system, we see from Section 7.7 and Problem 6 in Chapter 8 that asymptotic optimality can be achieved by separating source coding and channel coding. Recall from Section 5.3 that the goal of source coding is to represent the information source in (almost) fair bits<sup>1</sup>. Then the role of channel coding is to enable the transmission of fair bits through the channel essentially free of error with no reference to the meaning of these fair bits. Thus a theme in classical information theory for point-topoint communication is that fair bits can be drawn equivalence to a *commodity*.

It is intuitively appealing that this theme in classical information theory would continue to hold in network communication where the network consists of *noiseless* point-to-point communication channels. If so, in order to multi $cast^2$  information from a source node to possibly more than one sink node, we only need to compress the information at the source node into fair bits, organize them into data packets, and route the packets to the sink node through the intermediate nodes in the network. In the case when there are more than one sink node, the information needs to be replicated at certain intermediate nodes so that every sink node can receive a copy of the information. This method of transmitting information in a network is generally referred to as store-and-forward—bb or routing. As a matter of fact, almost all computer networks built in the last few decades are based on this paradigm, where routers are deployed at the intermediate nodes to switch a data packet from an input channel to an output channel without processing the data content. The delivery of data packets in a computer network resembles mail delivery in a postal system. We refer the readers to textbooks on *data communication* [195] and switching theory [157][203].

 $<sup>^1</sup>$  Fair bits refer to i.i.d. bits, each being distributed uniformly on  $\{0,1\}$ . See Section 5.3.

 $<sup>^2\,</sup>$  Multicast means to transmit information from a source node to a specified set of sink nodes.

### 402 17 Introduction

However, we will see very shortly that in network communication, it does not suffice to simply route and/or replicate information within the network. Specifically, coding generally needs to be employed at the intermediate nodes in order to achieve bandwidth optimality. This notion, called *network coding*, is the subject of discussion in the second part of this book.

# 17.1 The Butterfly Network

In this section, the advantage of network coding over routing is explained by means of a few simple examples. The application of network coding in wireless and satellite communication will be discussed in the next section.

We will use a finite directed graph to represent a point-to-point communication network. A node in the network corresponds to a vertex in the graph, while a communication channel in the network corresponds to an edge in the graph. We will not distinguish a node from a vertex, nor will we distinguish a channel from an edge. In the graph, a node is represented by a circle, with the exception that the unique source node, denoted by s (if exists), is represented by a square. Each edge is labeled by a positive integer called the *capacity*<sup>3</sup> or the *rate constraint*, which gives the maximum number of information symbols taken from some finite alphabet that can be transmitted over the channel per unit time. In this section, we assume that the information symbol is binary. When there is only one edge from node a to node b, we denote the edge by (a, b).

Example 17.1 (Butterfly Network I). Consider the network in Figure 17.1(a). In this network, two bits  $b_1$  and  $b_2$  are generated at source node s, and they are to be multicast to two sink nodes  $t_1$  and  $t_2$ . In Figure 17.1(b), we try to devise a routing scheme for this purpose. By symmetry, we send the two bits on different output channels at node s. Without loss of generality,  $b_1$  is sent on channel (s, 1) and  $b_2$  is sent on channel (s, 2). At nodes 1 and 2, the received bit is replicated and the copies are sent on the two output channels. At node 3, since both  $b_1$  and  $b_2$  are received but there is only one output channel, we have to choose one of the two bits to be sent on the output channel (3, 4). Suppose we send  $b_1$  as in Figure 17.1(b). Then the bit is replicated at node 4 and the two copies are sent to nodes  $t_1$  and  $t_2$ , respectively. At node  $t_2$ , both  $b_1$  and  $b_2$  are received. However, at node  $t_1$ , two copies of  $b_1$  are received and  $b_2$  cannot be recovered. Thus this routing scheme does not work. Similarly, if  $b_2$  instead of  $b_1$  is sent on channel (3, 4),  $b_1$  cannot be recovered at node  $t_2$ .

However, if network coding is allowed, it is actually possible to achieve our goal. Figure 17.1(c) shows a scheme which multicasts both  $b_1$  and  $b_2$  to nodes  $t_1$  and  $t_2$ , where '+' denotes *modulo* 2 addition. At node  $t_1$ ,  $b_1$  is received, and  $b_2$  can be recovered by adding  $b_1$  and  $b_1 + b_2$ , because

<sup>&</sup>lt;sup>3</sup> Here the term "capacity" is used in the sense of graph theory.

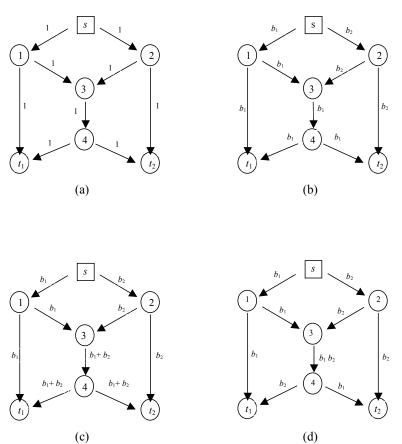


Fig. 17.1. Butterfly Network I.

$$b_1 + (b_1 + b_2) = (b_1 + b_1) + b_2 = 0 + b_2 = b_2.$$
 (17.1)

Similarly,  $b_2$  is received at node  $t_2$ , and  $b_1$  can be recovered by adding  $b_2$  and  $b_1 + b_2$ .

In this scheme,  $b_1$  and  $b_2$  are encoded into the bit  $b_1 + b_2$  which is then sent on channel (3, 4). If network coding is not allowed, in order to multicast both  $b_1$  and  $b_2$  to nodes  $t_1$  and  $t_2$ , at least one more bit has to be sent. Figure 17.1(d) shows such a scheme. In this scheme, however, the capacity of channel (3, 4) is exceeded by 1 bit. If the capacity of channel (3, 4) cannot be exceeded and network coding is not allowed, it can be shown that at most 1.5 bits can be multicast per unit time on the average (see Problem 3).

#### 404 17 Introduction

The above example shows the advantage of network coding over routing for a single multicast in a network. The next example shows the advantage of network coding over routing for multiple unicasts<sup>4</sup> in a network.

Example 17.2 (Butterfly Network II). In Figure 17.1, instead of both being generated at node s, suppose bit  $b_1$  is generated at node 1 and bit  $b_2$  is generated at node 2. Then we can remove node s and obtain the network in Figure 17.2(a). We again want to multicast  $b_1$  and  $b_2$  to both nodes  $t_1$  and  $t_2$ . Since this network is essentially the same as the previous one, Figure 17.2(b) shows the obvious network coding solution.

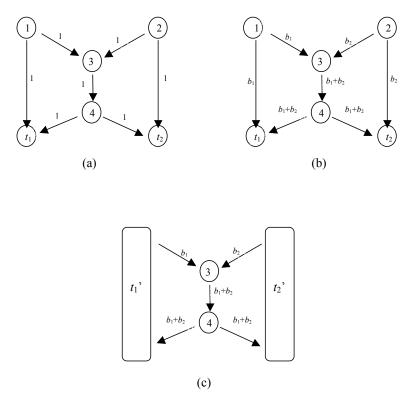


Fig. 17.2. Butterfly Network II.

There are two multicasts in this network. However, if we merge node 1 and node  $t_1$  into a new node  $t'_1$  and merge node 2 and node  $t_2$  into a new node  $t'_2$ , then we obtain the network and the corresponding network coding solution in Figure 17.2(c). In this new network, bits  $b_1$  and  $b_2$  are generated at nodes  $t'_1$  and  $t'_2$ , respectively, and the communication goal is to exchange the two

<sup>&</sup>lt;sup>4</sup> Unicast is the special case of multicast with one sink node.

bits through the network. In other words, the two multicasts in Figure 17.2(a) become two unicasts in Figure 17.2(c).

If network coding is not allowed, we need to route  $b_1$  from node  $t'_1$  to node  $t'_2$  and to route  $b_2$  from node  $t'_2$  to node  $t'_1$ . Since each of these routes has to go through node 3 and node 4, if  $b_1$  and  $b_2$  are routed simultaneously, the capacity of channel (3, 4) is exceeded. Therefore, we see the advantage of network coding over routing when there are multiple unicasts in the network.

For the network in Figure 17.2(b), the two sink nodes are required to recover both of the information sources, namely the bits  $b_1$  and  $b_2$ . Even though they are generated at two different source nodes 1 and 2, they can be regarded as being generated at a *super source node s* connecting to nodes  $t_1$  and  $t_2$  as in Figure 17.1(c). Precisely, the network (network code) in Figure 17.2(b) is obtained from the network (network code) in Figure 17.1(c) by removing node *s* and all its output channels. A further elaboration on this theme will be given in Example 19.26 in Chapter 19 when we discuss single-source linear network coding.

# 17.2 Wireless and Satellite Communications

In wireless communication, when a node broadcasts, different noisy versions of the signal is received by the neighboring nodes. Under certain conditions, with suitable channel coding, we can assume the existence of an error-free channel between the broadcast node and the neighboring nodes such that each of the latter receives exactly the same information. Such an abstraction, though generally suboptimal, provides very useful tools for communication systems design.

Our model for network communication can be used for modeling the above broadcast scenario by imposing the following constraints on the broadcast node:

- 1. all the output channels have the same capacity;
- 2. the same symbol is sent on each of the output channels.

We will refer to these constraints as the *broadcast constraint*. Figure 17.3(a) is an illustration of a broadcast node b with two neighboring nodes  $n_1$  and  $n_2$ , where the two output channels of node b have the same capacity.

In order to express the broadcast constraint in the usual graph-theoretic terminology, we need to establish the following simple fact about network coding.

**Proposition 17.3.** Network coding is not necessary at a node which has only one input channel and the capacity of each output channel is the same as that of the input channel.

406 17 Introduction

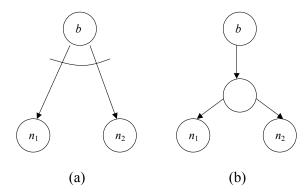


Fig. 17.3. A broadcast node b with two neighboring nodes  $n_1$  and  $n_2$ .

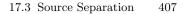
*Proof.* Consider a node in the network as prescribed and denote the symbol(s) received on the input channel by x. (There is more than one symbol in x if the input channel has capacity larger than 1.) Let a coding scheme be given, and denote the symbol sent on the *i*th output channel by  $g_i(x)$ .

We now show that one may assume without loss of generality that x is sent on all the output channels. If x instead of  $g_i(x)$  is sent on the *i*th output channel, then the receiving node can mimic the effect of receiving  $g_i(x)$ by applying the function  $g_i$  on x upon receiving it. In other words, any coding scheme that does not send x on all the output channels can readily be converted into one that does. This proves the proposition.  $\Box$ 

We now show that the broadcast constraint depicted in Figure 17.3(a) is logically equivalent to the usual graph representation in Figure 17.3(b). In this figure, the unlabeled node is a dummy node associated with the broadcast node which is inserted for the purpose of modeling the broadcast constraint, where the input channel and all the output channels of the dummy node have the same capacity as an output channel of the broadcast node b in Figure 17.3(a). Although no broadcast constraint is imposed on the dummy node in Figure 17.3(b), by Proposition 17.3, we may assume without loss of generality that the dummy node simply sends the symbol received on the input channel on each of the output channels. Then Figures 17.3(a) and (b) are logically equivalent to each other because a coding scheme for the former corresponds to a coding scheme for the latter, and vice versa.

Example 17.4 (A Wireless/Satellite System). Consider a communication system with two wireless nodes  $t'_1$  and  $t'_2$  that generate two bits  $b_1$  and  $b_2$ , respectively, and the two bits are to be exchanged through a relay node. Such a system can also be the model of a satellite communication system, where the relay node corresponds to a satellite, and the two nodes  $t'_1$  and  $t'_2$  correspond to ground stations that communicate with each other through the satellite.

We make the usual assumption that a wireless node cannot simultaneously



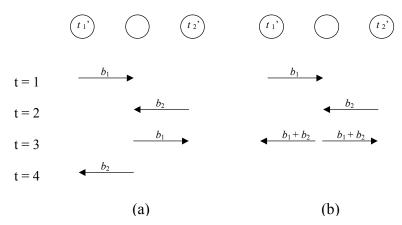


Fig. 17.4. A network coding application in wireless communication.

1. transmit and receive;

2. receive the transmission from more than one neighboring node.

A straightforward routing scheme which takes a total of 4 time units to complete is shown in Figure 17.4(a), with t being the time index.

By taking into account the broadcast nature of the relay node, the system can be modeled by the network in Figure 17.2(c), where node 3 corresponds to the relay node and node 4 corresponds to the associated dummy node. Then the network coding solution is shown in Figure 17.4(b), which takes a total of 3 time units to complete. In other words, a very simple coding scheme at the relay node can save 50 percent of the downlink bandwidth.

### **17.3 Source Separation**

In an error-free point-to-point communication system, suppose we want to transmit two information sources X and Y. If we compress the two sources separately, we need to transmit approximately H(X) + H(Y) bits. If we compress the two sources jointly, we need to transmit approximately H(X, Y) bits. If X and Y are independent, we have

$$H(X,Y) = H(X) + H(Y).$$
 (17.2)

In other words, if the information sources are independent, asymptotically there is no difference between coding them separately or jointly.

We will refer to coding independent information sources separately as source separation. Example 17.2 reveals the important fact that source separation is not necessary optimal in network communication, which is explained as follows. Let  $B_1$  and  $B_2$  be random bits generated at nodes  $t'_1$  and  $t'_2$ , respectively, where  $B_1$  and  $B_2$  are independent and each of them are distributed

#### 408 17 Introduction

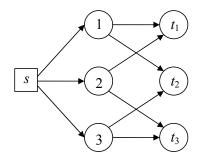
uniformly on  $\{0, 1\}$ . With  $B_2$  as side-information which is independent of  $B_1$ , node  $t'_2$  has to receive at least 1 bit in order to decode  $B_1$ . Since node  $t'_2$  can receive information only from node 4 which in turn can receive information only from node 3, any coding scheme that transmits  $B_1$  from node  $t'_1$  to node  $t'_2$ must send at least 1 bit on channel (3, 4). Similarly, any coding scheme that transmits  $B_2$  from node  $t'_2$  to node  $t'_1$  must send at least 1 bit on channel (3, 4). Therefore, any source separation solution must send at least 2 bits on channel (3, 4). Since the network coding solution in Figure 17.2(c) sends only 1 bit on channel (3, 4), we see that source separation is not optimal.

For a network coding problem with multiple information sources, since source separation does not guarantee optimality, the problem cannot always be decomposed into a number single-source problems. We will see that while single-source network coding has a relatively simple characterization, the characterization of multi-source network coding is much more involved.

### Problems

In the following problems, the rate constraint for an edge is in bits per unit time.

1. Consider the following network.

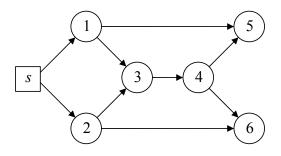


We want to multicast information to the sink nodes at the maximum rate without using network coding. Let  $B = \{b_1, b_2, \dots, b_\kappa\}$  be the set of bits to be multicast. Let  $B_i$  be the set of bits sent in edge (s, i), where  $|B_i| = 2$ , i = 1, 2, 3. At node *i*, the received bits are duplicated and sent in the two out-going edges. Thus two bits are sent in each edge in the network.

- a) Show that  $B = B_i \cup B_j$  for any  $1 \le i < j \le 3$ .
- b) Show that  $B_3 \cup (B_1 \cap B_2) = B$ .
- c) Show that  $|B_3 \cup (B_1 \cap B_2)| \le |B_3| + |B_1| + |B_2| |B_1 \cup B_2|$ .
- d) Determine the maximum value of  $\kappa$  and devise a network code which achieves this maximum value.
- e) What is the percentage of improvement if network coding is used?

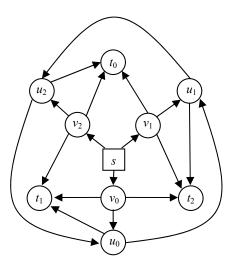
(Ahlswede et al. [7].)

2. Consider the following butterfly network.



Devise a network coding scheme which multicasts two bits  $b_1$  and  $b_2$  from node s to all the other nodes such that nodes 3, 5, and 6 receive  $b_1$  and  $b_2$  after 1 unit time and nodes 1, 2, and 4 receive  $b_1$  and  $b_2$  after 2 units of time. In other words, node i receives information at a rate equal to maxflow(s, i) for all  $i \neq s$ .

- 3. Determine the maximum rate at which information can be multicast to nodes 5 and 6 only in the network in Problem 2 if network coding is not used. Devise a network coding scheme which achieves this maximum rate.
- 4. Convolutional network code In the following network,  $maxflow(s, t_l) = 3$  for l = 1, 2, 3. The max-flow bound asserts that 3 bits can be multicast to



all the three sink nodes per unit time. We now describe a network coding scheme which achieve this. Let 3 bits  $b_0(k), b_1(k), b_2(k)$  be generated at node s at time  $k = 1, 2, \cdots$ , where we assume without loss of generality

410 17 Introduction

that  $b_l(k)$  is an element of the finite field GF(2). We adopt the convention that  $b_l(k) = 0$  for  $k \leq 0$ . At time  $k \geq 1$ , information transactions T1 to T11 occur in the following order:

- T1. *s* sends  $b_l(k)$  to  $v_l$ , l = 0, 1, 2
- T2.  $v_l$  sends  $b_l(k)$  to  $u_l$ ,  $t_{l\oplus 1}$ , and  $t_{l\oplus 2}$ , l = 0, 1, 2
- T3.  $u_0$  sends  $b_0(k) + b_1(k-1) + b_2(k-1)$  to  $u_1$
- T4.  $u_1$  sends  $b_0(k) + b_1(k-1) + b_2(k-1)$  to  $t_2$
- T5.  $u_1$  sends  $b_0(k) + b_1(k) + b_2(k-1)$  to  $u_2$
- T6.  $u_2$  sends  $b_0(k) + b_1(k) + b_2(k-1)$  to  $t_0$
- T7.  $u_2$  sends  $b_0(k) + b_1(k) + b_2(k)$  to  $u_0$
- T8.  $u_0$  sends  $b_0(k) + b_1(k) + b_2(k)$  to  $t_1$
- T9.  $t_2$  decodes  $b_2(k-1)$
- T10.  $t_0$  decodes  $b_0(k)$
- T11.  $t_1$  decodes  $b_1(k)$

where " $\oplus$ " denotes modulo 3 addition and "+" denotes modulo 2 addition.

- a) Show that the information transactions T1 to T11 can be performed at time k = 1.
- b) Show that T1 to T11 can be performed at any time  $k \geq 1$  by induction on k.
- c) Verify that at time k, nodes  $t_0$  and  $t_1$  can recover  $b_0(k')$ ,  $b_1(k')$ , and  $b_2(k')$  for all  $k' \leq k$ .
- d) Verify that at time k, node  $t_2$  can recover  $b_0(k')$  and  $b_1(k')$  for all  $k' \leq k$ , and  $b_2(k')$  for all  $k' \leq k 1$ . Note the unit time delay for  $t_2$  to recover  $b_2(k)$ .

(Ahlswede et al. [7].)

### **Historical Notes**

The concept of *network coding* was first introduced for satellite communication networks in Yeung and Zhang [368] and then fully developed in Ahlswede *et al.* [7], where in the latter the term "network coding" was coined. In this work, the advantage of network coding over store-and-forward was first demonstrated by the butterfly network, thus refuting the folklore that information transmission in a point-to-point network is equivalent to a commodity flow.

Prior to [368] and [7], network coding problems for special networks had been studied in the context of distributed source coding. The suboptimality of source separation was first demonstrated by Yeung [359]. Source separation was proved to be optimal for special networks by Hau [143], Roche *et al.* [279], and Yeung and Zhang [367]. Some other special cases of single-source network coding had been studied by Roche *et al.* [278], Rabin [268], Ayanoglu *et al.* [21], and Roche [277].

For a tutorial on the theory, we refer the reader to the unifying work by Yeung *et al.* [365]. Tutorials on the subject have also been written by Fragouli and Soljanin [111] and Chou and Wu [64] from the algorithm and application perspectives. We also refer the reader to the book by Ho and Lun [147]. For an update of the literature, the reader may visit the *Network Coding Homepage* [246].

# The Max-Flow Bound

In this chapter, we discuss an important bound for single-source network coding which has a strong connection with graph theory. This bound, called the max-flow min-cut bound, or simply the *max-flow bound*, gives a fundamental limit on the amount of information that can be multicast in the network.

The max-flow bound is established in a general setting where information can be transmitted within the network in some arbitrary manner. Toward this end, we first formally define a point-to-point network and a class of codes on such a network. In Chapters 19 and 20, we will prove the achievability of the max-flow bound by *linear* network coding<sup>1</sup>.

### **18.1** Point-to-Point Communication Networks

A point-to-point communication network is represented by a directed graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ , where  $\mathcal{V}$  is the set of nodes in the network and  $\mathcal{E}$  is the set of edges in  $\mathcal{G}$  which represent the point-to-point channels. Parallel edges between a pair of nodes is allowed<sup>2</sup>. We assume that  $\mathcal{G}$  is finite, i.e.,  $|\mathcal{E}| < \infty$  (and hence  $|\mathcal{V}| < \infty$ ). The unique *source node* in the network, where information is generated, is denoted by *s*. All the other nodes are referred to as *non-source* nodes. The sets of input channels and output channels of a node *i* are denoted by In(*i*) and Out(*i*), respectively.

For a channel e, let  $R_e$  be the *rate constraint*, i.e., the maximum number of information symbols taken from a finite alphabet that can be sent on the channel per unit time. As before, we also refer to  $R_e$  as the capacity of channel e in the sense of graph theory. Let

$$\mathbf{R} = [R_e : e \in \mathcal{E}] \tag{18.1}$$

<sup>&</sup>lt;sup>1</sup> A more specific form of the max-flow bound will be proved in Theorem 19.10 for linear network coding.

<sup>&</sup>lt;sup>2</sup> Such a graph is sometimes called a *multigraph*.

#### 414 18 The Max-Flow Bound

be the rate constraints for the graph  $\mathcal{G}$ . To simplify our discussion, we assume that  $R_e$  are positive integers for all  $e \in \mathcal{E}$ .

In the following, we introduce some notions in graph theory which will facilitate the characterization of a point-to-point network. Temporarily regard an edge in the graph  $\mathcal{G}$  as a water pipe and  $\mathcal{G}$  as a network of water pipes. Fix a node  $t \neq s$  and call it the *sink node*. Suppose water is generated at a constant rate at node s. We assume that the rate of water flow in each pipe does not exceed its capacity. We also assume that there is no leakage in the network, so that water is conserved at every node other than s and t in the sense that the total rate of water flowing into the node is equal to the total rate of water flowing out of the node. The water generated at node s is eventually drained at node t.

A flow

$$\mathbf{F} = [F_e : e \in \mathcal{E}] \tag{18.2}$$

in  $\mathcal{G}$  from node s to node t with respect to rate constraints **R** is a valid assignment of a nonnegative integer  $F_e$  to every edge  $e \in \mathcal{E}$  such that  $F_e$  is equal to the rate of water flow in edge e under all the assumptions in the last paragraph. The integer  $F_e$  is referred to as the value of **F** on edge e. Specifically, **F** is a flow in  $\mathcal{G}$  from node s to node t if for all  $e \in \mathcal{E}$ ,

$$0 \le F_e \le R_e,\tag{18.3}$$

and for all  $i \in \mathcal{V}$  except for s and t,

$$F_{+}(i) = F_{-}(i), \tag{18.4}$$

where

$$F_{+}(i) = \sum_{e \in \operatorname{In}(i)} F_{e} \tag{18.5}$$

and

$$F_{-}(i) = \sum_{e \in \operatorname{Out}(i)} F_e.$$
(18.6)

In the above,  $F_+(i)$  is the total flow into node *i* and  $F_-(i)$  is the total flow out of node *i*, and (18.4) is called the *conservation* conditions.

Since the conservation conditions require that the resultant flow out of any node other than s and t is zero, it is intuitively clear and not difficult to show that the resultant flow out of node s is equal to the resultant flow into node t. This common value is called the *value* of  $\mathbf{F}$ .  $\mathbf{F}$  is a *max-flow* from node s to node t in  $\mathcal{G}$  with respect to rate constraints  $\mathbf{R}$  if  $\mathbf{F}$  is a flow from node s to node t whose value is greater than or equal to the value of any other flow from node s to node t.

A cut between node s and node t is a subset  $\mathcal{U}$  of  $\mathcal{V}$  such that  $s \in \mathcal{U}$  and  $t \notin \mathcal{U}$ . Let

$$\mathcal{E}_{\mathcal{U}} = \{ e \in \mathcal{E} : e \in \operatorname{Out}(i) \cap \operatorname{In}(j) \text{ for some } i \in \mathcal{U} \text{ and } j \notin \mathcal{U} \}$$
(18.7)

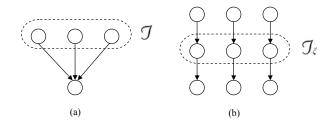


Fig. 18.1. Illustrations of the max-flow and the min-cut from the source node to (a) a collection of non-source node  $\mathcal{T}$  and (b) a collection of edges  $\xi$ .

be the set of edges across the cut  $\mathcal{U}$ . The *capacity* of the cut  $\mathcal{U}$  with respect to rate constraints **R** is defined as the sum of the capacities of all the edges across the cut, i.e.,

$$\sum_{e \in \mathcal{E}_{\mathcal{U}}} R_e. \tag{18.8}$$

A cut  $\mathcal{U}$  is a *min-cut* between node s and node t if it is a cut between node s and node t whose capacity is less than or equal to the capacity of any other cut between s and t.

A min-cut between node s and node t can be thought of as a *bottleneck* between node s and node t. Therefore, it is intuitively clear that the value of a max-flow from node s to node t cannot exceed the capacity of a min-cut between the two nodes. The following theorem, known as the *max-flow min-cut theorem*, states that the capacity of a min-cut is always achievable. This theorem will play a key role in the subsequent discussions.

**Theorem 18.1 (Max-Flow Min-Cut Theorem [104]).** Let  $\mathcal{G}$  be a graph with source node s, sink node t, and rate constraints  $\mathbf{R}$ . Then the value of a max-flow from node s to node t is equal to the capacity of a min-cut between the two nodes.

The notions of max-flow and min-cut can be generalized to a collection of non-source nodes  $\mathcal{T}$ . To define the max-flow and the min-cut from s to  $\mathcal{T}$ , we expand the graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  into  $\mathcal{G}' = (\mathcal{V}', \mathcal{E}')$  by installing a new node  $\tau$  which is connected from every node in  $\mathcal{T}$  by an edge. The capacity of an edge  $(t, \tau), t \in \mathcal{T}$ , is set to infinity. Intuitively, node  $\tau$  acts as a single sink node that collects all the flows into  $\mathcal{T}$ . Then the max-flow and the min-cut from node s to  $\mathcal{T}$  in graph  $\mathcal{G}$  are defined as the max-flow and the min-cut from node s to node  $\tau$  in graph  $\mathcal{G}'$ , respectively. This is illustrated in Figure 18.1(a).

The notions of max-flow and min-cut can be further generalized to a collection of edges  $\xi$ . For an edge  $e \in \xi$ , let the edge be from node  $v_e$  to node  $w_e$ . We modify the graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  to obtain the graph  $\tilde{\mathcal{G}} = (\tilde{\mathcal{V}}, \tilde{\mathcal{E}})$  by installing a new node  $t_e$  for each edge  $e \in \xi$  and replacing edge e by two new edges e' and e'', where e' is from node  $v_e$  to node  $t_e$  and e'' is from node  $t_e$  to node  $w_e$ . Let  $\mathcal{T}_{\xi}$  be the set of nodes  $t_e, e \in \xi$ . Then the max-flow and the min-cut between 416 18 The Max-Flow Bound

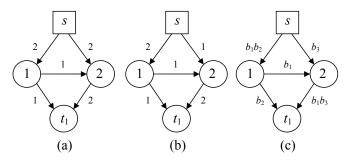


Fig. 18.2. A one-sink network.

node s and the collection of edges  $\xi$  in graph  $\mathcal{G}$  are defined as the max-flow and the min-cut between node s and the collection of nodes  $\mathcal{T}_{\xi}$  in graph  $\tilde{\mathcal{G}}$ , respectively. This is illustrated in Figure 18.1(b).

### 18.2 Examples Achieving the Max-Flow Bound

Let  $\omega$  be the rate at which information is multicast from source node s to sink nodes  $t_1, t_2, \dots, t_L$  in a network  $\mathcal{G}$  with rate constraints **R**. We are naturally interested in the maximum possible value of  $\omega$ . With a slight abuse of notation, we denote the value of a max-flow from source node s to a sink node  $t_l$  by maxflow( $t_l$ ). It is intuitive that

$$\omega \le \max flow(t_l) \tag{18.9}$$

for all  $l = 1, 2, \dots, L$ , i.e.,

$$\omega \le \min_{l} \max flow(t_l). \tag{18.10}$$

This is called the *max-flow bound*, which will be formally established in the next two sections. In this section, we first show by a few examples that the max-flow bound can be achieved. In these examples, the unit of information is the bit.

First, we consider the network in Figure 18.2 which has one sink node. Figure 18.2(a) shows the capacity of each edge. By identifying the min-cut to be  $\{s, 1, 2\}$  and applying the max-flow min-cut theorem, we see that

$$\max flow(t_1) = 3.$$
 (18.11)

Therefore the flow in Figure 18.2(b) is a max-flow. In Figure 18.2(c), we show how we can send three bits  $b_1, b_2$ , and  $b_3$  from node s to node  $t_1$  based on the max-flow in Figure 18.2(b). Evidently, the max-flow bound is achieved.

In fact, we can easily see that the max-flow bound can always be achieved when there is only one sink node in the network. In this case, we only need to

### 18.2 Examples Achieving the Max-Flow Bound 417

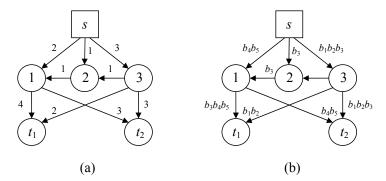


Fig. 18.3. A two-sink network without coding.

treat the information bits constituting the message as a commodity and route them through the network according to any fixed routing scheme. Eventually, all the bits will arrive at the sink node. Since the routing scheme is fixed, the sink node knows which bit is coming in from which edge, and the message can be recovered accordingly.

Next, we consider the network in Figure 18.3 which has two sink nodes. Figure 18.3(a) shows the capacity of each edge. It is easy to see that

$$\max flow(t_1) = 5 \tag{18.12}$$

and

$$\max flow(t_2) = 6.$$
 (18.13)

So the max-flow bound asserts that we cannot send more than 5 bits to both  $t_1$  and  $t_2$ . Figure 18.3(b) shows a scheme which sends 5 bits  $b_1, b_2, b_3, b_4$ , and  $b_5$  to  $t_1$  and  $t_2$  simultaneously. Therefore, the max-flow bound is achieved. In this scheme,  $b_1$  and  $b_2$  are replicated at node 3,  $b_3$  is replicated at node s, while  $b_4$  and  $b_5$  are replicated at node 1. Note that each bit is replicated exactly once in the network because two copies of each bit are needed to be sent to the two sink nodes.

We now revisit the butterfly network reproduced in Figure 18.4(a), which again has two sink nodes. It is easy to see that

$$\max flow(t_l) = 2 \tag{18.14}$$

for l = 1, 2. So the max-flow bound asserts that we cannot send more than 2 bits to both sink nodes  $t_1$  and  $t_2$ . We have already seen the network coding scheme in Figure 18.4(b) that achieves the max-flow bound. In this scheme, coding is required at node 3.

Finally, we consider the network in Figure 18.5 which has three sink nodes. Figure 18.5(a) shows the capacity of each edge. It is easy to see that

$$\max flow(t_l) = 2 \tag{18.15}$$

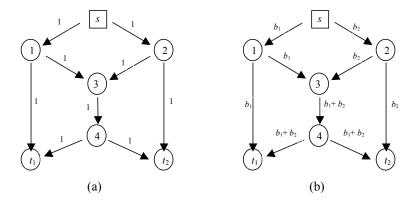


Fig. 18.4. Butterfly network I.

for all *l*. In Figure 18.5(b), we show how to multicast 2 bits  $b_1$  and  $b_2$  to all the sink nodes. Therefore, the max-flow bound is achieved. Again, it is necessary to code at the nodes in order to multicast the maximum number of bits to all the sink nodes.

The network in Figure 18.5 is of special interest in practice because it is a special case of the *diversity coding* scheme used in commercial *disk arrays*, which are a kind of fault-tolerant data storage system. For simplicity, assume the disk array has three disks which are represented by nodes 1, 2, and 3 in the network, and the information to be stored are the bits  $b_1$  and  $b_2$ . The information is encoded into three pieces, namely  $b_1$ ,  $b_2$ , and  $b_1 + b_2$ , which are stored on the disks represented by nodes 1, 2, and 3, respectively. In the system, there are three decoders, represented by sink nodes  $t_1$ ,  $t_2$ , and  $t_3$ , such that each of them has access to a distinct set of two disks. The idea is that when any one disk is out of order, the information can still be recovered from the remaining two disks. For example, if the disk represented by node 1 is out of order, then the information can be recovered by the decoder represented by

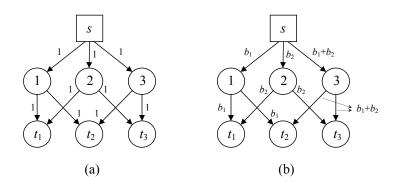


Fig. 18.5. A diversity coding scheme.

sink node  $t_3$  which has access to the disks represented by node 2 and node 3. When all the three disks are functioning, the information can be recovered by any decoder.

# 18.3 A Class of Network Codes

In this section, we introduce a general class of codes for the point-to-point network defined in Section 18.1. In the next section, the max-flow bound will be proved for this class of network codes.

Since the max-flow bound concerns only the values of max-flows from source node s to the sink nodes, we assume without loss of generality that there is no loop in the graph  $\mathcal{G}$ , i.e.,  $\operatorname{In}(i) \cap \operatorname{Out}(i) = \emptyset$  for all  $i \in \mathcal{V}$ , because such edges do not increase the value of a max-flow from node s to a sink node. For the same reason, we assume that there is no input edge at node s, i.e.,  $\operatorname{In}(s) = \emptyset$ .

We consider a block code of length n. Let X denote the information source and assume that x, the outcome of X, is obtained by selecting an index from a set  $\mathcal{X}$  according to the uniform distribution. The elements in  $\mathcal{X}$  are called messages. The information sent on an output channel of a node can depend only on the information previously received by that node. This constraint specifies the *causality* of any coding scheme on the network.

An  $(n, (\eta_e : e \in \mathcal{E}), \tau)$  network code on the graph  $\mathcal{G}$  that multicasts information from source node s to sink nodes  $t_1, t_2, \dots, t_L$ , where n is the block length, is defined by the components listed below; the construction of the code from these components will be described after their definitions are given.

1) A positive integer K.

2) Mappings

$$u: \{1, 2, \cdots, K\} \to \mathcal{V},\tag{18.16}$$

$$v: \{1, 2, \cdots, K\} \to \mathcal{V},\tag{18.17}$$

and

$$\hat{e}: \{1, 2, \cdots, K\} \to \mathcal{E},\tag{18.18}$$

such that  $\hat{e}(k) \in \text{Out}(u(k))$  and  $\hat{e}(k) \in \text{In}(v(k))$ . 3) Index sets  $A_k = \{1, 2, \dots, |A_k|\}, 1 \le k \le K$ , such that

$$\prod_{k\in T_e} |A_k| = \eta_e, \tag{18.19}$$

where

$$T_e = \{ 1 \le k \le K : \hat{e}(k) = e \}.$$
(18.20)

420 18 The Max-Flow Bound

4) (Encoding functions). If u(k) = s, then

$$f_k: \mathcal{X} \to A_k, \tag{18.21}$$

where

$$\mathcal{X} = \{1, 2, \cdots, \lceil 2^{n\tau} \rceil\}.$$
(18.22)

If 
$$u(k) \neq s$$
, if

$$Q_k = \{1 \le k' < k : v(k') = u(k)\}$$
(18.23)

is nonempty, then

$$f_k: \prod_{k' \in Q_k} A_{k'} \to A_k; \tag{18.24}$$

otherwise, let  $f_k$  be an arbitrary constant taken from  $A_k$ .

5) (Decoding functions). Mappings

$$g_l: \prod_{k' \in W_l} A_{k'} \to \mathcal{X} \tag{18.25}$$

for  $l = 1, 2, \cdots, L$ , where

$$W_l = \{1 \le k \le K : v(k) = t_l\}$$
(18.26)

such that for all  $l = 1, 2, \cdots, L$ ,

$$\tilde{g}_l(x) = x \tag{18.27}$$

for all  $x \in \mathcal{X}$ , where  $\tilde{g}_l$  is the function from  $\mathcal{X}$  to  $\mathcal{X}$  induced inductively by  $f_k, 1 \leq k \leq K$  and  $g_l$ , such that  $\tilde{g}_l(x)$  denotes the value of  $g_l$  as a function of x.

The quantity  $\tau$  is the rate of the information source X, which is also the rate at which information is multicast from the source node to all the sink nodes. The  $(n, (\eta_e : e \in \mathcal{E}), \tau)$  code is constructed from the above components as follows. At the beginning of a coding session, the value of X is available to node s. During the coding session, there are K transactions which take place in chronological order, where each transaction refers to a node sending information to another node. In the kth transaction, node u(k) encodes according to encoding function  $f_k$  and sends an index in  $A_k$  to node v(k). The domain of  $f_k$  is the set of all possible information that can be received by node u(k)just before the kth transaction, and we distinguish two cases. If u(k) = s, the domain of  $f_k$  is  $\mathcal{X}$ . If  $u(k) \neq s$ ,  $Q_k$  gives the time indices of all the previous transactions for which information was sent to node u(k), so the domain of  $f_k$  is  $\prod_{k' \in Q_k} A_{k'}$ . The set  $T_e$  gives the time indices of all the transactions for which information is sent on channel e, so  $\eta_e$  is the number of possible index tuples that can be sent on channel e during the coding session. Finally,  $W_l$  gives the indices of all the transactions for which information is sent to node  $t_l$ , and  $g_l$  is the decoding function at node  $t_l$  which recovers x with zero error.

# 18.4 Proof of the Max-Flow Bound

In this section, we state and prove the max-flow bound for the class of network codes defined in the last section.

**Definition 18.2.** For a graph  $\mathcal{G}$  with rate constraints  $\mathbf{R}$ , an information rate  $\omega \geq 0$  is asymptotically achievable if for any  $\epsilon > 0$ , there exists for sufficiently large n an  $(n, (\eta_e : e \in \mathcal{E}), \tau)$  network code on  $\mathcal{G}$  such that

$$n^{-1}\log_2\eta_e \le R_e + \epsilon \tag{18.28}$$

for all  $e \in \mathcal{E}$ , where  $n^{-1} \log_2 \eta_e$  is the average bit rate of the code on channel e, and

$$\tau \ge \omega - \epsilon. \tag{18.29}$$

For brevity, an asymptotically achievable information rate will be referred to as an achievable information rate.

**Remark** It follows from the above definition that if  $\omega \geq 0$  is achievable, then  $\omega'$  is also achievable for all  $0 \leq \omega' \leq \omega$ . Also, if  $\omega^{(k)}$  is achievable for all  $k \geq 1$ , then it can be shown that  $\omega = \lim_{k\to\infty} \omega^{(k)}$ , if exists, is also achievable. Therefore, the set of all achievable information rates is closed and fully characterized by the maximum value in the set.

**Theorem 18.3 (Max-Flow Bound).** For a graph  $\mathcal{G}$  with rate constraints **R**, if  $\omega$  is achievable, then

$$\omega \le \min_{l} \max flow(t_l). \tag{18.30}$$

*Proof.* It suffices to prove that for a graph  $\mathcal{G}$  with rate constraints  $\mathbf{R}$ , if for any  $\epsilon > 0$  there exists for sufficiently large n an  $(n, (\eta_e : e \in \mathcal{E}), \tau)$  code on  $\mathcal{G}$  such that

$$n^{-1}\log_2\eta_e \le R_e + \epsilon \tag{18.31}$$

for all  $e \in \mathcal{E}$  and

$$\tau \ge \omega - \epsilon, \tag{18.32}$$

then  $\omega$  satisfies (18.30).

Consider such a code for a fixed  $\epsilon$  and a sufficiently large n, and consider any  $l = 1, 2, \dots, L$  and any cut  $\mathcal{U}$  between node s and node  $t_l$ . Let

$$w_j(x) = (\tilde{f}_k(x) : k \in \bigcup_{e \in \text{In}(j)} T_e),$$
 (18.33)

where  $x \in \mathcal{X}$  and  $\tilde{f}_k$  is the function from  $\mathcal{X}$  to  $A_k$  induced inductively by  $f_{k'}, 1 \leq k' \leq k$ , such that  $\tilde{f}_k(x)$  denotes the value of  $f_k$  as a function of x. The tuple  $w_j(x)$  is all the information known by node j during the whole coding session when the message is x. Since  $\tilde{f}_k(x)$  is a function of the information

#### 42218 The Max-Flow Bound

previously received by node u(k), it can be shown by induction (see Problem 3) that  $w_{t_l}(x)$  is a function of  $\tilde{f}_k(x), k \in \bigcup_{e \in \mathcal{E}_{\mathcal{U}}} T_e$ , where  $\mathcal{E}_{\mathcal{U}}$  is the set of edges across the cut  $\mathcal{U}$  as previously defined in (18.7). Since x can be determined at node  $t_l$ , we have

$$H(X) \le H(X, w_{t_l}(X))$$
 (18.34)

$$=H(w_{t_l}(X)) \tag{18.35}$$

$$\stackrel{a)}{\leq} H\left(\tilde{f}_k(X), k \in \bigcup_{e \in \mathcal{E}_{\mathcal{U}}} T_e\right)$$
(18.36)

$$\stackrel{b)}{\leq} \sum_{e \in \mathcal{E}_{\mathcal{U}}} \sum_{k \in T_e} H(\tilde{f}_k(X)) \tag{18.37}$$

$$\stackrel{c)}{\leq} \sum_{e \in \mathcal{E}_{\mathcal{U}}} \sum_{k \in T_e} \log_2 |A_k| \tag{18.38}$$

$$=\sum_{e\in\mathcal{E}_{\mathcal{U}}}\log_2\left(\prod_{k\in T_e}|A_k|\right)$$
(18.39)

$$\stackrel{d)}{\leq} \sum_{e \in \mathcal{E}_{\mathcal{U}}} \log_2 \eta_e, \tag{18.40}$$

where

- a) follows because  $w_{t_l}(x)$  is a function of  $\tilde{f}_k(x), k \in \bigcup_{e \in \mathcal{E}_{\mathcal{U}}} T_e$ ; •
- b) follows from the independence bound for entropy (Theorem 2.39); •
- c) follows from (18.21) and Theorem 2.43; •
- d) follows from (18.19). •

Thus

$$\omega - \epsilon \le \tau \tag{18.41}$$
$$\le n^{-1} \log_2 \lceil 2^{n\tau} \rceil \tag{18.42}$$

$$\leq n^{-1} \log_2 |2^{n/2}| \tag{18.42}$$

$$= n^{-1} \log_2 |\mathcal{X}|$$
(18.43)  
=  $n^{-1} H(X)$ (18.44)

$$= n^{-1}H(X) (18.44)$$

$$\leq \sum_{e \in \mathcal{E}_{\mathcal{U}}} n^{-1} \log_2 \eta_e \tag{18.45}$$

$$\leq \sum_{e \in \mathcal{E}_{\mathcal{U}}} (R_e + \epsilon) \tag{18.46}$$

$$\leq \sum_{e \in \mathcal{E}_{\mathcal{U}}} R_e + |\mathcal{E}|\epsilon, \tag{18.47}$$

where (18.45) follows from (18.40). Minimizing the right hand side over all  $\mathcal{U}$ , we have

18.4 Proof of the Max-Flow Bound 423

$$\omega - \epsilon \le \min_{\mathcal{U}} \sum_{e \in \mathcal{E}_{\mathcal{U}}} R_e + |\mathcal{E}|\epsilon.$$
(18.48)

The first term on the right hand side is the capacity of a min-cut between node s and node  $t_l$ . By the max-flow min-cut theorem, it is equal to the value of a max-flow from node s to node  $t_l$ , i.e., maxflow $(t_l)$ . Letting  $\epsilon \to 0$ , we obtain

$$\omega \le \max flow(t_l). \tag{18.49}$$

Since this upper bound on  $\omega$  holds for all  $l = 1, 2, \dots, L$ ,

$$\omega \le \min_{l} \max flow(t_l). \tag{18.50}$$

The theorem is proved.  $\Box$ 

**Remark 1** In proving the max-flow bound, the time evolution and the causality of the network code have been taken into account.

**Remark 2** Even if we allow an arbitrarily small probability of decoding error in the usual Shannon sense, by modifying our proof by means of a standard application of Fano's inequality, it can be shown that it is still necessary for  $\omega$  to satisfy (18.50). The details are omitted here.

#### Problems

- 1. In a network, for a flow **F** from a source node s to a sink node t, show that  $F_+(s) = F_-(t)$  provided that the conservation conditions in (18.4) hold.
- 2. For the class of codes defined in Section 18.3, show that
  - a) If the rates  $\omega^{(k)}$  are achievable for all  $k \ge 1$ , then  $\omega = \lim_{k \to \infty} \omega^{(k)}$ , if exists, is also achievable.
  - b) The tuple  $w_{t_l}(x)$  is a function of  $\tilde{f}_k(x), k \in \bigcup_{e \in \mathcal{E}_{\mathcal{U}}} T_e$ , where  $\mathcal{U}$  is any cut between source node s and sink node  $t_l$ .
- 3. Prove the claim in the proof of Theorem 18.3 that for any cut  $\mathcal{U}$  between node s and node  $t_l, w_{t_l}(x)$  is a function of  $\tilde{f}_k(x), k \in \bigcup_{e \in \mathcal{E}_{\mathcal{U}}} T_e$ . Hint: Define

$$w_{j,\kappa}(x) = (f_k(x) : k \in \bigcup_{e \in \operatorname{In}(j)} T_e \text{ and } k \le \kappa)$$

and prove by induction on  $\kappa$  that for all  $1 \leq \kappa \leq K$ ,  $(w_{j,\kappa}(x) : j \notin \mathcal{U})$  is a function of  $(\tilde{f}_k(x) : k \in \bigcup_{e \in \mathcal{E}_{\mathcal{U}}} T_e$  and  $k \leq \kappa)$ .

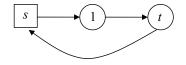
4. Probabilistic network code For a network code defined in Section 18.3, the kth transaction of the coding process is specified by a mapping  $f_k$ . Suppose instead of a mapping  $f_k$ , the kth transaction is specified by a transition probability matrix from the domain of  $f_k$  to the range of  $f_k$ . Also, instead of a mapping  $g_l$ , decoding at sink node  $t_l$  is specified by a transition probability matrix from the domain of  $g_l$  to the range of  $g_l$ ,  $1 \le l \le L$ .

#### 424 18 The Max-Flow Bound

Conditioning on the indices received by node u(k) during  $1 \leq k' < k$ , the index sent from node u(k) to node v(k) in the kth transaction is independent of all the previously generated random variables. Similarly, conditioning on all the indices received by sink node  $t_l$  during the whole coding session, the decoding at  $t_l$  is independent of all the previously generated random variables.

We refer to such a code as a probabilistic network code. Since a deterministic network code is a special case of a probabilistic network code, the latter can potentially multicast at a higher rate compared with the former. Prove that this is not possible.

5. Consider a probabilistic network code on the network below.



Let  $X = (X_1, X_2)$  be uniformly distributed on  $GF(2)^2$ , and Z be independent of X and uniformly distributed on GF(2). We use  $\tilde{F}_k$  to denote the index transmitted in the kth transaction and  $W_{t_l}$  to denote  $(\tilde{F}_k, k \in \bigcup_{e \in \text{In}(t_l)} T_e)$ . The probabilistic network code is specified by the following 5 transactions:

$$u(1) = s, \quad v(1) = 1, \quad \tilde{F}_1 = X_1,$$
  

$$u(2) = 1, \quad v(2) = t, \quad \tilde{F}_2 = X_1 + Z,$$
  

$$u(3) = t, \quad v(3) = s, \quad \tilde{F}_3 = X_1 + Z,$$
  

$$u(4) = s, \quad v(4) = 1, \quad \tilde{F}_4 = (X_1, X_2 + Z),$$
  

$$u(5) = 1, \quad v(5) = t, \quad \tilde{F}_5 = (X_1, X_2 + Z).$$

Note that the 4th transaction is possible because upon knowing  $X_1$  and  $X_1 + Z$ , Z can be determined.

- a) Determine  $W_t$ .
- b) Verify that X can be recovered from  $W_t$ .
- c) Show that  $X \to (\tilde{F}_1, \tilde{F}_4) \to W_t$  does not form a Markov chain. Here,  $\tilde{F}_1$  and  $\tilde{F}_4$  are all the random variables sent on edge (s, 1) during the coding session. Although node t receives all the information through the edge (s, 1), the Markov chain in c) does not hold. (Ahlswede *et al.* [7].)

# **Historical Notes**

The max-flow bound presented in this chapter was proved by Ahlswede et al. [7], where the point-to-point channels in the network are noiseless.

The max-flow bound can be established when the point-to-point channels in the network are discrete memoryless channels. Borade [44] proved the bound with the assumptions that the channels are independent of each other and that the transmissions in the channels are synchronous. Song *et al.* [309] proved the bound without the latter assumption. These results are network generalizations of the result by Shannon [295] asserting that the capacity of a discrete memoryless channel is not increased by feedback (see Section 7.6), and they imply the asymptotic optimality of separating network coding and channel coding under the corresponding assumptions.

In the last chapter, we have established the max-flow bound as the fundamental bound for multicasting a single information source in a point-to-point communication network. In the next two chapters, we will construct *linear network codes* that achieve the max-flow bound at various levels of generality.

A finite field is a system of symbols on which one can perform operations corresponding to the four operations in arithmetic for real numbers, namely addition, subtraction, multiplication, and division. The set of real numbers together with the associated operations are referred to as the field of real numbers, or simply the real field. Unlike the real field that has an infinite number of elements, a finite field has only a finite number of elements. For finite field theory, we refer the reader to [238]. For our discussions here, since we will not make use of the detailed structural properties of a finite field, the reader may by and large regard the algebra on a finite field and the algebra on the real field as the same.

In a linear network code, all the information symbols are regarded as elements of a finite field F called the *base field*. These include the symbols that comprise the information source as well as the symbols transmitted on the channels. For example, F is taken to be the binary field GF(2) when the information unit is the bit. Furthermore, encoding and decoding are based on linear algebra defined on the based field, so that efficient algorithms for encoding and decoding as well as for code construction can be obtained.

In this chapter, we consider acyclic networks, i.e., networks with no directed cycle. We study the network coding problem in which a message consisting of a finite block of symbols is multicast. We make the ideal assumption that the propagation delay in the network, which includes the processing delay at the nodes and the transmission delay over the channels, is zero. In a general setting, a pipeline of messages may be multicast, and the propagation delay may be non-negligible. If the network is acyclic, then the operations in the network can be so synchronized that sequential messages are processed independent of each other. In this way, the network coding problem is inde-

pendent of the propagation delay. Therefore, it suffices to study the network coding problem as described.

On the other hand, when a network contains directed cycles, the processing and transmission of sequential messages can convolve with together. Then the amount of delay incurred becomes part of the consideration in network coding. This will be discussed in the next chapter.

### **19.1 Acyclic Networks**

Denote a directed network by  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ , where  $\mathcal{V}$  and  $\mathcal{E}$  are the sets of nodes and channels, respectively. A pair of channels  $(d, e) \in \mathcal{E} \times \mathcal{E}$  is called an *adjacent pair* if there exists a node  $t \in \mathcal{V}$  such that  $d \in \text{In}(t)$  and  $e \in \text{Out}(t)$ . A *directed path* in  $\mathcal{G}$  is a sequence of channels

$$e_1, e_2, \cdots, e_m \tag{19.1}$$

such that  $(e_i, e_{i+1})$  is an adjacent pair for all  $1 \leq i < m$ . Let  $e_1 \in \text{Out}(t)$ and  $e_m \in \text{In}(t')$ . The sequence in (19.1) is called a directed path from  $e_1$  to  $e_m$ , or equivalently, a directed path from node t to node t'. If t = t', then the directed path is called a *directed cycle*. A directed network  $\mathcal{G}$  is *cyclic* if it contains a directed cycle, otherwise  $\mathcal{G}$  is *acyclic*.

Acyclic networks are easier to handle because the nodes in the network can be ordered in a way which allows encoding at the nodes to be carried out in a sequential and consistent manner. The following proposition and its proof describe such an order.

**Proposition 19.1.** If  $\mathcal{G}$  is a finite directed acyclic graph, then it is possible to order the nodes of  $\mathcal{G}$  in a sequence such that if there is an edge from node i to node j, then node i appears before node j in the sequence.

*Proof.* We partition the set  $\mathcal{V}$  into subsets  $\mathcal{V}_1, \mathcal{V}_2, \cdots$ , such that node i is in  $\mathcal{V}_k$  if and only if the length of a longest directed path ending at node i is equal to k. We first prove that if node i is in  $\mathcal{V}_{k'}$  and node j is in  $\mathcal{V}_k$  such that there exists a directed path from node i to node j, then k' < k. Since the length of a longest directed path ending at node i is equal to k' and there exists a directed path from node i to node j (with length at least equal to 1), there exists a directed path ending at node j with length equal to k' + 1. As node j is in  $\mathcal{V}_k$ , we have

$$k' + 1 \le k,\tag{19.2}$$

so that

$$k' < k. \tag{19.3}$$

Hence, by listing the nodes of  $\mathcal{G}$  in a sequence such that the nodes in  $\mathcal{V}_{k'}$  appear before the nodes in  $\mathcal{V}_k$  if k' < k, where the order of the nodes within

each  $\mathcal{V}_k$  is arbitrary, we obtain an order of the nodes of  $\mathcal{G}$  with the desired property.  $\Box$ 

Following the direction of the edges, we will refer to an order prescribed by Proposition 19.1 as an *upstream-to-downstream order*<sup>1</sup>. For a given acyclic network, such an order (not unique) is implicitly assumed. The nodes in the network encodes according to this order, referred to as the *encoding order*. Then whenever a node encodes, all the information needed would have already been received on the input channels of that node.

*Example 19.2.* Consider ordering the nodes in the butterfly network in Figure 17.1 by the sequence

$$s, 2, 1, 3, 4, t_2, t_1.$$
 (19.4)

It is easy to check that in this sequence, if there is a directed path from node i to node j, then node i appears before node j.

## 19.2 Linear Network Codes

In this section, we formulate a linear network code on an acyclic network  $\mathcal{G}$ . By allowing parallel channels between a pair of nodes, we assume without loss of generality that all the channels in the network have unit capacity, i.e., one symbol in the base field F can be transmitted on each channel. There exists a unique node s in  $\mathcal{G}$ , called the source node, where a message consisting of  $\omega$  symbols taken from the base field F is generated. To avoid trivially, we assume that every non-source node has at least one input channel.

As in Section 18.3, we assume that there is no loop in  $\mathcal{G}$ , and there is no input channel at node s. To facilitate our discussion, however, we let  $\operatorname{In}(s)$  be a set of  $\omega$  imaginary channels that terminate at node s but have no originating nodes. The reader may think of the  $\omega$  symbols forming the message as being received by source node s on these  $\omega$  imaginary channels. We emphasize that these imaginary channels are not part of the network, and the number of these channels is context dependent. Figure 19.1(a) illustrates the butterfly network with  $\omega = 2$  imaginary channels appended at source node s.

Two directed paths  $P_1$  and  $P_2$  in  $\mathcal{G}$  are *edge-disjoint* if the two paths do not share a common channel. It is not difficult to see from the conservation conditions in (18.4) that for a non-source node t, the maximum number of edge-disjoint paths from node s to node t is equal to maxflow(t).

The message generated at source node s, consisting of  $\omega$  symbols in the base field F, is represented by a row  $\omega$ -vector  $\mathbf{x} \in F^{\omega}$ . Based on the value of  $\mathbf{x}$ , source node s transmits a symbol over each output channel. Encoding at the nodes in the network is carried out according to a certain upstream-to-downstream order. At a node in the network, the ensemble of received symbols

<sup>&</sup>lt;sup>1</sup> Also called an ancestral order in graph theory.

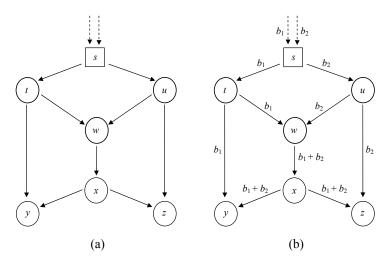


Fig. 19.1. (a) Two imaginary channels are appended to the source node of the butterfly network. (b) A 2-dimensional network code for the butterfly network.

is mapped to a symbol in F specific to each output channel, and the symbol is sent on that channel. The following definition of a network code formally describes this mechanism. Since the code so defined is not necessarily linear, the base field F can be regarded in this context as any finite alphabet.

**Definition 19.3 (Local Description of a Network Code).** An  $\omega$ -dimensional network code on an acyclic network over a base field F consists of a local encoding mapping

$$\tilde{k}_e: F^{|\mathrm{In}(t)|} \to F \tag{19.5}$$

for every channel e in the network, where  $e \in Out(t)$ .

With the encoding mechanism as described, the local encoding mappings derive recursively the symbols transmitted over all channels e, denoted by  $\tilde{f}_e(\mathbf{x})$ . The above definition of a network code does not explicitly give the values of  $\tilde{f}_e(\mathbf{x})$ , whose mathematical properties are at the focus of the present discussion. Therefore, we also present an equivalent definition below, which describes a network code in terms of both the local encoding mechanisms as well as the recursively derived values  $\tilde{f}_e(\mathbf{x})$ .

**Definition 19.4 (Global Description of a Network Code).** An  $\omega$ dimensional network code on an acyclic network over a base field F consists of a local encoding mapping

$$\tilde{k}_e: F^{|\mathrm{In}(t)|} \to F \tag{19.6}$$

and a global encoding mapping

19.2 Linear Network Codes 431

$$\tilde{f}_e: F^\omega \to F \tag{19.7}$$

for each channel e in the network, where  $e \in Out(t)$ , such that:

- (19.8) For every node t and every channel  $e \in \text{Out}(t)$ ,  $\hat{f}_e(\mathbf{x})$  is uniquely determined by  $(\tilde{f}_d(\mathbf{x}), d \in \text{In}(t))$  via the local encoding mapping  $\tilde{k}_e$ .
- (19.9) For the  $\omega$  imaginary channels e, the mappings  $\tilde{f}_e$  form the standard basis of the vector space  $F^{\omega}$ .

*Example 19.5.* Let  $\mathbf{x} = [b_1 b_2]$  denote a generic row vector in  $GF(2)^2$ . Figure 19.1(b) shows a 2-dimensional binary network code for the butterfly network with the following global encoding mappings:

$$f_e(x) = b_1$$
 for  $e = (o, s), (s, t), (t, w), (t, y)$  (19.10)

$$\hat{f}_e(x) = b_2$$
 for  $e = (o, s)', (s, u), (u, w), (u, z)$  (19.11)

$$\hat{f}_e(x) = b_1 + b_2$$
 for  $e = (w, x), (x, y), (x, z),$  (19.12)

where (o, s) and (o, s)' denote the two imaginary channels at node s. The corresponding local encoding mappings are

$$\tilde{k}_{(s,t)}(b_1, b_2) = b_1, \ \tilde{k}_{(s,u)}(b_1, b_2) = b_2,$$
(19.13)

$$k_{(t,w)}(b_1) = k_{(t,y)}(b_1) = b_1, (19.14)$$

$$\tilde{k}_{(u,w)}(b_2) = \tilde{k}_{(u,z)}(b_2) = b_2, \ \tilde{k}_{(w,x)}(b_1,b_2) = b_1 + b_2,$$
 (19.15)

etc.

When a global encoding mapping  $\tilde{f}_e$  is linear, it corresponds to a column  $\omega$ -vector  $\mathbf{f}_e$  such that  $\tilde{f}_e(x)$  is the product  $\mathbf{x} \cdot \mathbf{f}_e$ , where the row  $\omega$ -vector  $\mathbf{x}$  is the message generated at node s. Similarly, when a local encoding mapping  $\tilde{k}_e$ , where  $e \in \text{Out}(t)$ , is linear, it corresponds to a column |In(t)|-vector  $\mathbf{k}_e$  such that  $\tilde{k}_e(\mathbf{y}) = \mathbf{y} \cdot \mathbf{k}_e$ , where  $\mathbf{y} \in F^{|\text{In}(t)|}$  is the row vector representing the symbols received at node t. In an  $\omega$ -dimensional network code on an acyclic network, if all the local encoding mappings are linear, then so are the global encoding mappings since they are functional compositions of the local encoding mappings are all linear, then so are the local encoding mappings. We leave the proof as an exercise.

In the following, we formulate a linear network code as a network code whose local and global encoding mappings are all linear. Again, both the local and global descriptions are presented even though they are equivalent. The global description of a linear network code will be very useful when we construct such codes in Section 19.4.

**Definition 19.6 (Local Description of a Linear Network Code).** An  $\omega$ -dimensional linear network code on an acyclic network over a base field F consists of a scalar  $k_{d,e}$ , called the local encoding kernel, for every adjacent pair of channels (d, e) in the network. The  $|\text{In}(t)| \times |\text{Out}(t)|$  matrix

$$K_t = [k_{d,e}]_{d \in \operatorname{In}(t), e \in \operatorname{Out}(t)}$$
(19.16)

is called the local encoding kernel at node t.

Note that the matrix structure of  $K_t$  implicitly assumes an ordering among the channels.

**Definition 19.7 (Global Description of a Linear Network Code).** An  $\omega$ -dimensional linear network code on an acyclic network over a base field F consists of a scalar  $k_{d,e}$  for every adjacent pair of channels (d, e) in the network as well as a column  $\omega$ -vector  $\mathbf{f}_e$  for every channel e such that:

(19.17)  $\mathbf{f}_e = \sum_{d \in \text{In}(t)} k_{d,e} \mathbf{f}_d$ , where  $e \in \text{Out}(t)$ . (19.18) The vectors  $\mathbf{f}_e$  for the  $\omega$  imaginary channels  $e \in \text{In}(s)$  form the standard basis of the vector space  $F^{\omega}$ .

The vector  $\mathbf{f}_e$  is called the global encoding kernel for channel e.

We now explain how the global description above specifies the linear network code. Initially, source node s generates a message  $\mathbf{x}$  as a row  $\omega$ -vector. In view of (19.18), the symbols in  $\mathbf{x}$  are regarded as being received by source node s on the imaginary channels as  $\mathbf{x} \cdot \mathbf{f}_d$ ,  $d \in \text{In}(s)$ . Starting at source node s, any node t in the network receives the symbols  $\mathbf{x} \cdot \mathbf{f}_d$ ,  $d \in \text{In}(t)$ , from which it calculates the symbol  $\mathbf{x} \cdot \mathbf{f}_e$  for sending on each channel  $e \in \text{Out}(t)$  via the linear formula

$$\mathbf{x} \cdot \mathbf{f}_e = \mathbf{x} \sum_{d \in \text{In}(t)} k_{d,e} \, \mathbf{f}_d = \sum_{d \in \text{In}(t)} k_{d,e} (\mathbf{x} \cdot \mathbf{f}_d), \tag{19.19}$$

where the first equality follows from (19.17). In this way, the symbol  $\mathbf{x} \cdot \mathbf{f}_e$  is transmitted on any channel e (which may be an imaginary channel) in the network.

Given the local encoding kernels for all the channels in an acyclic network, the global encoding kernels can be calculated recursively in any upstream-todownstream order by (19.17), while (19.18) provides the boundary conditions.

**Remark** A partial analogy can be drawn between the global encoding kernels for the channels in a linear network code and the columns of a generator matrix of a linear block code in algebraic coding theory [209][37][341]. The former are indexed by the channels in the network, while the latter are indexed by "time." However, the global encoding kernels in a linear network code are constrained by the network topology via (19.17), while the columns in the generator matrix of a linear block code in general are not subject to any such constraint.

The following two examples illustrate the relation between the local encoding kernels and the global encoding kernels of a linear network code. The reader should understand these two examples thoroughly before proceeding to the next section.

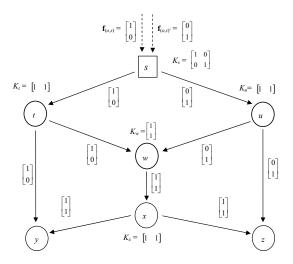


Fig. 19.2. The global and local encoding kernels for the 2-dimensional linear network code in Example 19.8.

*Example 19.8.* The network code in Figure 19.1(b) is in fact linear. Assume the alphabetical order among the channels  $(o, s), (o, s)', (s, t), \dots, (x, z)$ . Then the local encoding kernels at nodes are:

$$K_s = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, \quad K_t = K_u = K_x = \begin{bmatrix} 1 & 1 \end{bmatrix}, \quad K_w = \begin{bmatrix} 1 \\ 1 \end{bmatrix}.$$
(19.20)

The corresponding global encoding kernels are:

$$\mathbf{f}_{e} = \begin{cases} \begin{bmatrix} 1\\0 \end{bmatrix} \text{ for } e = (o, s), (s, t), (t, w), \text{ and } (t, y) \\ \begin{bmatrix} 0\\1 \end{bmatrix} \text{ for } e = (o, s)', (s, u), (u, w), \text{ and } (u, z) \\ \begin{bmatrix} 1\\1 \end{bmatrix} \text{ for } e = (w, x), (x, y), \text{ and } (x, z). \end{cases}$$
(19.21)

The local/global encoding kernels are summarized in Figure 19.2. In fact, they describe a 2-dimensional linear network code regardless of the choice of the base field.

*Example 19.9.* For a general 2-dimensional linear network code on the network in Figure 19.2, the local encoding kernels at the nodes can be expressed as

$$K_s = \begin{bmatrix} a & c \\ b & d \end{bmatrix}, \quad K_t = \begin{bmatrix} e & f \end{bmatrix}, \quad K_u = \begin{bmatrix} g & h \end{bmatrix}, \quad (19.22)$$

$$K_w = \begin{bmatrix} i \\ j \end{bmatrix}, \ K_x = \begin{bmatrix} k \ l \end{bmatrix}, \tag{19.23}$$

where  $a, b, c, \dots, l$ , the entries of the matrices, are indeterminates in the base field F. Starting with

$$\mathbf{f}_{(o,s)} = \begin{bmatrix} 1\\0 \end{bmatrix} \text{ and } \mathbf{f}_{(o,s)'} = \begin{bmatrix} 0\\1 \end{bmatrix}, \qquad (19.24)$$

we can obtain all the global encoding kernels below by applying (19.17) recursively:

$$\mathbf{f}_{(s,t)} = \begin{bmatrix} a \\ b \end{bmatrix}, \ \mathbf{f}_{(s,u)} = \begin{bmatrix} c \\ d \end{bmatrix}, \ \mathbf{f}_{(t,w)} = \begin{bmatrix} ae \\ be \end{bmatrix}, \ \mathbf{f}_{(t,y)} = \begin{bmatrix} af \\ bf \end{bmatrix},$$
(19.25)

$$\mathbf{f}_{(u,w)} = \begin{bmatrix} cg\\ dg \end{bmatrix}, \ \mathbf{f}_{(u,z)} = \begin{bmatrix} ch\\ dh \end{bmatrix}, \ \mathbf{f}_{(w,x)} = \begin{bmatrix} aei + cgj\\ bei + dgj \end{bmatrix},$$
(19.26)

$$\mathbf{f}_{(x,y)} = \begin{bmatrix} aeik + cgjk\\ beik + dgjk \end{bmatrix}, \ \mathbf{f}_{(x,z)} = \begin{bmatrix} aeil + cgjl\\ beil + dgjl \end{bmatrix}.$$
 (19.27)

For example,  $\mathbf{f}_{(w,x)}$  is obtained from  $\mathbf{f}_{(t,w)}$  and  $\mathbf{f}_{(u,w)}$  by

$$\mathbf{f}_{(w,x)} = k_{(t,w),(w,x)} \mathbf{f}_{(t,w)} + k_{(u,w),(w,x)} \mathbf{f}_{(u,w)}$$
(19.28)

$$= i \begin{bmatrix} ae\\be \end{bmatrix} + j \begin{bmatrix} cg\\dg \end{bmatrix}$$
(19.29)

$$= \begin{bmatrix} aei + cgj\\ bei + dgj \end{bmatrix}.$$
 (19.30)

The local/global encoding kernels of the general linear network code are summarized in Figure 19.3.

# 19.3 Desirable Properties of a Linear Network Code

We have proved in Section 18.4 that in a communication network represented by a graph  $\mathcal{G}$ , the rate at which information is transmitted from source node sto any node t cannot exceed maxflow(t), the value of a max-flow from node sto node t. For a collection of non-source nodes  $\mathcal{T}$ , denote by maxflow $(\mathcal{T})$  the value of a max-flow from node s to  $\mathcal{T}$ . Then it is readily seen that the rate at which information is transmitted from source node s to the collection of nodes  $\mathcal{T}$  cannot exceed maxflow $(\mathcal{T})$ .

In the sequel, we adopt the conventional notation  $\langle \cdot \rangle$  for the linear span of a set of vectors. For a node t, let

$$V_t = \langle \{ \mathbf{f}_e : e \in \mathrm{In}(t) \} \rangle \tag{19.31}$$

and for a collection  $\mathcal{T}$  of nodes, let

$$V_{\mathcal{T}} = \langle \cup_{t \in \mathcal{T}} V_t \rangle. \tag{19.32}$$

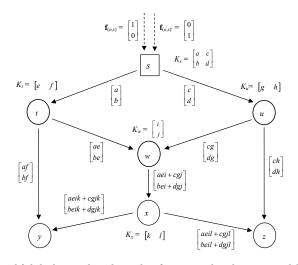


Fig. 19.3. Local/global encoding kernels of a general 2-dimensional linear network code.

For a collection  $\xi$  of channels, let

$$V_{\xi} = \langle \{ \mathbf{f}_e : e \in \xi \} \rangle, \tag{19.33}$$

with the convention  $V_{\emptyset} = \{0\}$ , where 0 denotes the zero column  $\omega$ -vector.

In the next theorem, we first establish a specific form of the max-flow bound which applies to linear network coding.

**Theorem 19.10 (Max-Flow Bound for Linear Network Coding).** For an  $\omega$ -dimensional linear network code on an acyclic network, for any collection  $\mathcal{T}$  of non-source nodes,

$$\dim(V_{\mathcal{T}}) \le \min\{\omega, \maxflow(\mathcal{T})\}.$$
(19.34)

*Proof.* Let the acyclic network be  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ . Consider a cut  $\mathcal{U}$  between source node s and a collection  $\mathcal{T}$  of non-source nodes, and let  $\mathcal{E}_{\mathcal{U}}$  be the set of edges across the cut  $\mathcal{U}$  as in (18.7). Then  $V_{\mathcal{T}}$  is a linear transformation of  $V_{\mathcal{E}_{\mathcal{U}}}$ , where

$$\dim(V_{\mathcal{T}}) \le \dim(V_{\mathcal{E}_{\mathcal{U}}}) \le |\mathcal{E}_{\mathcal{U}}| \tag{19.35}$$

Minimizing over all the cuts between s and  $\mathcal{T}$  and invoking the max-flow min-cut theorem, we have

$$\dim(V_{\mathcal{T}}) \le \max flow(\mathcal{T}). \tag{19.36}$$

On the other hand,  $V_{\mathcal{T}}$  is a linear transformation of the standard basis of  $F^{\omega}$  whose dimension is equal to  $\omega$ . Therefore,

$$\dim(V_{\mathcal{T}}) \le \omega. \tag{19.37}$$

Then the proof is completed by combining (19.36) and (19.37).  $\Box$ 

For a collection of channels  $\xi \subset \mathcal{E}$  (i.e., not including the imaginary channels), we denote by maxflow( $\xi$ ) the value of a max-flow from source node s to  $\xi$ . Theorem 19.10 has the following straightforward corollary.

**Corollary 19.11.** For an  $\omega$ -dimensional linear network code on an acyclic network, for any collection of channels  $\xi \subset \mathcal{E}$ ,

$$\dim(V_{\xi}) \le \min\{\omega, \maxflow(\xi)\}.$$
(19.38)

Whether the max-flow bound in Theorem 19.10 or Corollary 19.11 is achievable depends on the network topology, the dimension  $\omega$ , and the coding scheme. Three special classes of linear network codes are defined below by the achievement of this bound to three different extents.

**Definition 19.12.** An  $\omega$ -dimensional linear network code on an acyclic network qualifies as a linear multicast, a linear broadcast, or a linear dispersion, respectively, if the following hold:

(19.39) dim $(V_t) = \omega$  for every non-source node t with maxflow $(t) \ge \omega$ .

- (19.40)  $\dim(V_t) = \min\{\omega, \max(t)\}\$  for every non-source node t.
- (19.41) dim  $(V_{\mathcal{T}}) = \min\{\omega, \max flow(\mathcal{T})\}\$  for every collection  $\mathcal{T}$  of non-source nodes.

For a set  $\xi$  of channels, including possibly the imaginary channels, let

$$F_{\xi} = \left[ \mathbf{f}_e \right]_{e \in \xi} \tag{19.42}$$

be the  $\omega \times |\xi|$  matrix obtained by putting  $\mathbf{f}_e, e \in \xi$  in juxtaposition. For a node t, the symbols  $\mathbf{x} \cdot \mathbf{f}_e, e \in \text{In}(t)$  are received on the input channels. Equivalently, the row |In(t)|-vector

$$\mathbf{x} \cdot F_{\mathrm{In}(t)} \tag{19.43}$$

is received. Obviously, the message  $\mathbf{x}$ , consisting of  $\omega$  information units, can be uniquely determined at the node if and only if the rank of  $F_{\text{In}(t)}$  is equal to  $\omega$ , i.e.,

$$\dim(V_t) = \omega. \tag{19.44}$$

The same applies to a collection  $\mathcal{T}$  of non-source nodes.

For a linear multicast, a node t can decode the message **x** if and only if  $\max flow(t) \ge \omega$ . For a node t with  $\max flow(t) < \omega$ , nothing is guaranteed. An application of an  $\omega$ -dimensional linear multicast is for multicasting information at rate  $\omega$  to all (or some of) those non-source nodes with max-flow at least equal to  $\omega$ .

For a linear broadcast, like a linear multicast, a node t can decode the message **x** if and only if maxflow(t)  $\geq \omega$ . For a node t with maxflow(t)  $< \omega$ , the set of all received vectors, namely

$$\{\mathbf{x} \cdot F_{\mathrm{In}(t)} : \mathbf{x} \in F^{\omega}\},\tag{19.45}$$

form a vector subspace of  $F^{\omega}$  with dimension equal to maxflow(t), but there is no guarantee on which such subspace is actually received<sup>2</sup>. An application of linear broadcast is for multicasting information on a network at a variable rate (see Problem 13). A random version of linear broadcast (to be discussed in Section 19.4) is also useful for identifying the max-flow of a non-source in an unknown network topology [320].

For a linear dispersion, a collection  $\mathcal{T}$  of non-source nodes can decode the message  $\mathbf{x}$  if and only if maxflow $(\mathcal{T}) \geq \omega$ . If maxflow $(\mathcal{T}) < \omega$ , the collection  $\mathcal{T}$  receives a vector subspace with dimension equal to maxflow $(\mathcal{T})$ . Again, there is no guarantee on which such subspace is actually received. An application of linear dispersion is in a two-tier network system consisting of the backbone network and a number of local area networks (LANs), where each LAN is connected to one or more nodes on the backbone network. An information source with rate  $\omega$ , generated at a node s in the backbone network, is to be transmitted to every user on the LANs. With a linear dispersion on the backbone network, every user on a LAN can receive the information source as long as the LAN acquires through the backbone network an aggregated max-flow from node s at least equal to  $\omega$ . Moreover, new LANs can be established under the same criterion without modifying the linear dispersion on the backbone network.

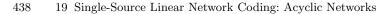
Note that for all the three classes of linear network codes in Definition 19.12, a sink node is not explicitly identified. Also, it is immediate from the definition that every linear dispersion is a linear broadcast, and every linear broadcast is a linear multicast. The example below shows that a linear broadcast is not necessarily a linear dispersion, a linear multicast is not necessarily a linear broadcast, and a linear network code is not necessarily a linear multicast.

*Example 19.13.* Figure 19.4(a) shows a 2-dimensional linear dispersion on an acyclic network with the global encoding kernels as prescribed. Figure 19.4(b) shows a 2-dimensional linear broadcast on the same network that is not a linear dispersion because

$$\max flow(\{t, u\}) = 2 = \omega, \tag{19.46}$$

while the global encoding kernels of the channels in  $\operatorname{In}(t) \cup \operatorname{In}(u)$  span only a 1dimensional subspace. Figure 19.4(c) shows a 2-dimensional linear multicast that is not a linear broadcast since node u receives no information at all. Finally, the 2-dimensional linear network code in Figure 19.4(d) is not a linear multicast.

<sup>&</sup>lt;sup>2</sup> Here  $F^{\omega}$  refers to the row vector space.



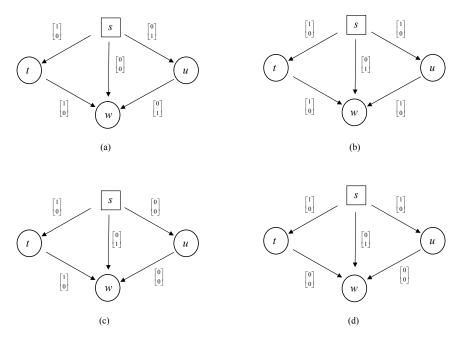


Fig. 19.4. (a) A 2-dimensional linear dispersion over an acyclic network. (b) A 2-dimensional linear broadcast that is not a linear dispersion. (c) A 2-dimensional linear multicast that is not a linear broadcast. (d) A 2-dimensional linear network code that is not a linear multicast.

Example 19.14. The linear network code in Example 19.8 meets all the criteria (19.39) through (19.41) in Definition 19.12. Thus it is a 2-dimensional linear dispersion, and hence also a linear broadcast and linear multicast, regardless of the choice of the base field. The same applies to the linear network code in Figure 19.4(a).

Example 19.15. The general linear network code in Example 19.9 meets the criterion (19.39) for a linear multicast when

- f<sub>(t,w)</sub> and f<sub>(u,w)</sub> are linearly independent;
  f<sub>(t,y)</sub> and f<sub>(x,y)</sub> are linearly independent;
  f<sub>(u,z)</sub> and f<sub>(x,z)</sub> are linearly independent.

Equivalently, the criterion says that e, f, g, h, k, l, ad - bc, abei + adgj - baei - baeibcgj, and daei + dcgj - cbei - cdgj are all nonzero. Example 19.8 has been the special case with

$$a = d = e = f = g = h = i = j = k = l = 1$$
 (19.47)

and

$$b = c = 0. (19.48)$$

19.3 Desirable Properties of a Linear Network Code 439

#### Transformation of a Linear Network Code

Consider an  $\omega$ -dimensional linear network code C on an acyclic network. Suppose source node s, instead of encoding the message  $\mathbf{x}$ , encodes

$$\mathbf{x}' = \mathbf{x}A,\tag{19.49}$$

where A is an *invertible*  $\omega \times \omega$  matrix. Then the symbol sent on a channel  $e \in \mathcal{E}$  is given by

$$\mathbf{x}' \cdot \mathbf{f}_e = (\mathbf{x}A) \cdot \mathbf{f}_e = \mathbf{x} \cdot (A \mathbf{f}_e). \tag{19.50}$$

This gives a new linear network code  $\mathcal{C}'$  with respect to the message  ${\bf x}$  with global encoding kernels

$$\mathbf{f}'_e = \begin{cases} A \, \mathbf{f}_e & \text{if } e \in \mathcal{E} \\ \mathbf{f}_e & \text{if } e \in \mathrm{In}(s). \end{cases}$$
(19.51)

Recall the definition of the matrix  $F_{\xi}$  in (19.42) for a set of channels  $\xi$ . Then (19.17) can be written in matrix form as

$$F_{\text{Out}(t)} = F_{\text{In}(t)} K_t \tag{19.52}$$

for all nodes t, where  $K_t$  is the local encoding kernel at node t. Similarly, letting

$$F'_{\xi} = \begin{bmatrix} \mathbf{f}'_e \end{bmatrix}_{e \in \xi}, \tag{19.53}$$

we obtain from (19.51) that

$$F'_{\operatorname{Out}(t)} = AF_{\operatorname{Out}(t)} \tag{19.54}$$

for all nodes t,

$$F'_{\mathrm{In}(t)} = AF_{\mathrm{In}(t)} \tag{19.55}$$

for all nodes  $t \neq s$ , and

$$F'_{\text{In}(s)} = F_{\text{In}(s)}.$$
 (19.56)

For a node  $t \neq s$ , from (19.54), (19.52), and (19.55),

$$F'_{\text{Out}(t)} = AF_{\text{Out}(t)} \tag{19.57}$$

$$=A(F_{\mathrm{In}(t)}K_t)\tag{19.58}$$

$$= (AF_{\mathrm{In}(t)})K_t \tag{19.59}$$

$$=F_{\mathrm{In}(t)}'K_t.\tag{19.60}$$

Since  $\mathbf{f}_e, e \in \text{In}(s)$  form the standard basis of  $F^{\omega}$ ,

$$F_{\text{In}(s)} = F'_{\text{In}(s)} = I,$$
 (19.61)

the  $\omega\times\omega$  identity matrix. It then follows from (19.54), (19.52), and (19.61) that

$$F'_{\operatorname{Out}(s)} = AF_{\operatorname{Out}(s)} \tag{19.62}$$

$$=A(F_{\mathrm{In}(s)}K_s)\tag{19.63}$$

$$=AK_s \tag{19.64}$$

$$=F'_{\text{In}(s)}(AK_s).$$
 (19.65)

Comparing (19.60) and (19.65) with (19.52), we see that the local encoding kernels of C' are given by

$$K'_t = \begin{cases} K_t & \text{if } t \neq s \\ AK_s & \text{if } t = s. \end{cases}$$
(19.66)

The network code C' is called the *transformation* of the network code C by the (invertible) matrix A. In view of Definition 19.12, the requirements of a linear multicast, a linear broadcast, and a linear dispersion are all in terms of the linear independence among the global encoding kernels. We leave it as an exercise for the reader to show that if a network code is a linear multicast, broadcast, or dispersion, then any transformation of it is also a linear multicast, broadcast, or dispersion, respectively.

Suppose C is an  $\omega$ -dimensional linear multicast and let C' be a transformation of C. When the network code C' is employed, the message  $\mathbf{x}$  can be decoded by any node t with maxflow $(t) \geq \omega$ , because from the foregoing C' is also a linear multicast. For the purpose of multicasting, there is no difference between C and C', and they can be regarded as equivalent.

If C is an  $\omega$ -dimensional linear broadcast and C' is a transformation of C, then C' is also an  $\omega$ -dimensional linear broadcast. However, C as a linear broadcast may deliver to a particular node t with maxflow $(t) < \omega$  a certain subset of symbols in the message  $\mathbf{x}$ , while C' may not be able to achieve the same. Then whether C and C' can be regarded as equivalent depends on the specific requirements of the application. As an example, the linear network code in Figure 19.1(b) delivers  $b_1$  to node t. However, taking a transformation of the network code with matrix

$$A = \begin{bmatrix} 1 & 0 \\ 1 & 1 \end{bmatrix}, \tag{19.67}$$

the resulting network code can no longer deliver  $b_1$  to node t, although nodes w, v, and z can continue to decode both  $b_1$  and  $b_2$ .

#### Implementation of a Linear Network Code

In implementation of a linear network code, be it a linear multicast, a linear broadcast, a linear dispersion, or any linear network code, in order that the code can be used as intended, the global encoding kernels  $\mathbf{f}_e, e \in \text{In}(t)$  must be known by each node t if node t is to recover any useful information from the symbols received on the input channels. These global encoding kernels can be made available ahead of time if the code is already decided. Alternatively, they

can be delivered through the input channels if multiple usage of the network is allowed.

One possible way to deliver the global encoding kernels to node t in a coding session of length n, where  $n > \omega$ , is as follows. At time  $k = 1, 2, \dots, \omega$ , the source node transmits the dummy message  $\mathbf{m}_k$ , a row  $\omega$ -vector with all the components equal to 0 except that the kth component is equal to 1. Note that

$$\begin{bmatrix} \mathbf{m}_1 \\ \mathbf{m}_2 \\ \vdots \\ \mathbf{m}_\omega \end{bmatrix} = I_\omega, \tag{19.68}$$

the  $\omega \times \omega$  identity matrix. At time  $k = \omega + i$ , where  $i = 1, 2, \dots, n - \omega$ , the source node transmits the message  $\mathbf{x}_i$ . Then throughout the coding session, node t receives

$$\begin{array}{c|c} \mathbf{m}_{1} \\ \mathbf{m}_{2} \\ \vdots \\ \mathbf{m}_{\omega} \\ \mathbf{x}_{1} \\ \mathbf{x}_{2} \\ \vdots \\ \mathbf{x}_{n-\omega} \end{array} \right| F_{\mathrm{In}(t)} = \begin{bmatrix} I_{\omega} \\ \mathbf{x}_{1} \\ \mathbf{x}_{2} \\ \vdots \\ \mathbf{x}_{n-\omega} \end{array} \right| F_{\mathrm{In}(t)} = \begin{bmatrix} F_{\mathrm{In}(t)} \\ \mathbf{x}_{1} \cdot F_{\mathrm{In}(t)} \\ \mathbf{x}_{2} \cdot F_{\mathrm{In}(t)} \\ \vdots \\ \mathbf{x}_{n-\omega} \cdot F_{\mathrm{In}(t)} \end{bmatrix}$$
(19.69)

on the input channels. In other words, the global encoding kernels of the input channels at node t are received at the beginning of the coding session. This applies to all the sink nodes in the network simultaneously because the  $\omega$  dummy messages do not depend on the particular node t. If  $F_{\text{In}(t)}$  has full rank, then node t can start to decode  $\mathbf{x}_1$  upon receiving  $\mathbf{x}_1 \cdot F_{\text{In}(t)}$ .

Since  $n - \omega$  messages are transmitted in a coding session of length n, the utilization of the network is equal to  $(n - \omega)/n$ , which tends to 1 as  $n \to \infty$ . That is, the overhead for delivering the global encoding kernels through the network is asymptotically negligible.

# **19.4** Existence and Construction

For a given acyclic network, the following three factors dictate the existence of an  $\omega$ -dimensional linear network code with desirable properties:

- the value of  $\omega$ ,
- the network topology,
- the choice of the base field *F*.

We begin with an example illustrating the third factor.

*Example 19.16.* On the network in Figure 19.5, a 2-dimensional ternary linear multicast can be constructed by the following local encoding kernels at the nodes:

$$K_s = \begin{bmatrix} 0 & 1 & 1 & 1 \\ 1 & 0 & 1 & 2 \end{bmatrix} \text{ and } K_{u_i} = \begin{bmatrix} 1 & 1 & 1 \end{bmatrix}$$
(19.70)

for  $1 \leq i \leq 4$ . On the other hand, we can prove the nonexistence of a 2dimensional binary linear multicast on this network as follows. Assuming the contrary that a 2-dimensional binary linear multicast exists, we will derive a contradiction. Let the global encoding kernel  $\mathbf{f}_{(s,u_i)} = [y_i \ z_i]^{\top}$  for  $1 \leq i \leq 4$ . Since maxflow $(t_k) = 2$  for all  $1 \leq k \leq 6$ , the global encoding kernels for the two input channels to each node  $t_k$  must be linearly independent. Thus, if node  $t_k$  is at the downstream of both nodes  $u_i$  and  $u_j$ , then the two vectors  $[y_i \ z_i]^{\top}$  and  $[y_i \ z_i]^{\top}$  must be linearly independent. As each node  $t_k$  is at the downstream of a different pair of nodes among  $u_1, u_2, u_3$ , and  $u_4$ , the four vectors  $[y_i \ z_i]^{\top}$ ,  $1 \leq i \leq 4$ , are pairwise linearly independent, and consequently, must be four distinct vectors in  $GF(2)^2$ . Then one of them must be  $[0 \ 0]^{\top}$  since there are only four vectors.

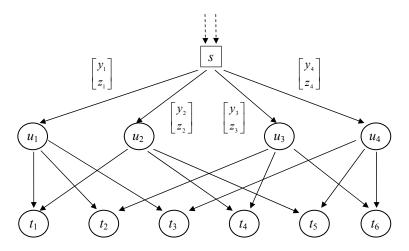


Fig. 19.5. A network with a 2-dimensional ternary linear multicast but without a 2-dimensional binary linear multicast.

In order for the linear network code to qualify as a linear multicast, a linear broadcast, or a linear dispersion, it is required that certain collections of global encoding kernels span the maximum possible dimensions. This is equivalent to certain polynomial functions taking nonzero values, where the indeterminates of these polynomials are the local encoding kernels. To fix ideas, take  $\omega = 3$ , consider a node t with two input channels, and put the

global encoding kernels of these two channels in juxtaposition to form a  $3 \times 2$  matrix. Then, this matrix attains the maximum possible rank of 2 if and only if there exists a  $2 \times 2$  submatrix with nonzero determinant.

According to the local description, a linear network code is specified by the local encoding kernels, and the global encoding kernels can be derived recursively in an upstream-to-downstream order. From Example 19.14, it is not hard to see that every component in a global encoding kernel is a polynomial function whose indeterminates are the local encoding kernels.

When a nonzero value of a polynomial function is required, it does not merely mean that at least one coefficient in the polynomial is nonzero. Rather, it means a way to choose scalar values for the indeterminates so that the polynomial function is evaluated to a nonzero scalar value.

When the base field is small, certain polynomial equations may be unavoidable. For instance, for any prime number p, the polynomial equation  $z^p - z = 0$  is satisfied for any  $z \in GF(p)$ . The nonexistence of a binary linear multicast in Example 19.16 can also trace its root to a set of polynomial equations that cannot be avoided simultaneously over GF(2).

However, when the base field is sufficiently large, every nonzero polynomial function can indeed be evaluated to a nonzero value with a proper choice of the values taken by the set of indeterminates involved. This is formally stated in the following elementary lemma, which will be instrumental in the proof of Theorem 19.20 asserting the existence of a linear multicast on an acyclic network when the base field is sufficiently large.

**Lemma 19.17.** Let  $g(z_1, z_2, \dots, z_n)$  be a nonzero polynomial with coefficients in a field F. If |F| is greater than the degree of g in every  $z_j$  for  $1 \le j \le n$ , then there exist  $a_1, a_2, \dots, a_n \in F$  such that

$$g(a_1, a_2, \cdots, a_n) \neq 0.$$
 (19.71)

*Proof.* The proof is by induction on n. For n = 0, g is a nonzero constant in F, and the lemma is obviously true. Assume that the lemma is true for n-1 for some  $n \ge 1$ . Express  $g(z_1, z_2, \dots, z_n)$  as a polynomial in  $z_n$  with coefficients in the polynomial ring  $F[z_1, z_2, \dots, z_{n-1}]$ , i.e.,

$$g(z_1, z_2, \cdots, z_n) = h(z_1, z_2, \cdots, z_{n-1}) z_n^{\ k} + \cdots,$$
(19.72)

where k is the degree of g in  $z_n$  and the leading coefficient  $h(z_1, z_2, \dots, z_{n-1})$ is a nonzero polynomial in  $F[z_1, z_2, \dots, z_{n-1}]$ . By the induction hypothesis, there exist  $a_1, a_2, \dots, a_{n-1} \in F$  such that  $h(a_1, a_2, \dots, a_{n-1}) \neq 0$ . Thus  $g(a_1, a_2, \dots, a_{n-1}, z)$  is a nonzero polynomial in z with degree k < |F|. Since this polynomial cannot have more than k roots in F and |F| > k, there exists  $a_n \in F$  such that

$$g(a_1, a_2, \cdots, a_{n-1}, a_n) \neq 0.$$
 (19.73)

**Corollary 19.18.** Let  $g(z_1, z_2, \dots, z_n)$  be a nonzero polynomial with coefficients in a field F with |F| > m, where m is the highest degree of g in  $z_i$  for

 $1 \leq j \leq n$ . Let  $a_1, a_2, \dots, a_n$  be chosen independently according to the uniform distribution on F. Then

$$\Pr\{g(a_1, a_2, \cdots, a_n) \neq 0\} \ge \left(1 - \frac{m}{|F|}\right)^n.$$
(19.74)

In particular,

$$\Pr\{g(a_1, a_2, \cdots, a_n) \neq 0\} \to 1 \tag{19.75}$$

as  $|F| \to \infty$ .

*Proof.* The first part of the corollary is proved by induction on n. For n = 0, g is a nonzero constant in F, and the proposition is obviously true. Assume that the proposition is true for n - 1 for some  $n \ge 1$ . From (19.72) and the induction hypothesis, we see that

$$\Pr\{g(z_1, z_2, \cdots, z_n) \neq 0\}$$
(19.76)

$$= \Pr\{h(z_1, z_2, \cdots, z_{n-1}) \neq 0\} \Pr\{g(z_1, z_2, \cdots, z_n) \neq 0|$$
(19.77)

$$h(z_1, z_2, \cdots, z_{n-1}) \neq 0\}$$
(19.78)

$$\geq \left(1 - \frac{m}{|F|}\right)^{n-1} \Pr\{g(z_1, z_2, \cdots, z_n) \neq 0|$$
(19.79)

$$h(z_1, z_2, \cdots, z_{n-1}) \neq 0\}$$
(19.80)

$$\geq \left(1 - \frac{m}{|F|}\right)^n \left(1 - \frac{m}{|F|}\right) \tag{19.81}$$

$$= \left(1 - \frac{m}{|F|}\right)^n. \tag{19.82}$$

This proves the first part of the corollary. As n is fixed, the lower bound above tends to 1 as  $|F| \to \infty$ . This completes the proof.  $\Box$ 

*Example 19.19.* Recall the 2-dimensional linear network code in Example 19.9 that is expressed in the 12 indeterminates  $a, b, c, \dots, l$ . Place the vectors  $\mathbf{f}_{(t,w)}$  and  $\mathbf{f}_{(u,w)}$  in juxtaposition into the  $2 \times 2$  matrix

$$L_w = \begin{bmatrix} ae \ cg \\ be \ dg \end{bmatrix},\tag{19.83}$$

the vectors  $\mathbf{f}_{(t,y)}$  and  $\mathbf{f}_{(x,y)}$  into the 2 × 2 matrix

$$L_y = \begin{bmatrix} af \ aeik + cgjk\\ bf \ beik + dgjk \end{bmatrix},$$
(19.84)

and the vectors  $\mathbf{f}_{(u,z)}$  and  $\mathbf{f}_{(x,z)}$  into the 2 × 2 matrix

$$L_Z = \begin{bmatrix} aeil + cgjl \ ch\\ beil + dgjl \ dh \end{bmatrix}.$$
 (19.85)

19.4 Existence and Construction 445

Clearly,

$$\det(L_w) \cdot \det(L_y) \cdot \det(L_z) \neq 0 \in F[a, b, c, \cdots, l].$$
(19.86)

Applying Lemma 19.17 to the polynomial on the left hand side above, we can set scalar values for the 12 indeterminates so that it is evaluated to a nonzero value in F when F is sufficiently large. This implies that the determinants on the left hand side of (19.86) are evaluated to nonzero values in F simultaneously. Thus these scalar values yield a 2-dimensional linear multicast. In fact,

$$\det(L_w) \cdot \det(L_y) \cdot \det(L_z) = 1 \tag{19.87}$$

when

$$b = c = 0$$
 (19.88)

and

$$a = d = e = f = \dots = l = 1.$$
 (19.89)

Therefore, the 2-dimensional linear network code depicted in Figure 19.2 is a linear multicast, and this fact is regardless of the choice of the base field F.

**Theorem 19.20.** There exists an  $\omega$ -dimensional linear multicast on an acyclic network for sufficiently large base field F.

*Proof.* For a directed path  $P = e_1, e_2, \cdots, e_m$ , define

$$K_P = \prod_{1 \le j < m} k_{e_j, e_{j+1}}.$$
 (19.90)

Calculating by (19.17) recursively from the upstream channels to the downstream channels, it is not hard to find that

$$\mathbf{f}_e = \sum_{d \in \text{In}(s)} \left( \Sigma_{P: \text{ a path from } d \text{ to } e} \ K_P \right) \ \mathbf{f}_d \tag{19.91}$$

for every channel e (see Example 19.23 below). Denote by F[\*] the polynomial ring over the field F with all the  $k_{d,e}$  as indeterminates, where the total number of such indeterminates is equal to  $\sum_t |\text{In}(t)| \cdot |\text{Out}(t)|$ . Thus, every component of every global encoding kernel belongs to F[\*]. The subsequent arguments in this proof actually depend on this fact alone but not on the exact form of (19.91).

Let t be a non-source node with  $\maxflow(t) \geq \omega$ . Then there exist  $\omega$  edgedisjoint paths from the  $\omega$  imaginary channels to  $\omega$  distinct channels in  $\operatorname{In}(t)$ . Put the global encoding kernels of these  $\omega$  channels in juxtaposition to form an  $\omega \times \omega$  matrix  $L_t$ . We will prove that

$$\det(L_t) = 1 \tag{19.92}$$

for properly set scalar values of the indeterminates.

Toward proving this claim, we set

$$k_{d,e} = 1$$
 (19.93)

for all adjacent pairs of channels (d,e) along any one of the  $\omega$  edge-disjoint paths, and set

$$k_{d,e} = 0$$
 (19.94)

otherwise. With such local encoding kernels, the symbols sent on the  $\omega$  imaginary channels at source node s are routed to node t via the  $\omega$  edge-disjoint paths. Thus the columns in  $L_t$  are simply the global encoding kernels of the imaginary channels, which form the standard basis of the space  $F^{\omega}$ . Then (19.92) follows, and the claim is proved.

Consequently,

$$\det(L_t) \neq 0 \in F[*],\tag{19.95}$$

i.e.,  $det(L_t)$  is a nonzero polynomial in the indeterminates  $k_{d,e}$ . Since this conclusion applies to every non-source node t with maxflow $(t) \ge \omega$ ,

$$\prod_{\text{smaxflow}(t) \ge \omega} \det(L_t) \neq 0 \in F[*].$$
(19.96)

Applying Lemma 19.17 to the above polynomial when the field F is sufficiently large, we can set scalar values in F for the indeterminates so that

$$\prod_{t:\max flow(t) > \omega} \det(L_t) \neq 0 \in F,$$
(19.97)

which in turns implies that

t

$$\det(L_t) \neq 0 \in F \tag{19.98}$$

for all t such that  $maxflow(t) \ge \omega$ . These scalar values then yield a linear network code that meets the requirement (19.39) for a linear multicast.  $\Box$ 

**Corollary 19.21.** There exists an  $\omega$ -dimensional linear broadcast on an acyclic network for sufficiently large base field F.

*Proof.* For every non-source node t in the given acyclic network, install a new node t' and  $\omega$  input channels to this new node, with  $\min\{\omega, \max flow(t)\}$  of them from node t and the remaining  $\omega - \min\{\omega, \max flow(t)\}$  from source node s. This constructs a new acyclic network. Now consider an  $\omega$ -dimensional linear multicast on the new network whose existence follows from Theorem 19.20. For every node t' as described above,  $\dim(V_{t'}) = \omega$  because  $\max flow(t') = \omega$ . Moreover, since  $|\operatorname{In}(t')| = \omega$ , the global encoding kernels  $\mathbf{f}_e, e \in \operatorname{In}(t')$  are linearly independent. Therefore,

$$\dim(\langle \{\mathbf{f}_e : e \in \mathrm{In}(t') \cap \mathrm{Out}(t)\} \rangle) = |\mathrm{In}(t') \cap \mathrm{Out}(t)|$$
(19.99)

$$= \min\{\omega, \maxflow(t)\}.$$
(19.100)

By (19.17),

19.4 Existence and Construction 447

$$\langle \{ \mathbf{f}_e : e \in \mathrm{In}(t') \cap \mathrm{Out}(t) \} \rangle \subset V_t.$$
 (19.101)

Therefore,

$$\dim(V_t) \ge \dim(\langle \{ \mathbf{f}_e : e \in \mathrm{In}(t') \cap \mathrm{Out}(t) \} \rangle)$$
(19.102)

$$= \min\{\omega, \max flow(t)\}.$$
(19.103)

Then by invoking Theorem 19.10, we conclude that

$$\dim(V_t) = \min\{\omega, \maxflow(t)\}.$$
(19.104)

In other words, an  $\omega$ -dimensional linear multicast on the new network incorporates an  $\omega$ -dimensional linear broadcast on the original network.  $\Box$ 

**Corollary 19.22.** There exists an  $\omega$ -dimensional linear dispersion on an acyclic network for sufficiently large base field F.

*Proof.* For every nonempty collection  $\mathcal{T}$  of non-source nodes in the given acyclic network, install a new node  $u_{\mathcal{T}}$  and  $\max(t)$  channels from every node  $t \in \mathcal{T}$  to this new node. This constructs a new acyclic network with

$$\max flow(u_{\mathcal{T}}) = \max flow(\mathcal{T}) \tag{19.105}$$

for every  $\mathcal{T}$ . Now consider an  $\omega$ -dimensional linear broadcast on the new network whose existence follows from Corollary 19.21. By (19.17),

$$V_{u_{\mathcal{T}}} \subset V_{\mathcal{T}}.\tag{19.106}$$

Then

$$\dim(V_{\mathcal{T}}) \ge \dim(V_{u_{\mathcal{T}}}) \tag{19.107}$$

$$= \min\{\omega, \maxflow(u_{\mathcal{T}})\}$$
(19.108)

$$= \min\{\omega, \maxflow(\mathcal{T})\}.$$
 (19.109)

By invoking Theorem 19.10, we conclude that

$$\dim(V_{\mathcal{T}}) = \min\{\omega, \maxflow(\mathcal{T})\}.$$
(19.110)

In other words, an  $\omega$ -dimensional linear broadcast on the new network incorporates an  $\omega$ -dimensional linear dispersion on the original network.  $\Box$ 

*Example 19.23.* We now illustrate the formula (19.91) in the proof of Theorem 19.20 with the 2-dimensional linear network code in Example 19.9 which is expressed in the 12 indeterminates  $a, b, c, \dots, l$ . The local encoding kernels at the nodes are

$$K_s = \begin{bmatrix} a & c \\ b & d \end{bmatrix}, \quad K_t = \begin{bmatrix} e & f \end{bmatrix}, \quad K_u = \begin{bmatrix} g & h \end{bmatrix}, \quad (19.111)$$

$$K_w = \begin{bmatrix} i \\ j \end{bmatrix}, \ K_x = \begin{bmatrix} k \ l \end{bmatrix}.$$
(19.112)

Starting with  $\mathbf{f}_{(o,s)} = \begin{bmatrix} 1 & 0 \end{bmatrix}^{\top}$  and  $\mathbf{f}_{(o,s)'} = \begin{bmatrix} 0 & 1 \end{bmatrix}^{\top}$ , we can calculate the global encoding kernels by the formula (19.91). Take  $\mathbf{f}_{(x,y)}$  as the example. There are two paths from (o, s) to (x, y) and two from (o, s)' to (x, y). For these paths,

$$K_{P} = \begin{cases} aeik & \text{for } P = (o, s), (s, t), (t, w), (w, x), (x, y) \\ beik & \text{for } P = (o, s)', (s, t), (t, w), (w, x), (x, y) \\ cgjk & \text{for } P = (o, s), (s, u), (u, w), (w, x), (x, y) \\ dgjk & \text{for } P = (o, s)', (s, u), (u, w), (w, x), (x, y). \end{cases}$$
(19.113)

Thus

$$\mathbf{f}_{(x,y)} = (aeik)\mathbf{f}_{(o,s)} + (beik)\mathbf{f}_{(o,s)'} + (cgjk)\mathbf{f}_{(o,s)} + (dgjk)\mathbf{f}_{(o,s)'} \quad (19.114)$$

$$= \begin{bmatrix} aeik + cgjk\\ beik + dgjk \end{bmatrix},$$
(19.115)

which is consistent with Example 19.9.

The proof of Theorem 19.20 provides an algorithm for constructing a linear multicast that uses Lemma 19.17 as a subroutine to search for scalars  $a_1, a_2, \dots, a_n \in F$  such that  $g(a_1, a_2, \dots, a_n) \neq 0$  whenever  $g(z_1, z_2, \dots, z_n)$ is a nonzero polynomial over a sufficiently large field F. The straightforward implementation of this subroutine is exhaustive search, which is generally computationally inefficient. Nevertheless, the proof of Theorem 19.20 renders a simple method to construct a linear multicast randomly.

**Corollary 19.24.** Consider an  $\omega$ -dimensional linear network code on an acyclic network. By choosing the local encoding kernels  $k_{d,e}$  for all adjacent pairs of channels (d, e) independently according to the uniform distribution on the base field F, a linear multicast can be constructed with probability tends to 1 as  $|F| \to \infty$ .

*Proof.* This follows directly from Corollary 19.18 and the proof of Theorem 19.20.  $\Box$ 

The technique described in the above theorem for constructing a linear network code is called *random network coding*. Random network coding has the advantage that the code construction can be done independent of the network topology, making it very useful when the network topology is unknown. A case study for an application of random network coding will be presented in Section 19.7.

While random network coding offers a simple construction and more flexibility, a much larger base field is usually required. In some applications, it is necessary to verify that the code randomly constructed indeed possesses the desired properties. Such a task can be computationally non-trivial. The next algorithm constructs a linear multicast deterministically in polynomial time. Unlike the algorithm given in the proof of Theorem 19.20 that assigns values to the local encoding kernels, this algorithm assigns values to the global encoding kernels.

Algorithm 19.25 (Jaggi-Sanders Algorithm). This algorithm constructs an  $\omega$ -dimensional linear multicast over a finite field F on an acyclic network when  $|F| > \eta$ , the number of non-source nodes t in the network with maxflow $(t) \ge \omega$ . Denote these  $\eta$  non-source nodes by  $t_1, t_2, \dots, t_{\eta}$ .

A sequence of channels  $e_1, e_2, \dots, e_l$  is called a path leading to a node  $t_q$  if  $e_1 \in \text{In}(s), e_l \in \text{In}(t_q)$ , and  $(e_j, e_{j+1})$  is an adjacent pair for all  $1 \leq j \leq l-1$ . For each  $q, 1 \leq q \leq \eta$ , there exist  $\omega$  edge-disjoint paths  $P_{q,1}, P_{q,2}, \dots, P_{q,\omega}$  leading to  $t_q$ . All together there are  $\eta\omega$  such paths. The following procedure assigns a global encoding kernel  $\mathbf{f}_e$  for every channel e in the network in an upstream-to-downstream order such that  $\dim(V_{t_q}) = \omega$  for  $1 \leq q \leq \eta$ .

{

// By definition, the global encoding kernels of the  $\omega$ // imaginary channels form the standard basis of  $F^{\omega}$ . for  $(q = 1; q \le \eta; q + +)$ for  $(i = 1; i \leq \omega; i + +)$  $e_{q,i}$  = the imaginary channel initiating path  $P_{q,i}$ ; // This initializes  $e_{q,i}$ . Subsequently,  $e_{q,i}$  will be // dynamically updated by moving down path  $P_{q,i}$ // until it finally becomes a channel in  $In(t_q)$ . for (every node t, in any upstream-to-downstream order) for (every channel  $e \in \operatorname{Out}(t)$ ) // With respect to this channel e, define a "pair" as a // pair (q, i) of indices such that channel e is on the // path  $P_{q,i}$ . Note that for each q, there exists at most // one pair (q, i). Thus the number of pairs is at least 0 // and at most  $\eta$ . Since the nodes t are chosen in // an upstream-to-downstream order, if (q, i) is a pair, // then  $e_{q,i} \in \text{In}(t)$  by induction, so that  $\mathbf{f}_{e_{q,i}} \in V_t$ . For // reasons to be explained in the algorithm verification  $\begin{array}{l} // \text{ below, } \mathbf{f}_{e_{q,i}} \not\in \langle \{\mathbf{f}_{e_{q,j}} : j \neq i\} \rangle, \text{ and therefore} \\ // | \mathbf{f}_{e_{q,i}} \in V_t \setminus \langle \{\mathbf{f}_{e_{q,j}} : j \neq i\} \rangle. \end{array}$ Choose a vector w in  $V_t$  such that  $w \notin \langle \{f_{e_{q,j}} : j \neq i\} \rangle$  for every pair (q, i); // To see the existence of such a vector w, let  $//\dim(V_t) = \nu$ . Then,  $\dim(V_t \cap \langle \{f_{e_{q,j}} : j \neq i\} \rangle) \leq$  $// \nu - 1$  for every pair (q, i) since  $// \mathbf{f}_{e_{q,i}} \in V_t \setminus \langle \{f_{e_{q,j}} : j \neq i\} \rangle$ . Thus

 $\begin{array}{c} // |V_t \cap (\cup_{(q,i):\ a \ pair} \langle \{f_{e_{q,j}} : j \neq i\} \rangle)| \\ // \leq \eta |F|^{\nu-1} < |F|^{\nu} = |V_t|. \\ \mathbf{f}_e = w; \\ // \text{ This is equivalent to choosing scalar values for local} \\ // \text{ encoding kernels } k_{d,e} \text{ for all } d \in \text{In}(t) \text{ such that} \\ // \sum_{d \in \text{In}(t)} k_{d,e} \mathbf{f}_d \notin \langle \{f_{e_{q,j}} : j \neq i\} \rangle \text{ for every pair } (q,i). \\ for (every \ pair \ (q,i)) \\ e_{q,i} = e; \\ \end{array} \right\}$ 

Algorithm Verification. For  $1 \leq q \leq \eta$  and  $1 \leq i \leq \omega$ , the channel  $e_{q,i}$  is on the path  $P_{q,i}$ . Initially  $e_{q,i}$  is an imaginary channel at source node s. Through dynamic updating, it moves downstream along the path until finally reaching a channel in  $\text{In}(t_q)$ .

Fix an index q, where  $1 \leq q \leq \eta$ . Initially, the vectors  $\mathbf{f}_{e_{q,1}}, \mathbf{f}_{e_{q,2}}, \cdots, \mathbf{f}_{e_{q,\omega}}$  are linearly independent because they form the standard basis of  $F^{\omega}$ . At the end, they need to span the vector space  $F^{\omega}$ . Therefore, in order for the eventually constructed linear network code to qualify as a linear multicast, it suffices to show the preservation of the linear independence among  $\mathbf{f}_{e_{q,1}}, \mathbf{f}_{e_{q,2}}, \cdots, \mathbf{f}_{e_{q,\omega}}$  throughout the algorithm.

We need to show the preservation in the generic step inside the "for loop" for each channel e in the algorithm. The algorithm defines a "pair" as a pair (q, i) of indices such that channel e is on path  $P_{q,i}$ . When no (q, i) is a pair for  $1 \leq i \leq \omega$ , the channels  $e_{q,1}, e_{q,2}, \dots, e_{q,\omega}$  are not changed in the generic step; neither are the vectors  $\mathbf{f}_{e_{q,1}}, \mathbf{f}_{e_{q,2}}, \dots, \mathbf{f}_{e_{q,\omega}}$ . So we only need to consider the scenario that a pair (q, i) exists for some i. The only change among the channels  $e_{q,1}, e_{q,2}, \dots, e_{q,\omega}$  is that  $e_{q,i}$  becomes e. Meanwhile, the only change among the vectors  $\mathbf{f}_{e_{q,1}}, \mathbf{f}_{e_{q,2}}, \dots, \mathbf{f}_{e_{q,\omega}}$  is that  $\mathbf{f}_{e_{q,i}}$  becomes a vector

$$w \notin \langle \{ \mathbf{f}_{e_{q,j}} : j \neq i \} \rangle. \tag{19.116}$$

This preserves the linear independence among  $\mathbf{f}_{e_{q,1}}, \mathbf{f}_{e_{q,2}}, \cdots, \mathbf{f}_{e_{q,\omega}}$  as desired.

Complexity Analysis. There are a total of  $|\mathcal{E}|$  channels in the network. In the algorithm, the generic step in the "for loop" for each channel *e* processes at most  $\eta$  pairs. Throughout the algorithm, at most  $|\mathcal{E}|\eta$  such collections of channels are processed. From this, it is not hard to implement the algorithm within a polynomial time in  $|\mathcal{E}|$  for a fixed  $\omega$ . The computational details can be found in [165].

**Remark 1** In the Jaggi-Sanders algorithm, all nodes t in the network with  $maxflow(t) \ge \omega$  serve as a sink node that receives the message **x**. The algorithm can easily be modified accordingly if only a subset of such nodes need

to serve as a sink node. In that case, the field size requirement is  $|F| > \eta'$ , where  $\eta'$  is the total number of sink nodes.

**Remark 2** It is not difficult to see from the lower bound on the required field size in the Jaggi-Sanders algorithm that if a field much larger than sufficient is used, then a linear multicast can be constructed with high probability by randomly choosing the global encoding kernels.

Example 19.26 (Multi-Source Multicast). Consider a network coding problem on an acyclic network  $\mathcal{G}$  with a set S of source nodes. At node  $s \in S$ , a message  $\mathbf{x}_s$  in the form of a row vector in  $F^{\omega_s}$  is generated. Let

$$\omega = \sum_{s \in S} \omega_s \tag{19.117}$$

be the total dimension of all the messages, and let

$$\mathbf{x} = (\mathbf{x}_s : s \in S) \tag{19.118}$$

be referred to as the message. Here, we do not impose the constraint that a node  $s \in S$  has no input channels.

Expand the network  $\mathcal{G}$  into a network  $\mathcal{G}'$  by installing a new node 0, and  $\omega_s$  channels from node 0 to node s for each  $s \in S$ . Denote the value of a max-flow from node 0 to node t in  $\mathcal{G}'$  by maxflow $_{\mathcal{G}'}(t)$ .

Suppose there exists a coding scheme on  $\mathcal{G}$  such that a node t can decode the message **x**. Such a coding scheme induces a coding scheme on  $\mathcal{G}'$  for which

- 1. the message  $\mathbf{x}$  is generated at node 0;
- 2. for all  $s \in S$ , the message  $\mathbf{x}_s$  is sent uncoded from node 0 to node s through the  $\omega_s$  channels from node 0 to node s.

Applying the max-flow bound to node t with respect to this coding scheme on  $\mathcal{G}'$ , we obtain

$$\max flow_{\mathcal{G}'}(t) \ge \omega. \tag{19.119}$$

Thus we have shown that if a node t in  $\mathcal{G}$  can decode the message  $\mathbf{x}$ , then (19.119) has to be satisfied.

We now show that for a sufficiently large base field F, there exists a coding scheme on  $\mathcal{G}$  such that a node t satisfying (19.119) can decode the message  $\mathbf{x}$ . Let  $\eta$  be the number of nodes in  $\mathcal{G}$  that satisfies (19.119). To avoid triviality, assume  $\eta \geq 1$ . By Theorem 19.20, there exists an  $\omega$ -dimensional linear multicast  $\mathcal{C}$  on  $\mathcal{G}'$  when the base field is sufficiently large. From the proof of Theorem 19.10, we see that for this linear multicast, the  $\omega \times \omega$  matrix  $F_{\text{Out}(0)}$ must be invertible, otherwise a node t satisfying (19.119) cannot possibly decode the message  $\mathbf{x}$ . Transforming  $\mathcal{C}$  by the matrix  $[F_{\text{Out}(0)}]^{-1}$ , we obtain from (19.54) a linear multicast  $\mathcal{C}'$  with

$$F'_{\text{Out}(0)} = \left[F_{\text{Out}(0)}\right]^{-1} F_{\text{Out}(0)} = I_{\omega}.$$
 (19.120)

Accordingly, for this linear multicast, the message  $\mathbf{x}_s$  is sent uncoded from node 0 to node s for all  $s \in S$ . Thus a coding scheme on  $\mathcal{G}$  with the message  $\mathbf{x}_s$ being generated at node s for all  $s \in S$  instead of being received from node 0 is naturally induced, and this coding scheme inherits from the linear multicast  $\mathcal{C}'$  that a node t satisfying (19.119) can decode the message  $\mathbf{x}$ .

Therefore, instead of tackling the multi-source multicast problem on  $\mathcal{G}$ , we can tackle the single-source multicast problem on  $\mathcal{G}'$ . This has already seen illustrated in Examples 17.1 and 17.2 for the butterfly network.

### 19.5 Generic Network Codes

In the last section, we have seen how to construct a linear multicast by the Jaggi-Sanders algorithm. In light of Corollaries 19.21 and 19.22, the same algorithm can be used for constructing a linear broadcast or a linear dispersion.

It is not difficult to see that if the Jaggi-Sanders algorithm is used for constructing a linear broadcast, then the computational complexity of the algorithm remains polynomial in  $|\mathcal{E}|$ , the total number of channels in the network. However, if the algorithm is used for constructing a linear dispersion, the computational complexity becomes exponential because the number of channels that need to be installed in constructing the new network in Corollary 19.22 grows exponentially with the number of channels in the original network.

In this section, we introduce a class of linear network codes called *generic network codes*. As we will see, if a linear network code is generic, then it is a linear dispersion, and hence also a linear broadcast and a linear multicast. Toward the end of the section, we will present a polynomial time algorithm that constructs a generic network code.

Imagine that in an  $\omega$ -dimensional linear network code, the base field F is replaced by the real field  $\mathbf{R}$ . Then arbitrary infinitesimal perturbation of the local encoding kernels would place the global encoding kernels at general positions with respect to one another in the space  $\mathbf{R}^{\omega}$ . General positions of the global encoding kernels maximize the dimensions of various linear spans by avoiding linear dependence in every conceivable way. The concepts of general positions and infinitesimal perturbation do not apply to the vector space  $F^{\omega}$  when F is a finite field. However, they can be emulated when F is sufficiently large with the effect of avoiding unnecessary linear dependence.

The following definitions of a generic network code captures the notion of placing the global encoding kernels in general positions. In the sequel, for a channel  $e_j \in \mathcal{E}$ , let  $e_j \in \text{Out}(t_j)$ , and for a collection of channels  $\xi = \{e_1, e_2, \dots, e_{|\xi|}\} \subset \mathcal{E}$ , let  $\xi_{\bar{k}} = \xi \setminus \{e_k\}$ .

**Definition 19.27 (Generic Network Code I).** An  $\omega$ -dimensional linear network code on an acyclic network is generic if for any  $m \geq 1$  and any

collection of channels  $\xi = \{e_1, e_2, \cdots, e_m\} \subset \mathcal{E}$ , if there exists  $1 \leq k \leq m$  such that

a) there is no directed path from  $t_k$  to  $t_j$  for  $j \neq k$ , b)  $V_{t_k} \not\subset V_{\xi_{\overline{k}}}$ ,

then  $\mathbf{f}_{e_k} \notin V_{\xi_{\bar{k}}}$ .

**Definition 19.28 (Generic Network Code II).** An  $\omega$ -dimensional linear network code on an acyclic network is generic if for any  $m \geq 1$  and any collection of channels  $\xi = \{e_1, e_2, \dots, e_m\} \subset \mathcal{E}$ , if there exists  $1 \leq k \leq m$  such that

a) there is no directed path from  $t_k$  to  $t_j$  for  $j \neq k$ , b)  $V_{t_k} \not\subset V_{\xi_{\bar{k}}}$ , c)  $\mathbf{f}_e, e \in \xi_{\bar{k}}$  are linearly independent<sup>3</sup>,

then  $\mathbf{f}_{e_k} \not\in V_{\xi_{\bar{k}}}$ .

In Definitions 19.27 and 19.28, if a) does not hold, then  $\mathbf{f}_{e_k} \notin V_{\xi_k}$  may not be possible at all as we now explain. Let  $\xi = \{e_1, e_2\}$  and k = 1. Suppose  $\operatorname{In}(t_2) = \{e_1\}$  so that a) is violated. Since node  $t_2$  has only  $e_1$  as the input channel,  $\mathbf{f}_{e_1}$  cannot possibly be linear independent of  $\mathbf{f}_{e_2}$ .

The only difference between Definitions 19.27 and 19.28 is the additional requirement c) in the latter. The equivalence between these two definitions of a generic network code can be seen as follows. It is obvious that if a linear network code satisfies Definition 19.27, then it also satisfies Definition 19.28. Conversely, suppose a linear network code satisfies Definition 19.28. Consider any collection of channels  $\xi = \{e_1, e_2, \dots, e_m\} \subset \mathcal{E}$  such that there exists  $1 \leq k \leq m$  satisfying a) and b) in Definition 19.27 but not necessarily c) in Definition 19.28. Then we can always find a subset  $\xi'_{\bar{k}}$  of  $\xi_{\bar{k}}$  such that  $\mathbf{f}_e, e \in \xi'_{\bar{k}}$  are linearly independent and  $V_{\xi'_{\bar{k}}} = V_{\xi_{\bar{k}}}$ . Upon letting  $\xi' = \{e_k\} \cup \xi'_{\bar{k}}$  and applying Definition 19.28 with  $\xi'$  in place of  $\xi$ , we have

$$\mathbf{f}_{e_k} \not\in V_{\xi_r'} = V_{\xi_{\bar{k}}},\tag{19.121}$$

so the network code also satisfies Definition 19.27. This shows that the two definitions of a generic network code are equivalent. Note that in Definition 19.28, if  $\xi$  satisfies all the prescribed conditions, then  $m \leq \omega$  because c) and  $\mathbf{f}_{e_k} \notin V_{\xi_{\bar{k}}}$  together imply that  $\mathbf{f}_e, e \in \xi$  are linearly independent. Definition 19.28, which has a slightly more complicated form compared with Definition 19.27, will be instrumental in the proof of Theorem 19.32 that establishes different characterizations of a generic network code.

 $<sup>^3</sup>$  We adopt the convention that an empty collection of vectors is linearly independent.

**Proposition 19.29.** For a generic network code, for any collection of m output channels at a node t, where  $1 \le m \le \dim(V_t)$ , the corresponding global encoding kernels are linearly independent.

*Proof.* Since the proposition becomes degenerate if  $\dim(V_t) = 0$ , we assume  $\dim(V_t) > 0$ . In Definition 19.27, let all the nodes  $t_j$  be equal to node t and let  $1 \leq m \leq \dim(V_t)$ . First note that there is no directed path from node t to itself because the network is acyclic. For m = 1,  $\xi_{\bar{1}} = \emptyset$  and  $V_{\xi_{\bar{1}}} = V_{\emptyset} = \emptyset$ . Since  $\dim(V_t) > 0$ , we have  $V_t \not\subset V_{\xi_{\bar{1}}}$ . Then  $\mathbf{f}_{e_1} \notin V_{\xi_{\bar{1}}}$ , which implies  $\mathbf{f}_{e_1} \neq 0$ . This proves the proposition for m = 1.

Assume that the proposition is true for m-1 some  $2 \le m \le \dim(V_t)$ . We now prove that the proposition is true for m. By the induction hypothesis,  $\mathbf{f}_{e_1}, \mathbf{f}_{e_2}, \dots, \mathbf{f}_{e_{m-1}}$  are linearly independent. Since

$$\dim(\langle \{\mathbf{f}_{e_1}, \mathbf{f}_{e_2}, \cdots, \mathbf{f}_{e_{m-1}}\}\rangle) = m - 1 < \dim(V_t), \tag{19.122}$$

we have

$$V_t \not\subset \langle \{\mathbf{f}_{e_1}, \mathbf{f}_{e_2}, \cdots, \mathbf{f}_{e_{m-1}} \} \rangle.$$
(19.123)

Then by Definition 19.27,

$$\mathbf{f}_{e_m} \not\in \langle \{\mathbf{f}_{e_1}, \mathbf{f}_{e_2}, \cdots, \mathbf{f}_{e_{m-1}}\} \rangle.$$
(19.124)

Hence,  $\mathbf{f}_{e_1}, \mathbf{f}_{e_2}, \cdots, \mathbf{f}_{e_m}$  are linearly independent. The proposition is proved.

**Corollary 19.30.** For a generic network code, if  $|\text{Out}(t)| \leq \dim(V_t)$  for a node t, then the global encoding kernels of all the output channels of t are linearly independent.

A linear dispersion on an acyclic network is not necessarily a generic network code. The following is a counterexample.

Example 19.31. The 2-dimensional linear dispersion on the network in Figure 19.6 is not generic because the global encoding kernels of two of the output channels from source node s are equal to  $[1 \ 1]^{\mathsf{T}}$ , a contradiction to Proposition 19.29. It can be shown, however, that a generic network code on an acyclic network  $\mathcal{G}$  can be constructed through a linear dispersion on an expanded network  $\mathcal{G}'$ . See Problem 12 for details.

Together with Example 19.13, the example above shows that the four classes of linear network codes we have discussed, namely linear multicast, linear broadcast, linear dispersion, and generic network code, achieve the max-flow bound to strictly increasing extents.

In the following theorem, we prove two characterizations of a generic network code, each can be regarded as an alternative definition of a generic network code. The reader should understand this theorem before proceeding further but may skip the proof at the first reading.

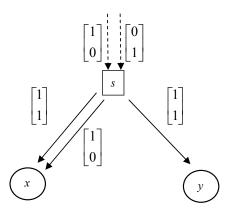


Fig. 19.6. A 2-dimensional linear dispersion that is not a generic network code.

**Theorem 19.32.** For an  $\omega$ -dimensional linear network code on an acyclic network, the following conditions are equivalent:

- 1) The network code is generic.
- 2) For any  $m \ge 1$  and any collection of channels  $\xi = \{e_1, e_2, \cdots, e_m\} \subset \mathcal{E}$ ,
- if  $V_{t_j} \not\subset V_{\xi_{\bar{j}}}$  for all  $1 \leq j \leq m$ , then  $\mathbf{f}_e, e \in \xi$  are linearly independent.
- 3) For any nonempty collection of channels  $\xi \subset \mathcal{E}$ , if

$$|\xi| = \min\{\omega, \maxflow(\xi)\}, \qquad (19.125)$$

then  $\mathbf{f}_e, e \in \xi$  are linearly independent.

*Proof.* Throughout this proof, we use Definition 19.28 as the definition of a generic network code. We will prove the theorem by showing that  $1) \Rightarrow 2) \Rightarrow 3) \Rightarrow 1$ .

We first show that 1)  $\Rightarrow$  2). Assume 1) holds. Consider any  $m \ge 1$  and any collection of channels  $\xi = \{e_1, e_2, \dots, e_m\} \subset \mathcal{E}$ , and assume  $V_{t_j} \not\subset V_{\xi_j}$  for all  $1 \le j \le m$ . We will show by induction on m that  $\mathbf{f}_e, e \in \xi$  are linearly independent. The claim is trivially true for m = 1. Assume the claim is true for m - 1 for some  $2 \le m \le \omega$ , and we will show that it is true for m.

Consider  $\xi = \{e_1, e_2, \dots, e_m\}$  and assume  $V_{t_j} \not\subset V_{\xi_j}$  for all  $1 \leq j \leq m$ . We first prove by contradiction that there exists at least one k such that there is no directed path from  $t_k$  to  $t_j$  for all  $j \neq k$ , where  $1 \leq j, k \leq m$ . Assume that for all k, there is at least one directed path from node  $t_k$  to node  $t_j$  for some  $j \neq k$ . Starting at any node  $t_k$ , by traversing such directed paths, we see that there exists a directed cycle in the network because the set  $\{t_k : 1 \leq k \leq m\}$  is finite. This leads to a contradiction because the network is acyclic, proving the existence of k as prescribed. Then apply Definition 19.28 to see that

$$\mathbf{f}_{e_k} \notin V_{\xi_{\bar{k}}} = \langle \{ \mathbf{f}_e : e \in \xi_{\bar{k}} \} \rangle.$$
(19.126)

Now for any  $j \neq k$ , since  $V_{t_j} \not\subset V_{\xi_{\bar{i}}}$  and  $V_{\xi_{\bar{k}} \setminus \{e_j\}} = V_{\xi_{\bar{i}} \setminus \{e_k\}} \subset V_{\xi_{\bar{i}}}$ , we have

$$V_{t_j} \not\subset V_{\xi_{\bar{k}} \setminus \{e_j\}}.\tag{19.127}$$

Then apply the induction hypothesis to  $\xi_{\bar{k}}$  to see that  $\mathbf{f}_e, e \in \xi_{\bar{k}}$  are linearly independent. It then follows from (19.126) that  $\mathbf{f}_e, e \in \xi$  are linearly independent. Thus  $1) \Rightarrow 2$ ).

We now show that 2)  $\Rightarrow$  3). Assume 2) holds and consider any collection of channel  $\xi = \{e_1, e_2, \dots, e_m\} \subset \mathcal{E}$  satisfying (19.125), where  $m \ge 1$ . Then

$$m = |\xi| = \min\{\omega, \maxflow(\xi)\}, \qquad (19.128)$$

which implies

$$\max flow(\xi) \ge m. \tag{19.129}$$

Therefore, there exist m edge-disjoint paths  $P_1, P_2, \dots, P_m$  from source node s to the channels in  $\xi$ , where the last channel on path  $P_j$  is  $e_j$ .

Denote the length of  $P_i$  by  $l_i$  and let

$$L = \sum_{j=1}^{m} l_j \tag{19.130}$$

be the total length of all the paths. We will prove the claim that  $\mathbf{f}_{e_1}, \mathbf{f}_{e_2}, \cdots, \mathbf{f}_{e_m}$ are linearly independent by induction on L. For the base case L = m, since  $m \leq \omega$  by (19.128), the claim is true by Proposition 19.29 with t = s. Assume that the claim is true for L - 1 for some  $L \geq m + 1$ , and we will prove that it is true for L. Let  $A = \{j : l_j > 1\}$  and for  $j \in A$ , let  $\xi'_j = \{e_1, e_2, \cdots, e_{j-1}, e'_j, e_{j+1}, \cdots, e_m\}$ , where  $e'_j$  is the channel preceding  $e_j$ on  $P_j$ . Then by the induction hypothesis,  $\mathbf{f}_e, e \in \xi'_j$  are linearly independent, which implies that

$$V_{t_i} \not\subset V_{\xi_{\overline{i}}}.\tag{19.131}$$

For  $j \notin A$ ,  $l_j = 1$ , i.e.,  $t_j = s$ . It follows from (19.128) that  $m \leq \omega$ . Then

$$V_{t_j} = V_s \not\subset V_{\xi_{\bar{j}}} \tag{19.132}$$

because dim $(V_{\xi_{\bar{j}}}) \leq |\xi_{\bar{j}}| = m - 1 < m \leq \omega$ . Therefore, (19.131) holds for all j, and hence by 2),  $\mathbf{f}_e, e \in \xi$  are linearly independent. Thus 2)  $\Rightarrow$  3).

Finally, we show that 3)  $\Rightarrow$  1). Assume 3) holds and consider any collection of channels  $\xi = \{e_1, e_2, \dots, e_m\} \subset \mathcal{E}$ , where  $1 \leq m \leq \omega$ , such that a) to c) in Definition 19.28 hold for some  $1 \leq k \leq m$ . Then either  $t_j = s$  for all  $1 \leq j \leq m$ , or  $t_k \neq s$ , because otherwise a) in Definition 19.28 is violated.

If  $t_j = s$  for all  $1 \le j \le m$ , then

$$m = |\xi| = \max \text{flow}(\xi). \tag{19.133}$$

Since  $m \leq \omega$ , we have

19.5 Generic Network Codes 457

$$|\xi| = \min\{\omega, \maxflow(\xi)\}.$$
(19.134)

Then  $\mathbf{f}_e, e \in \xi$  are linearly independently by 3), proving that  $\mathbf{f}_{e_k} \notin V_{\xi_{\bar{k}}}$ .

Otherwise,  $t_k \neq s$ . Following b) in Definition 19.28, there exists  $e'_k \in$ In $(t_k) \subset \mathcal{E}$  such that  $\mathbf{f}_{e'_k}$  and  $\mathbf{f}_e, e \in \xi_{\bar{k}}$  are linearly independent. Let  $\xi'_k = \{e_1, e_2, \dots, e_{k-1}, e'_k, e_{k+1}, \dots, e_m\}$ . By Corollary 19.11,

$$\max flow(\xi'_k) \ge \dim(V_{\xi'_k}) = m, \tag{19.135}$$

so  $e_1, e_2, \dots, e_{k-1}, e'_k, e_{k+1}, e_m$  can be traced back to source node *s* via some edge-disjoint paths  $P_1, P_2, \dots, P_{k-1}, P'_k, P_{k+1}, \dots, P_m$ , respectively. Let  $P_k$  be obtained by appending  $e_k$  to  $P'_k$ . Since there is no directed path from  $t_k$  to  $t_j$  and  $e_k \neq e_j$  for all  $j \neq k, P_1, P_2, \dots, P_{k-1}, P_k, P_{k+1}, \dots, P_m$  are edge-disjoint. Therefore,

$$\max flow(\xi) \ge m. \tag{19.136}$$

On the other hand,

$$\max \text{flow}(\xi) \le |\xi| = m. \tag{19.137}$$

Therefore,

$$m = |\xi| = \max \text{flow}(\xi), \tag{19.138}$$

i.e., (19.133). As before, we can further obtain (19.134). Then by 3),  $\mathbf{f}_e, e \in \xi$ , are linearly independent, and therefore  $\mathbf{f}_{e_k} \notin V_{\xi_{\bar{k}}}$ . Thus 3)  $\Rightarrow$  1).

Hence, the theorem is proved.  $\Box$ 

**Corollary 19.33.** An  $\omega$ -dimensional generic network code on an acyclic network is an  $\omega$ -dimensional linear dispersion on the same network.

*Proof.* Consider an  $\omega$ -dimensional generic network code on an acyclic network and let  $\mathcal{T}$  be any collection of non-source nodes. Let

$$m = \min\{\omega, \max \text{flow}(\mathcal{T})\}.$$
(19.139)

Since maxflow( $\mathcal{T}$ )  $\geq m$ , there exists m edge-disjoint paths  $P_1, P_2, \dots, P_m$  from source node s to  $\mathcal{T}$ . Let  $e_i$  be the last channel on path  $P_i$ , and let

$$\xi = \{e_1, e_2, \cdots, e_m\}.$$
(19.140)

Evidently,

$$\max flow(\xi) = m. \tag{19.141}$$

It follows from (19.139) that  $m \leq \omega$ . Therefore,

$$|\xi| = m = \max \text{flow}(\xi) = \min\{\omega, \max \text{flow}(\xi)\}.$$
(19.142)

By Theorem 19.32,  $\mathbf{f}_e, e \in \boldsymbol{\xi}$  are linearly independent. Then

$$\dim(V_{\mathcal{T}}) \ge \dim(V_{\xi}) = m = \min\{\omega, \maxflow(\mathcal{T})\}.$$
(19.143)

By Theorem 19.10, we conclude that

$$\dim(V_{\mathcal{T}}) = \min\{\omega, \maxflow(\mathcal{T})\}.$$
(19.144)

Hence, we have shown that a generic network code is a linear dispersion.  $\Box$ 

Theorem 19.32 renders the following important interpretation of a generic network code. Consider any linear network code and any collection of channels  $\xi \subset \mathcal{E}$ . If  $\mathbf{f}_e, e \in \xi$  are linearly independent, then

$$|\xi| = \dim(V_{\xi}). \tag{19.145}$$

By Corollary 19.11,

$$\dim(V_{\xi}) \le \min\{\omega, \maxflow(\xi)\}.$$
(19.146)

Therefore,

$$|\xi| \le \min\{\omega, \maxflow(\xi)\}.$$
(19.147)

On the other hand,

$$\max flow(\xi) \le |\xi|, \tag{19.148}$$

which implies

$$\min\{\omega, \maxflow(\xi)\} \le |\xi|. \tag{19.149}$$

Combining (19.147) and (19.149), we see that

$$|\xi| = \min\{\omega, \maxflow(\xi)\}$$
(19.150)

is a necessary condition for  $\mathbf{f}_e, e \in \xi$  to be linearly independent. For a generic network code, this is also a sufficient condition for  $\mathbf{f}_e, e \in \xi$  to be linearly independent. Thus for a generic network code, if a set of global encoding kernels can possibly be linearly independent, then it is linear independent. In this sense, a generic network code captures the notion of placing the global encoding kernels in general positions.

The condition 2) in Theorem 19.32 is the original definition of a generic network code given in [205]. Unlike 1) and 3), this condition is purely algebraic and does not depend upon the network topology. However, it does not suggest an algorithm for constructing such a code.

Motivated by Definition 19.28, we now present an algorithm for constructing a generic network code. The computational complexity of this algorithm is polynomial in  $|\mathcal{E}|$ , the total number of channels in the network.

Algorithm 19.34 (Construction of a Generic Network Code). This algorithm constructs an  $\omega$ -dimensional generic network code over a finite field F with  $|F| > \sum_{m=1}^{\omega} {|\mathcal{E}|-1 \choose m-1}$  by prescribing global encoding kernels that constitute a generic network code.

{ for (every node t, following an upstream-to-downstream order) ł for (every channel  $e \in Out(t)$ ) Choose a vector w in  $V_t$  such that  $w \notin V_{\zeta}$ , where  $\zeta$  is any collection of m-1 already processed channels, where  $1 \leq m \leq \omega$ , such that  $\mathbf{f}_e, e \in \zeta$  are linearly independent and  $V_t \not\subset V_{\zeta}$ ; // To see the existence of such a vector w, denote dim $(V_t)$ // by  $\nu$ . If  $\zeta$  is any collection of m-1 channels with  $V_t \not\subset V_{\zeta}$ , // then dim $(V_t \cap V_{\zeta}) \leq \nu - 1$ . There are at most  $\sum_{m=1}^{\omega} {|\mathcal{E}|^{-1} \choose m-1}$ // such collections  $\zeta$ . Thus //  $|V_t \cap (\cup_{\zeta} V_{\zeta})| \leq \sum_{m=1}^{\omega} {|\mathcal{E}|^{-1} \choose m-1} |F|^{\nu-1} < |F|^{\nu} = |V_t|.$  $\mathbf{f}_e = w;$ // This is equivalent to choosing scalar values for the local // encoding kernels  $k_{d,e}$  for all d such that  $\sum_{d\in In(t)} k_{d,e} \mathbf{f}_d$  $// \notin V_{\zeta}$  for every collection  $\zeta$  of channels as prescribed. } } }

Algorithm Verification. We will verify that the code constructed is indeed generic by way of Condition 3) in Theorem 19.32. Consider any collection of channels  $\xi = \{e_1, e_2, \dots, e_m\} \subset \mathcal{E}$  satisfying (19.125), where  $1 \leq m \leq \omega$ . Then there exist m edge-disjoint paths  $P_1, P_2, \dots, P_m$  from source node s to the channels in  $\xi$ , where the last channel on path  $P_j$  is  $e_j$ . Denote the length of  $P_j$  by  $l_j$  and let

$$L = \sum_{j=1}^{m} l_j \tag{19.151}$$

be the total length of all the paths. We will prove the claim that  $\mathbf{f}_e, e \in \xi$  are linearly independent by induction on L.

It is easy to verify that for any set of m channels in  $\operatorname{Out}(s)$ , the global encoding kernels assigned are linearly independent, so the base case L = mis verified. Assume the claim is true for L - 1 for some  $L \ge m + 1$ , and we will prove that it is true for L. Let  $e_k$  be the channel whose global encoding kernel is last assigned among all the channels in  $\xi$ . Note that  $P_k \ge 2$  since  $L \ge m + 1$  and the global encoding kernels are assigned by the algorithm in an upstream-to-downstream order. Then let  $e'_k$  be the channel preceding  $e_k$ on  $P_k$ , and let

$$\xi' = \{e_1, e_2, \cdots, e_{k-1}, e'_k, e_{k+1}, \cdots, e_m\}.$$
(19.152)

By the induction hypothesis,  $\mathbf{f}_{e}, e \in \xi'$  are linearly independent. Since  $\mathbf{f}_{e'_{k}}$  is linearly independent of  $\mathbf{f}_{e}$  for  $e \in \xi' \setminus \{e'_{k}\} = \xi_{\bar{k}}, V_{t_{k}} \not\subset V_{\xi_{\bar{k}}}$ . It then follows from the construction that  $\mathbf{f}_{e_{k}} \not\in V_{\xi_{\bar{k}}}$  because  $\xi_{\bar{k}}$  is one of the collections  $\zeta$ 

considered when  $\mathbf{f}_{e_k}$  is assigned. Hence,  $\mathbf{f}_e, e \in \xi$  are linearly independent, verifying that the network code constructed is generic.

Complexity Analysis. In the algorithm, the "for loop" for each channel e processes at most  $\sum_{m=1}^{\omega} \binom{|\mathcal{E}|-1}{m-1}$  collections of m-1 channels. The processing includes the detection of those collections  $\zeta$  as well as the computation of the set  $V_t \setminus (\bigcup_{\zeta} V_{\zeta})$ . This can be done, for instance, by Gauss elimination. Throughout the algorithm, the total number of collections of channels processed is at most  $|\mathcal{E}| \sum_{m=1}^{\omega} \binom{|\mathcal{E}|-1}{m-1}$ , a polynomial in  $|\mathcal{E}|$  of degree  $\omega$ . Thus for a fixed  $\omega$ , it is not hard to implement the algorithm within a polynomial time in  $|\mathcal{E}|$ . This is similar to the polynomial-time implementation of the Jaggi-Sanders algorithm for constructing a linear multicast.

Algorithm 19.34 constitutes a constructive proof for the next theorem.

**Theorem 19.35.** There exists an  $\omega$ -dimensional generic network code on an acyclic network for sufficiently large base field F.

By noting the lower bound on the required field size in Algorithm 19.34, a generic network code can be constructed with high probability by randomly choosing the global encoding kernels provided that the base field is much larger than sufficient.

### 19.6 Static Network Codes

In the discussion so far, a linear network code has been defined on a network with a fixed topology, where all the channels are assumed to be available at all times. In a real network, however, a channel may fail due to various reasons, for example, hardware failure, cable cut, or natural disasters. With the failure of some subset of channels, the communication capacity of the resulting network is generally reduced.

Consider the use of, for instance, an  $\omega$ -dimensional multicast on an acyclic network for multicasting a sequence of messages generated at the source node. When no channel failure occurs, a non-source node with the value of a maxflow at least equal to  $\omega$  would be able to receive the sequence of messages. In case of channel failures, if the value of a max-flow of that node in the resulting network is at least  $\omega$ , the sequence of messages in principle can still be received at that node. However, this would involve the deployment of a network code for the new network topology, which not only is cumbersome but also may cause a significant loss of data during the switchover.

In this section, we discuss a class of linear network codes called *static* network codes that can provide the network with maximum robustness in case of channel failures. To fix ideas, we first introduce some terminology. The status of the network is specified by a mapping  $\lambda : \mathcal{E} \to \{0, 1\}$  called a configuration. A channel being in the set

19.6 Static Network Codes 461

$$\lambda^{-1}(0) = \{ e \in \mathcal{E} : \lambda(e) = 0 \}$$
(19.153)

indicates the failure of that channel, and the subnetwork resulting from the deletion of all the channels in  $\lambda^{-1}(0)$  is called the  $\lambda$ -subnetwork. For the  $\lambda$ -subnetwork, the value of a max-flow from source node s to a non-source node t is denoted by maxflow<sub> $\lambda$ </sub>(t). Likewise, the value of a max-flow from source node s to a collection  $\mathcal{T}$  of non-source nodes is denoted by maxflow<sub> $\lambda$ </sub>( $\mathcal{T}$ ). It is easy to see that the total number of configurations is equal to  $2^{|\mathcal{E}|}$ .

**Definition 19.36.** Let  $\lambda$  be a configuration of the network. For an  $\omega$ -dimensional linear network code on the network, the  $\lambda$ -global encoding kernel of channel e, denoted by  $\mathbf{f}_{e,\lambda}$ , is the column  $\omega$ -vector calculated recursively in an upstream-to-downstream order by:

(19.154)  $\mathbf{f}_{e,\lambda} = \lambda(e) \sum_{d \in \text{In}(t)} k_{d,e} \mathbf{f}_{d,\lambda}$  for  $e \in \text{Out}(t)$ . (19.155) The  $\lambda$ -global encoding kernels of the  $\omega$  imaginary channels are independent of  $\lambda$  and form the standard basis of the space  $F^{\omega}$ .

Note that in the above definition, the local encoding kernels  $k_{d,e}$  are not changed with the configuration  $\lambda$ . Given the local encoding kernels, the  $\lambda$ -global encoding kernels can be calculated recursively by (19.154), while (19.155) serves as the boundary conditions. For a channel  $e \in \text{Out}(t)$  with  $\lambda(e) = 0$ , we see from (19.154) that

$$\mathbf{f}_{e,\lambda} = 0. \tag{19.156}$$

Equivalently, we can think of the local encoding kernels for all the adjacent pairs of channels  $(d, e), d \in \text{In}(t)$  as being set to 0. This can be seen by rewriting the formula in (19.154) as

$$\mathbf{f}_{e,\lambda} = \sum_{d \in \text{In}(t)} (\lambda(e)k_{d,e}) \, \mathbf{f}_{d,\lambda}.$$
(19.157)

As before, the message generated at source node s is denoted by a row  $\omega$ -vector **x**. When the prevailing configuration is  $\lambda$ , a node t receives the symbols  $\mathbf{x} \cdot \mathbf{f}_{d,\lambda}, d \in \text{In}(t)$ , from which it calculates the symbol  $\mathbf{x} \cdot \mathbf{f}_{e,\lambda}$  to be sent on each channel  $e \in \text{Out}(t)$  via

$$\mathbf{x} \cdot \mathbf{f}_{e,\lambda} = \mathbf{x} \left[ \lambda(e) \sum_{d \in \text{In}(t)} k_{d,e} \, \mathbf{f}_{d,\lambda} \right]$$
(19.158)

$$= \lambda(e) \sum_{d \in \text{In}(t)} k_{d,e}(\mathbf{x} \cdot \mathbf{f}_{d,\lambda}).$$
(19.159)

In particular, if  $\lambda(e) = 0$ , the zero symbol is sent on channel *e* regardless of the symbols received at node *t*.

In a real network, the zero symbol is not sent on a failed channel. Rather, whenever a symbol is not received on an input channel, the symbol is regarded by the receiving node as being the zero symbol.

For a configuration  $\lambda$  of the network, we let

$$V_{t,\lambda} = \langle \{ \mathbf{f}_{e,\lambda} : e \in \mathrm{In}(t) \} \rangle \tag{19.160}$$

for a node t,

$$V_{\mathcal{T},\lambda} = \langle \cup_{t \in \mathcal{T}} V_{t,\lambda} \rangle. \tag{19.161}$$

for a collection  $\mathcal{T}$  of nodes, and

$$V_{\xi,\lambda} = \langle \{ \mathbf{f}_{e,\lambda} : e \in \xi \} \rangle, \tag{19.162}$$

for a collection  $\xi$  of channels.

**Definition 19.37.** An  $\omega$ -dimensional linear network code on an acyclic network qualifies as a static linear multicast, a static linear broadcast, a static linear dispersion, or a static generic network code, respectively, if the following hold:

- (19.163) dim $(V_{t,\lambda}) = \omega$  for every configuration  $\lambda$  and every non-source node t with maxflow<sub> $\lambda$ </sub> $(t) \geq \omega$ .
- (19.164) dim $(V_{t,\lambda})$  = min $\{\omega, \max flow_{\lambda}(t)\}$  for every configuration  $\lambda$  and every non-source node t.
- (19.165) dim $(V_{\mathcal{T},\lambda})$  = min $\{\omega, \max flow_{\lambda}(\mathcal{T})\}\$  for every configuration  $\lambda$  and every collection  $\mathcal{T}$  of non-source nodes.
- (19.166) For any configuration  $\lambda$  and any nonempty collection of channels  $\xi \subset \mathcal{E}$ , if  $\xi = \min\{\omega, \max flow_{\lambda}(\xi)\}$ , then  $\mathbf{f}_{e,\lambda}, e \in \xi$  are linearly independent.

Here we have adopted Condition 3) in Theorem 19.32 for the purpose of defining a static generic network code. The qualifier "static" in the terms above stresses the fact that, while the configuration  $\lambda$  varies, the local encoding kernels remain unchanged. The advantage of using a static linear multicast, broadcast, or dispersion is that in case of channel failures, the local operation at every node in the network is affected only at the minimum level. Each receiving node in the network, however, needs to know the configuration  $\lambda$ before decoding can be done correctly. In implementation, this information can be provided by a separate signaling network.

For each class of static network codes in Definition 19.37, the requirement for its non-static version is applied to the  $\lambda$ -subnetwork for every configuration  $\lambda$ . Accordingly, a static linear multicast, a static linear broadcast, a static linear dispersion, and a static generic network code are increasingly stronger linear network codes as for the non-static versions.

*Example 19.38.* A 2-dimensional linear network code over GF(5) on the network in Figure 19.7 is prescribed by the local encoding kernels

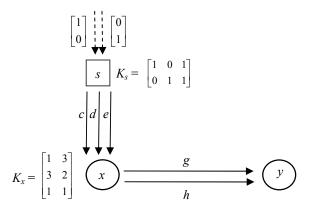


Fig. 19.7. A 2-dimensional GF(5)-valued static generic network code.

$$K_s = \begin{bmatrix} 1 & 0 & 1 \\ 0 & 1 & 1 \end{bmatrix} \tag{19.167}$$

and

$$K_x = \begin{bmatrix} 1 & 3 \\ 3 & 2 \\ 1 & 1 \end{bmatrix}.$$
 (19.168)

We claim that this is a static generic network code. Denote the three channels in In(x) by c, d, and e and the two channels in Out(x) by g and h. The vectors  $\mathbf{f}_{g,\lambda}$  and  $\mathbf{f}_{h,\lambda}$  for all possible configurations  $\lambda$  are tabulated in Table 19.1, from which it is straightforward to verify the condition (19.166).

The following is an example of a generic network code that does not qualify even as a static linear multicast.

*Example 19.39.* On the network in Figure 19.7, a 2-dimensional generic network code over GF(5) is prescribed by the local encoding kernels

$$K_s = \begin{bmatrix} 1 & 0 & 1 \\ 0 & 1 & 1 \end{bmatrix} \tag{19.169}$$

$\lambda(c)$	0		0		0		1		1		1		1	
$\lambda(d)$	0		1		1		0		0		1		1	
$\lambda(e)$	1		0		1		0		1		0		1	
$\mathbf{f}_{g,\lambda}$	$\lambda(g)$	1 1	$\lambda(g)$	$\begin{bmatrix} 0\\ 3 \end{bmatrix}$	$\lambda(g)$	$\begin{bmatrix} 1\\ 4 \end{bmatrix}$	$\lambda(g)$	$\begin{bmatrix} 1\\ 0 \end{bmatrix}$	$\lambda(g)$	$\begin{bmatrix} 2\\ 1 \end{bmatrix}$	$\lambda(g)$	$\begin{bmatrix} 1\\ 3 \end{bmatrix}$	$\lambda(g)$	$\begin{bmatrix} 2\\ 4 \end{bmatrix}$
$\mathbf{f}_{h,\lambda}$	$\lambda(h)$	1 1	$\lambda(h)$	$\begin{bmatrix} 0\\2 \end{bmatrix}$	$\lambda(h)$	$\begin{bmatrix} 1\\ 3 \end{bmatrix}$	$\lambda(h)$	$\begin{bmatrix} 3\\ 0 \end{bmatrix}$	$\lambda(h)$	$\begin{bmatrix} 4\\1 \end{bmatrix}$	$\lambda(h)$	$\begin{bmatrix} 3\\2 \end{bmatrix}$	$\lambda(h)$	$\begin{bmatrix} 4\\ 3 \end{bmatrix}$

**Table 19.1.** The vectors  $\mathbf{f}_{g,\lambda}$  and  $\mathbf{f}_{h,\lambda}$  for all possible configurations  $\lambda$  in Example 19.38.

and

$$K_x = \begin{bmatrix} 2 & 1 \\ 1 & 2 \\ 0 & 0 \end{bmatrix}.$$
 (19.170)

For a configuration  $\lambda$  such that

$$\lambda(c) = 0 \tag{19.171}$$

and

$$\lambda(d) = \lambda(e) = 1, \tag{19.172}$$

we have the  $\lambda$ -global encoding kernels

$$\mathbf{f}_{g,\lambda} = \begin{bmatrix} 0\\1 \end{bmatrix} \tag{19.173}$$

and

$$\mathbf{f}_{h,\lambda} = \begin{bmatrix} 0\\2 \end{bmatrix},\tag{19.174}$$

and therefore  $\dim(V_{y,\lambda}) = 1$ . On the other hand,  $\max flow_{\lambda}(y) = 2$ . Hence, this generic network code is not a static linear multicast.

Recall that in Algorithm 19.34 for constructing a generic network code, the key step chooses for a channel  $e \in \text{Out}(t)$  a vector in  $V_t$  to be the global encoding kernel  $\mathbf{f}_e$  such that

$$\mathbf{f}_e \notin V_{\zeta},\tag{19.175}$$

where  $\zeta$  is any collection of m-1 channels as prescribed with  $1 \leq m \leq \omega$ . This is equivalent to choosing scalar values for the local encoding kernels  $k_{d,e}$  for all  $d \in \text{In}(t)$  such that

$$\sum_{d\in \mathrm{In}(t)} k_{d,e} \, \mathbf{f}_d \notin V_{\zeta}. \tag{19.176}$$

Algorithm 19.34 is adapted below for the construction of a static generic network code.

Algorithm 19.40 (Construction of a Static Generic Network Code). This algorithm constructs an  $\omega$ -dimensional static generic network code over a finite field F on an acyclic network with  $|F| > 2^{|\mathcal{E}|} \sum_{m=1}^{\omega} {|\mathcal{E}|-1 \choose m-1}$ .

{ for (every node t, following an upstream-to-downstream order) for (every channel  $e \in Out(t)$ ) for (choose scalar values for  $k_{d,e}$  for all  $d \in In(t)$  such that for

any configuration  $\lambda$ ,  $\sum_{d \in \text{In}(t)} k_{d,e} \mathbf{f}_d \notin V_{\zeta,\lambda}$ , where  $\zeta$  is any collection of m-1 already processed channels such that  $\mathbf{f}_{e,\lambda}, e \in \zeta$  are linearly independent and  $V_{t,\lambda} \not\subset V_{\zeta,\lambda}$ ; // To see the existence of such values  $k_{d,e}$ , denote  $//\dim(V_{t,\lambda})$  by  $\nu$ . For any collection  $\zeta$  of channels // with  $V_{t,\lambda} \not\subset V_{\zeta,\lambda}$ , dim $(V_{t,\lambda} \cap V_{\zeta,\lambda}) < \nu$ . Consider // the linear mapping  $[k_{d,e}]_{d\in \text{In}(t)} \mapsto \sum_{d\in \text{In}(t)} k_{d,e} \mathbf{f}_{d,\lambda}$ // from  $F^{|\mathrm{In}(t)|}$  to  $F^{\omega}$ . The nullity of this linear // mapping is  $|In(t)| - \nu$ , so the pre-image of // the space  $(V_{t,\lambda} \cap V_{\zeta,\lambda})$  has dimension less than // |In(t)|. Thus the pre-image of  $\bigcup_{\lambda,\zeta} (V_{t,\lambda} \cap V_{\zeta,\lambda})$ // contains at most  $2^{|\mathcal{E}|} \sum_{m=1}^{\omega} {|\mathcal{E}|^{-1} \choose m-1} |F|^{|\text{In}(t)|-1}$ // elements, which are fewer than  $|F|^{|\text{In}(t)|}$  if 
$$\begin{split} // & |F| > 2^{|\mathcal{E}|} \sum_{m=1}^{\omega} \binom{|\mathcal{E}|-1}{m-1}, \\ for \; (every \; configuration \; \lambda) \end{split}$$
 $\mathbf{f}_{e,\lambda} = \lambda(e) \sum_{d \in \text{In}(t)} k_{d,e} \, \mathbf{f}_{d,\lambda};$ } }

Algorithm Verification. The explanation for the code constructed by Algorithm 19.40 being a static generic network code is exactly the same as that given for Algorithm 19.34. The details are omitted.

}

Algorithm 19.40 constitutes a constructive proof for the next theorem. By noting the lower bound on the required field size in the algorithm, we see that a generic network code can be constructed with high probability by randomly choosing the local encoding kernels provided that the base field is much larger than sufficient.

**Theorem 19.41.** There exist an  $\omega$ -dimensional static linear multicast, a static linear broadcast, a static linear dispersion, and a static generic network code on an acyclic network for sufficiently large base field F.

The requirements (19.163) through (19.166) in Definition 19.37 refer to all the  $2^{|\mathcal{E}|}$  possible configurations. Conceivably, a practical application may only need to deal with a certain collection  $\{\lambda_1, \lambda_2, \dots, \lambda_\kappa\}$  of configurations, where  $\kappa \ll 2^{|\mathcal{E}|}$ . Thus we may define, for instance, an  $\{\lambda_1, \lambda_2, \dots, \lambda_\kappa\}$ -static linear multicast and an  $\{\lambda_1, \lambda_2, \dots, \lambda_\kappa\}$ -static linear broadcast by replacing the conditions (19.19) and (19.20), respectively by

- (19.177) dim $(V_{t,\lambda}) = \omega$  for every configuration  $\lambda \in \{\lambda_1, \lambda_2, \dots, \lambda_\kappa\}$  and every non-source node t with maxflow $\lambda(t) \geq \omega$ .
- (19.178) dim $(V_{t,\lambda})$  = min $\{\omega, \max \text{flow}_{\lambda}(t)\}$  for every configuration  $\lambda \in \{\lambda_1, \lambda_2, \dots, \lambda_{\kappa}\}$  and every non-source node t.

Recall that Algorithm 19.34 is converted into Algorithm 19.40 by modifying the key step in the former. In a similar fashion, Algorithm 19.25 can be adapted for the construction of an  $\{\lambda_1, \lambda_2, \dots, \lambda_\kappa\}$ -static linear multicast or broadcast. This will lower the threshold on the sufficient size of the base field as well as the computational complexity. The details are left as an exercise.

## 19.7 Random Network Coding: A Case Study

We have seen in Corollary 19.24 that if the local encoding kernels of a linear network code are randomly chosen, a linear multicast can be obtained with high probability provided that the base field is sufficiently large. Since the code construction is independent of the network topology, the network code so constructed can be used when the network topology is unknown. In this section, we study an application of random network coding in *peer-to-peer* (P2P) networks. The system we will analyze is based on a prototype for large scale content distribution on such networks proposed in [121].

### 19.7.1 How the System Works

A file originally residing on a single server is to be distributed to a large number of users through a network. The server divides the file into k data blocks,  $B_1, B_2, \dots, B_k$ , and uploads coded versions of these blocks to different users according to some protocol. These users again help distributing the file by uploading blocks to other users in the network. By means of such repeated operations, a logical network called an *overlay network* is formed by the users as the process evolves. On this logical network, henceforth referred to as the network, information can be dispersed very rapidly, and the file is eventually delivered to every user in the network. Note that the topology of the network is not known ahead of time.

In the system, new users can join the network as a node at any time as long as the distribution process is active. Upon arrival, a new user will contact a designated node called the *tracker* that provides a subset of the other users already in the system, forming the set of neighboring nodes of the new user. Subsequent information flow in the network is possible only between neighboring nodes.

For the purpose of coding, the data blocks  $B_1, B_2, \dots, B_k$  are represented as symbols in a large finite field F referred to as the base field<sup>4</sup>. At the beginning of the distribution process, a Client A contacts the server and receives a number of *encoded blocks*. For example, the server uploads two encoded blocks  $E_1$  and  $E_2$  to Client A, where for i = 1, 2,

$$E_i = c_1^i B_1 + c_2^i B_2 + \dots + c_k^i B_k, (19.179)$$

<sup>&</sup>lt;sup>4</sup> In the system proposed in [121], the size of the base field is of the order  $2^{16}$ .

with  $c_j^i$ ,  $1 \leq j \leq k$  being chosen randomly from the base field F. Note that each  $E_1$  and  $E_2$  is some random linear combination of  $B_1, B_2, \dots, B_k$ .

In general, whenever a node needs to upload an encoded block to a neighboring node, the block is formed by taking a random linear combination of all the blocks possessed by that node. Continuing with the above example, when Client A needs to upload an encoded block  $E_3$  to a neighboring Client B, we have

$$E_3 = c_1^3 E_1 + c_2^3 E_2, (19.180)$$

where  $c_1^3$  and  $c_2^3$  are randomly chosen from *F*. Substituting (19.179) into (19.180), we obtain

$$E_3 = \sum_{j=1}^k (c_1^3 c_j^1 + c_2^3 c_j^2) B_j.$$
(19.181)

Thus  $E_3$  and in general every encoded block subsequently uploaded by a node in the network is some random linear combination of the data blocks  $B_1, B_2, \dots, B_k$ .

The exact strategy for downloading encoded blocks from the neighboring nodes so as to avoid receiving redundant information depends on the implementation. The main idea is that downloading from a neighboring node is necessary only if the neighboring node has at least one block not in the linear span of all the blocks possessed by that particular node. Upon receiving enough linearly independent encoded blocks, a node is able to decode the whole file.

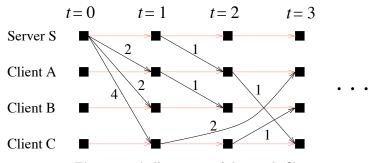
Compared with store-and-forward, the application of network coding as described in the above system can reduce the file download time because an encoded block uploaded by a node contains information about every block possessed by that node. Moreover, in case some nodes leave the system before the end of the distribution process, it is more likely that the remaining nodes have the necessary information to recover the whole file if network coding is used. In the following, we will give a quantitative analysis to substantiate these claimed advantages of network coding.

### 19.7.2 Model and Analysis

Let  $\mathcal{V}$  be the set of all the nodes in the system. In implementation, blocks of data are transmitted between neighboring nodes in an asynchronous manner, and possibly at different speeds. To simplify the analysis, we assume that every transmission from one node to a neighboring node is completed in an integral number of time units. Then we can unfold the network of nodes in discrete time into a graph  $\mathcal{G}^* = (\mathcal{V}^*, \mathcal{E}^*)$  with the node set

$$\mathcal{V}^* = \{ i_t : i \in \mathcal{V} \text{ and } t \ge 0 \},$$
 (19.182)

where node  $i_t \in \mathcal{V}^*$  corresponds to node  $i \in \mathcal{V}$  at time t. The edge set  $\mathcal{E}^*$  specified below is determined by the strategy adopted for the server as well



**Fig. 19.8.** A illustration of the graph  $\mathcal{G}^*$ .

as for all the other nodes in  $\mathcal{V}$  to request uploading of data blocks from the neighboring nodes. Specifically, there are two types of edges in  $\mathcal{E}^*$ :

- 1. There is an edge with capacity m from node  $i_t$  to node  $j_{t'}$ , where t < t', if m blocks are transmitted from node i to node j, starting at time t and ending at time t'.
- 2. For each  $i \in \mathcal{V}$  and  $t \geq 0$ , there is an edge with infinite capacity from node  $i_t$  to node  $i_{t+1}$ .

An edge of the second type models the assumption that the blocks, once possessed by a node, are retained in that node indefinitely over time. Without loss of generality, we may assume that all the blocks possessed by nodes  $i_l, l \leq t$ are transmitted uncoded on the edge from node  $i_t$  to node  $i_{t+1}$ .

An illustration of the graph  $\mathcal{G}^*$  up to t = 3 with  $\mathcal{V}$  consisting of the server S and three clients A, B, and C is given in Figure 19.8, where the edges with infinite capacities are lightened for clarity. Note that the graph  $\mathcal{G}^*$  is acyclic because each edge is pointed in the positive time direction and hence a cycle cannot be formed.

Denote the server S by node  $s \in \mathcal{V}$  and regard node  $s_0$  in  $\mathcal{G}^*$  as the source node generating the whole file consisting of k data blocks and multicasting it to all the other nodes in  $\mathcal{G}^*$  via random linear network coding, with the coefficients in the random linear combinations forming the encoded blocks being the local encoding kernels of the network code. Note that random network coding is applied on  $\mathcal{G}^*$ , not the logical network formed by the user nodes.

Also note that in order to simplify our description of the system, we have omitted the necessity of delivering the global encoding kernels to the nodes for the purpose of decoding. We refer the reader to the discussion toward the end of Section 19.3 for this implementation detail.

We are now ready to determine the time it takes for a particular node  $i \in \mathcal{V}$  to receive the whole file. Denote the value of a max-flow from node  $s_0$  to a node  $v \in \mathcal{G}^*$  other than  $s_0$  by maxflow(v). When the base field is sufficiently large, by Corollary 19.24, with probability close to 1, the network code generated randomly during the process is a linear multicast, so that those nodes  $i_t$  with

19.7 Random Network Coding: A Case Study 469

$$\max flow(i_t) \ge k \tag{19.183}$$

can receive the whole file. In other words, with high probability, the time it takes a node  $i \in \mathcal{V}$  to receive the whole file is equal to  $t^*$ , the minimum t that satisfies (19.183). Obviously, this is a lower bound on the time it takes a node  $i \in \mathcal{V}$  to receive the whole file, and it is achievable with high probability by the system under investigation. In the rare event that node i cannot decode at time  $t^*$ , it can eventually decode upon downloading some additional encoded blocks from the neighboring nodes.

When some nodes leave the system before the end of the distribution process, an important question is whether the remaining nodes have the necessary information to recover the whole file. To be specific, assume that a subset of users  $\mathcal{U}^c \subset \mathcal{V}$  leave the system after time t, and we want to know whether the users in  $\mathcal{U} = \mathcal{V} \setminus \mathcal{U}^c$  have sufficient information to recover the whole file. If they do, by further exchanging information among themselves, every user in  $\mathcal{U}$  can eventually receive the whole file (provided that no more nodes leave the system). Toward this end, again consider the graph  $\mathcal{G}^*$ . Let

$$\mathcal{U}_t = \{ u_t : u \in \mathcal{U} \} \tag{19.184}$$

and denote the value of a max-flow from node  $s_0$  to the set of nodes  $\mathcal{U}_t$  by maxflow( $\mathcal{U}_t$ ). If

$$\max flow(\mathcal{U}_t) \ge k, \tag{19.185}$$

then the users in  $\mathcal{U}$  with high probability would have the necessary information to recover the whole file. This is almost the best possible performance one can expect from such a system, because if

$$\max flow(\mathcal{U}_t) < k, \tag{19.186}$$

it is simply impossible for the users in  $\mathcal{U}$  to recover the whole file even if they are allowed to exchange information among themselves.

Thus we see that random network coding provides the system with both maximum bandwidth efficiency and maximum robustness. However, additional computational resource is required compared with store-and-forward. These are engineering tradeoffs in the design of such systems.

We conclude this section by an example demonstrating the advantage of random network coding when it is applied to packet networks with packet loss.

Example 19.42. The random network coding scheme discussed in this section can be applied to packet networks with packet loss. Consider the network depicted in Figure 19.9 consisting of three nodes, s, t, and u. Data packets are sent from node s to node u via node t. Let the packet loss rates of channels (s, t)and (t, u) be  $\gamma$ , i.e., a fraction  $\gamma$  of packets are lost during their transmission through the channel. Then the fraction of packets sent by node s that are eventually received at node u is  $(1 - \gamma)^2$ .

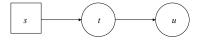


Fig. 19.9. A simple packet network.

To fix idea, assume the packet size is sufficiently large and one packet is sent on each channel per unit time. To remedy the problem of packet loss, a fountain code [49] can be employed at node s. This would allow data packets to be sent from node s to node u reliably at an effective rate equal to  $(1-\gamma)^2$ . On the other hand, by using the random network coding scheme we have discussed, data packets can be sent from node s to node u reliably at an effective rate equal to  $1-\gamma$ , which is strictly higher than  $(1-\gamma)^2$  whenever  $\gamma > 0$ . This can be proved by means of the analysis presented in this section. The details are left as an exercise.

While a fountain code can remedy the problem of packet loss between the source node and the sink node, it cannot prevent the packet loss rate from accumulating when packets are routed through the network. On the other hand, the use of random network coding allows information to be transmitted from the source node to the sink node at the maximum possible rate, namely the min-cut between the source node and the sink node after the packet loss in the channels has been taken into account.

### Problems

In the following, let  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  be the underlying directed acyclic network on which the linear network code is defined, and let s be the unique source node in the network.

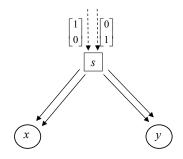
- 1. Show that in a network with the capacities of all the edges equal to 1, the number of edge-disjoint paths from source node s to a non-source node t is equal to maxflow(t).
- For the network code in Definitions 19.4 and 19.6, show that if the global encoding mappings are linear, then so are the local encoding mappings. (Yeung *et al.* [365].)
- Network transfer matrix Consider an ω-dimensional linear network code.
   a) Prove (19.91).
  - b) Fix an upstream-to-downstream order for the channels in the network and let K be the  $|\mathcal{E}| \times |\mathcal{E}|$  matrix with the (d, e)th element equal to  $k_{d,e}$  if (d, e) is an adjacent pair of channels and equal to 0 otherwise. Let A be the  $\omega \times |\mathcal{E}|$  matrix obtaining by appending  $|\mathcal{E}| - |\text{Out}(s)|$ columns of zeroes to  $K_s$ , and  $B_e$  be the  $|\mathcal{E}|$ -column vector with all the components equal to 0 except that the *e*th component is equal to 1. Show that

$$\mathbf{f}_e = A(I - K)^{-1}B_e$$

for all  $e \in \mathcal{E}$ . The matrix  $M = (I - K)^{-1}$  is called the *network transfer* matrix.

(Koetter and Médard [182].)

- 4. Apply Lemma 19.17 to obtain a lower bound on the field size for the existence of a 2-dimensional linear multicast on the butterfly network.
- 5. Show that  $\sum_{m=1}^{\omega} {\binom{|\mathcal{E}|-1}{m-1}}$  is a polynomial in  $|\mathcal{E}|$  of degree  $\omega$ . This is the lower bound on the required field size in Algorithm 19.34.
- 6. Verify that the network code in Example 19.38 is a generic network code.
- 7. Simplified characterization of a generic network code Consider an  $\omega$ dimensional generic network code on a network for which  $|\operatorname{Out}(s)| \geq \omega$ .
  - a) Show that Condition 3) in Theorem 19.32 can be modified to restricting the cardinality of  $\xi$  to  $\omega$ . Hint: If  $|\xi| < \omega$ , expand  $\xi$  by including a certain subset of the channels in Out(s).
  - b) Simplify Algorithm 19.34 and tighten the lower bound on the required field size accordingly.
  - (Tan *et al.* [315].)
- 8. For the network below, prove the non-existence of a two-dimensional binary generic network code.



- 9. Modify the Jaggi-Sanders algorithm for the construction of a static linear multicast.
- 10. Obtain a lower bound on the required field size and determine the computational complexity when Algorithm 19.40 is adapted for the construction of an  $\{\lambda_1, \lambda_2, \dots, \lambda_{\kappa}\}$ -static generic network code.
- 11. Show that a transformation of a static generic network code is also a static generic network code.
- 12. A generic network code as a linear dispersion Expand the network  $\mathcal{G}$  into a network  $\mathcal{G}' = (\mathcal{V}', \mathcal{E}')$  as follows. For an edge  $e \in \mathcal{E}$ , let the edge be from node  $v_e$  to node  $w_e$ . Install a new node  $t_e$  and replace edge e by two new edges e' and e'', where e' is from node  $v_e$  to node  $t_e$  and e'' is from node  $t_e$ to node  $w_e$ . Show that a linear dispersion on  $\mathcal{G}'$  is equivalent to a generic

network code on  $\mathcal{G}$ . Hint: Use Theorem 19.32. (Kwok and Yeung [196], Tan *et al.* [315].)

13. Multi-rate linear broadcast Consider a network on which an  $\omega$ -dimensional linear network code over a base field F is defined. For all  $e \in \mathcal{E}$ , let

$$\mathbf{f}_e' = \begin{bmatrix} I & \mathbf{b} \end{bmatrix} \mathbf{f}_e$$

where I is the  $(\omega - 1) \times (\omega - 1)$  identity matrix and **b** is an  $(\omega - 1)$ -column vector.

- a) Show that  $\mathbf{f}'_e, e \in \mathcal{E}$  constitute the global encoding kernels of an  $(\omega 1)$ -dimensional linear network code on the same network.
- b) Show that the  $(\omega 1)$ -dimensional linear network code in a) and the original  $\omega$ -dimensional linear network code have the same local encoding kernels for all the non-source nodes.

It was shown in Fong and Yeung [103] that an  $(\omega - 1)$ -dimensional linear broadcast can be constructed from any  $\omega$ -dimensional linear broadcast by choosing a suitable vector **b**, provided  $|F| \ge |\mathcal{V}|$ . As such, multi-rate linear multicast/broadcast can be supported on a network without changing the local encoding kernels of the non-source nodes.

- 14. Let a message  $\mathbf{x} \in F^{\omega}$  be generated at source node *s* in a network for which  $\max flow(t) \geq \omega$  for all non-source nodes *t*. Show that  $\mathbf{x}$  can be multicast to all the non-source nodes by store-and-forward. In other words, for this special case, network coding has no advantage over store-and-forward if complete information on the network topology is known ahead of time. This result is implied by a theorem on *directed spanning tree packing* by Edmonds [90] (see also Wu *et al.* [347]).
- 15. Let L be the length of the message  $\mathbf{x}$  generated at source node s, where L is divisible by maxflow(t) for all non-source nodes t. Allowing multiple usage of the network, devise a linear network coding scheme such that each non-source node t can receive  $\mathbf{x}$  in  $L/\max flow(t)$  units of time. Such a scheme enables each non-source node in the network to receive the message within the shortest possible time.
- 16. Consider distributing a message of 5 data blocks in a P2P network with 4 nodes, Server S and Clients A, B, and C, by the system discussed in Section 19.7. Assume each data block is sufficiently large. The following transmissions take place during the process.

From	Τo	Start Time	End Time	# Bloc
$\mathbf{S}$	А	0	1	2
$\mathbf{S}$	В	0	1	3
$\mathbf{S}$	$\mathbf{C}$	0	1	2
В	А	1	2	1
$\mathbf{C}$	В	1	3	2
$\mathbf{S}$	В	2	3	1
В	$\mathbf{C}$	2	3	2

From To Start Time End Time # Blocks

a) Which client is the first to receive the whole message?

- b) If Client B leaves the system after t = 3, do Clients A and C have sufficient information to reconstruct the whole message?
- c) Suppose the hard disk of Client B crashes at t = 1.5 and loses 2 blocks of data. Repeat b) by making the assumption that the transmissions by Client B starting at  $t \leq 1$  are not affected by the disk failure.
- 17. Prove the claim in Example 19.42 that by using random network coding, data packets can be sent from node s to node u at an effective rate equal to  $1 \gamma$ .

## **Historical Notes**

The achievability of the max-flow bound by linear network codes was proved by Li *et al.* [205] using a vector space approach and then by Koetter and Médard [182] using a matrix approach. These two approaches correspond respectively to the notions of global encoding kernel and local encoding kernel discussed here. Neither the construction in [205] for a generic network code nor the construction in [182] for a linear multicast are polynomial-time algorithms. Jaggi and Sanders *et al.* [165] obtained a polynomial-time algorithm for constructing a linear multicast by modifying the construction of a generic network code in [205]. A polynomial-time algorithm for constructing a generic network code was subsequently obtained in Yeung *et al.* [365].

In [182], static network code was introduced and its existence was proved. An explicit construction of such codes was given in [365].

The optimality of random network coding was proved in Ahlswede *et al.* [7]. Ho *et al.* [145] proved the optimality of random linear network coding and proposed the use of such codes on an unknown network topology. A tight upper bound on the probability of decoding error for random linear network coding has recently been obtained by Balli *et al.* [22].

Implementation issues of network coding were discussed in Chou *et al.* [65]. The application of random network coding in peer-to-peer networks discussed in Section 19.7 is due to Gkantsidis and Rodriguez [121].

Cai and Yeung have generalized single-source network coding on acyclic networks to network error correction [51][363][52] and secure network coding [50]. Network error correction subsumes classical algebraic coding, while secure network coding subsumes secret sharing in cryptography.

The presentation in this chapter is largely based on the tutorial paper by Yeung *et al.* [365]. The various characterizations of a generic network code is due to Tan *et al.* [315].

A directed network is *cyclic* if it contains at least one directed cycle. In Chapter 19, we have discussed network coding over an acyclic network, in which there exists an upstream-to-downstream order on the nodes. Following such an order, whenever a node encodes, all the information needed would have already been received on the input channels of that node. For a cyclic network, such an order of the nodes does not exists. This makes network coding over a cyclic network substantially different from that over an acyclic network.

### 20.1 Delay-Free Cyclic Networks

When we discussed network coding over an acyclic network in Chapter 19, we assume that there is no propagation delay in the network. Based on this assumption, a linear network code can be specified by either the local description in Definition 19.6 or the global description in Definition 19.7. The local and global descriptions of a linear network code are equivalent over an acyclic network because given the local encoding kernels, the global encoding kernels can be calculated recursively in any upstream-to-downstream order. In other words, the equation (19.17) has a unique solution for the global encoding kernels in terms of the local encoding kernels, while (19.18) serves as the boundary conditions.

If these descriptions are applied to a cyclic network, it is not clear whether for given local encoding kernels, there exists a unique solution for the global encoding kernels. In the following, we give one example with a unique solution, one with no solution, and one with multiple solutions.

*Example 20.1.* Consider the cyclic network in Figure 20.1. Let (s, t) precede (v, t) in the ordering among the channels. Similarly, let (s, t') precede (v, t'). Given the local encoding kernels

$$K_s = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, K_t = K_{t'} = \begin{bmatrix} 1 \\ 0 \end{bmatrix}, K_u = \begin{bmatrix} 1 \\ 1 \end{bmatrix}, K_v = \begin{bmatrix} 1 & 1 \end{bmatrix}, \quad (20.1)$$

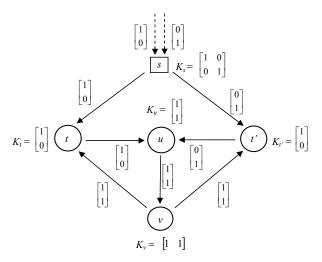


Fig. 20.1. A 2-dimensional linear broadcast on a cyclic network.

the equation (19.17) yields the following unique solution for the global encoding kernels:

$$\mathbf{f}_{(s,t)} = \mathbf{f}_{(t,u)} = \begin{bmatrix} 1\\0 \end{bmatrix}, \ \mathbf{f}_{(s,t')} = \mathbf{f}_{(t',u)} = \begin{bmatrix} 0\\1 \end{bmatrix}$$
(20.2)

$$\mathbf{f}_{(u,v)} = \mathbf{f}_{(v,t)} = \mathbf{f}_{(v,t')} = \begin{bmatrix} 1\\1 \end{bmatrix}.$$
 (20.3)

These global encoding kernels are shown in Figure 20.1, and they in fact define a 2-dimensional linear broadcast regardless of the choice of the base field. Note that

$$k_{(v,t),(t,u)} = 0 (20.4)$$

and

$$k_{(v,t'),(t',u)} = 0 (20.5)$$

in the prescribed local encoding kernels prevent information from looping in the directed cycles

$$(t, u), (u, v), (v, t)$$
 (20.6)

and

$$(t', u), (u, v), (v, t'),$$
 (20.7)

respectively.

*Example 20.2.* An arbitrarily prescribed set of local encoding kernels on a cyclic network is unlikely to be compatible with any global encoding kernels. In Figure 20.2(a), a local encoding kernel is prescribed at each node in a

cyclic network. Had a global encoding kernel  $\mathbf{f}_e$  existed for each channel e, the requirement (19.17) would imply the equations

$$\mathbf{f}_{(x,y)} = \begin{bmatrix} 1\\0 \end{bmatrix} + \mathbf{f}_{(w,x)}$$
(20.8)

$$\mathbf{f}_{(y,w)} = \begin{bmatrix} 0\\1 \end{bmatrix} + \mathbf{f}_{(x,y)} \tag{20.9}$$

$$\mathbf{f}_{(w,x)} = \mathbf{f}_{(y,w)},\tag{20.10}$$

which sum up to

$$\begin{bmatrix} 1\\0 \end{bmatrix} = \begin{bmatrix} 0\\1 \end{bmatrix},\tag{20.11}$$

a contradiction.

The nonexistence of compatible global encoding kernels can also be interpreted in terms of message transmission. Let the message  $\mathbf{x} = [a b]$  be a generic vector in  $F^2$ , where F denotes the base field. The symbol transmitted on channel e, given by  $\mathbf{x} \cdot \mathbf{f}_e$ , are shown in Figure 20.2(b). In particular, the symbols transmitted on channels (x, y), (y, w), and (w, x), namely p, q, and r, are related through

$$p = a + r \tag{20.12}$$

$$q = b + p \tag{20.13}$$

$$r = q. \tag{20.14}$$

These equalities imply that

$$a + b = 0,$$
 (20.15)

a contradiction to the independence between the two components a and b of a generic message.

*Example 20.3.* Let F be an extension field of  $GF(2)^1$ . Consider the same prescription of the local encoding kernels at the nodes as in Example 20.2 except that

$$K_S = \begin{bmatrix} 1 & 1 \\ 0 & 0 \end{bmatrix}. \tag{20.16}$$

The following three sets of global encoding kernels meet the requirement (19.17) in the definition of a linear network code:

$$\mathbf{f}_{(s,x)} = \mathbf{f}_{(s,y)} = \begin{bmatrix} 1\\0 \end{bmatrix}, \ \mathbf{f}_{(x,y)} = \begin{bmatrix} 0\\0 \end{bmatrix}, \ \mathbf{f}_{(y,w)} = \mathbf{f}_{(w,x)} = \begin{bmatrix} 1\\0 \end{bmatrix}; \qquad (20.17)$$

$$\mathbf{f}_{(s,x)} = \mathbf{f}_{(s,y)} = \begin{bmatrix} 1\\0\\ \end{bmatrix}, \ \mathbf{f}_{(x,y)} = \begin{bmatrix} 1\\0\\ \end{bmatrix}, \ \mathbf{f}_{(y,w)} = \mathbf{f}_{(w,x)} = \begin{bmatrix} 0\\0\\ \end{bmatrix}; \qquad (20.18)$$

$$\mathbf{f}_{(s,x)} = \mathbf{f}_{(s,y)} = \begin{bmatrix} 1\\0 \end{bmatrix}, \ \mathbf{f}_{(x,y)} = \begin{bmatrix} 0\\1 \end{bmatrix}, \ \mathbf{f}_{(y,w)} = \mathbf{f}_{(w,x)} = \begin{bmatrix} 1\\1 \end{bmatrix}.$$
(20.19)

 $^1$  In an extension field of GF(2), the arithmetic on the symbols 0 and 1 are modulo 2 arithmetic.

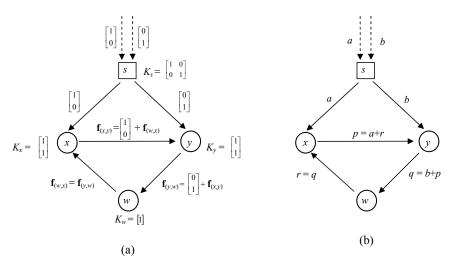


Fig. 20.2. An example of a cyclic network and local encoding kernels that do not render a solution for the global encoding kernels.

### 20.2 Convolutional Network Codes

In a real network, the propagation delay, which includes the processing delay at the nodes and the transmission delay over the channels, cannot be zero. For a cyclic network, this renders the implementation non-physical because the transmission on an output channel of a node can only depend on the information received on the input channels of that node. Besides, technical difficulties as described in the last section arise even with the ideal assumption that there is no propagation delay.

In this section, we introduce the *unit-delay network* as a model for network coding on a cyclic network  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ , where  $\mathcal{V}$  and  $\mathcal{E}$  are the sets of nodes and channels of the network, respectively. In this model, a symbol is transmitted on every channel in the network at every discrete time index, with the transmission delay equal to exactly one time unit. Intuitively, this assumption on the transmission delay over a channel ensures no information looping in the network even in the presence of a directed cycle. The results to be developed in this chapter, although discussed in the context of cyclic networks, apply equally well to acyclic networks.

As a time-multiplexed network in the combined space-time domain, a unitdelay network can be unfolded with respect to the time dimension into an indefinitely long network called a *trellis network*. Corresponding to a physical node t is a sequence of nodes  $t_0, t_1, t_2, \cdots$  in the trellis network, with the subscripts being the time indices. A channel  $e_j$  in the trellis network represents the transmission on the physical channel e between times j and j + 1. When the physical channel e is from node t to node u, the channel  $e_j$  in the trellis network is from node  $t_j$  to node  $u_{j+1}$ . Note that the trellis network is acyclic

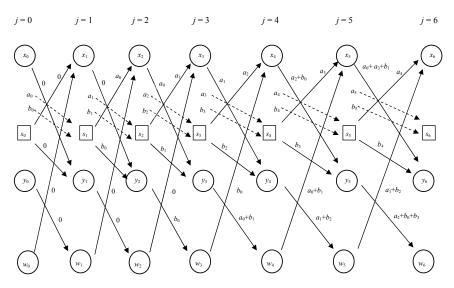


Fig. 20.3. The trellis network depicting a convolutional network code defined on the physical network in Figure 20.2.

regardless of the topology of the physical network, because all the channels are pointing in the forward time directions so that a directed cycle cannot be formed.

Example 20.4. Regard the network in Figure 20.2 as a unit-delay network. For each channel e in the network, the scalar values in the base field F transmitted on the channels  $e_j$ ,  $j \ge 0$  in the corresponding trellis network are determined by the local encoding kernels. This is illustrated in Figure 20.3. For instance, the channels  $(x, y)_j$ ,  $j \ge 0$  carry the scalar values

$$0, 0, a_0, a_1, a_2 + b_0, a_0 + a_3 + b_1, a_1 + a_4 + b_2, \cdots,$$
(20.20)

respectively. This constitutes an example of a convolutional network code to be formally defined in Definition 20.6.

Let  $c_j$  be the scalar value in F transmitted on a particular channel in the network at time j. A succinct mathematical expression for the sequence of scalars  $c_0, c_1, c_2, \cdots$  is the z-transform

$$\sum_{j=0}^{\infty} c_j z^j = c_0 + c_1 z + c_2 z^2 + \cdots, \qquad (20.21)$$

where the power j of the *dummy variable* z represents discrete time. The pipelining of scalars transmitted over a time-multiplexed channel can thus be regarded as the transmission of a power series over the channel. For example,

the transmission of a scalar value on the channel  $(x, y)_j$  for each  $j \ge 0$  in the trellis network in Figure 20.3 translates into the transmission of the power series

$$a_0 z^2 + a_1 z^3 + (a_2 + b_0) z^4 + (a_0 + a_3 + b_1) z^5 + (a_1 + a_4 + b_2) z^6 + \cdots$$
(20.22)

over the channel (x, y) in the network in Figure 20.2.

The z-transform in (20.21) is a power series in the dummy variable z, which is regarded as either a real or a complex number in the context of signal analysis. However, in the context of convolutional coding, the reader should not regard the z-transform in (20.21) as anything more than a representation of the sequence of scalars  $c_0, c_1, c_2, \cdots$ . Specifically, the dummy variable z is not associated with any value, and there is no notion of convergence. Such power series are called *formal power series*.

Given a field F, consider rational functions of a dummy variable z of the form

$$\frac{p(z)}{1+zq(z)},\tag{20.23}$$

where p(z) and q(z) are polynomials. The following properties of such a function are relevant to our subsequent discussion:

- 1. The denominator has a constant term, so the function can be expanded into a power series by long division (see Example 20.5).
- 2. If p(z) is not the zero polynomial, the inverse function, namely

$$\frac{1+zq(z)}{p(z)},$$
 (20.24)

exists.

Note that the rational function in (20.24) does not represent a power series if p(z) contains the factor z, or equivalently, does not contain a constant term.

The ring of power series over F is conventionally denoted by F[[z]]. Rational functions of the form (20.23) will be called *rational power series* which constitute a ring denoted by  $F\langle z \rangle$  [365]. It follows directly from the definitions that  $F\langle z \rangle$  is a subring of F[[z]]. We refer the reader to [112] for a comprehensive treatment of abstract algebra.

In the following, we illustrate the concepts of rational power series through a few simple examples.

Example 20.5. If z is a complex number, then we can write

$$\frac{1}{1-z} = 1 + z + z^2 + z^3 + \dots$$
 (20.25)

provided that |z| < 1, where we have interpreted the coefficients in the power series on the right hand side as real (or complex) numbers. If |z| > 1, the above expression is not meaningful because the power series diverges.

However, if we do not associate z with a value but regard the coefficients in the power series as elements in a commutative ring, we can always write

$$(1-z)(1+z+z^2+z^3+\cdots) = (1+z+z^2+z^3+\cdots) - (z+z^2+z^3+\cdots)$$
(20.26)

$$= 1.$$
 (20.27)

In this sense, we say that 1 - z is the reciprocal of the power series  $1 + z + z^2 + z^3 + \cdots$  and write

$$\frac{1}{1-z} = 1 + z + z^2 + z^3 + \cdots.$$
 (20.28)

We also say that  $1 + z + z^2 + z^3 + \cdots$  is the power series expansion of  $\frac{1}{1-z}$ . In fact, the power series on the right hand side can be readily obtained by dividing 1 by 1 - z using long division.

Alternatively, we can seek the inverse of 1 - z by considering the identity

$$(1-z)(a_0 + a_1z + a_2z^2 + \dots) = 1.$$
(20.29)

By equating the powers of z on both sides, we have

$$a_0 = 1$$
 (20.30)

$$-a_0 + a_1 = 0 \tag{20.31}$$

$$-a_1 + a_2 = 0 \tag{20.32}$$

Then by forward substitution, we immediately obtain

$$1 = a_0 = a_1 = a_2 = \cdots, \tag{20.34}$$

which gives exactly the power series obtained by long division. The reader can easily verify that long division indeed mimics the process of forward substitution.

For polynomials p(z) and q(z) where q(z) is not the zero polynomial, we can always expand the rational function  $\frac{p(z)}{q(z)}$  into a series. However, such a series is not always a power series. For example,

$$\frac{1}{z-z^2} = \frac{1}{z} \left[ \frac{1}{1-z} \right] \tag{20.35}$$

$$= \frac{1}{z}(1+z+z^2+\cdots)$$
(20.36)

$$= z^{-1} + 1 + z + z^2 + \cdots$$
 (20.37)

The above is not a power series because of the term involving a negative power of z. In fact, the identity

$$(z - z2)(a_0 + a_1 z + a_2 z2 + \dots) = 1$$
(20.38)

has no solution for  $a_0, a_1, a_2, \cdots$  since there is no constant term on the left hand side. Therefore,  $\frac{1}{z-z^2}$  indeed does not have a power series expansion.

From the above example, we see that  $\frac{p(z)}{q(z)}$  represents a rational power series if and only if q(z) has a nonzero constant term, or equivalently, does not contain the factor z.

**Definition 20.6 (Convolutional Network Code).** An  $\omega$ -dimensional convolutional network code on a unit-delay network over a base field F consists of an element  $k_{d,e}(z) \in F\langle z \rangle$  for every adjacent pair of channels (d, e) in the network as well as a column  $\omega$ -vector  $\mathbf{f}_e(z)$  over  $F\langle z \rangle$  for every channel e such that:

(20.39)  $\mathbf{f}_{e}(z) = z \sum_{d \in In(t)} k_{d,e}(z) \mathbf{f}_{d}(z)$  for  $e \in Out(t)$ .

(20.40) The vectors  $\mathbf{f}_{e}(z)$  for the imaginary channels  $e \in \mathrm{In}(s)$  consist of scalar components that form the standard basis of the vector space  $F^{\omega}$ .

The vector  $\mathbf{f}_e(z)$  is called the global encoding kernel for channel e, and  $k_{d,e}(z)$  is called the local encoding kernel for the adjacent pair of channels (d, e). The  $|\text{In}(t)| \times |\text{Out}(t)|$  matrix

$$K_t(z) = [k_{d,e}(z)]_{d \in In(t), e \in Out(t)}$$
(20.41)

is called the local encoding kernel at node t.

The constraint (20.39) is the time-multiplexed version of (19.17), with the factor z in the equation indicating a unit-time delay that represents the transmission delay over a channel. In the language of electronic circuit theory, for an adjacent pair of channels (d, e), the "gain" from channel d to channel e is given by  $zk_{d,e}(z)$ .

A convolutional network code over a unit-delay network can be viewed as a discrete-time *linear time-invariant* (LTI) system defined by the local encoding kernels, where the local encoding kernel  $k_{d,e}(z)$  specifies the impulse response of an LTI filter from channel d to channel e. The requirement that  $k_{d,e}(z)$  is a power series corresponds to the causality of the filter. The additional requirement that  $k_{d,e}(z)$  is rational ensures that the filter is implementable by a finite circuitry of shift registers. Intuitively, once the local encoding kernels are given, the global encoding kernels are uniquely determined. This is explained as follows. Write

$$\mathbf{f}_{e}(z) = \sum_{j=0}^{\infty} \mathbf{f}_{e,j} z^{j} = \mathbf{f}_{e,0} + \mathbf{f}_{e,1} z + \mathbf{f}_{e,2} z^{2} + \cdots$$
(20.42)

and

20.2 Convolutional Network Codes 483

$$k_{d,e}(z) = \sum_{j=0}^{\infty} k_{d,e,j} z^j = k_{d,e,0} + k_{d,e,1} z + k_{d,e,2} z^2 + \cdots, \qquad (20.43)$$

where  $\mathbf{f}_{e,j}$  is a column  $\omega$ -vector in  $F^{\omega}$  and  $k_{d,e,j}$  is a scalar in F. Then the equation in (20.39) can be written in time domain as the convolutional equation

$$\mathbf{f}_{e,j} = \sum_{d \in \text{In}(t)} \left( \sum_{u=0}^{j-1} k_{d,e,u} \, \mathbf{f}_{d,j-1-u} \right)$$
(20.44)

for  $j \ge 0$ , with the boundary conditions provided by (20.40):

- The vectors  $\mathbf{f}_{e,0}, e \in \text{In}(t)$  form the standard basis of the vector space  $F^{\omega}$ .
- The vectors  $\mathbf{f}_{e,j}, e \in \text{In}(t)$  are the zero vector for all  $j \ge 1$ .

For j = 0, the summation in (20.44) is empty, so that  $\mathbf{f}_{e,0}$  vanishes. For  $j \ge 0$ , the right hand side of (20.44) involves the vectors  $\mathbf{f}_{d,i}$  for only  $0 \le i \le j - 1$ . Thus the vectors  $\mathbf{f}_{e,j}$ ,  $j \ge 1$  can be calculated recursively via (20.44) with the boundary condition

$$\mathbf{f}_{d,0} = 0 \quad \text{for all } d \in \mathcal{E}. \tag{20.45}$$

Together with  $\mathbf{f}_{e,0} = 0$ , the global encoding kernel  $\mathbf{f}_e(z)$  is determined (cf. (20.42)). In other words, in a convolutional network code over a unit-delay network, the global encoding kernels are determined once the local encoding kernels are given. From (20.42), we see that the components of  $\mathbf{f}_e(z)$  are power series in z, so  $\mathbf{f}_e(z)$  is a column  $\omega$ -vector over F[[z]]. In Theorem 20.9, we will further establish that the components of the global encoding kernels are in fact rational functions in z, proving that  $\mathbf{f}_e(z)$  is indeed a column  $\omega$ -vector over  $f\langle z \rangle$  as required in Definition 20.6 for a convolutional network code.

*Example 20.7.* In Figure 20.2, denote the two imaginary channels by (o, s) and (o, s)'. A convolutional network code is specified by the prescription of a local encoding kernel at every node as shown in the figure:

$$K_s(z) = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, \ K_x(z) = K_y(z) = \begin{bmatrix} 1 \\ 1 \end{bmatrix}, \ K_w(z) = \begin{bmatrix} 1 \\ 1 \end{bmatrix},$$
(20.46)

and a global encoding kernel for every channel:

$$\mathbf{f}_{(o,s)}(z) = \begin{bmatrix} 1\\0 \end{bmatrix}, \ \mathbf{f}_{(o,s)'}(z) = \begin{bmatrix} 0\\1 \end{bmatrix}$$
(20.47)

$$\mathbf{f}_{(s,x)}(z) = z \begin{bmatrix} 1 & 0\\ 0 & 1 \end{bmatrix} \begin{bmatrix} 1\\ 0 \end{bmatrix} = \begin{bmatrix} z\\ 0 \end{bmatrix}$$
(20.48)

$$\mathbf{f}_{(s,y)}(z) = z \begin{bmatrix} 1 & 0\\ 0 & 1 \end{bmatrix} \begin{bmatrix} 0\\ 1 \end{bmatrix} = \begin{bmatrix} 0\\ z \end{bmatrix}$$
(20.49)

$$\mathbf{f}_{(x,y)}(z) = \begin{bmatrix} z^2/(1-z^3)\\ z^4/(1-z^3) \end{bmatrix}$$
(20.50)

$$\mathbf{f}_{(y,w)}(z) = \begin{bmatrix} z^3/(1-z^3) \\ z^2/(1-z^3) \end{bmatrix}$$
(20.51)

$$\mathbf{f}_{(w,x)}(z) = \begin{bmatrix} z^4/(1-z^3) \\ z^3/(1-z^3) \end{bmatrix},$$
(20.52)

where the last three global encoding kernels have been solved from the following equations:

$$\mathbf{f}_{(x,y)}(z) = z \left[ \mathbf{f}_{(s,x)}(z) \ \mathbf{f}_{(w,x)}(z) \right] \begin{bmatrix} 1\\1\\ 1 \end{bmatrix} = z^2 \begin{bmatrix} 1\\0\\ 1 \end{bmatrix} + z \ \mathbf{f}_{(w,x)}(z)$$
(20.53)

$$\mathbf{f}_{(y,w)}(z) = z \begin{bmatrix} \mathbf{f}_{(s,y)}(z) \ \mathbf{f}_{(x,y)}(z) \end{bmatrix} \begin{bmatrix} 1\\1 \end{bmatrix} = z^2 \begin{bmatrix} 0\\1 \end{bmatrix} + z \ \mathbf{f}_{(x,y)}(z)$$
(20.54)

$$\mathbf{f}_{(w,x)}(z) = z(\mathbf{f}_{(y,w)}(z)) \left[1\right] = z \, \mathbf{f}_{(y,w)}(z).$$
(20.55)

These local and global encoding kernels of a 2-dimensional convolutional network code are summarized in Figure 20.4.

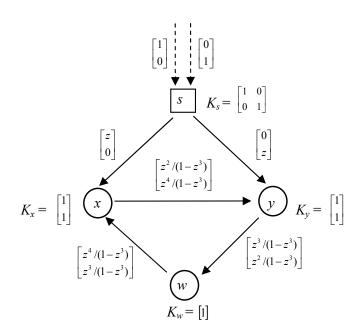


Fig. 20.4. The local and global encoding kernels of the convolutional network code in Example 20.7.

Represent the message generated at source node s at time j, where  $j \ge 0$ , by a row  $\omega$ -vector  $\mathbf{x}_j \in F^{\omega}$ . Equivalently, source node s generates the message pipeline represented by the z-transform

$$\mathbf{x}(z) = \sum_{j=0}^{\infty} \mathbf{x}_j z^j, \qquad (20.56)$$

which is a row  $\omega$ -vector over F[[z]], the ring of power series over F. Here,  $\mathbf{x}(z)$  is not necessarily rational.

Through a convolutional network code, each channel e carries the power series  $\mathbf{x}(z) \mathbf{f}_e(z)$ . Write

$$\mathbf{x}(z)\mathbf{f}_e(z) = \sum_{j=0}^{\infty} m_{e,j} z^j, \qquad (20.57)$$

where

$$m_{e,j} = \sum_{u=0}^{j} \mathbf{x}_u \, \mathbf{f}_{e,j-u}.$$
(20.58)

For  $e \in Out(t)$ , from the equation in (20.39), we obtain

$$\mathbf{x}(z) \mathbf{f}_e(z) = \mathbf{x}(z) \left[ z \sum_{d \in \text{In}(t)} k_{d,e}(z) \mathbf{f}_d(z) \right]$$
(20.59)

$$= z \sum_{d \in \operatorname{In}(t)} k_{d,e}(z) \left[ \mathbf{x}(z) \, \mathbf{f}_d(z) \right], \qquad (20.60)$$

or equivalently in time domain,

$$m_{e,j} = \sum_{d \in \text{In}(t)} \left( \sum_{u=0}^{j-1} k_{d,e,u} \, m_{d,j-1-u} \right).$$
(20.61)

The reader should compare (20.61) with (20.44). Note that the scalar values  $m_{e,j}$ ,  $j \ge 1$  can be calculated recursively via (20.61) with the boundary condition

$$m_{d,0} = 0 \quad \text{for all } d \in \mathcal{E}.$$
 (20.62)

Thus a node t calculates the scalar value  $m_{e,j}$  for transmitting on each output channel e at time j from the cumulative information it has received on all the input channels up to time j-1. The convolutional equation (20.61) can be implemented by a finite circuit of shift-registers in a causal manner because the local encoding kernels belong to  $F\langle z \rangle$ , the ring of rational power series over F (cf. Definition 20.6).

*Example 20.8.* Consider the convolutional network code in Example 20.7. Let source node s pipelines the message

$$\mathbf{x}(z) = \begin{bmatrix} \sum_{j=0}^{\infty} a_j z^j & \sum_{j=0}^{\infty} b_j z^j \end{bmatrix}.$$
 (20.63)

Then the five channels (s, x), (s, y), (x, y), (y, w), and (w, x) carry the following power series, respectively:

$$\mathbf{x}(z)\,\mathbf{f}_{(s,x)}(z) = \sum_{j=0}^{\infty} a_j z^{j+1}$$
(20.64)

$$\mathbf{x}(z)\,\mathbf{f}_{(s,y)}(z) = \sum_{j=0}^{\infty} b_j z^{j+1}$$
(20.65)

$$\mathbf{x}(z)\mathbf{f}_{(x,y)}(z) = \left(\sum_{j=0}^{\infty} a_j z^{j+2} + \sum_{j=0}^{\infty} b_j z^{j+4}\right) / (1-z^3)$$
(20.66)

$$= \left(\sum_{j=0}^{\infty} a_j z^{j+2} + \sum_{j=0}^{\infty} b_j z^{j+4}\right) \sum_{j=0}^{\infty} z^{3j}$$
(20.67)  
$$= a_0 z^2 + a_1 z^3 + (a_2 + b_0) z^4$$

$$+(a_0 + a_3 + b_1)z^5 + \cdots$$
(20.68)

$$\mathbf{x}(z)\,\mathbf{f}_{(y,w)}(z) = \left(\sum_{j=0}^{\infty} a_j z^{j+3} + \sum_{j=0}^{\infty} b_j z^{j+2}\right) / (1-z^3)$$
(20.69)

$$\mathbf{x}(z)\,\mathbf{f}_{(w,x)}(z) = \left(\sum_{j=0}^{\infty} a_j z^{j+4} + \sum_{j=0}^{\infty} b_j z^{j+3}\right) / (1-z^3).$$
(20.70)

At each time  $j \ge 0$ , the source generates a message  $\mathbf{x}_j = [a_j b_j]$ . Thus channel (s, x) carries the scalar 0 at time 0 and the scalar  $a_{j-1}$  at time  $j \ge 1$ . Similarly, channel (s, y) carries the scalar 0 at time 0 and the scalar  $b_{j-1}$  at time  $j \ge 1$ . For every channel e, write

$$\mathbf{x}(z) \mathbf{f}_e(z) = \sum_{j=0}^{\infty} m_{e,j} z^j$$
(20.71)

as in (20.57). The actual encoding process at node x is as follows. At time j, node x has received the sequence  $m_{d,0}, m_{d,1}, \dots, m_{d,j-1}$  for d = (s, x) and (w, x). Accordingly, at time  $j \ge 1$ , channel (x, y) transmits the scalar value

$$m_{(x,y),j} = \sum_{u=0}^{j-1} k_{(s,x),(x,y),u} m_{(s,x),j-1-u}$$

20.2 Convolutional Network Codes 487

$$+\sum_{u=0}^{j-1} k_{(w,x),(x,y),u} m_{(w,x),j-1-u}$$
(20.72)

$$= m_{(s,x),j-1} + m_{(w,x),j-1}.$$
(20.73)

Similarly, channels (y, w) and (w, x) transmit the scalar values

$$m_{(y,w),j} = m_{(s,y),j-1} + m_{(x,y),j-1}$$
(20.74)

and

$$m_{(w,x),j} = m_{(y,w),j-1},$$
 (20.75)

respectively. The values  $m_{(x,y),j}$ ,  $m_{(y,w),j}$ , and  $m_{(w,x),j}$  for  $j \ge 1$  can be calculated recursively by the above formulas with the boundary condition

$$m_{e,0} = 0 \quad \text{for all } e \in \mathcal{E},$$
 (20.76)

and they are shown in the trellis network in Figure 20.3 for small values of j. For instance, the channel (x, y) carries the scalar values

$$m_{(x,y),0} = 0, \ m_{(x,y),1} = 0, \ m_{(x,y),2} = a_0, \ m_{(x,y),3} = a_1, m_{(x,y),4} = a_2 + b_0, \ m_{(x,y),5} = a_0 + a_3 + b_1, \ \cdots.$$
(20.77)

The z-transform of this sequence is

$$x(z) \mathbf{f}_{(x,y)}(z) = \left(\sum_{j=0}^{\infty} a_j z^{j+2} + \sum_{j=0}^{\infty} b_j z^{j+4}\right) / (1-z^3), \quad (20.78)$$

as calculated in (20.68).

In the discussion following Definition 20.6, we have shown that once the local encoding kernels of a convolutional network code over a unit-delay network are given, the global encoding kernels are determined. The proof of the next theorem further provides a simple closed-form expression for the global encoding kernels  $\mathbf{f}_e(z)$ , from which it follows that the entries in  $\mathbf{f}_e(z)$  indeed belong to  $F\langle z \rangle$  as required in Definition 20.6.

**Theorem 20.9.** Let F be the base field and  $k_{d,e}(z) \in F\langle z \rangle$  be given for every adjacent pair of channels (d, e) on a unit-delay network. Then there exists a unique  $\omega$ -dimensional convolutional network code over F with  $k_{d,e}(z)$  as the local encoding kernel for every (d, e).

*Proof.* Let the unit-delay network be represented by a directed graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ . Let  $[k_{d,e}(z)]$  be the  $|\mathcal{E}| \times |\mathcal{E}|$  matrix in which both the rows and columns are indexed by  $\mathcal{E}$ , with the (d, e)th entry equal to the given  $k_{d,e}(z)$  if (d, e) is an adjacent pair of channels, and equal to zero otherwise. Denote the global encoding kernel of channel e by  $\mathbf{f}_e(z)$  if exists. Let  $[\mathbf{f}_e(z)]$  be the  $\omega \times |\mathcal{E}|$  matrix

obtained by putting the global encoding kernels  $\mathbf{f}_e(z), e \in \mathcal{E}$  in juxtaposition, Let  $H_s(z)$  be the  $\omega \times |\mathcal{E}|$  matrix obtained by appending  $|\mathcal{E}| - |\operatorname{Out}(s)|$  columns of zeroes to the local encoding kernel  $K_s(z)$ . The requirements (20.39) and (20.40) in Definition 20.6 can be written as

$$[\mathbf{f}_e(z)] = z[\mathbf{f}_e(z)] [k_{d,e}(z)] + zIH_s(z), \qquad (20.79)$$

where I in the above denotes the  $\omega \times \omega$  identity matrix representing the global encoding kernels  $\mathbf{f}_e(z)$ ,  $e \in \text{In}(s)$  in juxtaposition. Rearranging the terms in (20.79), we obtain

$$[\mathbf{f}_e(z)](I - z[k_{d,e}(z)]) = zH_s(z).$$
(20.80)

In the matrix  $z[k_{d,e}(z)]$ , the diagonal elements are equal to zero because (e, e) does not form an adjacent pair of channels for all  $e \in \mathcal{E}$ , while the non-zero off-diagonal elements all contain the factor z. Therefore,  $\det(I - z[k_{d,e}(z)])$  has the form

$$1 + zq(z),$$
 (20.81)

where  $q(z) \in F\langle z \rangle$ , so that it is invertible inside  $F\langle z \rangle$  because

$$[\det(I - z[k_{d,e}(z)])]^{-1} = \frac{1}{1 + zq(z)}$$
(20.82)

is a rational power series. It follows that

$$(I - z[k_{d,e}(z)])^{-1} (20.83)$$

exists and is a matrix over  $F\langle z \rangle$ . Then the unique solution for  $[\mathbf{f}_e(z)]$  in (20.80) is given by

$$[\mathbf{f}_e(z)] = zH_s(z)(I - z[k_{d,e}(z)])^{-1}.$$
(20.84)

With the two matrices  $[k_{d,e}(z)]$  and  $H_s(z)$  representing the given local encoding kernels and the matrix  $[\mathbf{f}_e(z)]$  representing the global encoding kernels, (20.84) is a closed-form expression for the global encoding kernels in terms of the local encoding kernels. In particular,  $[\mathbf{f}_e(z)]$  is a matrix over  $F\langle z \rangle$  because all the matrices on the right hand side of (20.84) are over  $F\langle z \rangle$ . Thus we conclude that all the components of the global encoding kernels are in  $F\langle z \rangle$ . Hence, the given local encoding kernels  $k_{d,e}(z)$  for all adjacent pairs (d, e) together with the associated global encoding kernels  $\mathbf{f}_e(z), e \in \mathrm{In}(s) \cup \mathcal{E}$  constitute a unique convolutional network code over the unit-delay network  $\mathcal{G}$ .  $\Box$ 

In view of Definition 19.7 for the global description of a linear network code over an acyclic network, Definition 20.6 can be regarded as the global description of a convolutional network code over a unit-delay network, while Theorem 20.9 renders a local description by specifying the local encoding kernels only.

### 20.3 Decoding of Convolutional Network Codes

For a node t, let

$$F_t(z) = [\mathbf{f}_e(z)]_{e \in \text{In}(t)}$$
 (20.85)

be the  $\omega \times |\text{In}(t)|$  matrix obtained by putting the global encoding kernels  $\mathbf{f}_e(z), e \in \text{In}(t)$  in juxtaposition. In the following, we define a *convolutional multicast*, the counterpart of a linear multicast defined in Chapter 19, for a unit-delay cyclic network. The existence of a convolutional multicast will also be established.

**Definition 20.10 (Convolutional Multicast).** An  $\omega$ -dimensional convolutional network code on a unit-delay network qualifies as an  $\omega$ -dimensional convolutional multicast if for every non-source node t with maxflow $(t) \geq \omega$ , there exists an  $|\text{In}(t)| \times \omega$  matrix  $D_t(z)$  over  $F\langle z \rangle$  and a positive integer  $\tau$  such that

$$F_t(z) D_t(z) = z^{\tau} I, \qquad (20.86)$$

where  $\tau > 0$  depends on node t and I is the  $\omega \times \omega$  identity matrix. The matrix  $D_t(z)$  are called the decoding kernel and the decoding delay at node t, respectively.

Source node s generates the message pipeline

$$\mathbf{x}(z) = \sum_{j=0}^{\infty} \mathbf{x}_j z^j, \qquad (20.87)$$

where  $\mathbf{x}_j$  is a row  $\omega$ -vector in  $F^{\omega}$  and  $\mathbf{x}(z)$  is a row  $\omega$ -vector over F[[z]]. Through the convolutional network code, a channel e carries the power series  $\mathbf{x}(z) \mathbf{f}_e(z)$ . The power series  $\mathbf{x}(z) \mathbf{f}_e(z)$  received by a node t from the input channels  $e \in \text{In}(t)$  form the row |In(t)|-vector  $\mathbf{x}(z) F_t(z)$  over F[[z]]. If the convolutional network code is a convolutional multicast, node t can use the decoding kernel  $D_t(z)$  to calculate

$$(\mathbf{x}(z) F_t(z)) D_t(z) = \mathbf{x}(z) (F_t(z) D_t(z))$$
(20.88)

$$= \mathbf{x}(z)(z^{\tau}I) \tag{20.89}$$

$$= z^{\tau} \mathbf{x}(z). \tag{20.90}$$

The row  $\omega$ -vector  $z^{\tau} \mathbf{x}(z)$  of power series represents the message pipeline generated by source node *s* delayed by  $\tau$  time units. Note that  $\tau > 0$  because the message pipeline  $\mathbf{x}(z)$  is delayed by one time unit at node *s*.

*Example 20.11.* Consider the network in Figure 20.4. Again let source node s pipelines the message

$$\mathbf{x}(z) = \begin{bmatrix} \sum_{j=0}^{\infty} a_j z^j & \sum_{j=0}^{\infty} b_j z^j \end{bmatrix}.$$
 (20.91)

For node x, we have

$$F_x(z) = \begin{bmatrix} z \ z^4 / (1 - z^3) \\ 0 \ z^3 / (1 - z^3) \end{bmatrix}.$$
 (20.92)

Let

$$D_x(z) = \begin{bmatrix} z^2 & -z^3 \\ 0 & 1-z^3 \end{bmatrix}.$$
 (20.93)

Then

$$F_x(z)D_t(z) = z^3 I_2 (20.94)$$

 $(I_2 \text{ is the } 2\times 2 \text{ identity matrix}).$  From channels (s,x) and (w,x), node x receives the row vector

$$\mathbf{x}(z)F_x(z) = \left[\sum_{j=0}^{\infty} a_j z^{j+1} \quad \sum_{j=0}^{\infty} \frac{a_j z^{j+4} + b_j z^{j+3}}{1 - z^3}\right]$$
(20.95)

and decodes the message pipeline as

$$z^{3} \mathbf{x}(z) = \left[\sum_{j=0}^{\infty} a_{j} z^{j+1} \sum_{j=0}^{\infty} \frac{a_{j} z^{j+4} + b_{j} z^{j+3}}{1 - z^{3}}\right] \begin{bmatrix} z^{2} & -z^{3} \\ 0 & 1 - z^{3} \end{bmatrix}.$$
 (20.96)

Decoding at node y is similar. Thus the 2-dimensional convolutional network code is a convolutional multicast.

Toward proving the existence of a convolutional multicast, we first observe that Lemma 19.17 can be strengthened as follows with essentially no change in the proof.

**Lemma 20.12.** Let  $g(y_1, y_2, \dots, y_m)$  be a nonzero polynomial with coefficients in a field  $\tilde{F}$ . For any subset  $\tilde{E}$  of  $\tilde{F}$ , if  $|\tilde{E}|$  is greater than the degree of g in every  $y_j$ , then there exist  $a_1, a_2, \dots, a_m \in \tilde{E}$  such that

$$g(a_1, a_2, \cdots, a_m) \neq 0.$$
 (20.97)

In the above lemma, the values  $a_1, a_2, \dots, a_m$  can be found by exhaustive search in  $\tilde{E}$  provided that  $\tilde{E}$  is finite. If  $\tilde{E}$  is infinite, simply replace  $\tilde{E}$  by a sufficiently large finite subset of  $\tilde{E}$ .

**Theorem 20.13.** There exists an  $\omega$ -dimensional convolutional multicast over any base field F. Furthermore, the local encoding kernels of the convolutional multicast can be chosen in any sufficiently large subset  $\Phi$  of  $F\langle z \rangle$ .

*Proof.* Recall the equation (20.84) in the proof of Theorem 20.10:

$$[\mathbf{f}_e(z)] = zH_s(z)(I - z[k_{d,e}(z)])^{-1}.$$
(20.98)

In this equation, the  $\omega \times |\mathcal{E}|$  matrix  $[\mathbf{f}_e(z)]$  on the left hand side represents the global encoding kernels, while the  $\omega \times |\mathcal{E}|$  matrix  $H_s(z)$  and the  $|\mathcal{E}| \times |\mathcal{E}|$  matrix  $[k_{d,e}(z)]$  on the right hand side represent the local encoding kernels. Analogous to the proof of Theorem 19.20, denote by  $(F\langle z \rangle)[*]$  the polynomial ring over  $F\langle z \rangle$  with all the  $k_{d,e}(z)$  as indeterminates.

Let t be a non-source node with maxflow $(t) \geq \omega$ . Then there exist  $\omega$  edgedisjoint paths from the  $\omega$  imaginary channels to  $\omega$  distinct channels in In(t). Put the global encoding kernels of these  $\omega$  channels in juxtaposition to form the  $\omega \times \omega$  matrix  $L_t(z)$  over  $(F\langle z \rangle)[*]$ . We will show that

$$\det(L_t(z)) \neq 0 \in (F\langle z \rangle)[*]. \tag{20.99}$$

Toward proving (20.99), it suffices to show that

$$\det(L_t(z)) \neq 0 \in F\langle z \rangle \tag{20.100}$$

when the determinant is evaluated at some particular values for the indeterminates  $k_{d,e}(z)$ . Analogous to the proof of Theorem 19.20, we set

$$k_{d,e}(z) = 1 \tag{20.101}$$

for all adjacent pairs of channels (d,e) along any one of the  $\omega$  edge-disjoint paths, and set

$$k_{d,e}(z) = 0 \tag{20.102}$$

otherwise. Then with a suitable indexing of the columns, the matrix  $L_t(z)$  becomes diagonal with all the diagonal entries being powers of z. Hence,  $\det(L_t(z))$  is equal to some positive power of z, proving (20.100) for this particular choice of the indeterminates  $k_{d,e}(x)$  and hence proving (20.99). As the conclusion (20.99) applies to every non-source node t with maxflow $(t) \geq \omega$ , it follows that

$$\prod_{t:\max flow(t) \ge \omega} \det(L_t(z)) \neq 0 \in (F\langle z \rangle)[*].$$
(20.103)

Let F(z) be the conventional notation for the field of rational functions in z over the given base field F. The ring  $F\langle z \rangle$  of rational power series is a subset of F(z). Then any subset  $\Phi$  of  $F\langle z \rangle$  is also a subset of F(z). Note that the ring  $F\langle z \rangle$  is infinite. Then for any sufficiently large subset  $\Phi$  of  $F\langle z \rangle$ , we can apply Lemma 20.12 to the polynomial in (20.103) with  $\tilde{F} = F(z)$  and  $\tilde{E} = \Phi$  to see that we can choose a value  $a_{d,e}(z) \in F\langle z \rangle$  for each of the indeterminates  $k_{d,e}(z)$  so that

$$\prod_{t:\max flow(t) > \omega} \det(L_t(z)) \neq 0 \in F\langle z \rangle.$$
(20.104)

when evaluated at  $k_{d,e}(z) = a_{d,e}(z)$  for all (d, e), which in turn implies that

$$\det(L_t(z)) \neq 0 \in F\langle z \rangle \tag{20.105}$$

for all nodes t such that  $\max(t) \ge \omega$ .

Henceforth, the local encoding kernel  $k_{d,e}(z)$  will be fixed at the appropriately chosen value  $a_{d,e}(z)$  for all (d, e) as prescribed above. Without loss

#### 492 20 Single-Source Linear Network Coding: Cyclic Networks

of generality, we assume that  $L_t(z)$  consists of the first  $\omega$  columns of  $F_t(z)$ . From (20.105), we can write

$$\det(L_t(z)) = z^{\tau} \left[ \frac{1 + zq(z)}{p(z)} \right],$$
(20.106)

where p(z) and q(z) are polynomials over F and p(z) is not the zero polynomial. Note that the right hand side of (20.106) is the general form for a nonzero rational function in z. In this particular context, since the columns of  $L_t(z)$  are global encoding kernels as prescribed by (20.84), each containing the factor z in the numerator, we see that  $\tau > 0$ .

Denote by  $J_t(z)$  the adjoint matrix<sup>2</sup> of  $L_t(z)$ . Take the  $\omega \times \omega$  matrix

$$\left[\frac{p(z)}{1+zq(z)}\right]J_t(z) \tag{20.107}$$

and append to it  $|In(t)| - \omega$  rows of zeroes to form an  $|In(t)| \times \omega$  matrix  $D_t(z)$ . Then

$$F_t(z)D_t(z) = \begin{bmatrix} L_t(z) & 0 \end{bmatrix} \begin{bmatrix} \begin{bmatrix} \frac{p(z)}{1+zq(z)} \end{bmatrix} J_t(z) \\ 0 \end{bmatrix}$$
(20.108)

$$= \left[\frac{p(z)}{1+zq(z)}\right] L_t(z)J_t(z)$$
(20.109)

$$= \left[\frac{p(z)}{1+zq(z)}\right] \det(L_t(z))I \tag{20.110}$$

$$= z^{\tau}I, \qquad (20.111)$$

where the last equality follows from (20.106). Hence, the matrix  $D_t(z)$  qualifies as a decoding kernel at node t in Definition 20.10. This proves the existence of the convolutional multicast as required.  $\Box$ 

The proof of Theorem 20.13 constitutes an algorithm for constructing a convolutional multicast. By noting the lower bound on the size of  $\tilde{E}$  in Lemma 20.12, a convolutional multicast can be constructed with high probability by randomly choosing the local encoding kernels in the subset  $\Phi$  of  $F\langle z \rangle$ provided that  $\Phi$  is much larger than sufficient.

Example 20.14. When the base field F is sufficiently large, Theorem 20.13 can be applied with  $\Phi = F$  so that the local encoding kernels of the convolutional multicast can be chosen to be scalars. This special case is the convolutional counterpart of Theorem 19.20 for the existence of a linear multicast over an acyclic network. In this case, the local encoding kernels can be found by exhaustive search over F.

<sup>&</sup>lt;sup>2</sup> For a matrix B whose entries are elements in a ring, denote by  $\operatorname{Adj}(B)$  the adjoint matrix of B. Then  $\operatorname{Adj}(B)B = B\operatorname{Adj}(B) = \det(A)I$ .

More generally, by virtue of Lemma 20.12, the same exhaustive search applies to any large enough subset  $\Phi$  of  $F\langle z \rangle$ . For example, F can be GF(2) and  $\Phi$  can be the set of all binary polynomials up to a sufficiently large degree.

## Problems

- 1. Show that the right hand side of (20.106) is the general form for a nonzero rational function in z.
- 2. A formal Laurent series over a field F has the form

$$a_{-m}z^{-m} + a_{-(m-1)}z^{-(m-1)} + \dots + a_{-1}z^{-1} + a_0 + a_1z + a_2z^2 + \dots,$$

where m is a nonnegative integer. Show that for any formal Laurent series f(z) over F, there exists a unique formal Laurent series g(z) over F such that f(z)g(z) = 1.

3. Verify the following series expansion:

$$\frac{1}{1-z} = -z^{-1} - z^{-2} - z^{-3} - \cdots$$

Can you obtain this series by long division?

4. Construct a finite circuit of shift-registers that implements a discrete-time LTI system with transfer function

$$\frac{a_0+a_1z+\cdots+a_nz^n}{b_0+b_1z+\cdots+b_nz^n},$$

where  $a_i$  and  $b_i$  are elements in a finite field and  $b_0 \neq 0$ .

- 5. Consider the convolutional network code in Figure 20.4.
  - a) Is it a convolutional multicast?
  - b) If your answer in a) is positive, give the decoding kernel at node y with minimum decoding delay.
  - c) Change  $K_s$  to  $\begin{bmatrix} 1 & 1 \\ 0 & 1 \end{bmatrix}$  and determine the corresponding global encoding kernels.
  - d) Instead of a convolutional multicast, can you construct a linear multicast on the network?

## **Historical Notes**

The asymptotic achievability of the max-flow bound for cyclic networks was proved by Ahlswede *et al.* [7], where an example of a convolutional network code achieving this bound was given. Li *et al.* [205] conjectured the existence of a convolutional multicast, which was subsequently proved by Koetter and

494 20 Single-Source Linear Network Coding: Cyclic Networks

Médard [182]. Construction and decoding of convolutional multicast have been studied by Erez and Feder [93][94], Fragouli and Soljanin [110], and Barbero and Ytrehus [23]. The formulation and treatment of convolutional codes here is based on Li and Yeung [204] (see also Yeung *et al.* [365]).

In Chapters 19 and 20, we have discussed single-source network coding in which an information source is multicast in a point-to-point communication network. The maximum rate at which information can be multicast has a simple characterization in terms of the maximum flows in the graph representing the network. In this chapter, we consider the more general multi-source network coding problem in which more than one *mutually independent* information sources are generated at possibly different nodes, and each of the information sources is multicast to a specific set of nodes.

The achievable information rate region of a multi-source network coding problem, which will be formally defined in Section 21.4, refers to the set of all possible rates at which multiple information sources can be multicast simultaneously on a network. In a single-source network coding problem, we are interested in characterizing the maximum rate at which information can be multicast from the source node to all the sink nodes. In a multi-source network coding problem, we are interested in characterizing the achievable information rate region.

As discussed in Section 17.3, source separation is not necessarily optimal for multi-source network coding. It is therefore *not* a simple extension of singlesource network coding. Unlike the single-source network coding problem which has an explicit solution, the multi-source network coding problem has not been completely solved. In this chapter, by making use of the tools we have developed for information inequalities in Chapter 13 to Chapter 15, we will develop an implicit characterization of the achievable information rate region for multi-source network coding on acyclic networks.

## 21.1 The Max-Flow Bounds

The max-flow bound, which fully characterizes the maximum rate of an information source that can be multicast in a network, plays a central role in single-source network coding. We now revisit this bound in the context of

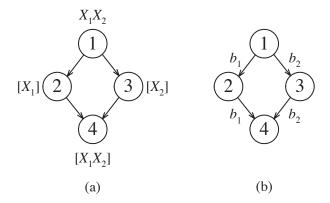


Fig. 21.1. A network which achieves the max-flow bound.

multi-source network coding. In the following discussion, the unit of information is the bit.

Consider the graph in Figure 21.1(a). The capacity of each edge is equal to 1. Two independent information sources  $X_1$  and  $X_2$  with rates  $\omega_1$  and  $\omega_2$ , respectively are generated at node 1. Suppose we want to multicast  $X_1$  to nodes 2 and 4 and multicast  $X_2$  to nodes 3 and 4. In the figure, an information source in square brackets is one which is to be received at that node.

It is easy to see that the values of a max-flow from node 1 to node 2, from node 1 to node 3, and from node 1 to node 4 are respectively 1, 1, and 2. At node 2 and node 3, information is received at rates  $\omega_1$  and  $\omega_2$ , respectively. At node 4, information is received at rate  $\omega_1 + \omega_2$  because  $X_1$  and  $X_2$  are independent. Applying the max-flow bound at nodes 2, 3, and 4, we have

$$\omega_1 \le 1 \tag{21.1}$$

$$\omega_2 \le 1 \tag{21.2}$$

and

$$\omega_1 + \omega_2 \le 2, \tag{21.3}$$

respectively. We refer to (21.1) to (21.3) as the max-flow bounds. Figure 21.2 is an illustration of all  $(\omega_1, \omega_2)$  which satisfy these bounds, where  $\omega_1$  and  $\omega_2$  are obviously nonnegative.

We now show that the rate pair (1, 1) is achievable. Let  $b_1$  be a bit generated by  $X_1$  and  $b_2$  be a bit generated by  $X_2$ . In the scheme in Figure 21.1(b),  $b_1$  is received at node 2,  $b_2$  is received at node 3, and both  $b_1$  and  $b_2$  are received at node 4. Thus the multicast requirements are satisfied, and the information rate pair (1, 1) is achievable. This implies that all  $(\omega_1, \omega_2)$  which satisfy the max-flow bounds are achievable because they are all inferior to (1, 1) (see Figure. 21.2). In this sense, we say that the max-flow bounds are achievable.

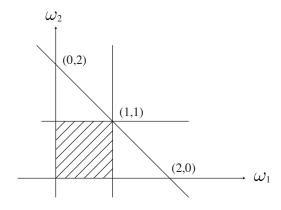


Fig. 21.2. The max-flow bounds for the network in Figure 21.1.

Suppose we now want to multicast  $X_1$  to nodes 2, 3, and 4 and multicast  $X_2$  to node 4 as illustrated in Figure 21.3. Applying the max-flow bound at either node 2 or node 3 gives

$$\omega_1 \le 1, \tag{21.4}$$

and applying the max-flow bound at node 4 gives

$$\omega_1 + \omega_2 \le 2. \tag{21.5}$$

Figure 21.4 is an illustration of all  $(\omega_1, \omega_2)$  which satisfy these bounds.

We now show that the information rate pair (1, 1) is not achievable. Suppose we need to send a bit  $b_1$  generated by  $X_1$  to nodes 2, 3, and 4 and send a bit  $b_2$  generated by  $X_2$  to node 4. Since  $b_1$  has to be recovered at node 2, the bit sent to node 2 must be an invertible transformation of  $b_1$ . This implies that the bit sent to node 2 cannot not depend on  $b_2$ . Similarly, the bit sent to node 3 also cannot depend on  $b_2$ . Therefore, it is impossible for node 4 to recover  $b_2$  because both the bits received at nodes 2 and 3 do not depend on

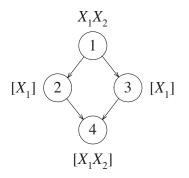


Fig. 21.3. A network which does not achieve the max-flow bounds.

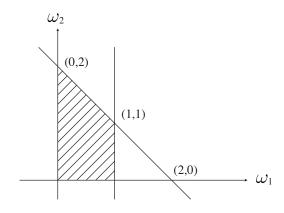


Fig. 21.4. The max-flow bounds for the network in Figure 21.3.

 $b_2$ . Thus the information rate pair (1, 1) is not achievable, which implies that the max-flow bounds (21.4) and (21.5) are not achievable.

From this example, we see that the max-flow bounds do not always fully characterize the achievable information rate region. We leave it as an exercise for the reader to show that for this example, source separation is in fact optimal.

## 21.2 Examples of Application

Multi-source network coding is a very rich model which encompasses many communication situations arising from fault-tolerant network communication, disk array, satellite communication, etc. In this section, we discuss some applications of the model.

### 21.2.1 Multilevel Diversity Coding

Let  $X_1, X_2, \dots, X_K$  be K information sources in decreasing order of importance. These information sources are encoded into pieces of information. There are a number of users, each of them having access to a certain subset of the information pieces. Each user belongs to a *level* between 1 and K, where a Level k user can decode  $X_1, X_2, \dots, X_k$ . This model, called *multilevel diver*sity coding, finds applications in fault-tolerant network communication, disk array, and distributed data retrieval.

Figure 21.5 shows a graph which represents a 3-level diversity coding system. The graph consists of three layers of nodes. The top layer consists of a node at which information sources  $X_1$ ,  $X_2$ , and  $X_3$  are generated. These information sources are encoded into three pieces, each of which is stored in a distinct node in the middle layer. A dummy node is associated with such a

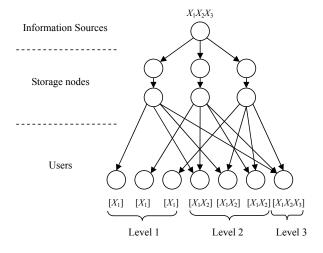


Fig. 21.5. A 3-level diversity coding system.

node to model the effect that the same information is retrieved every time the node is accessed (see the discussion in Section 17.2). The nodes in the bottom layer represent the users, each of them belonging to one of the three levels. Each of the three Level 1 users has access to a distinct node in the second layer (through the associated dummy node) and decodes  $X_1$ . Similarly, each of the three Level 2 users has access to a distinct set of two nodes in the second layer and decodes  $X_1$  and  $X_2$ . There is only one Level 3 user, who has access to all the three nodes in the second layer and decodes  $X_1$  and  $X_3$ .

The model represented by the graph in Figure 21.5 is called *symmetrical* 3-level diversity coding because the model is unchanged by permuting the nodes in the middle layer. By degenerating information sources  $X_1$  and  $X_3$ , the model is reduced to the diversity coding model discussed in Section 18.2.

In the following, we describe two applications of symmetrical multilevel diversity coding:

**Fault-Tolerant Network Communication** In a computer network, a data packet can be lost due to buffer overflow, false routing, breakdown of communication links, etc. Suppose the packet carries K messages,  $X_1, X_2, \dots, X_K$ , in decreasing order of importance. For improved reliability, the packet is encoded into K sub-packets, each of which is sent over a different channel. If any k sub-packets are received, then the messages  $X_1, X_2, \dots, X_k$  can be recovered.

**Disk Array** Consider a disk array which consists of K disks. The data to be stored in the disk array are segmented into K pieces,  $X_1, X_2, \dots, X_K$ , in decreasing order of importance. Then  $X_1, X_2, \dots, X_K$  are encoded into K pieces, each of which is stored on a separate disk. When any k out of the K disks are functioning, the data  $X_1, X_2, \dots, X_k$  can be recovered.

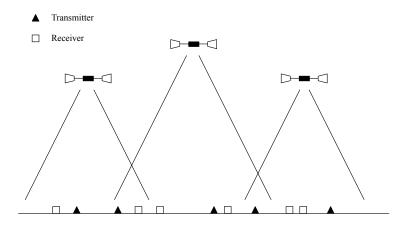


Fig. 21.6. A satellite communication network.

## 21.2.2 Satellite Communication Network

In a satellite communication network, a user is at any time covered by one or more satellites. A user can be a transmitter, a receiver, or both. Through the satellite network, each information source generated at a transmitter is multicast to a certain set of receivers. A transmitter can transmit to all the satellites within the line of sight, while a receiver can receive from all the satellites within the line of sight. Neighboring satellites may also communicate with each other. Figure 21.6 is an illustration of a satellite communication network.

The satellite communication network in Figure 21.6 can be represented by the graph in Figure 21.7 which consists of three layers of nodes. The top layer represents the transmitters, the middle layer consists of nodes representing the satellites as well as the associated dummy nodes modeling the broadcast nature of the satellites, and the bottom layer represents the receivers. If a satellite is within the line-of-sight of a transmitter, then the corresponding pair of nodes are connected by a directed edge. Likewise, if a receiver is within the line-of-sight of a satellite, then the corresponding pair nodes are connected by a directed edge. An edges between two nodes in the middle layer represent the communication links between two neighboring satellites. Each information source is multicast to a specified set of receiving nodes as shown.

## 21.3 A Network Code for Acyclic Networks

Let  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  denote an acyclic point-to-point communication network, where  $\mathcal{V}$  and  $\mathcal{E}$  are the set of nodes and the set of channels, respectively. We assume that each channel  $e \in \mathcal{E}$  is error-free with rate constraint  $R_e$ .

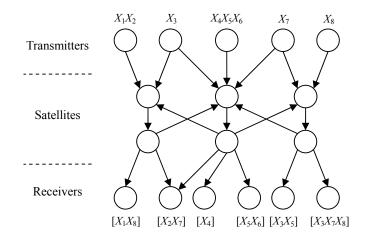


Fig. 21.7. A graph representing a satellite communication network.

As in our previous discussions, we let In(t) and Out(t) be the set of input channels and the set of output channels of node t, respectively.

Let  $S \subset V$  be the set of source nodes and  $T \subset V$  be the set of sink nodes. Without loss of generality, we assume  $\mathcal{G}$  has the structure that a source node has no input channel and a sink node has no output channel. Accordingly, Sand T are disjoint subsets of V.

An information source represented by a random variable  $X_s$  is generated at a source node  $s \in S$ , where  $X_s$  takes values in

$$\mathcal{X}_s = \{1, 2, \cdots, \lceil 2^{n\tau_s} \rceil\}$$
(21.6)

according to the uniform distribution, where  $\tau_s$  is the rate of the information source. The information sources  $X_s, s \in S$  are assumed to be mutually independent.

To simplify the notation, we will denote  $(X_s : s \in \mathcal{A})$  by  $X_{\mathcal{A}}$ ,  $\prod_{s \in \mathcal{A}} \mathcal{X}_s$  by  $\mathcal{X}_{\mathcal{A}}$ , etc. At a sink node  $t \in \mathcal{T}$ , the set of information sources  $X_{\beta(t)}$ , where  $\beta(t) \subset S$ , is received. We assume that each information source is received at at least one sink node, i.e., for every  $s \in \mathcal{S}$ ,  $s \in \beta(t)$  for some  $t \in \mathcal{T}$ . In the case when  $\beta(t) = S$  for all  $t \in \mathcal{T}$ , the problem is reduced to the single-source network coding problem.

**Definition 21.1.** An  $(n, (\eta_e : e \in \mathcal{E}), (\tau_s : s \in \mathcal{S}))$  block code of length n on a given communication network is defined by

1) for all source node  $s \in S$  and all channels  $e \in \text{Out}(s)$ , a local encoding function

$$k_e: \mathcal{X}_s \to \{0, 1, \cdots, \eta_e\}; \tag{21.7}$$

2) for all node  $i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T})$  and all channels  $e \in \text{Out}(i)$ , a local encoding function

$$k_e : \prod_{d \in \text{In}(i)} \{0, 1, \cdots, \eta_d\} \to \{0, 1, \cdots, \eta_e\};$$
(21.8)

3) for all sink node  $t \in \mathcal{T}$ , a decoding function

$$g_t: \prod_{d\in \mathrm{In}(t)} \{0, 1, \cdots, \eta_d\} \to \mathcal{X}_{\beta(t)}.$$
(21.9)

The nodes in  $\mathcal{V}$  are assumed to be ordered in an upstream-to-downstream manner as prescribed in Proposition 19.1. This defines a coding order among the nodes such that whenever a node encodes, all the information needed would have already been received on the input channels of that node.

For all sink node  $t \in \mathcal{T}$ , define

$$\Delta_t = \Pr\left\{\tilde{g}_t(X_{\mathcal{S}}) \neq X_{\beta(t)}\right\},\tag{21.10}$$

where  $\tilde{g}_t(X_S)$  denotes the value of  $g_t$  as a function of  $X_S$ .  $\Delta_t$  is the probability that the set of information sources  $X_{\beta(t)}$  is decoded correctly at sink node t.

Throughout this chapter, all the logarithms are in the base 2.

**Definition 21.2.** An information rate tuple  $\boldsymbol{\omega} = (\omega_s : s \in \mathcal{S})$ , where  $\boldsymbol{\omega} \geq 0$ (componentwise) is asymptotically achievable if for any  $\epsilon > 0$ , there exists for sufficient large n an  $(n, (\eta_e : e \in \mathcal{E}), (\tau_s : s \in \mathcal{S}))$  code such that

$$n^{-1}\log\eta_e \le R_e + \epsilon, \ e \in \mathcal{E} \tag{21.11}$$

$$\tau_s \ge \omega_s - \epsilon, \ s \in \mathcal{S} \tag{21.12}$$

$$\Delta_t \le \epsilon, \ t \in \mathcal{T}. \tag{21.13}$$

For brevity, an asymptotically achievable information rate tuple will be referred to as an achievable information rate tuple.

## 21.4 The Achievable Information Rate Region

In this section, we define the achievable information rate region and give a characterization of this region.

**Definition 21.3.** The achievable information rate region, denoted by  $\mathcal{R}$ , is the set of all achievable information rate tuple  $\boldsymbol{\omega}$ .

**Remark** It follows from the definition of the achievability of an information rate vector that if  $\boldsymbol{\omega}$  is achievable, then  $\boldsymbol{\omega}'$  is achievable for all  $0 \leq \boldsymbol{\omega}' \leq \boldsymbol{\omega}$ . Also, if  $\boldsymbol{\omega}^{(k)}, k \geq 1$  are achievable, then it can be proved by techniques similar to those in the proof of Theorem 8.12 that

$$\boldsymbol{\omega} = \lim_{k \to \infty} \boldsymbol{\omega}^{(k)} \tag{21.14}$$

is also achievable, i.e.,  $\mathcal{R}$  is closed. The details are omitted here.

Consider the set of all information rate tuples  $\omega$  such that there exist auxiliary random variables  $\{Y_s, s \in S\}$  and  $\{U_e, e \in \mathcal{E}\}$  which satisfy the following conditions:

$$H(Y_s) \ge \omega_s, \ s \in \mathcal{S} \tag{21.15}$$

$$H(Y_{\mathcal{S}}) = \sum_{s \in \mathcal{S}} H(Y_s) \tag{21.16}$$

$$H(U_{\text{Out}(s)}|Y_s) = 0, \ s \in \mathcal{S}$$

$$(21.17)$$

$$U_{\text{Out}(s)}|U_{\text{Out}(s)}| = 0, \ i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T})$$

$$(21.18)$$

$$H(U_{\text{Out}(i)}|U_{\text{In}(i)}) = 0, \ i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T})$$
(21.18)

$$H(U_e) \le R_e, \ e \in \mathcal{E} \tag{21.19}$$

$$H(Y_{\beta(t)}|U_{\text{In}(t)}) = 0, \ t \in \mathcal{T},$$
 (21.20)

where  $Y_{\mathcal{S}}$  denotes  $(Y_s : s \in \mathcal{S}), U_{\text{Out}(s)}$  denotes  $(U_e : e \in \text{Out}(s))$ , etc. Here,  $Y_s$  is an auxiliary random variable associated with the information source  $X_s$ , and  $U_e$  is an auxiliary random variable associated with the codeword sent on channel e. The interpretations of (21.15) to (21.20) are as follows. The inequality in (21.15) says that the entropy of  $Y_s$  is greater than or equal to  $\omega_s$ , the rate of the information source  $X_s$ . The equality in (21.16) says that  $Y_s, s \in \mathcal{S}$  are mutually independent, which corresponds to the assumption that the information sources  $X_s, s \in \mathcal{S}$  are mutually independent. The equality in (21.17) says that  $U_{Out(s)}$  is a function of  $Y_s$  for  $s \in S$ , and the equality in (21.18) says that  $U_{\text{Out}(i)}$  is a function of  $U_{\text{In}(i)}$  for  $i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T})$ . These correspond to the requirement that the codewords sent out by a source node sare functions of the information source  $X_s$ , and that the codewords sent out by a non-source node i are functions of the codewords received by node i. The inequality in (21.19) says that the entropy of  $U_e$  is less than or equal to  $R_e$ , the rate constraint for channel e. The equality in (21.20) says that  $Y_{\beta(t)}$  is a function of  $U_{\text{In}(t)}$  for  $t \in \mathcal{T}$ , which corresponds to the requirement that the information sources to be received at a sink node t can be decoded from the codewords received at node t.

For a given multi-source network coding problem, let

$$\mathcal{N} = \{Y_s : s \in \mathcal{S}; U_e : e \in \mathcal{E}\}$$
(21.21)

be a collection of discrete random variables whose joint distribution is unspecified, and let

$$\mathcal{Q}_{\mathcal{N}} = 2^{\mathcal{N}} \setminus \{\phi\} \tag{21.22}$$

with cardinality  $2^{|\mathcal{N}|} - 1$ . Let  $\mathcal{H}_{\mathcal{N}}$  be the  $|\mathcal{Q}_{\mathcal{N}}|$ -dimensional Euclidean space with the coordinates labeled by  $h_A, A \in \mathcal{Q}_N$ . A vector

$$\mathbf{h} = (h_A : A \in \mathcal{Q}_{\mathcal{N}}) \tag{21.23}$$

in  $\mathcal{H}_{\mathcal{N}}$  is said to be *finitely entropic* if there exists a joint distribution for all  $X \in \mathcal{N}$ , where  $|\mathcal{X}| < \infty$  for all  $X \in \mathcal{N}$  and

$$h_A = H(X : X \in A) \tag{21.24}$$

for all  $A \in \mathcal{Q}_{\mathcal{N}}$ . Note that  $\mathbf{h} \in \mathcal{H}_{\mathcal{N}}$  is entropic if it is finitely entropic, but not vice versa. We then define the region

$$\Gamma_{\mathcal{N}}^{**} = \{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : \mathbf{h} \text{ is finitely entropic} \}.$$
(21.25)

To simplify notation, for any nonempty  $A, A' \in \mathcal{Q}_{\mathcal{N}}$ , define

$$h_{A|A'} = h_{AA'} - h_{A'}, (21.26)$$

where we have used juxtaposition to denote the union of two sets. In using the above notation, we do not distinguish elements and singletons of  $\mathcal{N}$ , i.e., for a random variable  $Z \in \mathcal{N}$ ,  $h_Z$  is the same as  $h_{\{Z\}}$ . We now define the following regions in  $\mathcal{H}_{\mathcal{N}}$ :

$$\mathcal{C}_1 = \left\{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{Y_{\mathcal{S}}} = \sum_{s \in \mathcal{S}} h_{Y_s} \right\}$$
(21.27)

$$\mathcal{C}_2 = \left\{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{U_{\text{Out}(s)}|Y_s} = 0, s \in \mathcal{S} \right\}$$
(21.28)

$$\mathcal{C}_{3} = \left\{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{U_{\text{Out}(i)}|U_{\text{In}(i)}} = 0, i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T}) \right\}$$
(21.29)

$$\mathcal{C}_4 = \{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{U_e} \le R_e, e \in \mathcal{E} \}$$
(21.30)

$$\mathcal{C}_{5} = \left\{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{Y_{\beta(t)}|U_{\mathrm{In}(t)}} = 0, t \in \mathcal{T} \right\}.$$
(21.31)

Evidently, (21.27) to (21.31) are the regions in  $\mathcal{H}_{\mathcal{N}}$  corresponding to (21.16) to (21.20), respectively. We further denote  $\bigcap_{i \in \alpha} C_i$  by  $\mathcal{C}_{\alpha}$  for  $\alpha \subset \{1, 2, 3, 4, 5\}$ .

We now introduce a few notations. For a vector  $\mathbf{h} \in \mathcal{H}_{\mathcal{N}}$ , let  $\mathbf{h}_{Y_{\mathcal{S}}} = (h_{Y_s} : s \in \mathcal{S})$ . For a subset  $\mathcal{B}$  of  $\mathcal{H}_{\mathcal{N}}$ , let

- 1.  $\operatorname{proj}_{Y_{\mathcal{S}}}(\mathcal{B}) = {\mathbf{h}_{Y_{\mathcal{S}}} : \mathbf{h} \in \mathcal{B}}$  be the projection of the set  $\mathcal{B}$  on the coordinates  $h_{Y_s}, s \in \mathcal{S}$ ;
- 2.  $\Lambda(\mathcal{B}) = \{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : 0 \leq \mathbf{h} \leq \mathbf{h}' \text{ for some } \mathbf{h}' \in \mathcal{B} \};$
- 3.  $\operatorname{con}(\mathcal{B})$  be the convex hull of  $\mathcal{B}$ ;
- 4.  $\overline{\mathcal{B}}$  be the closure of  $\mathcal{B}$ .

Note that a vector  $\mathbf{h} \geq 0$  is in  $\Lambda(\mathcal{B})$  if and only if it is inferior to some vector  $\mathbf{h}'$  in  $\mathcal{B}$ . The following theorem gives a characterization of the achievable information rate region  $\mathcal{R}$  in terms of the region  $\Gamma_{\mathcal{N}}^{**}$ .

Definition 21.4. Define the region

$$\mathcal{R}' = \Lambda \left( \operatorname{proj}_{Y_{\mathcal{S}}} \left( \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_4 \cap \mathcal{C}_5 \right) \right).$$
(21.32)

Theorem 21.5.  $\mathcal{R} = \mathcal{R}'$ .

This theorem, which characterizes the achievable information rate region  $\mathcal{R}$ , will be proved in Sections 21.6 and 21.7. In the next section, we first discuss how more explicit inner and outer bounds on  $\mathcal{R}$  can be obtained.

## 21.5 Explicit Inner and Outer Bounds

Theorem 21.5 gives a characterization of the achievable information rate region  $\mathcal{R}$  in terms of the region  $\Gamma_{\mathcal{N}}^{**}$ . However, so far there exists no complete characterization of  $\Gamma_{\mathcal{N}}^{**}$ . Therefore, the region  $\mathcal{R}$  cannot be evaluated explicitly.

In the definition of  $\mathcal{R}'$  in (21.32), if  $\Gamma_{\mathcal{N}}^{**}$  is replaced by an inner bound (outer bound) on  $\Gamma_{\mathcal{N}}^{**}$ , then an inner bound (outer bound) on  $\mathcal{R}$  is obtained. The results in [235] and [233] which are beyond the scope of our discussion here, provide explicit constructions of inner bounds on  $\Gamma_{\mathcal{N}}^{**}$ .

We now discuss how an explicit outer bound on  $\mathcal{R}$  can be obtained. To facilitate our discussion, we further define

$$i_{A;A'} = h_A - h_{A|A'} \tag{21.33}$$

and

$$i_{A;A'|A''} = h_{A|A''} - h_{A|A'A''}$$
(21.34)

for  $A, A', A'' \in \mathcal{Q}_{\mathcal{N}}$ . Let  $\Gamma_{\mathcal{N}}$  be the set of  $\mathbf{h} \in \mathcal{H}_{\mathcal{N}}$  such that  $\mathbf{h}$  satisfies all the basic inequalities involving some or all of the random variables in  $\mathcal{N}$ , i.e., for all  $A, A', A'' \in \mathcal{Q}_{\mathcal{N}}$ ,

$$h_A \ge 0 \tag{21.35}$$

$$h_{A|A'} \ge 0 \tag{21.36}$$

$$i_{A;A'} \ge 0$$
 (21.37)

$$i_{A;A'|A''} \ge 0.$$
 (21.38)

We know from Section 14.2 that  $\overline{\Gamma}_{\mathcal{N}}^{**} \subset \Gamma_{\mathcal{N}}$ . Then upon replacing  $\overline{\Gamma}_{\mathcal{N}}^{**}$  by  $\Gamma_{\mathcal{N}}$  in (21.32), we obtain an outer bound on  $\mathcal{R}_{out}$ . This outer bound, called the *LP bound* (LP for linear programming), is given by

$$\mathcal{R}_{LP} = \Lambda \left( \operatorname{proj}_{Y_{\mathcal{S}}} \left( \overline{\operatorname{con}(\Gamma_{\mathcal{N}} \cap \mathcal{C}_{123})} \cap \mathcal{C}_{4} \cap \mathcal{C}_{5} \right) \right).$$
(21.39)

Since  $\mathcal{R}_{LP}$  involves only a finite number of linear constraints,  $\mathcal{R}_{LP}$  can be evaluated explicitly.

Using the technique in [368], it can be proved that  $\mathcal{R}_{LP}$  is tight for most special cases of multi-source network coding on an acyclic network for which the achievable information region is known. In addition to single-source network coding, these include the models described in [143] [359] [279] [367] [368] [308]. Since  $\mathcal{R}_{LP}$  encompasses all Shannon-type information inequalities and the converse proofs of the achievable information rate region for all these special cases do not involve non-Shannon-type inequalities, the tightness of  $\mathcal{R}_{LP}$ for all these cases is expected.

However, there exist multi-source network coding problems that requires non-Shannon-type inequalities for the characterization of the achievable information rate region [88] [59]. As new non-Shannon-type inequalities are discovered from time to time, improved outer bounds on  $\mathcal{R}$  can be obtained by incorporating these inequalities.

## 21.6 The Converse

In this section, we establish the converse part of Theorem 21.5, namely

$$\mathcal{R} \subset \Lambda \left( \operatorname{proj}_{Y_{\mathcal{S}}} \left( \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_{4} \cap \mathcal{C}_{5} \right) \right) = \mathcal{R}'.$$
(21.40)

Let  $\epsilon_k$  be a sequence such that  $0 < \epsilon_k < 1$  for all k and  $\epsilon_k$  monotonically decreases to 0 as  $k \to \infty$ . Consider an achievable information rate tuple  $\omega \in \mathcal{R}$ . Then for all k, for all sufficiently large n, there exists an

$$\left(n, (\eta_e^{(k)} : e \in \mathcal{E}), (\tau_s^{(k)} : s \in \mathcal{S})\right)$$
(21.41)

code satisfying

$$n^{-1}\log\eta_e^{(k)} \le R_e + \epsilon_k, \ e \in \mathcal{E}$$
(21.42)

$$\tau_s^{(k)} \ge \omega_s - \epsilon_k, \ s \in \mathcal{S} \tag{21.43}$$

$$\Delta_t^{(k)} \le \epsilon_k, \ t \in \mathcal{T},\tag{21.44}$$

where  $\Delta_t^{(k)}$  denotes the decoding error probability at sink node t (cf. (21.10)).

We now fix k to be any positive integer and temporarily suppress all the superscripts involving k. For all  $e \in \mathcal{E}$ , let  $U_e$  be the codeword sent on channel e and denote the alphabet of  $U_e$  by  $\mathcal{U}_e$ . The following lemma, whose proof will be deferred to the end of the section, is a consequence of Fano's inequality.

**Lemma 21.6.** For all n and k, for all  $t \in \mathcal{T}$ ,

$$H(X_{\beta(t)}|U_{\text{In}(t)}) \le n\phi_t(n,\epsilon_k), \qquad (21.45)$$

where

1.  $\phi_t(n, \epsilon_k)$  is bounded; 2.  $\phi_t(n, \epsilon_k) \to 0$  as  $n, k \to \infty$ ; 3.  $\phi_t(n, \epsilon_k)$  is monotonically decreasing in both n and k.

Since the information source  $X_s, s \in \mathcal{S}$  are mutually independent,

$$H(X_{\mathcal{S}}) = \sum_{s \in \mathcal{S}} H(X_s).$$
(21.46)

For any  $s \in \mathcal{S}$ , since  $U_{\text{Out}(s)}$  is a function of  $X_s$ ,

$$H(U_{\text{Out}(s)}|X_s) = 0.$$
 (21.47)

Similarly, for all  $i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T})$ ,

$$H(U_{\text{Out}(i)}|X_{\text{In}(i)}) = 0.$$
 (21.48)

For all  $e \in \mathcal{E}$ ,

$$H(U_e) \le \log |\mathcal{U}_e| \tag{21.49}$$

$$= \log(\eta_e + 1) \tag{21.50}$$

$$\leq n(R_e + 2\epsilon_k) \tag{21.51}$$

where (21.51) follows from (21.42) assuming that n is sufficiently large. For all  $t \in \mathcal{T}$ , from Lemma 21.6, we have

$$H(X_{\beta(t)}|U_{\mathrm{In}(t)}) \le n\phi_t(n,\epsilon_k).$$
(21.52)

For all  $s \in \mathcal{S}$ , from (21.43),

$$H(X_s) = \log |\mathcal{X}_s| = \log \lceil 2^{n\tau_s} \rceil \ge n\tau_s \ge n(\omega_s - \epsilon_k).$$
(21.53)

By letting  $Y_s = X_s$  for all  $s \in S$ , we then obtain from (21.46) to (21.48) and (21.51) to (21.53) that

$$H(Y_{\mathcal{S}}) = \sum_{s \in \mathcal{S}} H(Y_s) \tag{21.54}$$

$$H(U_{\text{Out}(s)}|Y_s) = 0, \ s \in \mathcal{S}$$

$$(21.55)$$

$$H(U_{\text{Out}(i)}|U_{\text{In}(i)}) = 0, \ i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T})$$
(21.56)

$$H(U_e) \le n(R_e + 2\epsilon_k), \ e \in \mathcal{E}$$
(21.57)

$$H(Y_{\beta(t)}|U_{\mathrm{In}(t)}) \le n\phi_t(n,\epsilon_k), \ t \in \mathcal{T}$$
(21.58)

$$H(Y_s) \ge n(\omega_s - \epsilon_k), \ s \in \mathcal{S}.$$
(21.59)

Now define the following two regions in  $\mathcal{H}_{\mathcal{N}}$ :

$$\mathcal{C}_{4,\epsilon_k}^n = \{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{U_e} \le n(R_e + 2\epsilon_k), e \in \mathcal{E} \}$$
(21.60)

$$\mathcal{C}^n_{5,\epsilon_k} = \{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{Y_{\beta(t)}|U_{\mathrm{In}(t)}} \le n\phi_t(n,\epsilon_k), t \in \mathcal{T} \}.$$
(21.61)

Note that all the auxiliary random variables  $Y_s, s \in S$  and  $U_e, e \in E$  have finite alphabets, because

$$|\mathcal{Y}_s| = |\mathcal{X}_s| = \lceil 2^{n\tau_s} \rceil < \infty \tag{21.62}$$

and

$$\log |\mathcal{U}_e| \le n(R_e + 2\epsilon_k) < \infty \tag{21.63}$$

(cf. (21.49) through (21.51)). Then we see from (21.54) to (21.59) that there exists

$$\mathbf{h}^{(k)} \in \Gamma_{\mathcal{N}}^{**} \tag{21.64}$$

such that

$$\mathbf{h}^{(k)} \in \mathcal{C}_{123} \cap \mathcal{C}^n_{4,\epsilon_k} \cap \mathcal{C}^n_{5,\epsilon_k} \tag{21.65}$$

and

$$h_{Y_s}^{(k)} \ge n(\omega_s - \epsilon_k) \tag{21.66}$$

for all  $s \in S$ . From (21.64) and (21.65), we obtain

$$\mathbf{h}^{(k)} \in \Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123} \cap \mathcal{C}_{4,\epsilon_k}^n \cap \mathcal{C}_{5,\epsilon_k}^n.$$
(21.67)

Upon dividing by n, (21.66) becomes

$$n^{-1}h_{Y_s}^{(k)} \ge \omega_s - \epsilon_k. \tag{21.68}$$

Since  $\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}$  contains the origin in  $\mathcal{H}_{\mathcal{N}}$ , we see that

$$n^{-1}\mathbf{h}^{(k)} \in \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_{4,\epsilon_k} \cap \mathcal{C}_{5,\epsilon_k},$$
(21.69)

where

$$\mathcal{C}_{4,\epsilon_k} = \{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{U_e} \le R_e + 2\epsilon_k, e \in \mathcal{E} \}$$
(21.70)

$$\mathcal{C}_{5,\epsilon_k} = \{ \mathbf{h} \in \mathcal{H}_{\mathcal{N}} : h_{Y_{\beta(t)}|U_{\mathrm{In}(t)}} \le \phi_t(n,\epsilon_k), t \in \mathcal{T} \}.$$
(21.71)

Note that the region  $C_{5,\epsilon_k}$  depends on n though it is not indicated explicitly. For all n and k, define the set

$$\mathcal{B}^{(n,k)} = \{ \mathbf{h} \in \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_{4,\epsilon_k} \cap \mathcal{C}_{5,\epsilon_k} : \\ h_{Y_s} \ge \omega_s - \epsilon_k \text{ for all } s \in \mathcal{S} \}.$$
(21.72)

**Lemma 21.7.** For all n and k, the set  $\mathcal{B}^{(n,k)}$  is compact<sup>1</sup>.

Again, the proof of this lemma is deferred to the end of the section. Now from Lemma 21.6,  $\phi_t(n, \epsilon_k)$  is monotonically decreasing in both n and k, so for all n and k,

$$\mathcal{B}^{(n+1,k)} \subset \mathcal{B}^{(n,k)} \tag{21.73}$$

and

$$\mathcal{B}^{(n,k+1)} \subset \mathcal{B}^{(n,k)}.$$
(21.74)

For any fixed k and all sufficiently large n, we see from (21.68) and (21.69) that  $\mathcal{B}^{(n,k)}$  is nonempty. Since  $\mathcal{B}^{(n,k)}$  is compact by Lemma 21.7,

$$\lim_{n \to \infty} \mathcal{B}^{(n,k)} = \bigcap_{n=1}^{\infty} \mathcal{B}^{(n,k)}$$
(21.75)

is both compact and nonempty. By the same argument, we conclude that

$$\lim_{k \to \infty} \lim_{n \to \infty} \mathcal{B}^{(n,k)} = \bigcap_{k=1}^{\infty} \bigcap_{n=1}^{\infty} \mathcal{B}^{(n,k)}$$
(21.76)

is also nonempty. Now the set

<sup>&</sup>lt;sup>1</sup> A subset of the Euclidean space is compact if and only if it is closed and bounded.

21.6 The Converse 509

$$\lim_{k \to \infty} \lim_{n \to \infty} \mathcal{B}^{(n,k)} \tag{21.77}$$

is equal to

$$\left\{ \mathbf{h} \in \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_4 \cap \mathcal{C}_5 : h_{Y_s} \ge \omega_s \text{ for all } s \in \mathcal{S} \right\}.$$
 (21.78)

Hence, there exists  $\mathbf{h}'$  satisfying

$$\mathbf{h}' \in \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_4 \cap \mathcal{C}_5$$
(21.79)

and

$$h'_{Y_s} \ge \omega_s, \ s \in \mathcal{S}. \tag{21.80}$$

Let  $\mathbf{r} = \operatorname{proj}_{Y_{\mathcal{S}}}(\mathbf{h}')$ . Then we have

$$\mathbf{r} \in \operatorname{proj}_{Y_{\mathcal{S}}} \left( \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_{4} \cap \mathcal{C}_{5} \right)$$
(21.81)

and

$$\mathbf{r} \ge \boldsymbol{\omega}$$
 (21.82)

componentwise. By (21.81) and (21.82), we finally conclude that

$$\boldsymbol{\omega} \in \Lambda\left(\operatorname{proj}_{Y_{\mathcal{S}}}\left(\overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_{4} \cap \mathcal{C}_{5}\right)\right).$$
(21.83)

This completes the proof of the converse part of Theorem 21.5.

Proof of Lemma 21.6. For any  $t \in \mathcal{T}$ , by Fano's inequality, we have

$$H(X_{\beta(t)}|U_{\text{In}(t)}) \le 1 + \Delta_t \log |\mathcal{X}_{\beta(t)}|$$
(21.84)

$$= 1 + \Delta_t H(X_{\beta(t)}) \tag{21.85}$$

$$\leq 1 + \epsilon_k H(X_{\beta(t)}), \tag{21.86}$$

where (21.85) follows because  $X_s$  distributes uniformly on  $\mathcal{X}_s$  and  $X_s$ ,  $s \in \mathcal{S}$  are mutually independent, and (21.86) follows from (21.44). Then

$$H(X_{\beta(t)}) = I(X_{\beta(t)}; U_{\text{In}(t)}) + H(X_{\beta(t)}|U_{\text{In}(t)})$$
(21.87)  
a)

$$\leq I(X_{\beta(t)}; U_{\text{In}(t)}) + 1 + \epsilon_k H(X_{\beta(t)})$$
(21.88)

$$\leq H(U_{\mathrm{In}(t)}) + 1 + \epsilon_k H(X_{\beta(t)}) \tag{21.89}$$

<sup>b)</sup>

$$\leq \left(\sum_{e \in \text{In}(t)} \log \eta_e\right) + 1 + \epsilon_k H(X_{\beta(t)}) \tag{21.90}$$

$$\stackrel{c)}{\leq} \left( \sum_{e \in \text{In}(t)} n(R_e + \epsilon_k) \right) + 1 + \epsilon_k H(X_{\beta(t)}), \quad (21.91)$$

where

- a) follows from (21.86);
- b) follows from Theorem 2.43;
- c) follows from (21.42).

Rearranging the terms in (21.91), we obtain

$$H(X_{\beta(t)}) \le \frac{n}{1 - \epsilon_k} \left( \sum_{e \in \operatorname{In}(t)} (R_e + \epsilon_k) + \frac{1}{n} \right).$$
 (21.92)

Substituting (21.92) into (21.86), we have

$$H(X_{\beta(t)}|U_{\mathrm{In}(t)}) < n \left[ \frac{1}{n} + \frac{\epsilon_k}{1 - \epsilon_k} \left( \sum_{e \in \mathrm{In}(t)} (R_e + \epsilon_k) + \frac{1}{n} \right) \right]$$
(21.93)  
=  $n\phi_t(n, \epsilon_k),$  (21.94)

$$= n\phi_t(n,\epsilon_k), \tag{21.94}$$

where

$$\phi_t(n,\epsilon_k) = \frac{1}{n} + \frac{\epsilon_k}{1-\epsilon_k} \left( \sum_{e \in \text{In}(t)} (R_e + \epsilon_k) + \frac{1}{n} \right).$$
(21.95)

Invoking the assumption that  $0 < \epsilon_k < 1$  for all k and  $\epsilon_k$  monotonically decreases to 0 as  $k \to \infty$ , it is evident that

- 1.  $\phi_t(n, \epsilon_k)$  is bounded for all n and k;
- 2.  $\phi_t(n, \epsilon_k) \to 0$  as  $n, k \to \infty$ ;

3.  $\phi_t(n, \epsilon_k)$  is monotonically nonincreasing in both n and k.

The lemma is proved.  $\Box$ 

Proof of Lemma 21.7. We need to show that the set  $\mathcal{B}^{(n,k)}$  is both closed and bounded. The closedness of  $\mathcal{B}^{(n,k)}$  is immediate from its definition. To establish the boundedness of  $\mathcal{B}^{(n,k)}$ , we need to show that for any  $\mathbf{h} \in \mathcal{B}^{(n,k)}$ , all the components of **h** are bounded. Consider any  $\mathbf{h} \in \mathcal{B}^{(n,k)}$ . Since

$$\mathcal{B}^{(n,k)} \subset \mathcal{C}_{4,\epsilon_k},\tag{21.96}$$

we see from (21.70) that  $h_{U_e}$  are bounded for all  $e \in \mathcal{E}$ . Since

$$\mathcal{B}^{(n,k)} \subset \mathcal{C}_{5,\epsilon_k},\tag{21.97}$$

we see from (21.71) that for every  $t \in \mathcal{T}$ ,

$$h_{Y_{\beta(t)}} \le h_{Y_{\beta(t)}U_{\text{In}(t)}}$$
 (21.98)  
= h (21.00)

$$=h_{Y_{\beta(t)}|U_{\text{In}(t)}} + h_{U_{\text{In}(t)}}$$
(21.99)

$$\leq \phi_t(n,\epsilon_k) + h_{U_{\text{In}(t)}} \tag{21.100}$$

$$\leq \phi_t(n,\epsilon_k) + \sum_{e \in \text{In}(t)} h_{U_e} \tag{21.101}$$

where (21.99) and the boundedness of  $\phi_t(n, \epsilon_k)$  follow from Lemma 21.6. This shows that  $h_{Y_{\beta(t)}}$  is bounded for all  $t \in \mathcal{T}$ .

In our model, for every  $s \in S$ , there exists at least one  $t \in T$  such that  $s \in \beta(t)$ . Then the boundedness of  $h_{Y_{\beta(t)}}$  for all  $t \in T$  implies the boundedness of  $h_{Y_s}$  for all  $s \in S$ . Finally, the boundedness of all the other components of **h** is established by invoking the independence bound for entropy. The lemma is proved.  $\Box$ 

## 21.7 Achievability

In this section, we establish the direct part of Theorem 21.5, namely,

$$\mathcal{R}' = \Lambda \left( \operatorname{proj}_{Y_{\mathcal{S}}} \left( \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_4 \cap \mathcal{C}_5 \right) \right) \subset \mathcal{R}.$$
(21.102)

Before we proceed, we first prove an alternative form of  $\mathcal{R}'$  that will be used in constructing the random code. For a subset  $\mathcal{B}$  of  $\mathcal{H}_{\mathcal{N}}$ , let

$$D(\mathcal{B}) = \{ \alpha \mathbf{h} : \mathbf{h} \in \mathcal{H}_{\mathcal{N}} \text{ and } 0 \le \alpha \le 1 \}.$$
(21.103)

Define the two subsets

$$\mathcal{A}_1 = \overline{\operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \tag{21.104}$$

and

$$\mathcal{A}_2 = \overline{D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \tag{21.105}$$

of  $\mathcal{H}_{\mathcal{N}}$ .

## Lemma 21.8. $A_1 = A_2$ .

*Proof.* Since the origin of  $\mathcal{H}_{\mathcal{N}}$  is in  $\Gamma_{\mathcal{N}}^{**}$ , it is also in  $\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}$  because  $\mathcal{C}_{123}$  is a linear subspace of  $\mathcal{H}_{\mathcal{N}}$ . Upon observing that for  $0 \leq \alpha \leq 1$ ,

$$\alpha \mathbf{h} = (1 - \alpha)\mathbf{0} + \alpha \mathbf{h} \tag{21.106}$$

is a convex combination of 0 and  $\mathbf{h}$ , we obtain

$$D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}) \subset \operatorname{con}(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}).$$
(21.107)

It follows that  $\mathcal{A}_2 \subset \mathcal{A}_1$ .

To prove that  $\mathcal{A}_1 \subset \mathcal{A}_2$ , it suffices to show that  $\mathcal{A}_2$  is convex because

- 1.  $(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}) \subset \mathcal{A}_2$ , where  $\mathcal{A}_2$  is closed;
- 2.  $\mathcal{A}_1$  is the smallest closed convex set containing  $\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}$ .

Toward this end, consider any  $\mathbf{h}_1, \mathbf{h}_2 \in \mathcal{A}_2$  and any  $0 \leq \lambda \leq 1$ . We will show that

$$\mathbf{h} = \lambda \mathbf{h}_1 + (1 - \lambda) \mathbf{h}_2 \in \mathcal{A}_2. \tag{21.108}$$

Here, we can assume without loss of generality that  $\mathbf{h}_1, \mathbf{h}_2 \neq 0$ , because otherwise (21.108) holds by the definition of  $\mathcal{A}_2$ . Since  $\mathbf{h}_1, \mathbf{h}_2 \in \mathcal{A}_2$ , there exist  $\mathbf{h}_1^k, \mathbf{h}_2^k \in D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})$  such that  $\mathbf{h}_1^k \to \mathbf{h}_1$  and  $\mathbf{h}_2^k \to \mathbf{h}_2$ . Since  $\mathbf{h}_1^k, \mathbf{h}_2^k \in$  $D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})$ ,

$$\mathbf{h}_1^k = \alpha_1^k \hat{\mathbf{h}}_1^k \tag{21.109}$$

and

$$\mathbf{h}_2^k = \alpha_2^k \hat{\mathbf{h}}_2^k, \qquad (21.110)$$

where  $\hat{\mathbf{h}}_{1}^{k}, \hat{\mathbf{h}}_{2}^{k} \in \Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}$  and  $0 < \alpha_{1}^{k}, \alpha_{2}^{k} \leq 1$ . Note that  $\alpha_{1}^{k}$  and  $\alpha_{2}^{k}$  are strictly positive because  $\mathbf{h}_{1}, \mathbf{h}_{2} \neq 0$ . Now let  $n_{1}^{k}$  and  $n_{2}^{k}$  be integer sequences such that

$$\frac{n_1^k \alpha_1^k}{n_2^k \alpha_2^k} \to \frac{\lambda}{1-\lambda},\tag{21.111}$$

and let

$$\hat{\mathbf{h}}^k = n_1^k \mathbf{h}_1^k + n_2^k \mathbf{h}_2^k.$$
(21.112)

By Lemma 15.3 (also Corollary 15.4),

$$\hat{\mathbf{h}}^k \in \Gamma_{\mathcal{N}}^{**}.\tag{21.113}$$

Furthermore, since  $\mathbf{h}_1^k, \mathbf{h}_2^k \in \mathcal{C}_{123}$  and  $\mathcal{C}_{123}$  is a linear subspace,  $\hat{\mathbf{h}}^k \in \mathcal{C}_{123}$ . Therefore,

$$\mathbf{\hat{h}}^k \in \Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}.$$
(21.114)

Let

$$\mathbf{h}^{k} = \frac{\hat{\mathbf{h}}^{k}}{n_{1}^{k}\alpha_{1}^{k} + n_{2}^{k}\alpha_{2}^{k}}.$$
(21.115)

Then

$$\mathbf{h}^{k} \in D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}) \subset \overline{D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} = \mathcal{A}_{2}.$$
 (21.116)

Substituting (21.112), (21.109), and (21.110) into (21.115), we obtain

$$\mathbf{h}^{k} = \frac{n_{1}^{k}\alpha_{1}^{k}}{n_{1}^{k}\alpha_{1}^{k} + n_{2}^{k}\alpha_{2}^{k}}\mathbf{h}_{1}^{k} + \frac{n_{2}^{k}\alpha_{2}^{k}}{n_{1}^{k}\alpha_{1}^{k} + n_{2}^{k}\alpha_{2}^{k}}\mathbf{h}_{2}^{k}.$$
 (21.117)

It can readily be seen from (21.111) that

$$\frac{n_1^k \alpha_1^k}{n_1^k \alpha_1^k + n_2^k \alpha_2^k} \to \lambda \tag{21.118}$$

and

$$\frac{n_2^k \alpha_2^k}{n_1^k \alpha_1^k + n_2^k \alpha_2^k} \to 1 - \lambda.$$
 (21.119)

Since  $\mathbf{h}_1^k \to \mathbf{h}_1$  and  $\mathbf{h}_2^k \to \mathbf{h}_2$ , we see from (21.108) that  $\mathbf{h}^k \to \mathbf{h}$ . Finally, since  $\mathbf{h}^k \in \mathcal{A}_2$  and  $\mathcal{A}_2$  is closed, we conclude that  $\mathbf{h} \in \mathcal{A}_2$ . Therefore,  $\mathcal{A}_2$  is convex, and hence  $\mathcal{A}_1 \subset \mathcal{A}_2$ . The lemma is proved.  $\Box$ 

21.7 Achievability 513

By virtue of this lemma, we can write

$$\mathcal{R}' = \Lambda \left( \operatorname{proj}_{Y_{\mathcal{S}}} \left( \overline{D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_4 \cap \mathcal{C}_5 \right) \right),$$
(21.120)

and we will establish  $\mathcal{R}' \subset \mathcal{R}$  by proving that

$$\Lambda\left(\operatorname{proj}_{Y_{\mathcal{S}}}\left(\overline{D(\Gamma_{\mathcal{N}}^{**}\cap\mathcal{C}_{123})}\cap\mathcal{C}_{4}\cap\mathcal{C}_{5}\right)\right)\subset\mathcal{R}.$$
(21.121)

By the remark following Definition 21.3, we only need to show the achievability of the region

$$\operatorname{proj}_{Y_{\mathcal{S}}}\left(\overline{D(\Gamma_{\mathcal{N}}^{**}\cap\mathcal{C}_{123})}\cap\mathcal{C}_{4}\cap\mathcal{C}_{5}\right).$$
(21.122)

Consider any  $\boldsymbol{\omega}$  in this region. Then there exists

$$\mathbf{h} \in \overline{D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})} \cap \mathcal{C}_4 \cap \mathcal{C}_5 \tag{21.123}$$

such that

$$\boldsymbol{\omega} = \operatorname{proj}_{Y_{\mathcal{S}}}(\mathbf{h}). \tag{21.124}$$

Since

$$\mathbf{h} \in \overline{D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123})},\tag{21.125}$$

there exist a sequence

$$\mathbf{h}^{(k)} \in D(\Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}) \tag{21.126}$$

such that

$$\mathbf{h} = \lim_{k \to \infty} \mathbf{h}^{(k)}.\tag{21.127}$$

Let

$$\boldsymbol{\omega}^{(k)} = \operatorname{proj}_{Y_{\mathcal{S}}}(\mathbf{h}^{(k)}). \tag{21.128}$$

It then follows from (21.127) that

$$\lim_{k \to \infty} \boldsymbol{\omega}^{(k)} = \boldsymbol{\omega}.$$
 (21.129)

By (21.126),

$$\mathbf{h}^{(k)} = \alpha^{(k)} \hat{\mathbf{h}}^{(k)} \tag{21.130}$$

where  $\hat{\mathbf{h}}^{(k)} \in \Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}$  and

$$0 \le \alpha^{(k)} \le 1.$$
 (21.131)

Note that  $\hat{\mathbf{h}}^{(k)}$  is an entropy function because it is in  $\Gamma_{\mathcal{N}}^{**}$ , but  $\mathbf{h}^{(k)}$  and  $\mathbf{h}$  are not necessarily entropy functions. Since  $\hat{\mathbf{h}}^{(k)} \in \Gamma_{\mathcal{N}}^{**} \cap \mathcal{C}_{123}$ , there exists a collection of random variables with finite alphabets

$$\mathcal{N}^{(k)} = \left\{ (Y_s^{(k)} : s \in \mathcal{S}), (U_e^{(k)} : e \in \mathcal{E}) \right\}$$
(21.132)

such that

$$\alpha^{(k)}H\left(Y_s^{(k)}\right) = \omega_s^{(k)}, \ s \in \mathcal{S}$$
(21.133)

$$H\left(Y_{\mathcal{S}}^{(k)}\right) = \sum_{s \in \mathcal{S}} H\left(Y_{s}^{(k)}\right)$$
(21.134)

$$H\left(U_{\text{Out}(s)}^{(k)} \middle| Y_s^{(k)}\right) = 0, \ s \in \mathcal{S}$$
(21.135)

$$H\left(U_{\text{Out}(i)}^{(k)} \middle| U_{\text{In}(i)}^{(k)}\right) = 0, \ i \in \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{T}),$$
(21.136)

where (21.133) is implied by (21.128). Furthermore, since  $\mathbf{h} \in \mathcal{C}_4 \cap \mathcal{C}_5$ , it follows from (21.127) and (21.130) that

$$\alpha^{(k)} H\left(U_e^{(k)}\right) \le R_e + \mu^{(k)}, \ e \in \mathcal{E}$$
(21.137)

$$\alpha^{(k)} H\left(\left. Y_{\beta(t)}^{(k)} \right| U_{\mathrm{In}(t)}^{(k)} \right) \le \gamma^{(k)}, \ t \in \mathcal{T},$$
(21.138)

where  $\mu^{(k)}, \gamma^{(k)} \to 0$  as  $k \to \infty$ . In the rest of the section, we will prove the achievability of  $\omega^{(k)}$  for all sufficiently large k. Then the closedness of  $\mathcal{R}$ implies the achievability of  $\omega$  by the remark following Definition 21.3.

#### 21.7.1 Random Code Construction

Fix k and  $\epsilon > 0$ , and let  $\delta$  be a small positive quantity to be specified later. We first construct a random

$$(n, (\eta_e^{(k)} : e \in \mathcal{E}), (\tau_s^{(k)} : s \in \mathcal{S}))$$
 (21.139)

code with

$$\eta_e^{(k)} \le 2^{n(\alpha^{(k)}H(U_e^{(k)}) + \psi_e^{(k)})} \tag{21.140}$$

for all  $e \in \mathcal{E}$  and

$$\omega_s^{(k)} - \frac{\epsilon}{2} \le \tau_s^{(k)} \le \omega_s^{(k)} - \frac{\epsilon}{3}, \qquad (21.141)$$

where  $\psi_e^{(k)} > 0$  and  $\psi_e^{(k)} \to 0$  as  $\delta \to 0$ , by the steps below. For the sake of simplicity, we temporarily suppress all the superscripts involving k.

1. Let

$$\hat{n} = \lceil n\alpha \rceil. \tag{21.142}$$

Here n is the block length of the random code we will construct, while  $\hat{n}$  is the length of a sequence of the typical sets that we will use for constructing the random code. For each source  $s \in S$ , let

$$\theta_s = \lceil 2^{n\tau_s} \rceil \tag{21.143}$$

and construct a codebook  $C_s$  by generating  $\theta_s$  codewords in  $\mathcal{Y}_s^{\hat{n}}$  randomly and independently according to  $p^{\hat{n}}(y_s)$ . Denote these sequences by  $\mathbf{Y}_s(1), \mathbf{Y}_s(2), \cdots, \mathbf{Y}_s(\theta_s)$ , and let  $\mathbf{Y}_s(0)$  be an arbitrary constant sequence in  $\mathcal{Y}_s^{\hat{n}}$ .

- 2. Reveal the codebook  $C_s, s \in S$  to all the nodes in the network.
- 3. At a source node  $s \in S$ , the information source  $X_s$  is generated according to the uniform distribution on

$$\mathcal{X}_s = \{1, 2, \cdots, \theta_s\}.$$
 (21.144)

4. Let  $T^{\hat{n}}_{[U_e]\delta}$  denote the set of strongly typical sequences<sup>2</sup> with respect to the distribution  $p(u_e)$ . Let

$$\zeta_e = |T^{\hat{n}}_{[U_e]\delta}|. \tag{21.145}$$

By the strong AEP and (21.142),

$$\zeta_e \le 2^{\hat{n}(H(U_e) + \psi_e/(2\alpha))} \le 2^{n(\alpha H(U_e) + \psi_e/2)}, \tag{21.146}$$

where  $\psi_e \to 0$  as  $\delta \to 0$ . For all channels  $e \in \mathcal{E}$ , choose an  $\eta_e$  satisfying

$$2^{n(\alpha H(U_e) + \psi_e/2)} \le \eta_e \le 2^{n(\alpha H(U_e) + \psi_e)}.$$
(21.147)

Denote the sequences in  $T^{\hat{n}}_{[U_e]\delta}$  by  $\mathbf{U}_e(1), \mathbf{U}_e(2), \cdots, \mathbf{U}_e(\zeta_e)$ , and let  $\mathbf{U}_e(0)$  be an arbitrary constant sequence in  $\mathcal{U}_e^{\hat{n}}$ .

a) Let the outcome of  $X_s$  be  $x_s$  for a source node s. For a channel  $e \in Out(s)$ , define the local encoding function

$$k_e: \mathcal{X}_s \to \{0, 1, \cdots, \eta_e\}$$
(21.148)

as follows. By (21.135), for each channel  $e \in \text{Out}(s)$ , there exists a function  $u_e$  such that

$$U_e = u_e(Y_s),$$
 (21.149)

i.e.,

$$\Pr\{U_e = u_e(y) | Y_s = y\} = 1 \tag{21.150}$$

for all  $y \in \mathcal{Y}_s$ . By the preservation property of strong typicality (Theorem 6.8), if

$$\mathbf{Y}_s(x_s) \in T^{\hat{n}}_{[Y_s]\delta},\tag{21.151}$$

then

$$u_e(\mathbf{Y}_s(x_s)) \in T^{\hat{n}}_{[U_e]\delta}, \qquad (21.152)$$

where in  $u_e(\mathbf{Y}_s(x_s))$ , the function  $u_e$  is applied to  $\mathbf{Y}_s(x_s)$  componentwise. If so, let  $k_e(x_s)$  be the index of  $u_e(\mathbf{Y}_s(x_s))$  in  $T^{\hat{n}}_{[U_e]\delta}$ , i.e.,

$$\mathbf{U}_e(k_e(x_s)) = u_e(\mathbf{Y}_s(x_s)).$$
 (21.153)

Otherwise, let  $k_e(x_s)$  be 0. Note that  $k_e$  is well-defined because

1

$$\zeta_e \le \eta_e \tag{21.154}$$

by (21.146) and (21.147).

 $<sup>^2</sup>$  Strong typicality applies because all the random variables in  $\mathcal{N}^{(k)}$  have finite alphabets.

b) Let  $C_e$  be the index sent on channel e. For a channel  $e \in \text{Out}(i)$ , where  $i \in \mathcal{V} \setminus (S \cup T)$ , define the local encoding function

$$k_e : \prod_{d \in \text{In}(i)} \{0, 1, \cdots, \eta_d\} \to \{0, 1, \cdots, \eta_e\}$$
(21.155)

as follows. By (21.136), there exists a function  $u_e$  such that

$$U_e = u_e(U_{\text{In}(i)}).$$
 (21.156)

With a slight abuse of notation, we write

$$\mathbf{U}_{\mathcal{E}'}(C_{\mathcal{E}'}) = (\mathbf{U}_d(C_d) : d \in \mathcal{E}')$$
(21.157)

for  $\mathcal{E}' \subset \mathcal{E}$ , and

$$\mathbf{Y}_{\mathcal{S}'}(x_{\mathcal{S}'}) = (\mathbf{Y}_s(x_s) : s \in \mathcal{S}')$$
(21.158)

for  $\mathcal{S}' \subset \mathcal{S}$ . By the preservation property of strong typicality, if

$$\mathbf{U}_{\mathrm{In}(i)}(C_{\mathrm{In}(i)}) \in T^{\dot{n}}_{[U_{\mathrm{In}(i)}]\delta},\tag{21.159}$$

then

$$u_e(\mathbf{U}_{\mathrm{In}(i)}(C_{\mathrm{In}(i)})) \in T^{\hat{n}}_{[U_e]\delta}.$$
(21.160)

If so, let  $k_e(C_{\text{In}(i)})$  be the index of  $u_e(\mathbf{U}_{\text{In}(i)}(C_{\text{In}(i)}))$  in  $T^{\hat{n}}_{[U_e]\delta}$ , i.e.,

$$\mathbf{U}_{e}(k_{e}(C_{\mathrm{In}(i)})) = u_{e}(\mathbf{U}_{\mathrm{In}(i)}(C_{\mathrm{In}(i)})).$$
 (21.161)

Otherwise, let  $k_e(C_{\text{In}(i)})$  be 0. Again,  $k_e$  is well-defined because (21.154) holds.

5. For a sink node  $t \in \mathcal{T}$ , define the decoding function

$$g_t: \prod_{d\in \mathrm{In}(t)} \{0, 1, \cdots, \eta_d\} \to \mathcal{X}_{\beta(t)}$$
(21.162)

as follows. If the received index  $C_d$  on channel d is nonzero for all  $d \in In(t)$ and there exists a unique tuple

$$x_{\beta(t)} \in \mathcal{X}_{\beta(t)} \tag{21.163}$$

such that

$$(\mathbf{Y}_{\beta(t)}(x_{\beta(t)}), \mathbf{U}_{\mathrm{In}(i)}(C_{\mathrm{In}(t)})) \in T^{\hat{n}}_{[U_{\mathrm{In}(t)}Y_{\beta(t)}]\delta},$$
(21.164)

then let  $g_t(C_{\text{In}(t)})$  be  $x_{\beta(t)}$ . Otherwise, declare a decoding error.

#### 21.7.2 Performance Analysis

Let us reinstate all the superscripts involving k that were suppressed when we described the construction of the random code. Our task is to show that for any sufficiently large k and any  $\epsilon > 0$ , the random code we have constructed satisfies the following when n is sufficiently large:

$$n^{-1}\log\eta_e^{(k)} \le R_e + \epsilon, \ e \in \mathcal{E}$$
(21.165)

$$\tau_s^{(k)} \ge \omega_s^{(k)} - \epsilon, \ s \in \mathcal{S}$$
(21.166)

$$\Delta_t^{(k)} \le \epsilon, \ t \in \mathcal{T}. \tag{21.167}$$

For  $e \in \mathcal{E}$ , consider

$$n^{-1}\log\eta_e^{(k)} \le \alpha^{(k)}H(U_e^{(k)}) + \psi_e^{(k)}$$
(21.168)

$$\leq R_e + \mu^{(k)} + \psi_e^{(k)}, \tag{21.169}$$

where the first inequality follows from the upper bound in (21.147) and the second inequality follows from (21.137). Since  $\mu^{(k)} \to 0$  as  $k \to \infty$ , we can let k be sufficiently large so that

$$\mu^{(k)} < \epsilon. \tag{21.170}$$

With k fixed, since  $\psi_e^{(k)} \to 0$  as  $\delta \to 0$ , by letting  $\delta$  be sufficiently small, we have

$$\mu^{(k)} + \psi_e^{(k)} \le \epsilon, \tag{21.171}$$

and (21.165) follows from (21.169). For  $s \in S$ , from the lower bound in (21.141), we have

$$\tau_s^{(k)} \ge \omega_s^{(k)} - \epsilon, \qquad (21.172)$$

proving (21.166).

The proof of (21.167), which is considerably more involved, will be organized into a few lemmas. For the sake of presentation, the proofs of these lemmas will be deferred to the end of the section.

For  $i \in \mathcal{S}$  and  $i \in \mathcal{V} \setminus (S \cup T)$ , the function  $u_e$ , where  $e \in \text{Out}(i)$ , has been defined in (21.149) and (21.156), respectively. Since the network is acyclic, we see by induction that all the auxiliary random variables  $U_e, e \in \mathcal{E}$  are functions of the auxiliary random variables  $Y_{\mathcal{S}}$ . Thus there exists a function  $\tilde{u}_e$  such that

$$U_e = \tilde{u}_e(Y_\mathcal{S}). \tag{21.173}$$

Equating the above with (21.149) and (21.156), we obtain

$$u_e(Y_s) = \tilde{u}_e(Y_S) \tag{21.174}$$

and

$$u_e(U_{\mathrm{In}(i)}) = \tilde{u}_e(Y_{\mathcal{S}}), \qquad (21.175)$$

respectively. These relations will be useful subsequently.

In the rest of the section, we will analyze the probabilities of decoding error for the random code we have constructed for a fixed k, namely  $\Delta_t^{(k)}$ , for  $t \in \mathcal{T}$ . With a slight abuse of notation, we write

$$\tilde{u}_{\mathcal{E}'}(\cdot) = (\tilde{u}_d(\cdot) : d \in \mathcal{E}') \tag{21.176}$$

for  $\mathcal{E}' \subset \mathcal{E}$ . Again, we temporarily suppress all the superscripts invoking k.

#### Lemma 21.9. Let

$$X_{\mathcal{S}} = x_{\mathcal{S}} \tag{21.177}$$

$$\mathbf{Y}_{\mathcal{S}}(x_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}} \in T^{\hat{n}}_{[Y_{\mathcal{S}}]\delta},\tag{21.178}$$

and for  $e \in \mathcal{E}$ , let  $C_e$  take the value  $c_e$ , which by the code construction is a function of  $x_{\mathcal{S}}$  and  $\mathbf{y}_{\mathcal{S}}$ . Then

$$\mathbf{U}_{\mathrm{In}(t)}(c_{\mathrm{In}(t)}) = \tilde{u}_{\mathrm{In}(t)}(\mathbf{y}_{\mathcal{S}}). \tag{21.179}$$

and

$$(\mathbf{y}_{\mathcal{S}}, \mathbf{U}_{\mathrm{In}(t)}(c_{\mathrm{In}(t)})) \in T^{\tilde{n}}_{[Y_{\mathcal{S}}U_{\mathrm{In}(t)}]\delta}$$
(21.180)

for all  $t \in \mathcal{T}$ .

Let

$$Err_{t} = \{g_{t}(C_{\text{In}(t)}) \neq X_{\beta(t)}\} = \{\tilde{g}_{t}(X_{\mathcal{S}}) \neq X_{\beta(t)}\}$$
(21.181)

be the event of a decoding error at sink node t, i.e.,

$$\Pr\{Err_t\} = \Delta_t \tag{21.182}$$

(cf. (21.10)). In the following, we will obtain an upper bound on  $Pr\{Err_t\}$ . Consider

$$\Pr\{Err_t\} = \sum_{x_{\beta(t)} \in \mathcal{X}_{\beta(t)}} \Pr\{Err_t | X_{\beta(t)} = x_{\beta(t)}\} \Pr\{X_{\beta(t)} = x_{\beta(t)}\}, \quad (21.183)$$

and for  $\mathcal{S}' \subset \mathcal{S}$ , let

~

$$1_{\mathcal{S}'} = \underbrace{(1, 1, \cdots, 1)}_{|\mathcal{S}'|}.$$
 (21.184)

Since  $\Pr\{Err_t|X_{\beta(t)} = x_{\beta(t)}\}$  are identical for all  $x_{\beta(t)}$  by symmetry in the code construction, from (21.183), we have

$$\Pr\{Err_t\} = \Pr\{Err_t | X_{\beta(t)} = 1_{\beta(t)}\} \sum_{x_{\beta(t)} \in \mathcal{X}_{\beta(t)}} \Pr\{X_{\beta(t)} = x_{\beta(t)}\}$$
(21.185)

$$= Pr\{Err_t | X_{\beta(t)} = 1_{\beta(t)}\}.$$
(21.186)

In other words, we can assume without loss of generality that  $X_{\beta(t)} = 1_{\beta(t)}$ . To facilitate our discussion, define the event

$$E_{\mathcal{S}} = \{ \mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) \in T^{\hat{n}}_{[Y_{\mathcal{S}}]\delta} \}$$
(21.187)

Following (21.186), we have

$$Pr\{Err_{t}\} = Pr\{Err_{t}|X_{\beta(t)} = 1_{\beta(t)}, E_{\mathcal{S}}\} Pr\{E_{\mathcal{S}}|X_{\beta(t)} = 1_{\beta(t)}\} + Pr\{Err_{t}|X_{\beta(t)} = 1_{\beta(t)}, E_{\mathcal{S}}^{c}\} Pr\{E_{\mathcal{S}}^{c}|X_{\beta(t)} = 1_{\beta(t)}\} = Pr\{Err_{t}|X_{\beta(t)} = 1_{\beta(t)}, E_{\mathcal{S}}\} Pr\{E_{\mathcal{S}}\}$$
(21.188)

$$+Pr\{Err_t|X_{\beta(t)} = 1_{\beta(t)}, E_{\mathcal{S}}^c\} \Pr\{E_{\mathcal{S}}^c\}$$
(21.189)

$$\leq Pr\{Err_t | X_{\beta(t)} = 1_{\beta(t)}, E_{\mathcal{S}}\} \cdot 1 + 1 \cdot \Pr\{E_{\mathcal{S}}^c\}$$
(21.190)

$$\leq Pr\{Err_t|X_{\beta(t)} = 1_{\beta(t)}, E_{\mathcal{S}}\} + \lambda, \qquad (21.191)$$

where the last inequality follows from the strong AEP and  $\lambda \to 0$  as  $\delta \to 0$ . Upon defining the event

$$E'_{\mathcal{S}} = \{X_{\beta(t)} = 1_{\beta(t)}\} \cap E_{\mathcal{S}}, \qquad (21.192)$$

we have

$$\Pr\{Err_t\} \le \Pr\{Err_t | E_{\mathcal{S}}'\} + \lambda.$$
(21.193)

We now further analyze the conditional probability in (21.193). For  $x_{\beta(t)} \in \mathcal{X}_{\beta(t)}$ , define the event

$$E_t(x_{\beta(t)}) = \left\{ (\mathbf{Y}_{\beta(t)}(x_{\beta(t)}), \mathbf{U}_{\mathrm{In}(t)}(C_{\mathrm{In}(t)})) \in T^{\hat{n}}_{[Y_{\beta(t)}U_{\mathrm{In}(t)}]\delta} \right\}.$$
 (21.194)

Since  $X_{\beta(t)} = 1_{\beta(t)}$ , decoding at sink node t is correct if the received indices  $C_{\text{In}(t)}$  is decoded to  $1_{\beta(t)}$ . This is the case if and only if  $E_t(1_{\beta(t)})$  occurs but  $E_t(x_{\beta(t)})$  does not occur for all  $x_{\beta(t)} \neq 1_{\beta(t)}$ . It follows that

$$Err_{t}^{c} = E_{t}(1_{\beta(t)}) \cap \left( \bigcap_{x_{\beta(t)} \neq 1_{\beta(t)}} E_{t}(x_{\beta(t)})^{c} \right), \qquad (21.195)$$

or

$$Err_{t} = E_{t}(1_{\beta(t)})^{c} \cup \left( \bigcup_{x_{\beta(t)} \neq 1_{\beta(t)}} E_{t}(x_{\beta(t)}) \right), \qquad (21.196)$$

which implies

$$\Pr\{Err_t | E'_{\mathcal{S}}\} = \Pr\{E_t(1_{\beta(t)})^c \cup \left(\bigcup_{x_{\beta(t)} \neq 1_{\beta(t)}} E_t(x_{\beta(t)})\right) | E'_{\mathcal{S}}\}.$$
 (21.197)

By the union bound, we have

$$\Pr\{Err_t | E'_{\mathcal{S}}\} \le \Pr\{E_t(1_{\beta(t)})^c | E'_{\mathcal{S}}\} + \sum_{x_{\beta(t)} \neq 1_{\beta(t)}} \Pr\{E_t(x_{\beta(t)}) | E'_{\mathcal{S}}\}$$
(21.198)

$$= \sum_{x_{\beta(t)} \neq 1_{\beta(t)}} \Pr\{E_t(x_{\beta(t)}) | E'_{\mathcal{S}}\},$$
(21.199)

where the last step follows because  $\Pr\{E_t(1_{\beta(t)})^c | E'_{\mathcal{S}}\}\$  in (21.198) vanishes by Lemma 21.9. The next two lemmas will be instrumental in obtaining an upper bound on  $\Pr\{E_t(x_{\beta(t)}) | E'_{\mathcal{S}}\}\$  in (21.199).

For any proper subset  $\Psi$  of  $\beta(t)$ , let

$$\Lambda_{\Psi} = \{ x_{\beta(t)} \neq 1_{\beta(t)} : x_s = 1 \text{ if and only if } s \in \Psi \}.$$
(21.200)

Note that  $\{\Lambda_{\Psi}\}$  is a partition of the set  $\mathcal{X}_{\beta(t)} \setminus \{1_{\beta(t)}\}$ . For  $x_{\beta(t)} \in \Lambda_{\Psi}$ ,  $x_{\beta(t)}$  and  $1_{\beta(t)}$  are identical for exactly the components indexed by  $\Psi$ .

**Lemma 21.10.** For  $x_{\beta(t)} \in \Lambda_{\Psi}$ , where  $\Psi$  is a proper subset of  $\beta(t)$ ,

$$\Pr\{E_t(x_{\beta(t)})|E'_{\mathcal{S}}\} \le 2^{-n\alpha(H(Y_{\beta(t)\setminus\Psi}) - H(Y_{\beta(t)}|U_{\mathrm{In}(t)}) - \varphi_t)},$$
(21.201)

where  $\varphi_t \to 0$  as  $n \to \infty$  and  $\delta \to 0$ .

Lemma 21.11. For all sufficiently large n,

$$|\Lambda_{\Psi}| \le 2^{n(\alpha H(Y_{\beta(t)\setminus\Psi}) - \epsilon/4)}.$$
(21.202)

We now reinstate all the superscript involving k that have been suppressed. By Lemma 21.10, Lemma 21.11, and (21.199),

$$\Pr\{Err_t | E'_{\mathcal{S}}\}$$

$$\leq \sum_{x_{\beta(t)} \neq 1_{\beta(t)}} \Pr\{E_t(x_{\beta(t)}) | E'_{\mathcal{S}}\}$$
(21.203)

$$\leq \sum_{\Psi} \sum_{x_{\beta(t)} \in \Psi} \Pr\{E_t(x_{\beta(t)}) | E_{\mathcal{S}}'\}$$
(21.204)

$$\leq 2^{|\mathcal{E}|} 2^{-n\alpha^{(k)}} \left[ \left( H(Y_{\beta(t)\setminus\Psi}^{(k)}) - H\left(Y_{\beta(t)}^{(k)}|U_{\mathrm{In}(t)}^{(k)}\right) - \varphi_t \right] \\ \cdot 2^{n \left[ \alpha^{(k)} H\left(Y_{\beta(t)\setminus\Psi}^{(k)}\right) - \epsilon/4 \right]}$$

$$(21.205)$$

$$= 2^{|\mathcal{E}|} 2^{-n \left[\epsilon/4 - \alpha^{(\kappa)} H\left(Y_{j(\kappa)}^{(\kappa)}\right) U_{\mathrm{In}(t)}^{(\kappa)}\right) - \alpha^{(\kappa)} \varphi_t\right]}$$
(21.206)

$$\leq 2^{|\mathcal{E}|} 2^{-n\left(\epsilon/4 - \gamma^{(k)} - \alpha^{(k)}\varphi_t\right)} \tag{21.207}$$

$$\leq 2^{|\mathcal{E}|} 2^{-n\left(\epsilon/4 - \gamma^{(k)} - \varphi_t\right)} \tag{21.208}$$

where (21.207) follows from (21.138) and (21.208) follows from (21.131). Then from (21.182), (21.193), and (21.208), we have

$$\Delta_t^{(k)} \le 2^{|\mathcal{E}|} 2^{-n(\epsilon/4 - \gamma^{(k)} - \varphi_t)} + \lambda$$
(21.209)

We now choose  $k, n, \text{ and } \delta$  to make the upper bound above smaller than any prescribed  $\epsilon > 0$ . Since  $\gamma^{(k)} \to 0$  as  $k \to \infty$ , we can let k to be sufficiently large so that

21.7 Achievability 521

$$\gamma^{(k)} < \epsilon/4. \tag{21.210}$$

Then with k fixed, since  $\varphi_t \to 0$  as  $n \to \infty$  and  $\delta \to 0$ , and  $\lambda \to 0$  as  $\delta \to 0$ , by letting n be sufficiently large n and  $\delta$  be sufficiently small, we have

1. 
$$\gamma^{(k)} + \varphi_t < \epsilon/4$$
, so that  $2^{|\mathcal{E}|} 2^{-n(\epsilon/4 - \gamma^{(k)} - \varphi_t)} \to 0$  as  $n \to \infty$ ;  
2.  $\Delta_t^{(k)} \le \epsilon$ .

This completes the proof of (21.167).

Hence, we have proved the achievability of  $\omega^{(k)}$  for all sufficiently large k. Then the closedness of  $\mathcal{R}$  implies the achievability of  $\boldsymbol{\omega} = \lim_{k \to \infty} \boldsymbol{\omega}^{(k)}$ , where  $\omega \in \mathcal{R}'$ . The achievability of  $\mathcal{R}'$  is established.

Proof of Lemma 21.9. We first prove that given

$$X_{\mathcal{S}} = x_{\mathcal{S}} \tag{21.211}$$

and

$$\mathbf{Y}_{\mathcal{S}}(x_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}} \in T^{\hat{n}}_{[Y_{\mathcal{S}}]\delta}, \qquad (21.212)$$

the following hold for all non-source nodes i (i.e.,  $i \in \mathcal{V} \setminus S$ ):

- $$\begin{split} &\text{i) } \mathbf{U}_{\mathrm{In}(i)}(c_{\mathrm{In}(i)}) \in T^{\hat{n}}_{[U_{\mathrm{In}(i)}]\delta}; \\ &\text{ii) } k_e(c_{\mathrm{In}(i)}) \neq 0, \, e \in \mathrm{Out}(i); \\ &\text{iii) } \mathbf{U}_e(k_e(c_{\mathrm{In}(i)})) = \tilde{u}_e(\mathbf{y}_{\mathcal{S}}), \, e \in \mathrm{Out}(i). \end{split}$$

Note that for  $i \in \mathcal{T}$ ,  $Out(i) = \emptyset$  in ii) and iii). By the consistency of strong typicality (Theorem 6.7),

$$\mathbf{y}_s \in T^n_{[Y_s]\delta} \tag{21.213}$$

for all  $s \in \mathcal{S}$ . Then according to the construction of the code, for all  $e \in \text{Out}(s)$ ,

$$k_e(x_s) \neq 0 \tag{21.214}$$

and

$$\mathbf{U}_{e}(k_{e}(x_{s})) = u_{e}(\mathbf{y}_{s}).$$
 (21.215)

We now prove i) to iii) by induction on the non-source nodes according to any given coding order. Let  $i_1$  be the first non-source node to encode. Since

$$\operatorname{In}(i_1) \subset \mathcal{S},\tag{21.216}$$

for all  $d \in \text{In}(i_1)$ ,  $d \in \text{Out}(s)$  for some  $s \in S$ . Then

$$\mathbf{U}_d(c_d) = \mathbf{U}_d(k_d(x_s)) \tag{21.217}$$

$$= u_d(\mathbf{y}_s) \tag{21.218}$$

$$=\tilde{u}_d(\mathbf{y}_{\mathcal{S}}),\tag{21.219}$$

where (21.218) and (21.219) follows from (21.215) and (21.174), respectively. Thus

$$\mathbf{U}_{\mathrm{In}(i_1)}(c_{\mathrm{In}(i_1)}) = \tilde{u}_{\mathrm{In}(i_1)}(\mathbf{y}_{\mathcal{S}}).$$
(21.220)

Since  $U_{\text{In}(i_1)}$  is a function of  $Y_S$ , in light of (21.178),

$$\mathbf{U}_{\mathrm{In}(i)}(c_{\mathrm{In}(i)}) \in T^{n}_{[U_{\mathrm{In}(i)}]\delta} \tag{21.221}$$

by the preservation property of strong typicality, proving i). According to the code construction, this also implies ii). Moreover,

$$\mathbf{U}_{e}(k_{e}(c_{\mathrm{In}(i_{1})})) = u_{e}(\mathbf{U}_{\mathrm{In}(i_{1})}(c_{\mathrm{In}(i_{1})}))$$
(21.222)

$$= \tilde{u}_e(\mathbf{y}_{\mathcal{S}}), \tag{21.223}$$

where the last equality is obtained by replacing in (21.175) the random variable  $U_d$  by the sequence  $\mathbf{U}_d(c_d)$  and the random variable  $Y_s$  by the sequence  $\mathbf{Y}_{\mathcal{S}}(x_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}$ , proving iii).

We now consider any non-source node i in the network. Assume that i) to iii) are true for all the nodes upstream to node i. For  $d \in \text{In}(i)$ , if  $d \in \text{Out}(s)$ for some  $s \in S$ , we have already proved in (21.219) that

$$\mathbf{U}_d(c_d) = \tilde{u}_d(\mathbf{y}_{\mathcal{S}}). \tag{21.224}$$

Otherwise,  $d \in Out(i')$ , where node i' is upstream to node i. Then

$$\mathbf{U}_d(c_d) = \mathbf{U}_d(k_d(c_{\mathrm{In}(i')})) \tag{21.225}$$

$$= u_d(\mathbf{U}_{\mathrm{In}(i')}(c_{\mathrm{In}(i')}))$$
(21.226)

$$=\tilde{u}_d(\mathbf{y}_{\mathcal{S}}),\tag{21.227}$$

In the above, (21.226) follows from ii) for node i' by the induction hypothesis and the code construction, and (21.227) follows from (21.175). Therefore, (21.224) is valid for all  $d \in \text{In}(i)$ . Hence,

$$\mathbf{U}_{\mathrm{In}(i)}(c_{\mathrm{In}(i)}) = \tilde{u}_{\mathrm{In}(i)}(\mathbf{y}_{\mathcal{S}}). \tag{21.228}$$

which is exactly the same as (21.220) except that  $i_1$  is replaced by i. Then by means of the same argument, we conclude that i), ii), and iii) hold for node i.

As (21.228) holds for any non-source node i, it holds for any sink node t. This proves (21.179) for all  $t \in \mathcal{T}$ . Furthermore, since  $U_{\text{In}(t)}$  is a function of  $Y_{\mathcal{S}}$ ,  $(Y_{\mathcal{S}}, U_{\text{In}(t)})$  is also a function of  $Y_{\mathcal{S}}$ . Then in view of (21.179), (21.180) follows from the preservation property of strong typicality. This completes the proof of the lemma.  $\Box$ 

Proof of Lemma 21.10. Consider

$$\Pr\{E_t(x_{\beta(t)})|E'_{\mathcal{S}}\}\$$

$$=\sum_{\mathbf{y}_{\mathcal{S}}\in T^{\hat{n}}_{[Y_{\mathcal{S}}]\delta}}\Pr\{E_t(x_{\beta(t)})|\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}})=\mathbf{y}_{\mathcal{S}},E'_{\mathcal{S}}\}\Pr\{\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}})=\mathbf{y}_{\mathcal{S}}|E'_{\mathcal{S}}\}.$$
(21.229)

To analyze  $\Pr\{E_t(x_{\beta(t)})|\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}, E'_{\mathcal{S}}\}\$  in the above summation, let us condition on the event  $\{\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}, E'_{\mathcal{S}}\}\$ , where  $\mathbf{y}_{\mathcal{S}} \in T^{\hat{n}}_{[Y_{\mathcal{S}}]\delta}$ . It then follows from (21.179) in Lemma 21.9 that

$$\mathbf{U}_{\mathrm{In}(t)}(c_{\mathrm{In}(t)}) = \tilde{u}_{\mathrm{In}(t)}(\mathbf{y}_{\mathcal{S}}). \tag{21.230}$$

Therefore, the event  $E_t(x_{\beta(t)})$  is equivalent to

$$(\mathbf{y}_{\Psi}, \mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi}), \tilde{u}_{\mathrm{In}(t)}(\mathbf{y}_{\mathcal{S}})) \in T^{\hat{n}}_{[Y_{\Psi}Y_{\beta(t)\setminus\Psi}U_{\mathrm{In}(t)}]\delta}$$
(21.231)

(cf. (21.194)), or

$$\mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi}) \in T^{\hat{n}}_{[Y_{\beta(t)\setminus\Psi}|Y_{\Psi}U_{\mathrm{In}(t)}]\delta}(\mathbf{y}_{\Psi}, \tilde{u}_{\mathrm{In}(t)}(\mathbf{y}_{\mathcal{S}})).$$
(21.232)

Thus

$$\Pr\{E_{t}(x_{\beta(t)})|\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}, E_{\mathcal{S}}'\} = \sum_{\mathbf{y}_{\beta(t)\setminus\Psi}\in T_{[Y_{\beta(t)\setminus\Psi}|Y_{\Psi}U_{\mathrm{In}(t)}]^{\delta}} (\mathbf{y}_{\Psi}, \tilde{u}_{\mathrm{In}(t)}(\mathbf{y}_{\mathcal{S}}))} \Pr\{\mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi}) = \mathbf{y}_{\beta(t)\setminus\Psi}|\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}, E_{\mathcal{S}}'\}.$$
 (21.233)

Since  $x_s \neq 1$  for  $s \in \beta(t) \setminus \Psi$ ,  $\mathbf{Y}_{\beta(t) \setminus \Psi}(x_{\beta(t) \setminus \Psi})$  is independent of the random sequences  $\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}})$  and the event  $E'_{\mathcal{S}}$  by construction. Therefore,

$$\Pr\{\mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi}) = \mathbf{y}_{\beta(t)\setminus\Psi}|\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}, E'_{\mathcal{S}}\} = \Pr\{\mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi}) = \mathbf{y}_{\beta(t)\setminus\Psi}\}.$$
(21.234)

By the consistency of strong typicality, if

$$\mathbf{y}_{\beta(t)\setminus\Psi} \in T^{\hat{n}}_{[Y_{\beta(t)\setminus\Psi}|Y_{\Psi}U_{\mathrm{In}(t)}]\delta}(\mathbf{y}_{\Psi}, \tilde{u}_{\mathrm{In}(t)}(\mathbf{y}_{\mathcal{S}})), \qquad (21.235)$$

then

$$\mathbf{y}_{\beta(t)\setminus\Psi} \in T^{\hat{n}}_{[Y_{\beta(t)\setminus\Psi}]\delta}.$$
(21.236)

Since  $\mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi})$  are generated i.i.d. according to the distribution of  $Y_{\beta(t)\setminus\Psi}$ , by the strong AEP,

$$\Pr\{\mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi}) = \mathbf{y}_{\beta(t)\setminus\Psi}\} \le 2^{-\hat{n}(H(Y_{\beta(t)\setminus\Psi})-\rho)}, \qquad (21.237)$$

where  $\rho \to 0$  as  $\delta \to 0$ . Combining (21.234) and (21.237), we have

$$\Pr\{\mathbf{Y}_{\beta(t)\setminus\Psi}(x_{\beta(t)\setminus\Psi}) = \mathbf{y}_{\beta(t)\setminus\Psi} | \mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}, E_{\mathcal{S}}'\} \le 2^{-\hat{n}(H(Y_{\beta(t)\setminus\Psi})-\rho)}.$$
(21.238)

By Theorem 6.10,

$$|T^{\hat{n}}_{[Y_{\beta(t)\setminus\Psi}|Y_{\Psi}U_{\mathrm{In}(t)}]\delta}(\mathbf{y}_{\Psi},\tilde{u}_{\mathrm{In}(t)}(\mathbf{y}_{\mathcal{S}}))| \leq 2^{\hat{n}(H(Y_{\beta(t)\setminus\Psi}|Y_{\Psi}U_{\mathrm{In}(t)})+\sigma)}, \quad (21.239)$$

where  $\sigma \to 0$  as  $\hat{n} \to \infty$  and  $\delta \to 0$ . It then follows from (21.233), (21.238), and (21.239) that

$\Pr\{E_t(x_{\beta(t)}) \mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}, E_{\mathcal{S}}'\}$	
$\leq 2^{\hat{n}(H(Y_{\beta(t)\setminus\Psi} Y_{\Psi}U_{\mathrm{In}(t)})+\sigma)}2^{-\hat{n}(H(Y_{\beta(t)\setminus\Psi})-\rho)}$	(21.240)
$= 2^{-\hat{n}(H(Y_{\beta(t)\setminus\Psi}) - H(Y_{\beta(t)\setminus\Psi} Y_{\Psi}U_{\mathrm{In}(t)}) - \sigma - \rho)}$	(21.241)
$\leq 2^{-\hat{n}(H(Y_{\beta(t)\setminus\Psi})-H(Y_{\beta(t)} U_{\mathrm{In}(t)})-\sigma-\rho)}$	(21.242)
$\leq 2^{-n\alpha(H(Y_{\beta(t)\setminus\Psi})-H(Y_{\beta(t)} U_{\mathrm{In}(t)})-\varphi_t)}$	(21.243)

where (21.242) is justified by

$$H(Y_{\beta(t)\setminus\Psi}|Y_{\Psi}U_{\mathrm{In}(t)}) \le H(Y_{\beta(t)\setminus\Psi}|Y_{\Psi}U_{\mathrm{In}(t)}) + H(Y_{\Psi}|U_{\mathrm{In}(t)})$$
(21.244)

$$=H(Y_{\beta(t),\backslash\Psi},Y_{\Psi}|U_{\mathrm{In}(t)})$$
(21.245)

$$= H(Y_{\beta(t)}|U_{\text{In}(t)}), \qquad (21.246)$$

(21.243) follows from (21.142), and  $\varphi_t \to 0$  as  $n \to \infty$  and  $\delta \to 0$ . In (21.229),

$$\Pr\{\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}} | E'_{\mathcal{S}} \}$$
  
= 
$$\Pr\{\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}} | X_{\beta(t)} = 1_{\beta(t)}, E_{\mathcal{S}} \}$$
(21.247)

$$= \Pr\{\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}|E_{\mathcal{S}}\}$$
(21.248)

$$= \Pr\{\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}} | \mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) \in T^{\hat{n}}_{[Y_{\mathcal{S}}]\delta}\}.$$
(21.249)

Hence, it follows from (21.229) and (21.243) that

$$\Pr\{E_{t}(x_{\beta(t)})|E_{\mathcal{S}}'\}$$

$$\leq 2^{-n\alpha(H(Y_{\beta(t)\setminus\Psi})-H(Y_{\beta(t)}|U_{\mathrm{In}(t)})-\varphi_{t})} .$$

$$\sum_{\mathbf{y}_{\mathcal{S}}\in T_{[Y_{\mathcal{S}}]\delta}^{\hat{n}}} \Pr\{\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) = \mathbf{y}_{\mathcal{S}}|\mathbf{Y}_{\mathcal{S}}(1_{\mathcal{S}}) \in T_{[Y_{\mathcal{S}}]\delta}^{\hat{n}}\}$$
(21.250)

$$= 2^{-n\alpha(H(Y_{\beta(t)}\setminus\psi) - H(Y_{\beta(t)}|U_{\mathrm{In}(t)}) - \varphi_t)} \cdot 1$$

$$(21.251)$$

$$= n\alpha(H(Y_{\beta(t)}\setminus\psi) - H(Y_{\beta(t)}|U_{\mathrm{In}(t)}) - \varphi_t) \cdot 1$$

$$(21.252)$$

$$=2^{-n\alpha(H(Y_{\beta(t)}\setminus\Psi)-H(Y_{\beta(t)}|U_{\text{In}(t)})-\varphi_t)}.$$
(21.252)

The lemma is proved.  $\Box$ 

Proof of Lemma 21.11. Let n be sufficiently large. Consider

$$|\Lambda_{\Psi}| = \prod_{s \in \beta(t) \setminus \Psi} |\mathcal{X}_s| \tag{21.253}$$

$$\stackrel{a)}{=} \prod_{s \in \beta(t) \setminus \Psi} \theta_s \tag{21.254}$$

21.7 Achievability 525

$$\stackrel{b)}{=} \prod_{s \in \beta(t) \setminus \Psi} \lceil 2^{n\tau_s} \rceil \tag{21.255}$$

$$\stackrel{(2)}{\leq} \prod_{s \in \beta(t) \setminus \Psi} \lceil 2^{n(\omega_s - \epsilon/3)} \rceil \tag{21.256}$$

$$\leq \prod_{s \in \beta(t) \setminus \Psi} 2^{n(\omega_s - \epsilon/4)} \tag{21.257}$$

$$\stackrel{d)}{=} \prod_{s \in \beta(t) \setminus W} 2^{n(\alpha H(Y_s) - \epsilon/4)} \tag{21.258}$$

$$=2^{n\sum_{s\in\beta(t)\setminus\Psi}(\alpha H(Y_s)-\epsilon/4)}$$
(21.259)

$$=2^{n\left\lfloor\alpha\sum_{s\in\beta(t)\setminus\Psi}H(Y_s)-(|\beta(t)|-|\Psi|)\epsilon/4\right\rfloor}$$
(21.260)

$$=2^{n\left[\alpha H(Y_{\beta(t)\setminus\Psi})-(|\beta(t)|-|\Psi|)\epsilon/4\right]}$$
(21.261)

$$\stackrel{e)}{\leq} 2^{n(\alpha H(Y_{\beta(t)\setminus\Psi})-\epsilon/4)},\tag{21.262}$$

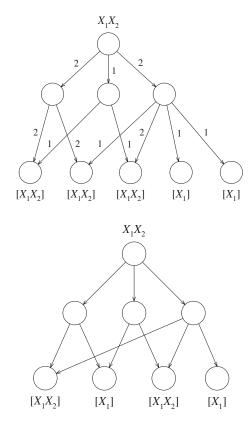
where

- a) follows from (21.144);
- b) follows from (21.143);
- c) follows from (21.141);
- d) follows from (21.133);
- e) follows because  $\Psi$  is a proper subset of  $\beta(t)$ .

The lemma is proved.  $\Box$ 

## Problems

- 1. Show that source separation is optimal for the networking problem depicted in Figure 21.3.
- 2. Consider the following network. ??
  - a) Let  $\omega_i$  be the rate of information source  $X_i$ . Determine and illustrate the max-flow bounds.
  - b) Are the max-flow bounds achievable?
  - c) Is source separation always optimal?
- 3. Repeat Problem 2 for the following network in which the capacities of all the edges are equal to 1.
- 4. Consider a disk array with 3 disks. Let  $X_1$ ,  $X_2$ , and  $X_3$  be 3 mutually independent pieces of information to be retrieved from the disk array, and let  $S_1$ ,  $S_2$ , and  $S_3$  be the data to be stored separately in the 3 disks. It is required that  $X_1$  can be retrieved from  $S_i$ ,  $i = 1, 2, 3, X_2$  can be retrieved from  $(S_i, S_j), 1 \le i < j \le 3$ , and  $X_3$  can be retrieved from  $(S_1, S_2, S_3)$ .



a) Prove that for i = 1, 2, 3,

$$H(S_i) = H(X_1) + H(S_i|X_1).$$

b) Prove that for  $1 \le i < j \le 3$ ,

 $H(S_i|X_1) + H(S_j|X_1) \ge H(X_2) + H(S_i, S_j|X_1, X_2).$ 

c) Prove that

$$H(S_1, S_2, S_3 | X_1, X_2) = H(X_3).$$

d) Prove that for i = 1, 2, 3,

$$H(S_i) \ge H(X_1).$$

e) Prove that

$$H(S_i) + H(S_j) \ge 2H(X_1) + H(X_2) + H(S_i, S_j | X_1, X_2).$$

f) Prove that

$$2S_i + S_{i\oplus 1} + S_{i\oplus 2} \ge 4H(X_1) + 2H(X_2) + H(X_3),$$

where i = 1, 2, 3 and

$$i \oplus j = \begin{cases} i+j & \text{if } i+j \leq 3\\ i+j-3 & \text{if } i+j > 3 \end{cases}$$

for  $1 \le i, j \le 3$ .

g) Prove that

$$H(S_1) + H(S_2) + H(S_3) \ge 3H(X_1) + \frac{3}{2}H(X_2) + H(X_3).$$

Parts d) to g) give constraints on  $H(S_1)$ ,  $H(S_2)$ , and  $H(S_3)$  in terms of  $H(X_1)$ ,  $H(X_2)$ , and  $H(X_3)$ . It was shown in Roche *et al.* [279] that these constraints are the tightest possible.

5. Generalize the setup in Problem 4 to K disks and show that

$$\sum_{i=1}^{K} H(S_i) \ge K \sum_{\alpha=1}^{K} \frac{H(X_{\alpha})}{\alpha}.$$

Hint: Use the inequalities in Problem 18 in Chapter 2 to prove that for  $s = 0, 1, \dots, K - 1$ ,

$$\sum_{i=1}^{K} H(S_i) \ge nK \sum_{\alpha=1}^{s} \frac{H(X_\alpha)}{\alpha} + \frac{K}{\binom{K}{s+1}}$$
$$\times \sum_{T:|T|=s+1} \frac{H(S_T|X_1, X_2, \cdots, X_s)}{s+1}$$

by induction on s, where T is a subset of  $\{1, 2, \dots, K\}$ .

- 6. Write out the achievable information rate region  $\mathcal{R}$  for the network in Problem 2.
- 7. By letting  $S = \{s\}$  and  $\beta(t) = \{s\}$  for all  $t \in \mathcal{T}$ , the multi-source network coding problem described in Section 21.3 becomes a single-source network coding problem. Write  $\boldsymbol{\omega} = \omega_s$ .
  - a) Write out the achievable information rate region  $\mathcal{R}$ .
  - b) Show that if  $\omega_s \in \mathcal{R}$ , then  $\omega_s \leq \max flow(t)$  for all  $t \in \mathcal{T}$ .
- 8. Show that if there exists an

$$(n, (\eta_{ij}: (i, j) \in \mathcal{E}), (\tau_s: s \in \mathcal{S}))$$

code which satisfies (21.11) and (21.13), then there always exists an

$$(n, (\eta_{ij}: (i, j) \in \mathcal{E}), (\tau'_s: s \in \mathcal{S}))$$

code which satisfies (21.11) and (21.13), where  $\tau'_s \leq \tau_s$  for all  $s \in S$ . Hint: use a random coding argument.

### **Historical Notes**

Multilevel diversity coding was studied by Yeung [359], where it was shown that source separation is not always optimal. Roche *et al.* [279] showed that source separation is optimal for symmetrical three-level diversity coding. This result was extended to any level by Yeung and Zhang [367] with a painstaking proof. Hau [143] studied all the one hundred configurations of a three-encoder diversity coding systems and found that source separation is optimal for eighty-six configurations.

Yeung and Zhang [368] introduced the distributed source coding model discussed in Section 21.2.2 which subsumes multilevel diversity coding. The region of all entropy functions previously introduced by Yeung [360] for studying information inequalities enabled them to obtain inner and outer bounds on the achievable information rate region for a variety of networks.

Distributed source coding is equivalent to multi-source network coding on a special class of acyclic networks. The inner and outer bounds on the achievable information rate region in [368] were generalized to arbitrary acyclic networks by Song *et al.* [308]. The gap between these bounds was finally closed by Yan *et al.* [351].

The insufficiency of specific forms of linear coding for multi-source network coding were demonstrated and discussed by Riis [274], Rasala Lehman and Lehman [270], and Medard *et al.* [242]. The insufficiency of very general forms of linear coding has been proved by Dougherty *et al.* [86].

Even though the achievable information rate region for multi-source network coding is characterized by all information inequalities (Shannon-type and non-Shannon-type), it is not clear whether there exists a multi-source network coding problem for which the characterization of the achievable information rate region necessarily involves non-Shannon-type inequalities. This important question was resolved by Dougherty *et al.* [88]. In this work, they constructed a multi-source network coding problem from matroids and demonstrated that a tighter outer bound on the achievable information rate region can be obtained by invoking the unconstrained non-Shannon-type inequality discovered by Zhang and Yeung [373]. Chan and Grant [59] recently proved that for every non-Shannon-type inequality that exists, there is a multi-source network coding problem for which the characterization of the achievable information rate region necessarily involves that particular inequality.

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a posterior distribution, 210 Abel, N.H., 380 Abelian group, 380, 392, 395, 396 Abrahams, J., 97, 529 Abramson, N., 77, 529 abstract algebra, 377, 480 Abu-Mostafa, Y.S., 529 acyclic network, 427-470, 475, 478, 492, 528, see also directed graph Aczél, J., 529 additive colored Gaussian noise, 281 additive noise, worst, 264, 283, 289 additive white Gaussian noise (AWGN), 273, 275 adjacent pair of channels, 428, 431, 432, 446, 449, 461, 470, 482, 488 adjoint matrix, 492 Ahlswede, R., 96, 176, 409, 410, 424, 425, 473, 493, 529 Ahmed, E., 539 algebraic coding, 432, 473 Algoet, P., 529 algorithm, 427, 448, 449, 449, 458, 458, **464** exponential-time, 452 polynomial-time, 449, 450, 452, 458, 460, 473 almost everywhere (a.e.), 241 almost perfect reconstruction, 102, 104 alternating optimization algorithm, 208-210, 213 convergence, 218 Amari, S., 47, 530

Anantharam, V., 530 ancestral order, 429, see upstream-todownstream order Anderson, J.B., 530 applied mathematics, 4 applied probability, 3 Argawal, A., 529 Arimoto, S., 163, 176, 223, 530, see also Blahut-Arimoto algorithms arithmetic, 427 arithmetic mean, 147 ascendant, 92 Ash, R.B., 530 asymptotic equipartition property (AEP) for continuous random variables, 242, 242 - 244typical sequence, 243 typical set, 243 strong, see also strong asymptotic equipartition property weak, see also weak asymptotic equipartition property asymptotically reliable communication, 136atom of a field, 50 weight of, 303 audio signal, 180 audio source, 372 autocorrelation function, 274, 279 auxiliary random variable, 360, 361, 373, 525

average distortion, 180, 181, 183, 197, 201.207 expected, 197 average input constraint, 262, 263, 288 average probability of error, 146, 175, 255Ayanoglu, E., 410, 530 backbone network, 437 Balkenhol, B., 96, 529 Balli, H., 473, 530 bandlimited signal, 277, 278, 280 orthonormal basis, 277, 278 bandpass filter, 275, 281 bandwidth, 275, 402, 469 downlink, 407 Barbero, Á.I., 494, 530 Barron, A.R., 530 base field, 429-434, 438, 441, 443, 445-448, 451, 452, 460, 465-468, 472, 476, 477, 479, 482, 487, 490 - 492basic inequalities, 27, 26-28, 57, 129, 304, 314, 329-349, 351, 362, 363, 371-375, 377, 392 Bassalygo, L.A., 530 Bayesian network, 11, 148, 341, 349 intersection, 11 BCH (Bose-Chaudhuri-Hocquenghem) code, 163 Beethoven's violin concerto, 1 Bell Telephone Laboratories, 2 Berger, T., VIII, 75, 132, 205, 223, 530, 535, 545 Berlekamp, E.R., 530, 542 Berrou, C., 162, 530 Berstel, J., 530 beta distribution, 249 Bhargava, V.K., 544 biased coin, 79 binary arbitrarily varying channel, 176 binary covering radius, 203 binary entropy function, 13, 32, 33, 193 binary erasure channel, 144, 168 binary symmetric channel (BSC), 133, 142, 172, 176 binomial formula, 298, 345, 346 bit, 3, 13, 416 Blackwell, D., 530

Blahut, R.E., 176, 205, 223, 530, 531, 544Blahut-Arimoto algorithms, VIII, 145, 196, 207-223 channel capacity, 210-214, 223 convergence, 222 rate-distortion function, 215-218, 223 Blakley, G.R., 531 block code, 102 linear, 432 block length, 102, 135, 147, 179, 207, 280, 419 Blundo, C., 76, 531 Bollobás, B., 531 Bondy, J.A., 306, 531 Borade, S., 425, 531 Bose, R.C., 531, see also BCH code bottleneck, 415 boundary condition, 432, 475, 483, 485 brain, 372 branching probabilities, 92 Breiman, L., 110, 530, 531, see also Shannon-McMillan-Breiman theorem broadcast, 405 constraint, 406 broadcast constraint, 405 Bruck, J., 538 BSC, see binary symmetric channel Burrows, M., 531 butterfly network, 402-405, 410, 417, 452, 471 Byers, J., 531 cable cut, 460 Cai, N., 115, 409, 410, 424, 425, 473, 493, 494, 528, 529, 531, 538, 543, 545Calderbank, R., 531 Capocelli, R.M., 76, 531 Cartesian product, 208 cascade of channels, 173 Cauchy distribution, 250 causality, 172, 419, 423, 482, 485 CDF, see cumulative distribution function Cesáro mean, 39 chain rule for conditional entropy, 21

conditional mutual information, 22 differential entropy, 240 entropy, 20, 29 mutual information, 21, 30 Chan, T.H., 12, 129, 373, 375, 397, 528, 531, 532 channel capacity computation of, VIII channel characteristics, 133 channel code, 133, 147, 171 probability of error, 134 rate, 136, 147 with feedback, 163, 174 without feedback, 145 channel coding, 401, 405 channel coding theorem, 132 for continuous memoryless channel, 255, 254-255 achievability, 259-263 converse, 256-259 random code, 259 for discrete memoryless channel, VIII, 3, 47, 147, 145-147, 176 achievability, 145, 153-160 converse, 145, 167 random code, 155 strong converse, 153, 176 channel failure, 460 channel with memory, 168, 172, 174, 175Charikar, M., 529 Chatin, G.J., 532 Chekuri, C., 532 Chernoff bound, 115 Chernoff, H., 532 child, 92 Chou, P.A., 411, 473, 532, 536 chronological order, 420 Chuang, I.L., 375, 540 Chung, K.L., 110, 532 cipher text, 69, 75 classical entropy, 375 closure, of a group, 378 CMC, see continuous memoryless channel code alphabet, 80 code tree, 84, 84-95, 97 pruning of, 93 codebook, 145, 183, 198, 254, 261

codeword, 102, 145, 183, 198, 254 coding session, 420, 441 transaction, 410, 420 coding theory, 162 coefficient of correlation, 249 coloring filter, 289 column space, 324 combinatorics, 125 commodity, 401, 417 commodity flow, 410 communication channel, 275, 402 capacity, 402 error-free, 405 receiver, 275 communication engineer, 3 communication engineering, 133, 230, 264communication system, 1, 3, 235, 251, 401, 406 design of, 405 discrete-time, 138 practical, 133, 163 Shannon's model, 2 communication theory, 3 commutative, 380, 381, 481 commutative group, see Abelian group compact disc, 1 compact set, 142, 221 complex conjugate, 274 complex number, 480 composite function, 381 compound source, 204 computational complexity, 450, 452, 458, 460, 466, 471procedure, 333, 336 resource, 469 computer communication, 162 computer science, 13 computer storage systems, 162 concavity, 43, 46, 66, 68, 218-221, 253, 286conditional branching distribution, 92 conditional distribution, 235 conditional entropy, 7, 14 conditional independence, IX, 8, 312, 326-327, 341, 349 elemental forms, 348 structure of, 12, 375

conditional mutual independence, 292 - 301conditional mutual information, 7, 16, 19, 239 conditional probability of error, 146, 255configuration, 460, 462, 463 constant sequence, 183 constant term, 480, 482 continuous memoryless channel (CMC), 252, 252-263, 288, 289 achievable rate, 255 average input constraint, 253 capacity, 253, 255 continuous partial derivatives, 208, 213, 220, 221 convergence, 480 in divergence, 26, 45 in probability, 99, 100, 106, 260 in variational distance, 18, 26, 45, 46 convex closure, 388, 391 convex cone, 356 convexity, 24, 43, 44, 67, 184, 187, 189, 197, 208, 212, 217, 316, 358 convolution, 276 convolutional code, 163 convolutional coding, 480 convolutional multicast, 489, 490, 492, 493convolutional network code, 409, 482, 477 - 493decoding, 488-493 decoding delay, 489, 493 decoding kernel, 489, 492, 493 multicast, 489 correlation matrix, 226, 227, 250, 283-285, 287 coset left, 381, 382-384 right, 381 Costello, Jr., D.J., 531, 538, 544 countable alphabet, 32-34, 40, 45, 108, 110, 130-132 covariance, 226 covariance matrix, 226-228, 248-250, 271Cover, T.M., VII, 223, 289, 529, 532, 533cross-correlation function, 274, 275, 278

cross-spectral density, 275 crossover probability, 133, 173, 176, 194 cryptography, 473 Csiszár, I., VIII, 47, 77, 132, 223, 328, 529, 532, 535, 543 cumulative distribution function (CDF), 225conditional, 225 joint, 225, 267 marginal, 225 cyclic network, 475-493, see also directed graph delay-free, 475-477 D-adic distribution, 86, 91 D-ary source code, 80 D-it, 13, 82, 94  $D_{max}, 182, 187, 189, 193$  $d_{max}, 197$ Dantzig, G.B., 533 Daróczy, Z., 529 data block, 466 aysnchronous transmission, 467 encoded, 466, 467 data communication, 163, 401 data packet, 401, 469, 473 data processing theorem, 31, 71, 152, 256, 339, 372 Davisson, L.D., 533 Dawid, A.P., 341, 533 De Santis, A., 76, 531 De Simone, R., 76, 531 decoder, 102, 155, 198 decoding function, 108, 145, 163, 183, 254, 420 decorrelation, 228, 271 deep space communication, 163 delay processing, 427, 478 propagation, 427, 475, 478 transmission, 427, 478, 482 Dembo, A., 410, 541 Dempster, A.P., 533 denominator, 480 dependency graph, 148-149, 165, 175, 176directed edge, 148 dotted edge, 148

parent node, 148

solid edge, 148 descendant, 85, 93 destination, 2 determinant, 227, 234, 443 deterministic distribution, 45 diagonal element, 228, 229, 272, 273, 283, 288, 488 diagonal matrix, 227, 229, 235, 250, 273, 288, 491 diagonalization, 227, 228, 229, 234, 271, 288differential entropy, 230, 225-250, 265, 375 conditional, 236, 251 joint, 234 scaling, 233, 234 translation, 232-234, 265 Diggavi, S.N., 289, 533 digital, 2 digital communication system, 3 directed cycle, 427, 428, 475, 476, 478, 479directed graph, 402, 413 acyclic, 427, 428 cut, 414 capacity of, 415 cyclic, 428, 475 directed cycle, 427, 428 directed path, 428 edge, 402, 413 max-flow, 414 min-cut, 415 node, 402, 413 non-source node, 413 rate constraint, 414, 415 sink node, 401, 414, 416 source node, 401, 413, 416 directed network, 428, 475, see also directed graph directed path, 428, 445, 455 longest, 428 directional derivative, 220, 221 discrete alphabet, 136, 138 discrete channel, 136-141 noise variable, 137, 138 discrete memoryless channel (DMC), 136-145, 163, 172, 174, 207, 251, 280, 425 achievable rate, 147

capacity, 3, 141, 133-177, 207 computation of, 145, 176, 210-214, 218, 223 feedback, 163-168, 172, 174, 425 generic noise variable, 140, 141, 252 symmetric, 173 discrete-time continuous channel, 251, 252noise variable, 251 discrete-time stochastic system, 138 disk array, 418, 498, 525 distortion measure, 179-205 average, 180 context dependent, 181 Hamming, 181, 192, 195 normalization, 181, 191 single-letter, 180 square-error, 181 distortion-rate function, 187 distributed source coding, 410, 528 divergence, 7, 23, 22-26, 43, 45, 46, 245, 245-246, 250, 368 convexity of, 44 divergence inequality, 24, 25, 44, 211, 245, 368 diversity coding, 418, 499 DMC, see discrete memoryless channel Dobrushin, R.L., 240, 530, 533 dominated convergence theorem, 238 double infimum, 209, 217 double supremum, 208, 209, 212, 213 Dougherty, R., 375, 528, 533 duality theorem, 335 dual, 335 primal, 335 Dueck, G., 533 dummy message, 441 dummy node, 406, 407 dyadic distribution, 86 dyadic expansion, 231 ear drum, 372 east-west direction, 209 eavesdropper, 70 edge-disjoint paths, 429, 445, 449, 456, 457, 459, 470, 491 Edmonds, J., 472, 533 efficient source coding, 104-105

Effros, M., 473, 528, 536, 537, 540

Egner, S., 473, 536 eigenvalue, 228, 229, 249 eigenvector, 228, 249 electronic circuit, 482 elemental inequalities, 331, 329-331, 335, 351, 352  $\alpha$ -inequalities, 343–347  $\beta$ -inequalities, 343–347 minimality of, 343-347 Elias, P., 533 EM algorithm, 223 emotion, 1 empirical differential entropy, 243 empirical distribution, 108, 130, 131 joint, 130, 201 empirical entropy, 100, 102, 111 encoder, 102, 155, 198 encoding function, 108, 145, 163, 183, 254, 420 Encyclopedia Britannica, 3, 531 energy constraint, 265 energy signal, 274 engineering, 3, 171 engineering tradeoff, 469 English, entropy rate of, 107 ensemble average, 106 entropic, 316, 355, 386 entropies, linear combination of, 12, 27, 77, 316, 331 entropy, 3, 7, 12, 18, 46, 265, 383, 407 concavity of, 66 relation with groups, 377-397 entropy bound, VIII, 82, 80-83, 86, 87, 91, 94 for prefix code, 91 entropy function, VII, 316, 351, 356, 358, 363, 372, 374, 383 continuity of, 46, 391 discontinuity of, 45, 46 group characterization, 386, 383-387 entropy rate, VIII, 7, 37, 37-40, 179 of English, 107 entropy space, 316, 319, 331, 351, 383 equivalence relation, 303 erasure probability, 144 Erez, E., 494, 533 ergodic, 106 ergodic stationary source, 106, 110, 168 entropy rate, 107

Estrin, D., 536 Euclidean distance, 212, 243, 359 Euclidean space, 316, 352 expected distortion, 182, 215 minimum, 183, 193 exponential distribution, 249 extreme direction, 355, 364, 370 facsimile, 1 fair bits, 104, 231, 401 almost fair bits, 104, 105 fair coin, 79 Fano's inequality, 33, 31-35, 47, 105, 152, 167, 171, 176, 423, 506, 509 simplified version, 34 tightness of, 45 Fano, R.M., 47, 176, 533 fault-tolerant data storage system, 371, 418 fault-tolerant network communication, 498 FCMI, see full conditional mutual independencies Feder, M., 494, 533 feedback, VIII, 136, 172, 175, 176 Feinstein, A., 176, 533 Feldman, J., 533 Feller, W., 534 ferromagnetic material, 306, 312 field, 493, see also finite field real, 427, 452 field size, 448, 449, 451, 458, 460, 464, 465, 471 field, in measure theory, 50 file, 466 finite alphabet, 18, 32, 34, 43, 46, 105, 108, 110, 120, 130, 131, 141, 155, 168, 180, 199, 207, 391, 402, 413 finite duration, 280 finite field, 427, 452, 493 algebra, 427 extension field, 477 finite group, 379, 377-397 finite resolution, 1 finite-dimensional maximization, 207 Fitingof, B.M., 534 fix-free code, 96 flow on a directed graph, 414

conservation conditions, 414

value of, 414 Fong, S.L., 472, 534 Ford, Jr., L.K., 534 Forney, Jr., G.D., 534 forward substitution, 481 fountain code, 470 Fourier transform, 274, 276, 278, 280 inverse, 274, 279 Fragouli, C., 410, 494, 532, 534 Fraleigh, J.B., 534 Freiling, C., 375, 528, 533 frequency, 274 frequency band, 275, 281 frequency component, 276 frequency of error, 181 frequency response, 281 frequency spectrum, 275 Frey, B.J., 538 Fu, F., 73, 534 Fujishige, S., 349, 534 Fulkerson, D.K., 534 full conditional independence, VIII full conditional mutual independencies, 292, 301-305, 341 axiomatization, 312 image of, **302**, 304 set-theoretic characterization, 312 functional dependence, 326 fundamental inequality, VIII, 23, 83, 241, 245 fundamental limits, 3 Gács, P., 532 Gallager, R.G., 162, 173, 176, 280, 289, 534, 542 Galois field, see finite field gamma distribution, 249  $\Gamma_n$ , 331–349, 351  $\Gamma_n^*, \, \mathbf{316}, \, 332, \, 351, \, 377$  $\overline{\Gamma}_n^*, \, 355$ group characterization of, 388-391 Gargano, L., 76, 531 Gauss elimination, 460 Gaussian channel. 264 bandlimited colored Gaussian channel, 280-283, 289 capacity, 283, 289 bandlimited white Gaussian channel,

273-281, 289

Index 553capacity, 276, 289 bandpass white Gaussian channel, 281, 289 capacity, 281 correlated Gaussian channels, 271-273, 287, 289 capacity, 273, 283, 288 noise variable, 271, 272 memoryless Gaussian channel, 264, 263-266, 276, 280, 282 capacity, 264-267, 288 parallel Gaussian channels, 266-272, 289capacity, 270, 273, 288 noise variable, 266 Gaussian distribution, 227, 233, 280 multivariate, 227, 228, 234, 248-250, 272, 279, 285, 375 zero-mean, 273, 279 Gaussian noise, 264, 289 independent, 289 process, 273, 279, 281 zero-mean, 267, 271, 273, 281, 283-287, 289 Ge, Y., 312, 534 generator matrix, 432 generic continuous channel, 252, 253, 264generic discrete channel, 139, 154, 207, 210generic message, 477 generic network code, 452, 452-460, 463, 471, 473 alternative definition, 454 construction, 458 simplified characterization, 471 static, 462, 463, 465, 471 construction, 464 transformation of, 471 geometric distribution, 37 Gersho, A., 534 Gitlin, R.D., 410, 530 Gkantsidis, C., 473, 534 Glavieux, A., 530 global encoding kernel, 432, 437, 440, 442, 443, 445, 446, 449, 451, 454, 458, 459, 464, 472, 473, 475-477, 482, 483, 487, 490, 492, 493

general positions, 452, 458

global Markov property, 306, 311 Goldman, S., 534 Goldsmith, A., 534 Golić, J.Dj., 359, 534 Golomb, S.W., 535 Govindan, R., 536 gradient, 220 Grant, A., 528, 532 graph theory, 306, 402, 405, 413, 414, 429graphical models, VIII, 312 Gray, R.M., 532, 534, 535 group, 378, 377-397 associativity, 378, 379-381, 383 axioms of, 378 closure, 380, 381, 383 identity, 378, 380-384 inverse, 378-381 order of, 377, 379, 381 group inequalities, 377-391, 395, 397 group theory, VII, IX, 132, 356 relation with information theory, 377 - 397group-characterizable entropy function, 386, 383-387 Guiasu, S., 535 Gutman, M., 533 Hadamard's inequality, 250 Hadamard, J., 535 Hajek, B.E., 535 half-space, 319 Hammer, D., 375, 535 Hamming ball, 203 Hamming code, 163 Hamming distance, 46, 176 Hamming distortion measure, 181 Hamming, R.V., 535 Han, T.S., 44, 47, 77, 327, 359, 535, 544 hard disk, 473 hardware failure, 460 Hardy, G.H., 46, 535 Harremoës, P., 26, 535 Hau, K.P., 410, 527, 528, 535, 542 Heegard, C., 535 heuristic argument, 276, 281 hiker, 209 Ho, S.-T., 471-473, 543 Ho, S.-W., 47, 108, 131, 132, 536

Ho, T., 411, 473, 528, 535, 536, 539, 540 Hocquenghem, A., 536, see also BCH code home entertainment systems, 162 Horibe, Y., 536 Hu, G.D., 77, 536 Huffman code, 86, 86–91 expected length, 88, 89 optimality of, 88 Huffman procedure, 86, 86–91 dummy symbols, 87 Huffman, D.A., 97, 536 Hui, J.Y., 536 human factor, 1 hypergraph, 312 hyperplane, 318, 322, 326, 355, 370 Hyvarinen, L.P., 536 I, C.-L., 410, 530 I-Measure, VIII, 56, 49-77, 150, 291-312, 351, 358 empty atom, 54 Markov chain, 58-65, 73, 309-311 Markov structures, 291-312 negativity of, 57-58 nonempty atom, 54 uniqueness, 56, 61 universal set, 51, 54 i.i.d. source, 102, 105, 109, 111, 180, 204, 207 bivariate, 120, 122 Ibinson, B., 536 identity matrix, 439, 472, 488, 489 Ihara, S., 250, 289, 536 image, 180 imaginary channel, 429, 436, 445, 449, 450, 482, 483, 491 imperfect secrecy theorem, 69 implication problem, IX, 326-327, 341-343, 349, 375 involves only FCMI's, 304, 326 impulse response, 275, 289, 482 inclusion-exclusion formula, 53 a variation of, 73 incomplete data, 223 incompressible, 104 independence bound for differential entropy, 242, 258 entropy, 28, 104

independence of random variables, 7-12 mutual, 8, 29, 30, 38, 42, 60, 75, 199, 272, 321, 340 pairwise, 8, 42, 57, 353 indeterminate, 434, 442-447, 491 inferior, 496, 504 infinite group, 379 infinitesimal perturbation, 452 Information Age, 4 information diagram, VIII, 59, 49-77, 337, 342, 362 Markov chain, 61-65, 73, 149, 309-311 information expressions, 313 canonical form, 316-319 alternative, 327 uniqueness, 317, 327 nonlinear, 327 symmetrical, 327 information identities, VII, 27, 77, 313 constrained, 321, 334-335 unconstrained, 319 information inequalities, VII, 27, 66, 313, 377, 391-395, 528 constrained, 320-321, 334-335 equivalence of, 322-325, 328 framework for, IX, 313-328 machine-proving, ITIP, 315, 337-340 non-Shannon-type, IX, 28, 351-375 Shannon-type, IX, 329-349 symmetrical, 348 unconstrained, 319, 333-334, 355, 378, 391 information looping, 476, 478 information rate-distortion function, 188, 198 continuity of, 198 properties of, 189 information source, 2, 37, 80, 179, 259, 401, 407, 420, 427 informational divergence, see divergence Ingleton inequality, 374, 396 Ingleton, A.W., 536 input distribution, 141, 154, 207, 210, 216, 259 strictly positive, 213, 223 input energy, 267 input power allocation, 268, 270, 273,

282, 288

input power constraint, 264, 266, 271, 272, 276, 280, 281, 288 Intanagonwiwat, C., 536 internal node, 84, 84-95 conditional entropy of, 92 inverse function, 480 invertible matrix, 273, 439, 488 invertible transformation, 318 Ising model, 306, 312 iterative algorithm, 176, 205, 207, 216, 223ITIP, IX, 337-340, 349, 359, 373, 375 efficient implementation, 343 Jacquet, P., 536 Jaggi, S., 473, 536, 537 Jaggi-Sanders algorithm, 449, 450-452, 460, 471 Jain, K., 472, 473, 532, 536, 545 Jaynes, E.T., 47, 537 Jelinek, F., 537 Jensen's inequality, 197, 286 Jensen, J.L.W.V., 537 Jerohin, V.D., 204, 537 Jewell, W.S., 543 Johnsen, O., 95, 537 Johnson, R.W., 542 joint entropy, 13, 315 joint source-channel coding, 170, 171 Jones, G.A., 537 Jones, J.M., 537 Kakihara, Y., 537 Karger, D., 528, 540 Karger, D.R., 473, 535, 536, 539 Karush, J., 97, 270, 537 Karush-Kuhn-Tucker (KKT) condition, 270Katabi, D., 537 Katona, G.O.H., 535, 543 Katti, S., 537 Kawabata, T., 61, 77, 312, 537 key, of a cryptosystem, 69, 75 Khachatrian, L., 96, 529 Khinchin, A.I., 537 Kieffer, J.C., 537, 545 Kindermann, R., 537 King, R., 532

Kischischang, F., 537

Kobayashi, K., 535 Koetter, R., 471, 473, 493, 535-537, 539 Kolmogorov complexity, VII, 375, 397 Kolmogorov, A.N., 250, 537 Körner, J., VIII, 77, 132, 328, 529, 532, 533Kraft inequality, VIII, 80, 82, 83, 85, 86, 90, 95-97 Kraft, L.G., 537 Kramer, G., 537 Kschischang, F.R., 538 Kuhn, H.W., 270, 538, see also Karush-Kuhn-Tucker (KKT) condition Kullback, S., 47, 538 Kullback-Leibler distance, see divergence Kung, S.-Y., 472, 545 Kurose. J.F., 538 Kwok, P.-W., 472, 538 Lagrange multipliers, 213, 268 Lagrange's theorem, 383 Laird, N.M., 533 Landau, H.J., 538 Langberg, M., 537, 538 Langdon, G.G., 538 Laplace distribution, 249 large scale content distribution, 466 lattice theory, 77 Lau, L.C., 538 Laurent series, formal, 493 Lauritzen, S.L., 538 laws of information theory, IX, 314, 371, 374Le Boudec, J.-Y., 534 leaf, 84, 84-95 Lebesgue measure, 245, 327, 355 Lebesgue-Stieltjes integration, 225 Lee, T.T., 312, 545 left-continuous, 253, 254, 261 Lehman, E., 528, 541 Leibler, R.A., 47, 538 Lempel, A., 546 Leong, B., 536 letter, 37 Leung, S.K., 532 Li, B., 538 Li, M., 538

Li, S.-Y.R., 409, 410, 424, 425, 473, 493, 494, 529, 538, 545 Li, Z., 538 Liang, X.-B., 538 Lieb, E.H., 375, 538 Lin, S., 538 Linden, N., 375, 536, 538 Linder, T., 97, 538 line of sight, 500 linear broadcast, 436, 437, 440, 442, 446, 452, 454, 476 multi-rate, 437, 472 static, 462, 465 linear code, 162 linear constraints, 321, 323 linear dispersion, 436, 437, 440, 442, 447, 452, 454, 457, 471 static, 462, 465 linear mapping nullity, 465 pre-image, 465 linear multicast, 436, 437, 440, 442, 445, 451, 452, 454, 460, 466, 468, 471, 473, 489, 492, 493 construction, see Jaggi-Sanders algorithm random, 448 static, 462, 465, 471 transformation of, 451 linear network code, 429-470, 473, 477 base field, 427 dimension, 431 global description, 432, 475 implementation of, 440-441, 468 overhead, 441 linear broadcast, 436 linear dispersion, 436 linear multicast, 436 local description, 431, 443, 475 transformation of, 439-440 linear network coding, 413 matrix approach, 473 vector space approach, 473 linear programming, IX, 329, 331-335, 337, 349 linear span, 434 linear subspace, 321, 364, 370 linear time-invariant (LTI) system, 482, 493

causality, 482 linear transformation, 435 linear transformation of random variables, 227-230, 249, 272 Littlewood, J.E., 46, 535 Lněnička, R., 375, 539 local area network (LAN), 437 local encoding kernel, **431**, 432, 442, 443, 446, 447, 463, 465, 466, 472, 473, 475-477, 479, 482, 483, 485, 487, 488, 490-492 local Markov property, 311 local redundancy, 94 local redundancy theorem, 94–95 Loeliger, H.-A., 538 log-optimal portfolio, 223 log-sum inequality, 24, 44, 222 logical network, 466 long division, 480, 481, 493 Longo, G., 530 lossless data compression, 3, 97 Lovász, L., 539 low-density parity-check (LDPC) code, 162Luby, M., 531 Lun, D.S., 536, 539 MacKay, D.J.C., 162, 539 majority vote, 135 majorization, 46 Makarychev, K., 374, 539 Makarychev, Y., 374, 539 Malkin, T., 533 Malvestuto, F.M., 312, 539 Mann, H.B., 530 Mansuripur, M., 539 mapping approach, 205 marginal distribution, 260, 360, 361 Marko, H., 175, 539 Markov chain, VIII, 7, 9, 30, 61, 62, 65, 66, 69, 71, 73, 139, 149, 167, 172, 174, 252, 256, 291, 292, 306, 309-311, 325, 337, 339, 360, 361, 372, 375 information diagram, 61-65, 73, 149, 309 - 311Markov graph, 306 Markov random field, VIII, 65, 292, 306-308, 312

hypergraph characterization of, 312 Markov star. 312 Markov structures, VIII, 291-312 Markov subchain, 10 Marshall, A.W., 539 Marton, K., 539 Massey, J.L., 70, 175, 539 Mathai, A.M., 539 MATLAB, 337 matroid, 374, 528 Matúš, F., 326, 359, 374, 375, 539, 540 Maurer, U.M., 531, 540, 544 max-flow, 414, 423, 434, 436, 437, 451, 460, 461, 468, 469 collection of edges, 415 collection of non-source nodes, 415 max-flow bound, VIII, 421, 413-423, 427, 451, 454, 473, 493, 495 for linear network coding, 435 max-flow bounds, 495-498, 525 max-flow min-cut theorem, 415, 423, 435maximal probability of error, 146, 154, 170, 176, 255 maximization, 269 maximum differential entropy, 246-248 maximum entropy, 35-37, 250 maximum likelihood decoding, 175 Mazo, J., 410, 530 McEliece, R.J., 540 McGill, W.J., 77, 540 McLaughlin, S.W., 544 McMillan, B., 97, 110, 540, see also Shannon-McMillan-Breiman theorem mean ergodic, 106 mean-square error, 181 meaningful information, 1 measure theory, 50, 106, 225, 250 membership table, 384 Merhav, N., 533 message, 417 message pipeline, 427, 485, 486, 489 message set, 133, 145, 146, 254, 255 method of types, 115, 132 microelectronics, 162 min-cut, 415, 423, 470 minimization, 333-335 minimum distance decoding, 176

Mittelholzer, T., 531, 544 Mitter, S., 543 Mitzenmacher. M., 531, 540 mixing random variable, 67, 258 modulo 2 addition, 379-380, 386, 387, 402, 410 modulo 2 arithmetic, 477 modulo 3 addition, 410 Mohan, S., 530 most likely sequence, 102 Moulin, P., 540 Moy, S.C., 540  $\mu^*$ , see *I*-Measure multi-dimensional direction, 209 multi-source multicast, 451 multi-source network coding, IX, 408, 495 - 528achievable information rate region, 495LP bound, 505 information rate region, 528 insufficiency of linear coding, 528 network code for acyclic network, 500 source separation, 525, 528 multicast, VIII, IX, 401, 402, 413, 427, 495 - 528multigraph, 413 multilevel diversity coding, 498-499, 528symmetrical, 499, 528 multiple descriptions, 73 multiple unicasts, 404 multiterminal source coding, 132, 205 Murty, U.S.R., 306, 531 mutual information, 7, 15, 215, 238, 239, 250, 259 between more than two random variables. 58 concavity of, 68, 253, 259, 288 convexity of, 67, 190 mutual typicality, 259-260 typical sequence, 260 typical set, 259 mutually independent information sources, 495-525 Médard, M., 471, 473, 494, 528, 535-537, 539, 540

nat, 13, 232 natural disasters, 460 neighborhood, 243 neighboring node, 405 nerve impulse, 372 network code deployment, 460 global description, 430 global encoding mapping, 430, 470 local description, 430 local encoding mapping, 430, 470 network coding, VII, VIII, 495-528 advantage of, 402-405, 410 source separation, 407–408, 410 Network Coding Homepage, 411 network communication, 401, 402, 405, 407 network error correction, 473 network topology, 441, 460, 472 unknown, 437, 448, 466, 473 network transfer matrix, 470 Nielsen, M.A., 375, 540 Nobel, A., 410, 541 noise energy, 264, 267 noise power, 264, 271, 275, 280 noise process, 275, 279, 281 noise source, 2 noise variable, 251, 266, 271, 272 noise vector, 271, 272, 283, 287-289 noisy channel, 3, 133, 160 noisy environment, 1 non-decreasing, 253 non-increasing, 187 non-Shannon-type inequalities, IX, 28, **314**, 315, 337, 351–375, 528 constrained, 364-370 unconstrained, 359-364, 378, 394, 528 nonlinear optimization, 207 nonnegative linear combination, 335, 336 nonnegative orthant, 316, 318, 331, 352, 359 normal distribution, see Gaussian distribution north-south direction, 209 null space, 323 numerator, 492 numerical computation, 145, 196,

207 - 223

Narayan, P., 115, 532

Nyquist, H., 540 O'Sullivan, J.A., 540 off-diagonal element, 488 Olkin, I., 539 Omura, J.K., 540, 544 Ooi, J.M., 540 optimal coding scheme, 3 order of a node, 85 ordinate, 215 Orlitsky, A., 540 Ornstein, D.S., 540 orthogonal complement, 323, 324 orthogonal matrix, 227, 230 orthogonal transformation, 228, 230 orthonormal basis, 277, 278, 280 orthonormal set, 278 orthonormal system, 228 output channel, 402 overlay network, 466 Oxley, J.G., 540 P2P, see peer-to-peer network packet loss, 469 rate, 469 packet network, 469 Papadimitriou, C.H., 540 Papoulis, A., 77, 541 parallel channels, 173, 349 parity check, 163 partition, 303 pdf, see probability density function Pearl, J., 541 peer-to-peer (P2P) network, 466, 473 client, 466 neighboring node, 466, 467, 469 server, 466 tracker. 466 Peile, R.E., 535 perceived distortion, 180 Perez, A., 541 perfect secrecy theorem, Shannon's, VIII, 70 permutation, 380 Perrin, D., 530 physical network, 479 physical system, 4 physics, VII Pierce, J.R., 541

Pinkston, J.T., 205, 541 Pinsker's inequality, 25, 26, 44, 46, 47, 115Pinsker, M.S., 47, 250, 289, 530, 541 Pippenger, N., 374, 375, 541 plain text, 69, 75 point-to-point channel, 133 noiseless, 401, 413, 425 capacity, 413 point-to-point communication network, VIII, 402, 413, 416, 427, 495 point-to-point communication system, 2,407 Pollak, H.O., 538, 543 Polya, G., 46, 535 polymatroid, 346, 349, 374 polynomial, 327, 442, 443, 471, 480, 481, 490-493 equation, 443 degree, 443 nonzero, 443, 446 root, 443 polynomial ring, 443, 445, 491 positive definite matrix, 227, 249, 284 positive semidefinite matrix, 227, 228, 229, 250 postal system, 401 power series, 479 expansion, 481 formal, 480 rational, 480, 482, 488, 491 ring of, 480, 485 ring of, 480, 485 power spectral density, 274, 275, 279, 281prefix code, VIII, 80, 83, 83-95 entropy bound, VIII existence of, 85 expected length, 93 random coding, 96 redundancy, VIII, 91-95 prefix-free code, see prefix code Preston, C., 312, 541 prime number, 443 probabilistic coding, 108, 172 probability density function (pdf), 225 bounded, 236 conditional, 226, 235, 236 bounded, 236, 238

joint, 225, 283-285, 287 probability distribution rational, 388 strictly positive, 7, 11, 210, 348 factorization of, 12 with zero masses, 7, 210, 312 probability of error, 33, 102, 134, 153, 181. 187 probability theory, VII, 326, 371 product measure, 240 product source, 204, 349 projection, 504 prolate spheroidal wave functions, 280 pyramid, 331, 332, 335 quantization, 239 quantized samples, 180 quantum information theory, 375 quantum mechanics, 375 quasi-uniform structure, 385 asymptotic, 128, 385 Rabin, M.O., 410, 541 Radon-Nikodym derivative, 240 random code, 155, 198, 262, 527 random coding error exponent, 176 random linear combination, 467, 468 random network coding, 448, 466-470, 473robustness, 469 random noise, 133 random variable, real, 35, 225-289 continuous, 225 discrete, 225 mixed, 225 second moment, 230, 280 rank function, 397 rank of a matrix, 323 full, 324, 325, 441 Rasala Lehman, A., 528, 541 rate constraint, 402, 413, 414, 416, 421 rate-distortion code, 183, 179-205, 207 rate-distortion function, 187, 183-188,

198, 204, 207 binary source, 192 forward channel description, 203 reverse channel description, 194 computation of, VIII, 196, 205, 215–218, 223

normalization, 205 product source, 204, 349 properties of, 187 Shannon lower bound, 204 rate-distortion pair, 184, 196 rate-distortion region, 184, 187, 215 rate-distortion theorem, VIII, 132, 189, 188-196, 205 achievability, 198-203 converse, 196-198 random code, 198 relation with source coding theorem, 195rate-distortion theory, VIII, 179-205 Rathie, P.N., 539 rational function, 480, 492 field of, 491 rational number, 185, 388 Ratnakar, N., 539 Ray-Chaudhuri, D.K., 531, see also BCH code Rayleigh's energy theorem, 277 reaching probability, 92, 93 real number, 427 receiver, 2 reciprocal, 481 rectangular lattice, 306, 312 reduced code tree, 89 reduced probability set, 89 redundancy of prefix code, VIII, 91–95 of uniquely decodable code, 83 Reed, I.S., 541 Reed-Solomon code, 163 relative entropy, see divergence relative frequency, 111, 120 relay node, 406, 407 Rényi, A., 541 repetition code, 135 replication of information, 417 reproduction alphabet, 180, 192, 199 reproduction sequence, 179-181, 186, 199reservoir, 271, 283 resultant flow, 414 Reza, F.M., 77, 541 right-continuous, 225 Riis, S., 528, 541

ring, 443, 445, 480, 492

commutative, 481 Rissanen, J., 541 Roche, J.R., 410, 527, 528, 541, 542 Rockafellar, R.T., 542 Rodriguez, P.R., 473, 534 Roman, S., 542 Romashchenko, A., 374, 375, 397, 535, 539.542 Rose, K., 205, 542 Ross, K.W., 538 routing, 401, 401, 402, 417 row space, 323, 324 Rubin, D.B., 533 Rudin, W., 542 Ruskai, M.B., 375, 538 Ruskey, F., 542 Russian, 77 Rustin, R., 533 sampling theorem, 276, 278 bandpass, 281 sampling time, 279 Sanders, P., 473, 536, see also Jaggi-Sanders algorithm Santhanam, N.P., 540 satellite communication, 402, 405-407, 410 satellite communication network, 498, 500Savari, S.A., 537, 542 Scholtz, R.A., 535 Schur-concave function, 46 science, 3 science of information, the, 3 secret key cryptosystem, 69, 75 secret sharing, 75, 76, 339, 473 access structure, 76 information-theoretic bounds, 75-76, 340 participants, 76 secure network coding, 473 security level of cryptosystem, 70 self-information, 16 semi-graphoid, 349 axioms of, 341 separation of network and channel coding, 425 source and channel coding, 136, 168-171, 204, 401

Servedio, R.A., 533 set function. 316 additive, 50, 73, 301 set identity, 53, 77, 297 set operations, 49, 50 set theory, VIII, 49 Shamai, S., 542, 544 Shamir, A., 542 Shannon code, 91 Shannon's information measures, VIII, 7, 12-18, 27, 49 continuity of, 18-20 discontinuity of, 20 elemental forms, 330, 348 irreducible, **329**, 347 linear combination of, 313 reducible, 329, 347 set-theoretic structure of, see *I*-Measure Shannon's papers, collection of, 4 Shannon, C.E., VII, 2, 46, 97, 102, 110, 176, 204, 205, 250, 289, 349, 374, 542Shannon-McMillan-Breiman theorem, VIII, 40, 106, 105-107, 169 Shannon-type identities constrained, 334-335 Shannon-type inequalities, IX, 314, 315, 329-349, 359, 528 constrained, 334-335 machine-proving, ITIP, 329, 337-340, 351unconstrained, 333-334 Shen, A., 375, 397, 535, 542 Shields, P.C., 542 shift-register, 482, 485, 493 Shore, J.E., 542 Shtarkov, Y.M., 544 Shunsuke, I., 542 sibling, 88 side-information, 205, 408 signal, 133 signal analysis, 274, 276, 480 signal-to-noise ratio, 266 signaling network, 462 signed measure, 50, 56, 57 Simmons, G.J., 539 Simonnard, M., 543 simplex method, 334, 335

optimality test, 334, 335 sinc function, 276, 279 single-input single-output system, 133, 136, 138 single-letter characterization, 207 single-letter distortion measure, 207 single-source network code a class of, 419 causality, 419, 423 single-source network coding, 410, 413, 495achievable information rate, 421 acyclic network, 427-470 cyclic network, 475-493 one sink node, 416 three sink nodes, 417 two sink nodes, 417 sink node, 401, 414, 416, 417, 451 Slepian, D., 205, 543 Slepian-Wolf coding, 205 Sloane, N.J.A., 531, 543 Snell, J., 537 Soljanin, E., 411, 494, 532, 534 Solomon, G., 541, see also Reed-Solomon code Song, L., 425, 528, 543 sound wave, 372 source code, 80, 171, 179 source coding theorem, VIII, 3, 102, 102-103, 110, 179, 187, 195 coding rate, 102 converse, 103 direct part, 102 general block code, 108 source node, 401, 413, 416 super, 405 source random variable, 204 source sequence, 179-181, 186, 198 space-time domain, 478 spanning tree packing, 472 Spitzer, F., 312, 543 Sprintson, A., 538 standard basis, 432, 435, 446, 450, 482 standard deviation, 248 static network code, 461, 460-466, 473 configuration, 460 generic, 462, see also generic network code, static

linear broadcast, 462, see also linear broadcast, static linear dispersion, 462, see also linear dispersion, static linear multicast, 462, see also linear multicast, static robustness, 460 stationary source, VIII, 38, 106 entropy rate, 7, 37-40 Steiglitz, K., 540 Stein, C., 533 Steinberg, Y., 544 still picture, 2 Stinson, D.R., 543 Stirling's approximation, 123 stock market, 223 store-and-forward, 401, 410, 467, 469, 472strong asymptotic equipartition property (AEP), 99, 112, 111-119, 130, 200 strong law of large numbers, 106 strong typicality, VIII, 111-132, 199, 242alternative definition, 130 consistency, 120, 154, 200 joint, 120-129 joint AEP, 122 joint typicality array, 127, 385 jointly typical sequence, 120 jointly typical set, 120 typical sequence, 111 typical set, 111 vs. weak typicality, 119 Studený, M., 77, 349, 374, 375, 539, 540, 543 sub-channel, 281 subcode, 186 subgroups, IX, 377-397 intersection of, 377, 383 membership table, 384 subnetwork, 461 subring, 480 substitution of symbols, 53 suffix code, 95 summit of a mountain, 209 support, 7, 13, 22, 35, 43, 108, 111, 180, **225**, 285, 289, 354, 386 supremum, 253, 267

switching theory, 401 symmetric group, 380, 390 symmetric matrix, 227, 228, 249, 284 synchronous transmission, 425 Szpankowski, W., 536, 543 Tan, M., 471-473, 543 Taneja, I.J., 543 tangent, 215 Tardos, G., 535, 543 Tarokh, V., 97, 538 Tatikonda, S., 543 Telatar, I.E., 543 telephone conversation, 163 telephone line, 1, 163 television broadcast channel. 2 ternary channel, 174 thermodynamics, 46 Thitimajshima, P., 530 Thomas, J.A., VII, 532 Thomasian, A.J., 530 time average, 106 time domain, 276, 483, 485 time-sharing, 185 Tjalkens, T.J., 544 Toledo, A.L., 543 Tolhuizen, L., 473, 536 Topsöe, F., 535, 543 transfer function, 493 transition matrix, 136, 172, 190, 207, 210, 215, 216 strictly positive, 217 transmitter, 2 trellis network, 478-480, 487 acyclicity of, 478 triangular inequality, 23, 43 Tse, D.N.C., 543, 546 Tucker, A.W., 270, 538, see also Karush-Kuhn-Tucker (KKT) condition Tunstall, B.P., 543 turbo code, 162 Tusnády, G., 223, 532 Type I atom, 307 Type II atom, 307 type of a sequence, 131

uncorrelated random variables, 228, 229, 234, 272, 279 undirected graph, 306 component, 306 cutset, 306 edge, 306 loop, 306 vertex, 306 unified asymptotic equipartition property (AEP), 131 unified typicality, 131, 132 consistency, 132 uniform distribution, 35, 37, 143, 145, 146, 155, 204, 231, 255, 258, 261, 354, 358, 386, 419, 444 union bound, 114, 156, 169, 187 unique solution, 475, 476 uniquely decodable code, VIII, 80, 83, 85-87, 91, 95, 97 expected length, 81 redundancy, 83 unit-delay network, 478, 478–493 universal source coding, 109 upstream-to-downstream order, 429, 432, 443, 449, 459, 464, 470, 475 Vaccaro, U., 76, 531 van der Lubbe, J.C.A., 544 van der Meulen, E.C., 544 van Dijk, M., 76, 544 variable-length channel code, 168 variance, 226, 229, 247, 279 variational distance, 18, 20, 25, 26, 43-46, 130, 391 vector space, 397 Vembu, S., 544 Venn diagram, 16, 50, 59, 77 Verdú, S., 47, 530, 535, 542, 544 Vereshchagin, N., 374, 539 Vereshchagin, N.K., 375, 397, 535, 542 video signal, 180 Viswanath, P., 543 Viswanathan, H., 530 Vitányi, P., 538 Viterbi, A.J., 544 von Neumann entropy, 375 strong subadditivity, 375 von Neumann, J., 544

uncertainty, 2, 13

Wald, A., 544 Wang, X., 543 water flow, 414 water leakage, 414 water pipe, 160, 414 water-filling, 271, 273, 283, 289 waveform channel, 251, 273, 275, 276, 278-281, 289 weak asymptotic equipartition property (AEP), VIII, 99, 99-110, 242, see also Shannon-McMillan-Breiman theorem weak independence, 75 weak law of large numbers, 99, 100, 135, 260, 262 weak typicality, VIII, 99-111, 119, 131, 132, 242 alternative definition, 108 typical sequence, 100, 100-110, 242 typical set, **100**, 100–110 Weaver, W.W., 542 Wegener, I., 529 Weingarten, H., 544 Welch, T.A., 544 Welsh, D.J.A., 536 Wheeler, D.J., 531 Whittaker, E.T., 544 Wicker, S.B., 535, 544 wide-sense stationary process, 274, 275, 288Widmer, J., 534 Willems, F.M.J., 544 Winter, A., 375, 536, 538 wired-line communication, 275 wireless communication, 162, 275, 402, 405 - 407

Wolf, J.K., 205, 543, see also Slepian-Wolf coding Wolfowitz, J., 132, 176, 529, 544 Woodard, P.M., 544 World Wide Web, 337 Wu, Y., 411, 472, 473, 532, 545 Wyner, A.D., 289, 543, 545 WYSIWYG, 66 Yan, X., 473, 528, 530, 545 Yan, Y.-O., 349, 369, 545 Yang, E.-h., 537, 545 Yang, S., 545 Ye, C., 96, 545 Ye, Z., 312, 534, 545 Yeung, R.W., 73, 75, 77, 95-97, 108, 115, 131, 132, 223, 312, 328, 349, 374, 409, 410, 424, 425, 471-473, 493, 494, 527-532, 534, 536-538, 542, 543, 545, 546 Ytrehus, Ø., 494, 530 z-transform, 479, 480, 485, 487 dummy variable, 479 Zamir, R., 73, 542 Zeger, K., 97, 375, 528, 533, 538 zero mean, 247, 248, 250, 267, 271 zero-error data compression, VIII, 79-97, 99zero-error reconstruction, 420 Zhang, J., 540 Zhang, Z., 341, 374, 410, 473, 494, 528, 530, 545, 546 Zhao, F., 539 Zheng, L., 546 Zigangirov, K.Sh., 546 Zimmerman, S., 97, 546 Ziv, J., 545, 546