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Victor Klee, Editor

COMPUTERS AND INTRACTABILITY

A Guide to the Theory of NP-Completeness

Michael R. Garey / David S. Johnson

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Contents

Preface	ix
1 Computers, Complexity, and Intractability	1
1.1 Introduction	1
1.2 Problems, Algorithms, and Complexity	4
1.3 Polynomial Time Algorithms and Intractable Problems	6
1.4 Provably Intractable Problems	11
1.5 NP-Complete Problems	13
1.6 An Outline of the Book	14
2 The Theory of NP-Completeness	17
2.1 Decision Problems, Languages, and Encoding Schemes	18
2.2 Deterministic Turing Machines and the Class P	23
2.3 Nondeterministic Computation and the Class NP	27
2.4 The Relationship Between P and NP	32
2.5 Polynomial Transformations and NP-Completeness	34
2.6 Cook's Theorem	38
3 Proving NP-Completeness Results	45
3.1 Six Basic NP-Complete Problems	46
3.1.1 3-SATISFIABILITY	48
3.1.2 3-DIMENSIONAL MATCHING	50
3.1.3 VERTEX COVER and CLIQUE	53
3.1.4 HAMILTONIAN CIRCUIT	56
3.1.5 PARTITION	60
3.2 Some Techniques for Proving NP-Completeness	63
3.2.1 Restriction	63
3.2.2 Local Replacement	66
3.2.3 Component Design	72
3.3 Some Suggested Exercises	74

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4 Using NP-Completeness to Analyze Problems	77
4.1 Analyzing Subproblems	80
4.2 Number Problems and Strong NP-Completeness	90
4.2.1 Some Additional Definitions	92
4.2.2 Proving Strong NP-Completeness Results	95
4.3 Time Complexity as a Function of Natural Parameters	106
5 NP-Hardness	109
5.1 Turing Reducibility and NP-Hard Problems	109
5.2 A Terminological History	118
6 Coping with NP-Complete Problems	121
6.1 Performance Guarantees for Approximation Algorithms	123
6.2 Applying NP-Completeness to Approximation Problems	137
6.3 Performance Guarantees and Behavior "In Practice"	148
7 Beyond NP-Completeness	153
7.1 The Structure of NP	154
7.2 The Polynomial Hierarchy	161
7.3 The Complexity of Enumeration Problems	167
7.4 Polynomial Space Completeness	170
7.5 Logarithmic Space	177
7.6 Proofs of Intractability and P vs. NP	181
Appendix: A List of NP-Complete Problems	187
A1 Graph Theory	190
A1.1 Covering and Partitioning	190
A1.2 Subgraphs and Supergraphs	194
A1.3 Vertex Ordering	199
A1.4 Iso- and Other Morphisms	202
A1.5 Miscellaneous	203
A2 Network Design	206
A2.1 Spanning Trees	206
A2.2 Cuts and Connectivity	209
A2.3 Routing Problems	211
A2.4 Flow Problems	214
A2.5 Miscellaneous	218
A3 Sets and Partitions	221
A3.1 Covering, Hitting, and Splitting	221
A3.2 Weighted Set Problems	223
A4 Storage and Retrieval	226
A4.1 Data Storage	226
A4.2 Compression and Representation	228
A4.3 Database Problems	232

A5 Sequencing and Scheduling	236
A5.1 Sequencing on One Processor	236
A5.2 Multiprocessor Scheduling	238
A5.3 Shop Scheduling	241
A5.4 Miscellaneous	243
A6 Mathematical Programming	245
A7 Algebra and Number Theory	249
A7.1 Divisibility Problems	249
A7.2 Solvability of Equations	250
A7.3 Miscellaneous	252
A8 Games and Puzzles	254
A9 Logic	259
A9.1 Propositional Logic	259
A9.2 Miscellaneous	261
A10 Automata and Language Theory	265
A10.1 Automata Theory	265
A10.2 Formal Languages	267
A11 Program Optimization	272
A11.1 Code Generation	272
A11.2 Programs and Schemes	275
A12 Miscellaneous	279
A13 Open Problems	285
Symbol Index	289
Reference and Author Index	291
Subject Index	327
Update for the Current Printing	339

Preface

Few technical terms have gained such rapid notoriety as the appellation "NP-complete." In the short time since its introduction in the early 1970's, this term has come to symbolize the abyss of inherent intractability that algorithm designers increasingly face as they seek to solve larger and more complex problems. A wide variety of commonly encountered problems from mathematics, computer science, and operations research are now known to be NP-complete, and the collection of such problems continues to grow almost daily. Indeed, the NP-complete problems are now so pervasive that it is important for anyone concerned with the computational aspects of these fields to be familiar with the meaning and implications of this concept.

This book is intended as a detailed guide to the theory of NP-completeness, emphasizing those concepts and techniques that seem to be most useful for applying the theory to practical problems. It can be viewed as consisting of three parts.

The first part, Chapters 1 through 5, covers the basic theory of NP-completeness. Chapter 1 presents a relatively low-level introduction to some of the central notions of computational complexity and discusses the significance of NP-completeness in this context. Chapters 2 through 5 provide the detailed definitions and proof techniques necessary for thoroughly understanding and applying the theory.

The second part, Chapters 6 and 7, provides an overview of two alternative directions for further study. Chapter 6 concentrates on the search for efficient "approximation" algorithms for NP-complete problems, an area whose development has seen considerable interplay with the theory of NP-completeness. Chapter 7 surveys a large number of theoretical topics in computational complexity, many of which have arisen as a consequence of previous work on NP-completeness. Both of these chapters (especially Chapter 7) are intended solely as introductions to these areas, with our expectation being that any reader wishing to pursue particular topics in more detail will do so by consulting the cited references.

The third and final part of the book is the Appendix, which contains an extensive list (more than 300 main entries, and several times this many results in total) of NP-complete and NP-hard problems. Annotations to the main entries discuss what is known about the complexity of subproblems and variants of the stated problems.

The book should be suitable for use as a supplementary text in courses on algorithm design, computational complexity, operations research, or combinatorial mathematics. It also can be used as a starting point for seminars on approximation algorithms or computational complexity at the graduate or advanced undergraduate level. The second author has used a preliminary draft as the basis for a graduate seminar on approximation algorithms, covering Chapters 1 through 5 in about five weeks and then pursuing the topics in Chapter 6, supplementing them extensively with additional material from the references. A seminar on computational complexity might proceed similarly, substituting Chapter 7 for Chapter 6 as the initial access point to the literature. It is also possible to cover both chapters in a combined seminar.

More generally, the book can serve both as a self-study text for anyone interested in learning about the subject of NP-completeness and as a reference book for researchers and practitioners who are concerned with algorithms and their complexity. The list of NP-complete problems in the Appendix can be used by anyone familiar with the central notions of NP-completeness, even without having read the material in the main text. The novice can gain such familiarity by skimming the material in Chapters 1 through 5, concentrating on the informal discussions of definitions and techniques, and returning to the more formal material only as needed for clarification. To aid those using the book as a reference, we have included a substantial number of terms in the Subject Index, and the extensive Reference and Author Index gives the sections where each reference is mentioned in the text.

We are indebted to a large number of people who have helped us greatly in preparing this book. Hal Gabow, Larry Landweber, and Bob Tarjan taught from preliminary versions of the book and provided us with valuable suggestions based on their experience. The following people read preliminary drafts of all or part of the book and made constructive comments: Al Aho, Shimon Even, Ron Graham, Harry Hunt, Victor Klee, Albert Meyer, Christos Papadimitriou, Henry Pollak, Sartaj Sahni, Ravi Sethi, Larry Stockmeyer, and Jeff Ullman. A large number of researchers, too numerous to mention here (but see the Reference and Author Index), responded to our call for NP-completeness results and contributed toward making our list of NP-complete problems as extensive as it is. Several of our colleagues at Bell Laboratories, especially Brian Kernighan, provided invaluable assistance with computer typesetting on the UNIX® system. Finally, special thanks go to Jeanette Reinbold, whose facility with translating our handwritten hieroglyphics into faultless input to the typesetting system made the task of writing this book so much easier.

Murray Hill, New Jersey
October, 1978

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COMPUTERS AND INTRACTABILITY

A Guide to the Theory of NP-Completeness

Computers, Complexity, and Intractability

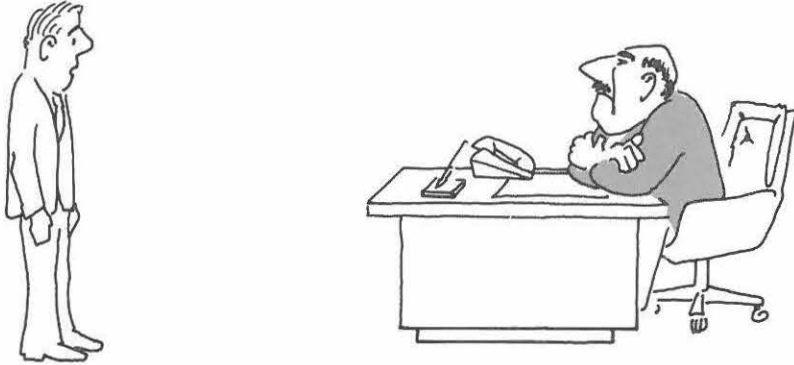
1.1 Introduction

The subject matter of this book is perhaps best introduced through the following, somewhat whimsical, example.

Suppose that you, like the authors, are employed in the halls of industry. One day your boss calls you into his office and confides that the company is about to enter the highly competitive “bandersnatch” market. For this reason, a good method is needed for determining whether or not any given set of specifications for a new bandersnatch component can be met and, if so, for constructing a design that meets them. Since you are the company’s chief algorithm designer, your charge is to find an efficient algorithm for doing this.

After consulting with the bandersnatch department to determine exactly what the problem is, you eagerly hurry back to your office, pull down your reference books, and plunge into the task with great enthusiasm. Some weeks later, your office filled with mountains of crumpled-up scratch paper, your enthusiasm has lessened considerably. So far you have not been able to come up with any algorithm substantially better than searching through all possible designs. This would not particularly endear you to your boss, since it would involve years of computation time for just one set of

specifications, and the bandersnatch department is already 13 components behind schedule. You certainly don't want to return to his office and report:



"I can't find an efficient algorithm, I guess I'm just too dumb."

To avoid serious damage to your position within the company, it would be much better if you could prove that the bandersnatch problem is *inherently* intractable, that no algorithm could possibly solve it quickly. You then could stride confidently into the boss's office and proclaim:



"I can't find an efficient algorithm, because no such algorithm is possible!"

Unfortunately, proving inherent intractability can be just as hard as finding efficient algorithms. Even the best theoreticians have been stymied in their attempts to obtain such proofs for commonly encountered hard problems. However, having read this book, you have discovered something

almost as good. The theory of NP-completeness provides many straightforward techniques for proving that a given problem is "just as hard" as a large number of other problems that are widely recognized as being difficult and that have been confounding the experts for years. Armed with these techniques, you might be able to prove that the bandersnatch problem is NP-complete and, hence, that it is equivalent to all these other hard problems. Then you could march into your boss's office and announce:



"I can't find an efficient algorithm, but neither can all these famous people."

At the very least, this would inform your boss that it would do no good to fire you and hire another expert on algorithms.

Of course, our own bosses would frown upon our writing this book if its sole purpose was to protect the jobs of algorithm designers. Indeed, discovering that a problem is NP-complete is usually just the beginning of work on that problem. The needs of the bandersnatch department won't disappear overnight simply because their problem is known to be NP-complete. However, the knowledge that it is NP-complete does provide valuable information about what lines of approach have the potential of being most productive. Certainly the search for an efficient, exact algorithm should be accorded low priority. It is now more appropriate to concentrate on other, less ambitious, approaches. For example, you might look for efficient algorithms that solve various special cases of the general problem. You might look for algorithms that, though not guaranteed to run quickly, seem likely to do so most of the time. Or you might even relax the problem somewhat, looking for a fast algorithm that merely finds designs that

meet *most* of the component specifications. In short, the primary application of the theory of NP-completeness is to assist algorithm designers in directing their problem-solving efforts toward those approaches that have the greatest likelihood of leading to useful algorithms.

In the first chapter of this “guide” to NP-completeness, we introduce many of the underlying concepts, discuss their applicability (as well as give some cautions), and outline the remainder of the book.

1.2 Problems, Algorithms, and Complexity

In order to elaborate on what is meant by “inherently intractable” problems and problems having “equivalent” difficulty, it is important that we first agree on the meaning of several more basic terms.

Let us begin with the notion of a problem. For our purposes, a *problem* will be a general question to be answered, usually possessing several *parameters*, or free variables, whose values are left unspecified. A problem is described by giving: (1) a general description of all its parameters, and (2) a statement of what properties the answer, or *solution*, is required to satisfy. An *instance* of a problem is obtained by specifying particular values for all the problem parameters.

As an example, consider the classical “traveling salesman problem.” The parameters of this problem consist of a finite set $C = \{c_1, c_2, \dots, c_m\}$ of “cities” and, for each pair of cities c_i, c_j in C , the “distance” $d(c_i, c_j)$ between them. A solution is an ordering $\langle c_{\pi(1)}, c_{\pi(2)}, \dots, c_{\pi(m)} \rangle$ of the given cities that minimizes

$$\left(\sum_{i=1}^{m-1} d(c_{\pi(i)}, c_{\pi(i+1)}) \right) + d(c_{\pi(m)}, c_{\pi(1)})$$

This expression gives the length of the “tour” that starts at $c_{\pi(1)}$, visits each city in sequence, and then returns directly to $c_{\pi(1)}$ from the last city $c_{\pi(m)}$.

One instance of the traveling salesman problem, illustrated in Figure 1.1, is given by $C = \{c_1, c_2, c_3, c_4\}$, $d(c_1, c_2) = 10$, $d(c_1, c_3) = 5$, $d(c_1, c_4) = 9$, $d(c_2, c_3) = 6$, $d(c_2, c_4) = 9$, and $d(c_3, c_4) = 3$. The ordering $\langle c_1, c_2, c_4, c_3 \rangle$ is a solution for this instance, as the corresponding tour has the minimum possible tour length of 27.

Algorithms are general, step-by-step procedures for solving problems. For concreteness, we can think of them simply as being computer programs, written in some precise computer language. An algorithm is said to *solve* a problem Π if that algorithm can be applied to any instance I of Π and is guaranteed always to produce a solution for that instance I . We emphasize that the term “solution” is intended here strictly in the sense introduced above, so that, in particular, an algorithm does not “solve” the traveling

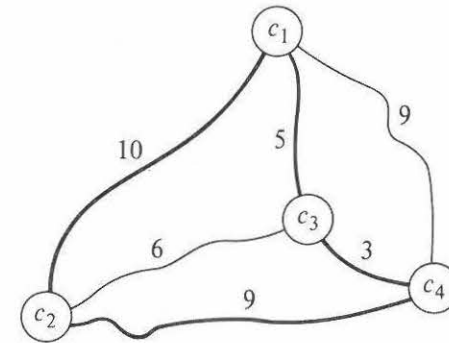


Figure 1.1 An instance of the traveling salesman problem and a tour of length 27, which is the minimum possible in this case.

salesman problem unless it always constructs an ordering that gives a minimum length tour.

In general, we are interested in finding the most “efficient” algorithm for solving a problem. In its broadest sense, the notion of efficiency involves all the various computing resources needed for executing an algorithm. However, by the “most efficient” algorithm one normally means the fastest. Since time requirements are often a dominant factor determining whether or not a particular algorithm is efficient enough to be useful in practice, we shall concentrate primarily on this single resource.

The time requirements of an algorithm are conveniently expressed in terms of a single variable, the “size” of a problem instance, which is intended to reflect the amount of input data needed to describe the instance. This is convenient because we would expect the relative difficulty of problem instances to vary roughly with their size. Often the size of a problem instance is measured in an informal way. For the traveling salesman problem, for example, the number of cities is commonly used for this purpose. However, an m -city problem instance includes, in addition to the labels of the m cities, a collection of $m(m-1)/2$ numbers defining the inter-city distances, and the sizes of these numbers also contribute to the amount of input data. If we are to deal with time requirements in a precise, mathematical manner, we must take care to define instance size in such a way that all these factors are taken into account.

To do this, observe that the description of a problem instance that we provide as input to the computer can be viewed as a single finite string of symbols chosen from a finite input alphabet. Although there are many different ways in which instances of a given problem might be described, let us assume that one particular way has been chosen in advance and that each problem has associated with it a fixed *encoding scheme*, which maps problem

instances into the strings describing them. The *input length* for an instance I of a problem Π is defined to be the number of symbols in the description of I obtained from the encoding scheme for Π . It is this number, the input length, that is used as the formal measure of instance size.

For example, instances of the traveling salesman problem might be described using the alphabet $\{c, [,], /, 0, 1, 2, 3, 4, 5, 6, 7, 8, 9\}$, with our previous example of a problem instance being encoded by the string " $c[1]c[2]c[3]c[4]//10/5/9//6/9//3$." More complicated instances would be encoded in analogous fashion. If this were the encoding scheme associated with the traveling salesman problem, then the input length for our example would be 32.

The *time complexity function* for an algorithm expresses its time requirements by giving, for each possible input length, the largest amount of time needed by the algorithm to solve a problem instance of that size. Of course, this function is not well-defined until one fixes the encoding scheme to be used for determining input length and the computer or computer model to be used for determining execution time. However, as we shall see, the particular choices made for these will have little effect on the broad distinctions made in the theory of NP-completeness. Hence, in what follows, the reader is advised merely to fix in mind a particular encoding scheme for each problem and a particular computer or computer model, and to think in terms of time complexity as determined from the corresponding input lengths and execution times.

1.3 Polynomial Time Algorithms and Intractable Problems

Different algorithms possess a wide variety of different time complexity functions, and the characterization of which of these are "efficient enough" and which are "too inefficient" will always depend on the situation at hand. However, computer scientists recognize a simple distinction that offers considerable insight into these matters. This is the distinction between polynomial time algorithms and exponential time algorithms.

Let us say that a function $f(n)$ is $O(g(n))$ whenever there exists a constant c such that $|f(n)| \leq c \cdot |g(n)|$ for all values of $n \geq 0$. A *polynomial time algorithm* is defined to be one whose time complexity function is $O(p(n))$ for some polynomial function p , where n is used to denote the input length. Any algorithm whose time complexity function cannot be so bounded is called an *exponential time algorithm* (although it should be noted that this definition includes certain non-polynomial time complexity functions, like $n^{\log n}$, which are not normally regarded as exponential functions).

The distinction between these two types of algorithms has particular significance when considering the solution of large problem instances. Figure 1.2 illustrates the differences in growth rates among several typical complexity functions of each type, where the functions express execution time

in terms of microseconds. Notice the much more explosive growth rates for the two exponential complexity functions.

Time complexity function	Size n					
	10	20	30	40	50	60
n	.00001 second	.00002 second	.00003 second	.00004 second	.00005 second	.00006 second
n^2	.0001 second	.0004 second	.0009 second	.0016 second	.0025 second	.0036 second
n^3	.001 second	.008 second	.027 second	.064 second	.125 second	.216 second
n^5	.1 second	3.2 seconds	24.3 seconds	1.7 minutes	5.2 minutes	13.0 minutes
2^n	.001 second	1.0 second	17.9 minutes	12.7 days	35.7 years	366 centuries
3^n	.059 second	58 minutes	6.5 years	3855 centuries	2×10^8 centuries	1.3×10^{13} centuries

Figure 1.2 Comparison of several polynomial and exponential time complexity functions.

Even more revealing is an examination of the effects of improved computer technology on algorithms having these time complexity functions. Figure 1.3 shows how the largest problem instance solvable in one hour would change if we had a computer 100 or 1000 times faster than our present machine. Observe that with the 2^n algorithm a thousand-fold increase in computing speed only adds 10 to the size of the largest problem instance we can solve in an hour, whereas with the n^5 algorithm this size almost quadruples.

These tables indicate some of the reasons why polynomial time algorithms are generally regarded as being much more desirable than exponential time algorithms. This view, and the distinction between the two types of algorithms, is central to our notion of inherent intractability and to the theory of NP-completeness.

The fundamental nature of this distinction was first discussed in [Cobham, 1964] and [Edmonds, 1965a]. Edmonds, in particular, equated poly-

Size of Largest Problem Instance
Solvable in 1 Hour

Time complexity function	With present computer	With computer 100 times faster	With computer 1000 times faster
n	N_1	$100 N_1$	$1000 N_1$
n^2	N_2	$10 N_2$	$31.6 N_2$
n^3	N_3	$4.64 N_3$	$10 N_3$
n^5	N_4	$2.5 N_4$	$3.98 N_4$
2^n	N_5	$N_5 + 6.64$	$N_5 + 9.97$
3^n	N_6	$N_6 + 4.19$	$N_6 + 6.29$

Figure 1.3 Effect of improved technology on several polynomial and exponential time algorithms.

nomial time algorithms with “good” algorithms and conjectured that certain integer programming problems might not be solvable by such “good” algorithms. This reflects the viewpoint that exponential time algorithms should not be considered “good” algorithms, and indeed this usually is the case. Most exponential time algorithms are merely variations on exhaustive search, whereas polynomial time algorithms generally are made possible only through the gain of some deeper insight into the structure of a problem. There is wide agreement that a problem has not been “well-solved” until a polynomial time algorithm is known for it. Hence, we shall refer to a problem as *intractable* if it is so hard that no polynomial time algorithm can possibly solve it.

Of course, this formal use of “intractable” should be viewed only as a rough approximation to its dictionary meaning. The distinction between “efficient” polynomial time algorithms and “inefficient” exponential time algorithms admits of many exceptions when the problem instances of interest have limited size. Even in Figure 1.2, the 2^n algorithm is faster than the n^5 algorithm for $n \leq 20$. More extreme examples can be constructed easily.

Furthermore, there are some exponential time algorithms that have been quite useful in practice. Time complexity as defined is a *worst-case* measure, and the fact that an algorithm has time complexity 2^n means only that at least one problem instance of size n requires that much time. Most problem instances might actually require far less time than that, a situation

that appears to hold for several well-known algorithms. The simplex algorithm for linear programming has been shown to have exponential time complexity [Klee and Minty, 1972], [Zadeh, 1973], but it has an impressive record of running quickly in practice. Likewise, branch-and-bound algorithms for the knapsack problem have been so successful that many consider it to be a “well-solved” problem, even though these algorithms, too, have exponential time complexity.

Unfortunately, examples like these are quite rare. Although exponential time algorithms are known for many problems, few of them are regarded as being very useful in practice. Even the successful exponential time algorithms mentioned above have not stopped researchers from continuing to search for polynomial time algorithms for solving those problems. In fact, the very success of these algorithms has led to the suspicion that they somehow capture a crucial property of the problems whose refinement could lead to still better methods. So far, little progress has been made toward explaining this success, and no methods are known for predicting in advance that a given exponential time algorithm will run quickly in practice.

On the other hand, the much more stringent bounds on execution time satisfied by polynomial time algorithms often permit such predictions to be made. Even though an algorithm having time complexity n^{100} or $10^{99}n^2$ might not be considered likely to run quickly in practice, the polynomially solvable problems that arise naturally tend to be solvable within polynomial time bounds that have degree 2 or 3 at worst and that do not involve extremely large coefficients. Algorithms satisfying such bounds *can* be considered to be “provably efficient,” and it is this much-desired property that makes polynomial time algorithms the preferred way to solve problems.

Our definition of “intractable” also provides a theoretical framework of considerable generality and power. The intractability of a problem turns out to be essentially independent of the particular encoding scheme and computer model used for determining time complexity.

Let us first consider encoding schemes. Suppose for example that we are dealing with a problem in which each instance is a graph $G = (V, E)$, where V is the set of vertices and E is the set of edges, each edge being an unordered pair of vertices. Such an instance might be described (see Figure 1.4) by simply listing all the vertices and edges, or by listing the rows of the adjacency matrix for the graph, or by listing for each vertex all the other vertices sharing a common edge with it (a “neighbor” list). Each of these encodings can give a different input length for the same graph. However, it is easy to verify (see Figure 1.5) that the input lengths they determine differ at most polynomially from one another, so that any algorithm having polynomial time complexity under one of these encoding schemes also will have polynomial time complexity under all the others. In fact, the standard encoding schemes used in practice for any particular problem always seem to differ at most polynomially from one another. It would be difficult to imagine a “reasonable” encoding scheme for a problem that differs more

than polynomially from the standard ones. Although what we mean here by “reasonable” cannot be formalized, the following two conditions capture much of the notion:

- (1) the encoding of an instance I should be concise and not “padded” with unnecessary information or symbols, and
- (2) numbers occurring in I should be represented in binary (or decimal, or octal, or in any fixed base other than 1).

If we restrict ourselves to encoding schemes satisfying these conditions, then the particular encoding scheme used should not affect the determination of whether a given problem is intractable.

Encoding Scheme	String	Length
Vertex list, Edge list	V[1]V[2]V[3]V[4](V[1]V[2])(V[2]V[3])	36
Neighbor list	(V[2])(V[1]V[3])(V[2])()	24
Adjacency matrix rows	0100/1010/0010/0000	19

Figure 1.4 Descriptions of the graph $G = (V, E)$ where $V = \{V_1, V_2, V_3, V_4\}$ and $E = \{\{V_1, V_2\}, \{V_2, V_3\}\}$, under three different encoding schemes.

Encoding Scheme	Lower Bound	Upper Bound
Vertex list, Edge list	$4v + 10e$	$4v + 10e + (v + 2e) \cdot \lceil \log_{10} v \rceil$
Neighbor list	$2v + 8e$	$2v + 8e + 2e \cdot \lceil \log_{10} v \rceil$
Adjacency matrix	$v^2 + v - 1$	$v^2 + v - 1$

Figure 1.5 General bounds on input lengths for the three encoding schemes of Figure 1.4 for graphs $G = (V, E)$ with $|V| = v$, $|E| = e$. Since $e < v^2$, these show that the input lengths differ at most polynomially from each other. ($\lceil x \rceil$ denotes the least integer not less than x .)

Similar comments can be made concerning the choice of computer models. All the realistic models of computers studied so far, such as one-tape Turing machines, multi-tape Turing machines, and random-access machines (RAMs), are equivalent with respect to polynomial time complexity (for example, see Figure 1.6). One would expect any other “reasonable” model to share in this equivalence. The notion of “reasonable” in-

tended here is essentially that there is a polynomial bound on the amount of work that can be done in a single unit of time. Thus, for example, a model having the capability of performing arbitrarily many operations in parallel would not be considered “reasonable,” and indeed no existing (or planned) computer has this capability. At any rate, so long as we restrict ourselves to the standard models of realistic computers, the class of intractable problems will be unaffected by the particular model used, and we can make our choice on the basis of convenience without sacrificing the applicability of our results.

Simulated machine B	Simulating machine A		
	1TM	kTM	RAM
1-Tape Turing Machine (1TM)	—	$O(T(n))$	$O(T(n)\log T(n))$
k-Tape Turing Machine (kTM)	$O(T^2(n))$	—	$O(T(n)\log T(n))$
Random Access Machine (RAM)	$O(T^3(n))$	$O(T^2(n))$	—

Figure 1.6 Time required by machine A to simulate the execution of an algorithm of time complexity $T(n)$ on Machine B (for example, see [Hopcroft and Ullman, 1969] and [Aho, Hopcroft, and Ullman, 1974]).

1.4 Provably Intractable Problems

Now that we have discussed the formal meaning of “intractable problem,” it is appropriate that we briefly survey the current state of knowledge about the existence of intractable problems.

It is useful to begin by distinguishing between two different causes of intractability allowed by our definition. The first, which is the one we usually have in mind, is that the problem is so difficult that an exponential amount of time is needed to discover a solution. The second is that the solution *itself* is required to be so extensive that it cannot be described with an expression having length bounded by a polynomial function of the input length.

This second cause occurs, for example, in the variant of the traveling salesman problem that includes a number B as an additional parameter and that asks for *all* tours having total length B or less. It is easy to construct instances of this problem in which exponentially many tours are shorter than the given bound, so that no polynomial time algorithm could possibly list them all.

Intractability of this sort is by no means insignificant, and it is important to recognize it when it occurs. However, in most cases its existence is

apparent from the problem definition. In fact, this type of intractability can be regarded as a signal that the problem is not defined realistically, because we are asking for more information than we could ever hope to use. Thus, from now on we shall restrict our attention to the first type of intractability. Accordingly, only problems for which the solution length is bounded by a polynomial function of the input length will be considered.

The earliest intractability results for such problems are the classical undecidability results of Alan Turing. Over 40 years ago, Turing demonstrated that certain problems are so hard that they are “undecidable,” in the sense that no algorithm at all can be given for solving them. He proved, for example, that it is impossible to specify *any* algorithm which, given an arbitrary computer program and an arbitrary input to that program, can decide whether or not the program will eventually halt when applied to that input [Turing, 1936]. A variety of other problems are now known to be undecidable, including the triviality problem for finitely presented groups [Rabin, 1958], Hilbert’s tenth problem (solvability of polynomial equations in integers) [Matijasevic, 1970], and several problems of “tiling the plane” [Berger, 1966]. Since these undecidable problems cannot be solved by *any* algorithm, much less a polynomial time algorithm, they indeed are intractable in an especially strong sense.

The first examples of intractable “decidable” problems were obtained in the early 1960’s, as part of work on complexity “hierarchies” by Hartmanis and Stearns [1965]. However, these results involved only “artificial” problems, specifically constructed to have the appropriate properties. It was not until the early 1970’s that Meyer and Stockmeyer [1972], Fischer and Rabin [1974], and others finally succeeded in proving some “natural” decidable problems to be intractable. These include a variety of previously studied problems from automata theory, formal language theory, and mathematical logic. In fact, the proofs show that these problems cannot be solved in polynomial time using even a “nondeterministic” computer model, which has the ability to pursue an unbounded number of independent computational sequences in parallel. We shall see that this “unreasonable” computer model plays an important role in the theory of NP-completeness, and its capabilities will be specified more fully in Chapter 2.

All the provably intractable problems known to date fall into the two categories we have just mentioned. They are either undecidable or “nondeterministically” intractable. However, most of the apparently intractable problems encountered in practice *are* decidable and *can* be solved in polynomial time with the aid of a nondeterministic computer. Thus, none of the proof techniques developed so far is powerful enough to verify the apparent intractability of these problems.

1.5 NP-Complete Problems

As theoreticians continue to seek more powerful methods for proving problems intractable, parallel efforts focus on learning more about the ways in which various problems are interrelated with respect to their difficulty. As we suggested earlier, the discovery of such relationships between problems often can provide information useful to algorithm designers.

The principal technique used for demonstrating that two problems are related is that of “reducing” one to the other, by giving a constructive transformation that maps any instance of the first problem into an equivalent instance of the second. Such a transformation provides the means for converting any algorithm that solves the second problem into a corresponding algorithm for solving the first problem.

Many simple examples of such reductions have been known for some time. For example, Dantzig [1960] reduced a number of combinatorial optimization problems to the general zero-one integer linear programming problem. Edmonds [1962] reduced the graph theoretic problems of “covering all edges with a minimum number of vertices” and “finding a maximum independent set of vertices” to the general “set covering problem.” Gimpel [1965] reduced the general set covering problem to the “prime implicant covering problem” of logic design. Dantzig, Blattner, and Rao [1966] described a “well-known” reduction from the traveling salesman problem to the “shortest path problem” with negative edge lengths allowed.

These early reductions, although rather isolated and limited in scope, foreshadow the kind of results proved in the theory of NP-completeness.

The foundations for the theory of NP-completeness were laid in a paper of Stephen Cook, presented in 1971, entitled “The Complexity of Theorem Proving Procedures” [Cook, 1971a]. In this brief but elegant paper Cook did several important things.

First, he emphasized the significance of “polynomial time reducibility,” that is, reductions for which the required transformation can be executed by a polynomial time algorithm. If we have a polynomial time reduction from one problem to another, this ensures that any polynomial time algorithm for the second problem can be converted into a corresponding polynomial time algorithm for the first problem.

Second, he focused attention on the class NP of decision problems that can be solved in polynomial time by a nondeterministic computer. (A decision problem is one whose solution is either “yes” or “no”.) Most of the apparently intractable problems encountered in practice, when phrased as decision problems, belong to this class.

Third, he proved that one particular problem in NP, called the “satisfiability” problem, has the property that every other problem in NP can be polynomially reduced to it. If the satisfiability problem can be solved with a polynomial time algorithm, then so can every problem in NP, and if any problem in NP is intractable, then the satisfiability problem also must

be intractable. Thus, in a sense, the satisfiability problem is the “hardest” problem in NP.

Finally, Cook suggested that other problems in NP might share with the satisfiability problem this property of being the “hardest” member of NP. He showed this to be the case for the problem “Does a given graph G contain a complete subgraph on a given number k of vertices?”

Subsequently, Richard Karp presented a collection of results [Karp, 1972] proving that indeed the decision problem versions of many well known combinatorial problems, including the traveling salesman problem, are just as “hard” as the satisfiability problem. Since then a wide variety of other problems have been proved equivalent in difficulty to these problems, and this equivalence class, consisting of the “hardest” problems in NP, has been given a name: the class of *NP-complete problems*.

Cook’s original ideas have turned out to be remarkably powerful. They have provided the means for combining many individual complexity questions into the single question: Are the NP-complete problems intractable? The lists included in the Appendix of this book contain literally hundreds of different problems now known to be NP-complete. As more and more problems of independent interest are shown to belong to this equivalence class, its importance is continually reinforced.

The question of whether or not the NP-complete problems are intractable is now considered to be one of the foremost open questions of contemporary mathematics and computer science. Despite the willingness of most researchers to conjecture that the NP-complete problems are all intractable, little progress has yet been made toward establishing either a proof or a disproof of this far-reaching conjecture. However, even without a proof that NP-completeness implies intractability, the knowledge that a problem is NP-complete suggests, at the very least, that a major breakthrough will be needed to solve it with a polynomial time algorithm.

1.6 An Outline of the Book

Although this book is intended mainly as a primer on how to determine whether or not any particular problem is NP-complete (either by looking it up in the lists we present or by proving it yourself), we shall also discuss some of the options available for dealing with a problem that is known to be NP-complete. A brief outline of subsequent chapters follows.

In Chapter 2, we present the formal underpinnings of NP-completeness and prove Cook’s theorem. The central definitions involve certain theoretical concepts, such as “languages” and “Turing machines,” which we develop in a straightforward manner, relating them to the notions of problems and computer models already discussed. This chapter should give the reader a good understanding of the technical meaning of NP-completeness.

Chapter 3 is devoted to methods for proving a problem NP-complete. A number of examples are presented to illustrate the usual structure of such proofs, and to indicate how one goes about generating one. In essence, one proves a new problem to be NP-complete by polynomially reducing a known NP-complete problem to it. We survey the known NP-complete problems that have been most useful for this purpose and demonstrate their use.

In Chapter 4, we examine the ways in which the theory of NP-completeness can be used for conducting a detailed analysis of the complexity of a problem, seeking to determine the “boundary” between those cases of the problem that are polynomially solvable and those that are NP-complete.

In Chapter 5, we show how the techniques used for proving NP-completeness can be generalized so that problems other than just decision problems can be proved to be “as hard as” the NP-complete problems. As an aid to reading the published literature on the theory of NP-completeness, we also provide a brief historical survey of the development of the main ideas and the varying terminology that has been used for discussing them.

In Chapter 6, we discuss several approaches for dealing with intractable problems, especially that of finding near-optimal solutions using fast algorithms. Examples of the successes and failures of each approach are described, and we illustrate how the theory of NP-completeness can be applied even here.

Chapter 7 is intended to acquaint the reader with some of the theoretical issues and ideas that have arisen in parallel with the theory of NP-completeness. Among other topics we discuss the polynomial hierarchy, #P-completeness, polynomial space completeness, and the “relativization” of the question of the intractability of the NP-complete problems.

The last third of the book consists of the Appendix, an extensive and annotated list of problems known to be NP-complete or harder. The list is divided into sections, each devoted to problems from a particular subject area, such as graph theory, scheduling, algebra and number theory, covering and partitioning, mathematical programming, program optimization, automata and language theory, and, of course, miscellaneous topics. The list includes references to related problems known to be solvable in polynomial time and to problems whose status remains open in that neither polynomial time algorithms nor NP-completeness proofs are known for them.