Undergraduate Texts in Mathematics

S. Axler F.W. Gehring K.A. Ribet

Kai Lai Chung Farid AitSahlia

Elementary Probability Theory

With Stochastic Processes and an Introduction to Mathematical Finance

Fourth Edition

With 57 Figures



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Preface to the Fourth Edition

In this edition two new chapters, 9 and 10, on mathematical finance are added. They are written by Dr. Farid AitSahlia, *ancien élève*, who has taught such a course and worked on the research staff of several industrial and financial institutions.

The new text begins with a meticulous account of the uncommon vocabulary and syntax of the financial world; its manifold options and actions, with consequent expectations and variations, in the marketplace. These are then expounded in clear, precise mathematical terms and treated by the methods of probability developed in the earlier chapters. Numerous graded and motivated examples and exercises are supplied to illustrate the applicability of the fundamental concepts and techniques to concrete financial problems. For the reader whose main interest is in finance, only a portion of the first eight chapters is a "prerequisite" for the study of the last two chapters. Further specific references may be scanned from the topics listed in the Index, then pursued in more detail.

I have taken this opportunity to fill a gap in Section 8.1 and to expand Appendix 3 to include a useful proposition on martingale stopped at an optional time. The latter notion plays a basic role in more advanced financial and other disciplines. However, the level of our compendium remains elementary, as befitting the title and scheme of this textbook. We have also included some up-to-date financial episodes to enliven, for the beginners, the stratified atmosphere of "strictly business". We are indebted to Ruth Williams, who read a draft of the new chapters with valuable suggestions for improvement; to Bernard Bru and Marc Barbut for information on the Pareto-Lévy laws originally designed for income distributions. It is hoped that a readable summary of this renowned work may be found in the new Appendix 4.

Kai Lai Chung August 3, 2002

Prologue to Introduction to Mathematical Finance

The two new chapters are self-contained introductions to the topics of mean-variance optimization and option pricing theory. The former covers a subject that is sometimes labeled "modern portfolio theory" and that is widely used by money managers employed by large financial institutions. To read this chapter, one only needs an elementary knowledge of probability concepts and a modest familiarity with calculus. Also included is an introductory discussion on stable laws in an applied context, an often neglected topic in elementary probability and finance texts. The latter chapter lays the foundations for option pricing theory, a subject that has fueled the development of finance into an advanced mathematical discipline as attested by the many recently published books on the subject. It is an initiation to martingale pricing theory, the mathematical expression of the so-called "arbitrage pricing theory", in the context of the binomial random walk. Despite its simplicity, this model captures the flavors of many advanced theoretical issues. It is often used in practice as a benchmark for the approximate pricing of complex financial instruments.

I would like to thank Professor Kai Lai Chung for inviting me to write the new material for the fourth edition. I would also like to thank my wife Unnur for her support during this rewarding experience.

> Farid AitSahlia November 1, 2002

1

Set

1.1. Sample sets

These days schoolchildren are taught about sets. A second grader* was asked to name "the set of girls in his class." This can be done by a complete list such as:

"Nancy, Florence, Sally, Judy, Ann, Barbara, ..."

A problem arises when there are duplicates. To distinguish between two Barbaras one must indicate their family names or call them B_1 and B_2 . The same member cannot be counted twice in a set.

The notion of a set is common in all mathematics. For instance, in geometry one talks about "the set of points which are equidistant from a given point." This is called a circle. In algebra one talks about "the set of integers which have no other divisors except 1 and itself." This is called the set of prime numbers. In calculus the domain of definition of a function is a set of numbers, e.g., the interval (a,b); so is the range of a function if you remember what it means.

In probability theory the notion of a set plays a more fundamental role. Furthermore we are interested in very general kinds of sets as well as specific concrete ones. To begin with the latter kind, consider the following examples:

- (a) a bushel of apples;
- (b) fifty-five cancer patients under a certain medical treatment;

^{*}My son Daniel.

- (c) all the students in a college;
- (d) all the oxygen molecules in a given container;
- (e) all possible outcomes when six dice are rolled;
- (f) all points on a target board.

Let us consider at the same time the following "smaller" sets:

- (a') the rotten apples in that bushel;
- (b') those patients who respond positively to the treatment;
- (c') the mathematics majors of that college;
- (d') those molecules that are traveling upwards;
- (e') those cases when the six dice show different faces;
- (f') the points in a little area called the "bull's-eye" on the board.

We shall set up a mathematical model for these and many more such examples that may come to mind, namely we shall abstract and generalize our intuitive notion of "a bunch of things." First we call the things points, then we call the bunch a space; we prefix them by the word "sample" to distinguish these terms from other usages, and also to allude to their statistical origin. Thus a sample point is the abstraction of an apple, a cancer patient, a student, a molecule, a possible chance outcome, or an ordinary geometrical point. The sample space consists of a number of sample points and is just a name for the totality or aggregate of them all. Any one of the examples (a)–(f) above can be taken to be a sample space, but so also may any one of the smaller sets in (a')–(f'). What we choose to call a space [a universe] is a relative matter.

Let us then fix a sample space to be denoted by Ω , the capital Greek letter *omega*. It may contain any number of points, possibly infinite but at least one. (As you have probably found out before, mathematics can be very pedantic!) Any of these points may be denoted by ω , the small Greek letter omega, to be distinguished from one another by various devices such as adding subscripts or dashes (as in the case of the two Barbaras if we do not know their family names), thus $\omega_1, \omega_2, \omega', \ldots$. Any partial collection of the points is a *subset* of Ω , and since we have fixed Ω we will just call it a set. In extreme cases a set may be Ω itself or the *empty set*, which has no point in it. You may be surprised to hear that the empty set is an important entity and is given a special symbol \emptyset . The number of points in a set S will be called its *size* and denoted by |S|; thus it is a nonnegative integer or ∞ . In particular $|\emptyset| = 0$.

A particular set S is well defined if it is possible to tell whether any given point *belongs to* it or not. These two cases are denoted respectively by

Thus a set is determined by a specified rule of membership. For instance, the sets in (a')–(f') are well defined up to the limitations of verbal descriptions. One can always quibble about the meaning of words such as "a rotten apple," or attempt to be funny by observing, for instance, that when dice are rolled on a pavement some of them may disappear into the sewer. Some people of a pseudo-philosophical turn of mind get a lot of mileage out of such *caveats*, but we will not indulge in them here. Now, one sure way of specifying a rule to determine a set is to enumerate all its members, namely to make a complete list as the second grader did. But this may be tedious if not impossible. For example, it will be shown in §3.1 that the size of the set in (e) is equal to $6^6 = 46656$. Can you give a quick guess as to how many pages of a book like this will be needed just to record all these possibilities of a mere throw of six dice? On the other hand, it can be described in a systematic and unmistakable way as the set of all ordered 6-tuples of the form below:

$$(s_1, s_2, s_3, s_4, s_5, s_6)$$

where each of the symbols s_j , $1 \le j \le 6$, may be any of the numbers 1, 2, 3, 4, 5, 6. This is a good illustration of mathematics being economy of thought (and printing space).

If every point of A belongs to B, then A is contained or included in B and is a subset of B, while B is a superset of A. We write this in one of the two ways below:

$$A \subset B$$
, $B \supset A$.

Two sets are *identical* if they contain exactly the same points, and then we write

$$A = B$$
.

Another way to say this is: A = B if and only if $A \subset B$ and $B \subset A$. This may sound unnecessarily roundabout to you, but is often the only way to check that two given sets are really identical. It is not always easy to identify two sets defined in different ways. Do you know for example that the set of even integers is identical with the set of all solutions x of the equation $\sin(\pi x/2) = 0$? We shall soon give some examples of showing the identity of sets by the roundabout method.

1.2. Operations with sets

We learn about sets by operating on them, just as we learn about numbers by operating on them. In the latter case we also say that we compute

with numbers: add, subtract, multiply, and so on. These operations performed on given numbers produce other numbers, which are called their sum, difference, product, etc. In the same way, operations performed on sets produce other sets with new names. We are now going to discuss some of these and the laws governing them.

Complement. The complement of a set A is denoted by A^c and is the set of points that do not belong to A. Remember we are talking only about points in a fixed Ω ! We write this symbolically as follows:

$$A^c = \{ \omega \mid \omega \notin A \},\$$

which reads: " A^c is the set of ω that does not belong to A." In particular $\Omega^c = \emptyset$ and $\emptyset^c = \Omega$. The operation has the property that if it is performed twice in succession on A, we get A back:

$$(A^c)^c = A. (1.2.1)$$

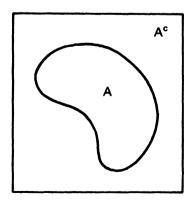
Union. The union $A \cup B$ of two sets A and B is the set of points that belong to at least one of them. In symbols:

$$A \cup B = \{ \omega \mid \omega \in A \text{ or } \omega \in B \}$$

where "or" means "and/or" in pedantic [legal] style and will always be used in this sense.

Intersection. The intersection $A \cap B$ of two sets A and B is the set of points that belong to both of them. In symbols:

$$A \cap B = \{ \omega \mid \omega \in A \text{ and } \omega \in B \}.$$



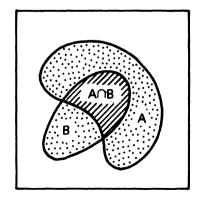


Figure 1

We hold the truth of the following laws as self-evident:

Commutative Law. $A \cup B = B \cup A$, $A \cap B = B \cap A$.

Associative Law.
$$(A \cup B) \cup C = A \cup (B \cup C),$$
 $(A \cap B) \cap C = A \cap (B \cap C).$

But observe that these relations are instances of identity of sets mentioned above, and are subject to proof. They should be compared, but not confused, with analogous laws for sum and product of numbers:

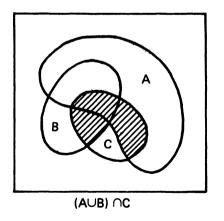
$$a+b=b+a, \quad a\times b=b\times a$$

$$(a+b)+c=a+(b+c), \quad (a\times b)\times c=a\times (b\times c).$$

Brackets are needed to indicate the order in which the operations are to be performed. Because of the associative laws, however, we can write

$$A \cup B \cup C$$
, $A \cap B \cap C \cap D$

without brackets. But a string of symbols like $A \cup B \cap C$ is ambiguous, therefore not defined; indeed $(A \cup B) \cap C$ is not identical with $A \cup (B \cap C)$. You should be able to settle this easily by a picture.



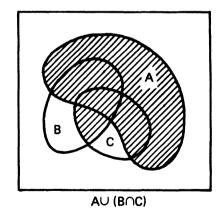


Figure 2

The next pair of distributive laws connects the two operations as follows:

$$(A \cup B) \cap C = (A \cap C) \cup (B \cap C); \tag{D_1}$$

$$(A \cap B) \cup C = (A \cup C) \cap (B \cup C). \tag{D_2}$$

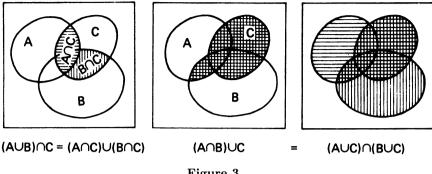


Figure 3

Several remarks are in order. First, the analogy with arithmetic carries over to (D_1) :

$$(a+b) \times c = (a \times c) + (b \times c);$$

but breaks down in (D_2) :

$$(a \times b) + c \neq (a + c) \times (b + c).$$

Of course, the alert reader will have observed that the analogy breaks down already at an earlier stage, for

$$A = A \cup A = A \cap A$$
:

but the only number a satisfying the relation a + a = a is 0; while there are exactly two numbers satisfying $a \times a = a$, namely 0 and 1.

Second, you have probably already discovered the use of diagrams to prove or disprove assertions about sets. It is also a good practice to see the truth of such formulas as (D_1) and (D_2) by well-chosen examples. Suppose then that

$$A = \text{inexpensive things}, B = \text{really good things},$$

$$C = \text{food [edible things]}.$$

Then $(A \cup B) \cap C$ means "(inexpensive or really good) food," while $(A \cap C) \cup (B \cap C)$ means "(inexpensive food) or (really good food)." So they are the same thing all right. This does not amount to a proof, as one swallow does not make a summer, but if one is convinced that whatever logical structure or thinking process involved above in no way depends on the precise nature of the three things A, B, and C, so much so that they can be *anything*, then one has in fact landed a general proof. Now it is interesting that the same example applied to (D_2) somehow does not make it equally obvious

(at least to the author). Why? Perhaps because some patterns of logic are in more common use in our everyday experience than others.

This last remark becomes more significant if one notices an obvious duality between the two distributive laws. Each can be obtained from the other by switching the two symbols \cup and \cap . Indeed each can be deduced from the other by making use of this duality (Exercise 11).

Finally, since (D_2) comes less naturally to the intuitive mind, we will avail ourselves of this opportunity to demonstrate the roundabout method of identifying sets mentioned above by giving a rigorous proof of the formula. According to this method, we must show: (i) each point on the left side of (D_2) belongs to the right side; (ii) each point on the right side of (D_2) belongs to the left side.

- (i) Suppose ω belongs to the left side of (D_2) , then it belongs either to $A \cap B$ or to C. If $\omega \in A \cap B$, then $\omega \in A$, hence $\omega \in A \cup C$; similarly $\omega \in B \cup C$. Therefore ω belongs to the right side of (D_2) . On the other hand, if $\omega \in C$, then $\omega \in A \cup C$ and $\omega \in B \cup C$ and we finish as before.
- (ii) Suppose ω belongs to the right side of (D_2) , then ω may or may not belong to C, and the trick is to consider these two alternatives. If $\omega \in C$, then it certainly belongs to the left side of (D_2) . On the other hand, if $\omega \notin C$, then since it belongs to $A \cup C$, it must belong to A; similarly it must belong to B. Hence it belongs to $A \cap B$, and so to the left side of (D_2) . Q.E.D.

1.3. Various relations

The three operations so far defined: complement, union, and intersection obey two more laws called *De Morgan's laws*:

$$(A \cup B)^c = A^c \cap B^c; \tag{C_1}$$

$$(A \cap B)^c = A^c \cup B^c. \tag{C_2}$$

They are dual in the same sense as (D_1) and (D_2) are. Let us check these by our previous example. If A = inexpensive, and B = really good, then clearly $(A \cup B)^c =$ not inexpensive nor really good, namely high-priced junk, which is the same as $A^c \cap B^c =$ inexpensive and not really good. Similarly we can check (C_2) .

Logically, we can deduce either (C_1) or (C_2) from the other; let us show it one way. Suppose then (C_1) is true, then since A and B are arbitrary sets we can substitute their complements and get

$$(A^c \cup B^c)^c = (A^c)^c \cap (B^c)^c = A \cap B \tag{1.3.1}$$

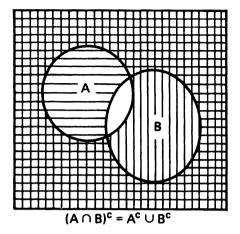


Figure 4

where we have also used (1.2.1) for the second equation. Now taking the complements of the first and third sets in (1.3.1) and using (1.2.1) again we get

$$A^c \cup B^c = (A \cap B)^c.$$

This is (C_2) . Q.E.D.

It follows from De Morgan's laws that if we have complementation, then either union or intersection can be expressed in terms of the other. Thus we have

$$A \cap B = (A^c \cup B^c)^c,$$

$$A \cup B = (A^c \cap B^c)^c;$$

and so there is redundancy among the three operations. On the other hand, it is impossible to express complementation by means of the other two although there is a magic symbol from which all three can be derived (Exercise 14). It is convenient to define some other operations, as we now do.

Difference. The set $A \setminus B$ is the set of points that belong to A and (but) not to B. In symbols:

$$A \setminus B = A \cap B^c = \{\omega \mid \omega \in A \text{ and } \omega \not \in B\}.$$

This operation is neither commutative nor associative. Let us find a *counterexample* to the associative law, namely, to find some A, B, C for which

$$(A \setminus B) \setminus C \neq A \setminus (B \setminus C). \tag{1.3.2}$$

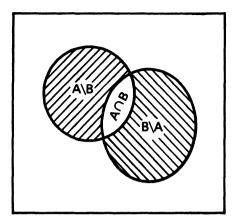


Figure 5

Note that in contrast to a proof of identity discussed above, a single instance of falsehood will destroy the identity. In looking for a counterexample one usually begins by specializing the situation to reduce the "unknowns." So try B = C. The left side of (1.3.2) becomes $A \setminus B$, while the right side becomes $A \setminus \emptyset = A$. Thus we need only make $A \setminus B \neq A$, and that is easy.

In case $A \supset B$ we write A - B for $A \setminus B$. Using this new symbol we have

$$A \setminus B = A - (A \cap B)$$

and

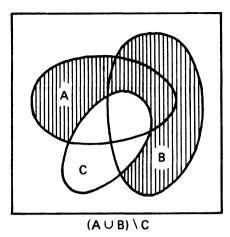
$$A^c = \Omega - A$$
.

The operation "-" has some resemblance to the arithmetic operation of subtracting, in particular $A - A = \emptyset$, but the analogy does not go very far. For instance, there is no analogue to (a + b) - c = a + (b - c).

Symmetric Difference. The set $A \triangle B$ is the set of points that belong to exactly one of the two sets A and B. In symbols:

$$A \triangle B = (A \cap B^c) \cup (A^c \cap B) = (A \setminus B) \cup (B \setminus A).$$

This operation is useful in advanced theory of sets. As its name indicates, it is symmetric with respect to A and B, which is the same as saying that it is commutative. Is it associative? Try some concrete examples or diagrams, which have succeeded so well before, and you will probably be as quickly confused as I am. But the question can be neatly resolved by a device to be introduced in $\S1.4$.



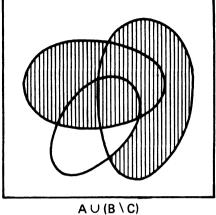


Figure 6

Having defined these operations, we should let our fancy run free for a few moments and imagine all kinds of sets that can be obtained by using them in succession in various combinations and permutations, such as

$$[(A \setminus C^c) \cap (B \cup C)^c]^c \cup (A^c \triangle B).$$

But remember we are talking about subsets of a fixed Ω , and if Ω is a finite set of a number of distinct subsets is certainly also finite, so there must be a tremendous amount of interrelationship among these sets that we can build up. The various laws discussed above are just some of the most basic ones, and a few more will be given among the exercises below.

An extremely important relation between sets will now be defined. Two sets A and B are said to be disjoint when they do not intersect, namely, have no point in common:

$$A \cap B = \emptyset$$
.

This is equivalent to either one of the following inclusion conditions:

$$A \subset B^c$$
: $B \subset A^c$.

Any number of sets are said to be disjoint when every pair of them is disjoint as just defined. Thus, "A, B, C are disjoint" means more than just $A \cap B \cap C = \emptyset$; it means

$$A \cap B = \emptyset$$
, $A \cap C = \emptyset$, $B \cap C = \emptyset$.

From here on we will omit the intersection symbol and write simply

$$AB$$
 for $A \cap B$

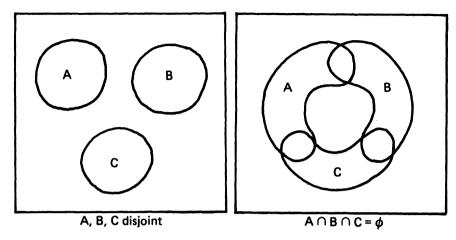


Figure 7

just as we write ab for $a \times b$. When A and B are disjoint we will write sometimes

$$A+B$$
 for $A\cup B$.

But be careful: not only does "+" mean addition for numbers but even when A and B are sets there are other usages of A+B such as their vectorial sum.

For any set A, we have the obvious decomposition:

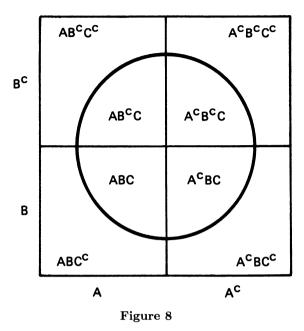
$$\Omega = A + A^c. \tag{1.3.3}$$

The way to think of this is: the set A gives a classification of all points ω in Ω according as ω belongs to A or to A^c . A college student may be classified according to whether he is a mathematics major or not, but he can also be classified according to whether he is a freshman or not, of voting age or not, has a car or not, ..., is a girl or not. Each two-way classification divides the sample space into two disjoint sets, and if several of these are superimposed on each other we get, e.g.,

$$\Omega = (A + A^{c})(B + B^{c}) = AB + AB^{c} + A^{c}B + A^{c}B^{c}, \qquad (1.3.4)$$

$$\Omega = (A + A^{c})(B + B^{c})(C + C^{c})
= ABC + ABC^{c} + AB^{c}C + AB^{c}C^{c} + A^{c}BC
+ A^{c}BC^{c} + A^{c}B^{c}C + A^{c}B^{c}C^{c}.$$
(1.3.5)

Let us call the pieces of such a decomposition the *atoms*. There are 2, 4, 8 atoms respectively above because 1, 2, 3 sets are considered. In general there



will be 2^n atoms if n sets are considered. Now these atoms have a remarkable property, which will be illustrated in the case (1.3.5), as follows: no matter how you operate on the three sets A, B, C, and no matter how many times you do it, the resulting set can always be written as the union of some of the atoms. Here are some examples:

$$A \cup B = ABC + ABC^c + AB^cC + AB^cC^c + A^cBC^c + A^cBC$$

$$(A \setminus B) \setminus C^c = AB^cC$$
$$(A \triangle B)C^c = AB^cC^c + A^cBC^c.$$

Can you see why?

Up to now we have considered only the union or intersection of a finite number of sets. There is no difficulty in extending this to an infinite number of sets. Suppose a finite or infinite sequence of sets A_n , n = 1, 2, ..., is given, then we can form their union and intersection as follows:

$$\bigcup_{n} A_{n} = \{ \omega \mid \omega \in A_{n} \text{ for at least one value of } n \};$$

$$\bigcap_{n} A_{n} = \{ \omega \mid \omega \in A_{n} \text{ for all values of } n \}.$$

When the sequence is infinite these may be regarded as obvious "set limits"

of finite unions or intersections, thus:

$$\bigcup_{n=1}^{\infty} A_n = \lim_{m \to \infty} \bigcup_{n=1}^{m} A_n; \quad \bigcap_{n=1}^{\infty} A_n = \lim_{m \to \infty} \bigcap_{n=1}^{m} A_n.$$

Observe that as m increases, $\bigcup_{n=1}^{m} A_n$ does not decrease while $\bigcap_{n=1}^{m} A_n$ does not increase, and we may say that the former swells up to $\bigcup_{n=1}^{\infty} A_n$, the latter shrinks down to $\bigcap_{n=1}^{\infty} A_n$.

The distributive laws and De Morgan's laws have obvious extensions to a finite or infinite sequence of sets. For instance,

$$\left(\bigcup_{n} A_{n}\right) \cap B = \bigcup_{n} (A_{n} \cap B), \tag{1.3.6}$$

$$\left(\bigcap_{n} A_{n}\right)^{c} = \bigcup_{n} A_{n}^{c}.\tag{1.3.7}$$

Really interesting new sets are produced by using both union and intersection an infinite number of times, and in succession. Here are the two most prominent ones:

$$\bigcap_{m=1}^{\infty} \left(\bigcup_{n=m}^{\infty} A_n \right); \quad \bigcup_{m=1}^{\infty} \left(\bigcap_{n=m}^{\infty} A_n \right).$$

These belong to a more advanced course (see [Chung 1, §4.2] of the References). They are shown here as a preview to arouse your curiosity.

1.4. Indicator*

The idea of classifying ω by means of a dichotomy: to be or not to be in A, which we discussed toward the end of §1.3, can be quantified into a useful device. This device will generalize to the fundamental notion of "random variable" in Chapter 4.

Imagine Ω to be a target board and A a certain marked area on the board as in Examples (f) and (f') above. Imagine that "pick a point ω in Ω " is done by shooting a dart at the target. Suppose a bell rings (or a bulb lights up) when the dart hits within the area A; otherwise it is a dud. This is the intuitive picture expressed below by a mathematical formula:

$$I_A(\omega) = \begin{cases} 1 & \text{if } \omega \in A, \\ 0 & \text{if } \omega \notin A. \end{cases}$$

^{*}This section may be omitted after the first three paragraphs.

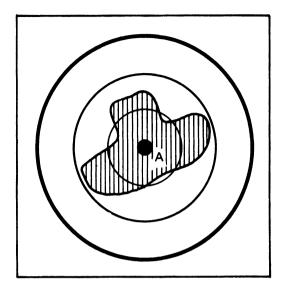


Figure 9

Thus the symbol I_A is a function that is defined on the whole sample space Ω and takes only the two values 0 and 1, corresponding to a dud and a ring. You may have learned in a calculus course the importance of distinguishing between a function (sometimes called a mapping) and one of its values. Here it is the function I_A that indicates the set A, hence it is called the indicator function, or briefly, *indicator* of A. Another set B has its indicator I_B . The two functions I_A and I_B are identical (what does that mean?) if and only if the two sets are identical.

To see how we can put indicators to work, let us figure out the indicators for some of the sets discussed before. We need two mathematical symbols \vee (cup) and \wedge (cap), which may be new to you. For any two real numbers a and b, they are defined as follows:

$$a \lor b = \text{maximum of } a \text{ and } b;$$

 $a \land b = \text{minimum of } a \text{ and } b.$ (1.4.1)

In case a = b, either one of them will serve as maximum as well as minimum. Now the salient properties of indicators are given by the formulas below:

$$I_{A\cap B}(\omega) = I_A(\omega) \wedge I_B(\omega) = I_A(\omega) \cdot I_B(\omega); \tag{1.4.2}$$

$$I_{A\cup B}(\omega) = I_A(\omega) \vee I_B(\omega).$$
 (1.4.3)

You should have no difficulty checking these equations, after all there are only two possible values 0 and 1 for each of these functions. Since the equations are true for every ω , they can be written more simply as equations

(identities) between functions:

$$I_{A \cap B} = I_A \wedge I_B = I_A \cdot I_B, \tag{1.4.4}$$

$$I_{A \cup B} = I_A \vee I_B. \tag{1.4.5}$$

Here for example the function $I_A \wedge I_B$ is that mapping that assigns to each ω the value $I_A(\omega) \wedge I_B(\omega)$, just as in calculus the function f + g is that mapping that assigns to each x the number f(x) + g(x).

After observing the product $I_A(\omega) \cdot I_B(\omega)$ at the end of (1.4.2) you may be wondering why we do not have the sum $I_A(\omega) + I_B(\omega)$ in (1.4.3). But if this were so we could get the value 2 here, which is impossible since the first member $I_{A\cup B}(\omega)$ cannot take this value. Nevertheless, shouldn't $I_A + I_B$ mean something? Consider target shooting again but this time mark out two overlapping areas A and B. Instead of bell-ringing, you get 1 penny if you hit within A, and also if you hit within B. What happens if you hit the intersection AB? That depends on the rule of the game. Perhaps you still get 1 penny, perhaps you get 2 pennies. Both rules are legitimate. In formula (1.4.3) it is the first rule that applies. If you want to apply the second rule, then you are no longer dealing with the set $A \cup B$ alone as in Figure 10a, but something like Figure 10b:

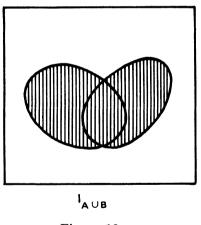


Figure 10a

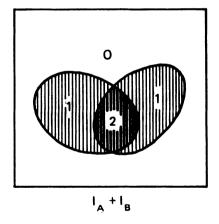


Figure 10b

This situation can be realized electrically by laying first a uniform charge over the area A, and then on top of this another charge over the area B, so that the resulting total charge is distributed as shown in Figure 10b. In this case the variable charge will be represented by the function $I_A + I_B$. Such a sum of indicators is a very special case of sum of random variables, which will occupy us in later chapters.

For the present let us return to formula (1.4.5) and note that if the two sets A and B are disjoint, then it indeed reduces to the sum of the indicators, because then at most one of the two indicators can take the

value 1, so that the maximum coincides with the sum, namely

$$0 \lor 0 = 0 + 0$$
, $0 \lor 1 = 0 + 1$, $1 \lor 0 = 1 + 0$.

Thus we have

$$I_{A+B} = I_A + I_B$$
 provided $A \cap B = \emptyset$. (1.4.6)

As a particular case, we have for any set A:

$$I_{\Omega} = I_A + I_{A^c}.$$

Now I_{Ω} is the constant function 1 (on Ω), hence we may rewrite the above as

$$I_{A^c} = 1 - I_A. (1.4.7)$$

We can now derive an interesting formula. Since $(A \cup B)^c = A^c B^c$, we get by applying (1.4.7), (1.4.4) and then (1.4.7) again:

$$I_{A \cup B} = 1 - I_{A^c B^c} = 1 - I_{A^c} I_{B^c} = 1 - (1 - I_A)(1 - I_B).$$

Multiplying out the product (we are dealing with numerical functions!) and transposing terms we obtain

$$I_{A \cup B} + I_{A \cap B} = I_A + I_B. \tag{1.4.8}$$

Finally we want to investigate $I_{A\triangle B}$. We need a bit of arithmetic (also called number theory) first. All integers can be classified as even or odd, depending on whether the remainder we get when we divide it by 2 is 0 or 1. Thus each integer may be identified with (or reduced to) 0 or 1, provided we are only interested in its *parity* and not its exact value. When integers are added or subtracted subject to this reduction, we say we are operating $modulo\ 2$. For instance:

$$5+7+8-1+3=1+1+0-1+1=2=0$$
, modulo 2

A famous case of this method of counting occurs when the maiden picks off the petals of some wild flower one by one and murmurs: "he loves me," "he loves me not" in turn. Now you should be able to verify the following equation for every ω :

$$I_{A\triangle B} = I_A(\omega) + I_B(\omega) - 2I_{AB}(\omega)$$

= $I_A(\omega) + I_B(\omega)$, modulo 2. (1.4.9)

We can now settle a question raised in §1.3 and establish without pain the identity:

$$(A \triangle B) \triangle C = A \triangle (B \triangle C). \tag{1.4.10}$$

Proof: Using (1.4.9) twice we have

$$I_{(A\triangle B)\triangle C} = I_{A\triangle B} + I_C = (I_A + I_B) + I_C, \quad \text{modulo 2.} \quad (1.4.11)$$

Now if you have understood the meaning of addition modulo 2 you should see at once that it is an associative operation (what does that mean, "modulo 2"?). Hence the last member of (1.4.11) is equal to

$$I_A + (I_B + I_C) = I_A + I_{B \triangle C} = I_{A \triangle (B \triangle C)},$$
 modulo 2.

We have therefore shown that the two sets in (1.4.10) have identical indicators, hence they are identical. Q.E.D.

We do not need this result below. We just want to show that a trick is sometimes neater than a picture!

Exercises

- 1. Why is the sequence of numbers $\{1, 2, 1, 2, 3\}$ not a set?
- 2. If two sets have the same size, are they then identical?
- 3. Can a set and a proper subset have the same size? (A *proper* subset is a subset that is not also a superset!)
- 4. If two sets have identical complements, then they are themselves identical. Show this in two ways: (i) by verbal definition, (ii) by using formula (1.2.1).
- 5. If A, B, C have the same meanings as in Section 1.2, what do the following sets mean:

$$A \cup (B \cap C); \quad (A \setminus B) \setminus C; \quad A \setminus (B \setminus C).$$

6. Show that

$$(A \cup B) \cap C \neq A \cup (B \cap C);$$

but also give some special cases where there is equality.

7. Using the atoms given in the decomposition (1.3.5), express

$$A \cup B \cup C$$
; $(A \cup B)(B \cup C)$; $A \setminus B$; $A \triangle B$;

the set of ω which belongs to exactly 1 [exactly 2; at least 2] of the sets A,B,C.

8. Show that $A \subset B$ if and only if AB = A; or $A \cup B = B$. (So the relation of inclusion can be defined through identity and the operations.)

- 9. Show that A and B are disjoint if and only if $A \setminus B = A$; or $A \cup B = A \triangle B$. (After No. 8 is done, this can be shown purely symbolically without going back to the verbal definitions of the sets.)
- 10. Show that there is a distributive law also for difference:

$$(A \setminus B) \cap C = (A \cap C) \setminus (B \cap C).$$

Is the dual

$$(A \cap B) \setminus C = (A \setminus C) \cap (B \setminus C)$$

also true?

- 11. Derive (D_2) from (D_1) by using (C_1) and (C_2) .
- *12. Show that

$$(A \cup B) \setminus (C \cup D) \subset (A \setminus C) \cup (B \setminus D).$$

*13. Let us define a new operation "/" as follows:

$$A/B = A^c \cup B$$
.

Show that

- (i) $(A / B) \cap (B / C) \subset A / C$;
- (ii) $(A / B) \cap (A / C) = A / BC$;
- (iii) $(A / B) \cap (B / A) = (A \triangle B)^c$.

In intuitive logic, "A/B" may be read as "A implies B." Use this to interpret the relations above.

- *14. If you like a "dirty trick" this one is for you. There is an operation between two sets A and B from which alone all the operations defined above can be derived. [Hint: It is sufficient to derive complement and union from it. Look for some combination that contains these two. It is not unique.]
 - 15. Show that $A \subset B$ if and only if $I_A \leq I_B$; and $A \cap B = \emptyset$ if and only if $I_A I_B = 0$.
 - 16. Think up some concrete schemes that illustrate formula (1.4.8).
 - 17. Give a direct proof of (1.4.8) by checking it for all ω . You may use the atoms in (1.3.4) if you want to be well organized.
 - 18. Show that for any real numbers a and b, we have

$$a + b = (a \lor b) + (a \land b).$$

Use this to prove (1.4.8) again.

19. Express $I_{A\setminus B}$ and I_{A-B} in terms of I_A and I_B .

- 20. Express $I_{A\cup B\cup C}$ as a polynomial of I_A,I_B,I_C . [Hint: Consider $1-I_{A\cup B\cup C}$.]
- *21. Show that

$$I_{ABC} = I_A + I_B + I_C - I_{A \cup B} - I_{A \cup C} - I_{B \cup C} + I_{A \cup B \cup C}.$$

You can verify this directly, but it is nicer to derive it from No. 20 by duality.

Probability

2.1. Examples of probability

We learned something about sets in Chapter 1; now we are going to measure them. The most primitive way of measuring is to count the number, so we will begin with such an example.

Example 1. In Example (a') of §1.1, suppose that the number of rotten apples is 28. This gives a measure to the set A described in (a'), called its size and denoted by |A|. But it does not tell anything about the total number of apples in the bushel, namely the size of the sample space Ω given in Example (a). If we buy a bushel of apples we are more likely to be concerned with the relative *proportion* of rotten ones in it rather than their absolute number. Suppose then the total number is 550. If we now use the letter P provisionarily for "proportion," we can write this as follows:

$$P(A) = \frac{|A|}{|\Omega|} = \frac{28}{550}. (2.1.1)$$

Suppose next that we consider the set B of unripe apples in the same bushel, whose number is 47. Then we have similarly

$$P(B) = \frac{|B|}{|\Omega|} = \frac{47}{550}.$$

It seems reasonable to suppose that an apple cannot be both rotten and unripe (this is really a matter of definition of the two adjectives); then the

two sets are disjoint so their members do not overlap. Hence the number of "rotten or unripe apples" is equal to the sum of the number of "rotten apples" and the number of "unripe apples": 28 + 47 = 75. This may be written in symbols as:

$$|A + B| = |A| + |B|. (2.1.2)$$

If we now divide through by $|\Omega|$, we obtain

$$P(A+B) = P(A) + P(B). (2.1.3)$$

On the other hand, if some apples can be rotten and unripe at the same time, such as when worms got into green ones, then the equation (2.1.2) must be replaced by an inequality:

$$|A \cup B| \le |A| + |B|,$$

which leads to

$$P(A \cup B) \le P(A) + P(B). \tag{2.1.4}$$

Now what is the excess of |A| + |B| over $|A \cup B|$? It is precisely the number of "rotten and unripe apples," that is, $|A \cap B|$. Thus

$$|A\cup B|+|A\cap B|=|A|+|B|,$$

which yields the pretty equation

$$P(A \cup B) + P(A \cap B) = P(A) + P(B).$$
 (2.1.5)

Example 2. A more sophisticated way of measuring a set is the area of a plane set as in Examples (f) and (f') of §1.1, or the volume of a solid. It is said that the measurement of land areas was the origin of geometry and trigonometry in ancient times. While the nomads were still counting on their fingers and toes as in Example 1, the Chinese and Egyptians, among other peoples, were subdividing their arable lands, measuring them in units and keeping accounts of them on stone tablets or papyrus. This unit varied a great deal from one civilization to another (who knows the conversion rate of an acre into mou's or hectares?). But again it is often the ratio of two areas that concerns us as in the case of a wild shot that hits the target board. The proportion of the area of a subset A to that of Ω may be written, if we denote the area by the symbol $|\cdot|$:

$$P(A) = \frac{|A|}{|\Omega|}. (2.1.6)$$

This means also that if we fix the unit so that the total area of Ω is 1 unit, then the area of A is equal to the fraction P(A) in this scale. Formula (2.1.6) looks just like formula (2.1.1) by the deliberate choice of notation in order to underline the similarity of the two situations. Furthermore, for two sets A and B the previous relations (2.1.3) to (2.1.5) hold equally well in their new interpretations.

Example 3. When a die is thrown there are six possible outcomes. If we compare the process of throwing a particular number [face] with that of picking a particular apple in Example 1, we are led to take $\Omega = \{1, 2, 3, 4, 5, 6\}$ and define

$$P({k}) = \frac{1}{6}, \quad k = 1, 2, 3, 4, 5, 6.$$
 (2.1.7)

Here we are treating the six outcomes as "equally likely," so that the same measure is assigned to all of them, just as we have done tacitly with the apples. This hypothesis is usually implied by saying that the die is "perfect." In reality, of course, no such die exists. For instance, the mere marking of the faces would destroy the perfect symmetry; and even if the die were a perfect cube, the outcome would still depend on the way it is thrown. Thus we must stipulate that this is done in a perfectly symmetrical way too, and so on. Such conditions can be approximately realized and constitute the basis of an assumption of equal likelihood on grounds of symmetry.

Now common sense demands an empirical interpretation of the "probability" given in (2.1.7). It should give a measure of what is likely to happen, and this is associated in the intuitive mind with the observable frequency of occurrence . Namely, if the die is thrown a number of times, how often will a particular face appear? More generally, let A be an event determined by the outcome; e.g., "to throw a number not less than 5 [or an odd number]." Let $N_n(A)$ denote the number of times the event A is observed in n throws; then the relative frequency of A in these trials is given by the ratio

$$Q_n(A) = \frac{N_n(A)}{n}. (2.1.8)$$

There is good reason to take this Q_n as a measure of A. Suppose B is another event such that A and B are incompatible or mutually exclusive in the sense that they cannot occur in the same trial. Clearly we have $N_n(A+B) = N_n(A) + N_n(B)$, and consequently

$$Q_n(A+B) = \frac{N_n(A+B)}{n}$$

$$= \frac{N_n(A) + N_n(B)}{n} = \frac{N_n(A)}{n} + \frac{N_n(B)}{n} = Q_n(A) + Q_n(B).$$
(2.1.9)

Similarly for any two events A and B in connection with the same game, not necessarily incompatible, the relations (2.1.4) and (2.1.5) hold with the P's there replaced by our present Q_n . Of course, this Q_n depends on n and will fluctuate, even wildly, as n increases. But if you let n go to infinity, will the sequence of ratios $Q_n(A)$ "settle down to a steady value"? Such a question can never be answered empirically, since by the very nature of a limit we cannot put an end to the trials. So it is a mathematical idealization to assume that such a limit does exist, and then write

$$Q(A) = \lim_{n \to \infty} Q_n(A). \tag{2.1.10}$$

We may call this the empirical limiting frequency of the event A. If you know how to operate with limits, then you can see easily that the relation (2.1.9) remains true "in the limit." Namely when we let $n \to \infty$ everywhere in that formula and use the definition (2.1.10), we obtain (2.1.3) with P replaced by Q. Similarly, (2.1.4) and (2.1.5) also hold in this context.

But the limit Q still depends on the actual sequence of trials that are carried out to determine its value. On the face of it, there is no guarantee whatever that another sequence of trials, even if it is carried out under the same circumstances, will yield the same value. Yet our intuition demands that a measure of the likelihood of an event such as A should tell something more than the mere record of one experiment. A viable theory built on the frequencies will have to assume that the Q defined above is in fact the same for all similar sequences of trials. Even with the hedge implicit in the word "similar," that is assuming a lot to begin with. Such an attempt has been made with limited success, and has a great appeal to common sense, but we will not pursue it here. Rather, we will use the definition in (2.1.7) which implies that if A is any subset of Ω and |A| its size, then

$$P(A) = \frac{|A|}{|\Omega|} = \frac{|A|}{6}.$$
 (2.1.11)

For example, if A is the event "to throw an odd number," then A is identified with the set $\{1,3,5\}$ and P(A) = 3/6 = 1/2.

It is a fundamental proposition in the theory of probability that under certain conditions (repeated independent trials with identical die), the limiting frequency in (2.1.10) will indeed exist and be equal to P(A) defined in (2.1.11), for "practically all" conceivable sequences of trials. This celebrated theorem, called the Law of Large Numbers, is considered to be the cornerstone of all empirical sciences. In a sense it justifies the intuitive foundation of probability as frequency discussed above. The precise statement and derivation will be given in Chapter 7. We have made this early announcement to quiet your feelings or misgivings about frequencies and to concentrate for the moment on sets and probabilities in the following sections.

2.2. Definition and illustrations

First of all, a probability is a number associated with or assigned to a set in order to measure it in some sense. Since we want to consider many sets at the same time (that is why we studied Chapter 1), and each of them will have a probability associated with it, this makes probability a "function of sets." You should have already learned in some mathematics course what a function means; in fact, this notion is used a little in Chapter 1. Nevertheless, let us review it in the familiar notation: a function f defined for some or all real numbers is a rule of association, by which we assign the number f(x) to the number x. It is sometimes written as $f(\cdot)$, or more painstakingly as follows:

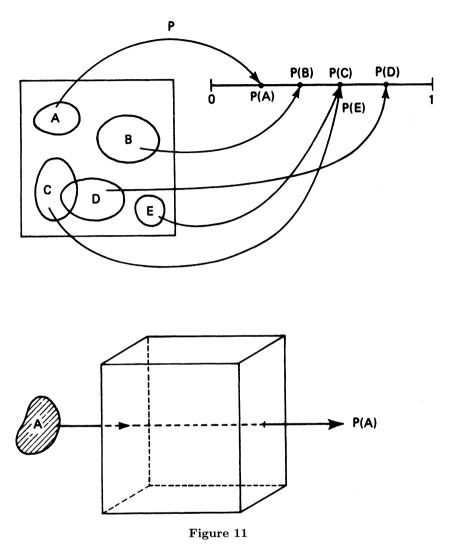
$$f: x \to f(x). \tag{2.2.1}$$

So when we say a probability is a function of sets we mean a similar association, except that x is replaced by a set S:

$$P: S \to P(S). \tag{2.2.2}$$

The value P(S) is still a number; indeed it will be a number between 0 and 1. We have not been really precise in (2.2.1), because we have not specified the set of x there for which it has a meaning. This set may be the interval (a,b) or the half-line $(0,\infty)$ or some more complicated set called the domain of f. Now what is the domain of our probability function P? It must be a set of sets or, to avoid the double usage, a family (class) of sets. As in Chapter 1 we are talking about subsets of a fixed sample space Ω . It would be nice if we could use the family of all subsets of Ω , but unexpected difficulties will arise in this case if no restriction is imposed on Ω . We might say that if Ω is too large, namely when it contains uncountably many points, then it has too many subsets, and it becomes impossible to assign a probability to each of them and still satisfy a basic rule [Axiom (ii*) ahead] governing the assignments. However, if Ω is a finite or countably infinite set, then no such trouble can arise and we may indeed assign a probability to each and all of its subsets. This will be shown at the beginning of §2.4. You are supposed to know what a finite set is (although it is by no means easy to give a logical definition, while it is mere tautology to say that "it has only a finite number of points"); let us review what a countably infinite set is. This notion will be of sufficient importance to us, even if it only lurks in the background most of the time.

A set is countably infinite when it can be put into 1-to-1 correspondence with the set of positive integers. This correspondence can then be exhibited by labeling the elements as $\{s_1, s_2, \ldots, s_n, \ldots\}$. There are, of course, many ways of doing this, for instance we can just let some of the elements swap labels (or places if they are thought of being laid out in a row). The set of positive rational numbers is countably infinite, hence they can be labeled



in some way as $\{r_1, r_2, \ldots, r_n, \ldots\}$, but don't think for a moment that you can do this by putting them in increasing order as you can with the positive integers $1 < 2 < \cdots < n < \cdots$. From now on we shall call a set countable when it is either finite or countably infinite. Otherwise it is called uncountable. For example, the set of all real numbers is uncountable. We shall deal with uncountable sets later, and we will review some properties of a countable set when we need them. For the present we will assume the sample space Ω to be countable in order to give the following definition in its simplest form, without a diverting complication. As a matter of fact, we could even assume Ω to be finite as in Examples (a) to (e) of §1.1, without losing the essence of the discussion below.

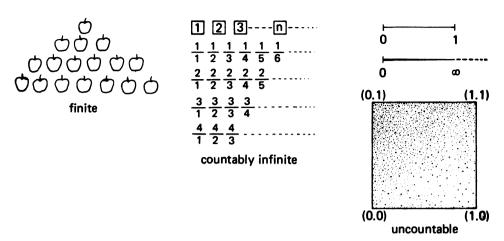


Figure 12

Definition. A probability measure on the sample space Ω is a function of subsets of Ω satisfying three axioms:

- (i) For every set $A \subset \Omega$, the value of the function is a nonnegative number: $P(A) \geq 0$.
- (ii) For any two disjoint sets A and B, the value of the function for their union A + B is equal to the sum of its value for A and its value for B:

$$P(A+B) = P(A) + P(B)$$
 provided $AB = \emptyset$.

(iii) The value of the function for Ω (as a subset) is equal to 1:

$$P(\Omega) = 1.$$

Observe that we have been extremely careful in distinguishing the function $P(\cdot)$ from its values such as P(A), P(B), P(A+B), $P(\Omega)$. Each of these is "a probability," but the function itself should properly be referred to as a "probability measure" as indicated.

Example 1 in §2.1 shows that the proportion P defined there is in fact a probability measure on the sample space, which is a bushel of 550 apples. It assigns a probability to every subset of these apples, and this assignment satisfies the three axioms above. In Example 2 if we take Ω to be all the land that belonged to the Pharaoh, it is unfortunately not a countable set. Nevertheless we can define the area for a very large class of subsets that are called "measurable," and if we restrict ourselves to these subsets only, the "area function" is a probability measure as shown in Example 2 where this restriction is ignored. Note that Axiom (iii) reduces to a convention: the decree of a unit. Now how can a land area not be measurable? While

this is a sophisticated mathematical question that we will not go into in this book, it is easy to think of practical reasons for the possibility: the piece of land may be too jagged, rough, or inaccessible (see Fig. 13).

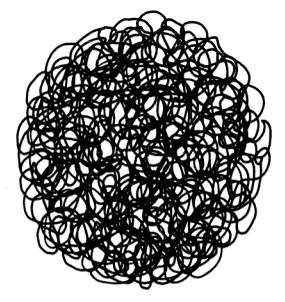


Figure 13

In Example 3 we have shown that the empirical relative frequency is a probability measure. But we will not use this definition in this book. Instead, we will use the first definition given at the beginning of Example 3, which is historically the earliest of its kind. The general formulation will now be given.

Example 4. A classical enunciation of probability runs as follows. The probability of an event is the ratio of the number of cases *favorable* to that event to the total number of cases, provided that all these are *equally likely*

To translate this into our language: the sample space is a finite set of possible cases: $\{\omega_1, \omega_2, \ldots, \omega_m\}$, each ω_i being a "case." An event A is a subset $\{\omega_{i_1}, \omega_{i_2}, \ldots, \omega_{i_n}\}$, each ω_{i_j} being a "favorable case." The probability of A is then the ratio

$$P(A) = \frac{|A|}{|\Omega|} = \frac{n}{m}.$$
(2.2.3)

As we see from the discussion in Example 1, this defines a probability measure P on Ω anyway, so that the stipulation above that the cases be equally likely is superfluous from the axiomatic point of view. Besides, what

does it really mean? It sounds like a bit of tautology, and how is one going to decide whether the cases are equally likely or not?

A celebrated example will illustrate this. Let two coins be tossed. D'Alembert (mathematician, philosopher, and encyclopedist, 1717–83) argued that there are three possible cases, namely:

(i) both heads, (ii) both tails, (iii) a head and a tail.

So he went on to conclude that the probability of "a head and a tail" is equal to 1/3. If he had figured that this *probability* should have something to do with the experimental *frequency* of the occurrence of the event, he might have changed his mind after tossing two coins more than a few times. (History does not record if he ever did that, but it is said that for centuries people believed that men had more teeth than women because Aristotle had said so, and apparently nobody bothered to look into a few mouths.) The three cases he considered are not equally likely. Case (iii) should be split into two:

- (iiia) first coin shows head and second coin shows tail.
- (iiib) first coin shows tail and second coin shows head.

It is the four cases (i), (ii), (iiia) and (iiib) that are equally likely by symmetry and on empirical evidence. This should be obvious if we toss the two coins one after the other rather than simultaneously. However, there is an important point to be made clear here. The two coins may be physically indistinguishable so that in so far as actual observation is concerned, D'Alembert's three cases are the only distinct patterns to be recognized. In the model of two coins they happen not to be equally likely on the basis of common sense and experimental evidence. But in an analogous model for certain microcosmic particles, called Bose–Einstein statistics (see Exercise 24 of Chapter 3), they are indeed assumed to be equally likely in order to explain some types of physical phenomena. Thus what we regard as "equally likely" is a matter outside the axiomatic formulation. To put it another way, if we use (2.2.3) as our definition of probability then we are in effect treating the ω 's as equally likely, in the sense that we count only their numbers and do not attach different weights to them.

Example 5. If six dice are rolled, what is the probability that all show different faces?

This is just Example (e) and (e'). It is stated elliptically on purpose to get you used to such problems. We have already mentioned that the total number of possible outcomes is equal to $6^6 = 46656$. They are supposed to be all "equally likely" although we never breathed a word about this assumption. Why, nobody can solve the problem as announced without such an assumption. Other data about the dice would have to be given before we

could begin—which is precisely the difficulty when similar problems arise in practice. Now if the dice are all perfect, and the mechanism by which they are rolled is also perfect, which excludes any collusion between the movements of the several dice, then our hypothesis of equal likelihood may be justified. Such conditions are taken for granted in a problem like this when nothing is said about the dice. The solution is then given by (2.2.3) with $n=6^6$ and m=6! (see Example 2 in §3.1 for these computations):

$$\frac{6!}{6^6} = \frac{720}{46656} = .015432$$

approximately.

Let us note that if the dice are not distinguishable from each other, then to the observer there is exactly one *pattern* in which the six dice show different faces. Similarly, the total number of different patterns when six dice are rolled is much smaller than 6⁶ (see Example 3 of §3.2). Yet when we count the possible outcomes we must think of the dice as distinguishable, as if they were painted in different colors. This is one of the vital points to grasp in the counting cases; see Chapter 3.

In some situations the equally likely cases must be searched out. This point will be illustrated by a famous historical problem called the "problem of points."

Example 6. Two players A and B play a series of games in which the probability of each winning a single game is equal to 1/2, irrespective [independent] of the outcomes of other games. For instance, they may play tennis in which they are equally matched, or simply play "heads or tails" by tossing an unbiased coin. Each player gains a "point" when he wins a game, and nothing when he loses. Suppose that they stop playing when A needs 2 more points and B needs 3 more points to win the stake. How should they divide it fairly?

It is clear that the winner will be decided in 4 more games. For in those 4 games either A will have won ≥ 2 points or B will have won ≥ 3 points, but not both. Let us enumerate all the possible outcomes of these 4 games using the letter A or B to denote the winner of each game:

These are equally likely cases on grounds of symmetry. There are* $\binom{4}{4}$ + $\binom{4}{3}$ + $\binom{4}{2}$ = 11 cases in which A wins the stake; and $\binom{4}{3}$ + $\binom{4}{4}$ = 5 cases

^{*}See (3.2.3) for notation used below.

in which B wins the stake. Hence the stake should be divided in the ratio 11:5. Suppose it is \$64000; then A gets \$44000, B gets \$20000. [We are taking the liberty of using the dollar as currency; the United States did not exist at the time when the problem was posed.]

This is Pascal's solution in a letter to Fermat dated August 24, 1654. [Blaise Pascal (1623–62); Pierre de Fermat (1601–65); both among the greatest mathematicians of all time.] Objection was raised by a learned contemporary (and repeated through the ages) that the enumeration above was not reasonable, because the series would have stopped as soon as the winner was decided and not have gone on through all 4 games in some cases. Thus the real possibilities are as follows:

But these are not equally likely cases. In modern terminology, if these 10 cases are regarded as constituting the sample space, then

$$P(AA) = \frac{1}{4}, \quad P(ABA) = P(BAA) = P(BBB) = \frac{1}{8},$$

$$P(ABBA) = P(BABA) = P(BBAA) = P(ABBB)$$

$$= P(BABB) = P(BBAB) = \frac{1}{16}$$

since A and B are independent events with probability 1/2 each (see §2.4). If we add up these probabilities we get of course

$$P(A \text{ wins the stake}) = \frac{1}{4} + \frac{1}{8} + \frac{1}{16} + \frac{1}{8} + \frac{1}{16} + \frac{1}{16} = \frac{11}{16},$$

$$P(B \text{ wins the stake}) = \frac{1}{16} + \frac{1}{16} + \frac{1}{16} + \frac{1}{8} = \frac{5}{16}.$$

Pascal did not quite explain his method this way, saying merely that "it is absolutely equal and indifferent to each whether they play in the natural way of the game, which is to finish as soon as one has his score, or whether they play the entire four games." A later letter by him seems to indicate that he fumbled on the same point in a similar problem with three players. The student should take heart that this kind of reasoning was not easy even for past masters.

2.3. Deductions from the axioms

In this section we will do some simple "axiomatics." That is to say, we shall deduce some properties of the probability measure from its definition, using, of course, the axioms but nothing else. In this respect the axioms of a mathematical theory are like the constitution of a government. Unless and until it is changed or amended, every *law* must be made to follow from it. In mathematics we have the added assurance that there are no divergent views as to how the constitution should be construed.

We record some consequences of the axioms in (iv) to (viii) below. First of all, let us show that a probability is indeed a number between 0 and 1.

(iv) For any set A, we have

$$P(A) \leq 1$$
.

This is easy, but you will see that in the course of deducing it we shall use all three axioms. Consider the complement A^c as well as A. These two sets are disjoint and their union is Ω :

$$A + A^c = \Omega. (2.3.1)$$

So far, this is just set theory, no probability theory yet. Now use Axiom (ii) on the left side of (2.3.1) and Axiom (iii) on the right:

$$P(A) + P(A^c) = P(\Omega) = 1.$$
 (2.3.2)

Finally use Axiom (i) for A^c to get

$$P(A) = 1 - P(A^c) \le 1.$$

Of course, the first inequality above is just Axiom (i). You might object to our slow pace above by pointing out that since A is contained in Ω , it is obvious that $P(A) \leq P(\Omega) = 1$. This reasoning is certainly correct, but we still have to pluck it from the axioms, and that is the point of the little proof above. We can also get it from the following more general proposition.

(v) For any two sets such that $A \subset B$, we have

$$P(A) \le P(B)$$
, and $P(B-A) = P(B) - P(A)$.

The proof is an imitation of the preceding one with B playing the role of Ω . We have

$$B = A + (B - A),$$

 $P(B) = P(A) + P(B - A) > P(A).$

The next proposition is such an immediate extension of Axiom (ii) that we could have adopted it instead as an axiom.

(vi) For any finite number of disjoint sets A_1, \ldots, A_n , we have

$$P(A_1 + \dots + A_n) = P(A_1) + \dots + P(A_n). \tag{2.3.3}$$

This property of the probability measure is called *finite additivity*. It is trivial if we recall what "disjoint" means and use (ii) a few times; or we may proceed by induction if we are meticulous. There is an important extension of (2.3.3) to a countable number of sets later, *not* obtainable by induction!

As already checked in several special cases, there is a generalization of Axiom (ii), hence also of (2.3.3), to sets that are not necessarily disjoint. You may find it trite, but it has the dignified name of *Boole's inequality*. Boole (1815–64) was a pioneer in the "laws of thought" and author of *Theories of Logic and Probabilities*.

(vii) For any finite number of arbitrary sets A_1, \ldots, A_n , we have

$$P(A_1 \cup \dots \cup A_n) \le P(A_1) + \dots + P(A_n). \tag{2.3.4}$$

Let us first show this when n = 2. For any two sets A and B, we can write their union as the sum of disjoint sets as follows:

$$A \cup B = A + A^c B. \tag{2.3.5}$$

Now we apply Axiom (ii) to get

$$P(A \cup B) = P(A) + P(A^{c}B).$$
 (2.3.6)

Since $A^cB \subset B$, we can apply (v) to get (2.3.4).

The general case follows easily by mathematical induction, and you should write it out as a good exercise on this method. You will find that you need the associative law for the union of sets as well as that for the addition of numbers.

The next question is the difference between the two sides of the inequality (2.3.4). The question is somewhat moot since it depends on what we want to use to express the difference. However, when n=2 there is a clear answer.

(viii) For any two sets A and B, we have

$$P(A \cup B) + P(A \cap B) = P(A) + P(B).$$
 (2.3.7)

This can be gotten from (2.3.6) by observing that $A^cB = B - AB$, so that we have by virtue of (v):

$$P(A \cup B) = P(A) + P(B - AB) = P(A) + P(B) - P(AB),$$

which is equivalent to (2.3.7). Another neat proof is given in Exercise 12.

We shall postpone a discussion of the general case until §6.2. In practice, the inequality is often more useful than the corresponding identity which is rather complicated.

We will not quit formula (2.3.7) without remarking on its striking resemblance to formula (1.4.8) of §1.4, which is repeated below for the sake of comparison:

$$I_{A \cup B} + I_{A \cap B} = I_A + I_B. \tag{2.3.8}$$

There is indeed a deep connection between the pair, as follows. The probability P(S) of each set S can be obtained from its indicator function I_S by a procedure (operation) called "taking expectation" or "integration." If we perform this on (2.3.8) term by term, their result is (2.3.7). This procedure is an essential part of probability theory and will be thoroughly discussed in Chapter 6. See Exercise 19 for a special case.

To conclude our axiomatics, we will now strengthen Axiom (ii) or its immediate consequence (vi), namely the finite additivity of P, into a new axiom.

(ii*) Axiom of countable additivity. For a countably infinite collection of disjoint sets A_k , $k=1,2,\ldots$, we have

$$P\left(\sum_{k=1}^{\infty} A_k\right) = \sum_{k=1}^{\infty} P(A_k). \tag{2.3.9}$$

This axiom includes (vi) as a particular case, for we need only put $A_k = \emptyset$ for k > n in (2.3.9) to obtain (2.3.3). The empty set is disjoint from any other set including itself and has probability zero (why?). If Ω is a finite set, then the new axiom reduces to the old one. But it is important to see why (2.3.9) cannot be deduced from (2.3.3) by letting $n \to \infty$. Let us try this by rewriting (2.3.3) as follows:

$$P\left(\sum_{k=1}^{n} A_k\right) = \sum_{k=1}^{n} P(A_k). \tag{2.3.10}$$

Since the left side above cannot exceed 1 for all n, the series on the right side must converge and we obtain

$$\lim_{n \to \infty} P\left(\sum_{k=1}^{n} A_k\right) = \lim_{n \to \infty} \sum_{k=1}^{n} P(A_k) = \sum_{k=1}^{\infty} P(A_k).$$
 (2.3.11)

Comparing this established result with the desired result (2.3.9), we see that the question boils down to

$$\lim_{n \to \infty} P\left(\sum_{k=1}^{n} A_k\right) = P\left(\sum_{k=1}^{\infty} A_k\right),\,$$

which can be exhibited more suggestively as

$$\lim_{n \to \infty} P\left(\sum_{k=1}^{n} A_k\right) = P\left(\lim_{n \to \infty} \sum_{k=1}^{n} A_k\right). \tag{2.3.12}$$

See end of $\S1.3$ (see Fig. 14).

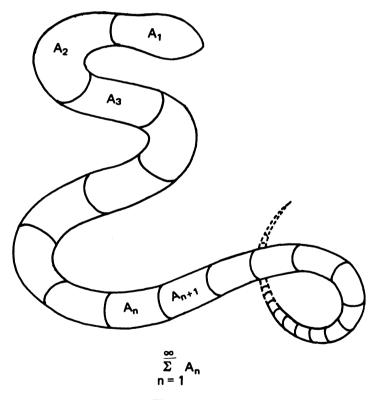


Figure 14

Thus it is a matter of interchanging the two operations "lim" and "P" in (2.3.12), or you may say, "taking the limit inside the probability relation." If you have had enough calculus you know this kind of interchange is often hard to justify and may be illegitimate or even invalid. The new axiom is created to secure it in the present case and has fundamental consequences in the theory of probability.

2.4. Independent events

From now on, a "probability measure" will satisfy Axioms (i), (ii*), and (iii). The subsets of Ω to which such a probability has been assigned will also be called an *event*.

We shall show how easy it is to *construct* probability measures for any countable space $\Omega = \{\omega_1, \omega_2, \dots, \omega_n, \dots\}$. To each sample point ω_n let us attach an arbitrary "weight" p_n subject only to the conditions

$$\forall n: \ p_n \ge 0, \ \sum_n p_n = 1.$$
 (2.4.1)

This means that the weights are positive or zero, and add up to 1 altogether. Now for any subset A of Ω , we define its probability to be the *sum of the weights of all the points in it.* In symbols, we put first

$$\forall n: \ P(\{\omega_n\}) = p_n; \tag{2.4.2}$$

and then for every $A \subset \Omega$:

$$P(A) = \sum_{\omega_n \in A} p_n = \sum_{\omega_n \in A} P(\{\omega_n\}).$$

We may write the last term above more neatly as

$$P(A) = \sum_{\omega \in A} P(\{\omega\}). \tag{2.4.3}$$

Thus P is a function defined for all subsets of Ω and it remains to check that it satisfies Axioms (i), (ii*), and (iii). This requires nothing but a bit of clearheaded thinking and is best done by yourself. Since the weights are quite arbitrary apart from the easy conditions in (2.4.1), you see that probability measures come "a dime a dozen" in a countable sample space. In fact, we can get them all by the above method of construction. For if any probability measure P is given, never mind how, we can define p_n to be $P(\{\omega_n\})$ as in (2.4.2), and then P(A) must be given as in (2.4.3), because of Axiom (ii*). Furthermore the p_n 's will satisfy (2.4.1) as a simple consequence of the axioms. In other words, any given P is necessarily of the type described by our construction.

In the very special case that Ω is finite and contains exactly m points, we may attach equal weights to all of them, so that

$$p_n = \frac{1}{m}, \quad n = 1, 2, \dots, m.$$

Then we are back to the "equally likely" situation in Example 4 of §2.2. But in general the p_n 's need not be equal, and when Ω is countably infinite

they cannot all be equal (why?). The preceding discussion shows the degree of arbitrariness involved in the general concept of a probability measure.

An important model of probability space is that of *repeated independent trials*: this is the model used when a coin is tossed, a die thrown, a card drawn from a deck (with replacement) several times. Alternately, we may toss several coins or throw several dice at the same time. Let us begin with an example.

Example 7. First toss a coin, then throw a die, finally draw a card from a deck of poker cards. Each trial produces an event; let

A = coin falls heads; B = die shows number 5 or 6;C = card drawn is a spade.

Assume that the coin is fair, the die is perfect, and the deck thoroughly shuffled. Furthermore assume that these three trials are carried out "independently" of each other, which means intuitively that the outcome of each trial does not influence that of the others. For instance, this condition is approximately fulfilled if the trials are done by different people in different places, or by the same person in different months! Then all possible joint outcomes may be regarded as equally likely. There are respectively 2, 6, and 52 possible cases for the individual trials, and the total number of cases for the whole set of trials is obtained by multiplying these numbers together: $2 \cdot 6 \cdot 52$ (as you will soon see it is better not to compute this product). This follows from a fundamental rule of counting, which is fully discussed in §3.1 and which you should read now if need be. [In general, many parts of this book may be read in different orders, back and forth.] The same rule yields the numbers of favorable cases to the events A, B, C, AB, AC, BC, ABC given below, where the symbol $|\dots|$ for size is used:

$$|A| = 1 \cdot 6 \cdot 52, \qquad |B| = 2 \cdot 2 \cdot 52, \qquad |C| = 2 \cdot 6 \cdot 13,$$

$$|AB| = 1 \cdot 2 \cdot 52, \quad |AC| = 1 \cdot 6 \cdot 13, \quad |BC| = 2 \cdot 2 \cdot 13,$$

$$|ABC| = 1 \cdot 2 \cdot 13.$$

Dividing these numbers by $|\Omega| = 2 \cdot 6 \cdot 52$, we obtain after quick cancellation of factors:

$$P(A) = \frac{1}{2},$$
 $P(B) = \frac{1}{3},$ $P(C) = \frac{1}{4},$ $P(AB) = \frac{1}{6},$ $P(AC) = \frac{1}{8},$ $P(BC) = \frac{1}{12},$ $P(ABC) = \frac{1}{24}.$

We see at a glance that the following set of equations holds:

$$P(AB) = P(A)P(B), \ P(AC) = P(A)P(C), \ P(BC) = P(B)P(C)$$
 (2.4.4)
 $P(ABC) = P(A)P(B)P(C).$

The reader is now asked to convince himself that this set of relations will also hold for any three events A, B, C such that A is determined by the coin, B by the die, and C by the card drawn alone. When this is the case we say that these trials are stochastically independent as well as the events so produced. The adverb "stochastically" is usually omitted for brevity.

The astute reader may observe that we have not formally defined the word "trial," and yet we are talking about independent trials! A logical construction of such objects is quite simple but perhaps a bit too abstract for casual introduction. It is known as "product space"; see Exercise 29. However, it takes less fuss to define "independent events" and we shall do so at once.

Two events A and B are said to be independent if we have P(AB) = P(A)P(B). Three events A, B, and C are said to be independent if the relations in (2.4.4) hold. Thus independence is a notion relative to a given probability measure (by contrast, the notion of disjointness, e.g., does not depend on any probability). More generally, the n events A_1, A_2, \ldots, A_n are independent if the intersection [joint occurrence] of any subset of them has as its probability the product of probabilities of the individual events. If you find this sentence too long and involved, you may prefer the following symbolism. For any subset (i_1, i_2, \ldots, i_k) of $(1, 2, \ldots, n)$, we have

$$P(A_{i_1} \cap A_{i_2} \cap \dots \cap A_{i_k}) = P(A_{i_1})P(A_{i_2}) \cdots P(A_{i_k}).$$
 (2.4.5)

Of course, here the indices i_1, \ldots, i_k are distinct and $1 \le k \le n$.

Further elaboration of the notion of independence is postponed to §5.5, because it will be better explained in terms of random variables. But we shall briefly describe a classical scheme—the grand daddy of repeated trials, and subject of intensive and extensive research by J. Bernoulli, De Moivre, Laplace, . . . , Borel,

Example 8. (The coin-tossing scheme). A coin is tossed repeatedly n times. The joint outcome may be recorded as a sequence of H's and T's, where H = "head," T = "tail." It is often convenient to quantify by putting H = 1, T = 0; or H = 1, T = -1; we shall adopt the first usage here. Then the result is a sequence of 0's and 1's consisting of n terms such as 110010110 with n = 9. Since there are 2 outcomes for each trial, there are 2^n possible joint outcomes. This is another application of the fundamental rule in §3.1. If all of these are assumed to be equally likely so that each particular joint outcome has probability $1/2^n$, then we can proceed as in Example 7 to verify that the trials are independent and the coin is fair. You will find this

a dull exercise, but it is recommended that you go through it in your head if not on paper. However, we will turn the table around here by assuming at the outset that the successive tosses do form independent trials. On the other hand, we do not assume the coin to be "fair," but only that the probabilities for head (H) and tail (T) remain constant throughout the trials. Empirically speaking, this is only approximately true since things do not really remain unchanged over long periods of time. Now we need a precise notation to record complicated statements, ordinary words being often awkward or ambiguous. Then let X_i denote the outcome of the ith trial and let ϵ_i denote 0 or 1 for each i, but of course varying with the subscript. Then our hypothesis above may be written as follows:

$$P(X_i = 1) = p;$$
 $P(X_i = 0) = 1 - p;$ $i = 1, 2, ..., n;$ (2.4.6)

where p is the probability of heads for each trial. For any particular, namely completely specified, sequence $(\epsilon_1, \epsilon_2, \dots, \epsilon_n)$ of 0's and 1's, the probability of the corresponding sequence of outcomes is equal to

$$P(X_1 = \epsilon_1, X_2 = \epsilon_2, \dots, X_n = \epsilon_n)$$

$$= P(X_1 = \epsilon_1) P(X_2 = \epsilon_2) \dots P(X_n = \epsilon_n)$$
(2.4.7)

as a consequence of independence. Now each factor on the right side above is equal to p or 1-p depending on whether the corresponding ϵ_i is 1 or 0. Suppose j of these are 1's and n-j are 0's; then the quantity in (2.4.7) is equal to

$$p^{j}(1-p)^{n-j}. (2.4.8)$$

Observe that for each sequence of trials, the number of heads is given by the sum $\sum_{i=1}^{n} X_i$. It is important to understand that the number in (2.4.8) is not the probability of obtaining j heads in n tosses, but rather that of obtaining a specific sequence of heads and tails in which there are j heads. In order to compute the former probability, we must count the total number of the latter sequences since all of them have the same probability given in (2.4.8). This number is equal to the binomial coefficient $\binom{n}{j}$; see §3.2 for a full discussion. Each one of these $\binom{n}{j}$ sequences corresponds to one possibility of obtaining j heads in n trials, and these possibilities are disjoint. Hence it follows from the additivity of P that we have

$$\begin{split} P\!\!\left(\sum_{i=1}^n X_i \! = \! j\right) &= P\!\!\left(\text{exactly } j \text{ heads in } n \text{ trials}\right) \\ &= \! \binom{n}{j} P\!\!\left(\text{any specified sequence of } n \text{ trials with exactly } j \text{ heads}\right) \\ &= \! \binom{n}{j} p^j (1-p)^{n-j}. \end{split}$$

This famous result is known as *Bernoulli's formula*. We shall return to it many times in the book.

2.5. Arithmetical density*

We study in this section a very instructive example taken from arithmetic.

Example 9. Let Ω be the first 120 natural numbers $\{1, 2, ..., 120\}$. For the probability measure P we use the proportion as in Example 1 of $\S 2.1$. Now consider the sets

$$A = \{ \omega \mid \omega \text{ is a multiple of } 3 \},$$

 $B = \{ \omega \mid \omega \text{ is a multiple of } 4 \}.$

Then every third number of Ω belongs to A, and every fourth to B. Hence we get the proportions

$$P(A) = 1/3, \quad P(B) = 1/4.$$

What does the set AB represent? It is the set of integers that are divisible by 3 and by 4. If you have not entirely forgotten your school arithmetic, you know this is just the set of multiples of $3 \cdot 4 = 12$. Hence P(AB) = 1/12. Now we can use (viii) to get $P(A \cup B)$:

$$P(A \cup B) = P(A) + P(B) - P(AB) = 1/3 + 1/4 - 1/12 = 1/2.$$
 (2.5.1)

What does this mean? $A \cup B$ is the set of those integers in Ω which are divisible by 3 or by 4 (or by both). We can count them one by one, but if you are smart you see that you don't have to do this drudgery. All you have to do is to count up to 12 (which is 10% of the whole population Ω), and check them off as shown:

$$1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12.$$

$$\checkmark \checkmark \qquad \checkmark \qquad \checkmark$$

Six are checked (one checked twice), hence the proportion of $A \cup B$ among these 12 is equal to 6/12 = 1/2 as given by (2.5.1).

An observant reader will have noticed that in the case above we have also

$$P(AB) = 1/12 = 1/3 \cdot 1/4 = P(A) \cdot P(B).$$

^{*}This section may be omitted.

This is true because the two numbers 3 and 4 happen to be *relatively prime*, namely they have no common divisor except 1. Suppose we consider another set:

$$C = \{\omega \mid \omega \text{ is a multiple of 6}\}.$$

Then P(C) = 1/6, but what is P(BC) now? The set BC consists of those integers that are divisible by both 4 and 6, namely divisible by their least common multiple (remember that?), which is 12 and not the product $4 \cdot 6 = 24$. Thus P(BC) = 1/12. Furthermore, because 12 is the least common multiple we can again stop counting at 12 in computing the proportion of the set $B \cup C$. An actual counting gives the answer 4/12 = 1/3, which may also be obtained from the formula (2.3.7):

$$P(B \cup C) = P(B) + P(C) - P(BC) = 1/4 + 1/6 - 1/12 = 1/3.$$
 (2.5.2)

This example illustrates a point that arose in the discussion in Example 3 of §2.1. Instead of talking about the proportion of the multiples of 3, say, we can talk about its frequency. Here no rolling of any fortuitous dice is needed. God has given us those natural numbers (a great mathematician Kronecker said so), and the multiples of 3 occur at perfectly regular periods with the frequency 1/3. In fact, if we use $N_n(A)$ to denote the number of natural numbers up to and including n which belong to the set A, it is a simple matter to show that

$$\lim_{n \to \infty} \frac{N_n(A)}{n} = \frac{1}{3}.$$

Let us call this P(A), the limiting frequency of A. Intuitively, it should represent the chance of picking a number divisible by 3, if we can reach into the whole bag of natural numbers as if they were so many indistinguishable balls in an urn. Of course, similar limits exist for the sets B, C, AB, BC, etc. and have the values computed above. But now with this infinite sample space of "all natural numbers," call it Ω^* , we can treat by the same method any set of the form

$$A_m = \{ \omega \mid \omega \text{ is divisible by } m \}, \tag{2.5.3}$$

where m is an arbitrary natural number. Why then did we not use this more natural and comprehensive model?

The answer may be a surprise for you. By our definition of probability measure given in §2.2, we should have required that every subset of Ω^* have a probability, provided that Ω^* is countable, which is the case here. Now take for instance the set that consists of the single number {1971} or, if you prefer, the set $Z = \{\text{all numbers from 1 to 1971}\}$. Its probability is given by $\lim_{n\to\infty} N_n(Z)/n$ according to the same rule that was applied to the set

A. But $N_n(Z)$ is equal to 1971 for all values of $n \geq 1971$; hence the limit above is equal to 0 and we conclude that every finite set has probability 0 by this rule. If P were to be countably additive as required by Axiom (ii*) in §2.3, then $P(\Omega^*)$ would be 0 rather than 1. This contradiction shows that P cannot be a probability measure on Ω^* . Yet it works perfectly well for sets such as A_m .

There is a way out of this paradoxical situation. We must abandon our previous requirement that the measure be defined for all subsets (of natural numbers). Let a finite number of the sets A_m be given, and let us consider the *composite sets* that can be obtained from these by the operations complementation, union, and intersection. Call this class of sets the class *generated by* the original sets. Then it is indeed possible to define P in the manner prescribed above for all sets in *this* class. A set that is not in the class has no probability at all. For example, the set Z does not belong to the class generated by A, B, C. Hence its probability is *not* defined, rather than zero. We may also say that the set Z is *nonmeasurable* in the context of Example 2 of §2.1. This saves the situation, but we will not pursue it further here except to give another example.

Example 10. What is the probability of the set of numbers divisible by 3, not divisible by 5, and divisible by 4 or 6?

Using the preceding notation, the set in question is $AD^c(B \cup C)$, where $D = A_5$. Using distributive law, we can write this as $AD^cB \cup AD^cC$. We also have

$$(AD^cB)(AD^cC) = AD^cBC = ABC - ABCD.$$

Hence by (v),

$$P(AD^cBC) = P(ABC) - P(ABCD) = \frac{1}{12} - \frac{1}{60} = \frac{1}{15}.$$

Similarly, we have

$$P(AD^{c}B) = P(AB) - P(ABD) = \frac{1}{12} - \frac{1}{60} = \frac{4}{60} = \frac{1}{15},$$

$$P(AD^{c}C) = P(AC) - P(ACD) = \frac{1}{6} - \frac{1}{30} = \frac{4}{30} = \frac{2}{15}.$$

Finally we obtain by (viii):

$$P(AD^{c}B \cup AD^{c}C) = P(AD^{c}B) + P(AD^{c}C) - P(AD^{c}BC)$$
$$= \frac{1}{15} + \frac{2}{15} - \frac{1}{15} = \frac{2}{15}.$$

You should check this using the space Ω in Example 9.

The problem can be simplified by a little initial arithmetic, because the set in question is seen to be that of numbers divisible by 2 or 3 and not by 5. Now our method will yield the answer more quickly.

Exercises

1. Consider Example 1 in §2.1. Suppose that each good apple costs $1 \not\in$ while a rotten one costs nothing. Denote the rotten ones by R, an arbitrary bunch from the bushel by S, and define

$$Q(S) = |S \setminus R|/|\Omega - R|.$$

Q is the relative value of S with respect to that of the bushel. Show that it is a probability measure.

2. Suppose that the land of a square kingdom is divided into three strips A, B, C of equal area and suppose the value per unit is in the ratio of 1:3:2. For any piece of (measurable) land S in this kingdom, the relative value with respect to that of the kingdom is then given by the formula:

$$V(S) = \frac{P(SA) + 3P(SB) + 2P(SC)}{2}$$

where P is as in Example 2 of §2.1. Show that V is a probability measure.

*3. Generalizing No. 2, let a_1, \ldots, a_n be arbitrary positive numbers and let $A_1 + \cdots + A_n = \Omega$ be an arbitrary partition. Let P be a probability measure on Ω and

$$Q(S) = [a_1 P(SA_1) + \dots + a_n P(SA_n)]/[a_1 P(A_1) + \dots + a_n P(A_n)]$$

for any subset of Ω . Show that P is a probability measure.

- 4. Let A and B denote two cups of coffee you drank at a lunch counter. Suppose the first cup of coffee costs 15¢, and a second cup costs 10¢. Using P to denote "price," write down a formula like Axiom (ii) but with an inequality (P is "subadditive").
- 5. Suppose that on a shirt sale each customer can buy two shirts at \$4 each, but the regular price is \$5. A customer bought 4 shirts S_1, \ldots, S_4 . Write down a formula like Axiom (ii) and contrast it with Exercise 3. Forget about sales tax! (P is "superadditive.")
- 6. Show that if P and Q are two probability measures defined on the same (countable) sample space, then aP + bQ is also a probability measure for any two nonnegative numbers a and b satisfying a + b = 1. Give a concrete illustration of such a *mixture*.

- *7. If P is a probability measure, show that the function P/2 satisfies Axioms (i) and (ii) but not (iii). The function P^2 satisfies (i) and (iii) but not necessarily (ii); give a counterexample to (ii) by using Example 1.
- *8. If A, B, C are arbitrary sets, show that
 - (a) $P(A \cap B \cap C) \leq P(A) \wedge P(B) \wedge P(C)$,
 - (b) $P(A \cup B \cup C) \ge P(A) \lor P(B) \lor P(C)$.
- *9. Prove that for any two sets A and B, we have

$$P(AB) \ge P(A) + P(B) - 1.$$

Give a concrete example of this inequality. [Hint: Use (2.3.4) with n=2 and De Morgan's laws.]

- 10. We have $A \cap A = A$, but when is $P(A) \cdot P(A) = P(A)$? Can P(A) = 0 but $A \neq \emptyset$?
- 11. Find an example where P(AB) < P(A)P(B).
- 12. Prove (2.3.7) by first showing that

$$(A \cup B) - A = B - (A \cap B).$$

- 13. Two groups share some members. Suppose that Group A has 123, Group B has 78 members, and the total membership in both groups is 184. How many members belong to both?
- 14. Groups A, B, C have 57, 49, 43 members, respectively. A and B have 13, A and C have 7, B and C have 4 members in common; and there is a lone guy who belongs to all three groups. Find the total number of people in all three groups.
- *15. Generalize Exercise 14 when the various numbers are arbitrary but, of course, subject to certain obvious inequalities. The resulting formula, divided by the total population (there may be any nonjoiners!), is the extension of (2.3.7) to n=3.
 - 16. Compute $P(A \triangle B)$ in terms of P(A), P(B), and P(AB); also in terms of P(A), P(B), and $P(A \cup B)$.
- *17. Using the notation (2.5.3) and the probability defined in that context, show that for any two m and n we have

$$P(A_m A_n) \ge P(A_m) P(A_n).$$

When is there equality above?

*18. Recall the computation of plane areas by double integration in calculus; for a nice figure such as a parallelogram, trapezoid, or circle, we have

area of
$$S = \iint_S 1 \, dx \, dy$$
.

Show that this can be written in terms of the indicator I_S as

$$A(S) = \iint I_S(x, y) \, dx dy,$$

where Ω is the whole plane and $I_S(x,y)$ is the value of the function I_S for (at) the point (x,y) (denoted by ω in §1.4). Show also that for two such figures S_1 and S_2 , we have

$$A(S_1) + A(S_2) = \iint (I_{S_1} + I_{S_2}),$$

where we have omitted some unnecessary symbols.

- *19. Now you can demonstrate the connection between (2.3.7) and (2.3.8) mentioned there, in the case of plane areas.
- 20. Find several examples of $\{p_n\}$ satisfying the conditions in (2.4.1); give at least two in which all $p_n > 0$.
- *21. Deduce from Axiom (ii*) the following two results. (a) If the sets A_n are nondecreasing, namely $A_n \subset A_{n+1}$ for all $n \geq 1$, and $A_\infty = \bigcup_n A_n$, then $P(A_\infty) = \lim_{n \to \infty} P(A_n)$. (b) If the sets A_n are nonincreasing, namely $A_n \supset A_{n+1}$ for all $n \geq 1$, and $A_\infty = \bigcap_n A_n$, then $P(A_\infty) = \lim_{n \to \infty} P(A_n)$. [Hint: For (a), consider $A_1 + (A_2 A_1) + (A_3 A_2) + \cdots$; for (b), dualize by complementation.]
 - 22. What is the probability (in the sense of Example 10) that a natural number picked at random is not divisible by any of the numbers 3, 4, 6 but is divisible by 2 or 5?
- *23. Show that if (m_1, \ldots, m_n) are co-prime positive integers, then the events $(A_{m_1}, \ldots, A_{m_n})$ defined in §2.5 are independent.
 - 24. What can you say about the event A if it is independent of itself? If the events A and B are disjoint and independent, what can you say of them?
 - 25. Show that if the two events (A, B) are independent, then so are (A, B^c) , (A^c, B) , and (A^c, B^c) . Generalize this result to three independent events.
 - 26. Show that if A, B, C are independent events, then A and $B \cup C$ are independent, and $A \setminus B$ and C are independent.
 - 27. Prove that

$$P(A \cup B \cup C) = P(A) + P(B) + P(C)$$
$$-P(AB) - P(AC) - P(BC) + P(ABC)$$

when A, B, C are independent by considering $P(A^cB^cC^c)$. [The formula remains true without the assumption of independence; see §6.2.]

- 28. Suppose five coins are tossed; the outcomes are independent but the probability of heads may be different for different coins. Write the probability of the specific sequence *HHTHT* and the probability of exactly three heads.
- *29. How would you build a mathematical model for arbitrary repeated trials, namely without the constraint of independence? In other words, describe a sample space suitable for recording such trials. What is the mathematical definition of an event that is determined by one of the trials alone, two of them, etc.? You do not need a probability measure. Now think how you would cleverly construct such a measure over the space in order to make the trials independent. The answer is given in, e.g., [Feller 1, §V.4], but you will understand it better if you first give it a try yourself.

Counting

3.1. Fundamental rule

The calculation of probabilities often leads to the counting of various possible cases. This has been indicated in Examples 4 and 5 of §2.2 and forms the backbone of the classical theory with its stock in trade the games of chance. But combinatorial techniques are also needed in all kinds of applications arising from sampling, ranking, partitioning, allocating, programming, and model building, to mention a few. In this chapter we shall treat the most elementary and basic types of problems and the methods of solving them.

The author has sometimes begun a discussion of "permutations and combinations" by asking in class the following question. If a man has three shirts and two ties, in how many ways can he dress up [put on one of each]? Only two numbers, 2 and 3, are involved, and it's anybody's guess that one must combine them in some way. Does one add: 2+3? or multiply: 2×3 ? (or perhaps make 2^3 or 3^2). The question was meant to be rhetorical but experience revealed an alarming number of wrong answers. So if we dwell on this a little longer than you deem necessary you will know why.

First of all, in a simple example like that, one can simply picture the various possibilities and count them mentally:

A commonly used tabulation is as follows:

T	1	2	3
1	11	21	31
2	12	22	32

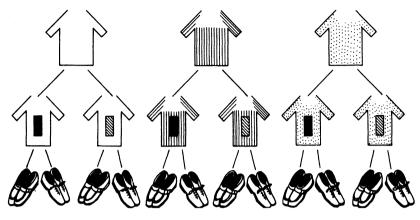


Figure 15

As mathematics is economy of thought we can schematize (program) this in a more concise way:

$$(s_1, t_1)(s_1, t_2)(s_2, t_1)(s_2, t_2)(s_3, t_1)(s_3, t_2),$$

and finally we see that it is enough just to write

$$(1,1)(1,2)(2,1)(2,2)(3,1)(3,2)$$
 (3.1.2)

by assigning the first slot to "shirt" and the second to "tie." Thus we have reached the mathematical method of naming the collection in (3.1.2). It is the set of all ordered couples (a, b) such that a = 1, 2, 3; b = 1, 2; and you see that the answer to my question is $3 \times 2 = 6$.

In general we can talk about ordered k-tuples (a_1, \ldots, a_k) where for each j from 1 to k, the symbol a_j indicates the assignment (choice) for the jth slot, and it may be denoted by a numeral between 1 and m_j . In the example above k = 2, $m_1 = 3$, $m_2 = 2$, and the collection of all (a_1, a_2) is what is enumerated in (3.1.2).

This symbolic way of doing things is extremely convenient. For instance, if the man has also two pairs of shoes, we simply extend each 2-tuple to a 3-tuple by adding a third slot into which we can put either "1" or "2." Thus each of the original 2-tuples in (3.1.2) splits into two 3-tuples, and so the total of 3-tuples will be $3 \times 2 \times 2 = 12$. This is the number of ways the man can choose a shirt, a tie, and a pair of shoes. You see it is all automated as on a computing machine. As a matter of fact, it is mathematical symbolism that taught the machines, not the other way around (at least, not yet).

The idea of splitting mentioned above lends well to visual imagination. It shows why 3 "shirts" *multiply* into 6 "shirt–ties" and 12 "shirt–tieshoes." Take a good look at it. Here is the general proposition:

Fundamental Rule. A number of multiple choices are to be made. There are m_1 possibilities for the first choice, m_2 for the second, m_3 for the third, etc. If these choices can be combined freely, then the total number of possibilities for the whole set of choices is equal to

$$m_1 \times m_2 \times m_3 \times \cdots$$
.

A formal proof would amount to repeating what is described above in more cut-and-dried terms, and is left to your own discretion. Let us point out, however, that "free combination" means in the example above that no matching of shirt and ties is required, etc.

Example 1. A menu in a restaurant reads like this:

Choice of one: Soup, Juice, Fruit Cocktail

Choice of one:
Beef Hash
Roast Ham
Fried Chicken
Spaghetti with Meatballs

Choice of one: Mashed Potatoes, Broccoli, Lima Beans

> Choice of one: Ice Cream, Apple Pie

> > Choice of one: Coffee, Tea, Milk

Suppose you take one of each "course" without substituting or skipping; how many options do you have? Or if you like the language nowadays employed in more momentous decisions of this sort, how many *scenarios* of a "complete 5-course dinner" (as advertised) can you make out of this menu? The total number of items you see on the menu is

$$3+4+3+2+3=15.$$

But you don't eat them all. On the other hand, the number of different dinners available is equal to

$$3 \times 4 \times 3 \times 2 \times 3 = 216$$
.

according to the fundamental rule. True, you eat only one dinner at a time, but it is quite possible for you to try all these 216 dinners if you have

catholic taste in food and patronize that restaurant often enough. More realistically and statistically significant: all these 216 dinners may be actually served to different customers over a period of time and perhaps even on a single day. This possibility forms the empirical basis of combinatorial counting and its relevance to computing probabilities.

Example 2. We can now solve the problem about Example (e) and (e') in §1.1: in how many ways can six dice appear when they are rolled? And in how many ways can they show all different faces?

Each die here represents a multiple choice of six possibilities. For the first problem these 6 choices can be freely combined so the rule applies directly to give the answer $6^6 = 46656$. For the second problem the choices cannot be freely combined since they are required to all be different. Off-hand the rule does not apply, but the reasoning behind it does. This is what counts in mathematics: not a blind reliance on a rule but a true understanding of its meaning. (Perhaps that is why "permutation and combination" are for many students harder stuff than algebra or calculus.) Look at the splitting diagram in Fig. 16.

The first die can show any face, but the second must show a different one. Hence, after the first choice has been made, there are five possibilities for the second choice. Which five depends on the first choice, but their number does not. So there are 6×5 possibilities for the first and second choices together. After these have been made, there are four possibilities left for the third, and so on. For the complete sequence of six choices we have, therefore, $6\cdot 5\cdot 4\cdot 3\cdot 2\cdot 1=6!=720$ possibilities. By the way, make sure by analyzing the diagram that the first die hasn't got preferential treatment. Besides, which is "first"?

Of course, we can re-enunciate a more general rule to cover the situation just discussed, but is it necessary once the principles are understood?

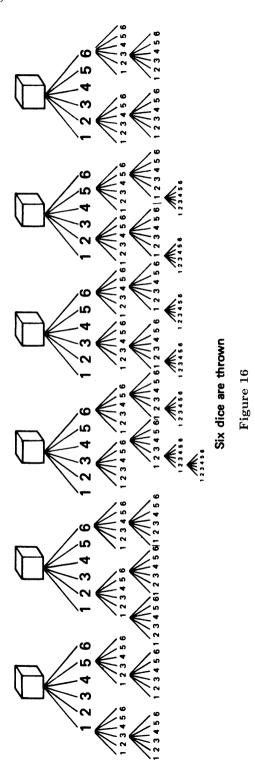
3.2. Diverse ways of sampling

Let us proceed to several standard methods of counting which constitute the essential elements in the majority of combinatorial problems. These can be conveniently studied either as sampling or as allocating problems. We begin with the former.

An urn contains m distinguishable balls marked 1 to m, from which n balls will be drawn under various specified conditions, and the number of all possible outcomes will be counted in each case.

I. Sampling with replacement and with ordering.

We draw n balls sequentially, each ball drawn being put back into the urn before the next drawing is made, and we record the numbers on the balls together with their order of appearance. Thus we are dealing with



ordered n-tuples (a_1, \ldots, a_n) in which each a_j can be any number from 1 to m. The fundamental rule applies directly and yields the answer m^n . This corresponds to the case of rolling six dice without restriction, but the analogy may be clearer if we think of the same dice being rolled six times in succession, so that each rolling corresponds to a drawing.

II. Sampling without replacement and with ordering.

We sample as in Case I, but after each ball is drawn it is left out of the urn. We are dealing with ordered n-tuples (a_1, \ldots, a_n) as above with the restriction that the a_j all be different. Clearly we must have $n \leq m$. The Fundamental Rule does not apply directly, but the splitting argument works as in Example 2 and yields the answer

$$m \cdot (m-1) \cdot (m-2) \cdots (m-n+1) = (m)_n.$$
 (3.2.1)

Observe that there are n factors on the left side of (3.2.1) and that the last factor is m - (n - 1) rather than m - n, why? We have introduced the symbol $(m)_n$ to denote this "continued product" on the left side of (3.2.1).

Case II has a very important subcase, which can be posed as a "permutation" problem.

IIa. Permutation of m distinguishable balls.

This is the case of II when m=n. Thus all m balls are drawn out one after another without being put back. The result is therefore just the m numbered balls appearing in some order, and the total number of such possibilities is the same as that of all possible arrangements (ordering, ranking, permutation) of the set $\{1, 2, \ldots, m\}$. This number is called the factorial of m and denoted by

$$m! = (m)_m = m(m-1) \cdots 2 \cdot 1$$

$$\begin{array}{c|cccc}
 & n! & & \\
\hline
 & 1 & & 1 \\
 & 2 & & 2 \\
 & 3 & & 6 \\
 & 4 & & 24 \\
 & 5 & & 120 \\
 & 6 & & 720 \\
 & 7 & & 5040 \\
 & 8 & & 40320 \\
 & 9 & & 362880 \\
\end{array}$$

III. Sampling without replacement and without ordering.

10

Here the balls are not put back and their ordering of appearance is not recorded; hence we might as well draw all n balls at one grab. We are therefore dealing with subsets of size n from a set (population) of size m.

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To count their number, we will compare with Case II where the balls are ranged in order. Now a bunch of n balls, if drawn one by one, can appear in n! different ways by Case IIa. Thus each unordered sample of size n produces n! ordered ones, and conversely every ordered sample of size n can be produced in this manner. For instance, if m=5, n=3, the subset $\{3,5,2\}$ can be drawn in 3!=6 ways as follows:

$$(2,3,5)(2,5,3)(3,2,5)(3,5,2)(5,2,3)(5,3,2).$$

In general we know from Case II that the total number of ordered samples of size n is $(m)_n$. Let us denote for one moment the unknown number of unordered samples of size n by x, then the argument above shows that

$$n! x = (m)_n$$
.

Solving for x, we get the desired answer, which will be denoted by

$$\binom{m}{n} = \frac{(m)_n}{n!}. (3.2.2)$$

If we multiply both numerator and denominator by (m-n)!, we see from (3.2.1) that

$${m \choose n} = \frac{(m)_n (m-n)!}{n! (m-n)!}$$

$$= \frac{m(m-1) \cdots (m-n+1)(m-n) \cdots 2 \cdot 1}{m! (m-n)!} = \frac{m!}{n! (m-n)!}.$$
(3.2.3)

When n=m, there is exactly one subset of size n, namely the whole set, hence the number in (3.2.3) must reduce to 1 if it is to maintain its significance. So we are obliged to set 0!=1. Under this convention, formula (3.2.3) holds for $0 \le n \le m$. The number $\binom{m}{n}$ is called a *binomial coefficient* and plays an important role in probability theory. Note that

$$\binom{m}{n} = \binom{m}{m-n},\tag{3.2.4}$$

which is immediate from (3.2.3). It is also obvious without this explicit evaluation from the interpretation of both sides as counting formulas (why?).

The argument used in Case III leads to a generalization of IIa:

IIIa. Permutation of m balls that are distinguishable by groups.

Suppose that there are m_1 balls of color no. 1, m_2 balls of color no. 2, ..., m_r balls of color no. r. Their colors are distinguishable, but balls of the same color are not. Of course, $m_1 + m_2 + \cdots + m_r = m$. How many distinguishable arrangements of these m balls are there?

For instance, if $m_1 = m_2 = 2$, m = 4, and the colors are black and white, there are 6 distinguishable arrangements as follows:

To answer the question in general, we compare with Case IIa where all balls are distinguishable. Suppose we mark the balls of color no. 1 from 1 to m_1 , the balls of color no. 2 from 1 to m_2 , and so forth. Then they become all distinguishable and so the total number of arrangements after the markings will be m! by Case IIa. Now the m_1 balls of color no. 1 can be arranged in $m_1!$ ways by their new marks, the m_2 balls of color no. 2 can be arranged in $m_2!$ ways by their new marks, etc. Each arrangement for one color can be freely combined with any arrangement for another color. Hence according to the fundamental rule, there are altogether

$$m_1! m_2! \dots m_r!$$

new arrangements produced by the various markings, for each original unmarked arrangement. It follows as in the discussion of Case III that the total number of distinguishable unmarked arrangements is equal to the quotient

$$\frac{m!}{m_1! \, m_2! \dots m_r!}.$$

This is called a multinomial coefficient. When r=2, it reduces to the binomial coefficient $\binom{m}{m_1} = \binom{m}{m_2}$.

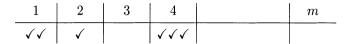
IV. Sampling with replacement and without ordering.

We draw n balls one after another, each ball being put back into the urn before the next drawing is made, but we record the numbers drawn with possible repetitions as a lot without paying attention to their order of appearance. This is a slightly more tricky situation, so we will begin by a numerical illustration. Take m=n=3; all the possibilities in this case are listed in the first column below:

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Do you see the organization principle used in making the list?

In general think of a "tally sheet" with numbers indicating the balls in the top line:



After each drawing we place a check under the number (of the ball) that is drawn. Thus at the end of all the drawings the total number of checks on the sheet will be n (which may be greater than m); there may be as many as that in an entry, and there may be blanks in some entries. Now economize by removing all dispensable parts of the tally sheet, so that column 1 in (3.2.5) becomes the skeleton in column 2. Check this over carefully to see that no information is lost in the simplified method of accounting. Finally, align the symbols \checkmark and | in column 2 to get column 3. Now forget about columns 1 and 2 and concentrate on column 3 for a while. Do you see how to reconstruct from each little cryptogram of "checks and bars" the original tally? Do you see that all possible ways of arranging 3 checks and 2 bars are listed there? Thus the total number is (by Case IIIa with m=5, $m_1=3$, $m_2=2$, or equivalently Case III with m=5, n=3) equal to $5!/3! \, 2! = 10$ as shown. This must therefore also be the number of all possible tally results.

In general each possible record of sampling under Case IV can be transformed by the same method into the problem of arranging n checks and m-1 bars (since m slots have m-1 lines dividing them) in all possible ways. You will have to draw some mental pictures to convince yourself that there is one-to-one correspondence between the two problems as in the particular case illustrated above. From IIIa we know the solution to the second problem is

$$\binom{m+n-1}{n} = \binom{m+n-1}{m-1}.$$
 (3.2.6)

Hence this is also the total number of outcomes when we sample under Case IV.

Example 3. D'Alembert's way of counting discussed in Example 4 of §2.2 is equivalent to sampling under Case IV with m=n=2. Tossing each coin corresponds to drawing a head or a tail if the results for two coins are tallied without regard to "which shows which" (or without ordering when the coins are tossed one after the other); then there are three possible outcomes:

$$\checkmark \checkmark | \qquad \checkmark | \checkmark \qquad | \checkmark \checkmark$$

HH HT = TH TT

Similarly, if six dice are rolled and the dice are not distinguishable, then the total number of recognizably distinct patterns is given by (3.2.6) with m = n = 6, namely

$$\binom{6+6-1}{6} = \binom{11}{6} = 462.$$

This is less than 1% of the number 46656 under Case 1.

We will now illustrate in a simple numerical case the different ways of counting in the four sampling procedures: m = 4, n = 2.

Case (I)				
	(1,1)	(1, 2)	(1,3)	(1,4)
	(2, 1)	(2, 2)	(2, 3)	(2, 4)
	(3, 1)	(3, 2)	(3, 3)	(3, 4)
ĺ	(4,1)	(4,2)	(4,3)	(4,4)

Case (II)			
	(1, 2)	(1,3)	(1,4)
(2,1)		(2, 3)	(2, 4)
(3,1)	(3, 2)		(3, 4)
(4,1)	(4, 2)	(4,3)	

			'	
	(1,1)	(1, 2)	(1,3)	(1,4)
1		(2, 2)	(2, 3)	(2, 4)
1			(3, 3)	(3, 4)
				(4,4)

Case (IV)

	(Case (III	.)
	(1, 2)	(1, 3)	(1, 4)
		(2, 3)	(2, 4)
ļ			(3,4)

3.3. Allocation models; binomial coefficients

A source of inspiration as well as frustration in combinatorics is that the same problem may appear in different guises and it may take an effort to recognize their true identity. Sampling under Case (IV) is a case in point; another will be discussed under (IIIb) ahead. People have different thinking habits and often prefer one way to another. But it may be worthwhile to learn some other ways as we learn foreign languages. In the above we have treated several basic counting methods as sampling problems. Another formulation, preferred by physicists and engineers, is "putting balls into boxes." Since these balls play a different role from those used above, we will call them tokens instead to simplify the translation later.

There are m boxes labeled from 1 to m and there are n tokens, numbered from 1 to n. The tokens are put into the boxes with or without the condition that no box can contain more than one token. We record the outcome of

the allocation (occupancy) by noting the number of tokens in each box, with or without noting the labels on the tokens. The four cases below then correspond respectively with the four cases of sampling discussed above.

- I'. Each box may contain any number of tokens, and the labels on the tokens are observed.
- II'. No box may contain more than one token, and the labels on the tokens are observed.
- III'. No box may contain more than one token, and the labels on the tokens are not observed.
- IV'. Each box may contain any number of tokens, and the labels on the tokens are not observed.

It will serve no useful purpose to prove that the corresponding problems are really identical, for this is the kind of mental exercise one must go through by oneself to be convinced. (Some teachers even go so far as to say that combinatorial thinking cannot be taught.) However, here are the key words in the translation from one to the other description.

$\operatorname{Sampling}$	Allocating	
Ball	Box	
Number of drawing	Number on token	
jth drawing gets ball no. k	jth token put into box no. k	

In some way the new formulation is more adaptable in that further conditions on the allocation can be imposed easily. For instance, one may require that no box be left empty when n > m, or specify "loads" in some boxes. Of course, these conditions can be translated into the other language, but they may then become less natural. Here is one important case of this sort, which is just Case IIIa in another guise.

IIIb. Partition into numbered groups.

Let a population of m objects be subdivided into r subpopulations or just "groups": m_1 into group no. 1, m_2 into group no. 2, ..., m_r into group no. r, where $m_1 + \cdots + m_r = m$ and all $m_i \ge 1$. This is a trivial paraphrasing of putting m tokens into r boxes so that m_i tokens are put into box no. j. It is important to observe that it is not the same as subdividing into r groups of sizes m_1, \ldots, m_r ; for the groups are numbered. A simple example will make this clear.

Example 4. In how many ways can four people split into two pairs?

The English language certainly does not make this question unambiguous, but offhand one would have to consider the following three ways as the answer:

$$(12)(34) \qquad (13)(24) \qquad (14)(23). \tag{3.3.1}$$

This is the correct interpretation if the two pairs are going to play chess or pingpong games and two equally good tables are available to both pairs. But now suppose the two pairs are going to play double tennis together and the "first" pair has the choice of side of the court, or will be the first to serve. It will then make a difference whether (12) precedes (34), or vice versa. So each case in (3.3.1) must be permuted by pairs into two orders, and the answer is then the following six ways:

$$(12)(34)$$
 $(34)(12)$ $(13)(24)$ $(24)(13)$ $(14)(23)$ $(23)(14)$.

This is the situation covered by the general problem of partition under IIIb, which will now be solved.

Think of putting tokens (people) into boxes (groups). According to sampling Case III, there are $\binom{m}{m_1}$ ways of choosing m_1 tokens to be put into box 1; after that, there are $\binom{m-m_1}{m_2}$ ways of choosing m_2 tokens from the remaining $m-m_1$ to be put into box 2; and so on. The fundamental rule does not apply but its modification used in sampling Case II does, and so the answer is

$${\binom{m}{m_1}} {\binom{m-m_1}{m_2}} {\binom{m-m_1-m_2}{m_3}} \cdots {\binom{m-m_1-m_2-\cdots-m_{r-1}}{m_r}}$$

$$= \frac{m!}{m_1! (m-m_1)!} \frac{(m-m_1)!}{m_2! (m-m_1-m_2)!} \frac{(m-m_1-m_2)!}{m_3! (m-m_1-m_2-m_3)!}$$

$$\cdots \frac{(m-m_1-\cdots-m_{r-1})!}{m_r! \ 0!}$$

$$= \frac{m!}{m_1! \ m_2! \cdots m_r!}.$$
(3.3.2)

Observe that there is no duplicate counting involved in the argument, even if some of the groups have the same size as in the tennis player example above. This is because we have given numbers to the boxes (groups). On the other hand, we are not arranging the numbered groups in order (as the words "ordered groups" employed by some authors would seem to imply). To clarify this essentially linguistic confusion, let us consider another simple example.

Example 5. Six mountain climbers decide to divide into three groups for the final assault on the peak. The groups will be of size 1, 2, 3, respectively, and all manners of deployment are considered. What is the total number of possible grouping and deploying?

The number of ways of splitting in G_1, G_2, G_3 , where the subscript denotes the size of group, is given by (3.3.2):

$$\frac{6!}{1!\,2!\,3!} = 60.$$

Having formed these three groups, there remains the decision of which group leads, which is in the middle, and which backs up. This is solved by Case IIa: 3! = 6. Now each grouping can be combined freely with each deploying; hence the fundamental rule gives the final answer: $60 \cdot 6 = 360$.

What happens when some of the groups have the same size? Think about the tennis players again.

Returning to (3.3.2), this is the same multinomial coefficient obtained as solution to the permutation problem IIIa. Here it appears as a combination type of problem since we have used sampling Case III repeatedly in its derivation. Thus it is futile to try to pin a label on a problem as permutation or combination. The majority of combinatorial problems involves a mixture of various ways of counting discussed above. We will now illustrate this with several worked problems.

In the remainder of this section we will establish some useful formulas connecting binomial coefficients. First of all, let us lay down the convention:

$$\binom{m}{n} = 0 \quad \text{if } m < n \text{ or if } n < 0.$$
 (3.3.3)

Next we show that

is what (3.3.4) says. Isn't this neat?

$$\binom{m}{n} = \binom{m-1}{n-1} + \binom{m-1}{n}, \qquad 0 \le n \le m. \tag{3.3.4}$$

Since we have the explicit evaluation of $\binom{m}{n}$ from (3.2.3), this can of course be verified at once. But here is a combinatorial argument without computation. Recall that $\binom{m}{n}$ is the number of different ways of choosing n objects out of m objects, which may be thought of as being done at one stroke. Now think of one of the objects as "special." This special one may or may not be included in the choice. If it is included, then the number of ways of choosing n-1 more objects out of the other m-1 objects is equal to $\binom{m-1}{n-1}$. If it is not included, then the number of ways of choosing all n objects from the other m-1 objects is equal to $\binom{m-1}{n}$. The sum of the two alternatives must then give the total number of choices, and this

As a consequence of (3.3.4), we can obtain $\binom{m}{n}$, $0 \le n \le m$, step by step as m increases, as follows:

For example, each number in the last row shown above is obtained by adding its two neighbors in the preceding row, where a vacancy may be regarded as zero:

$$1 = 0 + 1$$
, $7 = 1 + 6$, $21 = 6 + 15$, $35 = 15 + 20$, $35 = 20 + 15$, $21 = 15 + 6$, $7 = 6 + 1$, $1 = 1 + 0$.

Thus $\binom{7}{n} = \binom{6}{n-1} + \binom{6}{n}$ for $0 \le n \le 7$. The array in (3.3.5) is called *Pascal's triangle*, though apparently he was not the first one to have used it.

*Observe that we can split the last term $\binom{m-1}{n}$ in (3.3.4) as we split the first term $\binom{m}{n}$ by the same formula applied to m-1. Thus we obtain successively

$$\binom{m}{n} = \binom{m-1}{n-1} + \binom{m-1}{n} = \binom{m-1}{n-1} + \binom{m-2}{n-1} + \binom{m-2}{n}$$
$$= \binom{m-1}{n-1} + \binom{m-2}{n-1} + \binom{m-3}{n-1} + \binom{m-3}{n} = \cdots$$

The final result is

$$\binom{m}{n} = \binom{m-1}{n-1} + \binom{m-2}{n-1} + \dots + \binom{n}{n-1} + \binom{n}{n}$$

$$= \sum_{k=n-1}^{m-1} \binom{k}{n-1} = \sum_{k\leq m-1} \binom{k}{n-1}$$
(3.3.6)

^{*}The rest of the section may be omitted.

since the last term in the sum is $\binom{n}{n} = \binom{n-1}{n-1}$, and for k < n-1 the terms are zero by our convention (3.3.3).

Example. $35 = \binom{7}{4} = \binom{6}{3} + \binom{5}{3} + \binom{4}{3} + \binom{3}{3} = 20 + 10 + 4 + 1$. Look at Pascal's triangle to see where these numbers are located.

As an application, we can now give another solution to the counting problem for sampling under (IV) in §3.2. By the second formulation (IV') above, this is the number of ways of putting n indistinguishable [unnumbered] tokens into m labeled boxes without restriction. We know from (3.2.6) that it is equal to $\binom{m+n-1}{m-1}$, but the argument leading to this answer is pretty tricky. Suppose we were not smart enough to have figured it out that way, but have surmised the result by experimenting with small values of m and n. We can still establish the formula in general as follows. [Actually, that tricky argument was probably invented as an afterthought after the result had been surmised.]

We proceed by mathematical induction on the value of m. For m=1 clearly there is just one way of dumping all the tokens, no matter how many, into the one box, which checks with the formula since $\binom{1+n-1}{1-1} =$

 $\binom{n}{0}=1$. Now suppose that the formula holds true for any number of tokens when the number of boxes is equal to m-1. Introduce a new box; we may put any number of tokens into it. If we put j tokens into the new one, then we must put the remaining n-j tokens into the other m-1 boxes. According to the induction hypothesis, there are $\binom{m-2+n-j}{m-2}$ different ways of doing this. Summing over all possible values of j, we have

$$\sum_{j=0}^{n} {m-2+n-j \choose m-2} = \sum_{k=m-2}^{m+n-2} {k \choose m-2},$$

where we have changed the index of summation by setting m-2+n-j=k. The second sum above is equal to $\binom{m+n-1}{m-1}$ by (3.3.6), and the induction is complete.

Next, let us show that

$$\binom{n}{0} + \binom{n}{1} + \binom{n}{2} + \dots + \binom{n}{n} = \sum_{k=0}^{n} \binom{n}{k} = 2^{n};$$
 (3.3.7)

that is, the sum of the nth row in Pascal's triangle is equal to 2^n [the first row shown in (3.3.5) is the 0th]. If you know Newton's binomial theorem,

this can be shown from

$$(a+b)^n = \sum_{k=0}^n \binom{n}{k} a^k b^{n-k}$$
 (3.3.8)

by substituting a=b=1. But here is a combinatorial proof. The terms on the left side of (3.3.7) represent the various numbers of ways of choosing $0,1,2,\ldots,n$ objects out of n objects. Hence the sum is the total number of ways of choosing any subset [the empty set and the entire set both included] from a set of size n. Now in such a choice each object may or may not be included, and the inclusion or exclusion of each object may be freely combined with that of any other. Hence the fundamental rule yields the total number of choices as

$$\underbrace{2 \times 2 \times \cdots \times 2}_{n \text{ times}} = 2^n.$$

This is the number given on the right side of (3.3.7). It is the total number of distinct subsets of a set of size n.

Example. For n = 2, all choices from (a, b) are

$$\emptyset$$
, $\{a\}$, $\{b\}$, $\{ab\}$.

For n = 3, all choices from (a, b, c) are

$$\emptyset, \{a\}, \{b\}, \{c\}, \{ab\}, \{ac\}, \{bc\}, \{abc\}.$$

Finally, let $k \leq m$ be two positive integers. We show

$$\binom{m}{n} = \sum_{j=0}^{k} \binom{k}{j} \binom{m-k}{n-j}.$$
 (3.3.9)

Observe how the indices on top [at bottom] on the right side add up to the index on top [at bottom] on the left side; it is not necessary to indicate the precise range of j in the summation; we may let j range over all integers because the superfluous terms will automatically be zero by our convention (3.3.3). To see the truth of (3.3.9), we think of the m objects as being separated into two piles, one containing k objects and the other m-k. To choose n objects from the entire set, we may choose j objects from the first pile and n-j objects from the second pile, and combine them. By the fundamental rule, for each fixed value j the number of such combinations is equal to $\binom{k}{j}\binom{m-k}{n-j}$. So if we allow j to take all possible values and add up the results, we obtain the total number of choices which is equal

to $\binom{m}{n}$. You need not worry about "impossible" values for j when j > n or n - j > m - k, because the corresponding term will be zero by our convention.

Example.

$$\binom{7}{3} = \binom{3}{0} \binom{4}{3} + \binom{3}{1} \binom{4}{2} + \binom{3}{2} \binom{4}{1} + \binom{3}{3} \binom{4}{0},$$

$$\binom{7}{5} = \binom{3}{1} \binom{4}{4} + \binom{3}{2} \binom{4}{3} + \binom{3}{3} \binom{4}{2}.$$

In particular, if k = 1 in (3.3.9), we are back to (3.3.4). In this case our argument also reduces to the one used there.

An algebraic derivation of (3.3.9), together with its extension to the case where the upper indices are no longer positive integers, will be given in Chapter 6.

3.4. How to solve it

This section may be entitled "How to count." Many students find these problems hard, partly because they have been inured in other elementary mathematics courses to the cookbook variety of problems such as: "Solve $x^2 = 5x + 10 = 0$," "differentiate xe^{-x} " (maybe twice), etc. One can do such problems by memorizing certain rules without any independent thought. Of course, we have this kind of problem in "permutation and combination" too, and you will find some of these among the exercises. For instance, there is a famous formula to do the "round-table" problem: "In how many different ways can 8 people be seated at a round table?" If you learned it, you could solve the problem without knowing what the word "different" means. But a little variation may get you into deep trouble. The truth is, and that's also a truism: there is no substitute for true understanding. However, it is not easy to understand the principles without concrete applications, so the handful of examples below are selected to be the "test cases." More are given in the Exercises and you should have a lot of practice if you want to become an expert. Before we discuss the examples in detail, a few general tips will be offered to help you to do your own thing. They are necessarily very broad and rather slippery, but they may be of some help sometimes.

(a) If you don't *see* the problem well, try some particular (but not too particular) case with small numbers so you can see better. This will fix in your mind what is to be counted, and help you especially in spotting duplicates and omissions.

- (b) Break up the problem into pieces provided that they are simpler, cleaner, and easier to concentrate on. This can be done sometimes by fixing one of the "variables," and the number of similar pieces may be counted as a subproblem.
- (c) Don't try to argue step by step if you can see complications rising rapidly. Of all the negative advice I gave my classes this was the least heeded but probably the most rewarding. Counting step by step may seem easy for the first couple of steps, but do you see how to carry it through to the end?
- (d) Don't be turned off if there is ambiguity in the statement of the problem. This is a semantical hang-up, not a mathematical one. Try all interpretations if necessary. This may not be the best strategy in a quiz, but it's a fine thing to do if you want to learn the stuff. In any case, don't take advantage of the ambiguities of the English language or the oversight of your instructor to turn a reasonable problem into a trivial one. (See Exercise 13.)

Problem 1. (Quality Control). Suppose that in a bushel of 550 apples 2% are rotten ones. What is the probability that a "random sample" of 25 apples contains 2 rotten apples?

This is the principle behind testing the quality of products by random checking. If the probability turns out to be too small on the basis of the claimed percentage compared with that figured on some other suspected percentage, then the claim is in doubt. This problem can be done just as easily with arbitrary numbers so we will formulate it in the general case. Suppose there are k defective items in a lot of m products. What is the probability that a random sample of size n contains j defective items? The word "random" here signifies that all samples of size n, under Case III in §3.2, are considered equally likely. Hence the total number is $\binom{m}{n}$. How many of these contain exactly j defective items? To get such a sample we must choose any j out of the k defective items and combine it freely with k0 out of the k1 nondefective items. The first choice can be made in $\binom{k}{j}$ ways, the second in $\binom{m-k}{n-j}$ ways, by sampling under Case III. By the fundamental rule, the total number of samples of size k2 containing k3 defective items is equal to the product, and consequently the desired probability is the ratio

$$\binom{k}{j} \binom{m-k}{n-j} / \binom{m}{n}. \tag{3.4.1}$$

In the case of the apples, we have m = 550, k = 11, n = 25, j = 2; so the

probability is equal to

$$\binom{11}{2} \binom{539}{23} \bigg/ \binom{550}{25}.$$

This number is not easy to compute, but we will learn how to get a good approximation later. Numerical tables are also available.

If we sum the probabilities in (3.4.1) for all j, $0 \le j \le n$, the result ought to equal 1 since all possibilities are counted. We have therefore *proved* the formula

$$\sum_{j=0}^{k} \binom{k}{j} \binom{m-k}{n-j} = \binom{m}{n}$$

by a probabilistic argument. This is confirmed by (3.3.9); indeed a little reflection should convince you that the two arguments are really equivalent.

Problem 2. If a deck of poker cards is thoroughly shuffled, what is the probability that the four aces are found in a row?

There are 52 cards among which are 4 aces. A thorough shuffling signifies that all permutations of the cards are equally likely. For the whole deck, there are (52)! outcomes by Case IIa. In how many of these do the four aces stick together? Here we use tip (b) to break up the problem according to where the aces are found. Since they are supposed to appear in a row, we need only locate the first ace as we check the cards in the order they appear in the deck. This may be the top card, the next, and so on, until the 49th. Hence there are 49 positions for the 4 aces. After this has been fixed, the 4 aces can still permute among themselves in 4! ways, and so can the 48 nonaces. This may be regarded as a case of IIIa with r=2, $m_1=4$, $m_2=48$. The fundamental rule carries the day, and we get the answer

$$\frac{49 \cdot 4! (48)!}{(52)!} = \frac{24}{52 \cdot 51 \cdot 50}.$$

This problem is a case where my tip (a) may be helpful. Try four cards with two aces. The total number of permutations in which the aces stick together is only $3 \cdot 2! \cdot 2! = 12$, so you can list them all and look.

Problem 3. Fifteen new students are to be evenly distributed among three classes. Suppose that there are 3 whiz-kids among the 15. What is the probability that each class gets one? One class gets them all?

It should be clear that this is the partition problem discussed under Case IIIb, with m=15, $m_1=m_2=m_3=5$. Hence the total number of outcomes is given by

$$\frac{15!}{5! \, 5! \, 5!}$$

To count the number of these assignments in which each class gets one whiz-kid, we will first assign these three kids. This can be done in 3! ways by IIa. The other 12 students can be evenly distributed in the 3 classes by Case IIIb with m = 12, $m_1 = m_2 = m_3 = 4$. The fundamental rule applies, and we get the desired probability

$$3! \frac{12!}{4! \cdot 4! \cdot 4!} / \frac{15!}{5! \cdot 5! \cdot 5!} = \frac{6 \cdot 5^3}{15 \cdot 14 \cdot 13}.$$

Next, if one class gets them all, then there are three possibilities according to which class it is, and the rest is similar. So we just replace the numerator above by $3 \cdot (12!/5! \, 5! \, 2!)$ and obtain

$$3 \cdot \frac{12!}{5! \cdot 5! \cdot 2!} / \frac{15!}{5! \cdot 5! \cdot 5!} = \frac{5 \cdot 4 \cdot 3^2}{15 \cdot 14 \cdot 13}.$$

By the way, we can now get the probability of the remaining possibility, namely that the number of whiz-kids in the three classes be two, one, zero respectively.

Problem 4. Six dice are rolled. What is the probability of getting three pairs?

One can ask at once "which three pairs?" This means a choice of 3 numbers out of the 6 numbers from 1 to 6. The answer is given by sampling under Case III: $\binom{6}{3} = 20$. Now we can concentrate on one of these cases, say $\{2,3,5\}$, and figure out the probability of getting "a pair of 2, a pair of 3, and a pair of 5." This is surely more clear-cut, so my tip (b) should be used here. To count the number of ways 6 dice can show the pattern $\{2,2,3,3,5,5\}$, one way is to consider this as putting six labeled tokens (the dice as distinguishable) into three boxes marked $\boxed{2}$, $\boxed{3}$, $\boxed{5}$, with two going into each box. So the number is given by IIIb:

$$\frac{6!}{2! \, 2! \, 2!} = 90.$$

Another way is to think of the dice as six distinguishable performers standing in line waiting for cues to do their routine acts, with two each doing acts nos. 2, 3, 5, respectively, but who does which is up to Boss Chance. This then becomes a permutation problem under IIIa and gives of course the same number above. Finally, we multiply this by the number of choices of the 3 numbers to get

$$\binom{6}{3} \frac{6!}{2! \ 2! \ 2!} = 20 \cdot 90.$$

You may regard this multiplication as another application of the ubiquitous fundamental rule, but it really just means that 20 mutually exclusive categories are *added* together, each containing 90 cases. The desired probability is given by

$$\frac{20 \cdot 90}{6^6} = \frac{25}{648}.$$

This problem is a case where my negative tip (c) may save you some wasted time as I have seen students trying an argument as follows. If we want to end up with three pairs, the first two dice can be anything; the third die must be one of the two if they are different, and a new one if they are the same. The probability of the first two being different is 5/6, in which case the third die has probability 4/6; on the other hand, the probability of the first two being the same is 1/6, in which case the third has probability 5/6. Are you still with us? But what about the next step, and the next?

However, this kind of sequential analysis, based on conditional probabilities, will be discussed in Chapter 5. It works very well sometimes, as in the next problem.

Problem 5. (Birthdays). What is the probability that among n people there are at least two who have the same birthday? We are assuming that they "choose" their birthdays independently of one another, so that the result is as if they had drawn n balls marked from 1 to 365 (ignoring leap years) by sampling under Case I. All these outcomes are equally likely, and the total number is $(365)^n$. Now we must count those cases in which some of the balls drawn bear the same number. This sounds complicated but it is easy to figure out the "opposite event," namely when all n balls are different. This falls under Case II, and the number is $(365)_n$. Hence the desired probability is

$$p_n = 1 - \frac{(365)_n}{(365)^n}.$$

What comes as a surprise is the numerical fact that this probability exceeds 1/2 as soon as $n \ge 23$; see table below.* What would you have guessed?

^{*}Computation from n=2 to n=55 was done on a small calculator to five decimal places in a matter of minutes.

n	p_n
5	.03
10	.12
15	.25
20	.41
21	.44
22	.48
23	.51
24	.54
25	.57
30	.71
35	.81
40	.89
45	.94
50	.97
55	.99

One can do this problem by a naive argument that turns out to be correct. To get the probability that n people all have different birthdays, we order them in some way and consider each one's "choice," as in the case of the six dice that show different faces (Example 2, §3.1). The first person can have any day of the year for her birthday, hence probability 1; the second can have any but one, hence probability 364/365; the third any but two, hence probability 363/365; and so on. Thus the final probability is

$$\frac{365}{365} \frac{364}{365} \frac{363}{365} \cdots$$
 (*n* factors),

which is just another way of writing $(365)_n/(365)^n$. The intuitive idea of sequential conditional probabilities used here is equivalent to a splitting diagram described in Section 3.1, beginning with 365 cases, each of which splits into 364, then again into 363, etc. If one divides out by 365 at each stage, one gets the product above.

Problem 6. (Matching). Four cards numbered 1 to 4 are laid face down on a table and a person claiming clairvoyance will name them by his extrasensory power. If he is a faker and just guesses at random, what is the probability that he gets at least one right?

There is a neat solution to this famous problem by a formula to be established later in §6.2. But for a small number like 4, brute force will do and in the process we shall learn something new. Now the faker simply picks any one of the 4! permutations, and these are considered equally likely. Using tip (b), we will count the number of cases in which there is exactly 1 match. This means the other three cards are mismatched, and so

we must count the "no-match" cases for three cards. This can be done by enumerating all the 3! = 6 possible random guesses as tabulated below:

There are two cases of no-match: the 4th and 5th above. We obtain all cases in which there is exactly one match in four cards by fixing that one match and mismatch the three other cards. There are four choices for the card to be matched, and after this is chosen, there are two ways to mismatch the other three by the tabulation above. Hence by the modified fundamental rule there are $4 \cdot 2 = 8$ cases of exactly 1 match in 4 cards. Next, fix two matches and mismatch the other two. There is only one way to do the latter; hence the number of cases of exactly 2 matches in 4 cards is equal to that of choosing 2 cards (to be matched) out of 4, which is $\binom{4}{2} = 6$.

Finally, it is clear that if three cards match, the remaining one must also, and there is just one way of matching them all. The results are tabulated as follows:

Exact number of matches	Number of cases	Probability Probability
4	1	1/24
3	0	0
2	6	1/4
1	8	1/3
0	9	3/8

The last row above, for the number of cases of no-match, is obtained by subtracting the sum of the other cases from the total number:

$$24 - (1 + 6 + 8) = 9.$$

The probability of at least one match is 15/24 = 5/8; of at least two matches is 7/24.

You might propose to do the counting without any reasoning by listing all 24 cases for 4 cards, as we did for 3 cards. That is a fine thing to do, not only for your satisfaction but also to check the various cases against our reasoning above. But our step leading from 3 cards to 4 cards is meant to be an illustration of the empirical inductive method and can lead also from 4 to 5, etc. In fact, that is the way the computing machines do things. They are really not very smart, and always do things step by step, but they are organized and tremendously fast. In our case a little neat algebra does it better, and we can establish the following general formula for the number

of cases of at least one match for n cards:

$$n! \left(1 - \frac{1}{2!} + \frac{1}{3!} - + \dots + (-1)^{n-1} \frac{1}{n!}\right).$$

Problem 7. In how many ways can n balls be put into n numbered boxes so that exactly one box is empty? This problem is instructive as it illustrates several points made above. First of all, it is ambiguous whether the balls are distinguishable or not. Using my tip (d), we will treat both hypotheses.

Hypothesis 1. The balls are indistinguishable. Then it is clearly just a matter of picking the empty box and the one that must have two balls. This is a sampling problem under Case II, and the answer is $(n)_2 = n(n-1)$.

This easy solution would probably be acceptable granted the ambiguous wording, but we can learn more if we try the harder way too.

Hypothesis 2. The balls are distinguishable. Then after the choice of the two boxes as under Hypothesis 1 (call it step 1), we still have the problem as to which ball goes into which box. This is a problem of partition under Case IIIb with m = n, $m_1 = 2$, $m_2 = \cdots = m_{n-1} = 1$, the empty box being left out of consideration. Hence the answer is

$$\frac{n!}{2! \, 1! \cdots 1!} = \frac{n!}{2!}.\tag{3.4.2}$$

You don't have to know about that formula, since you can argue directly as follows. The question is how to put n numbered balls into n-1 numbered boxes with 2 balls going into a certain box (already chosen by step 1) and 1 ball each into all the rest. There are $\binom{n}{2}$ ways of choosing the two balls to go into that particular box, after that the remaining n-2 balls can go into the other n-2 boxes in (n-2)! ways. The product of these two numbers is the same as (3.4.2). Finally, the total number of ways under Hypothesis 2 is given by

$$n(n-1) \cdot \frac{n!}{2}.\tag{3.4.3}$$

We have argued in two steps above. One may be tempted to argue in three steps as follows. First choose the empty box; then choose n-1 balls and put them one each into the other n-1 boxes; finally throw the last ball into any one of the latter. The number of possible choices for each step is equal, respectively, to n, $(n)_{n-1}$ (by sampling under II), and n-1. If they are multiplied together, the result is

$$n \cdot n! \, (n-1), \tag{3.4.4}$$

which is twice as large as (3.4.3). Which is correct?

70 Counting

This is the kind of situation my tip (a) is meant to help. Take n=3 and suppose the empty box has been chosen, so the problem is to put balls 1, 2, 3 into boxes A and B. For the purpose of the illustration let A be square and B round. Choose two round balls to put into these two boxes; there are six cases as shown:

Now throw the last ball into one of the two boxes, so that each case above splits into two according to which box gets it:

13 2	23	$\boxed{12}\boxed{3}$	$\boxed{32} \boxed{1}$	21	31 2
1 23	2 (13)	1 32	3 (12)	2 (31)	3 (21)

You see what the trouble is. Each final case is counted twice because the box that gets two balls can get them in two orders! The trouble is the same in the general case, and so we must divide the number in formula (3.4.4) by 2 to eliminate double counting, which makes it come out as in formula (3.4.3). All is harmonious.

Exercises

[When probabilities are involved in the problems below, the equally likely cases should be "obvious" from the context. In case you demur, follow my tip (d).]

- 1. A girl decides to choose either a shirt or a tie for a birthday present. There are three shirts and two ties to choose from. How many choices does she have if she will get only one of them? If she may get both a shirt and a tie?
- 2. Three kinds of shirts are on sale. (a) If two men buy one shirt each, how many possibilities are there? (b) If two shirts are sold, how many possibilities are there?
- 3. As in No. 2 make up a good question with 3 shirts and 2 men, to which the answer is 2^3 , or $\binom{3+2-1}{3}$.
- 4. If on the menu shown in §3.1 there are three kinds of ice cream and two kinds of pie to choose from, how many different dinners are there? If we take into account that the customer may skip the vegetable or the dessert or both, how many different dinners are there?
- 5. How many different initials can be formed with two or three letters of the alphabet? How large must the alphabet be in order that 1 million people can be identified by 3-letter initials?
- 6. How many integers are there between 1 million and 10 million, in whose decimal form no two consecutive digits are the same?

- 7. In a "true or false" test there are 12 questions. If a student decides to check six of each at random, in how many ways can she do it?
- 8. In how many ways can four boys and four girls pair off? In how many ways can they stand in a row in alternating gender?
- 9. In how many ways can a committee of 3 be chosen from 20 people? In how many ways can a president, a secretary, and a treasurer be chosen?
- 10. If you have two dollars, two quarters, and three nickels, how many different sums can you pay without making change? Change the quarters into dimes and answer again.
- 11. Two screws are missing from a machine that has screws of three different sizes. If three screws of different sizes are sent over, what is the probability that they are what's needed?
- 12. There are two locks on the door and the keys are among the six different ones you carry in your pocket. In a hurry you dropped one somewhere. What is the probability that you can still open the door? What is the probability that the first two keys you try will open the door?
- 13. A die is rolled three times. What is the probability that you get a larger number each time? (I gave this simple problem in a test but inadvertently used the words "... that the numbers you obtain increase steadily." Think of a possible *mis* interpretation of the words!)
- *14. Three dice are rolled twice. What is the probability that they show the same numbers (a) if the dice are distinguishable, (b) if they are not. [Hint: divide into cases according to the pattern of the first throw: a pair, a triple, or all different; then match the second throw accordingly.]
 - 15. You walk into a party without knowing anyone there. There are six women and four men and you know there are four married couples. In how many ways can you guess who the couples are? What if you know there are exactly three couples?
 - 16. Four shoes are taken at random from five different pairs. What is the probability that there is at least one pair among them?
- 17. A California driver decides that he must switch lanes every minute to get ahead. If he is on a 4-lane divided highway and does this at random, what is the probability that he is back on his original lane after 4 minutes (assuming no collision)? [Hint: the answer depends on whether he starts in an outside or inside lane.]
- 18. In sampling under Case I or II of §3.2, what is the probability that in n drawings a particular ball is never drawn? Assume n < m.
- 19. You are told that of the four cards face down on the table, two are red and two are black. If you guess all four at random, what is the probability that you get 0, 2, 4 right?
- 20. An airport shuttle bus makes 4 scheduled stops for 15 passengers. What is the probability that all of them get off at the same stop? What is the probability that someone (at least one person) gets off at each stop?

- 21. Ten books are made into 2 piles. In how many ways can this be done if books as well as piles may or may not be distinguishable? Treat all four hypotheses, and require that neither pile be empty.
- 22. Ten different books are to be given to Daniel, Phillip, Paul, and John, who will get in the order given 3, 3, 2, 2 books, respectively. In how many ways can this be done? Since Paul and John screamed "no fair," it is decided that they draw lots to determine which two get 3 and which two get 2. How many ways are there now for a distribution? Finally, Marilda and Corinna also want a chance and so it is decided that the six kids should draw lots to determine which two get 3, which two get 2, and which two get none. Now how many ways are there? [There is real semantical difficulty in formulating these distinct problems in general. It is better to be verbose than concise in such a situation. Try putting tokens into boxes.]
- 23. In a draft lottery containing the 366 days of the year (including February 29), what is the probability that the first 180 days drawn (without replacement of course) are evenly distributed among the 12 months? What is the probability that the first 30 days drawn contain none from August or September? [Hint: first choose 15 days from each month.]
- 24. At a certain resort the travel bureau finds that tourists occupy the 20 hotels there as if they were so many lookalike tokens (fares) placed in numbered boxes. If this theory is correct, what is the probability that when the first batch of 30 tourists arrive, no hotel is left vacant? [This model is called the Bose-Einstein statistic in physics. If the tourists are treated as distinct persons, it is the older Boltzmann-Maxwell statistic; see [Feller 1, §II.5].]
- 25. One hundred trout are caught in a little lake and returned after they are tagged. Later another 100 are caught and found to contain 7 tagged ones. What is the probability of this if the lake contains n trout? [What is your best guess as to the true value of n? The latter is the kind of question asked in statistics.]
- 26. Program a one-to-one correspondence between the various possible cases under the two counting methods in IIIa and IIIb by taking m = 4, $m_1 = 2$, $m_2 = m_3 = 1$.
- *27. (For poker players only.) In a poker hand assume all "hands" are equally likely as under Sampling Case III. Compute the probability of (a) flush, (b) straight, (c) straight flush, (d) four of a kind, (e) full house.
- *28. Show that

$$\binom{2n}{n} = \sum_{k=0}^{\infty} \binom{n}{k}^2.$$

[Hint: apply (3.3.9).]

*29. The number of different ways in which a positive integer n can be written as the sum of positive integers not exceeding n is called (in number theory) the "partition number" of n. For example,

6 = 6	sextuple
= 5 + 1	quintuple
= 4 + 2	quadruple and pair
=4+1+1	quadruple
= 3 + 3	two triples
= 3 + 2 + 1	triple and pair
= 3 + 1 + 1 + 1	triple
= 2 + 2 + 2	three pairs
= 2 + 2 + 1 + 1	two pairs
=2+1+1+1+1	one pair
= 1 + 1 + 1 + 1 + 1 + 1	all different ("no same")

Thus the partition number of 6 is equal to 11; compare this with the numbers 46656 and 462 given in Examples 2 and 3. This may be called the total number of distinguishable "coincidence patterns" when six dice are thrown. We have indicated simpler (but vaguer) names for these patterns in the listing above. Compute their respective probabilities. [It came to me as a surprise that "two pairs" is more probable than "one pair" and has a probability exceeding 1/3. My suspicion of an error in the computation was allayed only after I had rolled 6 dice 100 times. It was an old custom in China to play this game over the New Year holidays, and so far as I can remember, "two pairs" were given a higher rank (prize) than "one pair." This is unfair according to their probabilities. Subsequently to my own experiment I found that Feller* had listed analogous probabilities for seven dice. His choice of this "random number" 7 in disregard or ignorance of a time-honored game had probably resulted in my overlooking his tabulation.]

*30. (Banach's match problem.) The Polish mathematician Banach kept two match boxes, one in each pocket. Each box contains n matches. Whenever he wanted a match he reached at random into one of his pockets. When he found that the box he picked was empty, what is the distribution of the number of matches left in the other box? [Hint: divide into two cases according to whether the left or right box is empty, but be careful about the case when both are empty.]

^{*}William Feller (1906–70), leading exponent of probability.

Random Variables

4.1. What is a random variable?

We have seen that the points of a sample space may be very concrete objects such as apples, molecules, and people. As such they possess various qualities, some of which may be measurable. An apple has its weight and volume; its juicy content can be scientifically measured; even its taste may be graded by expert tasters. A molecule has mass and velocity, from which we can compute its momentum and kinetic energy by formulas from physics. For a human being, there are physiological characteristics such as age, height, and weight. But there are many other numerical data attached to him (or her) like I.Q., number of years of schooling, number of brothers and sisters, annual income earned and taxes paid, and so on. We will examine some of these illustrations and then set up a mathematical description in general terms.

Example 1. Let Ω be a human population containing n individuals. These may be labeled as

$$\Omega = \{\omega_1, \omega_2, \dots, \omega_n\}. \tag{4.1.1}$$

If we are interested in their age distribution, let $A(\omega)$ denote the age of ω . Thus to each ω is associated a number $A(\omega)$ in some unit, such as "year." So the mapping

$$\omega \to A(\omega)$$

is a function with Ω as its domain of definition. The range is the set of integers but can be made more precise by fractions or decimals or spelled out as, e.g., "18 years, 5 months, and 1 day." There is no harm if we take all positive integers or all positive real numbers as the range, although only a very small portion of it will be needed. Accordingly, we say A is an integer-valued or real-valued function. Similarly, we may denote the height, weight, and income by the functions

$$\omega \to H(\omega),$$
 $\omega \to W(\omega),$
 $\omega \to I(\omega).$

In the last case I may take negative values! Now for some medical purposes, a linear combination of height and weight may be an appropriate measure:

$$\omega \to \lambda H(\omega) + \mu W(\omega)$$
,

where λ and μ are two numbers. This then is also a function of ω . Similarly, if ω is a "head of family," alias breadearner, the census bureau may want to compute the function:

$$\omega o rac{I(\omega)}{N(\omega)}$$

where $N(\omega)$ is the number of persons in his or her family, namely the number of mouths to be fed. The ratio above represents then the "income per capita" for the family.

Let us introduce some convenient symbolism to denote various sets of sample points derived from random variables. For example, the set of ω in Ω for which the age is between 20 and 40 will be denoted by

$$\{\omega \mid 20 \le A(\omega) \le 40\}$$

or more briefly when there is no danger of misunderstanding by

$$\{20 \leq A \leq 40\}.$$

The set of ω for which the height is between 65 and 75 (in inches) and the weight is between 120 and 180 (in pounds) can be denoted in several ways as follows:

$$\begin{split} \{\omega \mid 65 \leq H(\omega) \leq 75\} \bigcap \{\omega \mid 120 \leq W(\omega) \leq 180\} \\ &= \{\omega \mid 65 \leq H(\omega) \leq 75; 120 \leq W(\omega) \leq 180\} \\ &= \{65 \leq H \leq 75; 120 \leq W \leq 180\}. \end{split}$$

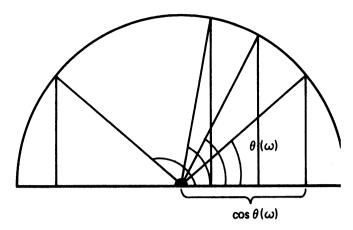


Figure 17

Example 2. Let Ω be gaseous molecules in a given container. We can still represent Ω as in (4.1.1) even though n is now a very large number such as 10^{25} . Let m= mass, v= velocity, M= momentum, E= kinetic energy. Then we have the corresponding functions:

$$\begin{split} \omega &\to m(\omega), \\ \omega &\to v(\omega), \\ \omega &\to M(\omega) = m(\omega)v(\omega), \\ \omega &\to E(\omega) = \frac{1}{2}m(\omega)v(\omega)^2. \end{split}$$

In experiments with gases actual measurements may be made of m and v, but the quantities of interest may be M or E, which can be derived from the formulas. Similarly, if θ is the angle of the velocity relative to the x-axis, $\omega \to \theta(\omega)$ is a function of ω and

$$\omega \to \cos \theta(\omega)$$

may be regarded as the composition of the function "cos" with the function " θ ." The set of all molecules moving toward the right is represented by

$$\{\omega \mid \cos \theta(\omega) > 0\}.$$

Example 3. Let Ω be the outcome space of throwing a die twice. Then it consists of $6^2 = 36$ points listed below:

(1,1)	(1, 2)	(1,3)	(1,4)	(1,5)	(1,6)
(2,1)	(2, 2)	(2, 3)	(2,4)	(2, 5)	(2, 6)
(3,1)	(3, 2)	(3, 3)	(3, 4)	(3, 5)	(3, 6)
(4,1)	(4,2)	(4,3)	(4,4)	(4, 5)	(4, 6)
(5,1)	(5, 2)	(5, 3)	(5, 4)	(5, 5)	(5, 6)
(6,1)	(6, 2)	(6, 3)	(6, 4)	(6, 5)	(6, 6)

Thus each ω is represented by an ordered pair or a two-dimensional *vector*:

$$\omega_k = (x_k, y_k), \quad k = 1, 2, \dots, 36,$$

where x_k and y_k take values from 1 to 6. The first *coordinate* x represents the outcome of the first throw, the second coordinate y the outcome of the second throw. These two coordinates are determined by the point ω , hence they are functions of ω :

$$\omega \to x(\omega), \quad \omega \to y(\omega).$$
 (4.1.2)

On the other hand, each ω is completely determined by its two coordinates, so much so that we may say that ω is the pair of them:

$$\omega \equiv (x(\omega), y(\omega)).$$

This turnabout is an important concept to grasp. For example, let the die be thrown n times and the results of the successive throws be denoted by

$$x_1(\omega), x_2(\omega), \ldots, x_n(\omega);$$

then not only is each $x_k(\omega), k = 1, 2, \dots, n$, a function of ω that may be called its *kth coordinate*, but the totality of these n functions in turn determines ω , and therefore ω is nothing more or less than the n-dimensional vector

$$\omega \equiv (x_1(\omega), x_2(\omega), \dots, x_n(\omega)). \tag{4.1.3}$$

In general each $x_k(\omega)$ represents a certain numerical characteristic of the sample ω , and although ω may possess many, many characteristics, in most questions only a certain set of them is taken into account. Then a representation like (4.1.3) is appropriate. For example, in a traditional beauty contest, only three bodily measurements given in inches are considered, such as (36, 29, 38). In such a contest (no "song and dance") each contestant is reduced to such an ordered triple:

contestant =
$$(x, y, z)$$
.

Another case of this kind is when a student takes a number of tests, say 4, which are graded on the usual percentage basis. Let the student be ω , his

score on the 4 tests be $x_1(\omega), x_2(\omega), x_3(\omega), x_4(\omega)$. For the grader (or the computing machine if all the tests can be machine processed), each ω is just the 4 numbers $(x_1(\omega), \ldots, x_4(\omega))$. Two students who have the same scores are not distinguished. Suppose the criterion for success is that the total should exceed 200; then the set of successful candidates is represented by

$$\{\omega \mid x_1(\omega) + x_2(\omega) + x_3(\omega) + x_4(\omega) > 200\}.$$

A variation is obtained if different weights $\lambda_1, \lambda_2, \lambda_3, \lambda_4$ are assigned to the 4 tests – then the criterion will depend on the linear combination $\lambda_1 x_1(\omega) + \cdots + \lambda_4 x_4(\omega)$. Another possible criterion for passing the tests is given by

$$\{\omega \mid \min(x_1(\omega), x_2(\omega), x_3(\omega), x_4(\omega)) > 35\}.$$

What does this mean in plain English?

4.2. How do random variables come about?

We can now give a general formulation for numerical characteristics of sample points. Assume first that Ω is a countable space. This assumption makes an essential simplification that will become apparent; other spaces will be discussed later.

Definition of Random Variable. A numerically valued function X of ω with domain Ω :

$$\omega \in \Omega: \quad \omega \to X(\omega)$$
 (4.2.1)

is called a random variable [on Ω].

The term "random variable" is well established and so we will use it in this book, but "chance variable" or "stochastic variable" would have been good too. The adjective "random" is just to remind us that we are dealing with a sample space and trying to describe certain things that are commonly called "random events" or "chance phenomena." What might be said to have an element of randomness in $X(\omega)$ is the sample point ω that is picked "at random," such as in a throw of dice or the polling of an individual from a population. Once ω is picked, $X(\omega)$ is thereby determined and there is nothing vague, indeterminate, or chancy about it anymore. For instance, after an apple ω is picked from a bushel, its weight $W(\omega)$ can be measured and may be considered as known. In this connection the term "variable" should also be understood in the broad sense as a "dependent variable," namely a function of ω , as discussed in §4.1. We can say that the sample point ω serves here as an "independent variable" in the same way the variable x in $\sin x$ does, but it is better not to use this language

since "independent" has a very different and more important meaning in probability theory (see §5.5).

Finally, it is a custom (not always observed) to use a capital letter to denote a random variable, such as X, Y, N, or S, but there is no reason why we cannot use small letters x or y as we did in the examples of $\S 4.1$.

Observe that random variables can be defined on a sample space before any probability is mentioned. Later we shall see that they acquire their probability distributions through a probability measure imposed on the space.

Starting with some random variables, we can at once make new ones by operating on them in various ways. Specific examples have already been given in §4.1. The general proposition may be stated as follows:

Proposition 1. If X and Y are random variables, then so are

$$X + Y$$
, $X - Y$, XY , X/Y $(Y \neq 0)$, $(4.2.2)$

and aX + bY where a and b are two numbers.

This is immediate from the general definition, since, e.g.,

$$\omega \to X(\omega) + Y(\omega)$$

is a function on Ω as well as X and Y. The situation is exactly the same as in calculus: if f and g are functions, then so are

$$f+g$$
, $f-g$, fg , f/g $(g \neq 0)$, $af+bg$.

The only difference is that in calculus these are functions of x, a real number, while here the functions in (4.2.2) are those of ω , a sample point. Also, as in calculus where a constant is regarded as a very special kind of function, so is a constant a very special kind of random variable. For example, it is quite possible that in a class in elementary school, all pupils are of the same age. Then the random variable $A(\omega)$ discussed in Example 1 of §4.1 is equal to a constant, say = 9 (years) in a fourth-grade class.

In calculus a function of a function is still a function such as $x \to \log(\sin x)$ or $x \to f(\varphi(x)) = (f \circ \varphi)(x)$. A function of a random variable is still a random variable such as the $\cos \theta$ in Example 2 of §4.1. More generally we can have a function of several random variables.

Proposition 2. If φ is a function of two (ordinary) variables and X and Y are random variables, then

$$\omega \to \varphi(X(\omega), Y(\omega))$$
 (4.2.3)

is also a random variable, which is denoted more concisely as $\varphi(X,Y)$.

A good example is the function $\varphi(x,y)=\sqrt{x^2+y^2}$. Suppose $X(\omega)$ and $Y(\omega)$ denote, respectively, the horizontal and vertical velocities of a gas molecule; then

$$\varphi(X,Y) = \sqrt{x^2 + Y^2}$$

will denote its absolute speed.

Let us note in passing that Proposition 2 contains Proposition 1 as a particular case. For instance, if we take $\varphi(x,y) = x+y$, then $\varphi(X,Y) = X+Y$. It also contains functions of a single random variable as a particular case such as f(X). Do you see why? Finally, an extension of Proposition 2 to more than two variables is obvious. A particularly important case is the sum of n random variables:

$$S_n(\omega) = X_1(\omega) + \dots + X_n(\omega) = \sum_{k=1}^n X_k(\omega). \tag{4.2.4}$$

For example, if X_1, \ldots, X_n denote the successive outcomes of a throw of a die, then S_n is the total obtained in n throws. We shall have much to do with these partial sums S_n .

We will now illustrate the uses of random variables in some everyday situations. Quite often the intuitive notion of some random quantity precedes that of a sample space. Indeed one can often talk about random variables X, Y, etc. without bothering to specify Ω . The rather formal (and formidable?) mathematical setup serves as a necessary logical backdrop, but it need not be dragged into the open on every occasion when the language of probability can be readily employed.

Example 4. The cost of manufacturing a certain book is \$3 per book up to 1000 copies, \$2 per copy between 1000 and 5000 copies, and \$1 per copy afterwards. In reality, of course, books are printed in round lots and not on demand "as you go." What we assume here is tantamount to selling all overstock at cost, with no loss of business due to understock. Suppose we print 1000 copies initially and price the book at \$5. What is "random" here is the number of copies that will be sold; call it X. It should be evident that once X is known, we can compute the profit or loss from the sales; call this Y. Thus Y is a function of X and is random only because X is so. The formula connecting Y with X is given below (see Fig. 18 on page 82):

$$Y = \begin{cases} 5X - 3000 & \text{if } X \le 1000, \\ 2000 + 3(X - 1000) & \text{if } 1000 < X \le 5000, \\ 14000 + 4(X - 5000) & \text{if } X > 5000. \end{cases}$$

What is the probability that the book is a financial loss? It is that of the

event represented by the set

$${5X - 3000 < 0} = {X < 600}.$$

What is the probability that the profit will be at least \$10000? It is that of the set

$$\{2000 + 3(X - 1000) \ge 10000\} \cup \{X > 5000\}$$
$$= \left\{X \ge \frac{8000}{3} + 1000\right\} \cup \{X > 5000\}$$
$$= \{X > 3667\}.$$

But what are these probabilities? They will depend on a knowledge of X. One can only guess at it in advance; so it is a random phenomenon. But after the sales are out, we shall know the exact value of X; just as after a die is cast we shall know the outcome. The various probabilities are called the distribution of X and will be discussed in $\S 4.3$ ahead.

What is the sample space here? Since the object of primary interest is X, we may very well take it as our sample point and call it ω instead to conform with our general notation. Then each ω is some positive integer and $\omega \to Y(\omega)$ is a random variable with Ω the space of positive integers. To pick an ω means in this case to hazard a guess (or make a hypothesis) on the number of sales, from which we can compute the profit by the preceding formula. There is nothing wrong about this model of a sample space, though it seems a bit superfluous.

A more instructive way of thinking is to consider each ω as representing "a possible sales record for the book." A publisher is sometimes interested in other information than the total number of sales. An important factor left out of consideration above is the time element involved in the sale. Surely it makes a difference whether 5000 copies are sold in 1 or 10 years. If the book is a college text like this one, it may be important to know how it does in different types of schools and in different regions of the country. If it is fiction or drama, it may mean a great deal (even only from the profit motive) to know what the critics say about it, though this would be in a promotions rather than sales record. All these things may be contained in a capsule, which is the sample point ω . You can imagine it to be a complete record of every bit of information pertaining to the book, of which $X(\omega)$ and $Y(\omega)$ represent only two facets. Then what is Ω ? It is the totality of all such conceivable records. This concept of a sample space may sound weird and is unwieldy (can we say that Ω is countable?), but it gives the appropriate picture when one speaks of, e.g., the path of a particle in Brownian motion or the evolution of a stochastic process (see Chapter 8). On the other hand, it also shows the expediency of working with some specific random variables rather than worrying about the whole universe.

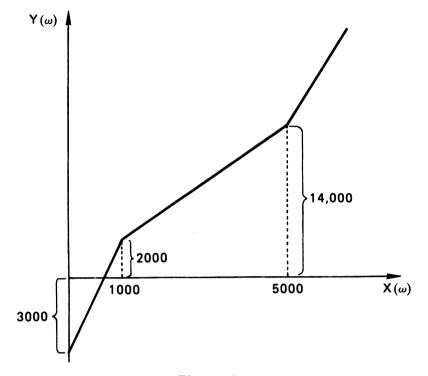


Figure 18

Example 5. An insurance company receives claims for indemnification from time to time. Both the times of arrival of such claims and their amounts are unknown in advance and determined by chance; ergo, random. The total amount of claims in one year, say, is of course also random, but clearly it will be determined as soon as we know the "when" and "how much" of the claims. Let the claims be numbered as they arrive and let S_n denote the date of the nth claim. Thus $S_3 = 33$ means the third claim arrives on February 2. So we have

$$1 < S_1 < S_2 < \cdots < S_n < \cdots$$

and there is equality whenever several claims arrive on the same day. Let the amount of the nth claim be C_n (in dollars). What is the total number of claims received in the year? It is given by N, where

$$N = \max\{n \mid S_n \le 365\}.$$

Obviously N is also random but it is determined by the sequence of S_n 's; in theory we need to know the entire sequence because N may be arbitrarily large. Knowing N and the sequence of C_n 's, we can determine the total

amount of claims in that year:

$$C_1 + \dots + C_N \tag{4.2.5}$$

in the notation of (4.2.4). Observe that in (4.2.5) not only is each term a random variable but also the number of terms. Of course the sum is also a random variable. It depends on the S_n 's as well as the C_n 's.

In this case we can easily imagine that the claims arrive at the office one after another and a complete record of them is kept in a ledger printed like a diary. Under some dates there may be no entry, under others there may be many in various different amounts. Such a ledger is kept over the years and will look quite different from one period of time to another. Another insurance company will have another ledger that may be similar in some respects and different in others. Each conceivable account kept in such a ledger may be considered as a sample point, and a reasonably large collection of them may serve as the sample space. For instance, an account in which one million claims arrive on the same day may be left out of the question, or a claim in the amount of 95 cents. In this way we can keep the image of a sample space within proper bounds of realism.

If we take such a view, other random variables come easily to mind. For example, we may denote by Y_k the total amount of claims on the kth day. This will be the number that is the sum of all the entries under the date, possibly zero. The total claims from the first day of the account to the nth day can then be represented by the sum

$$Z_n = \sum_{k=1}^n Y_k = Y_1 + Y_2 + \dots + Y_n. \tag{4.2.6}$$

The total claims over any period of time [s,t] can then be represented as

$$Z_t - Z_{s-1} = \sum_{k=s}^t Y_k = Y_s + Y_{s+1} + \dots + Y_t.$$
 (4.2.7)

We can plot the *accumulative* amount of claims Z_t against the time t by a graph of the kind in Figure 19.

There is a jump at t when the entry for the tth day is not empty, and the size of the rise is the total amount of claims on that day. Thus the successive rises correspond to the Y_k 's that are greater than 0. Clearly you can read off from such a graph the total claim over any given period of time, and also, e.g., the lengths of the "free periods" between the claims, but you cannot tell what the individual claims are when several arrive on the same day. If all the information you need can be gotten from such a graph, then you may regard each conceivable graph as a sample point. This will yield a somewhat narrower sample space than the one described above, but it will serve our purpose. From the mathematical point of view, the

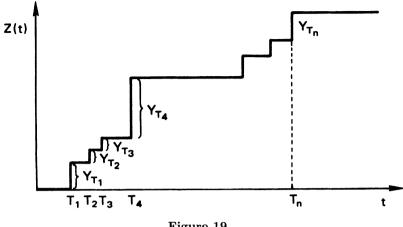


Figure 19

identification of a sample point with a graph (also called a sample curve, path, or trajectory) is very convenient, since a curve is a more precise (and familiar!) object than a ledger or some kind of sales-and-promotions record.

4.3. Distribution and expectation

In Chapter 2 we discussed the probabilities of sets of sample points. These sets are usually determined by the values of random variables. A typical example is

$$\{a \leq X \leq b\} = \{\omega \mid a \leq X(\omega) \leq b\}, \tag{4.3.1}$$

where X is a random variable, a and b are two constants. Particular cases of this have been indicated among the examples in §4.1. Since every subset of Ω has a probability assigned to it when Ω is countable, the set above has a probability, which will be denoted by

$$P(a \le X \le b). \tag{4.3.2}$$

More generally let A be a set of real numbers, alias a set of points on the real line $R^1 = (-\infty, +\infty)$; then we can write

$$P(X \in A) = P(\{\omega \mid X(\omega) \in A\}). \tag{4.3.3}$$

For instance, if A is the closed interval [a, b], then this is just the set in (4.3.2); but A may be the open interval (a, b), half-open interval (a, b] or [a,b), infinite intervals $(-\infty,b)$ or $(a,+\infty)$; the union of several intervals, or a set of integers say $\{m, m+1, \ldots, m+n\}$. An important case occurs when A reduces to a single point x; it is then called the singleton $\{x\}$.

The distinction between the point x and the set $\{x\}$ may seem academic. Anyway, the probability

$$P(X = x) = P(X \in \{x\}) \tag{4.3.4}$$

is "the probability that X takes (or assumes) the value x." If X is the age of a human population, $\{X=18\}$ is the subpopulation of 18-year-olds—a very important set!

Now the hypothesis that Ω is countable will play an essential simplifying role. Since X has Ω as domain of definition, it is clear that the range of X must be finite when Ω is finite, and at most countably infinite when Ω is so. Indeed, the *exact range* of X is just the set of real numbers below:

$$V_X = \bigcup_{\omega \in \Omega} \{X(\omega)\},\tag{4.3.5}$$

and many of these numbers may be the same – because the mapping $\omega \to X(\omega)$ is in general many-to-one, not necessarily one-to-one. In the extreme case when X is a constant random variable, the set V_X reduces to a single number. Let the distinct values in V_X be listed in any order as

$$\{v_1, v_2, \ldots, v_n, \ldots\}.$$

The sequence may be finite or infinite. Clearly if $x \notin V_X$, namely if x is not one of the values v_n , then P(X=x)=0. On the other hand, we do not forbid that some v_n may have zero probability. This means that some sample points may have probability zero. You may object: why don't we throw such nuisance points out of the sample space? Because it is often hard to know in advance which ones to throw out. It is easier to leave them in since they do no harm. [In an uncountable Ω , every single point ω may have probability zero! But we are not talking about this at present; see §4.5 ahead.]

Let us introduce the notation

$$p_n = P(X = v_n), \quad v_n \in V_x.$$
 (4.3.6)

It should be obvious that if we know all the p_n 's, then we can calculate all probabilities concerning the random variable X, alone. Thus the probabilities in (4.3.2) and (4.3.3) can be expressed in terms of the p_n 's as follows:

$$P(a \le X \le b) = \sum_{a \le v_n \le b} p_n; \quad P(X \in A) = \sum_{v_n \in A} p_n. \tag{4.3.7}$$

The first is a particular case of the second, and the last-written sum reads this way: "the sum of the p_n 's for which the corresponding v_n 's belong to A."

When A is the infinite interval $(-\infty, x]$ for any real number x, we can introduce a function of x as follows:

$$F_X(x) = P(X \le x) = \sum_{v_n \le x} p_n.$$
 (4.3.8)

This function $x \to F_X(x)$ defined on R^1 is called the distribution function of X. Its value at x "picks up" all the probabilities of values of X up to x (inclusive); for this reason the adjective "cumulative" is sometimes added to its name. For example, if X is the annual income (in \$'s) of a breadwinner, then $F_X(10000)$ is the probability of the income group earning anywhere up to \$10,000 and can theoretically include all those whose incomes are negative.

The distribution function F_X is determined by the v_n 's and p_n 's as shown in (4.3.8). Conversely, if we know F_X , namely we know $F_X(x)$ for all x, we can "recover" the v_n 's and p_n 's. We will not prove this fairly obvious assertion here. For the sake of convenience, we shall say that the two sets of numbers $\{v_n\}$ and $\{p_n\}$ determine the probability distribution of X, where any v_n for which $p_n = 0$ may be omitted. It is easy to see that if a < b, then

$$P(a < X \le b) = P(X \le b) - P(X \le a) = F_X(b) - F_X(a); \tag{4.3.9}$$

but how do we get $P(a \le X \le b)$, or P(X = x) from F_X ? (See Exercise 7 ahead.)

Now to return to the p_n 's, which are sometimes called the "elementary probabilities" for the random variable X. In general they have the following two properties:

(i)
$$\forall n$$
: $p_n \ge 0$;
(ii) $\sum_{n=0}^{\infty} p_n = 1$. (4.3.10)

Compare this with (2.4.1). The sum in (ii) may be over a finite or infinite sequence depending on whether V_X is a finite or infinite set. Property (i) is obvious, apart from the observation already made that some p_n may = 0. Property (ii) says that the values $\{v_n\}$ in V_X exhaust all possibilities for X, hence their probabilities must add up to that of the "whole universe." This is a fine way to say things, but let us learn to be more formal by converting the verbal argument into a symbolic proof. We begin with

$$\bigcup_{n} \{X = v_n\} = \Omega.$$

Since the v_n 's are distinct, the sets $\{X = v_n\}$ must be disjoint. Hence by

countable additivity (see §2.3) we have

$$\sum_{n} P(X = v_n) = P(\Omega) = 1.$$

This is Property (ii).

Before making further specialization on the random variables, let us formulate a fundamental new definition in its full generality. It is motivated by the intuitive notion of the average of a random quantity.

Definition of Mathematical Expectation. For a random variable X defined on a countable sample space Ω , its mathematical expectation is the number E(X) given by the formula

$$E(X) = \sum_{\omega \in \Omega} X(\omega) P(\{\omega\}), \tag{4.3.11}$$

provided that the series converges absolutely, namely

$$\sum_{\omega \in \Omega} |X(\omega)| P(\{\omega\}) < \infty. \tag{4.3.12}$$

In this case we say that the mathematical expectation of X exists. The process of "taking expectations" may be described in words as follows: take the value of X at each ω , multiply it by the probability of that point, and sum over all ω in Ω . If we think of $P(\{\omega\})$ as the weight attached to ω , then E(X) is the weighted average of the function X. Note that if we label the ω 's as $\{\omega_1, \omega_2, \ldots, \omega_n, \ldots\}$, then we have

$$E(X) = \sum_{n} X(\omega_n) P(\{\omega_n\}).$$

But we may as well use ω itself as label, and save a subscript, which explains the cryptic notation in (4.3.11).

Example 6. Let $\Omega = \{\omega_1, \dots, \omega_7\}$ be a parcel of land subdivided into seven "lots for sale." These lots have percentage areas and prices as follows:

Define $X(\omega)$ to be the price per acre of the lot ω . Then E(X) is the average price per acre of the whole parcel and is given by

$$(800)\frac{5}{100} + (900)\frac{10}{100} + (1000)\frac{10}{100} + (1200)\frac{10}{100} + (800)\frac{15}{100} + (900)\frac{20}{100} + (800)\frac{30}{100} = 890;$$

namely \$890 per acre. This can also be computed by first lumping together all acreage at the same price, and then summing over the various prices;

$$(800) \left(\frac{5}{100} + \frac{15}{100} + \frac{30}{100}\right) + (900) \left(\frac{10}{100} + \frac{20}{100}\right) + (1000) \frac{10}{100} + (1200) \frac{10}{100}$$
$$= (800) \frac{50}{100} + (900) \frac{30}{100} + (1000) \frac{10}{100} + (1200) \frac{10}{100} = 890.$$

The adjective in "mathematical expectation" is frequently omitted, and it is also variously known as "expected value," "mean (value)," or "first moment" (see §6.3 for the last term). In any case, do not expect the value E(X) when X is observed. For example, if you toss a fair coin to win \$1 or nothing depending on whether it falls heads or tails, you will never get the expected value \$.50! However, if you do this n times and n is large, then you can expect to get about n/2 dollars with a good probability. This is the implication of the law of large numbers, to be made precise in §7.6.

We shall now amplify the condition given in (4.3.12). Of course, it is automatically satisfied when Ω is a finite space, but it is essential when Ω is countably infinite because it allows us to calculate the expectation in any old way by rearranging and regrouping the terms in the series in (4.3.11), without fear of getting contradictory results. In other words, if the series is absolutely convergent, then it has a uniquely defined "sum" that in no way depends on how the terms are picked out and added together. The fact that contradictions can indeed arise if this condition is dropped may be a surprise to you. If so, you will do well to review your knowledge of the convergence and absolute convergence of a numerical series. This is a part of the calculus course which is often poorly learned (and taught), but will be essential for probability theory, not only in this connection but generally speaking. Can you, for instance, think of an example where the series in (4.3.11) converges but the one in (4.3.12) does not? [Remember that the p_n 's must satisfy the conditions in (4.3.10), though the v_n 's are quite arbitrary. So the question is a little harder than just to find an arbitrary nonabsolutely convergent series; but see Exercise 21. In such a case the expectation is *not* defined at all. The reason why we are being so strict is: absolutely convergent series can be manipulated in ways that nonabsolutely [conditionally] convergent series cannot be. Surely the definition of E(X) would not make sense if its value could be altered simply by shuffling around the various terms in the series in (4.3.11), which merely means that we enumerate the sample points in a different way. Yet this can happen without the condition (4.3.12)!

Let us explicitly state a general method of calculating E(X) which is often expedient. Suppose the sample space Ω can be decomposed into disjoint sets A_n :

$$\Omega = \bigcup_{n} A_n \tag{4.3.13}$$

in such a way that X takes the same value on each A_n . Thus we may write

$$X(\omega) = a_n \quad \text{for} \quad \omega \in A_n,$$
 (4.3.14)

where the a_n 's need not all be different. We then have

$$E(X) = \sum_{n} P(A_n)a_n = \sum_{n} P(X = a_n)a_n.$$
 (4.3.15)

This is obtained by regrouping the ω 's in (4.3.11) first into the subsets A_n , and then summing over all n. In particular, if $(v_1, v_2, \ldots, v_n, \ldots)$ is the range of X, and we group the sample points ω according to the values of $X(\omega)$, i.e., putting

$$A_n = \{ \omega \mid X(\omega) = v_n \}, \quad P(A_n) = p_n,$$

then we get

$$E(X) = \sum_{n} p_n v_n, (4.3.16)$$

where the series will automatically converge absolutely because of (4.3.12). In this form it is clear that the expectation of X is determined by its probability distribution.

Finally, it is worthwhile to point out that formula (4.3.11) contains an expression for the expectation of any function of X:

$$E(\varphi(X)) = \sum_{\omega \in \Omega} \varphi(X(\omega)) P(\{\omega\})$$

with a proviso like (4.3.12). For by Proposition 2 or rather a simpler analogue, $\varphi(X)$ is also a random variable. It follows that we have

$$E(\varphi(X)) = \sum_{n} p_n \varphi(v_n), \qquad (4.3.17)$$

where the v_n 's are as in (4.3.16), but note that the $\varphi(v_n)$'s need not be distinct. Thus the expectation of $\varphi(X)$ is already determined by the probability distribution of X (and of course also by the function φ), without the intervention of the probability distribution of $\varphi(X)$ itself. This is most convenient in calculations. In particular, for $\varphi(x) = x^r$ we get the rth moment of X:

$$E(X^r) = \sum_{n} p_n v_n^r; (4.3.18)$$

4.4. Integer-valued random variables

In this section we consider random variables that take only nonnegative integer values. In this case it is convenient to consider the range to be the entire set of such numbers:

$$N^0 = \{0, 1, 2, \dots, n, \dots\}$$

since we can assign probability zero to those that are not needed. Thus we have, as specialization of (4.3.6), (4.3.8), and (4.3.11),

$$p_n = P(X = n), n \in \mathbb{N}^0,$$

$$F_X(x) = \sum_{0 \le n \le x} p_n,$$

$$E(X) = \sum_{n=0}^{\infty} n p_n.$$
(4.4.1)

Since all terms in the last-written series are nonnegative, there is no difference between convergence and absolute convergence. Furthermore, since such a series either converges to a finite sum or diverges to $+\infty$, we may even allow $E(X) = +\infty$ in the latter case. This is in contrast to our general definition in the last section but is a convenient extension.

In many problems there is a practical justification to consider the random variables to take only integer values, provided a suitably small unit of measurement is chosen. For example, monetary values can be expressed in cents rather than dollars, or one tenth of a cent if need be; if "inch" is not a small enough unit for lengths we can use one hundredth or one thousandth of an inch. There is a unit called angstrom (Å), which is equal to 10^{-7} of a millimeter, used to measure electromagnetic wavelengths. For practical purposes, of course, incommensurable magnitudes (irrational ratios) do not exist; at one time π was legally defined to be 3.14 in some state of the United States! But one can go too far in this kind of justification!

We proceed to give some examples of (4.4.1).

Example 7. Suppose L is a positive integer, and

$$p_n = \frac{1}{L}, \quad 1 \le n \le L.$$
 (4.4.2)

Then automatically all other p_n 's must be zero because $\sum_{n=1}^{L} p_n = L \cdot \frac{1}{L} = 1$ and the conditions in (4.3.10) must be satisfied. Next, we have

$$E(X) = \frac{1}{L} \sum_{n=1}^{L} n = \frac{1}{L} \cdot \frac{L(L+1)}{2} = \frac{L+1}{2}.$$

The preceding sum is done by a formula for *arithmetical progression* which you have probably learned in school.

We say in this case that X has a uniform distribution over the set $\{1,2,\ldots,L\}$. In the language of Chapter 3, the L possible cases $\{X=1\}, \{X=2\},\ldots, \{X=L\}$ are all equally likely. The expected value of X is equal to the arithmetical mean [average] of the L possible values. Here is an illustration of its meaning. Suppose you draw at random a token X from a box containing 100 tokens valued at $1 \not\in$ to $100 \not\in$. Then your expected prize is given by $E(X) = 50.5 \not\in$. Does this sound reasonable to you?

Example 8. Suppose you toss a perfect coin repeatedly until a head turns up. Let X denote the number of tosses it takes until this happens, so that $\{X = n\}$ means n - 1 tails before the first head. It follows from the discussion in Example 8 of §2.4 that

$$p_n = P(T=n) = \frac{1}{2^n} \tag{4.4.3}$$

because the favorable outcome is just the specific sequence $TT \cdots TH$.

What is the expectation of X? According to (4.4.1), it is given by the formula

$$\sum_{n=1}^{\infty} \frac{n}{2^n} = ? \tag{4.4.4}$$

Let us learn how to sum this series, though properly speaking this does not belong to this course. We begin with the *fountainhead* of many of such series:

$$\frac{1}{1-x} = 1 + x + x^2 + \dots + x^n + \dots = \sum_{n=0}^{\infty} x^n \quad \text{for} \quad |x| < 1.$$
 (4.4.5)

This is a geometric series of the simplest kind which you surely have seen. Now differentiate it term by term:

$$\frac{1}{(1-x)^2} = 1 + 2x + 3x^2 + \dots + nx^{n-1} + \dots = \sum_{n=0}^{\infty} (n+1)x^n \quad \text{for } |x| < 1.$$
(4.4.6)

This is valid because the radius of convergence of the power series in (4.4.5) is equal to 1, so such manipulations are legitimate for |x| < 1. [Absolute and uniform convergence of the power series are involved here.] If we substitute

x = 1/2 in (4.4.6), we obtain

$$4 = \sum_{n=0}^{\infty} (n+1) \left(\frac{1}{2}\right)^n. \tag{4.4.7}$$

There is still some difference between (4.4.4) and the preceding series, so a little algebraic manipulation is needed. One way is to split up the terms above:

$$4 = \sum_{n=0}^{\infty} \frac{n}{2^n} + \sum_{n=0}^{\infty} \frac{1}{2^n} = \sum_{n=1}^{\infty} \frac{n}{2^n} + 2,$$

where we have summed the second series by substituting x = 1/2 into (4.4.5). Thus the answer to (4.4.4) is equal to 2. Another way to manipulate the formula is to change the index of summation: $n + 1 = \nu$. Then we have

$$4 = \sum_{n=0}^{\infty} \frac{n+1}{2^n} = \sum_{\nu=1}^{\infty} \frac{\nu}{2^{\nu-1}} = 2 \sum_{\nu=1}^{\infty} \frac{\nu}{2^{\nu}},$$

which of course yields the same answer. Both techniques are very useful!

The expectation E(X) = 2 seems eminently fair on intuitive grounds. For if the probability of your obtaining a head is 1/2 on one toss, then two tosses should get you $2 \cdot 1/2 = 1$ head, on the average. This plausible argument [which was actually given in a test paper by a smart student] can be made rigorous, but the necessary reasoning involved is far more sophisticated than you might think. It is a case of Wald's equation* or martingale theorem [for the advanced reader].

Let us at once generalize this problem to the case of a biased coin, with probability p for head and q = 1 - p tail. Then (4.4.3) becomes

$$p_n = \underbrace{(q \cdots q)}_{n-1 \text{ times}} p = q^{n-1} p, \tag{4.4.8}$$

and (4.4.4) becomes

$$\sum_{n=1}^{\infty} nq^{n-1}p = p\sum_{n=0}^{\infty} (n+1)q^n = \frac{p}{(1-q)^2} = \frac{p}{p^2} = \frac{1}{p}.$$
 (4.4.9)

The random variable X is called the *waiting time*, for heads to fall, or more generally for a "success." The distribution $\{q^{n-1}p; n=1,2,\dots\}$ will be called the *geometrical distribution with success probability* p.

^{*}Named after Abraham Wald (1902-50), leading U.S. statistician.

Example 9. A perfect coin is tossed n times. Let S_n denote the number of heads obtained. In the notation of §2.4, we have $S_n = X_1 + \cdots + X_n$. We know from §3.2 that

$$p_k = P(S_n = k) = \frac{1}{2^n} \binom{n}{k}, \quad 0 \le k \le n.$$
 (4.4.10)

If we believe in probability, then we know $\sum_{k=0}^{\infty} p_k = 1$ from (4.3.10). Hence

$$\sum_{k=0}^{n} \frac{1}{2^n} \binom{n}{k} = 1 \quad \text{or} \quad \sum_{k=0}^{n} \binom{n}{k} = 2^n. \tag{4.4.11}$$

This has been shown in (3.3.7) and can also be obtained from (4.4.13) below by putting x = 1 there, but we have done it by an argument based on probability. Next we have

$$E(S_n) = \sum_{k=0}^{n} \frac{k}{2^n} \binom{n}{k}.$$
 (4.4.12)

Here again we must sum a series, a finite one. We will do it in two different ways, both useful for other calculations. First by direct manipulation, the series may be rewritten as

$$\sum_{k=0}^{n} \frac{k}{2^{n}} \frac{n!}{k! (n-k)!} = \frac{n}{2^{n}} \sum_{k=1}^{n} \frac{(n-1)!}{(k-1)! (n-k)!} = \frac{n}{2^{n}} \sum_{k=1}^{n} \binom{n-1}{k-1}.$$

What we have done above is to cancel k from k!, split off n from n!, and omit a zero term for k = 0. Now change the index of summation by putting k - 1 = j (we have done this kind of thing in Example 8):

$$\frac{n}{2^n} \sum_{k=1}^n \binom{n-1}{k-1} = \frac{n}{2^n} \sum_{i=0}^{n-1} \binom{n-1}{j} = \frac{n}{2^n} \cdot 2^{n-1} = \frac{n}{2},$$

where the step before the last is obtained by using (4.4.11) with n replaced by n-1. Hence the answer is n/2.

This method is highly recommended if you enjoy playing with combinatorial formulas such as the binomial coefficients. But most of you will probably find the next method easier because it is more like a cookbook recipe. Start with Newton's binomial theorem in the form

$$(1+x)^n = \sum_{k=0}^n \binom{n}{k} x^k. \tag{4.4.13}$$

Observe that this is just an expression of a polynomial in x and is a special case of Taylor's series, just as the series in (4.4.5) and (4.4.6) are. Now differentiate to get

$$n(1+x)^{n-1} = \sum_{k=0}^{n} \binom{n}{k} kx^{k-1}.$$
 (4.4.14)

Substitute x = 1:

$$n2^{n-1} = \sum_{k=0}^{n} \binom{n}{k} k;$$

divide through by 2^n , and get the answer n/2 again for (4.4.12). So the expected number of heads in n tosses is n/2. Once more, what could be more reasonable since heads are expected half of the time!

We can generalize this problem to a biased coin, too. Then (4.4.10) becomes

$$P(S_n = k) = \binom{n}{k} p^k q^{n-k}, \quad 0 \le k \le n.$$
 (4.4.15)

There is a preview of the above formula in §2.4. We now see that it gives the probability distribution of the random variable $S_n = \sum_{i=1}^n X_i$. It is called the *binomial distribution* B(n; p). The random variable X_i here as well as its distribution are often referred to as *Bernoullian*; and when p = 1/2, the adjective *symmetric* is added. Next, (4.4.12) becomes

$$\sum_{k=0}^{n} \binom{n}{k} k p^k q^{n-k} = np. (4.4.16)$$

Both methods used above still work. The second is quicker: setting x = p/q in (4.4.14), since p + q = 1 we obtain

$$n\left(1+\frac{p}{q}\right)^{n-1} = \frac{n}{q^{n-1}} = \sum_{k=0}^{n} \binom{n}{k} k \left(\frac{p}{q}\right)^{k-1};$$

multiplying through by pq^{n-1} , we establish (4.4.16).

Example 10. In Problem 1 of §3.4, if we denote by X the number of defective items, then P(X = j) is given by the formula in (3.4.1). This is called the *hypergeometric distribution*.

4.5. Random variables with densities

In the preceding sections we have given a quite rigorous discussion of random variables that take only a countable set of values. But even at an elementary level there are many important questions in which we must consider random variables not subject to such a restriction. This means that we need a sample space that is not countable. Technical questions of "measurability" then arise which cannot be treated satisfactorily without more advanced mathematics. As we have mentioned in Chapter 2, this kind of difficulty stems from the impossibility of assigning a probability to every subset of the sample space when it is uncountable. The matter is resolved by confining ourselves to sample sets belonging to an adequate class called a Borel field; see Appendix 1. Without going into this here we will take up a particular but very important situation that covers most applications and requires little mathematical abstraction. This is the case of a random variable with a "density."

Consider a function f defined on $R^1 = (-\infty, +\infty)$:

$$u \to f(u)$$

and satisfying two conditions:

(i)
$$\forall u$$
: $f(u) \ge 0$;
(ii) $\int_{-\infty}^{\infty} f(u) du = 1$. (4.5.1)

Such a function is called a *density function* on R^1 . The integral in (ii) is the Riemann integral taught in calculus. You may recall that if f is continuous or just *piecewise continuous*, then the definite integral

$$\int_a^b f(u) \, du$$

exists for any interval [a, b]. But in order that the "improper integral" over the infinite range $(-\infty, +\infty)$ should exist, further conditions are needed to make sure that f(u) is pretty small for large |u|. In general, such a function is said to be "integrable over R^1 ." The requirement that the total integral be equal to 1 is less serious than it might appear, because if

$$\int_{-\infty}^{\infty} f(u) \, du = M < \infty,$$

we can just divide through by M and use f/M instead of f. Here are some possible pictures of density functions, some smooth, some not so.

You see what a variety they can be. The only constraints are that the curve should not lie below the x-axis anywhere, that the area under the curve

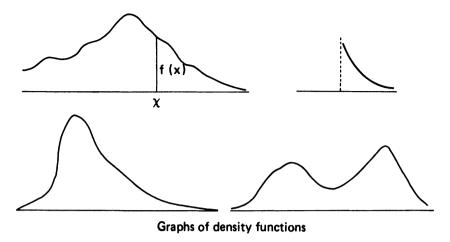


Figure 20

should have a meaning, and the total area should be equal to 1. You may agree that this is not asking for too much.

We can now define a class of random variables on a general sample space as follows. As in §4.2, X is a function on $\Omega: \omega \to X(\omega)$, but its probabilities are prescribed by means of a density function f so that for any interval [a,b] we have

$$P(a \le X \le b) = \int_{a}^{b} f(u) \, du. \tag{4.5.2}$$

More generally, if A is the union of intervals not necessarily disjoint and some of which may be infinite, we have

$$P(X \in A) = \int_{A} f(u) du. \tag{4.5.3}$$

Such a random variable is said to have a density, and its density function is f. [In some books this is called a "continuous" random variable, whereas the kind discussed in $\S 2$ is called "discrete." Both adjectives are slightly misleading so we will not use them here.]

If A is a finite union of intervals, then it can be split up into disjoint ones, some of which may abut on each other, such as

$$A = \bigcup_{j=1}^{k} [a_j, b_j],$$

and then the right-hand side of (4.5.3) may be written as

$$\int_{A} f(u) \, du = \sum_{j=1}^{k} \int_{a_{j}}^{b_{j}} f(u) \, du.$$

This is a property of integrals which is geometrically obvious when you consider them as areas. Next if $A = (-\infty, x]$, then we can write

$$F(x) = P(X \le x) = \int_{-\infty}^{x} f(u) \, du; \tag{4.5.4}$$

compare with (4.3.8). This formula defines the distribution function F of X as a primitive [indefinite integral] of f. It follows from the fundamental theorem of calculus that if f is continuous, then f is the derivative of F:

$$F'(x) = f(x). (4.5.5)$$

Thus in this case the two functions f and F mutually determine each other. If f is not continuous everywhere, (4.5.5) is still true for every x at which f is continuous. These things are proved in calculus.

Let us observe that in the definition above of a random variable with a density, it is *implied* that the sets $\{a \leq X \leq b\}$ and $\{X \in A\}$ have probabilities assigned to them; in fact, they are specified in (4.5.2) and (4.5.3) by means of the density function. This is a subtle point in the wording that should be brought out but will not be elaborated on. [Otherwise we shall be getting into the difficulties we are trying to circumvent here. But see Appendix 1.] Rather, let us remark on the close resemblance between the formulas above and the corresponding ones in §4.3. This will be amplified by a definition of mathematical expectation in the present case and listed below for comparison.

	Countable case	Density case
Range	$v_n, n=1,2,\ldots$	$-\infty < u < +\infty$
element of probability	p_n	f(u) du = dF(u)
$P(a \le X \le b)$	$\sum_{a \le v_n \le b} p_n$	$\int_a^b f(u) du$
$P(X \le x) = F(x)$	$\sum_{v_n \leq x} p_n$	$\int_{-\infty}^{x} f(u) du$
E(X)	$\sum_{n} p_n v_n$	$\int_{-\infty}^{\infty} u f(u) du$
proviso	$ \sum_{n} p_n v_n < \infty $	$\int_{-\infty}^{\infty} u f(u) du < \infty$

More generally, the analogue of (4.3.17) is

$$E(\varphi(X)) = \int_{-\infty}^{\infty} \varphi(u) f(u) du. \tag{4.5.6}$$

You may ignore the second item in the density case above involving a differential if you don't know what it means.

Further insight into the analogy is gained by looking at the following picture:

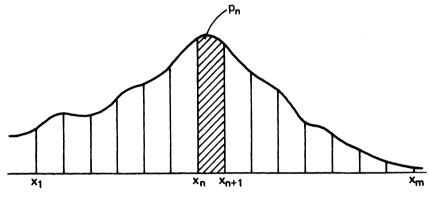


Figure 21

The curve is the graph of a density function f. We have divided the x-axis into m+1 pieces, not necessarily equal and not necessarily small, and denote the area under the curve between x_n and x_{n+1} by p_n , thus:

$$p_n = \int_{x_n}^{x_{n+1}} f(u) du, \quad 0 \le n \le m,$$

where $x_0 = -\infty, x_{m+1} = +\infty$. It is clear that we have

$$\forall n: \quad p_n \ge 0; \quad \sum_n p_n = 1.$$

Hence the numbers p_n satisfy the conditions in (4.3.10). Instead of a finite partition we may have a countable one by suitable labeling such as $\ldots, p_{-2}, p_{-1}, p_0, p_1, \ldots$. Thus we can derive a set of "elementary probabilities" from a density function, in infinitely many ways. This process may be called *discretization*. If X has the density f, we may consider a random variable Y such that

$$P(Y=x_n)=p_n,$$

where we may replace x_n by any other number in the subinterval $[x_n, x_{n+1}]$. Now if f is continuous and the partition is sufficiently fine, namely if the pieces are sufficiently small, then it is geometrically evident that Y is in some sense a discrete approximation of X. For instance,

$$E(Y) = \sum_{n} p_n x_n$$

will be an approximation of $E(X) = \int_{-\infty}^{\infty} uf(u) du$. Remember the *Riemann sums* defined in calculus to lead to a Riemann integral? There the strips with curved tops in Figure 21 are replaced by flat tops (rectangles), but the ideas involved are quite similar. From a practical point of view, it is the discrete approximations that can really be measured, whereas the continuous density is only a mathematical idealization. We shall return to this in a moment.

Having dwelled on the similarity of the two cases of random variable, we will pause to stress a fundamental difference between them. If X has a density, then by (4.5.2) with a = b = x, we have

$$P(X = x) = \int_{x}^{x} f(u) du = 0.$$
 (4.5.7)

Geometrically speaking, this merely states the trivial fact that a line segment has zero area. Since x is arbitrary in (4.5.7), it follows that X takes any preassigned value with probability zero. This is in direct contrast to a random variable taking a countable set of values, for then it must take some of these values with positive probability. It seems paradoxical that on the one hand, $X(\omega)$ must be some number for every ω , and on the other hand any given number has probability zero. The following simple concrete example should clarify this point.

Example 11. Spin a needle on a circular dial. When it stops it points at a random angle θ (measured from the horizontal, say). Under normal conditions it is reasonable to suppose that θ is uniformly distributed between 0° and 360° (cf. Example 7 of §4.4). This means it has the following density function:

$$f(u) = \begin{cases} \frac{1}{360} & \text{for } 0 \le u \le 360, \\ 0 & \text{otherwise.} \end{cases}$$

Thus for any $\theta_1 < \theta_2$ we have

$$P(\theta_1 \le \theta \le \theta_2) = \int_{\theta_1}^{\theta_2} \frac{1}{360} \, du = \frac{\theta_2 - \theta_1}{360}. \tag{4.5.8}$$

This formula says that the probability of the needle pointing between any two directions is proportional to the angle between them. If the angle $\theta_2 - \theta_1$ shrinks to zero, then so does the probability. Hence in the limit the probability of the needle pointing exactly at θ is equal to zero. From an empirical point of view, this event does not really make sense because the needle itself must have a width. So in the end it is the mathematical fiction or idealization of a "line without width" that is the root of the paradox.

There is a deeper way of looking at this situation which is very rich. It should be clear that instead of spinning a needle we may just as well "pick a number at random" from the interval [0,1]. This can be done by bending the circle into a line segment and changing the unit. Now every point in [0,1] can be represented by a decimal such as

$$.141592653589793\cdots. (4.5.9)$$

There is no real difference if the decimal terminates because then we just have all digits equal to 0 from a certain place on, and 0 is no different from any other digit. Thus, to pick a number in [0,1] amounts to picking all its decimal digits one after another. That is the kind of thing a computing machine churns out. Now the chance of picking any prescribed digit, say the first digit "1" above, is equal to 1/10 and the successive pickings from totally independent trials (see §2.4). Hence the chance of picking the 15 digits shown in (4.5.9) is equal to

$$\underbrace{\frac{1}{10} \cdot \frac{1}{10} \cdots \frac{1}{10}}_{15 \text{ times}} = \left(\frac{1}{10}\right)^{15}.$$

If we remember that 10^9 is 1 billion, this probability is already so small that according to Emile Borel [1871–1956; great French mathematician and one of the founders of modern probability theory], it is terrestrially negligible and should be equated to zero! But we have only gone 15 digits in the decimals of the number $\pi-3$, so there can be no question whatsoever of picking this number itself and yet if you can imagine going on forever, you will end up with some number which is just as impossible a priori as this $\pi-3$. So here again we are up against a mathematical fiction—the real number system.

We may generalize this example as follows. Let [a, b] be any finite, non-degenerate interval in \mathbb{R}^1 and put

$$f(u) = \begin{cases} \frac{1}{b-a} & \text{for } a \le u \le b, \\ 0 & \text{otherwise.} \end{cases}$$

This is a density function, and the corresponding distribution is called *the* uniform distribution on [a, b]. We can write the latter explicitly:

$$F(x) = \frac{[(a \lor x) \land b] - a}{b - a}$$

if you have a taste for such tricky formulas.

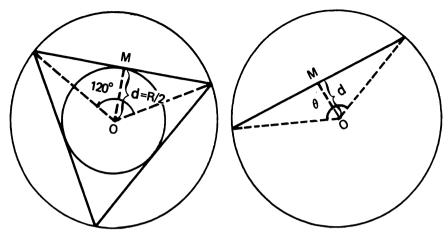


Figure 22

Example 12. A chord is drawn at random in a circle. What is the probability that its length exceeds that of a side of an inscribed equilateral triangle?

Let us draw such a triangle in a circle with center 0 and radius R, and make the following observations. The side is at distance R/2 from 0; its midpoint is on a concentric circle of radius R/2; it subtends an angle of 120 degrees at 0. You ought to know how to compute the length of the side, but this will not be needed. Let us denote by A the desired event that a random chord be longer than that side. Now the length of any chord is determined by any one of the three quantities: its distance d from 0; the location of its midpoint M; the angle θ it subtends at 0. We are going to assume in turn that each of these has a uniform distribution over its range and compute the probability of A under each assumption.

- (1) Suppose that d is uniformly distributed in [0, R]. This is a plausible assumption if we move a ruler parallel to itself with constant speed from a tangential position toward the center, stopping somewhere to intersect the circle in a chord. It is geometrically obvious that the event A will occur if and only if d < R/2. Hence P(A) = 1/2.
- (2) Suppose that M is uniformly distributed over the disk D formed by the given circle. This is a plausible assumption if a tiny dart is thrown at D and a chord is then drawn perpendicular to the line joining the hitting point to 0. Let D' denote the concentric disk of radius R/2. Then the event A will occur if and only if M falls within D'. Hence $P(A) = P(M \in D') = (\text{area of } D')/(\text{area of } D) = 1/4$.
- (3) Suppose that θ is uniformly distributed between 0 and 360 degrees. This is plausible if one endpoint of the chord is arbitrarily fixed and the other is obtained by rotating a radius at constant speed to stop somewhere

on the circle. Then it is clear from the picture that A will occur if and only if θ is between 120 and 240 degrees. Hence P(A) = (240 - 120)/360 = 1/3.

Thus the answer to the problem is 1/2, 1/4, or 1/3 according to the different hypotheses made. It follows that these hypotheses are not compatible with one another. Other hypotheses are possible and may lead to still other answers. Can you think of a good one? This problem was known as Bertrand's paradox in the earlier days of discussions of probability theory. But of course the paradox is due only to the fact that the problem is not well posed without specifying the underlying nature of the randomness. It is not surprising that the different ways of randomization should yield different probabilities, which can be verified experimentally by the mechanical procedures described. Here is a facile analogy. Suppose that you are asked how long it takes to go from your dormitory to the classroom without specifying whether we are talking about "walking," "biking," or "driving" time. Would you call it paradoxical that there are different answers to the question?

We end this section with some other simple examples of random variables with densities. Another important case, the normal distribution, will be discussed in Chapter 6.

Example 13. Suppose you station yourself at a spot on a relatively serene country road and watch the cars that pass by that spot. With your stopwatch you can clock the time before the first car passes. This is a random variable T called the *waiting time*. Under certain circumstances it is a reasonable hypothesis that T has the density function below with a certain $\lambda > 0$:

$$f(u) = \lambda e^{-\lambda u}, \quad u \ge 0. \tag{4.5.10}$$

It goes without saying that f(u) = 0 for u < 0. The corresponding distribution function is called the *exponential distribution with parameter* λ , obtained by integrating f as in (4.5.4):

$$F(x) = \int_{-\infty}^{x} f(u) du = \int_{0}^{x} \lambda e^{-\lambda u} du = 1 - e^{-\lambda x}.$$

In particular, if we put $x = +\infty$, or better, let $x \to \infty$ in the above, we see that f satisfies the conditions in (4.5.1), so it is indeed a density function. We have

$$P(T \le x) = F(x) = 1 - e^{-\lambda x};$$
 (4.5.11)

but in this case it is often more convenient to use the *tail probability*:

$$P(T > x) = 1 - F(x) = e^{-\lambda x}.$$
 (4.5.12)

This can be obtained directly from (4.5.3) with $A = (x, \infty)$; thus

$$P(T \in (x, \infty)) = \int_{(x, \infty)} \lambda e^{-\lambda u} du = \int_{x}^{\infty} \lambda e^{-\lambda u} du = e^{-\lambda x}.$$

For every given x, say 5 (seconds), the probability $e^{-5\lambda}$ in (4.5.12) decreases as λ increases. This means your waiting time tends to be shorter if λ is larger. On a busy highway λ will indeed be large. The expected waiting-time is given by

$$E(T) = \int_0^\infty u\lambda e^{-\lambda u} du = \frac{1}{\lambda} \int_0^\infty t e^{-t} dt = \frac{1}{\lambda}.$$
 (4.5.13)

[Can you compute the integral above using "integration by parts" without recourse to a table?] This result supports our preceding observation that T tends on the average to be smaller when λ is larger.

The exponential distribution is a very useful model for various types of waiting time problems such as telephone calls, service times, splitting of radioactive particles, etc.; see §7.2.

Example 14. Suppose in a problem involving the random variable T above, what we really want to measure is its logarithm (to the base e):

$$S = \log T. \tag{4.5.14}$$

This is also a random variable (cf. Proposition 2 in §4.2); it is negative if T > 1, zero if T = 1, and positive if T > 1. What are its probabilities? We may be interested in $P(a \le S \le b)$, but it is clear that we need only find $P(S \le x)$, namely the distribution function F_S of S. Now the function

$$x \to \log x$$

is monotone and its inverse is

$$x \to e^x$$

so that

$$S \le x \Leftrightarrow \log T \le x \Leftrightarrow T \le e^x$$
.

Hence by (4.5.11)

$$F_S(x) = P\{S \le x\} = P\{T \le e^x\} = 1 - e^{-\lambda e^x}.$$

The density function f_S is obtained by differentiating:

$$f_S(x) = F_S'(x) = \lambda e^x e^{-\lambda e^x} = \lambda e^{x - \lambda e^x}.$$

This looks formidable, but you see it is easily derived.

Example 15. A certain river floods every year. Suppose the low-water mark is set at 1, and the high-water mark Y has the distribution function

$$F(y) = P(Y \le y) = 1 - \frac{1}{y^2}, \quad 1 \le y < \infty. \tag{4.5.15}$$

Observe that F(1)=0, that F(y) increases with y, and that $F(y)\to 1$ as $y\to\infty$. This is as it should be from the meaning of $P(Y\leq y)$. To get the density function we differentiate:

$$f(y) = F'(y) = \frac{2}{y^3}, \quad 1 \le y < \infty.$$
 (4.5.16)

It is not necessary to check that $\int_{-\infty}^{\infty} f(y) dy = 1$, because this is equivalent to $\lim_{y\to\infty} F(y) = 1$. The expected value of Y is given by

$$E(Y) = \int_{1}^{\infty} u \cdot \frac{2}{u^3} du = \int_{1}^{\infty} \frac{2}{u^2} du = 2.$$

Thus the maximum of Y is twice that of the minimum, on the average.

What happens if we set the low-water mark at 0 instead of 1 and use a unit of measuring the height which is 1/10 of that used above? This means we set

$$Z = 10(Y - 1). (4.5.17)$$

As in Example 13 we have

$$Z \le z \Leftrightarrow 10(Y-1) \le z \Leftrightarrow Y \le 1 + \frac{z}{10}, \quad 0 \le z < \infty.$$

From this we can compute

$$F_Z(z) = 1 - \frac{100}{(10+z)^2},$$

 $f_Z(z) = \frac{200}{(10+z)^3}.$

The calculation of E(Z) from f_Z is tedious but easy. The answer is E(Z) = 10, and comparing with E(Y) = 2 we see that

$$E(Z) = 10(E(Y) - 1). (4.5.18)$$

Thus the *means* of Y and Z are connected by the same linear relation as the random variables themselves. Does this seem obvious to you? The general proposition will be discussed in $\S6.1$.

4.6. General case

The most general random variable is a function X defined on the sample space Ω such that for any real x, the probability $P(X \leq x)$ is defined.

To be frank, this statement has put the cart before the horse. What comes first is a probability measure P defined on a class of subsets of Ω . This class is called the sample Borel field or probability field and is denoted by \mathfrak{F} . Now if a function X has the property that for every x, the set $\{\omega \mid$ $X(\omega) < x$ belongs to the class \mathfrak{F} , then it is called a random variable. [We must refer to Appendix 1 for a full description of this concept; but the rest of this section should be intelligible without the formalities.] In other words, an arbitrary function must pass a test to become a member of the club. The new idea here is that P is defined only for subsets in \mathfrak{F} , not necessarily for all subsets of Ω . If it happens to be defined for all subsets, then of course the test described above becomes a nominal one and every function is automatically a random variable. This is the situation for a countable space Ω discussed in §4.1. In general, as we have hinted several times before, it is impossible to define a probability measure on all subsets of Ω , and so we must settle for a certain class \mathfrak{F} . Since only sets in \mathfrak{F} have probabilities assigned to them, and since we wish to discuss sample sets of the sort "X <x," we are obliged to require that these belong to \mathfrak{F} . Thus the necessity of such a test is easy to understand. What may be a little surprising is that this test is all we need. Namely, once we have made this requirement, we can then go on to discuss the probabilities of a whole variety of sample sets such as $\{a \leq X \leq b\}$, $\{X = x\}$, $\{X \text{ takes a rational value}\}$, or some crazy thing like $\{e^x > X^2 + 1\}.$

Next, we define for every real x:

$$F(x) = P(X \le x) \tag{4.6.1}$$

or equivalently for a < b:

$$F(b) - F(a) = P(a < X \le b);$$

and call the function F the distribution function of X. This has been done in previous cases, but we no longer have the special representative in (4.3.8) or (4.5.4):

$$F(x) = \sum_{\nu_n < x} p_n, \quad F(x) = \int_{-\infty}^x f(u) du$$

in terms of elementary probability or a density function. As a matter of fact, the general F turns out to be a mixture of these two kinds together with a weirder kind (the *singular* type). But we can operate quite well with the F as defined by (4.6.1) without further specification. The mathematical

equipment required to handle the general case, however, is somewhat more advanced (at the level of a course like "Fundamental concepts of analysis"). So we cannot go into this but will just mention two easy facts about F:

- (i) F is monotone nondecreasing: namely $x \le x' \Rightarrow F(x) \le F(x')$;
- (ii) F has limits 0 and 1 at $-\infty$ and $+\infty$, respectively:

$$F(-\infty) = \lim_{x \to -\infty} F(x) = 0, \quad F(+\infty) = \lim_{x \to +\infty} F(x) = 1.$$

Property (i) holds because if $x \le x'$, then $\{X \le x\} \subset \{X \le x'\}$. Property (ii) is intuitively obvious because the event $\{X \le x\}$ becomes impossible as $x \to -\infty$, and certain as $x \to +\infty$. This argument may satisfy you, but the rigorous proofs are a bit more sophisticated and depend on the countable additivity of P (see §2.3). Let us note that the existence of the limits in (ii) follows from the monotonicity in (i) and a fundamental theorem in calculus: a bounded monotone sequence of real numbers has a limit.

The rest of the section is devoted to a brief discussion of some basic notions concerning random vectors. This material may be postponed until it is needed in Chapter 6.

For simplicity of notation we will consider only two random variables X and Y, but the extension to any finite number is straightforward. We first consider the case where X and Y are countably valued. Let X take the values $\{x_i\}$, Y take the values $\{y_j\}$, and put

$$P(X = x_i, Y = y_j) = p(x_i, y_j). (4.6.2)$$

When x_i and y_j range over all possible values, the set of "elementary probabilities" above gives the *joint probability distribution* of the random vector (X, Y). To get the probability distribution of X alone, we let y_j range over all possible values in (4.6.2); thus

$$P(X = x_i) = \sum_{y_i} p(x_i, y_j) = p(x_i, *), \tag{4.6.3}$$

where the last quantity is defined by the middle sum. When x ranges over all possible values, the set of $p(x_i, *)$ gives the marginal distribution of X. The marginal distribution of Y is similarly defined. Let us observe that these marginal distributions do not in general determine the joint distribution.

Just as we can express the expectation of any function of X by means of its probability distribution [see (4.3.17)], we can do the same for any function of (X,Y) as follows:

$$E(\varphi(X,Y)) = \sum_{x_i} \sum_{y_j} \varphi(x_i, y_j) p(x_i, y_j). \tag{4.6.4}$$

It is instructive to see that this results from a rearrangement of terms in the definition of the expectation of $\varphi(X,Y)$ as *one* random variable as in (4.3.11):

$$E(\varphi(X,Y)) = \sum_{\omega} \varphi(X(\omega), Y(\omega)) P(\omega).$$

Next, we consider the density case extending the situation in $\S 4.5$. The random vector (X,Y) is said to have a joint density function f in case

$$P(X \le x, Y \le y) = \int_{-\infty}^{x} \int_{-\infty}^{y} f(u, v) \, du \, dv \tag{4.6.5}$$

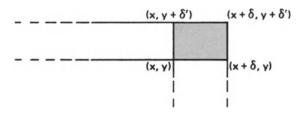
for all (x, y). It then follows that for any "reasonable" subset S of the Cartesian plane (called a *Borel set*), we have

$$P((X,Y) \in S) = \iint_{S} f(u,v) \, du \, dv. \tag{4.6.6}$$

For example, S may be polygons, disks, ellipses, and unions of such shapes. Note that (4.6.6) contains (4.6.5) as a very particular case and we can, at a pinch, accept the more comprehensive condition (4.6.6) as the *definition* of f as density for (X,Y). However, here is a heuristic argument from (4.6.5) to (4.6.6). Let us denote by R(x,y) the infinite rectangle in the plane with sides parallel to the coordinate axes and lying to the southwest of the point (x,y). The picture below shows that for any $\delta > 0$ and $\delta' > 0$:

$$R(x + \delta, y + \delta') - R(x + \delta, y) - R(x, y + \delta') + R(x, y)$$

is the shaded rectangle



It follows that if we manipulate the relation (4.6.5) in the same way, we get

$$P(x \le X \le x + \delta, y \le Y \le y + \delta') = \int_x^{x + \delta} \int_y^{y + \delta'} f(u, v) \, du \, dv.$$

This means (4.6.6) is true for the shaded rectangle. By varying x, y as well as δ, δ' , we see that the formula is true for any rectangle of this shape. Now

any reasonable figure can be approximated from inside and outside by a number of such small rectangles (even just squares)—a fact known already to the ancient Greeks. Hence in the limit we can get (4.6.6) as asserted.

The curious reader may wonder why a similar argument was not given earlier for the case of one random variable (4.5.3)? The answer is: heuristically speaking, there are hardly any sets in R^1 other than intervals, points, and their unions! Things are pretty tight in one dimension and our geometric intuition does not work well. This is one reason why classical measure theory is a sophisticated business.

The joint density function f satisfies the following conditions:

- (i) $f(u,v) \ge 0$ for all (u,v);
- (ii) $\int_{-\infty}^{\infty} \int_{-\infty}^{\infty} f(u, v) du dv = 1.$

Of course, (ii) implies that f is integrable over the whole plane. Frequently we also assume that f is continuous. Now the formulas analogous to (4.6.3) are

$$P(X \le x) = \int_{-\infty}^{x} f(u, *) du, \quad \text{where} \quad f(u, *) = \int_{-\infty}^{\infty} f(u, v) dv,$$

$$P(Y \le y) = \int_{-\infty}^{y} f(*, v) dv, \quad \text{where} \quad f(*, v) = \int_{-\infty}^{\infty} f(u, v) du.$$

$$(4.6.7)$$

The functions $u \to f(u,*)$ and $v \to f(*,v)$ are respectively called the marginal density functions of X and Y. They are derived from the joint density function after "integrating out" the variable that is not in question.

The formula corresponding to (4.6.4) becomes in the density case: for any "reasonable" [Borel] function φ :

$$E(\varphi(X,Y)) = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} \varphi(u,v) f(u,v) \, du \, dv. \tag{4.6.8}$$

The class of reasonable functions includes all bounded continuous functions in (u, v), indicators of reasonable sets, and functions that are continuous except across some smooth boundaries, for which the integral above exists, etc.

In the most general case the *joint distribution function* F of (X,Y) is defined by

$$F(x,y) = P(X \le x, Y \le y) \quad \text{for all } (x,y). \tag{4.6.9}$$

If we denote $\lim_{y\to\infty} F(x,y)$ by $F(x,\infty)$, we have

$$F(x,\infty) = P(X \le x, Y < \infty) = P(X \le x)$$

since " $Y < \infty$ " puts no restriction on Y. Thus $x \to F(x, \infty)$ is the marginal distribution function of X. The marginal distribution function of Y is similarly defined.

Although these general concepts form the background whenever several random variables are discussed, explicit use of them will be rare in this book.

Exercises

1. If X is a random variable [on a countable sample space], is it true that

$$X + X = 2X, X - X = 0$$
?

Explain in detail.

2. Let $\Omega = \{\omega_1, \omega_2, \omega_3\}$, $P(\omega_1) = P(\omega_2) = P(\omega_3) = 1/3$, and define X, Y, and Z as follows:

$$X(\omega_1) = 1, X(\omega_2) = 2, X(\omega_3) = 3;$$

 $Y(\omega_1) = 2, Y(\omega_2) = 3, mY(\omega_3) = 1;$
 $Z(\omega_1) = 3, Z(\omega_2) = 1, Z(\omega_3) = 2.$

Show that these three random variables have the same probability distribution. Find the probability distributions of X + Y, Y + Z, and Z + X.

3. In No. 2 find the probability distribution of

$$X + Y - Z, \sqrt{(X^2 + Y^2)Z}, \frac{Z}{|X - Y|}.$$

- 4. Take Ω to be a set of five real numbers. Define a probability measure and a random variable X on it that takes the values 1, 2, 3, 4, 5 with probabilities 1/10, 1/10, 1/5, 1/5, 2/5 respectively; another random variable Y that takes the value $\sqrt{2}, \sqrt{3}, \pi$ with probabilities 1/5, 3/10, 1/2. Find the probability distribution of XY. [Hint: the answer depends on your choice and is not unique.]
- 5. Generalize No. 4 by constructing Ω, P, X so that X takes the values v_1, v_2, \ldots, v_n with probabilities p_1, p_2, \ldots, p_n where the p_n 's satisfy (4.3.10).
- 6. In Example 3 of §4.1, what do the following sets mean?

$${X + Y = 7}, {X + T \le 7}, {X \lor Y > 4}, {X \ne Y}$$

List all the ω 's in each set.

*7. Let X be integer-valued and let F be its distribution function. Show that for every x and a < b

$$P(X = x) = \lim_{\epsilon \downarrow 0} [F(x + \epsilon) - F(x - \epsilon)],$$

$$P(a < X < b) = \lim_{\epsilon \downarrow 0} [F(b - \epsilon) - F(a + \epsilon)].$$

[The results are true for any random variable but require more advanced proofs even when Ω is countable.]

8. In Example 4 of §4.2, suppose that

$$B = 5000 + X'$$

where X' is uniformly distributed over the set of integers from 1 to 5000. What does this hypothesis mean? Find the probability distribution and mean of Y under this hypothesis.

9. As in No. 8 but now suppose that

$$X = 4000 + X'$$

where X' is uniformly distributed from 1 to 10000.

*10. As in No. 8 but now suppose that

$$X = 3000 + X'$$

and X' is the exponential distribution with mean 7000. Find E(Y).

11. Let $\lambda > 0$ and define f as follows:

$$f(u) = \begin{cases} \frac{1}{2} \lambda e^{-\lambda u} & \text{if } u \ge 0; \\ \frac{1}{2} \lambda e^{+\lambda u} & \text{if } u < 0. \end{cases}$$

This f is called bilateral exponential. If X has density f, find the density of |X|. [Hint: begin with the distribution function.]

- 12. If X is a positive random variable with density f, find the density of $+\sqrt{X}$. Apply this to the distribution of the side length of a square when its area is uniformly distributed in [a, b].
- 13. If X has density f, find the density of (i) aX + b where a and b are constants; (ii) X^2 .
- 14. Prove (4.4.5) in two ways: (a) by multiplying out $(1-x)(1+x+\cdots+x^n)$; (b) by using Taylor's series.
- 15. Suppose that

$$p_n = cq^{n-1}p, \quad 1 \le n \le m,$$

where c is a constant and m is a positive integer; cf. (4.4.8). Determine c so that $\sum_{n=1}^{m} p_n = 1$. (This scheme corresponds to the waiting time for a success when it is supposed to occur within m trials.)

- 16. A perfect coin is tossed n times. Let Y_n denote the number of heads obtained minus the number of tails. Find the probability distribution of Y_n and its mean. [Hint: there is a simple relation between Y_n and the S_n in Example 9 of $\S4.4$.]
- 17. Refer to Problem 1 in $\S 3.4$. Suppose there are 11 rotten apples in a bushel of 550, and 25 apples are picked at random. Find the probability distribution of the number X of rotten apples among those picked.
- *18. Generalize No. 17 to arbitrary numbers and find the mean of X. [Hint: this requires some expertise in combinatorics but becomes trivial after $\S 6.1$.]
 - 19. Let

$$P(X = n) - p_n = \frac{1}{n(n+1)}, \quad n \ge 1.$$

Is this a probability distribution for X? Find $P(X \ge m)$ for any m and E(X).

20. If all the books in a library have been upset and a monkey is hired to put them all back on the shelves, it can be shown that a good approximation for the probability of having exactly n books put back in their original places is

$$\frac{e^{-1}}{n!}, \quad n \ge 0.$$

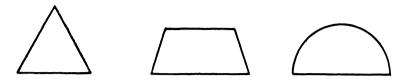
Find the expected number of books returned to their original places. [This oft-quoted illustration is a variant on the matching problem discussed in Problem 6 of §3.4.]

- 21. Find an example in which the series $\sum_{n} p_n v_n$ in (4.3.11) converges but not absolutely. [Hint: there is really nothing hard about this: choose $p_n = 1/2^n$ say, and now choose v_n so that $p_n v_n$ is the general term of any nonabsolutely convergent series you know.]
- 22. If f and g are two density functions, show that $\lambda f + \mu g$ is also a density function, where $\lambda + \mu = 1, \lambda \geq 0, \mu \geq 0$.
- 23. Find the probability that a random chord drawn in a circle is longer than the radius. As in Example 11 of $\S4.5$, work this out under the three different hypotheses discussed there.
- 24. Let

$$f(u) = ue^{-u}, \quad u \ge 0.$$

Show that f is a density function. Find $\int_0^\infty u f(u) du$.

25. In the figure below an equilateral triangle, a trapezoid, and a semidisk are shown:



Determine numerical constants for the sides and radius to make these the graphs of density functions.

- 26. Suppose a target is a disk of radius 10 feet and that the probability of hitting within any concentric disk is proportional to the area of the disk. Let R denote the distance of the bullet from the center. Find the distribution function, density function, and mean of R.
- 27. Agent 009 was trapped between two narrow abysmal walls. He swung his gun around in a vertical circle touching the walls as shown in Fig. 23, and fired a wild [random] shot. Assume that the angle his pistol makes with the horizontal is uniformly distributed between 0° and 90°. Find the distribution of the height where the bullet landed and its mean.
- 28. [St. Petersburg paradox] You play a game with your pal by tossing a perfect coin repeatedly and betting on the waiting time X until a head is tossed up. You agree to pay him $2^X \not\in$ when the value of X is known, namely $2^n \not\in$ if X = n. If you figure that a fair price for him to pay you in advance in order to win this random prize should be equal to the mathematical expectation $E(2^X)$, how much should he pay? How much, honestly, would you accept to play this game? [If you do not see any paradox in this, then you do not agree with such illustrious mathematicians as Daniel Bernoulli, D'Alembert, Poisson, Borel, to name only a few. For a brief account see [Keynes]. Feller believed that the paradox would go away if more advanced mathematics were used to reformulate the problem. You will have to decide for yourself whether it is not more interesting as a philosophical and psychological challenge. [See, however, Appendix 3 ahead.]
- 29. One objection to the scheme in No. 28 is that "time must have a stop." So suppose that only m tosses at most are allowed and your pal gets nothing if a head does not show up in m tosses. Try m=10 and m=100. What is now a fair price for him to pay? And do you feel more comfortable after this change of rule? In this case Feller's explanation melts away but the psychological element remains.
- *30. A number of μ is called the *median* of the random variable X iff $P(X \geq \mu) \geq 1/2$ and $P(X \leq \mu) \geq 1/2$. Show that such a number always exists but need not be unique. Here is a practical example. After n examination papers have been graded, they are arranged in descending order. There is one in the middle if n is odd, two if n is

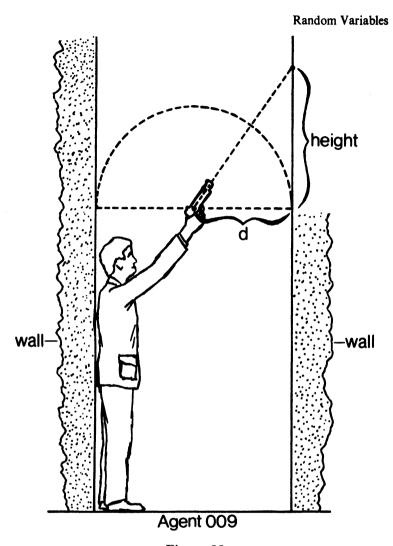


Figure 23

even, corresponding to the median(s). Explain the probability model used.

- 31. An urn contains n tickets numbered from 1 to n. Two tickets are drawn (without replacement). Let X denote the smaller, Y the larger of the two numbers so obtained. Describe the joint distribution of (X,Y), and the marginal ones. Find the distribution of Y-X from the joint distribution.
- 32. Pick two numbers at random from [0,1]. Define X and Y as No. 31 and answer the same question. [Hint: draw the picture and compute areas.]

Appendix 1

Borel Fields and General Random Variables

When the sample space Ω is uncountable, it may not be possible to define a probability measure for all its subsets, as we did for a countable Ω in §2.4. We must restrict the measure to sets of a certain family which must, however, be comprehensive enough to allow the usual operations with sets. Specifically, we require the family \mathfrak{F} to have two properties:

- (a) if a set A belongs to \mathfrak{F} , then its complement $A^c = \Omega A$ also belongs to \mathfrak{F} ;
- (b) if a countable number of sets A_1, A_2, \ldots all belong to \mathfrak{F} , then their union $\bigcup_n A_n$ also belongs to \mathfrak{F} .

It follows from De Morgan's laws that the union in (b) may be replaced by the intersection $\bigcap_n A_n$ as well. Thus if we operate on the members of the family with the three basic operations mentioned above, for a countable number of times, in any manner or order (see, e.g., (1.3.1)), the result is still a member of the family. In this sense the family is said to be *closed under* these operations, and so also under other derived operations such as differences. Such a family of subsets of Ω is called a *Borel field* on Ω . In general there are many such fields, for example the family of all subsets which is certainly a Borel field but may be too large to have a probability defined on it; or the family of two sets $\{\emptyset, \Omega\}$ or four sets $\{\emptyset, A, A^c, \Omega\}$ with a fixed set A, which are too small for most purposes. Now suppose that a reasonable Borel field \mathfrak{F} has been chosen and a probability measure P has been defined on it; then we have a *probability triple* $(\Omega, \mathfrak{F}, P)$ with which

we can begin our work. The sets in $\mathfrak F$ are said to be measurable and they alone have probabilities.

Let X be a real-valued function defined on Ω . Then X is called a random variable iff for any real number x, we have

$$\{\omega \mid X(\omega) \le x\} \in \mathfrak{F}.$$
 (A.1.1)

Hence $P\{X \leq x\}$ is defined, and as a function of x it is the distribution function F given in (4.6.1). Furthermore if a < b, then the set

$$\{a < X \le b\} = \{X \le b\} - \{X \le a\} \tag{A.1.2}$$

belongs to \mathfrak{F} since \mathfrak{F} is closed under difference. Thus its probability is defined and is in fact given by F(b) - F(a).

When Ω is countable and we take \mathfrak{F} to be the Borel field of all the subsets of Ω , then of course the condition (A.1.1) is satisfied for any function X. Thus in this case an arbitrary function on Ω is a random variable, as defined in §4.2. In general, the condition in (A.1.1) is imposed mainly because we wish to define the *mathematical expectation* by a procedure that requires such a condition. Specifically, if X is a bounded random variable, then it has an expectation given by the formula below:

$$E(X) = \lim_{\delta \downarrow 0} \sum_{n = -\infty}^{\infty} n \, \delta P\{n \, \delta < X \le (n+1) \, \delta\},\tag{A.1.3}$$

where the probabilities in the sum are well defined by the remark about (A.1.2). The existence of the limit in (A.1.3), and the consequent properties of the expectation that extend those discussed in Chapters 5 and 6, are part of a general theory known as that of *Lebesgue integration* [Henri Lebesgue (1875-1941), co-founder with Borel of the modern school of measure and integration]. We must refer the reader to standard treatments of the subject except to exhibit E(X) as an integral as follows:

$$E(X) = \int_{\Omega} X(\omega) P(d\omega);$$

cf. the discrete analogue (4.3.11) in a countable Ω .

Conditioning and Independence

5.1. Examples of conditioning

We have seen that the probability of a set A is its weighted proportion relative to the sample space Ω . When Ω is finite and all sample points have the same weight (therefore equally likely), then

$$P(A) = \frac{|A|}{|\Omega|}$$

as in Example 4 of §2.2. When Ω is countable and each point ω has the weight $P(\omega) = P(\{\omega\})$ attached to it, then

$$P(A) = \frac{\sum_{\omega \in A} P(\omega)}{\sum_{\omega \in \Omega} P(\omega)}$$
 (5.1.1)

from (2.4.3), since the denominator above is equal to 1. In many questions we are interested in the proportional weight of one set A relative to another set S. More accurately stated, this means the proportional weight of the part of A in S, namely the intersection $A \cap S$, or AS, relative to S. The formula analogous to (5.1.1) is then

$$\frac{\sum_{\omega \in AS} P(\omega)}{\sum_{\omega \in S} P(\omega)}.$$
 (5.1.2)

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Thus we are switching our attention from Ω to S as a new universe, and considering a new proportion or probability with respect to it. We introduce the notation

$$P(A \mid S) = \frac{P(AS)}{P(S)} \tag{5.1.3}$$

and call it the conditional probability of A relative to S. Other phrases such as "given S," "knowing S," or "under the hypothesis [of] S" may also be used to describe this relativity. Of course, if P(S)=0, then the ratio in (5.1.3) becomes the "indeterminate" 0/0, which has neither meaning nor utility; so whenever we write a conditional probability such as $P(A\mid S)$ we shall impose the proviso that P(S)>0 even if this is not explicitly mentioned. Observe that the ratio in (5.1.3) reduces to that in (5.1.2) when Ω is countable, but is meaningful in the general context where the probabilities of A and S are defined. The following preliminary examples will illustrate the various possible motivations and interpretations of the new concept.

Example 1. All students on a certain college campus are polled as to their reaction to a certain presidential candidate. Let D denote those who favor him. Now the student population Ω may be cross-classified in various ways, for instance according to sex, age, race, etc. Let

$$A = \text{female}, B = \text{black}, C = \text{of voting age}.$$

Then Ω is partitioned as in (1.3.5) into 8 subdivisions ABC, ABC^c , ..., $A^cB^cC^c$. Their respective numbers will be known if a complete poll is made, and the set D will in general cut across the various divisions. For instance,

$$P(D \mid A^cBC) = \frac{P(A^cBCD)}{P(A^cBC)}$$

denotes the proportion of male black students of voting age who favor the candidate;

$$P(D^c \mid A^c C) = \frac{P(A^c C D^c)}{P(A^c C)}$$

denotes the proportion of male students of voting age who do not favor the the candidate, etc.

Example 2. A perfect die is thrown twice. Given [knowing] that the total obtained is 7, what is the probability that the first point obtained is k, $1 \le k \le 6$?

Look at the list in Example 3 of §4.1. The outcomes with total equal to 7 are those on the "second diagonal," and their number is 6. Among

these there is one case in which the first throw is k. Hence the conditional probability is equal to 1/6. In symbols, let X_1 and X_2 respectively denote the point obtained in the first and second throw. Then we have as a case of (5.1.3), with $A = \{X_1 = k\}, S = \{X_1 + X_2 = 7\}$:

$$P\{X_1 = k \mid X_1 + X_2 = 7\} = \frac{P\{X_1 = k; X_1 + X_2 = 7\}}{P\{X_1 + X_2 = 7\}} = \frac{1}{6}.$$

The fact that this turns out to be the same as the unconditional probability $P\{X_1 = k\}$ is an accident due to the lucky choice of the number 7. It is the only value of the total that allows all six possibilities for each throw. As other examples, we have

$$\begin{split} &P\{X_1=k\mid X_1+X_2=6\}=\frac{1}{5},\quad 1\leq k\leq 5,\\ &P\{X_1=k\mid X_1+X_2=9\}=\frac{1}{4},\quad 3\leq k\leq 6. \end{split}$$

Here it should be obvious that the conditional probabilities will be the same if X_1 and X_2 are interchanged. Why?

Next, we ask the apparently simpler question: given $X_1 = 4$, what is the probability that $X_2 = k$? You may jump to the answer that this must be 1/6 since the second throw is not affected by the first, so the conditional probability $P\{X_2 = k \mid X_1 = 4\}$ must be the same as the unconditional one $P\{X_2 = k\}$. This is certainly correct provided we use the *independence* between the two trials (see §2.4). For the present we can use (5.1.3) to get

$$P\{X_2 = k \mid X_1 = 4\} = \frac{P\{X_1 = 4; X_2 = k\}}{P\{X_1 = 4\}} = \frac{\frac{1}{36}}{\frac{1}{6}} = \frac{1}{6}.$$
 (5.1.4)

Finally, we have

$$P\{X_1 + X_2 = 7 \mid X_1 = 4\} = \frac{P\{X_1 = 4; X_1 + X_2 = 7\}}{p\{X_1 = 4\}}.$$
 (5.1.5)

Without looking at the list of outcomes, we observe that the event $\{X_1 = 4; X_1 + X_2 = 7\}$ is exactly the same as $\{X_1 = 4; X_2 = 7 - 4 = 3\}$; so in effect (5.1.5) is a case of (5.1.4). This argument may seem awfully devious at this juncture, but is an essential feature of a random walk (see Chapter 8).

Example 3. Consider the waiting time X in Example 8 of §4.4, for a biased coin. Knowing that it has fallen tails three times, what is the probability that it will fall heads within the next two trials?

This is the conditional probability

$$P(X \le 5 \mid X \ge 4) = \frac{P(4 \le X \le 5)}{P(X \ge 4)}.$$
 (5.1.6)

We know that

$$P(X = n) = q^{n-1}p, \quad n = 1, 2, \dots;$$
 (5.1.7)

from which we can calculate

$$P(X \ge 4) = \sum_{n=4}^{\infty} q^{n-1} p = \frac{q^3 p}{1-q} = q^3$$
 (5.1.8)

(how do we sum the series?) Again from (5.1.7),

$$P(4 \le X \le 5) = q^3 p + q^4 p.$$

Thus the answer to (5.1.6) is p + qp. Now we have also from (5.1.7) the probability that the coin falls heads (at least once in two trials):

$$P(1 \le X \le 2) = p + qp.$$

Comparing these two results, we conclude that the three previous failures do not affect the future waiting time. This may seem obvious to you a priori, but it is a consequence of independence of the successive trials. By the way, many veteran gamblers at the roulette game believe that "if reds have appeared so many times in a row, then it is smart to bet on the black on the next spin because in the long run red and black should balance out." On the other hand, you might argue (with Lord Keynes* on your side) that if red has appeared say 10 times in a row, in the absence of other evidence, if would be a natural presumption that the roulette wheel or the croupier is biased toward the red, namely p > 1/2 in the above, and therefore the smart money should be on it. See Example 8 in §5.2 below for a similar discussion.

Example 4. We shall bring out an analogy between the geometrical distribution given in (5.1.7) [see also (4.4.8)] and the exponential distribution in (4.5.11). If X has the former distribution, then for any nonnegative integer n we have

$$P(X > n) = q^n. (5.1.9)$$

This can be shown by summing a geometrical series as in (5.1.8), but it is obvious if we remember that "X > n" means that the first n tosses all show tails. It now follows from (5.1.9) that for any nonnegative integers m and n, we have

$$P(X > n + m \mid X > m) = \frac{P(X > n + m)}{P(X > m)} = \frac{q^{m+n}}{q^m}$$
 (5.1.10)
= $q^n = P(X > n)$.

^{*}John Maynard Keynes [1883-1946], English economist and writer.

Now let T denote the waiting time in Example 12 of §4.5; then we have analogously for any nonnegative real values of s and t:

$$P(T > s + t \mid T > s) = \frac{P(T > s + t)}{P(T > s)} = \frac{e^{-\lambda(s+t)}}{e^{-\lambda s}}$$
 (5.1.11)
= $e^{-\lambda t} = P(T > t)$.

This may be announced as follows: if we have already spent some time in waiting, the distribution of further waiting time is the same as that of the initial waiting time as if we have waited in vain! A suggestive way of saying this is that the random variable T has no memory. This turns out to be a fundamental property of the exponential distribution which is not shared by any other and is basic for the theory of Markov processes. Note that although the geometrical distribution is a discrete analogue as shown in (5.1.10), strictly speaking it does not have the "memoryless" property because (5.1.10) may become false when n and m are not integers: take, e.g., n = m = 1/2.

Example 5. Consider all families with two children and assume that boys and girls are equally likely. Thus the sample space may be denoted schematically by 4 points:

$$\Omega = \{(bb), (bg), (gb), (gg)\}$$

where b = boy, g = girl; the order in each pair is the order of birth; and the 4 points have probability 1/4 each. We may of course instead use a space of 4N points, where N is a large number, in which the four possibilities have equal numbers. This will be a more realistic population model, but the arithmetic below will be the same.

If a family is chosen at random from Ω and found to have a boy in it, what is the probability that it has another boy, namely that it is of the type (b,b)? A quickie answer might be 1/2 if you jumped to the conclusion from the equal likelihood of the sexes. This is a mistake induced by a misplaced "relative clause" for the conditional probability in question. Here is the detailed explanation.

Let us put

$$A = \{\omega \mid \text{there is a boy in } \omega\},$$

$$B = \{\omega \mid \text{there are two boys in } \omega\}.$$

Then $B \subset A$ and so AB = B, thus

$$P(B \mid A) = \frac{P(B)}{P(A)} = \frac{\frac{1}{4}}{\frac{3}{4}} = \frac{1}{3}.$$

This is the correct answer to the question. But now let us ask a similar sounding but really different question. If a child is chosen at random from these families and is found to be a boy, what is the probability that the other child in his family is also a boy? This time the appropriate representation of the sample space should be

$$\tilde{\Omega} = \{g_a, g_b, b_a, b_b\},\$$

where the sample points are not families but the children of these families, and $g_g = a$ girl who has a sister, $g_b = a$ girl who has a brother, etc. Now we have

$$\begin{split} \tilde{A} &= \{\tilde{\omega} \mid \tilde{\omega} \text{ is a boy}\}, \\ \tilde{B} &= \{\tilde{\omega} \mid \tilde{\omega} \text{ has a brother}\}, \end{split}$$

so that

$$\tilde{A}\tilde{B} = \{\tilde{\omega} \mid \tilde{\omega} = b_b\}.$$

Therefore,

$$P(\tilde{B} \mid \tilde{A}) = \frac{P(\tilde{A}\tilde{B})}{P(\tilde{A})} = \frac{\frac{1}{4}}{\frac{1}{2}} = \frac{1}{2}.$$

This is a wonderful and by no means artificial illustration of the importance of understanding "what we are sampling" in statistics.

5.2. Basic formulas

Generally speaking, most problems of probability have to do with several events or random variables and it is their mutual relation or joint action that must be investigated. In a sense all probabilities are conditional because nothing happens in a vacuum. We omit the stipulation of conditions that are implicit or taken for granted, or if we feel that they are irrelevant to the situation in hand. For instance, when a coin is tossed we usually ignore the possibility that it will stand on its edge, and do not even specify whether it is Canadian or American. The probability that a certain candidate will win an election is certainly conditioned on his surviving the campaign—an assumption that has turned out to be premature in recent American history.

Let us begin by a few simple but fundamental propositions involving conditional probabilities: **Proposition 1.** For arbitrary events A_1, A_2, \ldots, A_n , we have

$$P(A_1 A_2 \dots A_n) = P(A_1) P(A_2 \mid A_1) P(A_3 \mid A_1 A_2) \dots P(A_n \mid A_1 A_2 \dots A_{n-1})$$
(5.2.1)

provided $P(A_1 A_2 \dots A_{n-1}) > 0$.

Proof: Under the proviso, all conditional probabilities in (5.2.1) are well defined since

$$P(A_1) \ge P(A_1 A_2) \ge \dots \ge P(A_1 A_2 \dots A_{n-1}) > 0.$$

Now the right side of (5.2.1) is explicitly

$$\frac{P(A_1)}{P(\Omega)} \frac{P(A_1 A_2)}{P(A_1)} \frac{P(A_1 A_2 A_3)}{P(A_1 A_2)} \cdots \frac{P(A_1 A_2 \cdots A_n)}{P(A_1 A_2 \cdots A_{n-1})},$$

which reduces to the left side by successive cancellations. Q.E.D.

By contrast with the additivity formula (2.3.3) for a disjoint union, the formula (5.2.1) may be called the *general multiplicative formula* for the probability of an intersection. But observe how the conditioning events are also "multiplied" step by step. A much simpler formula has been given in §2.4 for independent events. As an important application of (5.2.1), suppose the random variable $X_1, X_2, \ldots, X_n, \ldots$ are all countably valued; this is surely the case when Ω is countable. Now for arbitrary possible values $x_1, x_2, \ldots, x_n, \ldots$, we put

$$A_k = \{X_k = x_k\}, \quad k = 1, 2, \dots,$$

and obtain

$$P\{X_1 = x_1; X_2 = x_2; \dots; X_n = x_n\}$$

$$= P\{X_1 = x_1\} P\{X_2 = x_2 \mid X_1 = x_1\} P\{X_3 = x_3 \mid X_1 = x_1, X_2 = x_2\}$$

$$\cdots P\{X_n = x_n \mid X_1 = x_1, \dots, X_{n-1} = x_{n-1}\}.$$
(5.2.2)

The first term above is called the *joint probability* of X_1, X_2, \ldots, X_n ; so the formula expresses this by successive conditional probabilities. Special cases of this will be discussed later.

Proposition 2. Suppose that

$$\Omega = \sum_{n} A_n$$

is a partition of the sample space into disjoint sets. Then for any set B we have

$$P(B) = \sum_{n} P(A_n)P(B \mid A_n).$$
 (5.2.3)

Proof: First we write

$$B = \Omega B = \left(\sum_{n} A_{n}\right) B = \sum_{n} A_{n} B$$

by simple set theory, in particular (1.3.6); then we deduce

$$P(B) = P\left(\sum_{n} A_n B\right) = \sum_{n} P(A_n B)$$

by countable additivity of P. Finally we substitute

$$P(A_n B) = P(A_n)P(B \mid A_n)$$

from the definition (5.1.3). This establishes (5.2.3); note that if $P(A_n) = 0$ for some n, the corresponding term in the sum there may be taken to be 0 even though $P(B \mid A_n)$ is undefined. Q.E.D.

From now on we shall adopt the convention that $x \cdot 0 = 0$ if x is undefined, in order to avoid repetition of such remarks as in the preceding sentence.

The formula (5.2.3) will be referred to as that of total probability. Here is a useful interpretation. Suppose that the event B may occur under a number of mutually exclusive circumstances (or "causes"). Then the formula shows how its "total probability" is compounded from the probabilities of the various circumstances, and the corresponding conditional probabilities figured under the respective hypotheses.

Suppose X and Y are two integer-valued random variables and k is an integer. If we apply (5.2.3) to the sets

$$A_n = \{X = n\}, B = \{Y = k\},\$$

we obtain

$$P(Y = k) = \sum_{n} P(X = n)P(Y = k \mid X = n)$$
 (5.2.4)

where the sum is over all integers n, and if P(X = n) = 0 the corresponding term may be taken to be 0. It is easy to generalize the formula when Xtakes values in any countable range, and when "Y = k" is replaced by, e.g., " $a \leq Y \leq b$ " for a more general random variable, not necessarily taking integer values.

Proposition 3. Under the assumption and notation of Proposition 2, we have also

$$P(A_n \mid B) = \frac{P(A_n)P(B \mid A_n)}{\sum_{n} P(A_n)P(B \mid A_n)}$$
 (5.2.5)

provided P(B) > 0.

Proof: The denominator above is equal to P(B) by Proposition 2, so the equation may be multiplied out to read

$$P(B)P(A_n \mid B) = P(A_n)P(B \mid A_n).$$

This is true since both sides are equal to $P(A_nB)$. Q.E.D.

This simple proposition with an easy proof is very famous under the name of Bayes' Theorem, published in 1763. It is supposed to yield an "inverse probability," or probability of the "cause" A, on the basis of the observed "effect" B. Whereas $P(A_n)$ is the a priori, $P(A_n \mid B)$ is the a posteriori probability of the cause A_n . Numerous applications were made in all areas of natural phenomena and human behavior. For instance, if B is a "body" and the A_n 's are the several suspects of the murder, then the theorem will help the jury or court to decide the whodunit. [Jurisprudence was in fact a major field of early speculations on probability. If B is an earthquake and the A_n 's are the different physical theories to explain it, then the theorem will help the scientists to choose between them. Laplace [1749– 1827; one of the great mathematicians of all time who wrote a monumental treatise on probability around 1815 used the theorem to estimate the probability that "the sun will also rise tomorrow" (see Example 9 ahead). In modern times Bayes lent his name to a school of statistics. For our discussion here let us merely comment that Bayes has certainly hit upon a remarkable turnaround for conditional probabilities, but the practical utility of his formula is limited by our usual lack of knowledge on the various a priori probabilities.

The following simple examples are given to illustrate the three propositions above. Others will appear in the course of our work.

Example 6. We have actually seen several examples of Proposition 1 before in Chapter 3. Let us reexamine them using the new notion.

What is the probability of throwing six perfect die and getting six different faces? [See Example 2 of §3.1.] Number the dice from 1 to 6, and

put

 $A_1 =$ any face for Die 1,

 $A_2 = \text{Die } 2 \text{ shows a different face from Die } 1$,

 $A_3 = \text{Die } 3 \text{ shows a different face from Die 1 and Die 2},$

etc. Then we have, assuming that the dice act independently,

$$P(A_1) = 1$$
, $P(A_2 \mid A_1) = \frac{5}{6}$, $P(A_3 \mid A_1 A_2) = \frac{4}{6}$, ..., $P(A_6 \mid A_1 A_2 \cdots A_5) = \frac{1}{6}$.

Hence an application of Proposition 1 gives

$$P(A_1 A_2 \cdots A_6) = \frac{6}{6} \cdot \frac{5}{6} \cdot \frac{4}{6} \cdot \frac{3}{6} \cdot \frac{2}{6} \cdot \frac{1}{6} = \frac{6!}{6^6}.$$

The birthday problem [Problem 5 of §3.4] is now seen to be practically the same problem, where the number 6 above is replaced by 365. The sequential method mentioned there is just another case of Proposition 1.

Example 7. The family dog is missing after the picnic. Three hypotheses are suggested:

- (A) it has gone home;
- (B) it is still worrying that big bone in the picnic area;
- (C) it has wandered off into the woods.

The *a priori* probabilities, which are assessed from the habits of the dog, are estimated respectively to be $\frac{1}{4}$, $\frac{1}{2}$, $\frac{1}{4}$. A child each is sent back to the picnic ground and the edge of the woods to look for the dog. If it is in the former area, it is a cinch (90%) that it will be found; if it is in the latter, the chance is only a toss-up (50%). What is the probability that the dog will be found in the park?

Let A, B, C be the hypotheses above, and let D = "dog will be found in the park." Then we have the following data:

$$P(A) = \frac{1}{4}, \qquad P(B) = \frac{1}{2}, \qquad P(C) = \frac{1}{4};$$

$$P(D \mid A) = 0, \ P(D \mid B) = \frac{90}{100}, \quad P(D \mid C) = \frac{50}{100}.$$

Hence by (5.2.3),

$$P(D) = P(A)P(D \mid A) + P(B)P(D \mid B) + P(C)P(D \mid C)$$
$$= \frac{1}{4} \cdot 0 + \frac{1}{2} \cdot \frac{90}{100} + \frac{1}{4} \cdot \frac{50}{100} = \frac{115}{200}.$$

What is the probability that the dog will be found at home? Call this D', and assume that $P(D' \mid A) = 1$, namely that if it is home it will be there to greet the family. Clearly $P(D' \mid B) = P(D' \mid C) = 0$ and so

$$P(D') = P(A)P(D' \mid A) + P(B)P(D' \mid B) + P(C)P(D' \mid C)$$
$$= \frac{1}{4} \cdot 1 + \frac{1}{2} \cdot 0 + \frac{1}{4} \cdot 0 = \frac{1}{4}.$$

What is the probability that the dog is "lost"? It is

$$1 - P(D) - P(D') = \frac{35}{200}.$$

Example 8. Urn one contains 2 black and 3 red balls; urn two contains 3 black and 2 red balls. We toss an unbiased coin to decide on the urn to draw from but we do not know which is which. Suppose the first ball drawn is black and it is put back; what is the probability that the second ball drawn from the same urn is also black?

Call the two urns U_1 and U_2 ; the a priori probability that either one is chosen by the coin-tossing is 1/2:

$$P(U_1) = \frac{1}{2}, \quad P(U_2) = \frac{1}{2}.$$

Denote the event that the first ball is black by B_1 , that the second ball is black by B_2 . We have by (5.2.5)

$$P(U_1 \mid B_1) = \frac{\frac{1}{2} \cdot \frac{2}{5}}{\frac{1}{2} \cdot \frac{2}{5} + \frac{1}{2} \cdot \frac{3}{5}} = \frac{2}{5}, \quad P(U_2 \mid B_1) = \frac{3}{5}.$$

Note that the two probabilities must add up to 1 (why?) so we need only compute one of them. Note also that the two a posteriori probabilities are directly proportional to the probabilities $P(B_1 \mid U_1)$ and $P(B_1 \mid U_2)$. That is, the black ball drawn is more likely to have come from the urn that is more likely to yield a black ball, and in the proper ratio. Now use (5.2.3) to compute the probability that the second ball is also black. Here $A_1 = "B_1$ is from U_1 ," $A_2 = "B_1$ is from U_2 " are the two alternative hypotheses. Since the second drawing is conditioned on B_1 , the probabilities of the hypotheses are really conditional ones:

$$P(A_1) = P(U_1 \mid B_1) = \frac{2}{5}, \quad P(A_2) = P(U_2 \mid B_1) = \frac{3}{5}.$$

On the other hand, it is obvious that

$$P(B_2 \mid A_1) = \frac{2}{5}, \quad P(B_2 \mid A_2) = \frac{3}{5}.$$

Hence we obtain the conditional probability

$$P(B_2 \mid B_1) = \frac{2}{5} \cdot \frac{2}{5} + \frac{3}{5} \cdot \frac{3}{5} = \frac{13}{25}.$$

Compare this with

$$P(B_2) = P(U_1)P(B_2 \mid U_1) + P(U_2)P(B_2 \mid U_2) = \frac{1}{2} \cdot \frac{2}{5} + \frac{1}{2} \cdot \frac{3}{5} = \frac{1}{2}.$$

Note that $P(B_2 \mid U_1) = P(B_1 \mid U_1)$ – why? We see that the knowledge of the first ball drawn being black has strengthened the probability of drawing a second black ball, because it has increased the likelihood that we have picked the urn with more black balls. To proceed one more step, given that the first two balls drawn are both black and put back, what is the probability of drawing a third black ball from the same urn? We have in notation similar to the above:

$$P(U_1 \mid B_1 B_2) = \frac{\frac{1}{2} \left(\frac{2}{5}\right)^2}{\frac{1}{2} \cdot \left(\frac{2}{5}\right)^2 + \frac{1}{2} \left(\frac{3}{5}\right)^2} = \frac{4}{13}; \quad P(U_2 \mid B_1 B_2) = \frac{9}{13};$$

$$P(B_3 \mid B_1 B_2) = \frac{4}{13} \cdot \frac{2}{5} + \frac{9}{13} \cdot \frac{3}{5} = \frac{35}{65}.$$

This is greater than 13/25, so a further strengthening has occurred. Now it is easy to see that we can extend the result to any number of drawings. Thus,

$$P(U_1 \mid B_1 B_2 \cdots B_n) = \frac{\frac{1}{2} \left(\frac{2}{5}\right)^n}{\frac{1}{2} \left(\frac{2}{5}\right)^n + \frac{1}{2} \left(\frac{3}{5}\right)^n} = \frac{1}{1 + \left(\frac{3}{2}\right)^n},$$

where we have divided the denominator by the numerator in the middle term. It follows that as n becomes larger and larger, the a posteriori probability of U_1 becomes smaller and smaller. In fact it decreases to zero and consequently the a posteriori probability of U_2 increases to 1 in the limit. Thus we have

$$\lim_{n \to \infty} P(B_{n+1} \mid B_1 B_2 \cdots B_n) = \frac{3}{5} = P(B_1 \mid U_2).$$

This simple example has important implications on the empirical view-point of probability. Replace the two urns above by a coin that may be biased (as all real coins are). Assume that the probability p of heads is either 2/5 or 3/5 but we do not know which is the true value. The two possibilities are then two alternative hypotheses between which we must decide. If they both have a priori probability 1/2, then we are in the situation of the two urns. The outcome of each toss will affect our empirical estimate of the value of p. Suppose for some reason we believe that p = 2/5.

Then if the coin falls heads 10 times in a row, can we still maintain that p=2/5 and give probability $(2/5)^{10}$ to this rare event? Or shall we concede that really p = 3/5 so that the same event will have probability $(3/5)^{10}$? This is very small but still $(3/2)^{10}$ larger than the other. In certain problems of probability theory it is customary to consider the value of p as fixed and base the rest of our calculations on it. So the query is what reason do we have to maintain such a fixed stance in the face of damaging evidence given by observed outcomes? Keynes made a point of this criticism on the foundations of probability. From the axiomatic point of view, as followed in this book, a simple answer is this: our formulas are correct for each arbitrary value of p, but axioms of course do not tell us what this value is, or even whether it makes sense to assign any value at all. The latter may be the case when one talks about the probability of the existence of some "big living creatures somewhere in outer space." [It used to be the moon!] In other words, mathematics proper being a deductive science, the problem of evaluating, estimating, or testing the value of p lies outside its eminent domain. Of course, it is of the utmost importance in practice, and statistics was invented to cope with this kind of problem. But it need not concern us too much here. [The author had the authority of Dr. Albert Einstein on this point, while on a chance stroll on Mercer Street in Princeton, N.J., sometime in 1946 or 1947. Here is the gist of what he said: in any branch of science which has applications, there is always a gap, which needs a bridge between theory and practice. This is so for instance in geometry or mechanics; and probability is no exception.

The preceding example has a natural extension when the unknown p may take values in a finite or infinite range. Perhaps the most celebrated illustration is Laplace's law of succession below.

Example 9. Suppose that the sun has risen n times in succession; what is the probability that it will rise once more?

It is assumed that the a priori probability for a sunrise on any day is a constant whose value is unknown to us. Due to our total ignorance it will be assumed to take all possible values in [0,1] with equal likelihood. That is to say, this probability will be treated as a random variable ξ that is uniformly distributed over [0,1]. Thus ξ has the density function f such that f(p) = 1 for $0 \le p \le 1$. This can be written heuristically as

$$P(p \le \xi \le p + dp) = dp, \quad 0 \le p \le 1.$$
 (5.2.6)

See the discussion in Example 10 in §4.5. Now if the true value of ξ is p, then under this hypothesis the probability of n successive sunrises is equal to p^n , because they are assumed to be independent events. Let S^n denote the event that "the sun rises n times in succession"; then we may write

heuristically

$$P(S^n \mid \xi = p) = p^n.$$
 (5.2.7)

The analogue to (5.2.3) should then be

$$P(S^n) = \sum_{0 \le p \le 1} P(\xi = p) P(S^n \mid \xi = p).$$
 (5.2.8)

This is of course meaningless as it stands, but if we pass from the sum into an integral and use (5.2.6), the result is

$$P(S^n) = \int_0^1 P(S^n \mid \xi = p) \, dp = \int_0^1 p^n \, dp = \frac{1}{n+1}.$$
 (5.2.9)

This continuous version of (5.2.3) is in fact valid, although its derivation above is not quite so. Accepting the formula and applying it for both n and n+1, then taking the ratio, we obtain

$$P(S^{n+1} \mid S^n) = \frac{P(S^n S^{n+1})}{P(S^n)} = \frac{P(S^{n+1})}{P(S^n)} = \frac{\frac{1}{n+2}}{\frac{1}{n+1}} = \frac{n+1}{n+2}.$$
 (5.2.10)

This is Laplace's answer to the sunrise problem.

In modern parlance, Laplace used an "urn model" to study successive sunrise as a random process. A sunrise is assimilated to the drawing of a black ball from an urn of unknown composition. The various possible compositions are assimilated to so many different urns containing various proportions of black balls. Finally, the choice of the true value of the proportion is assimilated to the picking of a random number in [0, 1]. Clearly, these are weighty assumptions calling forth serious objections at several levels. Is sunrise a random phenomenon or is it deterministic? Assuming that it can be treated as random, is the preceding simple urn model adequate to its description? Assuming that the model is appropriate in principle, why should the a priori distribution of the true probability be uniformly distributed, and if not how could we otherwise assess it?

Leaving these great questions aside, let us return for a moment to (5.2.7). Since $P(\xi=p)=0$ for every p (see §4.5 for a relevant discussion), the so-called conditional probability in that formula is *not* defined by (5.1.3). Yet it makes good sense from the interpretation given before (5.2.7). In fact, it can be made completely legitimate by a more advanced theory [Radon–Nikodym derivative]. Once this is done, the final step (5.2.9) follows without the intervention of the heuristic (5.2.8). Although a full explanation of these matters lies beyond the depth of this textbook, it seems

proper to mention it here as a natural extension of the notion of conditional probability. A purely discrete approach to Laplace's formula is also possible, but the calculations are harder (see Exercise 35 below).

We end this section by introducing the notion of conditional expectation. In a countable sample space consider a random variable Y with range $\{y_k\}$ and an event S with P(S) > 0. Suppose that the expectation of Y exists; then its conditional expectation relative to S is defined to be

$$E(Y \mid S) = \sum_{k} y_{k} P(Y = y_{k} \mid S).$$
 (5.2.11)

Thus, we simply replace in the formula $E(Y) = \sum_k y_k P(Y = y_k)$ the probabilities by conditional ones. The series in (5.2.11) converges absolutely because the last-written series does so. In particular, if X is another random variable with range $\{x_j\}$, then we may take $S = \{X = x_j\}$ to obtain $E(Y \mid X = x_j)$. On the other hand, we have as in (5.2.4)

$$P(Y = y_k) = \sum_{j} P(X = x_j) P(Y = y_k \mid X = x_j).$$

Multiplying through by y_k , summing over k, and rearranging the double series, we obtain

$$E(Y) = \sum_{j} P(X = x_j) E(Y \mid X = x_j).$$
 (5.2.12)

The rearrangement is justified by absolute convergence.

The next two sections contain somewhat special material. The reader may read the beginnings of $\S\S5.3$ and 5.4 up to the statements of Theorems 1 and 3 to see what they are about, but postpone the rest and go to $\S5.5$.

*5.3. Sequential sampling

In this section we study an urn model in some detail. It is among the simplest schemes that can be handled by elementary methods. Yet it presents rich ideas involving conditioning which are important in both theory and practice.

An urn contains b black balls and r red balls. One ball is drawn at a time without replacement. Let $X_n=1$ or 0 depending on whether the nth ball drawn is black or red. Each sample point ω is then just the sequence $\{X_1(\omega), X_2(\omega), \ldots, X_{b+r}(\omega)\}$, briefly $\{X_n, 1 \leq n \leq b+r\}$; see the discussion around (4.1.3). Such a sequence is called a *stochastic process*, which is a fancy name for any family of random variables. [According to the dictionary, "stochastic" comes from a Greek word meaning "to aim at."] Here

the family is the finite sequence indexed by n from 1 to b+r. This index n may be regarded as a *time parameter* as if one drawing is made per unit time. In this way we can speak of the gradual evolution of the process as time goes on by observing the successive X_n 's.

You may have noticed that our model is nothing but sampling without replacement and with ordering, discussed in §3.2. You are right but our viewpoint has changed and the elaborate description above is meant to indicate this. Not only do we want to know, e.g., how many black balls are drawn after so many drawings, as we would previously, but now we want also to know how the sequential drawings affect each other, how the composition of the urn changes with time, etc. In other words, we want to investigate the mutual dependence of the X_n 's, and that's where conditional probabilities come in. Let us begin with the easiest kind of question.

Problem. A ball is drawn from the urn and discarded. Without knowing its color, what is the probability that a second ball drawn is black?

For simplicity let us write the events $\{X_n = 1\}$ as B_n and $\{X_n = 0\}$ as $R_n = B_n^c$. We then have from Proposition 2 of §5.2,

$$P(B_2) = P(B_1)P(B_2 \mid B_1) + P(B_1^c)P(B_2 \mid B_1^c).$$
 (5.3.1)

Clearly we have

$$P(B_1) = \frac{b}{b+r}, \quad P(B_1^c) = \frac{r}{b+r},$$
 (5.3.2)

whereas

$$P(B_2 \mid B_1) = \frac{b-1}{b+r-1}, \quad P(B_1 \mid B_1^c) = \frac{b}{b+r-1}$$

since there are b+r-1 balls left in the urn after the first drawing, and among these are b-1 or b black balls according to whether the first ball drawn is or is not black. Substituting into (5.3.1) we obtain

$$P(B_2) = \frac{b}{b+r} \frac{b-1}{b+r-1} + \frac{r}{b+r} \frac{b}{b+r-1} = \frac{b(b+r-1)}{(b+r)(b+r-1)} = \frac{b}{b+r}.$$

Thus $P(B_2) = P(B_1)$; namely if we take into account both possibilities for the color of the first ball, then the probabilities for the second ball are the same as if no ball had been drawn (and left out) before. Is this surprising or not? Anyone with curiosity would want to know whether this result is an accident or has a theory behind it. An easy way to test this is to try another step or two: suppose 2 or 3 balls have been drawn but their colors not noted; what then is the probability that the next ball will be black? You should carry out the simple computations by all means. The general result can be stated succinctly as follows.

Theorem 1. We have for each n

$$P(B_n) = \frac{b}{b+r}, \quad 1 \le n \le b+r.$$
 (5.3.3)

It is essential to pause here and remark on the economy of this mathematical formulation, in contrast to the verbose verbal description above. The condition that "we do not know" the colors of the n-1 balls previously drawn is observed as it were in silence, namely by the absence of conditioning for the probability $P(B_n)$. What should we have if we know the colors? It would be something like $P(B_2 \mid B_1)$ or $P(B_3 \mid B_1B_2^c)$ or $P(B_4 \mid B_1B_2^cB_3)$. These are trivial to compute (why?); but we can also have something like $P(B_4 \mid B_2)$ or $P(B_4 \mid B_1B_3^c)$, which is slightly less trivial. See Exercise 33.

There are many different ways to prove the beautiful theorem above; each method has some merit and is useful elsewhere. We will give two now, a third one in a tremendously more general form (Theorem 5 in §5.4) later. But there are others and perhaps you can think of one later. The first method may be the toughest for you; if so skip it and go at once to the second.*

First Method. This may be called "direct confrontation" or "brute force" and employs heavy (though standard) weaponry from the combinatory arsenal. Its merit lies in that it is bound to work provided that we have guessed the answer in advance, as we can in the present case after a few trials. In other words, it is a sort of experimental verification. We introduce a new random variable $Y_n =$ the number of black balls drawn in the first n drawings. This gives the proportion of black balls when the n+1st drawing is made since the total number of balls then is equal to b+r-n, regardless of the outcomes of the previous n drawings. Thus we have

$$P(B_{n+1} \mid Y_n = j) = \frac{b-j}{b+r-n}, \quad 0 \le j \le b.$$
 (5.3.4)

On the other hand, the probability $P(Y_n = j)$ can be computed as in Problem 1 of §3.4, with m = b + r, k = b in (3.4.1):

$$P(Y_n = j) = \frac{\binom{b}{j} \binom{r}{n-j}}{\binom{b+r}{n}}.$$
 (5.3.5)

^{*}A third method is to make mathematical induction on n.

We now apply (5.2.4):

$$P(B_{n+1}) = \sum_{j=0}^{b} P(Y_n = j) P(B_{n+1} \mid Y_n = j)$$

$$= \sum_{j=0}^{b} \frac{\binom{b}{j} \binom{r}{n-j}}{\binom{b+r}{n}} \frac{b-j}{b+r-n}.$$
(5.3.6)

This will surely give the answer, but how in the world are we going to compute a sum like that? Actually it is not so hard, and there are excellent mathematicians who make a career out of doing such (and much harder) things. The beauty of this kind of computation is that it's got to unravel if our guess is correct. This faith lends us strength. Just write out the several binomial coefficients above explicitly, cancelling and inserting factors with a view to regrouping them into new binomial coefficients:

$$\frac{b!}{j! (b-j)!} \frac{r!}{(n-j)! (r-n+j)!} \frac{n! (b+r-n)!}{(b+r)!} \frac{b-j}{b+r-n}$$

$$= \frac{b! r!}{(b+r)!} \frac{(b+r-n-1)!}{(r-n+j)! (b-j-1)!} \frac{n!}{j! (n-j)!}$$

$$= \frac{1}{\binom{b+r}{b}} \binom{b+r-n-1}{b-j-1} \binom{n}{j}.$$

Hence

$$P(B_{n+1}) = \frac{1}{\binom{b+r}{b}} \sum_{j=0}^{b-1} \binom{n}{j} \binom{b+r-1-n}{b-1-j},$$
 (5.3.7)

where the term corresponding to i = b has been omitted since it yields zero in (5.3.6). The new sum (5.3.7) is a well-known identity for binomial coefficients and is equal to $\binom{b+r-1}{b-1}$; see (3.3.9). Thus

$$P(B_{n+1}) = {b+r-1 \choose b-1} / {b+r \choose b} = \frac{b}{b+r}$$

as asserted in (5.3.3).

Second Method. This is purely combinatorial and can be worked out as an example in §3.2. Its merit is simplicity; but it cannot be easily generalized to apply to the next urn model we shall consider.

Consider the successive outcomes in n+1 drawings: $X_1(\omega), X_2(\omega), \ldots$, $X_n(\omega), X_{n+1}(\omega)$. Each $X_j(\omega)$ is 1 or 0 depending on the particular ω ; even the numbers of 1's and 0's among them depend on ω when n+1 < b+r. Two different outcome sequences such as 0011 and 0101 will not have the same probability in general. But now let us put numerals on the balls, say 1 to b for the black ones and b+1 to b+r for the red ones, so that all balls become distinguishable. We are then in the case of sampling without replacement and with ordering discussed in §3.2. The total number of possibilities with the new labeling is given by (3.2.1) with b+r for m and m+1 for m: $(b+r)_{m+1}$. These are now all equally likely! We are interested in the cases where the m+1st ball is black; how many are there for these? There are b choices for the n+1st ball, and after this is chosen there are $(b+r-1)_n$ ways of arranging the first n balls, by another application of (3.2.1). Hence by the fundamental rule in §3.1, the number of cases where the n+1st ball is black is equal to $b(b+r-1)_n$. Now the classical ratio formula for probability applies to yield the answer

$$P(B_{n+1}) = \frac{b(b+r-1)_n}{(b+r)_{n+1}} = \frac{b}{b+r}.$$

Undoubtedly this argument is easier to follow after it is explained, and there is little computation. But it takes a bit of perception to hit upon the counting method. Poisson [1781–1840; French mathematician for whom a distribution, a process, a limit theorem, and an integral were named, among other things] gave this solution, but his explanation is briefer than ours. We state his general result as follows.

Theorem 2 (Poisson's Theorem). Suppose in an urn containing b black and r red balls, n balls have been drawn first and discarded without their colors being noted. If m balls are drawn next, the probability that there are k black balls among them is the same as if we had drawn these m balls at the outset [without having discarded the n balls previously drawn].

Briefly stated: the probabilities are not affected by the preliminary drawing so long as we are in the dark as to what those outcomes are. Obviously if we know the colors of the balls discarded, the probabilities will be affected in general. To quote [Keynes, p. 349]: "This is an exceedingly good example ... that a probability cannot be influenced by the occurrence of a material event but only by such knowledge as we may have, respecting the occurrence of the event."

Here is Poisson's quick argument: if n+m balls are drawn out, the probability of a combination made up of n black and red balls in given proportions followed by m balls of which k are black and m-k are red must be the same as that of a similar combination in which the m balls precede the n balls. Hence the probability of k black balls in m drawings given that n balls have already been drawn out must be equal to the probability of the same result when no balls have been previously drawn out.

Is this totally convincing to you? The more explicit combinatorial argument given above for the case m=1 can be easily generalized to settle any doubt. The doubt is quite justified despite Poisson's authority. As we may learn from Chapter 3, in these combinatorial arguments one must do one's own thinking.

***5.4.** Pólva's urn scheme

To pursue the discussion in the preceding section a step further, we will study a famous generalization due to G. Pólya [1887–1986; professor at Stanford University, one of the most eminent analysts of modern times who also made major contributions to probability and combinatorial theories and their applications]. As before the urn contains b black and r red balls to begin with, but after a ball is drawn each time, it is returned to the urn and c balls of the same color are added to the urn, where c is an integer, and when c < 0 adding c balls means subtracting -c balls. This may be done whether we observe the color of the ball drawn or not; in the latter case, e.g., we may suppose that it is performed by an automaton. If c=0 this is just sampling with replacement, while if c=-1 we are in the situation studied in §5.3. In general if c is negative the process has to stop after a number of drawings, but if c is zero or positive it can be continued forever. This scheme can be further generalized (you know generalization is a mathematician's bug!) if after each drawing we add to the urn not only c balls of the color drawn but also d balls of the other color. But we will not consider this, and furthermore we will restrict ourselves to the case $c \ge -1$, referring to the scheme as Pólya's urn model. Pólya actually invented this model to study a problem arising in medicine; see the last paragraph of this section.

Problem. What is the probability that in Pólya's model the first three balls drawn have colors $\{b, b, r\}$ in this order? or $\{b, r, b\}$? or $\{r, b, b\}$?

An easy application of Proposition 1 in §5.2 yields, in the notation introduced in §5.3,

$$P(B_1B_2R_3) = P(B_1)P(B_2 \mid B_1)P(R_3 \mid B_1B_2)$$

$$= \frac{b}{b+r} \frac{b+c}{b+r+c} \frac{r}{b+r+2c}.$$
(5.4.1)

Similarly,

$$P(B_1 R_2 B_3) = \frac{b}{b+r} \frac{r}{b+r+c} \frac{b+c}{b+r+2c},$$

$$P(R_1 B_1 B_2) = \frac{r}{b+r} \frac{b}{b+r+c} \frac{b+c}{b+r+2c}.$$

Thus they are all the same, namely the probability of drawing 2 black and 1 red balls in three drawings does not depend on the *order* in which they are drawn. It follows that the probability of drawing 2 black and 1 red in the first three drawings is equal to three times the number on the right side of (5.4.1).

The general result is given below.

Theorem 3. The probability of drawing (from the beginning) any specified sequence of k black balls and n - k red balls is equal to

$$\frac{b(b+c)\cdots(b+(k-1)c)r(r+c)\cdots(r+(n-k-1)c)}{(b+r)(b+r+c)(b+r+2c)\cdots(b+r+(n-1)c)},$$
 (5.4.2)

for all $n \ge 1$ if $c \ge 0$; and for $0 \le n \le b+r$ if c = -1.

Proof: This is really an easy application of Proposition 1 in §5.2, but in a scrambled way. We have shown it above in the case k=2 and n=3. If you will try a few more cases with say n = 4, k = 2 or n = 5, k = 3, you will probably see how it goes in the general case more quickly than it can be explained in words. The point is: at the mth drawing, where $1 \leq m \leq n$, the denominator of the term corresponding to $P(A_m \mid A_1 A_2 \cdots A_{m-1})$ in (5.2.1) is b+r+(m-1)c, because a total of (m-1)c balls has been added to the urn by this time, no matter what balls have been drawn. Now at the first time when a black ball is drawn, there are b black balls in the urn; at the second time a black ball is drawn, the number of black balls in the urn is b+c, because one black ball has been previously drawn so c black balls have been added to the urn. This is true no matter at what time (which drawing) the second black ball is drawn. Similarly, when the third black ball is drawn, there will be b+2c black balls in the urn, and so on. This explains the k factors involving b in the numerator of (5.4.2). Now consider the red balls: at the first time a red ball is drawn, there are r red ones in the urn; at the second time a red ball is drawn, there are r + c red ones in the urn, because c red balls have been added after the first red one is drawn, and so on. This explains the n-k factors involving r (= red) in the numerator of (5.4.2). The whole thing there is therefore obtained by multiplying the successive ratios as the conditional probabilities in (5.2.1), and the exact order in which the factors in the numerator occur is determined by the specific order of blacks and reds in the given sequence. However, their product is the same so long as n and k are fixed. This establishes (5.4.2).

For instance, if the specified sequence is RBRRB, then the exact order in the numerator should be rb(r+c)(r+2c)(b+c).

Now suppose that only the number of black balls is given [specified!] but not the exact sequence; then we have the next result.

Theorem 4. The probability of drawing (from the beginning) k black balls in n drawings is equal to the number in (5.4.2) multiplied by $\binom{n}{k}$. In terms of generalized binomial coefficients [see (5.4.4) below], it is equal to

$$\frac{\binom{-\frac{b}{c}}{k}\binom{-\frac{r}{c}}{n-k}}{\binom{-\frac{b+r}{c}}{c}}.$$
 (5.4.3)

This is an extension of the hypergeometric distribution; see p. 90.

Proof: There are $\binom{n}{k}$ ways of permuting k black and n-k red balls; see §3.2. According to (5.4.2), every specified sequence of drawing k black and n-k red balls has the same probability. These various permutations correspond to disjoint events. Hence the probability stated in the theorem is just the sum of $\binom{n}{k}$ probabilities, each of which is equal to the number given in (5.4.2). It remains to express this probability by (5.4.3), which requires only a bit of algebra. Let us note that if a is a positive real number and j is a positive integer, then by definition

$${\binom{-a}{j}} = \frac{(-a)(-a-1)\cdots(-a-j+1)}{j!}$$

$$= (-1)^{j} \frac{a(a+1)\cdots(a+j-1)}{j!}.$$
(5.4.4)

Thus if we divide every factor in (5.4.2) by c, and write

$$\beta = \frac{b}{c}, \quad \gamma = \frac{r}{c}$$

for simplicity, then use (5.4.4), we obtain

$$\frac{\beta(\beta+1)\cdots(\beta+k-1)\gamma(\gamma+1)\cdots(\gamma+n-k-1)}{(\beta+\gamma)(\beta+\gamma+1)\cdots(\beta+\gamma+n-1)}$$

$$=\frac{(-1)^k k! \binom{-\beta}{k} (-1)^{n-k} (n-k)! \binom{-\gamma}{n-k}}{(-1)^n n! \binom{-\beta-\gamma}{n}} = \frac{\binom{-\beta}{k} \binom{-\gamma}{n-k}}{\binom{-\beta-\gamma}{n} \binom{n}{k}}.$$

After multiplying by $\binom{n}{k}$ we get (5.4.3) as asserted.

We can now give a far-reaching generalization of Theorems 1 and 2 in §5.3. Furthermore the result will fall out of the fundamental formula (5.4.2) like a ripe fruit. Only a bit of terminology and notation is in the way.

Recalling the definition of X_n in §5.3, we can record (5.4.2) as giving the *joint distribution* of the n random variable $\{X_1, X_2, \ldots, X_n\}$. Let us introduce the *hedge symbol* "①" to denote either "0" or "1" and use

subscripts (indices) to allow an arbitrary choice for each subscript, independently of each other. On the other hand, two such symbols with the same subscript must of course denote the same choice throughout a discussion. For instance, $\{ \bigcirc_1, \bigcirc_2, \bigcirc_3, \bigcirc_4 \}$ may mean $\{1,1,0,1\}$ or $\{0,1,0,1\}$, but then $\{ \bigcirc_1, \bigcirc_3 \}$ must mean $\{1,0\}$ in the first case and $\{0,0\}$ in the second. Theorem 3 can be stated as follows: if k of the ①'s below are 1's and n-k of them are 0's, then

$$P(X_1 = \bigcirc_1, X_2 = \bigcirc_2, \dots, X_n = \bigcirc_n)$$
 (5.4.5)

is given by the expression in (5.4.2). There are altogether 2^n possible choices for the ①'s in (5.4.5) [why?], and if we visualize all the resulting values corresponding to these choices, the set of 2^n probabilities determines the joint distribution of $\{X_1, X_2, \ldots, X_n\}$. Now suppose $\{n_1, n_2, \ldots, n_s\}$ is a subset of $\{1, 2, \ldots, n\}$; the joint distribution of $\{X_{n_1}, \ldots, X_{n_s}\}$ is determined by

$$P(X_{n_1} = \bigcirc_{n_1}, \dots, X_{n_s} = \bigcirc_{n_s})$$
 (5.4.6)

when the latter ①'s range over all the 2^s possible choices. This is called a marginal distribution with reference to that of the larger set $\{X_1, \ldots, X_n\}$.

We need more notation! Let $\{n'_1, \ldots, n'_t\}$ be the complementary set of $\{n_1, \ldots, n_s\}$ with respect to $\{1, \ldots, n\}$, namely those indices left over after the latter set has been taken out. Of course, t = n - s and the union $\{n_1, \ldots, n_s, n'_1, \ldots, n'_t\}$ is just some permutation of $\{1, \ldots, n\}$. Now we can write down the following formula expressing a marginal probability by means of joint probabilities of a larger set:

$$P(X_{n_1} = \textcircled{1}_1, \dots, X_{n_s} = \textcircled{1}_s)$$

$$= \sum_{\textcircled{1}'_1, \dots, \textcircled{1}'_t} P(X_{n_1} = \textcircled{1}_1, \dots, X_{n_s} = \textcircled{1}_s,$$

$$X_{n'_1} = \textcircled{1}'_1, \dots, X_{n'_t} = \textcircled{1}'_t),$$

$$(5.4.7)$$

where $\{ \textcircled{0}'_1, \dots, \textcircled{0}'_t \}$ is another set of hedge symbols and the sum is over all the 2^t choices for them. This formula follows from the obvious set relation

$$\{X_{n_1} = \textcircled{1}_1, \dots, X_{n_s} = \textcircled{1}_s\}$$

$$= \sum_{\textcircled{1}'_1, \dots, \textcircled{1}'_t} \{X_{n_1} = \textcircled{1}_1, \dots, X_{n_s} = \textcircled{1}_s X_{n_1'} = \textcircled{1}'_1, \dots, X_{n_t'} = \textcircled{1}'_t\}$$

and the additivity of P. [Clearly a similar relation holds when the X's take other values than 0 or 1, in which case the ①'s must be replaced by all possible values.]

We now come to the *pièce de résistance* of this discussion. It will sorely test your readiness to digest a general and abstract argument. If you can't

swallow it now, you need not be upset – but do come back and try it again later.

Theorem 5. The joint distribution of any s of the random variables $\{X_1, X_2, \ldots, X_n, \ldots\}$ is the same, no matter which s of them is in question.

As noted above, the sequence of X_n 's is infinite if $c \ge 0$, whereas $n \le b + r$ if c = -1.

Proof: What does the theorem say? Fix s and let X_{n_1}, \ldots, X_{n_s} be any set of s random variables chosen from the entire sequence. To discuss its joint distribution, we must consider all possible choices of values for these s random variables. So we need a notation for an arbitrary choice of that kind, call it $\mathfrak{O}_1, \ldots, \mathfrak{O}_s$. Now let us write

$$P(X_{n_1} = \bigcirc_1, X_{n_2} = \bigcirc_2, \dots, X_{n_s} = \bigcirc_s).$$

We must show that this has the same value no matter what $\{n_1, \ldots, n_s\}$ is, namely that it has the same value as

$$P(X_{m_1} = \bigcirc_1, X_{m_2} = \bigcirc_2, \dots, X_{m_s} = \bigcirc_s),$$

where $\{m_1, \ldots, m_s\}$ is any other subset of size s. The two sets $\{n_1, \ldots, n_s\}$ and $\{m_1, \ldots, m_s\}$ may very well be overlapping, such as $\{1, 3, 4\}$ and $\{3, 2, 1\}$. Note also that we have never said that the indices must be in increasing order!

Let the maximum of the indices used above be n. As before let t=n-s, and

$$\{n'_1, \dots, n'_t\} = \{1, \dots, n\} - \{n_1, \dots, n_s\},$$

$$\{m'_1, \dots, m'_t\} = \{1, \dots, n\} - \{m_1, \dots, m_s\}.$$

Next, let $\mathbb{O}'_1, \ldots, \mathbb{O}'_t$ be an arbitrary choice of t hedge symbols. We claim then

$$P(X_{n_1} = \textcircled{1}_1, \dots, X_{n_s} = \textcircled{1}_s, X_{n'_1} = \textcircled{1}'_1, \dots, X_{n'_t} = \textcircled{1}'_t)$$

$$= P(X_{m_1} = \textcircled{1}_1, \dots, X_{m_s} = \textcircled{1}_s, X_{m'_1} = \textcircled{1}_1, \dots, X_{m_t} = \textcircled{1}'_t).$$
(5.4.8)

If you can read this symbolism you will see that it is just a consequence of (5.4.2)! For both $(n_1, \ldots, n_s, n'_1, \ldots, n'_t)$ and $(m_1, \ldots, m_s, m'_1, \ldots, m'_t)$ are permutations of the whole set $(1, \ldots, n)$, whereas the set of hedge symbols $(\widehat{0}_1, \ldots, \widehat{0}_s, \widehat{0}'_1, \ldots, \widehat{0}'_t)$ are the same on both sides of (5.4.8). So the equation merely repeats the assertion of Theorem 3 that any two

specified sequences having the same number of black balls must have the same probability, irrespective of the permutations.

Finally, keeping $\bigcirc_1, \ldots, \bigcirc_s$ fixed but letting $\bigcirc_1', \ldots, \bigcirc_t'$ vary over all 2^t possible choices, we get 2^t equations of the form (5.4.8). Take their sum and use (5.4.7) once as written and another time when the n's are replaced by the m's. We get

$$P(X_{n_1} = \bigcirc_1, \dots, X_{n_s} = \bigcirc_s) = P(X_{m_1} = \bigcirc_1, \dots, X_{m_s} = \bigcirc_s),$$

as we set out to show. Q.E.D.

There is really nothing hard or tricky about this proof. "It's just the *notation*!", as some would say.

A sequence of random variables $\{X_n; n=1,2,...\}$ having the property given in Theorem 5 is said to be "permutable" or "exchangeable." It follows in particular that any *block* of given length s, such as $X_{s_0+1}, X_{s_0+2}, \ldots, X_{s_0+s}$, where s_0 is any nonnegative integer (and $s_0 + s \le b + r$ if c = -1), has the same distribution. Since the index is usually interpreted as the *time parameter*, the distribution of such a block may be said to be "invariant under a time shift." A sequence of random variables having this property is said to be "[strictly] stationary." This kind of process is widely used as a model in electrical oscillations, economic time series, queuing problems, etc.

Pólya's scheme may be considered as a model for a fortuitous happening [a "random event" in the everyday usage] whose likelihood tends to increase with each occurrence and decrease with each non occurrence. The drawing of a black ball from his urn is such an event. Pólya himself cited as an example the spread of an epidemic in which each victim produces many more new germs and so increases the chances of further contamination. To quote him directly (my translation from the French original), "In reducing this fact to its simplest terms and adding to it a certain symmetry, propitious for mathematical treatment, we are led to the urn scheme." The added symmetry refers to the adding of red balls when a red ball is drawn, which would mean that each nonvictim also increases the chances of other nonvictims. This half of the hypothesis for the urn model does not seem to be warranted and is slipped in without comment by several authors who discussed it. Professor Pólya's candor, in admitting it as a mathematical expediency, should be reassuring to scientists who invented elaborate mathematical theories to deal with crude realities such as hens pecking (mathematical psychology) and beetles crawling (mathematical biology).

5.5. Independence and relevance

An extreme and extremely important case of conditioning occurs when the condition has no effect on the probability. This intuitive notion is common

experience in tossing a coin or throwing a die several times, or drawing a ball several times from an urn with replacement. The knowledge of the outcome of the previous trials should not change the "virgin" probabilities of the next trial and in this sense the trials are intuitively independent of each other. We have already defined independent events in §2.4; observe that the defining relations in (2.4.5) are just special cases of (5.2.1) when all conditional probabilities are replaced by unconditional ones. The same replacement (5.2.2) will now lead to the fundamental definition below.

Definition of Independent Random Variables. The countably valued random variables X_1, \ldots, X_n are said to be independent iff for any real numbers x_1, \ldots, x_n , we have

$$P(X_1 = x_1, \dots, X_n = x_n) = P(X_1 = x_1) \dots P(X_n = x_n).$$
 (5.5.1)

This equation is trivial if one of the factors on the right is equal to zero; hence we may restrict the x's above to the countable set of all possible values of all the X's.

The deceptively simple condition (5.5.1) actually contains much more than meets the eye. To see this let us deduce at once a major extension of (5.5.1) in which single values x_i are replaced by arbitrary sets S_i . Let X_1, \ldots, X_n be independent random variables in Propositions 4 to 6 below.

Proposition 4. We have for arbitrary countable sets S_1, \ldots, S_n :

$$P(X_1 \in S_1, \dots, X_n \in S_n) = P(X_1 \in S_1) \dots P(X_n \in S_n).$$
 (5.5.2)

Proof: The left member of (5.5.2) is equal to

$$\sum_{x_1 \in S_1} \cdots \sum_{x_n \in S_n} P(X_1 = x_1, \dots, X_n = x_n)$$

$$= \sum_{x_1 \in S_1} \cdots \sum_{x_n \in S_n} P(X_1 = x_1) \dots P(X_n = x_n)$$

$$= \left\{ \sum_{x_1 \in S_1} P(X_1 = x_1) \right\} \cdots \left\{ \sum_{x_n \in S_n} P(X_n = x_n) \right\},$$

which is equal to the right member of (5.5.2) by simple algebra (which you should spell out if you have any doubt).

Note that independence of a set of random variables as defined above is a property of the set as a whole. Such a property is not necessarily inherited by a subset; can you think of an easy counterexample? However, as a consequence of Proposition 4, any subset of (X_1, \ldots, X_n) is indeed also a set of independent random variables. To see, e.g., (X_1, X_2, X_3) is

such a set when n > 3 above, we take $S_i = R^1$ for i > 3 and replace the other S_i 's by x_i in (5.5.2).

Next, the condition (5.5.2) will be further strengthened into its most useful form.

Proposition 5. The events

$$\{X_1 \in S_1\}, \dots, \{X_n \in S_n\}$$
 (5.5.3)

are independent.

Proof: It is important to recall that the definition of independent events requires not only the relation (5.5.2), but also similar relations for all subsets of (X_1, \ldots, X_n) . However, these also hold because the subsets are also sets of independent random variables, as just shown.

Before going further let us check that the notion of independent events defined in §2.4 is a special case of independent random variables defined in this section. With the arbitrary events $\{A_j, 1 \leq j \leq n\}$ we associate their indicators I_{A_i} (see §1.4), where

$$I_{A_i}(\omega) = \begin{cases} 1 & \text{if } \omega \in A_j, \\ 0 & \text{if } \omega \in A_j^c; \end{cases} \quad 1 \leq j \leq n.$$

These are random variables [at least in a countable sample space]. Each takes only the two values 0 or 1, and we have

$${I_{A_i} = 1} = A_j, \quad {I_{A_i} = 0} = A_j^c.$$

Now if we apply the condition (5.5.1) of independence to the random variables I_{A_1}, \ldots, I_{A_n} , they reduce exactly to the conditions

$$P(\tilde{A}_1 \cdots \tilde{A}_n) = P(\tilde{A}_1) \cdots P(\tilde{A}_n), \tag{5.5.4}$$

where each \tilde{A}_j may be A_j or A_j^c but, of course, must be the same on both sides. Now it can be shown (Exercise 36 ahead) that the condition (5.5.4) for all possible choices of \tilde{A}_j is exactly equivalent to the condition (2.4.5). Hence the independence of the events A_1, \ldots, A_n is equivalent to the independence of their indicators.

The study of independent random variables will be a central theme in any introduction to probability theory. Historically and empirically, they are known as independent trials. We have given an informal discussion of this concept in §2.4. Now it can be formulated in terms of random variables

as follows: a sequence of independent trials is just a sequence of independent random variables (X_1, \ldots, X_n) where X_i represents the outcome of the *i*th trial. Simple illustrations are given in Examples 7 and 8 of §2.4, where in Example 7 the missing random variables are easily supplied. Incidentally, these examples establish the *existence* of independent random variables so that we are assured that our theorems such as the propositions in this section are not vacuities. Actually we can even construct independent random variables with arbitrarily given distributions (see [Chung 1; Chapter 3]). [It may amuse you to know that mathematicians have been known to define and study objects that later turn out to be nonexistent!] This remark will be relevant in later chapters; for the moment we shall add one more general proposition to broaden the horizon.

Proposition 6. Let $\varphi_1, \ldots, \varphi_n$ be an arbitrary real-valued function on $(-\infty, \infty)$; then the random variables

$$\varphi_1(X_1), \dots, \varphi_n(X_n) \tag{5.5.5}$$

are independent.

Proof: Let us omit the subscripts on X and φ and ask the question: for a given real number y, what are the values of x such that

$$\varphi(x) = y$$
 and $X = x$?

The set of such values must be countable since X is countably valued; call it S; of course, it depends on y, φ and X. Then $\{\varphi(X) = y\}$ means exactly the same thing as $\{X \in S\}$. Hence for arbitrary y_1, \ldots, y_n , the events

$$\{\varphi_1(X_1)=y_1\},\ldots,\{\varphi_n(X_n)=y_n\}$$

are just those in (5.5.3) for certain sets S_1, \ldots, S_n specified above. So Proposition 6 follows from Proposition 5.

This proposition will be put to good use in Chapter 6. Actually there is a more general result as follows. If we separate the random variables X_1, \ldots, X_n into any number of blocks, and take a function of those in each block, then the resulting random variables are independent. The proof is not so different from the special case given above and will be omitted.

As for general random variables, they are defined to be independent iff for any real numbers x_1, \ldots, x_n , the events

$$\{X_1 \le x_1\}, \dots, \{X_n \le x_n\}$$
 (5.5.6)

are independent. In particular,

$$P(X_1 \le x_1, \dots, X_n \le x_n) = P(X_1 \le x_1) \dots P(X_n \le x_n).$$
 (5.5.7)

In terms of the joint distribution function F for the random vector (X_1, \ldots, X_n) discussed in §4.6, the preceding equation may be written as

$$F(x_1, \dots, x_n) = F_1(x_1) \dots F_n(x_n),$$
 (5.5.8)

where F_j is the marginal distribution of $X_j, 1 \leq j \leq n$. Thus in case of independence the marginal distributions determine the joint distribution.

It can be shown that as a consequence of the definition, events such as those in (5.5.3) are also independent, provided that the sets S_1, \ldots, S_n are reasonable [Borel]. In particular, if there is a joint density function f, then we have

$$P(X_1 \in S_1, \dots, X_n \in S_n) = \left\{ \int_{S_1} f_1(u) \, du \right\} \dots \left\{ \int_{S_n} f_n(u) \, du \right\}$$
$$= \int_{S_1} \dots \int_{S_n} f_1(u_1) \dots f_n(u_n) \, du_1 \dots du_n,$$

where f_1, \ldots, f_n are the marginal densities. But the probability in the first member above is also equal to

$$\int_{S_1} \cdots \int_{S_n} f(u_1, \ldots, u_n) du_1 \ldots du_n$$

as in (4.6.6). Comparison of these two expressions yields the equation

$$f(u_1, \dots, u_n) = f_1(u_1) \dots f_n(u_n).$$
 (5.5.9)

This is the form that (5.5.8) takes in the density case.

Thus we see that stochastic independence makes it possible to factorize a joint probability, distribution, or density. In the next chapter we shall see that it enables us to factorize mathematical expectation, generating function and other transforms.

Numerous results and applications of independent random variables will be given in Chapters 6 and 7. In fact, the main body of classical probability theory is concerned with them. So much so that in his epoch-making monograph Foundations of the Theory of Probability, Kolmogorov [1903–87; leading Russian mathematician and one of the founders of modern probability theory] said: "Thus one comes to perceive, in the concept of independence, at least the first germ of the true nature of problems in probability theory." Here we will content ourselves with two simple examples.

Example 10. A letter from Pascal to Fermat (dated Wednesday, 29th July, 1654), contains, among many other mathematical problems, the following passage:

M. de Méré told me that he had found a fallacy in the theory of numbers, for this reason: If one undertakes to get a six with one die, the advantage in getting it in 4 throws is as 671 is to 625. If one undertakes to throw 2 sixes with two dice, there is a disadvantage in undertaking it in 24 throws. And nevertheless 24 is to 36 (which is the number of pairings of the faces of two dice) as 4 is to 6 (which is the number of faces of one die). This is what made him so indignant and made him say to one and all that the propositions were not consistent and Arithmetic was self-contradictory: but you will very easily see that what I say is correct, understanding the principles as you do.

This famous problem, one of the first recorded in the history of probability and which challenged the intellectual giants of the time, can now be solved by a beginner.

To throw a six with 1 die in 4 throws means to obtain the point "six" at least once in 4 trials. Define $X_n, 1 \le n \le 4$, as follows:

$$P(X_n = k) = \frac{1}{6}, \quad k = 1, 2, \dots, 6,$$

and assume that X_1, X_2, X_3, X_4 are independent. Put $A_n = \{X_n = 6\}$; then the event in question is $A_1 \cup A_2 \cup A_3 \cup A_4$. It is easier to calculate the probability of its complement, which is identical to $A_1^c A_2^c A_3^c A_4^c$. The trials are assumed to be independent and the dice unbiased. We have as a case of (5.5.4),

$$P(A_1^c A_2^c A_3^c A_4^c) = P(A_1^c) P(A_2^c) P(A_3^c) P(A_4^c) = \left(\frac{5}{6}\right)^4;$$

hence

$$P(A_1 \cup A_2 \cup A_3 \cup A_4) = 1 - \left(\frac{5}{6}\right)^4 = 1 - \frac{625}{1296} = \frac{671}{1296}.$$

This last number is approximately equal to 0.5177. Since 1296-671=625, the "odds" are as 671 to 625, as stated by Pascal.

Next consider two dice; let (X'_n, X''_n) denote the outcome obtained in the *n*th throw of the pair, and let

$$B_n = \{X'_n = 6; X''_n = 6\}.$$

Then $P(B_n^c) = 35/36$, and

$$P(B_1^c B_2^c \cdots B_{24}^c) = \left(\frac{35}{36}\right)^{24},$$

$$P(B_1 \cup B_2 \cup \cdots \cup B_{24}) = 1 - \left(\frac{35}{36}\right)^{24}.$$

This last number is approximately equal to 0.4914, which confirms the disadvantage.

One must give great credit to de Méré for his sharp observation and long experience at gaming tables to discern the narrow inequality

$$P(A_1 \cup A_2 \cup A_3 \cup A_4) > \frac{1}{2} > P(B_1 \cup B_2 \cup \cdots \cup B_{24}).$$

His arithmetic went wrong because of a fallacious "linear hypothesis." [According to some historians the problem did not originate with de Méré.]

Example 11. If two points are picked at random from the interval [0,1], what is the probability that the distance between them is less than 1/2?

By now you should be able to interpret this kind of cryptogram. It means: if X and Y are two independent random variables each of which is uniformly distributed in [0,1], find the probability P(|X-Y|<1/2). Under the hypotheses the random vector (X,Y) is uniformly distributed over the unit square U (see Fig. 24); namely for any reasonable subset S of U, we have

$$P\{(X,Y) \in S\} = \iint\limits_{S} du \, dv.$$

This is seen from the discussion after (4.6.6); in fact the f(u,v) there is equal to $f_1(u)f_2(v)$ by (5.5.9) and both f_1 and f_2 are equal to 1 in [0,1] and 0 outside. For the present problem S is the set of points (u,v) in U satisfying the inequality |u-v|<1/2. You can evaluate the double integral above over this set if you are good at calculus, but it is a lot easier to do this geometrically as follows. Draw two lines u-v=1/2 and u-v=-1/2; then S is the area bounded by these lines and the sides of the square. The complementary area U-S is the union of two triangles each of area $1/2(1/2)^2=1/8$. Hence we have

area of
$$S = 1 - 2 \cdot \frac{1}{8} = \frac{3}{4}$$
,

and this is the required probability.

Example 12. Suppose X_1, X_2, \ldots, X_n are independent random variables with distributions F_1, F_2, \ldots, F_n as in (4.5.4). Let

$$M = \max(X_1, X_2, \dots, X_n),$$

$$m = \min(X_1, X_2, \dots, X_n).$$

Find the distribution functions of M and m.

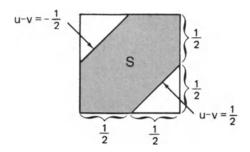


Figure 24

Using (5.5.7), we have for each x

$$F_{\max}(x) = P(M \le x) = P(X_1 \le x; X_2 \le x; \dots; X_n \le x)$$

$$= P(X_1 \le x) P(X_2 \le x) \dots P(X_n \le x)$$

$$= F_1(x) F_2(x) \dots F_n(x).$$

In particular, if all the F's are the same,

$$F_{\max}(x) = F(x)^n$$
.

As for the minimum, it is convenient to introduce the "tail distribution" G_j corresponding to each F_j as follows:

$$G_j(x) = P\{X_j > x\} = 1 - F_j(x).$$

Then we have, using the analogue of (5.5.2) this time with $S_j = (x_j, \infty)$,

$$G_{\min}(x) = P(m > x) = P(X_1 > x; X_2 > x; \dots; X_n > x)$$

$$= P(X_1 > x)P(X_2 > x) \cdots P(X_n > x)$$

$$= G_1(x)G_2(x) \cdots G_n(x).$$

Hence

$$F_{\min}(x) = 1 - G_1(x)G_2(x) \cdots G_n(x).$$

If all the F's are the same, this becomes

$$G_{\min}(x) = G(x)^n$$
, $F_{\min}(x) = 1 - G(x)^n$.

Here is a concrete illustration. Suppose a town depends on three reservoirs for its water supply, and suppose that its daily draws from them are independent and have exponential densities $\lambda_1 e^{-\lambda_1 x}$, $\lambda_2 e^{-\lambda_2 x}$, $\lambda_3 e^{-\lambda_3 x}$, respectively. Suppose each reservoir can supply a maximum of N gallons per

day to that town. What is the probability that on a specified day the town will run out of water?

Call the draws X_1, X_2, X_3 on that day; the probability in question is by (4.5.12)

$$P(X_1 > N; X_2 > N; X_3 > N) = e^{-\lambda_1 N} e^{-\lambda_2 N} e^{-\lambda_3 N} = e^{-(\lambda_1 + \lambda_2 + \lambda_3)N}.$$

* The rest of the section is devoted to a brief study of a logical notion that is broader than pairwise independence. This notation is inherent in statistical comparison of empirical data, operational evaluation of alternative policies, etc. Some writers even base the philosophical foundation of statistics on such a qualitative notion.

An event A is said to be favorable to another event B iff

$$P(AB) \ge P(A)P(B). \tag{5.5.10}$$

This will be denoted symbolically by $A \parallel B$. It is thus a binary relation between two events which includes pairwise independence as a special case. An excellent example is furnished by the divisibility by any two positive integers; see §2.4 and Exercise 17 in Chapter 2.

It is clear from (5.5.10) that the relation \parallel is symmetric; it is also reflexive since $P(A) \geq P(A)^2$ for any A. But it is not transitive, namely $A \parallel B$ and $B \parallel C$ do not imply $A \parallel C$. In fact, we will show by an example that even the stronger relation of pairwise independence is not transitive.

Example 13. Consider families with two children as in Example 5 of §5.1: $\Omega = \{(bb), (bg), (gb), (gg)\}$. Let such a family be chosen at random and consider the three events below:

A =first child is a boy;

B =the two children are of different sex;

C =the first child is a girl.

Then

$$AB = \{(bg)\}, BC = \{(gb)\}, AC = \emptyset.$$

A trivial computation then shows that P(AB) = P(A)P(B), P(BC) = P(B)P(C), but $P(AC) = 0 \neq P(A)P(C)$. Thus the pairs $\{A, B\}$ and $\{B, C\}$ are independent but the pair $\{A, C\}$ is not.

A slight modification will show that pairwise independence does not imply total independence for three events. Let Then

$$AD = \{(bb)\}, BD = \{(gb)\}, ABD = \emptyset;$$

and so
$$P(ABD) = 0 \neq P(A)P(B)P(D) = 1/8$$
.

Not so long ago one could still find textbooks on probability and statistics in which total independence was confused with pairwise independence. It is easy on hindsight to think of everyday analogues of the counterexamples above. For instance, if A is friendly to B, and B is friendly to C, why should it follow that A is friendly to C? Again, if every two of three people A, B, C get along well, it is not necessarily the case that all three of them do.

These commonplace illustrations should tell us something about the use and misuse of "intuition." Pushing a bit further, let us record a few more nonsequitors below (" \Rightarrow " reads "does not imply"):

$$A \parallel C \quad \text{and} \quad B \parallel C \Rightarrow (A \cap B) \parallel C;$$

$$A \parallel B \quad \text{and} \quad A \parallel C \Rightarrow A \parallel (B \cap C);$$

$$A \parallel C \quad \text{and} \quad B \parallel C \Rightarrow (A \cup B) \parallel C;$$

$$A \parallel B \quad \text{and} \quad A \parallel C \Rightarrow A \parallel (B \cup C).$$

$$(5.5.11)$$

You may try some verbal explanations for these; rigorous but artificial examples are also very easy to construct; see Exercise 15.

The great caution needed in making conditional evaluation is no academic matter, for much statistical analysis of experimental data depends on a critical understanding of the basic principles involved. The following illustration is taken from Colin R. Blyth, "On Simpson's paradox and the sure-thing principle," *Journal of American Statistical Association*, Vol. 67 (1972) pp. 364–366.

Example 14. A doctor has the following data on the effect of a new treatment. Because it involved extensive follow-up treatment after discharge, he could handle only a few out-of-town patients and had to work mostly with patients residing in the city.

	City residents		NonCity residents	
	Treated	Untreated	Treated	Untreated
Alive Dead	1000 9000	50 950	95 5	5000 5000

5000

5000

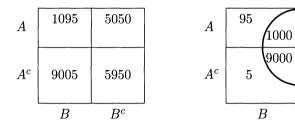
 B^c

 $\frac{50}{950}$

Let

$$A =$$
 alive,
 $B =$ treated,
 $C =$ cityresidents.

The sample space may be partitioned first according to A and B; then according to A, B, and C. The results are shown in the diagrams:



The various conditional probabilities, namely the classified proportions, are as follows:

$$\begin{split} P(A\mid B) &= \frac{1095}{10100} = \text{about } 10\%; \qquad P(A\mid BC) = \frac{1000}{10000}; \\ P(A\mid B^c) &= \frac{5050}{11000} = \text{about } 50\%; \qquad P(A\mid B^cC) = \frac{50}{1000}; \\ P(A\mid BC^c) &= \frac{95}{100}; \\ P(A\mid B^cC^c) &= \frac{5000}{10000}. \end{split}$$

Thus if the results (a matter of life or death) are judged from the conditional probabilities in the left column, the treatment seems to be a disaster since it had decreased the chance of survival five times! But now look at the right column, for city residents and noncity residents separately:

$$P(A \mid BC) = 10\%;$$
 $P(A \mid B^{c}C) = 5\%;$
 $P(A \mid BC^{c}) = 95\%;$ $P(A \mid B^{c}C^{c}) = 50\%.$

In both cases the chance of survival is doubled by the treatment.

The explanation is this: for some reason (such as air pollution), the C patients are much less likely to recover than the C^c patients, and most of those treated were C patients. Naturally, a treatment is going to show a poor recovery rate when used on the most seriously ill of the patients.

The arithmetical puzzle is easily solved by the following explicit formulas:

$$\begin{split} P(A \mid B) &= \frac{P(AB)}{P(B)} = \frac{P(ABC) + P(ABC^c)}{P(B)} \\ &= \frac{P(ABC)}{P(BC)} \frac{P(BC)}{P(B)} + \frac{P(ABC^c)}{P(BC^c)} \frac{P(BC^c)}{P(B)} \\ &= P(A \mid BC)P(C \mid B) + P(A \mid BC^c)P(C^c \mid B) \\ &= \frac{1000}{10000} \frac{10000}{10100} + \frac{95}{100} \frac{100}{10100}; \\ P(A \mid B^c) &= P(A \mid B^cC)P(C \mid B^c) + P(A \mid B^cC^c)P(C^c \mid B^c) \\ &= \frac{50}{1000} \frac{1000}{11000} + \frac{5000}{10000} \frac{10000}{11000}. \end{split}$$

It is those "hidden coefficients" $P(C \mid B), P(C^c \mid B), P(C \mid B^c), P(C^c \mid B^c)$ that have caused a reverse. A little parable will clarify the arithmetic involved. Suppose in two families both husbands and wives work. Husband of family 1 earns more than husband of family 2, wife of family 1 earns more than wife of family 2. For a certain good cause [or fun] both husband and wife of family 2 contribute half their monthly income; but in family 1 the husband contributes only 5% of his income, letting the wife contribute 95% of hers. Can you see why the poorer couple give more to the cause [or spend more on the vacation]?

This example should be compared with a simpler analogue in Exercise 11, where there is no paradox and intuition is a sure thing.

*5.6. Genetical models

This section treats an application to genetics. The probabilistic model discussed here is among the simplest and the most successful in empirical sciences.

Hereditary characters in *diploid* organisms such as human beings are carried by genes, which appear in pairs. In the simplest case each genre of a pair can assume two forms called *alleles*: A and a. For instance, A may be "blue-eyed" and a "brown-eyed" in a human being; or A may be "red blossom" and a "white blossom" in garden peas, which were the original subject of experiment by Mendel [1822–84]. We have then three *genotypes*:

there being no difference between Aa and aA [nature does not order the pair]. In some characters, A may be dominant whereas a recessive so that Aa cannot be distinguished from AA in appearance so far as the character

in question is concerned; in others Aa may be intermediate such as shades of green for eye color or pink for pea blossom. The reproductive cells, called gametes, are formed by splitting the gene pairs and have only one gene of each pair. At mating each parent therefore transmits one of the genes of the pair to the offspring through the gamete. The pure type AA or aa can of course transmit only A or a, whereas the mixed type Aa can transmit either A or a but not both. Now let us fix a gene pair and suppose that the parental genotypes AA, Aa, aa are in the proportions

$$u: 2v: w$$
 where $u > 0, v > 0, w > 0, u + 2v + w = 1.$

[The factor 2 in 2v is introduced to simplify the algebra below.] The total pool of these three genotypes is very large and the mating couples are formed "at random" from this pool. At each mating, each parent transmits one of the pair of genes to the offspring with probability 1/2, independently of each other, and independently of all other mating couples. Under these circumstances $random\ mating$ is said to take place. For example, if peas are well mixed in a garden, these conditions hold approximately; on the other hand, if the pea patches are segregated according to blossom colors then the mating will not be quite random.

The stochastic model can be described as follows. Two urns contain a very large number of coins of three types: with A on both sides, with A on one side and a on the other, and with a on both sides. Their proportions are as u:2v:w for each urn. One coin is chosen from each urn in such a way that all coins are equally likely. The two chosen coins are then tossed and the two uppermost faces determine the genotype of the offspring. What is the probability that it be AA, Aa, or aa? In a more empirical vein and using the frequency interpretation, we may repeat the process a large number of times to get an actual sample of the distribution of the types. Strictly speaking, the coins must be replaced each time so that the probability of each type remains constant in the repeated trials.

Let us tabulate the cases in which an offspring of type AA will result from the mating. Clearly this is possible only if there are at least two A genes available between the parents. Hence the possibilities are given in the first and second columns below.

Type of male	Type of female	Probability of mating of the couple	Probability of producing offspring AA from the couple	Probability of offspring AA
AA AA Aa Aa	AA Aa AA Aa	$u \cdot u = u^{2}$ $u \cdot 2v = 2uv$ $2v \cdot u = 2uv$ $2v \cdot 2v = 4v^{2}$	1 1/2 1/2 1/4	u^2 uv uv v^2

In the third column we give the probability of mating between the two designated genotypes in the first two entries of the same row; in the fourth column we give the conditional probability for the offspring to be of type AA given the parental types; in the fifth column the product of the probabilities in the third and fourth entries of the same row. By Proposition 2 of §5.2, the total probability for the offspring to be of type AA is given by adding the entries in the fifth column. Thus

$$P(\text{offspring is } AA) = u^2 + uv + uv + v^2 = (u + v)^2.$$

From symmetry, replacing u by w, we get

$$P(\text{offspring is } aa) = (v + w)^2.$$

Finally, we list all cases in which an offspring of type Aa can be produced, in a similar tabulation as the preceding one.

Type of male	Type of female	Probability of mating of the couple	Probability of producing offspring Aa from the couple	Probability of offspring Aa
AA Aa AA aa Aa	Aa AA aa AA aa	$u \cdot 2v = 2uv$ $2v \cdot u = 2uv$ $u \cdot w = uw$ $w \cdot u = uw$ $2v \cdot w = 2vw$	1/2 $1/2$ 1 1 $1/2$	uv uv uw uw vw
aa Aa	Aa Aa	$w \cdot 2v = 2vw$ $2v \cdot 2v = 4v^2$	$\frac{1/2}{1/2}$	$egin{array}{c} vw \ 2v^2 \end{array}$

By adding up the last column, we hence obtain:

$$P(\text{offspring is } Aa) = 2(uv + uw + vw + v^2) = 2(u+v)(v+w).$$

Let us put

$$p = u + v, \quad q = v + w$$
 (5.6.1)

so that p > 0, q > 0, p + q = 1. Let us also denote by $P_n(\cdots)$ the probability of the genotypes for offspring of the *n*th generation. Then the results obtained above are as follows:

$$P_1(AA) = p^2$$
, $P_1(Aa) = 2pq$, $P_1(aa) = q^2$. (5.6.2)

These give the proportions of the parental genotypes for the second generation. Hence in order to obtain P_2 , we need only substitute p^2 for u, pq for v, and q^2 for w in the two preceding formulas. Thus

$$P_2(AA) = (p^2 + pq)^2 = p^2,$$

 $P_2(Aa) = 2(p^2 + pq)(pq + q^2) = 2pq,$
 $P_2(aa) = (pq + q^2)^2 = q^2.$

Lo and behold: P_2 is the same as P_1 ! Does this mean that P_3 is also the same as P_1 , etc.? This is true, but only after the observation below. We have shown that $P_1 = P_2$ for an arbitrary P_0 [in fact, even the nit-picking conditions u > 0, v > 0, w > 0 may be omitted]. Moving over one generation, therefore, $P_2 = P_3$, even although P_1 may not be the same as P_0 . The rest is smooth sailing, and the result is known as the Hardy-Weinberg theorem. (G.H. Hardy [1877–1947] was a leading English mathematician whose main contributions were to number theory and classical analysis.)

Theorem. Under random mating for one pair of genes, the distribution of the genotypes becomes stationary from the first generation on, no matter what the original distribution is.

Let us assign the numerical values 2, 1, 0 to the three types AA, Aa, aa according to the number of A genes in the pair; and let us denote by X_n the random variable that represents the numerical genotype of the nth generation. Then the theorem says that for $n \ge 1$:

$$P(X_n = 2) = p^2$$
, $P(X_n = 1) = 2pq$, $P(X_n = 0) = q^2$. (5.6.3)

The distribution of X_n is stationary in the sense that these probabilities do not depend on n. Actually it can be shown that the process $\{X_n, n \geq 1\}$ is strictly stationary in the sense described in §5.4, because it is also a Markov chain; see Exercise 40.

The result embodied in (5.6.2) may be reinterpreted by an even simpler model than the one discussed above. Instead of gene pairs we may consider a pool of gametes, namely after the splitting of the pairs into individual genes. Then the A genes and a genes are originally in the proportion

$$(2u + 2v) : (2v + 2w) = p : q$$

because there are two A genes in the type AA, etc. Now we can think of these gametes as so many little tokens marked A or a in an urn, and assimilate the birth of an offspring to the random drawing (with replacement)

of two of the gametes to form a pair. Then the probabilities of drawing AA, Aa, aa are, respectively,

$$p \cdot p = p^2$$
, $p \cdot q + q \cdot p = 2pq$, $q \cdot q = q^2$.

This is the same result as recorded in (5.6.2).

The new model is not the same as the old one, but it leads to the same conclusion. It is tempting to try to identify the two models on hindsight, but the only logical way of doing so is to go through both cases as we have done. A priori or prima facie, they are not equivalent. Consider, for instance, the case of fishes: the females lay billions of eggs first and then the males come along and fertilize them with sperm. The partners may never meet. In this circumstance the second model fits the picture better, especially if we use two urns for eggs and sperm separately. [There are in fact creatures in which sex is not differentiated and which suits the one-urn model.] Such a model may be called the *spawning model*, in contrast to the *mating model* described earlier. In more complicated cases where more than one pair of genes is involved, the two models need not yield the same result.

Example 15. It is known in human genetics that certain "bad" genes cause crippling defects or disease. If a is such a gene, the genotype aa will not survive to adulthood. A person of genotype Aa is a carrier but appears normal because a is a recessive character. Suppose the probability of a carrier among the general population is p, irrespective of sex. Now if a person has an affected brother or sister who died in childhood, then he has a history in the family and cannot be treated genetically as a member of the general population. The probability of his being a carrier is a conditional one to be computed as follows. Both his parents must be carriers, namely of genotype Aa, for otherwise they could not have produced a child of genotype aa. Since each gene is transmitted with probability 1/2, the probabilities of their child to be AA, Aa, aa are 1/4, 1/2, 1/4, respectively. Since the person in question has survived he cannot be aa, and so the probability that he is AA or Aa is given by

$$P(AA \mid AA \cup Aa) = \frac{1}{3}, \quad P(Aa \mid AA \cup Aa) = \frac{2}{3}.$$

If he marries a woman who is not known to have a history of that kind in the family, then she is of genotype AA or Aa with probability 1-p or p as for the general population. The probabilities for the genotypes of their children are listed below.

Male	Female	Probability of the combination	Probability of producing AA	Probability of producing Aa	Probability of producing aa
\overline{AA}	AA	$\frac{1}{3}(1-p)$	1	0	0
AA	Aa	$\frac{1}{3}p$	$rac{1}{2}$	$\frac{1}{2}$	0
Aa	AA	$\frac{2}{3}(1-p)$	$rac{1}{2}$	$\frac{1}{2}$	0
Aa	Aa	$\frac{2}{3}p$	$\frac{1}{4}$	$\frac{1}{2}$	$\frac{1}{4}$

A simple computation gives the following distribution of the genotypes for the offspring:

$$P_1(AA) = \frac{2}{3} - \frac{p}{3};$$

$$P_1(Aa) = \frac{1}{3} + \frac{p}{6};$$

$$P_1(aa) = \frac{p}{6}.$$

The probability of a surviving child being a carrier is therefore

$$P_1(Aa \mid AA \cup Aa) = \frac{2+p}{6-p}.$$

If p is negligible, this is about 1/3. Hence from the surviving child's point of view, his having an affected uncle or aunt is only half as bad a hereditary risk as his father's having an affected sibling. One can now go on computing the chances for his children, and so on—exercises galore left to the reader.

In concluding this example, which concerns a serious human condition, it is proper to stress that the simple mathematical theory should be regarded only as a rough approximation since other genetical factors have been ignored in the discussion.

Exercises

- 1. Based on the data given in Example 14 of §5.5, what is the probability that (a) A living patient resides in the city? (b) a living treated patient lives outside the city?
- 2. All the screws in a machine come from the same factory but it is as likely to be from Factory A as from Factory B. The percentage of defective screws is 5% from A and 1% from B. Two screws are inspected;

if the first is found to be good, what is the probability that the second is also good?

- 3. There are two kinds of tubes in an electronic gadget. It will cease to function if and only if one of each kind is defective. The probability that there is a defective tube of the first kind is .1; the probability that there is a defective tube of the second kind is .2. It is known that two tubes are defective, what is the probability that the gadget still works?
- 4. Given that a throw of three unbiased dice shows different faces, what is the probability that (a) at least one is a six; (b) the total is eight?
- 5. Consider families with three children and assume that each child is equally likely to be a boy or a girl. If such a family is picked at random and the eldest child is found to be a boy, what is the probability that the other two are girls? The same question if a randomly chosen child from the family turns out to be a boy.
- 6. Instead of picking a family as in No. 5, suppose now a child is picked at random from all children of such families. If he is a boy, what is the probability that he has two sisters?
- 7. Pick a family as in No. 5, and then pick two children at random from this family. If they are found to both be girls, what is the probability they have a brother?
- 8. Suppose that the probability that both twins are boys is α , and that both are girls β ; suppose also that when the twins are of different sexes the probability of the first born being a girl is 1/2. If the first born of twins is a girl, what is the probability that the second is also a girl?
- 9. Three marksmen hit the target with probabilities 1/2, 2/3, 3/4, respectively. They shoot simultaneously and there are two hits. Who missed? Find the probabilities.
- 10. On a flight from Urbana to Paris my luggage did not arrive with me. It had been transferred three times and the probabilities that the transfer was not done in time were estimated to be 4/10, 2/10, 1/10, respectively, in the order of transfer. What is the probability that the first airline goofed?
- 11. Prove the "sure-thing principle": if

$$P(A \mid C) \ge P(B \mid C),$$

$$P(A \mid C^c) \ge P(B \mid C^c),$$

then P(A) > P(B).

12. Show that if $A \parallel B$, then

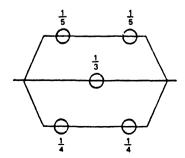
$$A^c \parallel B^c$$
, $A \not\parallel B^c$, $A^c \not\parallel B$.

- 13. Show that if $A \cap B = \emptyset$, then
 - (i) $A \parallel C$ and $B \parallel C \Rightarrow (A \cup B) \parallel C$;
 - (ii) $C \parallel A$ and $C \parallel B \Rightarrow C \parallel (A \cup B)$;
 - (iii) A and C are independent, B and C are independent $\Rightarrow A \cup B$ and C are independent.
- 14. Suppose P(H) > 0. Show that the set function

$$S \to P(S \mid H)$$
 for $S \in \Omega$ (countable)

is a probability measure.

- *15. Construct examples for all the assertions in (5.5.11). [Hint: a systematic but tedious way to do this is to assign p_1, \ldots, p_s to the eight atoms $ABC, \ldots, A^cB^cC^c$ (see (1.3.5)) and express the desired inequalities by means of them. The labor can be reduced by preliminary simple choices among the p's, such as making some of them zero and others equal. One can also hit upon examples by using various simple properties of a small set of integers; see an article by the author: "On mutually favorable events," Annals of Mathematical Statistics, Vol. 13 (1942), pp. 338–349.]
 - 16. Suppose that A_j , $1 \le j \le 5$, are independent events. Show that
 - (i) $(A_1 \cup A_2)A_3$ and $A_4^c \cup A_5^c$ are independent,
 - (ii) $A_1 \cup A_2, A_3 \cap A_3$ and A_5^c are independent.
 - 17. Suppose that in a certain casino there are three kinds of slot machines in equal numbers with payoff frequencies 1/3, 1/2, 2/3, respectively. One of these machines paid off twice in four cranks; what is the probability of a payoff on the next crank?
 - 18. A person takes four tests in succession. The probability of his passing the first test is p, that of his passing each succeeding test is p or p/2 depending on whether he passes or fails the preceding one. He qualifies provided he passes at least three tests. What is his chance of qualifying?
 - 19. An electric circuit looks as in the figure where the numbers indicate the probabilities of failure for the various links, which are all independent. What is the probability that the circuit is in operation?



- 20. It rains half of the time in a certain city, and the weather forecast is correct 2/3 of the time. Mr. Milquetoast goes out every day and is quite worried about the rain. So he will take his umbrella if the forecast is rain, but he will also take it 1/3 of the time even if the forecast is no rain. Find
 - (a) the probability of his being caught in rain without an umbrella,
 - (b) the probability of his carrying an umbrella without rain.
 - These are the two kinds of errors defined by Neyman and Pearson in their statistical history.* [Hint: compute the probability of "rain; forecast no rain; no umbrella," etc.]
- 21. Telegraphic signals "dot" and "dash" are sent in the proportion 3:4. Due to conditions causing very erratic transmission, a dot becomes a dash with probability 1/4, whereas a dash becomes a dot with probability 1/3. If a dot is received, what is the probability that it is sent as a dot?
- 22. A says B told him that C had lied. If each of these persons tells the truth with probability p, what is the probability that C indeed lied? Believe it or not, this kind of question was taken seriously one time under the name of "credibility of the testimony of witnesses." In the popular phrasing given above it is grossly ambiguous and takes a lot of words to explain the intended meaning. To cover one case in detail, suppose all three lied. Then B will tell A that C has told the truth, because B is supposed to know whether C has lied or not but decides to tell a lie himself: A will say that B told him that C had lied, since he wants to lie about what B told him, without knowing what C did. This is just one of the eight possible cases but the others can be similarly interpreted. A much clearer formulation is the model of transmission of signals used in No. 21. C transmits – or + according to whether he lies or not; then B transmits the message from C incorrectly or correctly if he lies or not; then A transmits the message from B in a similar manner. There will be no semantic impasse even if we go on this way to any number of witnesses. The question is: if "-" is received at the end of line, what is the probability that it is sent as such initially?
- 23. A particle starts from the origin and moves on the line 1 unit to the right or left with probability 1/2 each, the successive movements being independent. Let Y_n denote its position after n moves. Find the following probabilities:
 - (a) $P(Y_n \ge 0 \text{ for } 1 \le n \le 4);$
 - (b) $P(|Y_n| \le 2 \text{ for } 1 \le n \le 4);$
 - (c) $P(Y_n \ge 0 \text{ for } 1 \le n \le 4 \mid Y_4 = 0)$.

^{*}The reader is supposed to translate the verbal descriptions so that answers are obtainable. In the real world, predictions and decisions are made on even vaguer grounds.

24. In No. 23, show that if j < k < n, we have

$$P(Y_n = c \mid Y_j = a, Y_k = b) = P(Y_n = c \mid Y_k = b) = P(Y_{n-k} = c - b),$$

where a, b, c are any integers. Illustrate with j = 4, k = 6, n = 10, a = 2, b = 4, c = 6.

- 25. First throw an unbiased die, then throw as many unbiased coins as the point shown on the die.
 - (a) What is the probability of obtaining k heads?
 - (b) If 3 heads are obtained, what is the probability that the die showed n?
- 26. In a nuclear reaction a certain particle may split into 2 or 3 particles, or not split at all. The probabilities for these possibilities are p_2 , p_3 , and p_1 . The new particles behave in the same way and independently of each other as well as of the preceding reaction. Find the distribution of the total number of particles after two reactions.
- 27. An unbiased die is thrown n times; let M and m denote the maximum and minimum points obtained. Find P(m=2, M=5). [Hint: begin with $P(m \ge 2, M \le 5)$.]
- 28. Let X and Y be independent random variables with the same probability distribution $\{p_n, n \geq 1\}$. Find $P(X \leq Y)$ and P(X = Y).

In Problems 29–32, consider two numbers picked at random in [0,1].

- 29. If the smaller one is less than 1/4, what is the probability that the larger one is greater than 3/4?
- *30. Given that the smaller one is less than x, find the distribution of the larger one. [Hint: consider $P(\min < x, \max < y)$ and the two cases $x \le y$ and x > y.]
- *31. The two points picked divide [0,1] into three segments. What is the probability that these segments can be used to form a triangle? [Hint: this is the case if and only if the sum of lengths of any two is greater than the length of the third segment. Call the points X and Y and treat the case X < Y first.]
 - 32. Prove that the lengths of the three segments mentioned above have the same distributions. [Hint: consider the distribution of the smaller value picked, that of the difference between the two values, and use a symmetrical argument for the difference between 1 and the larger value.]
 - 33. In Pólya's urn scheme find:
 - (a) $P(R_3 \mid B_1 R_2)$;
 - (b) $P(R_3 \mid R_1R_2);$
 - (c) $P(R_3 \mid R_2)$;

- (d) $P(R_1 \mid R_2R_3)$;
- (e) $P(R_1 \mid R_2)$;
- (f) $P(R_1 | R_3)$.
- 34. Consider two urns U_i containing r_i red and b_i black balls respectively, i=1,2. A ball is drawn at random from U_1 and put into U_2 , then a ball is drawn at random from U_2 and put into urn U_1 . After this what is the probability of drawing a red ball from U_1 ? Show that if $b_1 = r_1$, $b_2 = r_2$, then this probability is the same as if no transfers have been made.
- *35. Assume that the a priori probabilities p in the sunrise problem (Example 9 of §5.2) can only take the values $k/100, 1 \le k \le 100$, with probability 1/100 each. Find $P(S^{n+1} \mid S^n)$. Replace 100 by N and let $N \to \infty$; what is the limit?
- *36. Prove that the events A_1, \ldots, A_n are independent if and only if

$$P(\tilde{A}_1 \cdots \tilde{A}_n) = P(\tilde{A}_1) \cdots P(\tilde{A}_n),$$

where each \tilde{A}_j may be A_j or A_j^c . [Hint: to deduce these equations from independence, use induction on n and also induction on k in $P(A_1^c \cdots A_k^c A_{k+1} \cdots A_n)$; the converse is easy by induction on n.]

- *37. Spell out a proof of Theorem 2 in §5.3. [Hint: label all balls and show that any particular sequence of balls has the same probability of occupying any given positions if all balls are drawn in order.]
 - 38. Verify Theorem 5 of §5.4 directly for the pairs (X_1, X_2) , (X_1, X_3) , and (X_2, X_3) .
 - 39. Assume that the three genotypes AA, Aa, aa are in the proportion $p^2: 2pq: q^2$, where p+q=1. If two parents chosen at random from the population have an offspring of type AA, what is the probability that another child of theirs is also of type AA? Same question with AA replaced by Aa.
 - 40. Let X_1 and X_2 denote the genotype of a female parent and her child. Assuming that the unknown genotype of the male parent is distributed as in Problem No. 39 and using the notation of (5.6.3), find the nine conditional probabilities below:

$$P{X_2 = k \mid X_1 = j}, \quad j = 0, 1, 2; k = 0, 1, 2.$$

These are called the $transition\ probabilities$ of a $Markov\ chain$; see $\S 8.3.$

*41. Prove that if the function φ defined on $[0,\infty)$ is nonincreasing and satisfies the Cauchy functional equation

$$\varphi(s+t) = \varphi(s)\varphi(t), \quad s \ge 0, t \ge 0;$$

then $\varphi(t) = e^{-\lambda t}$ for some $\lambda \ge 0$. Hence a positive random variable T has the property

$$P(T > s + t \mid T > s) = P(T > t), s \ge 0, t \ge 0$$

if and only if it has an exponential distribution. [Hint: $\varphi(0) = 1$: $\varphi(1/n) = \alpha^{1/n}$, where $\alpha = \varphi(1), \varphi(m/n) = \alpha^{m/n}$; if $m/n \le t < (m+1)/n$, then $\alpha^{(m+1)/n} \le \varphi(t) \le \alpha^{m/n}$; hence the general conclusion follows by letting $n \to \infty$.]

42. A needle of unit length is thrown onto a table that is marked with parallel lines at a fixed distance d from one another, where d>1. Let the distance from the midpoint of the needle to the nearest line be x, and let the angle between the needle and the perpendicular from its midpoint to the nearest line be θ . It is assumed that x and θ are independent random variables, each of which is uniformly distributed over its range. What is the probability that the needle intersects a line? This is known as Buffon's problem and its solution suggests an empirical [Monte Carlo] method of determining the value of π .

Mean, Variance, and Transforms

6.1. Basic properties of expectation

The mathematical expectation of a random variable, defined in §4.3, is one of the foremost notions in probability theory. It will be seen to play the same role as integration in calculus—and we know "integral calculus" is at least half of all calculus. Recall its meaning as a probabilistically weighted average [in a countable sample space] and rewrite (4.3.11) more simply as

$$E(X) = \sum_{\omega} X(\omega) P(\omega). \tag{6.1.1}$$

If we substitute |X| for X above, we see that the proviso (4.3.12) may be written as

$$E(|X|) < \infty. \tag{6.1.2}$$

We shall say that the random variable X is summable when (6.1.2) is satisfied. In this case we say also that "X has a finite expectation (or mean)" or "its expectation exists." The last expression is actually a little vague because we generally allow E(X) to be defined and equal to $+\infty$ when for instance $X \geq 0$ and the series in (6.1.1) diverges. See Exercises 27 and 28 of Chapter 4. We shall say so explicitly when this is the case.

It is clear that if X is bounded, namely when there exists a number M such that

$$|X(\omega)| \le M$$
 for all $\omega \in \Omega$,

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then X is summable and in fact

$$E(|X|) = \sum_{\omega} |X(\omega)|P(\omega) \le M \sum_{\omega} P(\omega) = M.$$

In particular, if Ω is finite then every random variable is bounded (this does not mean all of them are bounded by the same number). Thus the class of random variables having a finite expectation is quite large. For this class the mapping

$$X \to E(X) \tag{6.1.3}$$

assigns a number to a random variable. For instance, if X is the height of students in a school, then E(X) is their average height; if X is the income of wage earners, then E(X) is their average income; if X is the number of vehicles passing through a toll bridge in a day, then E(X) is the average daily traffic, etc.

If A is an event, then its indicator I_A (see §1.4) is a random variable, and we have

$$E(I_A) = P(A).$$

In this way the notion of mathematical expectation is seen to extend that of a probability measure.

Recall that if X and Y are random variables, then so is X + Y (Proposition 1 of §4.2). If X and Y both have finite expectations, it is intuitively clear what the expectation of X + Y should be. Thanks to the intrinsic nature of our definition, it is easy to prove the theorem.

Theorem 1. If X and Y are summable, then so is X + Y and we have

$$E(X+Y) = E(X) + E(Y).$$
 (6.1.4)

Proof: Applying the definition (6.1.1) to X + Y, we have

$$\begin{split} E(X+Y) &= \sum_{\omega} (X(\omega) + Y(\omega)) P(\omega) \\ &= \sum_{\omega} X(\omega) P(\omega) + \sum_{\omega} Y(\omega) P(\omega) = E(X) + E(Y). \end{split}$$

This is the end of the matter. You may wonder wherever do we need the condition (6.1.2)? The answer is: we want the defining series for E(X+Y) to converge absolutely, as explained in §4.3. This is indeed the case because

$$\sum_{\omega} |X(\omega) + Y(\omega)|P(\omega) \le \sum_{\omega} (|X(\omega)| + |Y(\omega)|)P(\omega)$$
$$= \sum_{\omega} |X(\omega)|P(\omega) + \sum_{\omega} |Y(\omega)|P(\omega) < \infty.$$

Innocuous or ingenious as Theorem 1 may appear, it embodies the most fundamental property of E. There is a pair of pale sisters as follows:

$$E(a) = a, \qquad E(aX) = aE(X) \tag{6.1.5}$$

for any constant a; and combining (6.1.4) and (6.1.5) we obtain

$$E(aX + bY) = aE(X) + bE(Y) \tag{6.1.6}$$

for any two constants a and b. This property makes the operation in (6.1.3) a "linear operator." This is a big name in mathematics; you may have heard of it in linear algebra or differential equations.

An easy extension of (6.1.4) by mathematical induction yields: if X_1, X_2, \ldots, X_n are summable random variables, then so is their sum and

$$E(X_1 + \dots + X_n) = E(X_1) + \dots + E(X_n). \tag{6.1.7}$$

Before we take up other properties of E, let us apply this to some interesting problems.

Example 1. A raffle lottery contains 100 tickets, of which there is one ticket bearing the prize \$10000, the rest being all zero. If I buy two tickets, what is my expected gain?

If I have only one ticket, my gain is represented by the random variable X, which takes the value 10000 on exactly one ω and 0 on all the rest. The tickets are assumed to be equally likely to win the prize; hence

$$X = \begin{cases} 10000 & \text{with probability } \frac{1}{100}, \\ 0 & \text{with probability } \frac{99}{100}, \end{cases}$$

and

$$E(X) = 10000 \cdot \frac{1}{100} + 0 \cdot \frac{99}{100} = 100.$$

Thus my expected gain is \$100. This is trivial, but now if I have two tickets I know very well only one of them can possibly win, so there is definite interference [dependence] between the two random variables represented by the tickets. Will this affect my expectation? Thinking a bit more deeply: if I am not the first person to have bought the tickets, perhaps by the time I get mine someone else has already taken the prize, albeit unknown to all. Will it then make a difference whether I get tickets early or late? Well, these questions have already been answered by the urn model discussed in

§5.3. We need only assimilate the tickets to 100 balls of which exactly 1 is black. Then if we define the outcome of the nth drawing by X_n , we know from Theorem 1 there that X_n has the same probability distribution as the X shown above, and so also the same expectation. For n=2 this was computed directly and easily without recourse to the general theorem. It follows that no matter what j and k are, namely for any two tickets drawn anytime, the expected value of both together is equal to

$$E(X_i + X_k) = E(X_i) + E(X_k) = 2E(X) = 200.$$

More generally, my expected gain is directly proportional to the number of tickets bought—a very fair answer, but is it so obvious in advance? In particular, if I buy all 100 tickets I stand to win 100 E(X) = 10000 dollars. This may sound dumb but it checks out.

To go one step further, let us consider two lotteries of exactly the same kind. Instead of buying two tickets X and Y from the same lottery, I may choose to buy one from each lottery. Now I have a chance to win \$20000. Does this make the scheme more advantageous to me? Yet Theorem 1 says that my expected gain is \$200 in either case. How is this accounted for? To answer this question you should figure out the distribution of X+Y under each scheme and compute E(X+Y) directly from it. You will learn a lot by comparing the results.

Example 2. (Coupon collecting problem). There are N coupons marked 1 to N in a bag. We draw one coupon after another with replacement. Suppose we wish to collect r different coupons; what is the expected number of drawings to get them? This is the problem faced by schoolchildren who collect baseball star cards; or by people who can win a sewing machine if they have a complete set of coupons that come in some detergent boxes. In the latter case the coupons may well be stacked against them if certain crucial ones are made very scarce. Here of course we consider the fair case in which all coupons are equally likely and the successive drawings are independent.

The problem may be regarded as one of waiting time, namely: we wait for the rth new arrival. Let X_1, X_2, \ldots denote the successive waiting times for a new coupon. Thus $X_1 = 1$ since the first is always new. Now X_2 is the waiting time for any coupon that is different from the first one drawn. Since at each drawing there are N coupons and all but one of them will be new, this reduces to Example 8 of §4.4 with success probability p = N - 1/N; hence

$$E(X_2) = \frac{N}{N-1}.$$

After these two different coupons have been collected, the waiting time for

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the third new one is similar with success probability p = N - 2/N; hence

$$E(X_3) = \frac{N}{N-2}.$$

Continuing this argument, we obtain for $1 \le r \le N$

$$E(X_1 + \dots + X_r) = \frac{N}{N} + \frac{N}{N-1} + \dots + \frac{N}{N-r+1}$$
$$= N\left(\frac{1}{N-r+1} + \dots + \frac{1}{N}\right).$$

In particular, if r = N, then

$$E(X_1 + \dots + X_N) = N\left(1 + \frac{1}{2} + \dots + \frac{1}{N}\right);$$
 (6.1.8)

and if N is even and r = N/2,

$$E(X_1 + \dots + X_{N/2}) = N\left(\frac{1}{\frac{N}{2} + 1} + \dots + \frac{1}{N}\right).$$
 (6.1.9)

Now there is a famous formula in mathematical analysis which says that

$$1 + \frac{1}{2} + \dots + \frac{1}{N} = \log N + C + \epsilon_N, \tag{6.1.10}$$

where the "log" is the natural logarithm to the base e, C is the Euler's constant = .5772... [nobody in the world knows whether it is a rational number or not], and ϵ_N tends to zero as N goes to infinity. For most purposes, the more crude asymptotic formula is sufficient:

$$\lim_{N \to \infty} \frac{1}{\log N} \left(1 + \frac{1}{2} + \dots + \frac{1}{N} \right) = 1.$$
 (6.1.11)

If we use this in (6.1.8) and (6.1.9), we see that for large values of N, the quantities there are roughly equal to $N \log N$ and $N \log 2 = \text{about } 7/10N$ respectively [how does one get $\log 2$ in the second estimate?]. This means: whereas it takes somewhat more drawings than half the number of coupons to collect half of them, it takes "infinitely" more drawings to collect all of them. The last few items are the hardest to get even if the game is not rigged.

A terrifying though not so realistic application is to the effects of aerial bombing in warfare. The results of the strikes are pretty much randomized under certain circumstances such as camouflage, decoy, foul weather, and intense enemy fire. Suppose there are 100 targets to be destroyed but each

strike hits one of them at random, perhaps repeatedly. Then it takes "on the average" about 100 log 100 or about 460 strikes to hit all targets at least once. Thus if the enemy has a large number of retaliatory launching sites, it will be very hard to knock them all out without accurate military intelligence. The conclusion should serve as a mathematical deterrent to the preemptive strike theory.

6.2. The density case

To return to saner matters, we will extend Theorem 1 to random variables in an arbitrary sample space. Actually the result is true in general, provided the mathematical expectation is properly defined. An inkling of this may be given by writing it as an abstract integral as follows:

$$E(X) = \int_{\Omega} X(\omega) d\omega, \qquad (6.2.1)$$

where " $d\omega$ " denotes the probability of the "element at ω ," as is commonly done for an area or volume element in multidimensional calculus—the so-called "differential." In this form (6.1.4) becomes

$$\int_{\Omega} (X(\omega) + Y(\omega)) d\omega = \int_{\Omega} X(\omega) d\omega + \int_{\Omega} Y(\omega) d\omega, \tag{6.2.2}$$

which is in complete analogy with the familiar formula in calculus:

$$\int_{I} (f(x) + g(x)) \, dx = \int_{I} f(x) \, dx + \int_{I} g(x) \, dx,$$

where I is an interval, say [0,1]. Do you remember anything of the proof of the last equation? It is established by going back to the definition of [Riemann] integrals through approximation by [Riemann] sums. For the probabilistic integral in (6.2.1) a similar procedure is followed. It is defined to be the limit of mathematical expectations of approximate discrete random variables [alluded to in §4.5]. These latter expectations are given by (6.1.1), and Theorem 1 is applicable to them. The general result (6.2.2) then follows by passing to the limit.

We cannot spell out the details of this proper approach in this text because it requires some measure theory, but there is a somewhat sneaky way to get Theorem 1 in the case where (X, Y) has a joint density as discussed in §4.6. Using the notation there, in particular (4.6.7), we have

$$E(X) = \int_{-\infty}^{\infty} u f(u, *) du, \quad E(Y) = \int_{-\infty}^{\infty} v f(*, v) dv.$$
 (6.2.3)

On the other hand, if we substitute $\varphi(x,y) = x + y$ in (4.6.8), we have

$$E(X+Y) = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} (u+v)f(u,v) \, du \, dv.$$
 (6.2.4)

Now this double integral can be split and evaluated by iterated integration:

$$\int_{-\infty}^{\infty} u \, du \left[\int_{-\infty}^{\infty} f(u, v) \, dv \right] + \int_{-\infty}^{\infty} v \, dv \left[\int_{-\infty}^{\infty} f(u, v) \, du \right]$$
$$= \int_{-\infty}^{\infty} u f(u, *) \, du + \int_{-\infty}^{\infty} v f(*, v) \, dv.$$

Comparison with (6.2.3) establishes (6.1.4).

The key to this method is the formula (6.2.4) whose proof was not given. The usual demonstration runs likes this. "Now look here: if X takes the value u and Y takes the value v, then X+Y takes the value u+v, and the probability that X=u and Y=v is $f(u,v)\,du\,dv$. See?" This kind of talk must be qualified as hand-waving or brow-beating. But it is a fact that applied scientists find such "demonstrations" quite convincing and one should go along until a second look becomes necessary, if ever. For the present the reader is advised to work out Exercise 40, which is the discrete analogue of the density argument above and is perfectly rigorous. These methods will be used again in the next section.

Example 3. Recall Example 5 in $\S 4.2$. The S_n 's being the successive times when the claims arrive, let us put

$$S_1 = T_1, S_2 - S_1 = T_2, \dots, S_n - S_{n-1} = T_n, \dots$$

Thus the T_n 's are the interarrival times. They are significant not only for our example of insurance claims, but also for various other models such as the "idle periods" for sporadically operated machinery, or "gaps" in a traffic pattern when the T's measure distance instead of time. In many applications it is these T's that are subject to statistical analysis. In the simplest case we may assume them to be exponentially distributed as in Example 12 of §4.5. If the density is $\lambda e^{-\lambda t}$ for all T_n , then $E(T_n) = 1/\lambda$. Since

$$S_n = T_1 + \cdots + T_n$$

we have by Theorem 1 in the density case

$$E(S_n) = E(T_1) + \dots + E(T_n) = \frac{n}{\lambda}.$$

Observe that there is no assumption about the independence of the T's, so that mutual influence between them is allowed. For example, several

accidents may be due to the same cause such as a 20-car smash-up on the freeway. Furthermore, the T's may have different λ 's due, e.g., to diurnal or seasonal changes. If $E(T_n) = 1/\lambda_n$ then

$$E(S_n) = \frac{1}{\lambda_1} + \dots + \frac{1}{\lambda_n}.$$

We conclude this section by solving a problem left over from §1.4 and Exercise 20 in Chapter 1; cf. also Problem 6 in §3.4.

Poincaré's Formula. For arbitrary events A_1, \ldots, A_n we have

$$P\left(\bigcup_{j=1}^{n} A_{j}\right) = \sum_{j} P(A_{j}) - \sum_{j,k} P(A_{j}A_{k}) + \sum_{j,k,l} P(A_{j}A_{k}A_{l}) \qquad (6.2.5)$$
$$- + \dots + (-1)^{n-1} P(A_{1} \dots A_{n}),$$

where the indices in each sum are distinct and range from 1 to n.

Proof: Let $\alpha_j = I_{A_j}$ be the indicator of A_j . Then the indicator of $A_1^c \cdots A_n^c$ is $\prod_{j=1}^n (1-\alpha_j)$; hence that of its complement is given by

$$I_{A_1} \cup \dots \cup A_n = 1 - \prod_{j=1}^n (1 - \alpha_j) = \sum_j \alpha_j - \sum_{j,k} \alpha_j \alpha_k$$
$$+ \sum_{j,k,l} \alpha_j \alpha_k \alpha_l - \dots + (-1)^{n-1} \alpha_1 \dots \alpha_n.$$

Now the expectation of an indicator random variable is just the probability of the corresponding event:

$$E(I_A) = P(A).$$

If we take the expectation of every term in the expansion above, and use (6.1.7) on the sums and differences, we obtain (6.2.5). [Henri Poincaré [1854–1912] was called the last universalist of mathematicians; his contributions to probability theory is largely philosophical and pedagogical. The formula above is a version of the "inclusion-exclusion principle" attributed to Sylvester.]

Example 4. (Matching problem, or *problem of rencontre*). Two sets of cards both numbered 1 to n are randomly matched. What is the probability of at least one match?

Solution. Let A_j be the event that the jth cards are matched, regardless of the others. There are n! permutations of the second set against the first set, which may be considered as laid out in natural order. If the jth cards match, that leaves (n-1)! permutations for the remaining cards, hence

$$P(A_j) = \frac{(n-1)!}{n!} = \frac{1}{n}.$$
(6.2.6)

Similarly if the jth and kth cards are both matched, where $j \neq k$, that leaves (n-2)! permutations for the remaining cards, hence

$$P(A_j A_k) = \frac{(n-2)!}{n!} = \frac{1}{n(n-1)};$$
(6.2.7)

next if j, k, l are all distinct, then

$$P(A_j A_k A_l) = \frac{(n-3)!}{n!} = \frac{1}{n(n-1)(n-2)},$$

and so on. Now there are $\binom{n}{1}$ terms in the first sum on the right side of (6.2.5), $\binom{n}{2}$ terms in the second, $\binom{n}{3}$ in the third, etc. Hence altogether the right side is equal to

$$\binom{n}{1}\frac{1}{n} - \binom{n}{2}\frac{1}{n(n-1)} + \binom{n}{3}\frac{1}{n(n-1)(n-2)} - + \dots + (-1)^{n-1}\frac{1}{n!}$$
$$= 1 - \frac{1}{2!} + \frac{1}{3!} - + \dots + (-1)^{n-1}\frac{1}{n!}.$$

Everybody knows (?) that

$$1 - e^{-1} = 1 - \frac{1}{2!} + \frac{1}{3!} - \dots + (-1)^{n-1} \frac{1}{n!} + \dots = \sum_{n=1}^{\infty} \frac{(-1)^{n-1}}{n!}.$$

This series converges very rapidly; in fact, it is easy to see that

$$\left|1 - e^{-1} - \left(1 - \frac{1}{2!} + \frac{1}{3!} - + \dots + (-1)^{n-1} \frac{1}{n!}\right)\right| \le \frac{1}{(n+1)!}$$

Hence as soon as $n \ge 4$, $1/(n+1)! \le 1/5! = 1/120$, and the probability of at least one match differs from $1 - e^{-1} \approx .63$ by less than .01. In other words, the probability of no match is about .63 for all $n \ge 4$.

What about the expected number of matches? The random number of matches is given by

$$N = I_{A_1} + \dots + I_{A_n}. \tag{6.2.8}$$

[Why? Think this one through thoroughly and remember that the A's are neither disjoint nor independent.] Hence its expectation is, by another application of Theorem 1,

$$E(N) = \sum_{j=1}^{n} E(I_{A_j}) = \sum_{j=1}^{n} P(A_j) = n \cdot \frac{1}{n} = 1,$$

namely exactly 1 for all n. This is neat, but it must be considered as a numerical accident.

6.3. Multiplication theorem; variance and covariance

We have indicated that Theorem 1 is really a general result in integration theory that is widely used in many branches of mathematics. In contrast, the next result requires stochastic independence and is special to probability theory.

Theorem 2. If X and Y are independent summable random variables, then

$$E(XY) = E(X)E(Y). (6.3.1)$$

Proof: We will prove this first when Ω is countable. Then both X and Y have a countable range. Let $\{x_j\}$ denote all the *distinct* values taken by X, similarly $\{y_k\}$ for Y. Next, let

$$A_{jk} = \{ w \mid X(\omega) = x_j, Y(\omega) = y_k \},\$$

namely A_{jk} is the sample subset for which $X = x_j$ and $Y = y_k$. Then the sets A_{jk} , as (j,k) range over all pairs of indices, are disjoint [why?] and their union is the whole space:

$$\Omega = \sum_{j} \sum_{k} A_{jk}.$$

The random variable XY takes the value $x_j y_k$ on the set A_{jk} , but some of these values may be equal, e.g., for $x_j = 2, y_k = 3$ and $x_j = 3, y_k = 2$. Nevertheless we get the expectation of XY by multiplying the probability of each A_{jk} with its value on the set, as follows:

$$E(XY) = \sum_{j} \sum_{k} x_{j} y_{k} P(A_{jk}). \tag{6.3.2}$$

This is a case of (4.3.15) and amounts merely to a grouping of terms in the defining series $\sum_{\omega} X(\omega)Y(\omega)P(\omega)$. Now by the assumption of independence,

$$P(A_{jk}) = P(X = x_j)P(Y = y_k).$$

Substituting this into (6.3.2), we see that the double sum splits into simple sums as follows:

$$\sum_{j} \sum_{k} x_j y_k P(X = x_j) P(Y = y_k)$$

$$= \left\{ \sum_{j} x_j P(X = x_j) \right\} \left\{ \sum_{k} y_k P(Y = y_k) \right\} = E(X) E(Y).$$

Here again the reassembling is justified by absolute convergence of the double series in (6.3.2).

Next, we prove the theorem when (X, Y) has a joint density function f, by a method similar to that used in §6.2. Analogously to (6.2.4), we have

$$E(XY) = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} uv f(u, v) du dv.$$

Since we have by (5.5.9), using the notation of §4.6,

$$f(u,v) = f(u,*)f(*,v),$$

the double integral can be split as follows:

$$\int_{-\infty}^{\infty} uf(u,*) du \int_{-\infty}^{\infty} vf(*,v) dv = E(X)E(Y).$$

Strictly speaking, we should have applied the calculations first to |X| and |Y|. These are also independent by Proposition 6 of §5.5, and we get

$$E(|XY|) = E(|X|)E(|Y|) < \infty.$$

Hence XY is summable and the manipulations above on the double series and double integral are valid. [These fussy details often distract from the main argument but are a necessary price to pay for mathematical rigor. The instructor as well as the reader is free to overlook some of these at his or her own discretion.]

The extension to any finite number of independent summable random variables is immediate:

$$E(X_1 \cdots X_n) = E(X_1) \cdots E(X_n). \tag{6.3.3}$$

This can be done directly or by induction. In the latter case we need that X_1X_2 is independent of X_3 , etc. This is true and was mentioned in §5.5—another fussy detail.

In the particular case of Theorem 2 where each X_j is the indicator of an event A_j , (6.3.3) reduces to

$$P(A_1 \cdots A_n) = P(A_1) \cdots P(A_n).$$

This makes it crystal clear that Theorem 2 cannot hold without restriction on the dependence. Contrast this with the corresponding case of (6.1.7):

$$E(I_{A_1} + \dots + I_{A_n}) = P(A_1) + \dots + P(A_n).$$

Here there is no condition whatsoever on the events such as their being disjoint, and the left member is to be emphatically distinguished from $P(A_1 \cup \cdots \cup A_n)$ or any other probability. This is the kind of confusion that has pestered pioneers as well as beginners. It is known as *Cardano's paradox*. [Cardano (1501–76) wrote the earliest book on games of chance.]

Example 5. Iron bars in the shape of slim cylinders are test-measured. Suppose the average length is 10 inches and average area of ends is 1 square inch. The average error made in the measurement of the length is .005 inch, that in the measurement of the area is .01 square inch. What is the average error made in estimating their weights?

Since weight is a constant times volume, it is sufficient to consider the latter: V = LA where L = length, A = area of ends. Let the errors be ΔL and ΔA , respectively; then the error in V is given by

$$\Delta V = (L + \Delta L)(A + \Delta A) - LA = L \Delta A + A \Delta L + \Delta L \Delta A.$$

Assuming independence between the measurements, we have

$$E(\Delta V) = E(L)E(\Delta A) + E(A)E(\Delta L) + E(\Delta A)E(\Delta L)$$
$$= 10 \cdot \frac{1}{100} + 1 \cdot \frac{1}{200} + \frac{1}{100} \cdot \frac{1}{200}$$

= .105 cubic inch if the last term is ignored.

Definition of Moment. For positive integer r, the mathematical expectation $E(X^r)$ is called the rth moment $[moment\ of\ order\ r]$ of X. Thus if X^r has a finite expectation, we say that X has a finite rth moment. For r=1, of course, the first moment is just the expectation or mean.

The case r=2 is of special importance. Since $X^2\geq 0$, we shall call $E(X^2)$ the second moment of X whether it is finite or equal to $+\infty$ depending on whether the defining series [in a countable Ω] $\sum_{\omega} X^2(\omega) P(\omega)$ converges or diverges.

When the mean E(X) is finite, it is often convenient to consider

$$X^0 = X - E(X) (6.3.4)$$

instead of X because its first moment is equal to zero. We shall say X^0 is obtained from X by *centering*.

Definition of Variance and Standard Deviation. The second moment of X^0 is called the *variance* of X and denoted by $\sigma^2(X)$; its positive square root $\sigma(X)$ is called the *standard deviation* of X.

There is an important relation among E(X), $E(X^2)$, and $\sigma^2(X)$, as follows.

Theorem 3. If $E(X^2)$ is finite, then so is E(|X|). We then have

$$\sigma^{2}(X) = E(X^{2}) - E(X)^{2}; \tag{6.3.5}$$

consequently,

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$$E(|X|)^2 \le E(X^2). \tag{6.3.6}$$

Proof: Since

$$X^{2} - 2|X| + 1 = (|X| - 1)^{2} \ge 0,$$

we must have (why?) $E(X^2 - 2|X| + 1) \ge 0$, and therefore $E(X^2) + 1 \ge 2E(|X|)$ by Theorem 1 and (6.1.5). This proves the first assertion of the theorem. Next we have

$$\sigma^{2}(X) = E\{(X - E(X))^{2}\} = E\{X^{2} - 2E(X)X + E(X)^{2}\}$$
$$= E(X^{2}) - 2E(X)E(X) + E(X)^{2} = E(X^{2}) - E(X)^{2}.$$

Since $\sigma^2(X) \geq 0$ from the first equation above, we obtain (6.3.6) by substituting |X| for X in (6.3.5).

What is the meaning of $\sigma(X)$? To begin with, X^0 is the deviation of X from its mean and can take both positive and negative values. If we are only interested in its magnitude then the mean absolute deviation is $E(|X^0|) = E(|X - E(X)|)$. This can actually be used but it is difficult for calculations. So we consider instead the mean square deviation $E(|X^0|^2)$, which is the variance. But then we should cancel out the squaring by extracting the root afterward, which gives us the standard deviation $+\sqrt{E(|X^0|^2)}$. This then is a gauge of the average deviation of a random variable [sample value] from its mean. The smaller it is, the better the random values cluster around its average and the population is well centered or concentrated. The true significance will be seen later when we discuss the convergence to a normal distribution in Chapter 7.

Observe that X and X + c have the same variance for any constant c; in particular, this is the case for X and X^0 . The next result resembles Theorem 1, but only in appearance.

Theorem 4. If X and Y are independent and both have finite variances, then

$$\sigma^{2}(X+Y) = \sigma^{2}(X) + \sigma^{2}(Y). \tag{6.3.7}$$

Proof: By the preceding remark, we may suppose that X and Y both have mean zero. Then X + Y also has mean zero and the variances in (6.3.7) are the same as second moments. Now

$$E(XY) = E(X)E(Y) = 0$$

by Theorem 2, and

$$E\{(X+Y)^2\} = E\{X^2 + 2XY + Y^2\}$$

= $E(X^2) + 2E(XY) + E(Y^2) = E(X^2) + E(Y^2)$

by Theorem 1, and this is the desired result.

The extension of Theorem 4 to any finite number of independent random variables is immediate. However, there is a general formula for the second moment without the assumption of independence, which is often useful. We begin with the algebraic identity:

$$(X_1 + \dots + X_n)^2 = \sum_{j=1}^n X_j^2 + 2 \sum_{1 \le j < k \le n} X_j X_k.$$

Taking expectations of both sides and using Theorem 1, we obtain

$$E\{(X_1 + \dots + X_n)^2\} = \sum_{j=1}^n E(X_j^2) + 2\sum_{1 \le j \le k \le n} E(X_j X_k).$$
 (6.3.8)

When the X's are centered and assumed to be independent, then all the mixed terms in the second sum above vanish and the result is the extension of Theorem 4 already mentioned.

Let us introduce two real indeterminants [dummy variables] ξ and η and consider the identity:

$$E\{(X\xi + Y\eta)^2\} = E(X^2)\xi^2 + 2E(XY)\xi\eta + E(Y^2)\eta^2.$$

The right member is a quadratic form in (ξ, η) and it is never negative because the left member is the expectation of a random variable which

does not take negative values. A well-known result in college algebra says that the coefficients of such a quadratic form $a\xi^2 + 2b\xi\eta + c\eta^2$ must satisfy the inequality $b^2 \leq ac$. Hence in the present case

$$E(XY)^2 \le E(X^2)E(Y^2).$$
 (6.3.9)

This is called the Cauchy-Schwarz inequality.

If X and Y both have finite variances, then the quantity

$$E(X^{0}Y^{0}) = E\{(X - E(X))(Y - E(Y))\}$$

$$= E\{XY - XE(Y) - YE(X) + E(X)E(Y)\}$$

$$= E(XY) - E(X)E(Y) - E(Y)E(X) + E(X)E(Y)$$

$$= E(XY) - E(X)E(Y)$$

is called the *covariance* of X and Y and denoted by Cov(X,Y); the quantity

$$\rho(X,Y) = \frac{\operatorname{Cov}(X,Y)}{\sigma(X)\sigma(Y)}$$

is called the *coefficient of correlation* between X and Y, provided of course the denominator does not vanish. If it is equal to zero, then X and Y are said to be *uncorrelated*. This is implied by independence but is in general a weaker property. As a consequence of (6.3.9), we always have $-1 \le \rho(X,Y) \le 1$. The sign as well as the absolute value of ρ gives a sort of gauge of the mutual dependence between the random variable. * See also Exercise 30 of Chapter 7.

Example 6. The most classical application of the preceding results is to the case of Bernoullian random variables (see Example 9 of §4.4). These are independent with the same probability distribution as follows:

$$X = \begin{cases} 1 & \text{with probability } p; \\ 0 & \text{with probability } q = 1 - p. \end{cases}$$
 (6.3.10)

We have encountered them in coin tossing (Example 8 of §2.4), but the scheme can be used in any repeated trials in which there are only two outcomes: success (X=1) and failure (X=0). For instance, Example 1 in §6.1 is the case where p=1/100 and the monetary unit is "ten grand." The chances of either "cure" or "death" in a major surgical operation is another illustration.

^{*}The mathematician Emil Artin told me the following story in 1947. "Everybody knows that probability and statistics are the same thing, and statistics is nothing but correlation. Now the correlation is just the cosine of an angle. Thus, all is trivial."

The mean and variance of X are easy to compute:

$$E(X) = p, \quad \sigma^{2}(X) = p - p^{2} = pq.$$

Let $\{X_n, n \geq 1\}$ denote Bernoullian random variables and write

$$S_n = X_1 + \dots + X_n \tag{6.3.11}$$

for the nth partial sum. It represents the total number of successes in n trials. By Theorem 1,

$$E(S_n) = E(X_1) + \dots + E(X_n) = np.$$
 (6.3.12)

This would be true even without independence. Next by Theorem 3,

$$\sigma^2(S_n) = \sigma^2(X_1) + \dots + \sigma^2(X_n) = npq.$$
 (6.3.13)

The ease with which these results are obtained shows a great technical advance. Recall that (6.3.12) has been established in (4.4.16), via the binomial distribution of S_n and a tricky computation. A similar method is available for (6.3.13) and the reader is strongly advised to carry it out for practice and comparison. But how much simpler is our new approach, going from the mean and variance of the individual summands to those of the sum without the intervention of probability distributions. In more complicated cases the latter will be very hard if not impossible to get. That explains why we are devoting several sections to the discussion of mean and variance which often suffice for theoretical as well as practical purposes.

Example 7. Returning to the matching problem in §6.2, let us now compute the standard deviation of the number of matches. The I_{A_j} 's in (6.2.8) are not independent, but formula (6.3.8) is applicable and yields

$$E(N^2) = \sum_{j=1}^{n} E(I_{A_j}^2) + 2 \sum_{1 \le j < k \le n} E(I_{A_j} I_{A_k}).$$

Clearly,

$$E(I_{A_j}^2) = P(A_j) = \frac{1}{n},$$

$$E(I_{A_j}I_{A_k}) = P(A_jA_k) = \frac{1}{n(n-1)}$$

by (6.2.6) and (6.2.7). Substituting into the above, we obtain

$$E(N^2) = n \cdot \frac{1}{n} + 2\binom{n}{2} \frac{1}{n(n-1)} = 1 + 1 = 2.$$

Hence

$$\sigma^2(N) = E(N^2) - E(N)^2 = 2 - 1 = 1.$$

Rarely an interesting general problem produces such simple numerical answers.

6.4. Multinomial distribution

A good illustration of the various notions and techniques developed in the preceding sections is the *multinomial distribution*. This is a natural generalization of the binomial distribution and serves as a model for repeated trials in which there are a number of possible outcomes instead of just "success or failure." We begin with the algebraic formula called the *multinomial theorem*:

$$(x_1 + \dots + x_r)^n = \sum_{n_1 \mid \dots \mid n_r \mid} \frac{n!}{n_1! \cdots n_r!} x_1^{n_1} \cdots x_r^{n_r}, \tag{6.4.1}$$

where the sum ranges over all ordered r-tuples of integers n_1, \ldots, n_r satisfying the following conditions:

$$n_1 \ge 0, \dots, n_r \ge 0, \quad n_1 + \dots + n_r = n.$$
 (6.4.2)

When r=2 this reduces to the binomial theorem. For then there are n+1 ordered couples

$$(0,n),(1,n-1),\ldots,(k,n-k),\ldots,(n,0)$$

with the corresponding coefficients

$$\frac{n!}{0! \, n!}, \frac{n!}{1! \, (n-1)!}, \dots, \frac{n!}{k! \, (n-k)!}, \dots, \frac{n!}{n! \, 0!},$$

i.e.,

$$\binom{n}{0}, \binom{n}{1}, \dots, \binom{n}{k}, \dots, \binom{n}{n}$$

Hence the sum can be written explicitly as

$$\binom{n}{0}x^0y^n + \binom{n}{1}x^1y^{n-1} + \dots + \binom{n}{k}x^ky^{n-k} + \dots + \binom{n}{n}x^ny^0$$
$$= \sum_{k=0}^n \binom{n}{k}x^ky^{n-k}.$$

In the general case the n identical factors $(x_1 + \cdots + x_r)$ on the left side of (6.4.1) are multiplied out and the terms collected on the right. Each term is of the form $x_1^{n_1} \cdots x_r^{n_r}$ with the exponents n_j satisfying (6.4.2). Such a term appears $n!/(n_1! \cdots n_r!)$ times because this is the number of ways of permuting n objects (the n factors) that belong to r different varieties (the r), such that r of them belong to the r th variety. [You see some combinatorics are in the nature of things and cannot be avoided even if you have skipped most of Chapter 3.]

A concrete model for the multinomial distribution may be described as follows. An urn contains balls of r different colors in the proportions

$$p_1:\cdots:p_r$$
, where $p_1+\cdots+p_r=1$.

We draw n balls one after another with replacement. Assume independence of the successive drawings, which is simulated by a thorough shake-up of the urn after each drawing. What is the probability that so many of the balls drawn are of each color?

Let X_1, \ldots, X_n be independent random variables all having the same distribution as the X below:

$$X = \begin{cases} 1 & \text{with probability } p_1, \\ 2 & \text{with probability } p_2, \\ \vdots & \\ r & \text{with probability } p_r. \end{cases}$$

$$(6.4.3)$$

What is the joint distribution of (X_1, \ldots, X_n) , namely

$$P(X_1 = x_1, \dots, X_n = x_n) \tag{6.4.4}$$

for all possible choices of x_1, \ldots, x_n from 1 to r? Here the numerical values correspond to labels for the colors. Such a quantification is not necessary but sometimes convenient. It also leads to questions that are not intended for the color scheme, such that the probability " $X_1 + \cdots + X_n = m$." But suppose we change the balls to lottery tickets bearing different monetary prizes, or to people having various ages or incomes, then the numerical formulation in (6.4.3) is pertinent. What about negative or fractional values for the X's? This can be accommodated by a linear transformation (cf. Example 15 in §4.5) provided all possible values are commensurable, say ordinary terminating decimals. For example, if the values are in three decimal places and range from -10 up, then we can use

$$X' = 10000 + 1000X$$

instead of X. The value -9.995 becomes 10000 - 9955 = 5 in the new scale. In a superpragmatic sense, we might even argue that the multinomial distribution is all we need for sampling in independent trials. For in

reality we shall never be able to distinguish between (say) $10^{10^{10}}$ different varieties of anything. But this kind of finitist attitude would destroy a lot of mathematics.

Let us evaluate (6.4.4). It is equal to $P(X_1 = x_1) \cdots P(X_n = x_n)$ by independence, and each factor is one of the p's in (6.4.3). To get an explicit expression we must know how many of the x_j 's are 1 or 2 or \cdots or r? Suppose n_1 of them are equal to 1, n_2 of them equal to $2, \ldots, n_r$ of them equal to r. Then these n_j 's satisfy (6.4.2) and the probability in question is equal to $p_1^{n_1} \cdots p_r^{n_r}$. It is convenient to introduce new random variables $N_j, 1 \leq j \leq r$, as follows:

$$N_i = \text{number of } X$$
's among (X_1, \ldots, X_n) that take the value j .

Each N_j takes a value from 0 to n, but the random variables N_1, \ldots, N_r cannot be independent since they are subject to the obvious restriction

$$N_1 + \dots + N_r = n. \tag{6.4.5}$$

However, their joint distribution can be written down:

$$P(N_1 = n_1, \dots, N_r = n_r) = \frac{n!}{n_1! \cdots n_r!} p_1^{n_1} \cdots p_r^{n_r}.$$
 (6.4.6)

The argument here is exactly the same as that given at the beginning of this section for (6.4.1), but we will repeat it. For any particular, or completely specified, sequence (X_1, \ldots, X_n) satisfying the conditions $N_1 = n_1, \ldots, N_r = n_r$, we have just shown that the probability is given by $p_1^{n_1} \cdots p_r^{n_r}$. But there are $n!/(n_1! \cdots n_r!)$ different particular sequences satisfying the same conditions, obtained by permuting the n factors of which n_1 factors are p_1, n_2 factors are p_2 , etc. To nail this down in a simple numerical example, let $n = 4, r = 3, n_1 = 2, n_2 = n_3 = 1$. This means in 4 drawings there are 2 balls of color 1, and 1 ball each of color 2 and 3. All the possible particular sequences are listed below:

$$\begin{array}{cccc} (1,1,2,3) & (1,1,3,2) \\ (1,2,1,3) & (1,3,1,2) \\ (1,2,3,1) & (1,3,2,1) \\ (2,1,1,3) & (3,1,1,2) \\ (2,1,3,1) & (3,1,2,1) \\ (2,3,1,1) & (3,2,1,1) \end{array}$$

Their number is 12 = 4!/(2! 1! 1!) and the associated probability is $12p_1^2p_2p_3$.

Formula (6.4.6), in which the indices n_j range over all possible integer values subject to (6.4.2), is called the *multinomial distribution* for the random variables (N_1, \ldots, N_r) . Specifically, it may be denoted by

 $M(n; r; p_1, \ldots, p_{r-1}, p_r)$ subject to $p_1 + \cdots + p_r = 1$. For the binomial case r = 2, this is often written as B(n; p), see Example 9 of §4.4.

If we divide (6.4.1) through by its left member, and put

$$p_j = \frac{x_j}{x_1 + \dots + x_r}, \quad 1 \le j \le r,$$

we obtain

$$1 = \sum \frac{n!}{n_1! \cdots n_r!} p_1^{n_1} \cdots p_r^{n_r}. \tag{6.4.7}$$

This yields a check that the sum of all the terms of the multinomial distribution is indeed equal to 1. Conversely, since (6.4.7) is a consequence of its probabilistic interpretation, we can deduce (6.4.1) from it, at least when $x_j \geq 0$, by writing $(p_1 + \cdots + p_r)^n$ for the left member in (6.4.7). This is another illustration of the way probability theory can add a new meaning to an algebraic formula; cf. the last part of §3.3.

Marginal distributions (see §4.6) of (N_1, \ldots, N_r) can be derived by a simple argument without computation. If we are interested in N_1 alone, then we can lump the r-1 other varieties as "not 1" with probability $1-p_1$. Thus the multinomial distribution collapses into a binomial one $B(n; p_1)$, namely:

$$P(N_1 = n_1) = \frac{n!}{n!(n - n_1)!} p_1^{n_1} (1 - p_1)^{n - n_1}.$$

From this we can deduce the mean and variance of N_1 as in Example 6 of §6.3. In general,

$$E(N_j) = np_j, \quad \sigma^2(N_j) = np_j(1 - p_j), \quad 1 \le j \le r.$$
 (6.4.8)

Next, if we are interested in the pair (N_1, N_2) , a similar lumping yields $M(n; 3; p_1, p_2, p_3)$, namely:

$$P(N_1 = n_1, N_2 = n_2)$$

$$= \frac{n!}{n_1! \, n_2! \, (n - n_1 - n_2)!} p_1^{n_1} p_2^{n_2} (1 - p_1 - p_2)^{n - n_1 - n_2}.$$
(6.4.9)

We can now express $E(N_1N_2)$ by using (4.3.15) or (6.3.2) [without independence]:

$$E(N_1 N_2) = \sum_{n_1 n_2} n_1 n_2 P(N_1 = n_1, N_2 = n_2)$$

$$= \sum_{n_1 n_2 n_3 n_1} \frac{n!}{n_1 n_2 p_1^{n_1} p_2^{n_2} p_3^{n_3}},$$

where $n_3 = n - n_1 - n_2$, $p_3 = 1 - p_1 - p_2$, and the sum ranges as in (6.4.2) with r = 3. The multiple sum above can be evaluated by generalizing the device used in Example 9 of §4.4 for the binomial case. Take the indicated second partial derivative below:

$$\frac{\partial^2}{\partial x_1 \partial x_2} (x_1 + x_2 + x_3)^n = n(n-1)(x_1 + x_2 + x_3)^{n-2}$$

$$= \sum \frac{n!}{n_1! \, n_2! \, n_3!} n_1 n_2 x_1^{n_1 - 1} x_2^{n_2 - 1} x_3^{n_3}.$$

Multiply through by x_1x_2 and then put $x_1 = p_1, x_2 = p_2, x_3 = p_3$. The result is $n(n-1)p_1p_2$ on one side and the desired multiple sum on the other. Hence we have in general for $j \neq k$:

$$E(N_{j}N_{k}) = n(n-1)p_{j}p_{k};$$

$$Cov(N_{j}, N_{k}) = E(N_{j}N_{k}) - E(N_{j})E(N_{k})$$

$$= n(n-1)p_{j}p_{k} - (np_{j})(np_{k}) = -np_{j}p_{k}.$$
(6.4.10)

It is fun to check out the formula (6.3.8), recalling (6.4.5):

$$n^{2} = E\{(N_{1} + \dots + N_{r})^{2}\}\$$

$$= \sum_{j=1}^{r} \{n(n-1)p_{j}^{2} + np_{j}\} + 2\sum_{1 \leq j < k \leq r} n(n-1)p_{j}p_{k}$$

$$= n(n-1) \left(\sum_{j=1}^{r} p_{j}\right)^{2} + n\sum_{j=1}^{r} p_{j} = n(n-1) + n = n^{2}.$$

There is another method of calculating $E(N_jN_k)$, similar to the first method in Example 6, §6.3. Let j be fixed and

$$\xi(x) = \begin{cases} 1 & \text{if } x = j, \\ 0 & \text{if } x \neq j. \end{cases}$$

As a function ξ of the real variable x, it is just the indicator of the singleton $\{j\}$. Now introduce the random variable

$$\xi_{\nu} = \xi(X_{\nu}) = \begin{cases} 1 & \text{if } X_{\nu} = j, \\ 0 & \text{if } X_{\nu} \neq j. \end{cases}$$
 (6.4.11)

namely, ξ_{ν} is the indicator for the vent $\{X_{\nu} = j\}$. In other words, the ξ_{ν} 's count just those X_{ν} 's taking the value j, so that $N_j = \xi_1 + \cdots + \xi_n$. Now

we have

$$E(\xi_{\nu}) = P(X_{\nu} = j) = p_j,$$

$$\sigma^2(\xi_{\nu}) = E(\xi_{\nu}^2) - E(\xi_{\nu})^2 = p_j - p_j^2 = p_j(1 - p_j).$$

Finally, the random variables ξ_1, \ldots, ξ_n are independent since X_1, \ldots, X_n are, by Proposition 6 of §5.5. Hence by Theorems 1 and 4,

$$E(N_j) = E(\xi_1) + \dots + E(\xi_n) = np_j,$$

$$\sigma^2(N_j) = \sigma^2(\xi_1) + \dots + \sigma^2(\xi_n) = np_j(1 - p_j).$$
(6.4.12)

Next, let $k \neq j$ and define η and η_{ν} in the same way as ξ and ξ_{ν} are defined, but with k in place of j. Consider now for $1 \leq \nu \leq n, 1 \leq \nu' \leq n$,

$$E(\xi_{\nu}\eta_{\nu'}) = P(X_{\nu} = j, X_{\nu'} = k) = \begin{cases} p_{j}p_{k} & \text{if } \nu \neq \nu', \\ 0 & \text{if } \nu = \nu'. \end{cases}$$
 (6.4.13)

Finally, we calculate

$$E(N_{j}N_{k}) = E\left\{\left(\sum_{\nu=1}^{n} \xi_{\nu}\right) \left(\sum_{\nu'=1}^{n} \eta_{\nu'}\right)\right\} = E\left\{\sum_{\nu=1}^{n} \xi_{\nu} \eta_{\nu} + \sum_{1 \le \nu \ne \nu' \le n} \xi_{\nu} \eta_{\nu'}\right\}$$

$$= \sum_{\nu=1}^{n} E(\xi_{\nu} \eta_{\nu}) + \sum_{1 \le \nu \ne \nu' \le n} E(\xi_{\nu} \eta_{\nu'})$$

$$= n(n-1)p_{j}p_{k}$$

by (6.4.13) because there are $(n)_2 = n(n-1)$ terms in the last written sum. This is, of course, the same result as in (6.4.10).

We conclude with a simple numerical illustration of a general problem mentioned above.

Example 8. Three identical dice are thrown. What is the probability of obtaining a total of 9? The dice are not supposed to be symmetrical, and the probability of turning up face j is equal to p_j , $1 \le j \le 6$; same for all three dice.

Let us list the possible cases in terms of the X's and the N's, respectively:

X_1	X_2	X_3	N_1	N_2	N_3	N_4	N_5	N_6	Permutation number	Probability
1	2	6	1	1				1	6	$6p_1p_3p_6$
1	3	5	1		1		1		6	$6p_1p_3p_5$
1	4	4	1			2			3	$3p_1p_4^2$
2	2	5		2			1		3	$3p_2^2p_5$
2	3	4		1	1	1			6	$6p_2p_3p_4$
3	3	3			3				1	p_3^3

Hence

$$P(X_1 + X_2 + X_3 = 9) = 6(p_1p_2p_6 + p_1p_3p_5 + p_2p_3p_4) + 3(p_1p_4^2 + p_2^2p_5) + p_3^3.$$

If all the p's are equal to 1/6, then this is equal to

$$\frac{6+6+3+3+6+1}{6^3} = \frac{25}{216}.$$

The numerator 25 is equal to the sum

$$\sum \frac{3!}{n_1! \cdots n_6!},$$

where the n_j 's satisfy the conditions

$$n_1 + n_2 + n_3 + n_4 + n_5 + n_6 = 3,$$

$$n_1 + 2n_2 + 3n_3 + 4n_4 + 5n_5 + 6n_6 = 9.$$

There are six solutions tabulated above as possible values of the N_i 's.

In the general context of the X's discussed above, the probability $P(X_1 + \cdots + X_n = m)$ is obtained by summing the right side of (6.4.6) over all (n_1, \ldots, n_r) satisfying both (6.4.2) and

$$1n_1 + 2n_2 + \dots + rn_r = m.$$

It is obvious that we need a computing machine to handle such explicit formulas. Fortunately in most problems we are interested in cruder results such that

$$P(a_n \le X_1 + \dots + X_n \le b_n)$$

for large values of n. The relevant asymptotic results and limit theorems will be the subject matter of Chapter 7. One kind of machinery needed for this purpose will be developed in the next section.

6.5. Generating function and the like

A powerful mathematical device, a true gimmick, is the generating function invented by the great prolific mathematician Euler [1707–83] to study the partition problem in number theory. Let X be a random variable taking only nonnegative integer values with the probability distribution given by

$$P(X = j) = a_j, \quad j = 0, 1, 2, \dots$$
 (6.5.1)

The idea is to put all the information contained above in a compact capsule. For this purpose a dummy variable z is introduced and the following power series in z set up:

$$g(z) = a_0 + a_1 z + a_2 z^2 + \dots = \sum_{j=0}^{\infty} a_j z^j.$$
 (6.5.2)

This is called the *generating function* associated with the sequence of numbers $\{a_j, j \geq 0\}$. In the present case we may also call it the generating function of the random variable X with the probability distribution (6.5.1). Thus g is a function of z that will be regarded as a real variable, although in some more advanced applications it is advantageous to consider z as a complex variable. Remembering that $\sum_j a_j = 1$, it is easy to see that the power series in (6.5.2) converges for $|z| \leq 1$. In fact, it is dominated as follows:

$$|g(z)| \le \sum_{j} |a_j||z|^j \le \sum_{j} a_j = 1$$
, for $|z| \le 1$.

[It is hoped that your knowledge about power series goes beyond the "ratio test." The above estimate is more direct and says a lot more.] Now a theorem in calculus asserts that we can differentiate the series term by term to get the derivatives of g, so long as we restrict its domain of validity to |z| < 1:

$$g'(z) = a_1 + 2a_2z + 3a_3z^2 + \dots = \sum_{n=1}^{\infty} na_n z^{n-1},$$

$$g''(z) = 2a_2 + 6a_3z + \dots = \sum_{n=2}^{\infty} n(n-1)a_n z^{n-2}.$$
(6.5.3)

In general we have

$$g^{(j)}(z) = \sum_{n=j}^{\infty} n(n-1)\cdots(n-j+1)a_n z^{n-j} = \sum_{n=j}^{\infty} \binom{n}{j} j! \, a_n z^{n-j}. \quad (6.5.4)$$

If we set z = 0 above, all the terms vanish except the constant term:

$$g^{(j)}(0) = j! a_j \quad \text{or} \quad a_j = \frac{1}{j!} g^{(j)}(0).$$
 (6.5.5)

This shows that we can recover all the a_j 's from g. Therefore, not only does the probability distribution determine the generating function, but also vice versa. So there is no loss of information in the capsule. In particular, putting z = 1 in g' and g'' we get by (4.3.18)

$$g'(1) = \sum_{n=0}^{\infty} na_n = E(X), \quad g''(1) = \sum_{n=0}^{\infty} n^2 a_n - \sum_{n=0}^{\infty} na_n = E(X^2) - E(X);$$

provided that the series converge, in which case (6.5.3) holds for z=1.* Thus

$$E(X) = g'(1), \quad E(X^2) = g'(1) + g''(1).$$
 (6.5.6)

In practice, the following qualitative statement, which is a corollary of the above, is often sufficient.

Theorem 5. The probability distribution of a nonnegative integer-valued random variable is uniquely determined by its generating function.

Let Y be a random variable having the probability distribution $\{b_k, k \geq 0\}$ where $b_k = P(Y = k)$, and let h be its generating function:

$$h(z) = \sum_{k=0}^{\infty} b_k z^k.$$

Suppose that g(z) = h(z) for all |z| < 1; then the theorem asserts that $a_k = b_k$ for all $k \ge 0$. Consequently X and Y have the same distribution, and this is what we mean by "unique determination." The explicit formula (6.5.4) of course implies this, but there is a simpler argument as follows. Since

$$\sum_{k=0}^{\infty} a_k z^k = \sum_{k=0}^{\infty} b_k z^k, \quad |z| < 1;$$

we get at once $a_0 = b_0$ by setting z = 0 in the equation. After removing these terms we can cancel a factor z on both sides, and then get $a_1 = b_1$ by again setting z = 0. Repetition of this process establishes the theorem. You ought to realize that we have just reproduced a terrible proof of a standard

^{*}This is an Abelian theorem; cf. the discussion after (8.4.17).

result that used to be given in some calculus text books! Can you tell what is wrong there and how to make it correct?

We proceed to discuss a salient property of generating functions when they are multiplied together. Using the notation above we have

$$g(z)h(z) = \left(\sum_{j=0}^{\infty} a_j z^j\right) \left(\sum_{k=0}^{\infty} b_k z^k\right) = \sum_j \sum_k a_j b_k z^{j+k}.$$

We will rearrange the terms of this double series into a power series in the usual form. Then

$$g(z)h(z) = \sum_{l=0}^{\infty} c_l z^l,$$
 (6.5.7)

where

$$c_l = \sum_{j+k=l} a_j b_k = \sum_{j=0}^l a_j b_{l-j}.$$
 (6.5.8)

The sequence $\{c_j\}$ is called the *convolution* of the two sequences $\{a_j\}$ and $\{b_j\}$. What does c_l stand for? Suppose that the random variables X and Y are independent. Then we have, by (6.5.8),

$$c_{l} = \sum_{j=0}^{l} P(X = j)P(Y = l - j)$$
$$= \sum_{j=0}^{l} P(X = j, Y = l - j) = P(X + Y = l).$$

The last equation above is obtained by the rules in §5.2, as follows. Given that X = j, we have X + Y = l if and only if Y = l - j; hence by Proposition 2 of §5.2 [cf. (5.2.4)],

$$P(X + Y = l) = \sum_{j=0}^{\infty} P(X = j)P(X + Y = l \mid X = j)$$

$$= \sum_{j=0}^{\infty} P(X = j)P(Y = l - j \mid X = j)$$

$$= \sum_{j=0}^{l} P(X = j)P(Y = l - j),$$

because Y is independent of X and it does not take negative values. In other words, we have shown that for all $l \ge 0$,

$$P(X+Y=l)=c_l,$$

so that $\{c_l, l \geq 0\}$ is the probability distribution of the random variable X+Y. Therefore, by definition its generating function is given by the power series in (6.5.7) and so equal to the product of the generating functions of X and of Y. After an easy induction, we can state the result as follows.

Theorem 6. If the random variables X_1, \ldots, X_n are independent and have g_1, \ldots, g_n as their generating functions, then the generating function of the sum $X_1 + \cdots + X_n$ is given by the product $g_1 \cdots g_n$.

This theorem is of great importance since it gives a method to study sums of independent random variables via generating functions, as we shall see in Chapter 7. In some cases the product of the generating functions takes a simple form and then we can deduce its probability distribution by looking at its power series, or equivalently by using (6.5.5). The examples ahead will demonstrate this method.

For future reference let us take note that given a sequence of real numbers $\{a_j\}$, we can define the power series g as in (6.5.2). This will be called the generating function associated with the sequence. If the series has a nonzero radius of convergence, then the preceding analysis can be carried over to this case without the probability interpretations. In particular, the convolution of two such sequences can be defined as in (6.5.8), and (6.5.7) is still valid. In §8.4 below we shall use generating functions whose coefficients are probabilities, such that the series may diverge for z = 1.

Example 9. For the Bernoullian random variables X_1, \ldots, X_n (Example 6 of §6.3), the common generating function is

$$g(z) = q + pz$$

since $a_0 = q, a_1 = p$ in (6.5.1). Hence the generating function of $S_n = X_1 + \cdots + X_n$, where the X's are independent, is given by the nth power of g:

$$g(z)^n = (q + pz)^n.$$

Its power series is therefore known from the binomial theorem, namely,

$$g(z)^n = \sum_{k=0}^n \binom{n}{k} q^{n-k} p^k z^k.$$

On the other hand, by definition of a generating function, we have

$$g(z)^n = \sum_{k=0}^{\infty} P(S_n = k) z^k.$$

Comparison of the last two expressions shows that

$$P(S_n = k) = \binom{n}{k} p^k q^{n-k}, \quad 0 \le k \le n; \quad P(X_n = k) = 0, \quad k > n.$$

This is the Bernoulli formula we learned sometime ago, but the derivation is new and it is machine-processed.

Example 10. For the waiting-time distribution (§4.4), we have $p_j = q^{j-1}p, j > 1$; hence

$$g(z) = \sum_{j=1}^{\infty} q^{j-1pz^{j}} = \frac{p}{q} \sum_{j=1}^{\infty} (qz)^{i} = \frac{p}{q} \frac{qz}{1-qz} = \frac{pz}{1-qz}.$$
 (6.5.9)

Let $S_n = T_1 + \cdots + T_n$ where the T's are independent and each has the g in (6.5.9) as generating function. Then S_n is the waiting time for the nth success. Its generating function is given by g^n , and this can be expanded into a power series by using the binomial series and (5.4.4):

$$g(z)^{n} = \left(\frac{pz}{1 - qz}\right)^{n} = p^{n}z^{n} \sum_{j=0}^{\infty} {\binom{-n}{j}} (-qz)^{j}$$

$$= \sum_{j=0}^{\infty} \frac{n(n+1)\cdots(n+j-1)}{j!} p^{n}q^{j}z^{n+j} = \sum_{j=0}^{\infty} {\binom{n+j-1}{n-1}} p^{n}q^{j}z^{n+j}$$

$$= \sum_{k=n}^{\infty} {\binom{k-1}{n-1}} p^{n}q^{k-n}z^{k}.$$

Hence we obtain for $j \geq 0$,

$$P(S_n = n + j) = \binom{n+j-1}{j} p^n q^j = \binom{-n}{j} p^n (-q)^j.$$

The probability distribution given by $\left\{\binom{-n}{j}p^n(-q)^j, j \geq 0\right\}$ is called the negative binomial distribution of order n. The discussion above shows that its generating function is given by

$$\left(\frac{g(z)}{z}\right)^n = \left(\frac{p}{1 - qz}\right)^n. \tag{6.5.10}$$

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Now g(z)/z is the generating function of $T_1 - 1$ (why?), which represents the number of failures before the first success. Hence the generating function in (6.5.10) is that of the random variable $S_n - n$, which is the total number of failures before the nth success.

Example 11. For the dice problem at the end of §6.4, we have $p_j = 1/6$ for $1 \le j \le 6$ if the dice are symmetrical. Hence the associated generating function is given by

$$g(z) = \frac{1}{6}(z + z^2 + z^3 + z^4 + z^5 + z^6) = \frac{z(1 - z^6)}{6(1 - z)}.$$

The generating function of the total points obtained by throwing three dice is just g^3 . This can be expanded into a power series as follows:

$$g(z)^{3} = \frac{z^{3}}{6^{3}} \frac{(1-z^{6})^{3}}{(1-z)^{3}} = \frac{z^{3}}{6^{3}} (1-3z^{6}+3z^{12}-z^{18})(1-z)^{-3}$$
$$= \frac{z^{3}}{6^{3}} (1-3z^{6}+3z^{12}-z^{18}) \sum_{k=0}^{\infty} {k+2 \choose 2} z^{k}.$$
 (6.5.11)

The coefficient of z^9 is easily found by inspection, since there are only two ways of forming it from the product above:

$$\frac{1}{6^3} \left\{ 1 \cdot \binom{6+2}{2} - 3 \cdot \binom{0+2}{2} \right\} = \frac{28-3}{6^3} = \frac{25}{6^3}.$$

You may not be overly impressed by the speed of this new method, as compared with a combinatorial counting done in §6.4, but you should observe how the machinery works:

Step 1°: Code the probabilities $\{P(X=j), j \geq 0\}$ into a generating function g.

Step 2°: Process the function by raising it to nth power g^n .

Step 3°: Decode the probabilities $\{P(S_n = k), k \geq 0\}$ from g^n by expanding it into a power series.

A characteristic feature of machine process is that parts of it can be performed mechanically such as the manipulations in (6.5.10). We need not keep track of what we are doing at every stage: plug something in, push a few buttons or crank some knobs, and out comes the product. To carry this gimmickry one step further, we will now exhibit the generating function of X in a form Euler would not have recognized [the concept of a random variable came late, not much before 1930]:

$$g(z) = E(z^X),$$
 (6.5.12)

namely the mathematical expectation of z^X . Let us first recall that for each z, the function $\omega \to z^{X(\omega)}$ is indeed a random variable. For countable Ω this is a special case of Proposition 2 in §4.2. When X takes the value j, z^X takes the value z^j ; hence by (4.3.15) the expectation of z^X may be expressed as $\sum_{j=0}^{\infty} P(X=j)z^j$, which is g(z).

An immediate payoff is a new and smoother proof of Theorem 6 based on different principles. The generating function of $X_1 + \cdots + X_n$ is, by what has just been said, equal to

$$E(z^{X_1 + \dots + X_n}) = E(z^{X_1} z^{X_2 \dots} z^{X_n})$$

by the law of exponentiation. Now the random variables $z^{X_1}, z^{X_2}, \ldots, z^{X_n}$ are independent by Proposition 6 of §5.5; hence by Theorem 2 of §6.3

$$E(z^{X_1}z^{X_2}\cdots z^{X_n}) = E(z^{X_1})E(z^{X_2})\cdots E(z^{X_n}).$$
(6.5.13)

Since $E(z^{X_j}) = g_j(z)$ for each j, this completes the proof of Theorem 6.

Another advantage of the expression $E(z^X)$ is that it leads to extensions. If X can take arbitrary real values, this expression still has a meaning. For simplicity let us consider only $0 < z \le 1$. Every such z can be represented as $e^{-\lambda}$ with $0 \le \lambda < \infty$; in fact, the correspondence $z = e^{-\lambda}$ is one-to-one; see Figure 25.

Now consider the new expression after such a change of variable:

$$E(e^{-\lambda X}), \quad 0 \le \lambda < \infty.$$
 (6.5.14)

If X has the probability distribution in (6.5.1), then

$$E(e^{-\lambda X}) = \sum_{j=0}^{\infty} a_j e^{-j\lambda},$$

which is of course just our previous g(z) with $z = e^{-\lambda}$. More generally if X takes the values $\{x_i\}$ with probabilities $\{p_i\}$, then

$$E(e^{-\lambda X}) = \sum_{j} p_{j} e^{-\lambda x_{j}}$$
(6.5.15)

provided that the series converges absolutely. This is the case if all the values $x_j \geq 0$ because then $e^{-\lambda x_j} \leq 1$ and the series is dominated by $\sum_j p_j = 1$. Finally, if X has the density function f, then by (4.5.6) with $\varphi(u) = e^{-\lambda u}$:

$$E(e^{-\lambda X}) = \int_{-\infty}^{\infty} e^{-\lambda u} f(u) du, \qquad (6.5.16)$$

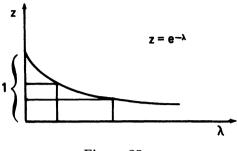


Figure 25

provided that the integral converges. This is the case if f(u) = 0 for u < 0, namely when X does not take negative values. We have therefore extended the notion of a generating function through (6.5.14) to a large class of random variables. This new gimmick is called the *Laplace transform* of X. In the analytic form given on the right side of (6.5.16) it is widely used in operational calculus, differential equations, and engineering applications.

If we replace the negative real parameter $-\lambda$ in (6.5.14) by the purely imaginary $i\theta$, where $i=\sqrt{-1}$ and θ is real, we get the Fourier transform $E(e^{i\theta X})$; in probability theory it is also known as the characteristic function of X. Let us recall De Moivre's formula (which used to be taught in high school trigonometry courses) for real u:

$$e^{iu} = \cos u + i \sin u$$
;

and its consequence

$$|e^{iu}|^2 = (\cos u)^2 + (\sin u)^2 = 1.$$

This implies that for any real random variable X, we have $|e^{i\theta X}| = 1$; hence the function φ :

$$\varphi(\theta) = E(e^{i\theta X}), \quad -\infty < \theta < \infty,$$
 (6.5.17)

is always defined; in fact, $|\varphi(\theta)| \leq 1$ for all θ . Herein lies the superiority of this new transform over the others discussed above, that cannot be defined sometimes because the associated series or integral does not converge. On the other hand, we pay the price of having to deal with complex variables and functions which lie beyond the scope of an elementary text. Nevertheless, we will invoke both the Laplace and Fourier transforms in Chapter 7, and for future reference let us record the following theorem.

Theorem 7. Theorems 5 and 6 remain true when the generating function is replaced by the Laplace transform (for nonnegative random variables) or the Fourier transform (for arbitrary random variables).

In the case of Theorem 6, this is immediate from (6.5.13) if the variable z there is replaced by $e^{-\lambda}$ or $e^{i\theta}$. For Theorem 5 the analogues lie deeper and require more advanced analysis (see [Chung 1, Chapter 6]). The reader is asked to accept their truth by analogy from the discussion above leading from $E(z^X)$ to $E(e^{-\lambda X})$ and $E(e^{i\theta X})$. After all, analogy is a time-honored method of learning.

Exercises

- 1. The Massachusetts state lottery has 1 million tickets. There is one first prize of \$50000; 9 second prizes of \$2500 each; 90 third prizes of \$250 each; 900 fourth prizes of \$25 each. What is the expected value of one ticket? Five tickets?
- 2. Suppose in the lottery above only 80% of the tickets are sold. What is the expected total to be paid out in prizes? If each ticket is sold at 50¢, what is the expected profit for the state?
- 3. Five residential blocks are polled for racial mixture. The number of houses having black or white owners listed below:

	1	2	3	4	5
Black	3	2	4	3	4
White	10	10	9	11	10

If two houses are picked at random from each block, what is the expected number of black-owned ones among them?

- 4. Six dice are thrown once. Find the mean and variance of the total points. Same question if the dice are thrown n times.
- 5. A lot of 1000 screws contain 1% with major defects and 5% with minor defects. If 50 screws are picked at random and inspected, what are the expected numbers of major and minor defectives?
- 6. In a bridge hand what is the expected number of spades? Of different suits? [Hint: for the second question let $X_j = 1$ or 0 depending on whether the jth suit is represented in the hand or not; consider $E(X_1 + X_2 + X_3 + X_4)$.]
- 7. An airport bus deposits 25 passengers at 7 stops. Assume that each passenger is as likely to get off at any stop as another and that they act independently. The bus stops only if someone wants to get off. What is the probability that nobody gets off at the third stop? What is the expected number of stops it will make? [Hint: let $X_j = 1$ or 0 according as someone gets off at the jth stop or not.]
- *8. Given 500 persons picked at random, (a) what is the probability that more than one of them have January 1 as birthday? (b) What is the expected number among them who have this birthday? (c) What is the

- expected number of days of the year that are birthdays of at least one of these persons? (d) What is the expected number of days of the year that are birthdays of more than one of these persons? Ignore leap years for simplicity. [Hint: for (b), (c), (d), proceed as in No. 7.]
- *9. Problems 6, 7, and 8 are different versions of occupancy problems, which may be formulated generally as follows. Put n unnumbered tokens into m numbered boxes (see §3.3). What is the expected number of boxes that get exactly [or at least] k tokens? One can also ask for instance: what is the expected number of tokens that do not share its box with any other token? Answer these questions and rephrase them in the language of Problem 6, 7, or 8.
- *10. Using the occupancy model above, find the distribution of the tokens in the boxes, namely, the probabilities that exactly n_j tokens go into the jth box, $1 \le j \le m$. Describe this distribution in the language of Problem 6, 7, or 8.
 - 11. An automatic machine produces a defective item with probability 2%. When this happens an adjustment is made. Find the average number of good items produced between adjustments.
 - 12. Exactly one of six similar-looking keys is known to open a certain door. If you try them one after another, how many do you expect to have tried before the door is opened?
 - 13. One hundred electric bulbs are tested. If the probability of failure is p for each bulb, what are the mean and standard deviation of the number of failures? Assume stochastic independence of the bulbs.
- *14. Fifty persons queue up for chest X-ray examinations. Suppose there are four "positive" cases among them. What is the expected number of "negative" cases before the first positive case is spotted? [Hint: think of the four as partitioning walls for the others. Thus the problem is equivalent to finding the expected number of tokens in the first box under (IV') of §3.3.]
 - 15. There are N coupons numbered 1 to N in a bag. Draw one after another with replacement. (a) What is the expected number of drawings until the first coupon drawn is drawn again? (b)* What is the expected number of drawings until the first time a duplication occurs? [Hint: for (b) compute first the probability of no duplication in n drawings.]
- *16. In the problem above, what is the expected maximum coupon number in n drawings? The same question if the coupons are drawn without replacement. [Hint: find $P(\text{maximum} \leq k)$.]
 - 17. In Pólya's urn scheme with $c \ge -1$ (see §5.4):
 - (a) What is the expected number of red balls in n drawings?
 - (b) What is the expected number of red balls in the urn after the nth drawing (and putting back c balls)?

18. If $p_n \geq 0$ and $r_n = \sum_{k=n}^{\infty} p_k$ show that

$$\sum_{n=1}^{\infty} n p_n = \sum_{n=1}^{\infty} r_n$$

whether both series converge or diverge to $+\infty$. Hence if X is a random variable taking nonnegative integer values, we have

$$E(X) = \sum_{n=1}^{\infty} P(X \ge n).$$
 (6.6.1)

[Hint: Write $p_n = r_n - r_{n+1}$, rearrange the series (called Abel's method of summation in some calculus textbooks).]

- 19. Apply the formula (6.6.1) to compute the mean waiting time discussed in Example 8 of §4.4. Note that $P(X \ge n) = q^{n-1}, n \ge 1$.
- 20. Let X_1, \ldots, X_m be independent nonnegative integer-valued random variables all having the same distribution $\{p_n, n \geq 0\}$; and $r_n = \sum_{k=n}^{\infty} p_k$. Show that

$$E\{\min(X_1,\ldots,X_m)\} = \sum_{n=1}^{\infty} r_n^m.$$

[Hint: use No. 18.]

21. Let X be a nonnegative random variable with density function f. Show that if $r(u) = \int_{u}^{\infty} f(t) dt$, then

$$E(X) = \int_0^\infty P(X \ge u) \, du = \int_0^\infty r(u) \, du. \tag{6.6.2}$$

[Hint: this is the analogue of No. 18. Calculation with integrals is smoother than with sums.]

- 22. Apply formula (6.6.2) to an X with the exponential density $\lambda e^{-\lambda t}$.
- 23. The duration T of a certain type of telephone call is found to satisfy the relation

$$P(T > t) = ae^{-\lambda t} + (1 - a)e^{-\mu t}, \quad t > 0,$$

where $0 \le a \le 1, \lambda > 0, \mu > 0$ are constants determined statistically. Find the mean and variance of T. [Hint: for the mean a quick method is to use No. 21.]

24. Suppose that the "life" of an electronic device has the exponential density $\lambda e^{-\lambda t}$ in hours. Knowing that it has been in use for n hours, how much longer can it be expected to last? Compare this with its initial life expectancy. Do you see any contradiction?

- 25. Let five devices described above be tested simultaneously. (a) How long can you expect before one of them fails? (b) How long can you expect before all of them fall?
- 26. The average error committed in measuring the diameter of a circular disk is .2% and the area of the disk is computed from this measurement. What is the average percentage error in the area if we ignore the square of the percentage error in the diameter?
- 27. Express the mean and variance of aX + b in terms of those of X, where a and b are two constants. Apply this to the conversion of temperature from Centigrade to Fahrenheit:

$$F = \frac{9}{5}C + 32.$$

- 28. A gambler figures that he can always beat the house by doubling his bet each time to recoup any losses. Namely he will quit as soon as he wins, otherwise he will keep doubling his ante until he wins. The only drawback to this winning system is that he may be forced to quit when he runs out of funds. Suppose that he has a capital of \$150 and begins with a dollar bet, and suppose he has an even chance to win each time. What is the probability that he will quit winning, and how much will he have won? What is the probability that he will quit because he does not have enough left to double his last ante, and how much will he have lost in this case? What is his overall mathematical expectation by using this system? The same question if he will bet all his remaining capital when he can no longer double.
- 29. Pick n points at random in [0,1]. Find the expected value of the maximum, minimum, and range (= maximum minus minimum).
- 30. Consider n independent events A_j with $P(A_j) = p_j, 1 \leq j \leq n$. Let N denote the (random) number of occurrences among them. Find the generating function of N and compute E(N) from it.
- 31. Let $\{p_j, j \geq 0\}$ be a probability distribution and

$$u_k = \sum_{j=0}^k p_j,$$

$$g(z) = \sum_{k=0}^{\infty} u_k z^k.$$

Show that the power series converges for |z| < 1. As an example let S_n be as in Example 9 of §6.5, and $p_j = P\{S_n = j\}$. What is the meaning of u_k ? Find its generating function g.

32. It is also possible to define the generating function of a random variable that takes positive and negative values. To take a simple case, if

$$P(X = k) = p_k, \quad k = 0, \pm 1, \pm 2, \dots, \pm N,$$

then

$$g(z) = \sum_{k=-N}^{+N} p_k z^k$$

is a rational function of z, namely the quotient of two polynomials. Find g when $p_k = 1/(2N+1)$ above, which corresponds to the uniform distribution over the set of integers $\{-N, -(N-1), \ldots, -1, 0, +1, \ldots, N-1, N\}$. Compute the mean from g' for a check.

33. Let $\{X_j, 1 \leq j \leq n\}$ be independent random variables such that

$$X_j = \left\{ egin{array}{ll} 1 & \mbox{with probability } rac{1}{4}, \\ 0 & \mbox{with probability } rac{1}{2}, \\ -1 & \mbox{with probability } rac{1}{4}; \end{array}
ight.$$

and $S_n = \sum_{j=1}^n X_j$. Find the generating function of S_n in the sense of No. 32, and compute $P(S_n = 0)$ from it. As a concrete application, suppose A and B toss an unbiased coin n times each. What is the probability that they score the same number of heads? [This problem can also be solved without using generating function, by using formula (3.3.9).]

- *34. In the coupon-collecting problem of No. 15, let T denote the number of drawings until a complete set of coupons is collected. Find the generating function of T. Compute the mean from it for a beautiful check with (6.1.8). [Hint: Let T_j be the waiting time between collecting the (j-1)th and the jth new card; then it has a geometric distribution with $p_j = (N-j+1)/N$. The T_j 's are independent.]
 - 35. Let X and g be as in (6.5.1) and (6.5.2). Derive explicit formulas for the first four moments of X in terms of g and its derivatives.
 - 36. Denote the Laplace transform of X in (6.5.16) by $L(\lambda)$. Express the nth moment of X in terms of L and its derivatives.
 - 37. Find the Laplace transform corresponding to the density function f given below.
 - (a) f(u) = 1/c in (0, c), c > 0.
 - (b) $f(u) = 2u/c^2$ in (0, c), c > 0.

- (c) $f(u) = (\lambda^n u^{n-1}/(n-1)!) e^{-\lambda u}$ in $[0, \infty), \lambda > 0, n \ge 1$. [First verify that this is a density function! The corresponding distribution is called the *gamma distribution* $\Gamma(n; \lambda)$.]
- 38. Let $S_n = T_1 + \cdots + T_n$, where the T_j 's are independent random variables all having the density $\lambda e^{-\lambda t}$. Find the Laplace transform of S_n . Compare with the result in No. 37(c). We can now use Theorem 7 to identify the distribution of S_n .
- *39. Consider a population of N taxpayers paying various amounts of taxes, of which the mean is m and variance is σ^2 . If n of these are selected at random, show that the mean and variance of their total taxes are equal to

$$nm$$
 and $\frac{N-n}{N-1}n\sigma^2$,

respectively. [Hint: denote the amounts by X_1, \ldots, X_n and use (6.3.8). Some algebra may be saved by noting that $E(X_jX_k)$ does not depend on n, so it can be determined when n = N, but this trick is by no means necessary.]

- 40. Prove Theorem 1 by the method used in the proof of Theorem 2. Do the density case as well.
- 41. Let $a(\cdot)$ and $b(\cdot)$ be two probability density functions and define their convolution $c(\cdot)$ as follows:

$$c(v) = \int_{-\infty}^{\infty} a(u)b(v-u) du, \quad -\infty < v < \infty;$$

cf. (6.5.8). Show that $c(\cdot)$ is also a probability density function, often denoted by a*b.

- *42. If $a(u) = \lambda e^{-\lambda u}$ for $u \ge 0$, find the convolution of $a(\cdot)$ with itself. Find by induction the *n*-fold convolution $\underbrace{a*a*\cdots*a}_{n \text{ times}}$. [Hint: the result is given in No. 37(c).]
 - 43. Prove Theorem 4 for nonnegative integer-valued random variables by using generating functions. [Hint: express the variance by generating functions as in No. 35 and then use Theorem 6.]
- 44. Prove the analogues of Theorem 6 for Laplace and Fourier transforms.
- 45. Consider a sequence of independent trials each having probability p for success and q for failure. Show that the probability that the nth success is preceded by exactly j failures is equal to

$$\binom{n+j-1}{j}p^nq^j.$$

*46. Prove the formula

$$\sum_{j+k=l} \binom{m+j-1}{j} \binom{n+k-1}{k} = \binom{m+n+l-1}{l}$$

where the sum ranges over all $j \ge 0$ and $k \ge 0$ such that j + k = l. [Hint: this may be more recognizable in the form

$$\sum_{j+k=l} {\binom{-m}{j}} {\binom{-n}{k}} = {\binom{-m-n}{l}};$$

cf. (3.3.9). Use
$$(1-z)^{-m}(1-z)^{-n} = (1-z)^{-m-n}$$
.

*47. The general case of the problem of points (Example 6 of §2.2) is as follows. Two players play a series of independent games in which A has probability p, B has probability q = 1 - p of winning each game. Suppose that A needs m and B needs n more games to win the series. Show that the probability that A will win is given by either one of the expressions below:

(i)

$$\sum_{k=m}^{m+n-1} {m+n-1 \choose k} p^k q^{m+n-1-k};$$

(ii)

$$\sum_{k=0}^{n-1} {m+k-1 \choose k} p^m q^k.$$

The solutions were first given by Montmort (1678–1719). [Hint: solution (i) follows at once from Bernoulli's formula by an obvious interpretation. This is based on the idea (see Example 6 of §2.2) to complete m+n-1 games even if A wins before the end. Solution (ii) is based on the more natural idea of terminating the series as soon as A wins m games before B wins n games. Suppose this happens after exactly m+k games, then A must win the last game and also m-1 among the first m+k-1 games, and $k \le n-1$.]

*48. Prove directly that the two expressions (i) and (ii) given in No. 47 are equal. [Hint: one can do this by induction on n, for fixed m; but a more interesting method is suggested by comparison of the two ideas

involved in the solutions. This leads to the expansion of (ii) into

$$\begin{split} &\sum_{k=0}^{n-1} \binom{m+k-1}{k} p^m q^k (p+q)^{n-1-k} \\ &= \sum_{k=0}^{n-1} \binom{m+k-1}{k} p^m q^k \sum_{j=0}^{n-k-1} \binom{n-1-k}{j} p^{n-1-k-j} q^j \\ &= \sum_{l=0}^{n-1} p^{m+n-1-l} q^l \sum_{j+k=l} \binom{m+k-1}{k} \binom{n-k-1}{j}; \end{split}$$

now use No. 46. Note that the equality relates a binomial distribution to a negative binomial distribution.]

Poisson and Normal Distributions

7.1. Models for Poisson distribution

The Poisson distribution is of great importance in theory and in practice. It has the added virtue of being a simple mathematical object. We could have introduced it at an earlier stage of the book, and the reader was alerted to this in §4.4. However, the belated entrance will give it more prominence, as well as a more thorough discussion than would be possible without the benefit of the last two chapters.

Fix a real positive number α and consider the probability distribution $\{a_k, k \in \mathbb{N}^0\}$, where \mathbb{N}^0 is the set of all nonnegative integers, given by

$$a_k = \frac{e^{-\alpha}}{k!} \alpha^k. (7.1.1)$$

We must first verify that

$$\sum_{k=0}^{\infty} a_k = e^{-\alpha} \sum_{k=0}^{\infty} \frac{\alpha^k}{k!} = e^{-\alpha} \cdot e^{\alpha} = 1,$$

where we have used the Taylor series of e^{α} . Let us compute its mean as

well:

$$\begin{split} \sum_{k=0}^{\infty} k a_k &= e^{-\alpha} \sum_{k=0}^{\infty} k \frac{\alpha^k}{k!} = e^{-\alpha} \alpha \sum_{k=1}^{\infty} \frac{\alpha^{k-1}}{(k-1)!} \\ &= e^{-\alpha} \alpha \sum_{k=0}^{\infty} \frac{\alpha^k}{k!} = e^{-\alpha} \alpha e^{+\alpha} = \alpha. \end{split}$$

[This little summation has been spelled out since I have found that students often do not learn such problems of "infinite series" from their calculus course.] Thus the parameter α has a very specific meaning indeed. We shall call the distribution in (7.1.1) the *Poisson distribution with parameter* α . It will be denoted by $\pi(\alpha)$, and the term with subscript k by $\pi_k(\alpha)$. Thus if X is a random variable having this distribution, then

$$P(X = k) = \pi_k(\alpha) = \frac{e^{-\alpha}}{k!} \alpha^k, k \in \mathbb{N}^0;$$

and

$$E(X) = \alpha. \tag{7.1.2}$$

Next, let us find the generating function g as defined in §6.5. We have, using Taylor's series for $e^{\alpha z}$ this time:

$$g(z) = \sum_{k=0}^{\infty} a_k z^k = e^{-\alpha} \sum_{k=0}^{\infty} \frac{\alpha^k}{k!} z^k = e^{-\alpha} e^{\alpha z} = e^{\alpha(z-1)}.$$
 (7.1.3)

This is a simple function and can be put to good use in calculations. If we differentiate it twice, we get

$$q'(z) = \alpha e^{\alpha(z-1)}, q''(z) = \alpha^2 e^{\alpha(z-1)}.$$

Hence by (6.5.6),

$$E(X) = g'(1) = \alpha,$$

$$E(X^{2}) = g'(1) + g''(1) = \alpha + \alpha^{2},$$

$$\sigma^{2}(X) = \alpha.$$
(7.1.4)

So the variance as well as the mean are equal to the parameter α (see below for an explanation).

Mathematically, the Poisson distribution can be derived in a number of significant ways. One of these is a limiting scheme via the binomial distribution. This is known historically as Poisson's limit law and will be discussed

first. Another way, that of adding exponentially distributed random variables, is the main topic of the next section.

Recall the binomial distribution B(n; p) in §4.4 and write

$$B_k(n;p) = \binom{n}{k} p^k (1-p)^{n-k}, \quad 0 \le k \le n.$$
 (7.1.5)

We shall allow p to vary with n; this means only that we put $p = p_n$ in the above. Specifically, we take

$$p_n = \frac{\alpha}{n}, \quad n \ge 1. \tag{7.1.6}$$

We are therefore considering the sequence of binomial distributions $B(n; \alpha/n)$, a typical term of which is given by

$$B_k\left(n; \frac{\alpha}{n}\right) = \binom{n}{k} \left(\frac{\alpha}{n}\right)^k \left(1 - \frac{\alpha}{n}\right)^{n-k}, \quad 0 \le k \le n. \tag{7.1.7}$$

For brevity let us denote this by $b_k(n)$. Now fix k and let n go to infinity. It turns out that $b_k(n)$ converges for every k and can be calculated as follows. To begin at the beginning, take k = 0: then we have

$$\lim_{n \to \infty} b_0(n) = \lim_{n \to \infty} \left(1 - \frac{\alpha}{n} \right)^n = e^{-\alpha}.$$
 (7.1.8)

This is one of the fundamental formulas for the exponential function which you ought to remember from calculus. An easy way to see it is to take natural logarithm and use the Taylor series $\log (1-x) = -\sum_{n=1}^{\infty} x^n/n$:

$$\log\left(1 - \frac{\alpha}{n}\right)^n = n\log\left(1 - \frac{\alpha}{n}\right) = n\left\{-\frac{\alpha}{n} - \frac{\alpha^2}{2n^2} - \cdots\right\}$$

$$= -\alpha - \frac{\alpha^2}{2n} - \cdots.$$
(7.1.9)

When $n \to \infty$ the last-written quantity converges to $-\alpha$, which is $\log e^{-\alpha}$. Hence (7.1.8) may be verified by taking logarithms and expanding into power series, a method very much in use in applied mathematics. A rigorous proof must show that the three dots at the end of (7.1.9) above can indeed be overlooked; see Exercise 18.

To proceed, we take the ratio of consecutive terms in (7.1.7):

$$\frac{b_{k+1}(n)}{b_k(n)} = \frac{n-k}{k+1} \left(\frac{\alpha}{n}\right) \left(1 - \frac{\alpha}{n}\right)^{-1} = \frac{\alpha}{k+1} \left[\left(\frac{n-k}{n}\right) \left(1 - \frac{\alpha}{n}\right)^{-1} \right].$$

The two factors within the square brackets above both converge to 1 as $n \to \infty$; hence

$$\lim_{n \to \infty} \frac{b_{k+1}(n)}{b_k(n)} = \frac{\alpha}{k+1}.$$
 (7.1.10)

Starting with (7.1.8), and using (7.1.10) for $k = 0, 1, 2, \ldots$, we obtain

$$\lim_{n \to \infty} b_1(n) = \frac{\alpha}{1} \lim_{n \to \infty} b_0(n) = \alpha e^{-\alpha},$$

$$\lim_{n \to \infty} b_2(n) = \frac{\alpha}{2} \lim_{n \to \infty} b_1(n) = \frac{\alpha^2}{1 \cdot 2} e^{-\alpha},$$

$$\lim_{n \to \infty} b_k(n) = \frac{\alpha}{k} \lim_{n \to \infty} b_{k-1}(n) = \frac{\alpha^k}{1 \cdot 2 \cdots k} e^{-\alpha}.$$

These limit values are the successive terms of $\pi(\alpha)$. Therefore we have proved Poisson's theorem in its simplest form as follows.

Poisson's limit law:

$$\lim_{n\to\infty} B_k\left(n; \frac{\alpha}{n}\right) = \pi_k(\alpha), \quad k \in \mathbb{N}^0.$$

This result remains true if the α/n on the left side above is replaced by α_n/n , where $\lim_n \alpha_n = \alpha$. In other words, instead of taking $p_n = \alpha/n$ as we did in (7.1.6), so that $np_n = \alpha$, we may take $p_n = \alpha_n/n$, so that $np_n = \alpha_n$ and

$$\lim_{n \to \infty} n p_n = \lim_n \alpha_n = \alpha. \tag{7.1.11}$$

The derivation is similar to the above except that (7.1.8) is replaced by the stronger result below: if $\lim_{n\to\infty} \alpha_n = \alpha$, then

$$\lim_{n \to \infty} \left(1 - \frac{\alpha_n}{n} \right)^n = e^{-\alpha}. \tag{7.1.12}$$

With this improvement, we can now enunciate the theorem in a more pragmatic form as follows. A binomial probability $B_k(n;p)$, when n is large compared with np which is nearly α , may be approximated by $\pi_k(\alpha)$, for modest values of k. Recall that np is the mean of B(n;p) (see §4.4); it is no surprise that its approximate value α should also be the mean of the approximate Poisson distribution, as we have seen under (7.1.2). Similarly, the variance of B(n;p) is $npq = n(\alpha/n) [1 - (\alpha/n)]$ for $p = \frac{\alpha}{n}$; as $n \to \infty$ the limit is also α as remarked under (7.1.4).

The mathematical introduction of the Poisson distribution is thus done. The limiting passage from the binomial scheme is quite elementary, in contrast to what will be done in §7.3 below. But does the condition (7.1.6),

or the more relaxed (7.1.11), make sense in any real situation? The astonishing thing here is that a great variety of natural and manmade random phenomena are found to fit the pattern nicely. We give four examples to illustrate the ways in which the scheme works to a greater or lesser degree.

Example 1. Consider a *rare* event, namely one with small probability p of occurrence. For instance, if one bets on a single number at roulette, the probability of winning is equal to $1/37 \approx .027$, assuming that the 36 numbers and one "zero" are equally likely. [The roulette wheels in Monte Carlo have a single "zero," but those in Las Vegas have "double zeros."] If we do this 37 times, we can "expect" to win once. (Which theorem says this?) But we can also compute the probabilities that we win no time, once, twice, etc. The exact answers are of course given by the first three terms of B(37; 1/37):

$$\left(1 - \frac{1}{37}\right)^{37},$$

$$\left(\frac{37}{1}\right) \left(1 - \frac{1}{37}\right)^{36} \frac{1}{37} = \left(1 - \frac{1}{37}\right)^{36},$$

$$\left(\frac{37}{2}\right) \left(1 - \frac{1}{37}\right)^{35} \frac{1}{37^2} = \frac{36}{2 \times 37} \left(1 - \frac{1}{37}\right)^{35}.$$

If we set

$$c = \left(1 - \frac{1}{37}\right)^{37} \approx .363,$$

then the three numbers above are

$$c, \quad \frac{37}{36}c, \quad \frac{37}{36} \times \frac{1}{2}c.$$

Hence if we use the approximation $e^{-1} \approx .368$ for c, committing thereby an error of 1.5%; and furthermore "confound" 37/36 with 1, committing another error of 3%, but in the opposite direction, we get the first three terms of $\pi(1)$, namely:

$$e^{-1}$$
, e^{-1} , $\frac{1}{2}e^{-1}$.

Further errors will be compounded if we go on, but some may balance others. We may also choose to bet, say, 111 times $(111 = 37 \times 3)$ on a single number, and vary it from time to time as gamblers usually do at a

roulette table. The same sort of approximation will then yield

$$\left(1 - \frac{1}{37}\right)^{111} = c^3 \approx e^{-3},$$

$$\frac{111}{37} \left(1 - \frac{1}{37}\right)^{110} = \frac{37}{36} \times 3c^3 \approx 3e^{-3},$$

$$\frac{111 \times 110}{2} \left(1 - \frac{1}{37}\right)^{109} = \frac{111 \times 110}{36 \times 36} \frac{1}{2}c^3 \approx \frac{9}{2}e^{-3},$$

etc. Here of course c^3 is a worse approximation of e^{-3} than c is of e^{-1} . Anyway it should be clear that we are simply engaged in more or less crude but handy numerical approximations, without going to any limit. For no matter how small p is, so long as it is fixed as in this example, np will of course go to infinity with n, and the limiting scheme discussed above will be wide of the mark when n is large enough. Nevertheless a reasonably good approximation can be obtained for values of n and p such that np is relatively small compared with n. It is just a case of pure and simple numerical approximation, but many such applications have been made to various rare events. In fact, the Poisson law was very popular at one time under the name of "the law of small numbers." Well-kept statistical data such as the number of Prussian cavalry men killed each year by a kick from a horse, or the number of child suicides in Prussia, were cited as typical examples of this remarkable distribution (see [Keynes]).

Example 2. Consider the card-matching problem in §6.2. If a person who claims ESP (extrasensory perception) is a fake and is merely trying to match the cards at random, will his average score be better or worse when the number of cards is increased? Intuitively, two opposite effects are apparent. On one hand, there will be more cards to score; on the other, it will be harder to score each. As it turns out (see §6.2) these two effects balance each other so nicely that the expected number is equal to 1 irrespective of the number of cards! Here is an ideal setting for (7.1.6) with $\alpha = 1$. In fact, we can make it conform exactly to the previous scheme by allowing duplication in the guessing. That is, if we think of a deck of n cards laid face down on the table, we are allowed to guess them one by one with total forgetfulness. Then we can guess each card to be any one of the n cards, with equal probability 1/n, and independently of all other guesses. The probability of exactly k matches is then given by (7.1.7) with $\alpha = 1$, and so the Poisson approximation $\pi_k(1)$ applies if n is large.

This kind of matching corresponds to sampling with replacement. It is not a realistic model when two decks of cards are matched against each other. There is then mutual dependence between the various guesses and the binomial distribution above of course does not apply. But it can be shown that when n is large the effect of dependence is small, as follows. Let the probability of "no match" be q_n when there are n cards to be matched. We see in Example 4 of §6.2 that

$$q_n \approx e^{-1}$$

is an excellent approximation even for moderate values of n. Now an easy combinatorial argument (Exercise 19) shows that the probability of exactly k matches is equal to

$$\binom{n}{k} \frac{1}{(n)_k} q_{n-k} = \frac{1}{k!} q_{n-k}.$$
 (7.1.13)

Hence for fixed k, this converges to $(1/k!)e^{-1} = \pi_k(1)$.

Example 3. The Poisson law in a spatial distribution is typified by the counting of particles in a sort of "homogeneous chaos." For instance, we may count the number of virus particles with a square grid under the microscope. Suppose that the average number per small square is μ and that there are N squares in the grid. The virus moves freely about in such a way that its distribution over the grid may be approximated by the "tokens in boxes" model described under (I') in §3.3. Namely, there are μN particles to be placed into the N squares, and each particle can go into any of the squares with probability 1/N, independently of each other. Then the probability of finding exactly k particles in a given square is given by the binomial distribution:

$$B_k\left(\mu N; \frac{1}{N}\right) = \binom{\mu N}{k} \left(\frac{1}{N}\right)^k \left(1 - \frac{1}{N}\right)^{\mu N - k}.$$

Now we should imagine that the virus specimen under examination is part of a much larger specimen with the same average spatial proportion μ . In practice, this assumption is reasonably correct when, for example, a little blood is drawn from a sick body. It is then legitimate to approximate the above probability by $\pi_k(\mu)$ when N is large. The point here is that the small squares in which the counts are made remain fixed in size, but the homogeneity of space permits a limiting passage when the number of such squares is multiplied.

A grim example of the spatial scheme is furnished by the counting of flying-bomb hits on the south of London during World War II. The area was divided into N=576 squares each of 1/4 square mile, and μ was found statistically to be about .930. The table below shows the actual counts N_k and the Poisson approximations $\pi_k(\mu)$ with $\mu=.9323$. The close fit in this case might be explained by the deliberate randomness of the attacks which

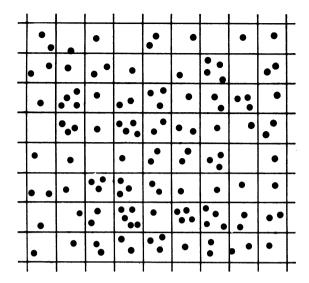


Figure 26

justified the binomial model above.

k	0	1	2	3	4	≥ 5
$\overline{N_k}$	229	211	93	35	7	1
$N\pi_k$	226.74	211.39	98.54	30.62	7.14	1.59

Example 4. In a large class of applications, time plays the role of space in the preceding example. If random occurrences are distributed over a period of time in such a way that their number per unit time may be supposed to be fairly constant over the period, then the Poisson scheme will operate with time acting as the medium for the homogeneous chaos. One could repeat the multiplication argument in Example 3 with time substituting for space, but here it is perhaps more plausible to subdivide the time. Suppose, for example, some cosmic ray impacts are registered on a geiger counter at the average rate of α per second. Then the probability of a register in a small time interval δ is given by $\alpha \, \delta + \mathrm{o}(\delta)$, where the "little-o" term represents an error term that is of smaller order of magnitude than δ , or roughly, "very small." Now divide the time interval [0,t] into N equal parts, so that the probability of a counter register in each subinterval is

$$\frac{\alpha t}{N} + \mathrm{o}\left(\frac{t}{N}\right),$$

with $\delta = t/N$ above. Of course, $\alpha t/N$ is much smaller than 1 when t is fixed and N is large. Let us first assume that for large enough values of N, the probability of more than one register in any small subinterval may

be neglected, so that we may suppose that the number of impacts received in each of the N subintervals is either 0 or 1. These numbers can then be treated as Bernoullian random variables taking the values 0 and 1 with probabilities $1-(\alpha t/N)$ and $\alpha t/N$, respectively. Finally we assume that they are independent of each other. This assumption can be justified on empirical grounds; for a deeper analysis in terms of the Poisson process, see the next section. Under these assumptions it is now clear that the probability of receiving exactly k impacts in the entire period [0,t] is given by the binomial $B_k(N;\alpha t/N)$; in fact, the total number registered in [0,t] is just the sum of N independent Bernoullian random variables described above. (See Example 9 of §4.4.) Since N is at our disposal and may be made arbitrarily large, in the limit we get $\pi_k(\alpha t)$. Thus is this case the validity of the Poisson scheme may be attributed to the infinite subdivisibility of time. The basic assumption concerning the independence of actions in disjoint subintervals will be justified in Theorem 2 of the following section.

*7.2. Poisson process

For a deeper understanding of the Poisson distribution we will construct a model in which it takes its proper place. The model is known as the *Poisson process* and is a fundamental stochastic process.

Consider a sequence of independent positive random variables all of which have the exponential density $\alpha e^{-\alpha t}$, $\alpha > 0$; see Example 3 of §6.2. Let them be denoted by T_1, T_2, \ldots so that for each j,

$$P(T_j \le t) = 1 - e^{-\alpha t}, \quad P(T_j > t) = e^{-\alpha t}, \quad t \ge 0.$$
 (7.2.1)

Since they are independent, we have for any nonnegative t_1, \ldots, t_n

$$P(T_1 > t_1, \dots, T_n > t_n) = P(T_1 > t_1) \cdots P(T_n > t_n)$$

= $e^{-\alpha(t_1 + \dots + t_n)}$

This determines the joint distribution of the T_j 's although we have given the "tail probabilities" for obvious simplicity. Examples of such random variables have been discussed before. For instance, they may be the *interarrival* times between vehicles in a traffic flow, or between claims received by an insurance company (see Example 5 of §4.2). They can also be the durations of successive telephone calls, or sojourn times of atoms at a specific energy level. Since

$$E(T_j) = \frac{1}{\alpha},\tag{7.2.2}$$

it is clear that the smaller α is, the longer the average *inter-arrival*, or waiting, or holding time. For instance, if T is the interarrival time between

automobiles at a checkpoint, then the corresponding α must be much larger on a Los Angeles freeway than in a Nevada desert. In this particular case α is also known as the *intensity of the flow*, in the sense that heavier traffic means a higher intensity, as every driver knows from his or her nerves.

Now let us put $S_0 = 0$ and for $n \ge 1$:

$$S_n = T_1 + \dots + T_n. (7.2.3)$$

Then by definition S_n is the waiting time until the nth arrival; and the event $\{S_n \leq t\}$ means that the nth arrival occurs before time t. [We shall use the preposition "before" loosely to mean "before or at" (time t). The difference can often be overlooked in continuous-time models but must be observed in discrete time.] Equivalently, this means "the total number of arrivals in the time interval [0,t]" is at least n. This kind of dual point of view is very useful, so we will denote the number just introduced by N(t). We can then record the assertion as follows:

$$\{N(t) \ge n\} = \{S_n \le t\}. \tag{7.2.4}$$

Like S_n , N(t) is also a random variable: $N(t,\omega)$ with the ω omitted from the notation as in $T_j(\omega)$. If you still remember our general discussion of random variables as functions of a sample point ω , now is a good time to review the situation. What is ω here? Just as in the examples of §4.2, each ω may be regarded as a possible record of the traffic flow or insurance claims or telephone service or nuclear transition. More precisely, N(t) is determined by the whole sequence $\{T_j, j \geq 1\}$ and depends on ω through the T_j 's. In fact, taking the difference of both sides in the equations (7.2.4) for n and n+1, we obtain

$$\{N(t) = n\} = \{S_n \le t\} - \{S_{n+1} \le t\} = \{S_n \le t < S_{n+1}\}. \tag{7.2.5}$$

The meaning of this new equation is clear from a direct interpretation: there are exactly n arrivals in [0,t] if and only if the nth arrival occurs before t but the (n+1)st occurs after t. For each value of t, the probability distribution of the random variable N(t) is therefore given by

$$P\{N(t) = n\} = P\{S_n \le t\} - P\{S_{n+1} \le t\}, \quad n \in \mathbb{N}^0.$$
 (7.2.6)

Observe the use of our convention $S_0 = 0$ in the above. We proceed to show that this is the Poisson distribution $\pi(\alpha t)$.

We shall calculate the probability $P\{S_n \leq t\}$ via the Laplace transform of S_n (see §6.5). The first step is to find the Laplace transform $L(\lambda)$ of each T_j , which is defined since $T_j \geq 0$. By (6.5.16) with $f(u) = \alpha e^{-\alpha u}$, we have

$$L(\lambda) = \int_0^\infty e^{-\lambda u} \alpha e^{-\alpha u} du = \frac{\alpha}{\alpha + \lambda}.$$
 (7.2.7)

Since the T_j 's are independent, an application of Theorem 7 of §6.5 yields the Laplace transform of S_n :

$$L(\lambda)^n = \left(\frac{\alpha}{\alpha + \lambda}\right)^n. \tag{7.2.8}$$

To get the distribution or density function of S_n from its Laplace transform is called an inversion problem; and there are tables of common Laplace transforms from which you can look up the *inverse*, namely the distribution or density associated with it. In the present case the answer has been indicated in Exercise 38 of Chapter 6. However, here is a trick that leads to it quickly. The basic formula is

$$\int_{0}^{\infty} e^{-xt} dt = \frac{1}{x}, \quad x > 0.$$
 (7.2.9)

Differentiating both sides n-1 times, which is easy to do, we obtain

$$\int_0^\infty (-t)^{n-1} e^{-xt} dt = \frac{(-1)^{n-1} (n-1)!}{x^n},$$

or

$$\frac{1}{(n-1)!} \int_0^\infty t^{n-1} e^{-xt} dt = \frac{1}{x^n}.$$
 (7.2.10)

Substituting $\alpha + \lambda$ for x in the above and multiplying both sides by α^n , we deduce

$$\int_0^\infty \frac{\alpha^n}{(n-1)!} u^{n-1} e^{-\alpha u} e^{-\lambda u} du = \left(\frac{\alpha}{\alpha+\lambda}\right)^n.$$

Thus if we put

$$f_n(u) = \frac{\alpha^n}{(n-1)!} u^{n-1} e^{-\alpha u}, \qquad (7.2.11)$$

we see that f_n is the density function for the Laplace transform in (7.2.8), namely that of S_n .* Hence we can rewrite the right side of (7.2.6) explicitly as

$$\int_0^t f_n(u) \, du - \int_0^t f_{n+1}(u) \, du. \tag{7.2.12}$$

^{*}Another derivation of this is contained in Exercise 42 of Chapter 6.

To simplify this we integrate the first integral by parts as indicated below:

$$\begin{split} \frac{\alpha^n}{(n-1)!} \int_0^t e^{-\alpha u} u^{n-1} \, du &= \frac{\alpha^n}{(n-1)!} \left\{ \frac{u^n}{n} e^{-\alpha u} \Big|_0^t + \int_0^t \frac{u^n}{n} e^{-\alpha u} \alpha \, du \right\} \\ &= \frac{\alpha^n}{n!} t^n e^{-\alpha t} + \frac{\alpha^{n+1}}{n!} \int_0^t u^n e^{-\alpha u} \, du. \end{split}$$

But the last-written integral is just the second integral in (7.2.12); hence the difference there is precisely $(\alpha^n/n!)t^ne^{-\alpha t} = \pi_n(\alpha t)$. For fixed n and α , this is the density function of the gamma distribution $\Gamma(n;\alpha)$; see p. 200. Let us record this as a theorem.

Theorem 1. The total number of arrivals in a time interval of length t has the Poisson distribution $\pi(\alpha t)$, for each t > 0.

The reader should observe that the theorem asserts more than has been proved. For in our formulation above we have implicitly chosen an initial instant from which time is measured, namely the zero time for the first arrival time T_1 . Thus the result was proved only for the total number of arrivals in the interval [0, t]. Now let us denote the number of arrivals in an arbitrary time interval [s, s+t] by N(s, s+t). Then it is obvious that

$$N(s, s + t) = N(s + t) - N(s)$$

in our previous notation, and N(0)=0. But we have yet to show that the distribution of N(s,s+t) is the same as N(0,t). The question becomes: if we start counting arrivals from time s on, will the same pattern of flow hold as from time 0 on? The answer is "yes" but it involves an essential property of the exponential distribution of the T_j 's. Intuitively speaking, if a waiting time such as T_j is broken somewhere in between, its duration after the break follows the original exponential distribution regardless of how long it has already endured before the break. This property is sometimes referred to as "lack of memory" and can be put in symbols: for any $s \geq 0$ and $t \geq 0$, we have

$$P(T > t + s \mid T > s) = P(T > t) = e^{-\alpha t};$$
 (7.2.13)

see Example 4 of $\S 5.1$. There is a converse: if a nonnegative random variable T satisfies the first equation in (7.2.13), then it must have an exponential distribution; see Exercise 41 of Chapter 5. Thus the lack of memory is characteristic of an exponential interarrival time.

We can now argue that the pattern of flow from time s on is the same as from 0 on. For the given instant s breaks one of the inter-arrival times, say T_k , into two stretches as shown below:

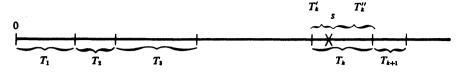


Figure 27

According to the above, the second stretch T''_k of the broken T_k has the same distribution as T_k , and it is *clearly* independent of all the succeeding T_{k+1}, T_{k+2}, \ldots . [The clarity is intuitive enough, but a formal proof takes some doing and is omitted.] Hence the new shifted interarrival times from s onward:

$$T_k'', T_{k+1}, T_{k+2}, \dots$$
 (7.2.14)

follow the same probability pattern as the original interarrival times beginning at 0:

$$T_1, T_2, T_3, \dots$$
 (7.2.15)

Therefore our previous analysis applies to the shifted flow as well as the original one. In particular, the number of arrivals in [s, s+t] must have the same distribution as that in [0, t]. This is the assertion of Theorem 1.

The fact that N(s, s+t) has the same distribution for all s is referred to as the *time-homogeneity* of the flow. Let us remember that this is shown under the assumption that the intensity α is constant for all time. In practice such an assumption is tenable only over specified periods of time. For example, in the case of traffic flow on a given highway, it may be assumed for the rush hour or from 2 a.m. to 3 a.m. with different values of α . However, for longer periods of time such as one day, an average value of α over 24 hours may be used. This may again vary from year to year, even week to week.

So far we have studied the number of arrivals in one period of time, of arbitrary length and origin. For a more complete analysis of the flow we must consider several such periods and their mutual dependence. In other words, we want to find the joint distribution of

$$N(s_1, s_1 + t_1), N(s_2, s_2 + t_2), N(s_3, s_3 + t_3), \dots,$$
 (7.2.16)

etc. The answer is given in the next theorem.

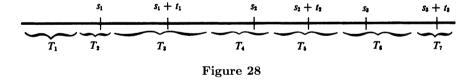
Theorem 2. If the intervals $(s_1, s_1+t_1), (s_2, s_2+t_2), \ldots$ are disjoint, then the random variables in (7.2.16) are independent and have the Poisson distributions $\pi(\alpha t_1), \pi(\alpha t_2) \ldots$

It is reasonable and correct to think that if we know the joint action of N over any arbitrary finite set of disjoint time intervals, then we know all about it in principle. Hence with Theorem 2 we shall be in full control of the process in question.

The proof of Theorem 2 depends again on the lack-of-memory property of the T_j 's. We will indicate the main idea here without going into formal details. Going back to the sequence in (7.2.14), where we put $s = s_2$, we now make the further observation that all the random variables there are not only independent of one another, but also of all those that precede s, namely:

$$T_1, \dots, T_{k-1}, T'_k.$$
 (7.2.17)

The fact that the two broken stretches T_k' and T_k'' are independent is a consequence of (7.2.13), whereas the independence of all the rest should be intuitively obvious because they have not been distributed by the break at s. [Again, it takes some work to justify the intuition.] Now the "past history" of the flow up to time s is determined by the sequence in (7.2.17), while its "future development" after s is determined by the sequence in (7.2.14). Therefore, relative to the "present" s, past and future are independent. In particular, $N(s_1, s_1 + t_1)$, which is part of the past, must be independent of $N(s_2, s_2 + t_2), N(s_3, s_3 + t_3), \ldots$, which are all part of the future. Repeating this argument for $s = s_3, s_4, \ldots$, the assertion of Theorem 2 follows.



We are now ready to give a general definition for the "flow" we have been discussing all along.

Definition of Poisson Process. A family of random variables $\{X(t)\}$, indexed by the continuous variable t ranging over $[0, \infty)$, is called a *Poisson process with parameter* (or *mean*) α iff it satisfies the following conditions:

- (i) X(0) = 0;
- (ii) the increments $X(s_i + t_i) X(s_i)$, over an arbitrary finite set of disjoint intervals $(s_i, s_i + t_i)$, are independent random variables;
- (iii) for each $s \ge 0$, $t \ge 0$, X(x+t) X(s) has the Poisson distribution $\pi(\alpha t)$.

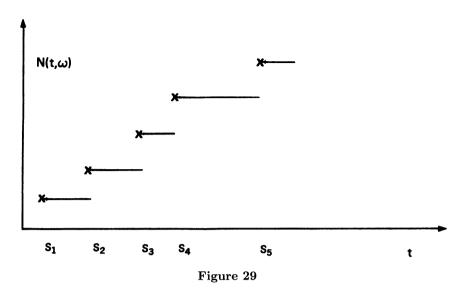
According to Theorems 1 and 2 above, the family $\{N(t), t \geq 0\}$ satisfies these conditions and therefore forms a Poisson process. Conversely, it can be shown that every Poisson process is representable as the N(t) above.

The concept of a stochastic process has already been mentioned in §§5.3–5.4, in connection with Pólya's urn model. The sequence $\{X_n, n \geq 1\}$ in Theorem 5 of §5.4 may well be called a Pólya process. In principle a stochastic process is just any family of random variables; but this is putting matters in an esoteric way. What is involved here goes back to the foundations of probability theory discussed in Chapters 2, 4, and 5. There are a sample space Ω with points ω , a probability measure P defined for certain sets of ω , a family of functions $\omega \to X_t(\omega)$ called random variables, and the process is concerned with the joint action or behavior of this family: the marginal and joint distributions, the conditional probabilities, the expectations, and so forth. Everything we have discussed (and are going to discuss) may be regarded as questions in stochastic processes, for in its full generality the term encompasses any random variable or sample set (via its indicator). But in its customary usage we mean a rather numerous and well-organized family governed by significant and useful laws. The preceding characterization of a Poisson process is a good example of this description.

As defined, $\omega \to N(t,\omega)$ is a random variable for each t, with the Poisson distribution $\pi(\alpha t)$. There is a dual point of view that is equally important in the study of a process, and that is the function $t \to N(t,\omega)$ for each ω . Such a function is called a sample function (path or trajectory). For example, in the case of telephone calls, to choose a sample point ω may mean to pick a day's record of the actual counts at a switchboard over a 24-hour period. This of course varies from day to day so the function $t \to N(t,\omega)$ gives only a sample (denoted by ω) of the telephone service. Its graph may look like Figure 29. The points of jumps are the successive arrival times $S_n(\omega)$, each jump being equal to 1, and the horizontal stretches indicate the interarrival times. So the sample function is a monotonically nondecreasing function that increases only by jumps of size 1 and is flat between jumps. Such a graphic is typical of the sample function of a Poisson process. If the flow is intense, then the points of jumps are crowded together.

The sequence $\{S_n, n \geq 1\}$ defined in (7.2.3) is also a stochastic process, indexed by n. A sample function $n \to S_n(\omega)$ for this process is an increasing sequence of positive numbers $\{S_1(\omega), S_2(\omega), \ldots, S_n(\omega), \ldots\}$. Hence it is often called a sample sequence. There is a reciprocal relation between this and the sample function $N(t,\omega)$ above. If we interchange the two coordinate axes, which can be done by turning the page 90°, and look at Figure 29 through the light from the back, we get the graph of $n \to S_n(\omega)$. Ignore the now vertical stretches except the lower endpoints, which indicate the values of S_n .

The following examples illustrate some of the properties of the Poisson distribution and process.



Example 5. Consider the number of arrivals in two disjoint time intervals: $X_1 = N(s_1, s_1 + t_1)$ and $X_2 = N(s_2, s_2 + t_2)$ as in (7.2.16). What is the probability that the total number $X_1 + X_2$ is equal to n?

By Theorem 2, X_1 and X_2 are independent random variables with the distributions $\pi(\alpha t_1)$ and $\pi(\alpha t_2)$, respectively. Hence

$$P(X_1 + X_2 = n) = \sum_{j+k=n} P(X_1 = j) P(X_2 = k)$$

$$= \sum_{j+k=n} \frac{e^{-\alpha t_1} (\alpha t_1)^j}{j!} \frac{e^{-\alpha t_2} (\alpha t_2)^k}{k!}$$

$$= \frac{e^{-\alpha (t_1 + t_2)}}{n!} \sum_{j=0}^n \binom{n}{j} (\alpha t_1)^j (\alpha t_2)^{n-j}$$

$$= \frac{e^{-\alpha (t_1 + t_2)}}{n!} (\alpha t_1 + \alpha t_2)^n = \pi_n (\alpha t_1 + \alpha t_2).$$

Namely, $X_1 + X_2$ is also Poissonian with parameter $\alpha t_1 + \alpha t_2$. The general proposition is as follows.

Theorem 3. Let X_j be independent random variables with Poisson distributions $\pi(\alpha_j), 1 \leq j \leq n$. Then $X_1 + \cdots + X_n$ has Poisson distribution $\pi(\alpha_1 + \cdots + \alpha_n)$.

This follows from an easy induction, but we can also make speedy use of generating functions. If we denote the generating function of X_i by g_{x_i} ,

then by Theorem 6 of §6.5:

$$g_{X_1 + \dots + X_n}(z) = g_{X_1}(z)g_{X_2}(z) \cdots g_{X_n}(z)$$

$$= e^{\alpha_1(z-1)}e^{\alpha_2(z-1)} \cdots e^{\alpha_n(z-1)}$$

$$= e^{(\alpha_1 + \dots + \alpha_n)(z-1)}.$$

Thus $X_1 + \cdots + X_n$ has the generating function associated with $\pi(\alpha_1 + \cdots + \alpha_n)$, and so by the uniqueness stated in Theorem 7 of §6.5 it has the latter as distribution.

Example 6. At a crossroad of America we watch cars zooming by bearing license plates of various states. Assume that the arrival process is Poissonian with intensity α , and that the probabilities of each car being from the states of California, Nevada ,and Arizona are, respectively, $p_1 = 1/25$, $p_2 = 1/100$, $p_3 = 1/80$. In a unit period of time what is the number of cars counted with these license plates?

We are assuming that if n cars are counted the distribution of various license plates follows a multinomial distribution $M(n; 50; p_1, \ldots, p_{50})$ where the first three p's are given. Now the number of cars passing in the period of time is a random variable N such that

$$P(N = n) = \frac{e^{-\alpha}}{n!} \alpha^n, \quad n = 0, 1, 2, \dots$$

Among these N cars, the number bearing the kth state license is also a random variable N_k ; of course,

$$N_1 + N_2 + \cdots + N_{50} = N.$$

The problem is to compute

$$P(N_1 = n_1, N_2 = n_2, N_3 = n_3)$$

for arbitrary n_1, n_2, n_3 . Let $q = 1 - p_1 - p_2 - p_3$; this is the probability of a license plate not being from one of the three indicated states. For $n \geq n_1 + n_2 + n_3$, the conditional probability under the hypothesis that N = n is given by the multinomial; hence

$$P(N_1 = n_1, N_2 = n_2, N_3 = n_3 \mid N = n) = \frac{n! \, p_1^{n_1} p_2^{n_2} p_3^{n_3} q^k}{n_1! \, n_2! \, n_3! \, k!},$$

where $k = n - n_1 - n_2 - n_3$. Using the formula for "total probability"

(5.2.3), we get

$$P(N_1 = n_1, N_2 = n_2, N_3 = n_3)$$

$$= \sum_{n} P(N = n)P(N_1 = n_1, N_2 = n_2, N_3 = n_3 \mid N = n)$$

$$= \sum_{k=0}^{\infty} \frac{e^{-\alpha}}{n!} \alpha^n \frac{n!}{n_1! n_2! n_3!} \frac{p_1^{n_1} p_2^{n_2} p_3^{n_3} q^k}{k!}.$$

Since $n_1 + n_2 + n_3$ is fixed and $n \ge n_1 + n_2 + n_3$, the summation above reduces to that with k ranging over all nonnegative integers. Now write in the above

$$e^{-\alpha} = e^{-\alpha(p_1 + p_2 + p_3)}e^{-\alpha q}, \quad \alpha^n = \alpha^{n_1 + n_2 + n_3}\alpha^k.$$

and take out the factors that do not involve the index of summation k. The result is

$$\frac{e^{-\alpha(p_1+p_2+p_3)}}{n_1! \, n_2! \, n_3!} (\alpha p_1)^{n_1} (\alpha p_2)^{n_2} (\alpha p_3)^{n_3} \sum_{k=0}^{\infty} \frac{e^{-\alpha q}}{k!} (\alpha q)^k$$
$$= \pi_{n_1}(\alpha p_1) \pi_{n_2}(\alpha p_2) \pi_{n_3}(\alpha p_3)$$

since the last-written sum equals 1. Thus the random variables N_1 , N_2 , N_3 are independent (why?) and have the Poisson distributions $\pi(\alpha p_1)$, $\pi(\alpha p_2)$, $\pi(\alpha p_3)$.

The substitution of the "fixed number n" in the multinomial $M(n;r;p_1,\ldots,p_r)$ by the "random number N" having a Poisson distribution is called in statistical methodology "randomized sampling." In the example here the difference is illustrated by either counting a fixed number of cars or counting whatever number of cars in a chosen time interval, or by some other selection method that allows a chance variation of the number. Which way of counting is more appropriate will in general depend on the circumstances and the information sought.

There is, of course, a general proposition behind the above example that may be stated as follows.

Theorem 4. Under randomized sampling from a multinomial population $M(n; r; p_1, ..., p_r)$ where the total number sampled is a Poisson random variable N with mean α , the numbers $N_1, ..., N_r$ of the various varieties obtained by the sampling become independent Poisson variables with means $\alpha p_1, ..., \alpha p_r$.

As an illustration of the finer structure of the Poisson process we will derive a result concerning the location of the jumps of its sample functions.

Let us begin with the remark that although (almost) all sample functions have infinitely many jumps in $(0, \infty)$, the probability that a jump occurs at any prescribed instant of time is equal to zero. For if t > 0 is fixed, then as $\delta \downarrow 0$ we have

$$P\{N(t+\delta) - N(t-\delta) \ge 1\} = 1 - \pi_0(\alpha, 2\delta) = 1 - e^{-2\alpha \delta} \to 0.$$

In particular, the number of jumps in an interval (t_1, t_2) has the same distribution whether the endpoints t_1 and t_2 are included or not. As before let us write $N(t_1, t_2)$ for this number. Now suppose N(0, t) = n for a given t, where $n \geq 1$, and consider an arbitrary subinterval (t_1, t_2) of (0, t). We have for $0 \leq j \leq n$

$$P\{N(t_1, t_2) = j; N(0, t) = n\}$$

$$= P\{N(t_1, t_2) = j; N(0, t_1) + N(t_2, t) = n - j\}.$$

Let $t_2 - t_1 = s$; then the sum of the lengths of the two intervals $(0, t_1)$ and (t_2, t) is equal to t - s. By property (ii) and Theorem 3 the random variable $N(0, t_1) + N(t_2, t)$ has the distribution $\pi(\alpha(t - s))$ and is independent of $N(t_1, t_2)$. Hence the probability above is equal to

$$e^{-\alpha s} \frac{(\alpha s)^j}{j!} e^{-\alpha(t-s)} \frac{(\alpha t - \alpha s)^{n-j}}{(n-j)!}.$$

Dividing this by $P\{N(0,t)=n\}=e^{-\alpha t}(\alpha t)^n/n!$, we obtain the conditional probability:

$$P\{N(t_1, t_2) = j \mid N(0, t) = n\} = \binom{n}{j} \left(\frac{s}{t}\right)^j \left(1 - \frac{s}{t}\right)^{n-j}.$$
 (7.2.18)

This is the binomial probability $B_i(n; s/t)$.

Now consider an arbitrary partition of (0,t) into a finite number of subintervals I_1, \ldots, I_l of lengths s_1, \ldots, s_l so that $s_1 + \cdots + s_l = t$. Let n_1, \ldots, n_l be arbitrary nonnegative integers with $n_1 + \cdots + n_l = n$. If we denote by $N(I_k)$ the number of jumps of the process in the interval I_k , then we have a calculation similar to the above:

$$P\{N(I_k) = n_k, 1 \le k \le l \mid N(0, t) = n\}$$

$$= \prod_{k=1}^{l} \frac{e^{-\alpha s_k} (\alpha s_k)^{n_k}}{n_k!} \left(e^{-\alpha t} \frac{(\alpha t)^n}{n!} \right)^{-1}$$

$$= \frac{n!}{n_1! \cdots n_l!} \prod_{k=1}^{l} \left(\frac{s_k}{t} \right)^{n_k}.$$
(7.2.19)

This is the multinomial distribution discussed in §6.4.

Let us pick n points at random in (0,t) and arrange them in nondecreasing order $0 < \xi_1 \le \xi_2 \le \cdots \le \xi_n < t$. Using the notation above let $\tilde{N}(I_k)$ denote the number of these points lying in I_k . It is not hard to see that the n-dimensional distribution of (ξ_1,\ldots,ξ_n) is uniquely determined by the distribution of $(\tilde{N}(I_1),\ldots,\tilde{N}(I_l))$ for all possible partitions of (0,t); for a rigorous proof of this fact see Exercise 26. In particular, if the n points are picked independently of one another and each is uniformly distributed in (0,t), then it follows from the discussion in §6.4 that the probability $P\{\tilde{N}(I_k)=n_k,1\le k\le l\}$ is given by the last term in (7.2.19). Therefore, under the hypothesis that there are exactly n jumps of the Poisson process in (0,t), the conditional distribution of the n points of jump is the same as if they were picked in the manner just described. This has been described as a sort of "homogeneous chaos."

7.3. From binomial to normal

From the point of view of approximating the binomial distribution B(n;p) for large values of n, the case discussed in §7.1 leading to the Poisson distribution is abnormal, because p has to be so small that np remains constant, or nearly so. The fact that many random phenomena follow this law rather nicely was not known in the early history of probability. One must remember that not only had radioactivity yet to be discovered, but neither the telephone nor automobile traffic existed as modern problems. On the other hand, counting heads by tossing coins or points by rolling dice, and the measurement of all kinds of physical and biological quantities were already done extensively. These led to binomial and multinomial distributions, and since computing machines were not available it became imperative to find manageable formulas for the probabilities. The normal way to approximate the binomial distribution

$$B_k(n;p) = \binom{n}{k} p^k (1-p)^{n-k}, \quad 0 \le k \le n, \tag{7.3.1}$$

is for a fixed value of p and large values of n. To illustrate by the simplest kind of example, suppose an unbiased coin is tossed 100 times; what is the probability of obtaining exactly 50 heads? The answer

$$\binom{100}{50} \frac{1}{2^{100}} = \frac{100!}{50! \, 50!} \, \frac{1}{2^{100}}$$

gives little satisfaction as we have no idea of the magnitude of this probability. Without some advanced mathematics (which will be developed presently), who can guess whether this is near 1/2 or 1/10 or 1/50?

Now it is evident that the key to such combinatorial formulas is the factorial n!, which just crops up everywhere. Take a look back at Chapter 3.

So the problem is to find a handy formula: another function $\chi(n)$ of n which is a good approximation for n! but of a simpler structure for computations. But what is "good"? Since n! increases very rapidly with n (see the short table in §3.2), it would be hopeless to make the difference $|n! - \chi(n)|$ small. [Does it really make a difference to have a million dollars or a million and three?] What counts is the ratio $n!/\chi(n)$, which should be close to 1. For two positive functions ψ and χ of the integer variable n, there is a standard notation:

$$\psi(n) \sim \chi(n)$$
 which means $\lim_{n \to \infty} \frac{\psi(n)}{\chi(n)} = 1.$ (7.3.2)

We say also that $\psi(n)$ and $\chi(n)$ are asymptotically equal (or equivalent) as $n \to \infty$. If so we also have

$$\lim_{n \to \infty} \frac{|\psi(n) - \chi(n)|}{\chi(n)} = 0$$

provided $\chi(n) > 0$ for large n; thus the difference $|\psi(n) - \chi(n)|$ is negligible in comparison with $\chi(n)$ or $\psi(n)$, though it may be large indeed in absolute terms. Here is a trivial example that you should have retained from a calculus course (under the misleading heading "indeterminate form"):

$$\psi(n) = 2n^2 + 10n - 100, \quad \chi(n) = 2n^2.$$

More generally, a polynomial in n is asymptotically equal to its highest term. Here, of course, we are dealing with something far more difficult: to find a simple enough $\chi(n)$ such that

$$\lim_{n \to \infty} \frac{n!}{\chi(n)} = 1.$$

Such a χ is given by *Stirling's formula* (see Appendix 2):

$$\chi(n) = \left(\frac{n}{e}\right)^n \sqrt{2\pi n} = n^{n+(1/2)} e^{-n} \sqrt{2\pi},\tag{7.3.3}$$

or more precisely

$$n! = \left(\frac{n}{e}\right)^n \sqrt{2\pi n} \, e^{\omega(n)}, \text{ where } \frac{1}{12(n+\frac{1}{2})} < \omega(n) < \frac{1}{12n}.$$
 (7.3.4)

You may think $\chi(n)$ is uglier looking than n!, but it is much easier to compute because powers are easy to compute. Here we will apply it at once to the little problem above. It does not pay to get involved in numericals at the beginning, so we will consider

$$\binom{2n}{n} \frac{1}{2^{2n}} = \frac{(2n)!}{n! \, n!} \, \frac{1}{2^{2n}}.\tag{7.3.5}$$

Substituting $\chi(n)$ and $\chi(2n)$ for n! and (2n)!, respectively, we see that this is asymptotically equal to

$$\frac{\left(\frac{2n}{e}\right)^{2n}\sqrt{4\pi n}}{\left(\frac{n}{e}\right)^{2n}2\pi n}\frac{1}{2^{2n}} = \frac{1}{\sqrt{\pi n}}.$$

In particular for n = 50, we get the desired answer $1/\sqrt{50\pi} = .08$ approximately. Try to do this by using logarithms on

$$\frac{(100)_{50}}{50!} \; \frac{1}{2^{100}}$$

and you will appreciate Stirling's formula more. We proceed at once to the slightly more general

$$\binom{2n}{n+k} \frac{1}{2^{2n}} = \frac{(2n)!}{(n+k)!(n-k)!} \frac{1}{2^{2n}},\tag{7.3.6}$$

where k is fixed. A similar application of (7.3.3) yields

$$\frac{\left(\frac{2n}{e}\right)^{2n}\sqrt{4\pi n} \cdot \frac{1}{2^{2n}}}{\left(\frac{n+k}{e}\right)^{n+k}\sqrt{2\pi(n+k)}\left(\frac{n-k}{e}\right)^{n-k}\sqrt{2\pi(n-k)}} \\
= \frac{n^{2n}}{(n+k)^{n+k}(n-k)^{n-k}}\sqrt{\frac{n}{\pi(n+k)(n-k)}} \\
= \left(\frac{n}{n+k}\right)^{n+k}\left(\frac{n}{n-k}\right)^{n-k}\sqrt{\frac{n}{\pi(n^2-k^2)}}.$$

Clearly the last-written factor is asymptotically equal to $1/\sqrt{\pi n}$. As for the two preceding ones, it follows from (7.1.8) that

$$\lim_{n \to \infty} \left(\frac{n}{n+k} \right)^{n+k} = \lim_{n \to \infty} \left(1 - \frac{k}{n+k} \right)^{n+k} = e^{-k},$$

$$\lim_{n \to \infty} \left(\frac{n}{n-k} \right)^{n-k} = \lim_{n \to \infty} \left(1 + \frac{k}{n-k} \right)^{n-k} = e^{k}.$$
(7.3.7)

Hence the asymptotic value of (7.3.6) is

$$e^{-k}e^k\frac{1}{\sqrt{\pi n}} = \frac{1}{\sqrt{\pi n}},$$

exactly as in (7.3.5), which is the particular case k = 0.

As a consequence, for any fixed number l, we have

$$\lim_{n \to \infty} \sum_{k=-l}^{l} {2n \choose n+k} \frac{1}{2^{2n}} = 0 \tag{7.3.8}$$

because each term in the sum has limit 0 as $n \to \infty$ as just shown, and there is only a fixed number of terms. Now if we remember Pascal's triangle (3.3.5), the binomial coefficients $\binom{2n}{n+k}$, $-n \le k \le n$, assume their maximum value $\binom{2n}{n}$ for the middle term k=0 and decrease as |k| increases (see Exercise 6). According to (7.3.8), the sum of a fixed number of terms centered around the middle term approaches zero; hence a fortiori the sum of any fixed number of terms will also approach zero, namely for any fixed a and b with a < b, we have

$$\lim_{n \to \infty} \sum_{i=a}^{b} {2n \choose j} \frac{1}{2^{2n}} = 0.$$

Finally, this result remains true if we replace 2n by 2n + 1 above, because the ratio of corresponding terms

$$\binom{2n+1}{j} \frac{1}{2^{2n+1}} / \binom{2n}{j} \frac{1}{2^{2n}} = \frac{2n+1}{2n+1-j} \cdot \frac{1}{2}$$

approaches 1/2, which does not affect the zero limit. Now let us return to the probability meaning of the terms, and denote as usual by S_n the number of heads obtained in n tosses of the coin. The result then asserts that for any fixed numbers a and b, we have

$$\lim_{n \to \infty} P(a \le S_n \le b) = 0. \tag{7.3.9}$$

Observe that there are n+1 possible values for S_n , whereas if the range [a,b] is fixed irrespective of n, it will constitute a negligible fraction of n when n is large. Thus the result (7.3.9) is hardly surprising, though certainly disappointing.

It is clear that in order to "catch" a sufficient number of possible values of S_n to yield a nonzero limit probability, the range allowed must increase to infinity with n. Since we saw that the terms near the middle are of the order of magnitude $1/\sqrt{n}$, it is plausible that the number of terms needed will be of the order of magnitude \sqrt{n} . More precisely, it turns out that for each l,

$$P\left(\frac{n}{2} - l\sqrt{n} \le S_n \le \frac{n}{2} + l\sqrt{n}\right) = \sum_{|j - \frac{n}{2}| \le l\sqrt{n}} \binom{n}{j} \frac{1}{2^n}$$
 (7.3.10)

will have a limit strictly between 0 and 1 as $n \to \infty$. Here the range for S_n is centered around the middle value n/2 and contains about $2l\sqrt{n}$ terms. When n is large this is still only a very small fraction of n, but it increases just rapidly enough to serve our purpose. The choice of \sqrt{n} rather than say $n^{1/3}$ or $n^{3/5}$ is crucial and is determined by a rather deep bit of mathematical analysis, which we proceed to explain.

Up to here we have considered the case p=1/2 in (7.3.1), in order to bring out the essential features in the simplest case. However, this simplification would obscure the role of np and npq in the general formula below. The reader is advised to carry out the following calculations in the easier case p=q=1/2 to obtain some practice and confidence in such calculations.

Theorem 5. Suppose 0 ; put <math>q = 1 - p, and

$$x_{nk} = \frac{k - np}{\sqrt{npq}}, \quad 0 \le k \le n. \tag{7.3.11}$$

Clearly x_{nk} depends on both n and k, but it will be written as x_k below.

Let A be an arbitrary but fixed positive constant. Then in the range of k such that

$$|x_k| \le A,\tag{7.3.12}$$

we have

$$\binom{n}{k} p^k q^{n-k} \sim \frac{1}{\sqrt{2\pi npq}} e^{-x_k^2/2}.$$
 (7.3.13)

The convergence is uniform with respect to k in the range specified above.

Proof: We have from (7.3.11)

$$k = np + \sqrt{npq} x_k, \quad n - k = nq - \sqrt{npq} x_k. \tag{7.3.14}$$

Hence in the range indicated in (7.3.12),

$$k \sim np, \quad n - k \sim nq.$$
 (7.3.15)

Using Stirling's formula (7.3.3), we may write the left member of (7.3.13) as

$$\frac{\left(\frac{n}{e}\right)^{n}\sqrt{2\pi n}\,p^{k}q^{n-k}}{\left(\frac{k}{e}\right)^{k}\sqrt{2\pi k}\left(\frac{n-k}{e}\right)^{n-k}\sqrt{2\pi(n-k)}}$$

$$=\sqrt{\frac{n}{2\pi k(n-k)}}\,\varphi(n,k) \sim \frac{1}{\sqrt{2\pi npq}}\varphi(n,k)$$
(7.3.16)

by (7.3.15), where

$$\varphi(n,k) = \left(\frac{np}{k}\right)^k \left(\frac{nq}{n-k}\right)^{n-k}.$$

Taking logarithms and using the Taylor series

$$\log(1+x) = x - \frac{x^2}{2} + \dots + (-1)^{n-1} \frac{x^n}{n} + \dots, \quad |x| < 1,$$

we have by (7.3.14)

$$\log\left(\frac{np}{k}\right)^{k} = k\log\left(1 - \frac{\sqrt{npq}\,x_{k}}{k}\right)$$

$$= k\left(-\frac{\sqrt{npq}\,x_{k}}{k} - \frac{npqx_{k}^{2}}{2k^{2}} - \cdots\right),$$

$$\log\left(\frac{nq}{n-k}\right)^{n-k} = (n-k)\log\left(1 + \frac{\sqrt{npq}\,x_{k}}{n-k}\right)$$

$$= (n-k)\left(\frac{\sqrt{npq}\,x_{k}}{n-k} - \frac{npqx_{k}^{2}}{2(n-k)^{2}} + \cdots\right),$$

$$(7.3.17)$$

provided that

$$\left| \frac{\sqrt{npq} x_k}{k} \right| < 1 \quad \text{and} \quad \left| \frac{\sqrt{npq} x_k}{n-k} \right| < 1.$$
 (7.3.17')

These conditions are satisfied for sufficiently large values of n, in view of (7.3.12) and (7.3.15). Adding the two series expansions above whereupon the first terms cancel out each other obligingly, ignoring the dots but using " \sim " instead of "=," we obtain

$$\log \varphi(n,k) \sim -\frac{npqx_k^2}{2k} - \frac{npqx_k^2}{2(n-k)} = -\frac{n^2pqx_k^2}{2k(n-k)}.$$

In Appendix 2 we will give a rigorous demonstration of this relation. Using (7.3.15) again, we see that

$$\log \varphi(n,k) \sim -\frac{n^2 p q x_k^2}{2n p n q} = -\frac{x_k^2}{2}.$$
 (7.3.18)

In view of (7.3.12) [why do we need this reminder?], this is equivalent to

$$\varphi(n,k) \sim e^{-x_k^2/2}$$
.

Going back to (7.3.16), we obtain (7.3.13).

Theorem 6 (De Moivre-Laplace Theorem). For any two constants a and b, $-\infty < a < b < +\infty$, we have

$$\lim_{n \to \infty} P\left(a < \frac{S_n - np}{\sqrt{npq}} \le b\right) = \frac{1}{\sqrt{2\pi}} \int_a^b e^{-x^2/2} \, dx. \tag{7.3.19}$$

Proof: Let k denote a possible value of S_n so that $S_n = k$ means $(S_n - np)/\sqrt{npq} = x_k$ by the transformation (7.3.11). Hence the probability on the left side of (7.3.19) is just

$$\sum_{a < x_k \le b} P(S_n = k) = \sum_{a < x_k \le b} \binom{n}{k} p^k q^{n-k}.$$

Substituting for each term its asymptotic value given in (7.3.13), and observing from (7.3.11) that

$$x_{k+1} - x_k = \frac{1}{\sqrt{npq}},$$

we obtain

$$\frac{1}{\sqrt{2\pi}} \sum_{a < x_k \le b} e^{-x_k^2/2} (x_{k+1} - x_k). \tag{7.3.20}$$

The correspondence between k and x_k is one-to-one and when k varies from 0 to n, x_k varies in the interval $[-\sqrt{np/q}, \sqrt{nq/p}]$, not continuously but by an increment $x_{k+1}-x_k=1/\sqrt{npq}$. For large enough n the interval contains the given (a,b] and the points x_k falling inside (a,b] form a partition of it into equal subintervals of length $1/\sqrt{npq}$. Suppose the smallest and greatest values of k satisfying the condition $a < x_k \le b$ are j and k, then we have

$$x_{j-1} \le a < x_j < x_{j+1} < \dots < x_{l-1} < x_l \le b < x_{l+1}$$

and the sum in (7.3.20) may be written as follows:

$$\sum_{k=i}^{l} \varphi(x_k)(x_{k+1} - x_k); \quad \text{where } \varphi(x) = \frac{1}{\sqrt{2\pi}} e^{-x^2/2}.$$
 (7.3.21)

This is a Riemann sum for the definite integral $\int_a^b \varphi(x) dx$, although in standard textbook treatments of Riemann integration the endpoints a and b are usually included as points of partition. But this makes no difference as $n \to \infty$ and the partition becomes finer, so the sum above converges to the integral as shown in (7.3.19).

The result in (7.3.19) is called the *De Moivre-Laplace theorem* [Abraham De Moivre (1667–1754), considered as successor to Newton, gave this

result in his *Doctrine of Chances* (1714). Apparently he had priority over Stirling (1692–1770) for the formula named after the latter. Laplace extended it and realized its importance in his monumental *Théorie Analytique des Probabilités* (1812)]. It was the first known particular case of the central limit theorem to be discussed in the next section. It solves the problem of approximation stated at the beginning of the section. The right member of (7.3.20) involves a new probability distribution to be discussed in the next section. Simple examples of application will be given at the end of §7.5 and among the exercises.

7.4. Normal distribution

The probability distribution with the φ in (7.3.21) as density function will now be formally introduced:

$$\Phi(x) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{x} e^{-u^2/2} du, \quad \varphi(x) = \frac{1}{\sqrt{2\pi}} e^{-x^2/2}.$$

It is called the *normal distribution*, also the *Laplace-Gauss distribution*; and sometimes the prefix *unit* is attached to distinguish it from a whole family of normal distributions derived by a linear transformation of the variable x; see below. But we have yet to show that φ is a true probability density as defined in §4.5, namely that

$$\int_{-\infty}^{\infty} \varphi(x) \, dx = 1. \tag{7.4.1}$$

A heuristic proof of this fact may be obtained by setting $a=-\infty, b=+\infty$ in (7.3.19), whereupon the probability on the left side certainly becomes 1. Why is this not rigorous? Because two (or three) passages to limit are involved here that are not necessarily interchangeable. Actually the argument can be justified (see Appendix 2); but it may be more important that you should convince yourself that a justification is needed at all. This is an instance where advanced mathematics separates from the elementary kind we are doing mostly in this book.

A direct proof of (7.4.1) is also very instructive; although it is given in most calculus texts we will reproduce it for its sheer ingenuity. The trick is to consider the square of the integral in (7.4.1) and convert it to a double

integral:

$$\left(\int_{-\infty}^{\infty} \varphi(x) \, dx\right) \left(\int_{-\infty}^{\infty} \varphi(y) \, dy\right)$$

$$= \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} \varphi(x) \varphi(y) \, dx \, dy$$

$$= \frac{1}{2\pi} \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} \exp\left(-\frac{1}{2}(x^2 + y^2)\right) \, dx \, dy.$$

We can then use polar coordinates:

$$\rho^2 = x^2 + y^2$$
, $dx dy = \rho d\rho d\theta$

to evaluate it:

$$\begin{split} \frac{1}{2\pi} \int_0^{2\pi} \int_0^{\infty} \exp\left(-\frac{1}{2}\rho^2\right) \rho \, d\rho \, d\theta &= \frac{1}{2\pi} \int_0^{2\pi} -\exp\left(-\frac{1}{2}\rho^2\right) \Big|_0^{\infty} \, d\theta \\ &= \frac{1}{2\pi} \int_0^{2\pi} 1 \, d\theta = 1. \end{split}$$

This establishes (7.4.1) if we take the positive square root.

The normal density φ has many remarkable analytical properties; in fact, Gauss determined it by selecting a few of them as characteristics of a "law of errors." [Carl Friedrich Gauss (1777–1855) ranked as one of the greatest of all mathematicians, also did fundamental work in physics, astronomy, and geodesy. His major contribution to probability was through his theory of errors of observations, known as the method of least squares.] Let us observe first that it is a symmetric function of x, namely $\varphi(x) = \varphi(-x)$, from which the convenient formula follows:

$$\int_{-x}^{x} \varphi(u) \, du = \Phi(x) - \Phi(-x) = 2\Phi(x) - 1. \tag{7.4.2}$$

Next, φ has derivatives of all orders, and each derivative is the product of φ by a polynomial called a *Hermite polynomial*. The existence of all derivatives makes the curve $x \to \varphi(x)$ very smooth, and it is usually described as "bell-shaped."* Furthermore as $|x| \to \infty$, $\varphi(x)$ decreases to 0 very rapidly. The following estimate of the tail of Φ is often useful:

$$1 - \Phi(x) = \int_{x}^{\infty} \varphi(u) \, du \le \frac{\varphi(x)}{x} = \frac{e^{-x^{2}/2}}{\sqrt{2\pi} \, x}, \qquad x > 0.$$

^{*}See the graph attached to the table of $\Phi(x)$ on p. 394.

To see this, note that $-\varphi'(u) = u\varphi(u)$, hence

$$\int_x^\infty 1 \cdot \varphi(u) \, du \le \int_x^\infty \frac{u}{x} \varphi(u) \, du = \frac{-1}{x} \int_x^\infty \varphi'(u) \, du = \frac{-1}{x} \varphi(u) \Big|_x^\infty = \frac{\varphi(x)}{x},$$

another neat trick. It follows that not only Φ has moments of all orders, but even the integral

$$M(\theta) = \int_{-\infty}^{\infty} e^{\theta x} \varphi(x) \, dx = \int_{-\infty}^{\infty} \exp\left(\theta x - \frac{x^2}{2}\right) \, dx \tag{7.4.3}$$

is finite for every real θ , because $e^{-x^2/2}$ decreases much faster than $e^{|\theta x|}$ increases as $|x| \to \infty$. As a function of θ , M is called the *moment-generating function* of φ or Φ . Note that if we replace θ by the purely imaginary $i\theta$, then $M(i\theta)$ becomes the characteristic function or Fourier transform of Φ [see (6.5.17)]. The reason why we did not introduce the moment-generating function in §6.5 is because the integral in (7.4.3) rarely exists if φ is replaced by an arbitrary density function, but for the normal φ , $M(\theta)$ is cleaner than $M(i\theta)$ and serves as well. Let us now calculate $M(\theta)$. This is done by completing a square in the exponent in (7.4.3):

$$\theta x - \frac{x^2}{2} = \frac{\theta^2}{2} - \frac{(x-\theta)^2}{2}.$$

Now we have

$$M(\theta) = e^{\theta^2/2} \int_{-\infty}^{\infty} \varphi(x - \theta) \, dx = e^{\theta^2/2}.$$
 (7.4.4)

From this we can derive all the moments of Φ by successive differentiation of M with respect to θ , as in the case of a generating function discussed in §6.5. More directly, we may expand the $e^{\theta x}$ in (7.4.3) into its Taylor series in θ and compare the result with the Taylor series of $e^{\theta^2/2}$ in (7.4.4):

$$\int_{-\infty}^{\infty} \left\{ 1 + \theta x + \frac{(\theta x)^2}{2!} + \dots + \frac{(\theta x)^n}{n!} + \dots \right\} \varphi(x) dx$$
$$= 1 + \frac{\theta^2}{2} + \frac{1}{2!} \left(\frac{\theta^2}{2}\right)^2 + \dots + \frac{1}{n!} \left(\frac{\theta^2}{2}\right)^n + \dots$$

If we denote the *n*th moment by $m^{(n)}$:

$$m^{(n)} = \int_{-\infty}^{\infty} x^n \varphi(x) \, dx,$$

the above equation becomes

$$\sum_{n=0}^{\infty} \frac{m^{(n)}}{n!} \theta^n = \sum_{n=0}^{\infty} \frac{1}{2^n n!} \theta^{2n}.$$

It follows from the uniqueness of power series expansion (cf. §6.5) that the corresponding coefficients on both sides must be equal: thus for $n \ge 1$,

$$m^{(2n-1)} = 0,$$

$$m^{(2n)} = \frac{(2n)!}{2^n n!}.$$
(7.4.5)

Of course, the vanishing of all moments of odd order is an immediate consequence of the symmetry of φ .

In general, for any real m and $\sigma^2 > 0$, a random variable X is said to have a normal distribution $N(m, \sigma^2)$ iff the reduced variable

$$X^* = \frac{X - m}{\sigma}$$

has Φ as its distribution function. In particular, for m=0 and $\sigma^2=1$, N(0,1) is just the unit normal Φ . The density function of $N(m,\sigma^2)$ is

$$\frac{1}{\sqrt{2\pi}\,\sigma}\exp\left(-\frac{(x-m)^2}{2\sigma^2}\right) = \frac{1}{\sigma}\varphi\left(\frac{x-m}{\sigma}\right). \tag{7.4.6}$$

This follows from a general proposition (see Exercise 13 of Chapter 4). The moment-generating function M_X of X is most conveniently calculated through that of X^* as follows:

$$M_X(\theta) = E(e^{\theta(m+\sigma X^*)}) = e^{m\theta} E(e^{(\sigma\theta)X^*})$$

= $e^{m\theta} M(\sigma\theta) = e^{m\theta+\sigma^2\theta^2/2}$. (7.4.7)

A basic property of the *normal family* is given below. Cf. the analogous Theorem 3 in §7.2 for the Poisson family.

Theorem 7. Let X_j be independent random variables with normal distributions $N(m_j, \sigma_j^2), 1 \leq j \leq n$. Then $X_1 + \cdots + X_n$ has the normal distribution $N(\sum_{j=1}^n m_j, \sum_{j=1}^n \sigma_j^2)$.

Proof: It is sufficient to prove this for n=2, since the general case follows by induction. This is easily done by means of the moment-generating function. We have by the product theorem as in Theorem 6 of §6.5

$$\begin{split} M_{X_1+X_2}(\theta) &= M_{X_1}(\theta) M_{X_2}(\theta) = e^{m_1\theta + \sigma_1^2\theta^2/2} e^{m_2\theta + \sigma_2^2\theta^2/2} \\ &= e^{(m_1+m_2)\theta + (\sigma_1^2 + \sigma_2^2)\theta^2/2}. \end{split}$$

which is the moment-generating function of $N(m_1+m_2, \sigma_1^2+\sigma_2^2)$ by (7.4.7). Hence X_1+X_2 has this normal distribution since it is uniquely determined by the moment-generating function. [We did not prove this assertion, but see the end of §6.5.]

*7.5. Central limit theorem

We will now return to the De Moivre-Laplace Theorem 6 and give it a more general formulation. Recall that

$$S_n = X_1 + \dots + X_n, \quad n \ge 1,$$
 (7.5.1)

where the X_j 's are independent Bernoullian random variables. We know that for every j:

$$E(X_i) = p, \quad \sigma^2(X_i) = pq;$$

and for every n:

$$E(S_n) = np, \quad \sigma^2(S_n) = npq;$$

see Example 6 of §6.3. Put

$$X_j^* = \frac{X_j - E(X_j)}{\sigma(X_j)}; \quad S_n^* = \frac{S_n - E(S_n)}{\sigma(S_n)} = \frac{1}{\sqrt{n}} \sum_{j=1}^n X_j^*.$$
 (7.5.2)

The S_n^* 's are the random variables appearing in the left member of (7.3.19) and are sometimes called the *normalized* or *normed sums*. We have for every j and n:

$$E(X_j^*) = 0, \quad \sigma^2(X_j^*) = 1,$$

 $E(S_n^*) = 0, \quad \sigma^2(S_n^*) = 1.$ (7.5.3)

The linear transformation from X_j to X_j^* or S_n to S_n^* amounts to a change of origin and scale in the measurement of a random quantity in order to reduce its mean to zero and variance to 1 as shown in (7.5.3). Each S_n^* is a random variable taking the set of values

$$x_{n,k} = \frac{k - np}{\sqrt{npq}}, \quad k = 0, 1, \dots, n.$$

This is just the x_{nk} in (7.3.11). The probability distribution of S_n^* is given by

$$P(S_n^* = x_{n,k}) = \binom{n}{k} p^k q^{n-k}, \quad 0 \le k \le n.$$

It is more convenient to use the corresponding distribution function; call it F_n so that

$$P(S_n^* \le x) = F_n(x), \quad -\infty < x < \infty.$$

Finally, if I is the finite interval (a, b], and F is any distribution function, we shall write

$$F(I) = F(b) - F(a).$$

[By now you should understand why we used (a, b] rather than (a, b) or [a, b]. It makes no difference if F is continuous, but the F_n 's above are not continuous. Of course in the limit the difference disappears in the present case, but it cannot be ignored generally.] After these elaborate preparations, we can rewrite the De Moivre-Laplace formula in the elegant form below: for any finite interval I,

$$\lim_{n \to \infty} F_n(I) = \Phi(I). \tag{7.5.4}$$

Thus we see that we are dealing with the convergence of a sequence of distribution functions to a given distribution function in a certain sense.

In this formulation the subject is capable of a tremendous generalization. The sequence of distribution functions need not be those of normalized sums, the given limit need not be the normal distribution nor even specified in advance, and the sense of convergence need not be that specified above. For example, the Poisson limit theorem discussed in §7.1 can be viewed as a particular instance. The subject matter has been intensively studied in the last 40 years and is still undergoing further evolutions. [For some reference books in English, see [Feller 2], [Chung 1].] Here we must limit ourselves to one such generalization, the so-called central limit theorem in its classical setting, which is about the simplest kind of extension of Theorem 6 in §7.3. Even so we shall need a powerful tool from more advanced theory that we can use but not fully explain. This extension consists in replacing the Bernoullian variables above by rather arbitrary ones, as we proceed to describe.

Let $\{X_j, j \geq 1\}$ be a sequence of independent and identically distributed random variables. The phrase "identically distributed" means they have a common distribution, which need not be specified. But it is assumed that the mean and variance of each X_j are finite and denoted by m and σ^2 , respectively, where $0 < \sigma^2 < \infty$. Define S_n and S_n^* exactly as before, then

$$E(S_n) = nm, \quad \sigma^2(S_n) = n\sigma^2,$$
 (7.5.5)

and (7.5.3) holds as before. Again let F_n denote the distribution of the normalized sum S_n^* . Then Theorem 8 below asserts that (7.5.4) remains true under the liberalized conditions for the X_j 's. To mention just some simple cases, each X_j may now be a "die-rolling" instead of a "coin-tossing" random variable to which Theorem 6 is applicable; or it may be uniformly distributed ("point-picking" variable); or again it may be exponentially distributed ("telephone-ringing" variable). Think of some other varieties if you wish.

Theorem 8. For the sums S_n under the generalized conditions spelled out above, we have for any a < b

$$\lim_{n \to \infty} P\left(a < \frac{S_n - nm}{\sqrt{n}\,\sigma} \le b\right) = \frac{1}{\sqrt{2\pi}} \int_a^b e^{-x^2/2} \, dx. \tag{7.5.6}$$

Proof: The powerful tool alluded to earlier is that of the characteristic function discussed in §6.5. [We could not have used the moment-generating function since it may not exist for S_n .] For the unit normal distribution Φ , its characteristic function g can be obtained by substituting $i\theta$ for θ in (7.4.4):

$$g(\theta) = e^{-\theta^2/2}. (7.5.7)$$

With each arbitrary distribution F_n there is also associated its characteristic function g_n that is in general expressible by means of F_n as a *Stieltjes integral*. This is beyond the scope of this book but luckily we can bypass it in the following treatment by using the associated random variables. [Evidently we will leave the reader to find out what may be concealed!] We can now state the following result.

Theorem 9. If we have for every θ

$$\lim_{n \to \infty} g_n(\theta) = g(\theta) = e^{-\theta^2/2},$$
(7.5.8)

then we have for every x:

$$\lim_{n \to \infty} F_n(x) = \Phi(x) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^x e^{-u^2/2} du; \tag{7.5.9}$$

in particular, (7.5.4) is true.

Although we shall not prove this (see [Chung 1, Chapter 6]), let us at least probe its significance. According to Theorem 7 of §6.5, each g_n uniquely determines F_n , and g determines Φ . The present theorem carries this correspondence between distribution function and its transform (characteristic function) one step further; for it says that the *limit* of the sequence $\{g_n\}$ also determines the *limit* of the sequence $\{F_n\}$. Hence it has been called the "continuity theorem" for the transform. In the case of the normal Φ above the result is due to Pólya; the general case is due to Paul Lévy (1886–1972) and Harald Cramér (1893–1985); both pioneers of modern probability theory.

Next we need a little lemma about characteristic functions.

Lemma. If X has mean 0 and variance 1, then its characteristic function h has the following Taylor expansion at $\theta = 0$:

$$h(\theta) = 1 - \frac{\theta^2}{2}(1 + \epsilon(\theta)),$$
 (7.5.10)

where ϵ is a function depending on h such that $\lim_{\theta \to 0} \epsilon(\theta) = 0$.

Proof: According to a useful form of Taylor's theorem (look it up in your calculus book): if h has a second derivative at $\theta = 0$, then we have

$$h(\theta) = h(0) + h'(0)\theta + \frac{h''(0)}{2}\theta^2(1 + \epsilon(\theta)). \tag{7.5.11}$$

From

$$h(\theta) = E(e^{i\theta X})$$

we obtain by formal differentiation:

$$h'(\theta) = E(e^{i\theta X}iX), \quad h''(\theta) = E(e^{i\theta X}(iX)^2);$$

hence

$$h'(0) = E(iX) = 0, \quad h''(0) = E(-X^2) = -1.$$

Substituting into (7.5.11), we get (7.5.10).

Theorem 8 can now be proved by a routine calculation. Consider the characteristic function of S_n^* :

$$E(e^{i\theta S_n^*}) = E(e^{i\theta(X_1^* + \dots + X_n^*)/\sqrt{n}}).$$

Since the X_j^* 's are independent and identically distributed as well as the X_j 's, by the analogue of Theorem 6 of §6.5, the right member above is equal to

$$E(e^{i\theta X_1^*/\sqrt{n}})^n = h\left(\frac{\theta}{\sqrt{n}}\right)^n, \tag{7.5.12}$$

where h denotes the characteristic function of X_1^* . It follows from the lemma that

$$h\left(\frac{\theta}{\sqrt{n}}\right) = 1 - \frac{\theta^2}{2n}\left(1 + \epsilon\left(\frac{\theta}{\sqrt{n}}\right)\right) \tag{7.5.13}$$

where θ is fixed and $n \to \infty$. Consequently we have

$$\begin{split} \lim_{n \to \infty} E(e^{i\theta S_n^*}) &= \lim_{n \to \infty} \left[1 - \frac{\theta^2}{2n} \left(1 + \epsilon \left(\frac{\theta}{\sqrt{n}} \right) \right) \right]^n \\ &= e^{-\theta^2/2} \end{split}$$

by an application of (7.1.12). This means the characteristic functions of S_n^* converge to that of the unit normal; therefore by Theorem 9, the distribution F_n converges to Φ in the sense of (7.5.9), from which (7.5.6) follows.

The name "central limit theorem" is used generally to designate a convergence theorem in which the normal distribution appears as the limit. More particularly it applies to sums of random variables as in Theorem 8. Historically these variables arose as errors of observations of chance fluctuations, so that the result is the all-embracing assertion that under "normal" conditions they all obey the same normal law, also known as the "error function." For this reason it had been regarded by some as a law of nature! Even in this narrow context Theorem 8 can be generalized in several directions: the assumptions of a finite second moment, of a common distribution, and of strict independence can all be relaxed. Finally, if the normal conditions are radically altered, then the central limit theorem will no longer apply, and random phenomena abound in which the limit distribution is no longer normal. The Poisson case discussed in §7.1 may be considered as one such example, but there are other laws closely related to the normal that are called "stable" and "infinitely divisible" laws. See [Chung 1, Chapter 7 for a discussion of the various possibilities mentioned here.

It should be stressed that the central limit theorem as stated in Theorems 6 and 8 is of the form (7.5.4), without giving an estimate of the "error" $F_n(I) - \Phi(I)$. In other words, it asserts convergence without indicating any "speed of convergence." This renders the result useless in accurate numerical computations. However, under specified conditions it is possible to obtain bounds for the error. For example in the De Moivre–Laplace case (7.3.19) we can show that the error does not exceed C/\sqrt{n} , where C is a numerical constant involving p but not a or b; see [Chung 1, §7.4] for a more general result. In crude, quick-and-dirty applications the error is simply ignored, as will be done below.

In contrast to the mathematical developments, simple practical applications that form the backbone of "large sample theory" in statistics are usually of the cookbook variety. The great limit theorem embodied in (7.5.6) is turned into a rough approximate formula that may be written as follows:

$$P(x_1\sigma\sqrt{n} < S_n - mn < x_2\sigma\sqrt{n}) \approx \Phi(x_2) - \Phi(x_1).$$

In many situations we are interested in a symmetric spread around the

mean, i.e., $x_1 = -x_2$. Then the above becomes by (7.4.2)

$$P(|S_n - mn| < x\sigma\sqrt{n}) \approx 2\Phi(x) - 1. \tag{7.5.14}$$

Extensive tabulations of the values of Φ and its inverse function Φ^{-1} are available; a short table is appended at the end of the book. The following example illustrates the routine applications of the central limit theorem.

Example 7. A physical quantity is measured many times for accuracy. Each measurement is subject to a random error. It is judged reasonable to assume that it is uniformly distributed between -1 and +1 in a conveniently chosen unit. Now if we take the arithmetical mean [average] of n measurements, what is the probability that it differs from the true value by less than a fraction δ of the unit?

Let the true value be denoted by m and the actual measurements obtained by $X_i, 1 \le j \le n$. Then the hypothesis says that

$$X_j = m + \xi_j,$$

where ξ_j is a random variable that has the uniform distribution in [-1, +1]. Thus

$$E(\xi_j) = \int_{-1}^{+1} \frac{x}{2} dx = 0, \quad \sigma^2(\xi_j) = E(\xi_j^2) = \int_{-1}^{+1} \frac{1}{2} x^2 dx = \frac{1}{3},$$

$$E(X_j) = m, \qquad \sigma^2(X_j) = \frac{1}{3}.$$

In our notation above, we want to compute the approximate value of $P\{|S_n - mn| < \delta n\}$. This probability must be put into the form given in (7.5.6), and the limit relation there becomes by (7.5.14)

$$P\left\{\left|\frac{S_n-mn}{\sqrt{n/3}}\right|<\delta\sqrt{3n}\right\}pprox 2\Phi(\delta\sqrt{3n})-1.$$

For instance, if n = 25 and $\delta = 1/5$, then the result is equal to

$$2\Phi(\sqrt{3}) - 1 \approx 2\Phi(1.73) - 1 \approx .92$$

from the Table on p. 394. Thus, if 25 measurements are taken, then we are 92% sure that their average is within one fifth of a unit from the true value.

Often the question is turned around: how many measurements should we take in order for the probability to exceed α (the "significance level") and for the average to differ from the true value by at most δ ? This means we must find the value x_{α} such that

$$2\Phi(x_{\alpha}) - 1 = \alpha$$
 or $\Phi(x_{\alpha}) = \frac{1+\alpha}{2}$,

and then choose n to make

$$\delta\sqrt{3n} > x_{\alpha}.$$

For instance, if $\alpha = .95$ and $\delta = 1/5$, then the table shows that $x_{\alpha} \approx 1.96$; hence

$$n > \frac{x_{\alpha}^2}{3\delta^2} \approx 32.$$

Thus, seven or eight more measurements should increase our degree of confidence from 92% to 95%. Whether this is worthwhile may depend on the cost of doing the additional work as well as the significance of the enhanced probability.

It is clear that there are three variables involved in questions of this kind, namely: δ , α , and n. If two of them are fixed, we can solve for the third. Thus if n=25 is fixed because the measurements are found in recorded data and not repeatable, and our credulity demands a high degree of confidence α , say 99%, then we must compromise on the coefficient of accuracy δ . We leave this as an exercise.

Admittedly these practical applications of the great theorem are dull stuff, but so are, e.g., Newton's laws of motion on the quotidian level.

7.6. Law of large numbers

In this section we collect two results related to the central limit theorem: the law of large numbers and Chebyshev's inequality.

The celebrated law of large numbers can be deduced from Theorem 8 as an easy consequence.

Theorem 10. Under the same conditions as in Theorem 8, we have for a fixed but arbitrary constant c > 0,

$$\lim_{n \to \infty} P\left(\left| \frac{S_n}{n} - m \right| < c \right) = 1. \tag{7.6.1}$$

Proof: Since c is fixed, for any positive constant l, we have

$$l\sigma\sqrt{n} < cn \tag{7.6.2}$$

for all sufficiently large values of n. Hence the event

$$\left\{ \left| \frac{S_n - mn}{\sigma \sqrt{n}} \right| < l \right\} \quad \text{certainly implies} \quad \left\{ \left| \frac{S_n - mn}{n} \right| < c \right\},$$

and so

$$P\left(\left|\frac{S_n - mn}{\sigma\sqrt{n}}\right| < l\right) \le P\left(\left|\frac{S_n - mn}{n}\right| < c\right) \tag{7.6.3}$$

for large n. According to (7.5.6) with a = -l, b = +l, the left member above converges to

$$\frac{1}{\sqrt{2\pi}} \int_{-l}^{l} e^{-x^2/2} dx$$

as $n \to \infty$. Given any $\delta > 0$ we can first choose l so large that the value of the integral above exceeds $1 - \delta$, then choose n so large that (7.6.3) holds. It follows that

$$P\left(\left|\frac{S_n}{n} - m\right| < c\right) > 1 - \delta \tag{7.6.4}$$

for all sufficiently large n, and this is what (7.6.1) says.

Briefly stated, the law of large numbers is a corollary to the central limit theorem because any large multiple of \sqrt{n} is negligible in comparison with any small multiple of n.

In the Bernoullian case the result was first proved by Jakob Bernoulli as a crowning achievement. [Jakob or Jacques Bernoulli (1654–1705), Swiss mathematician and physicist, author of the first treatise on probability: Ars conjectandi (1713), which contains this theorem. His proof depends on direct calculations with binomial coefficients without, of course, the benefit of such formulas as Stirling's. In a sense the De Moivre-Laplace Theorem 6 was a sequel to it. By presenting it in reverse to the historical development it is made to look like a trivial corollary. As a matter of fact, the law of large numbers is a more fundamental but also more primitive limit theorem. It holds true under much broader conditions than the central limit theorem. For instance, in the setting of Theorem 8, it is sufficient to assume that the common mean of X_i is finite, without any assumption on the second moment. Since the assertion of the law concerns only the mean, such an extension is significant and was first proved by A.Ya. Khintchine [1894– 1959, one of the most important of the school of Russian probabilists. In fact, it can be proved by the method used in the proof of Theorem 8 above, except that it requires an essential extension of Theorem 9 which will take us out of our depth here. (See Theorem 6.4.3 of [Chung 1].) Instead we will give an extension of Theorem 10 in another direction, when the random variables $\{X_i\}$ are not necessarily identically distributed. This is easy via another celebrated but simple result known as Chebyshev's inequality. [P.L. Chebyshev (1821–94) together with A.A. Markov (1856–1922) and A. M. Ljapunov (1857–1918) were founders of the Russian school of probability.

Theorem 11. Suppose the random variable X has a finite second moment. Then for any constant c > 0 we have

$$P(|X| \ge c) \le \frac{E(X^2)}{c^2}. (7.6.5)$$

Proof: We will carry out the proof for a countably valued X and leave the analogous proof for the density case as an exercise. The idea of the proof is the same for a general random variable.

Suppose that X takes the values v_i with probabilities p_i , as in §4.3. Then we have

$$E(X^2) = \sum_{j} p_j v_j^2. (7.6.6)$$

If we consider only those values v_j satisfying the inequality $|v_j| \geq c$ and denote by A the corresponding set of indices j, namely $A = \{j \mid |v_j| \geq c\}$, then of course $v_j^2 \geq c^2$ for $j \in A$, whereas

$$P(|X| \ge c) = \sum_{j \in A} p_j.$$

Hence if we sum the index j only over the partial set A, we have

$$E(X^2) \ge \sum_{j \in A} p_j v_j^2 \ge \sum_{j \in A} p_j c^2 = c^2 \sum_{j \in A} p_j = c^2 P(|X| \ge c),$$

which is (7.6.5).

We can now state an extended form of the law of large numbers as follows.

Theorem 12. Let $\{X_j, j \geq 1\}$ be a sequence of independent random variables such that for each j,

$$E(X_j) = m_j, \quad \sigma^2(X_j) = \sigma_j^2; \tag{7.6.7}$$

and furthermore suppose there exists a constant $M < \infty$ such that for all j,

$$\sigma_j^2 \le M. \tag{7.6.8}$$

Then we have for each fixed c > 0,

$$\lim_{n \to \infty} P\left(\left| \frac{X_1 + \dots + X_n}{n} - \frac{m_1 + \dots + m_n}{n} \right| < c \right) = 1. \tag{7.6.9}$$

Proof: If we write $X_j^0 = X_j - m_j$, $S_n^0 = \sum_{j=1}^n X_j^0$, then the expression between the bars above is just S_n^0/n . Of course, $E(S_n^0) = 0$, whereas

$$E((S_n^0)^2) = \sigma^2(S_n^0) = \sum_{j=1}^n \sigma^2(X_j^0) = \sum_{j=1}^n \sigma_j^2.$$

This string of equalities follows easily from the properties of variances and you ought to have no trouble recognizing them now at a glance. [If you still do, then you should look up the places in preceding chapters where they are discussed.] Now the condition in (7.6.8) implies that

$$E((S_n^0))^2) \le Mn, \quad E\left(\left(\frac{S_n^0}{n}\right)^2\right) \le \frac{M}{n}. \tag{7.6.10}$$

It remains to apply Theorem 11 to $X = S_n^0/n$ to obtain

$$P\left(\left|\frac{S_n^0}{n}\right| \ge c\right) \le \frac{E((S_n^0/n)^2)}{c^2} \le \frac{M}{c^2n}.$$
 (7.6.11)

Hence the probability above converges to zero as $n \to \infty$, which is equivalent to the assertion in (7.6.9).

Actually the proof yields more: it gives an estimate on the "speed of convergence." Namely, given M, c, and δ we can tell how large n must be in order for the probability in (7.6.9) to exceed $1 - \delta$. Note also that Theorem 10 is a particular case of Theorem 12 because there all the σ_j^2 's are equal and we may take $M = \sigma_1^2$.

Perhaps the reader will agree that the above derivations of Theorems 11 and 12 are relatively simple doings compared with the fireworks in §§7.3–7.4. Looking back, we may find it surprising that it took two centuries before the right proof of Bernoulli's theorem was discovered by Chebyshev. It is an instance of the triumph of an idea, a new way of thinking, but even Chebyshev himself buried his inequality among laborious and unnecessary details. The cleaning up as shown above was done by later authors. Let us observe that the method of proof is applicable to any sum S_n of random variables, whether they are independent or not, provided that the crucial estimates in (7.6.10) are valid.

We turn now to the meaning of the law of large numbers. This is best explained in the simplest Bernoullian scheme where each X_j takes the values 1 and 0 with probabilities p and q = 1 - p, as in Theorems 5 and 6 above. In this case $S_n^0 = S_n - np$ and $E((S_n^0)^2) = \sigma^2(S_n) = npq$, so that (7.6.11) becomes

$$P\left(\left|\frac{S_n}{n} - p\right| \ge c\right) \le \frac{pq}{c^2 n} \le \frac{1}{4c^2 n};\tag{7.6.12}$$

note that $p(1-p) \leq 1/4$ for $0 \leq p \leq 1$. In terms of coin tossing with p as the probability of a head, S_n/n represents the relative frequency of heads in n independent tosses; cf. Example 3 of §2.1. This is, of course, a random number varying from one experiment to another. It will clarify things if we reinstate the long-absent sample point ω and write explicitly

$$\frac{S_n(\omega)}{n} = \frac{\text{relative frequency of heads in } n \text{ tosses associated}}{\text{with the experiment denoted by } \omega.}$$

Thus each ω is conceived as the record of the outcomes of a sequence of coin tossing called briefly an *experiment*, so that different ω 's correspond to different experiments. The sequence of tossing should be considered as *infinite* [indefinite, unending] in order to allow arbitrarily large values of n in computing the frequency [but we are not necessarily assuming the existence of a limiting frequency; see below]. Symbolically, ω may be regarded as a concise representation of the sequence of successive outcomes:

$$\omega = \{X_1(\omega), X_2(\omega), \dots, X_n(\omega), \dots\};$$

see the discussion in §4.1. We can talk about the probability of certain sets of ω ; indeed we have written the probability of the set

$$\Lambda_n(c) = \left\{ \omega \left| \left| \frac{S_n(\omega)}{n} - p \right| \ge c \right\} \right.$$

in (7.6.12); see also, e.g., (7.3.10) and (7.3.19), where similar probabilities are evaluated. In the precise form given in (7.6.12), if the value c as well as any given positive number ϵ is assigned, we can determine how large n needs to be to make $P(\Lambda_n(c)) \leq \epsilon$, namely:

$$\frac{1}{4c^2n} \le \epsilon \quad \text{or} \quad n \ge \frac{1}{4c^2\epsilon}. \tag{7.6.13}$$

This bound for n is larger than necessary because Chebyshev's inequality is a crude estimate in general. A sharper bound can be culled from Theorem 6 as follows. Rewrite the probability in (7.6.12) and use the approximation given in (7.3.19) simplified by (7.4.2):

$$P\left(\left|\frac{S_n - np}{\sqrt{npq}}\right| \ge \sqrt{\frac{n}{pq}}c\right) \approx 2\left[1 - \Phi\left(\sqrt{\frac{n}{pq}}c\right)\right].$$
 (7.6.14)

This is *not* an asymptotic relation as it pretends to be, because the constants a and b must be fixed in Theorem 6, whereas here $\pm c\sqrt{n}/\sqrt{pq}$ vary with n (and increase too rapidly with n). It can be shown that the difference of the two sides above is of the form A/\sqrt{n} , where A is a numerical

constant depending on p, and this error term should not be ignored. But we do so below. Now put

$$\eta = \sqrt{\frac{n}{pq}} \, c;$$

our problem is to find the value of η to make

$$2[1 - \Phi(\eta)] \le \epsilon \quad \text{or} \quad \Phi(\eta) \ge 1 - \frac{\epsilon}{2};$$
 (7.6.15)

then solve for n from η . This can be done by looking up a table of values of Φ ; a short one is appended at the end of this book.

Example 8. Suppose c = 2% and $\epsilon = 5\%$. Then (7.6.15) becomes

$$\Phi(\eta) \ge 1 - \frac{5}{200} = .975.$$

From the table we see that this is satisfied if $\eta \geq 1.96$. Thus

$$n \ge \frac{(1.96)^2 pq}{c^2} = \frac{(1.96)^2 \times 10000}{4} pq.$$

The last term depends on p, but $p(1-p) \leq 1/4$ for all p, as already noted, and so $n \geq 10000 \cdot (1/4) = 2500$ will do. For comparison, the bound given in (7.6.13) requires $n \geq 12500$; but that estimate has been rigorously established whereas the normal approximation is a rough-and-dirty one. We conclude that if the coin is tossed more than 2500 times, then we can be 95% sure that relative frequency of heads computed from the actual experiment will differ from the true p by no more than 2%.

Such a result can be applied in two ways (both envisioned by Bernoulli): (i) if we consider p as known, then we can make a prediction on the outcome of the experiment; (ii) if we regard p as unknown, then we can make an estimate of its value by performing an actual experiment. The second application has been called a problem of "inverse probability" and is the origin of the so-called Monte Carlo method. Here is a numerical example. In an actual experiment 10000 tosses were made and the total number of heads obtained is 4979; see [Feller 1, p. 21] for details. The computation above shows that we can be 95% sure that

$$\left| p - \frac{4979}{10000} \right| \le \frac{2}{100}$$
 or $.4779 \le p \le .5179$.

Returning to the general situation in Theorem 10, we will state the law of large numbers in the following form reminiscent of the definition of

an ordinary limit. For any $\epsilon > 0$, there exists an $n_0(\epsilon)$ such that for all $n \geq n_0(\epsilon)$ we have

$$P\left(\left|\frac{S_n}{n} - m\right| < \epsilon\right) > 1 - \epsilon. \tag{7.6.16}$$

We have taken both c and δ in (7.6.4) to be ϵ without loss of generality [see Exercise 22]. If we interpret this as in the preceding example as an assertion concerning the proximity of the theoretical mean m to the empirical average S_n/n , the double hedge [margin of error] implied by the two ϵ 's in (7.6.16) seems inevitable. For in any experiment one can neither be 100% sure nor 100% accurate, otherwise the phenomenon would not be random. Nevertheless, mathematicians are idealists and long for perfection. What cannot be realized in the empirical world may be achieved in a purely mathematical scheme. Such a possibility was uncovered by Borel in 1909, who created a new chapter in probability by his discovery described below. In the Bernoullian case, his famous result may be stated as follows:

$$P\left(\lim_{n\to\infty}\frac{S_n}{n}=p\right)=1.^*\tag{7.6.17}$$

This is known as a "strong law of large numbers," which is an essential improvement on Bernoulli's "weak law of large numbers." It asserts the existence of a limiting frequency equal to the theoretical probability p, for all sample points ω except possibly a set of probability zero (but not necessarily an empty set). Thus the limit in (2.1.10) indeed exists, but only for almost all ω , so that the empirical theory of frequencies beloved by the applied scientist is justifiable through a sophisticated theorem. The difference between this and Bernoulli's weaker theorem:

$$\forall \epsilon > 0$$
: $\lim_{n \to \infty} P\left(\left|\frac{S_n}{n} - p\right| < \epsilon\right) = 1$,

is subtle and cannot be adequately explained without measure theory. The astute reader may observe that although we claim 100% certainty and accuracy in (7.6.17), the limiting frequency is not an empirically observable thing—so that the cynic might say that what we are sure of is only an *ideal*, whereas the sophist could retort that we shall never be caught wanting! Even so a probabilistic certainty does not mean absolute certainty in the deterministic sense. There is an analogue of this distinction in the second law of thermodynamics (which comes from statistical mechanics). According to that law, e.g., when a hot body is in contact with a cold body, it is *logically possible* that heat will flow from the cold to the hot,

^{*}For a discussion of Borel's theorem and related topics, see Chapter 5 of [Chung 1].

but the probability of this happening is zero. A similar exception is permitted in Borel's theorem. For instance, if a coin is tossed indefinitely, it is logically possible that it's heads every single time. Such an event constitutes an exception to the assertion in (7.6.17), but its probability is equal to $\lim_{n\to\infty} p^n = 0$.

The strong law of large numbers is the foundation of a mathematical theory of probability based on the concept of frequency; see §2.1. It makes better sense than the weak one and is indispensable for certain theoretical investigations. [In statistical mechanics it is known in an extended form under the name *ergodic theorem*.] But the dyed-in-the-wool empiricist, as well as a radical school of logicians called intuitionists, may regard it as an idealistic fiction. It is amusing to quote two eminent authors on the subject:

Feller: "[the weak law of large numbers] is of very limited interest and should be replaced by the more precise and more useful strong law of large numbers." (p. 152 of [Feller 1])

van der Waerden: "[the strong law of large numbers] scarcely plays a role in mathematical statistics." (p. 98 of *Mathematische Statistik*, 3rd ed., Springer-Verlag, 1971)

Let us end this discussion by keeping in mind the gap between observable phenomena in the real world and the theoretical models used to study them; see Einstein's remark on p. 129. The law of large numbers, weak or strong, is a mathematical theorem deduced from axioms. Its applicability to true-life experiences such as the tossing of a penny or nickel is necessarily limited and imperfect. The various examples given above to interpret and illustrate the theorems should be viewed with this basic understanding.

Exercises

- 1. Suppose that a book of 300 pages contains 200 misprints. Use Poisson approximation to write down the probability that there is more than one misprint on a particular page.
- 2. In a school where 4% of the children write with their left hands, what is the probability that there are no left-handed children in a class of 25?
- 3. Six dice are thrown 200 times by the players. Estimate the probability of obtaining "six different faces" k times, where k = 0, 1, 2, 3, 4, 5.
- 4. A home bakery made 100 loaves of raisin bread using 2000 raisins. Write down the probability that the loaf you bought contains 20 to 30 raisins.
- 5. It is estimated that on a certain island of 15 square miles there are 20 giant tortoises of one species and 30 of another species left. An

- ecological survey team spotted 2 of them in an area of 1 square mile, but neglected to record which species. Use Poisson distribution to find the probabilities of the various possibilities.
- 6. Find the maximum term or terms in the binomial distribution $B_k(n; p)$, $0 \le k \le n$. Show that the terms increase up to the maximum and then decrease. [Hint: take ratios of consecutive terms.]
- 7. Find the maximum term or terms in the Poisson distribution $\pi_k(\alpha)$, $0 < k < \infty$. Show the same behavior of the terms as in No. 6.
- 8. Let X be a random variable such that $P(X = c + kh) = \pi_k(\alpha)$, where c is a real and h is a positive number. Find the Laplace transform of X.
- 9. Find the convolution of two sequences given by Poisson distributions $\{\pi_k(\alpha)\}\$ and $\{\pi_k(\beta)\}\$.
- *10. If X_{α} has the Poisson distribution $\pi(\alpha)$, then

$$\lim_{\alpha \to \infty} P\left\{ \frac{X_{\alpha} - \alpha}{\sqrt{\alpha}} \le u \right\} = \Phi(u)$$

for every u. [Hint: use the Laplace transform $E(e^{-\lambda(X_{\alpha}-\alpha)/\sqrt{\alpha}})$, show that as $\alpha \to \infty$ it converges to $e^{\lambda^2/2}$, and invoke the analogue of Theorem 9 of §7.5.]

- 11. Assume that the distance between cars going in one direction on a certain highway is exponentially distributed with mean value of 100 meters. What is the probability that in a stretch of 5 kilometers there are between 50 to 60 cars?
- 12. On a certain highway the flow of traffic may be assumed to be Poissonian with intensity equal to 30 cars per minute. Write down the probability that it takes more than N seconds for n consecutive cars to pass by an observation post. [Hint: use (7.2.11).]
- 13. A perfect die is rolled 100 times. Find the probability that the sum of all points obtained is between 330 and 380.
- 14. It is desired to find the probability p that a certain thumbtack will fall on its flat head when tossed. How many trials are needed in order that we may be 95% sure that the observed relative frequency differs from p by less than p/10? [Hint: try it a number of times to get a rough bound for p.]
- 15. Two movie theaters compete for 1000 customers. Suppose that each customer chooses one of the two with "total indifference" and independently of other customers. How many seats should each theater have so that the probability of turning away any customer for lack of seats is less than 1%?
- 16. A sufficient number of voters are polled to determine the percentage in favor of a certain candidate. Assuming that an unknown proportion *p* of the voters favor him and they act independently of one another,

how many should be polled to predict the value of p within 4.5% with 95% confidence? [This is the so-called four percent margin of error in predicting elections, presumably because <.045 becomes \leq .04 by the rule of rounding decimals.]

- 17. Write $\Phi((a,b))$ for $\Phi(b) \Phi(a)$, where a < b and Φ is the unit normal distribution. Show that $\Phi((0,2)) > \Phi((1,3))$ and generalize to any two intervals of the same length. [Hint: $e^{-x^2/2}$ decreases as |x| increases.]
- 18. Complete the proof of (7.1.8) and then use the same method to prove (7.1.12). [Hint: $|\log(1-x)+x| \leq (1/2) \sum_{n=2}^{\infty} |x|^n = x^2/(2(1-|x|))$; hence if $|x| \leq 1/2$ this is bounded by x^2 .]
- 19. Prove (7.1.13).
- 20. Prove Chebyshev's inequality when X has a density. [Hint: $\sigma^2(X) = \int_{-\infty}^{\infty} (x-m)^2 f(x) dx \ge \int_{|x-m|>c} (x-m)^2 f(x) dx$.]
- 21. Prove the following analogue of Chebyshev's inequality where the absolute first moment is used in place of the second moment:

$$P(|X - m| > c) \le \frac{1}{c}E(|X - m|).$$

*22. Show that $\lim_{n\to\infty} P(|X_n| > \epsilon) = 0$ for every ϵ if and only if given any ϵ , there exists $n_0(\epsilon)$ such that

$$P(|X_n| > \epsilon) < \epsilon \quad \text{for} \quad n > n_0(\epsilon).$$

This is also equivalent to: given any δ and ϵ , there exists $n_0(\delta, \epsilon)$ such that

$$P(|X_n| > \epsilon) < \delta$$
 for $n > n_0(\delta, \epsilon)$.

[Hint: consider $\epsilon' = \delta \wedge \epsilon$ and apply the first form.]

- 23. If X has the distribution Φ , show that |X| has the distribution Ψ , where $\Psi = 2\Phi 1$; Ψ is called the "positive normal distribution."
- 24. If X has the distribution Φ , find the density function of X^2 and the corresponding distribution. This is known as the "chi-square distribution" in statistics. [Hint: differentiate $P(X^2 < x) = 2/\sqrt{2\pi} \int_0^{\sqrt{x}} e^{-u^2/2} du$.]
- *25. Use No. 24 to show that

$$\int_0^\infty x^{-1/2} e^{-x} \, dx = \sqrt{\pi}.$$

The integral is equal to $\Gamma(1/2)$, where Γ is the gamma function defined by $\Gamma(\alpha) = \int_0^\infty x^{\alpha-1} e^{-x} dx$ for $\alpha > 0$. [Hint: consider $E(X^2)$ in No. 24.]

*26. Let $\{\xi_k, 1 \leq k \leq n\}$ be n random variables satisfying $0 < \xi_1 \leq \xi_2 \leq \cdots \leq \xi_n \leq t$; let $(0,t] = \bigcup_{k=1}^l I_k$ be an arbitrary partition of (0,t] into subintervals $I_k = (x_{k-1}, x_k]$, where $x_0 = 0$; and let $\tilde{N}(I_k)$ denote the

number of ξ 's belonging to I_k . How can we express the event $\{\xi_k \leq x_k; 1 \leq k \leq l\}$ by means of $\tilde{N}(I_k), 1 \leq k \leq l$? Here, of course, $0 < x_1 < x_2 < \cdots < x_n \leq t$. Now suppose that $x_k, 1 \leq k \leq l$, are arbitrary and answer the question again. [Hint: try n=2 and 3 to see what is going on; relabel the x_k in the second part.]

- *27. Let $\{X(t), t \geq 0\}$ be a Poisson process with parameter α . For a fixed t > 0 define $\delta(t)$ to be the distance from t to the last jump before t if there is one, and to be t otherwise. Define $\delta'(t)$ to be the distance from t to the next jump after t. Find the distributions of $\delta(t)$ and $\delta'(t)$. [Hint: if $u < t, P\{\delta(t) > u\} = P\{N(t u, t) = 0\}$; for all u > 0, $P\{\delta'(t) > u\} = P\{N(t, t + u) = 0\}$.]
- *28. Let $\tau(t) = \delta(t) + \delta'(t)$ as in No. 27. This is the length of the between-jump interval containing the given time t. For each ω , this is one of the random variables T_k described in §7.2. Does $\tau(t)$ have the same exponential distribution as all the T_k 's? [This is a nice example where logic must take precedence over "intuition," and it is often referred to as a paradox. The answer should be easy from No. 27. For further discussion at a level slightly more advanced than this book, see Chung, "The Poisson process as renewal process," Periodica Mathematica Hungarica, Vol. 2 (1972), pp. 41–48.]
- 29. Use Chebyshev's inequality to show that if X and Y are two arbitrary random variables satisfying $E\{(X-Y)^2\}=0$, then we have P(X=Y)=1, namely X and Y are almost surely identical. [Hint: $P(|X-Y|>\epsilon)=0$ for any $\epsilon>0$.]
- 30. Recall the coefficient of correlation $\rho(X,Y)$ from §6.3. Show that if $\rho(X,Y)=1$, then the two "normalized" random variables

$$\tilde{X} = \frac{X - E(X)}{\sigma(X)}, \quad \tilde{Y} = \frac{Y - E(Y)}{\sigma(Y)}$$

are almost surely identical. What if $\rho(X,Y)=-1$? [Hint: compute $E\{(\tilde{X}-\tilde{Y})^2\}$ and use No. 29.]

Appendix 2

Stirling's Formula and De Moivre–Laplace's Theorem

In this appendix we complete some details in the proof of Theorem 5, establish Stirling's formula (7.3.3), and relate it to the normal integral (7.4.1). We begin with an estimate.

Lemma. If $|x| \leq 2/3$, then

$$\log(1+x) = x - \frac{x^2}{2} + \theta(x),$$

where $|\theta(x)| \leq |x|^3$.

Proof: We have by Taylor's series for $\log(1+x)$:

$$\log(1+x) = x - \frac{x^2}{2} + \sum_{n=3}^{\infty} (-1)^{n-1} \frac{x^n}{n}.$$

Hence $\theta(x)$ is equal to the series above and

$$\theta(x) \le \sum_{n=3}^{\infty} \frac{|x|^n}{n} \le \frac{1}{3} \sum_{n=3}^{\infty} |x|^n = \frac{|x|^3}{3(1-|x|)}.$$

For $|x| \leq 2/3, 3(1-|x|) \geq 1$ and the lemma follows. The choice of the constant 2/3 is a matter of convenience; a similar estimate holds for any constant < 1.

We will use the lemma first to complete the proof of Theorem 5, by showing that the omitted terms in the two series expansions in (7.3.17)

may indeed be ignored as $n \to \infty$. When n is sufficiently large the two quantities in (7.3.17') will be $\leq 2/3$. Consequently the lemma is applicable and the contribution from the "tails" of the two series, represented by dots there, is bounded by

$$k \left| \frac{\sqrt{npq} x_k}{k} \right|^3 + (n-k) \left| \frac{\sqrt{npq} x_k}{n-k} \right|^3.$$

Since pq < 1 and $|x_k| \leq A$, this does not exceed

$$\frac{n^{3/2}}{k^2}A^3 + \frac{n^{3/2}}{(n-k)^2}A^3,$$

which clearly tends to zero as $n \to \infty$, by (7.3.15). Therefore the tails vanish in the limit, and we are led to (7.3.18) as shown there.

Next we shall prove, as a major step toward Stirling's formula, the relation below:

$$\lim_{n \to \infty} \left\{ \log n! - \left(n + \frac{1}{2} \right) \log n + n \right\} = C, \tag{A.2.1}$$

where C is a constant to be determined later. Let d_n denote the quantity between the braces in (A.2.1). Then a simple computation gives

$$d_n - d_{n+1} = \left(n + \frac{1}{2}\right) \log\left(1 + \frac{1}{n}\right) - 1.$$

Using the notation in the lemma, we write this as

$$\left(n+\frac{1}{2}\right)\left(\frac{1}{n}-\frac{1}{2n^2}+\theta\left(\frac{1}{n}\right)\right)-1=\left(n+\frac{1}{2}\right)\theta\left(\frac{1}{n}\right)-\frac{1}{4n^2},$$

and consequently by the lemma with $x = 1/n, n \ge 2$:

$$|d_n - d_{n+1}| \le \left(n + \frac{1}{2}\right) \frac{1}{n^3} + \frac{1}{4n^2} = \frac{2n+1}{2n^3} + \frac{1}{4n^2}.$$

Therefore the series $\sum_{n} |d_n - d_{n+1}|$ converges by the comparison test. Now recall that an absolutely convergent series is convergent, which means the partial sum tends to a finite limit, say C_1 . Thus we have

$$\lim_{N \to \infty} \sum_{n=1}^{N} (d_n - d_{n+1}) = C_1;$$

but the sum above telescopes into $d_1 - d_{N+1}$, and so

$$\lim_{N \to \infty} d_{N+1} = d_1 - C_1,$$

and we have proved the assertion in (A.2.1) with $C = d_1 - C_1$. It follows that

$$\lim_{n \to \infty} \frac{n! e^n}{n^{n+(1/2)}} = e^C,$$

or if $K = e^C$:

$$n! \sim K n^{n+(1/2)} e^{-n}$$
. (A.2.2)

If we compare this with (7.3.3) we see that it remains to prove that $K=\sqrt{2\pi}$ to obtain Stirling's formula. But observe that even without this evaluation of the constant K, the calculations in Theorems 5 and 6 of §7.3 are valid provided we replace $\sqrt{2\pi}$ by K everywhere. In particular, formula (7.3.19) with a=-b becomes

$$\lim_{n \to \infty} P\left(\left| \frac{S_n - np}{\sqrt{npq}} \right| \le b \right) = \frac{1}{K} \int_{-b}^{b} e^{-x^2/2} dx. \tag{A.2.3}$$

On the other hand, we may apply Theorem 11 (Chebyshev's inequality) with $X = (S_n - np)/\sqrt{npq}$, E(X) = 0, $E(X^2) = 1$, to obtain the inequality

$$P\left(\left|\frac{S_n - np}{\sqrt{npq}}\right| \le b\right) \ge 1 - \frac{1}{b^2}.\tag{A.2.4}$$

Combining the last two relations and remembering that a probability cannot exceed 1, we obtain,

$$1 - \frac{1}{b^2} \le \frac{1}{K} \int_{-b}^b e^{-x^2/2} \, dx \le 1.$$

Letting $b \to \infty$, we conclude that

$$K = \int_{-\infty}^{\infty} e^{-x^2/2} \, dx. \tag{A.2.5}$$

Since the integral above has the value $\sqrt{2\pi}$ by (7.4.1), we have proved that $K = \sqrt{2\pi}$.

Another way of evaluating K is via the Wallis's product formula given in many calculus texts (see e.g., Courant-John, Introduction to Calculus and Analysis, Vol. 1, New York: Interscience Publishers, 1965). If this is done, then the argument above gives (A.2.5) with $K = \sqrt{2\pi}$, so that the formula for the normal integral (7.4.1) follows. This justifies the heuristic argument mentioned under (7.4.1) and shows the intimate relation between the two results named in the title of this appendix.

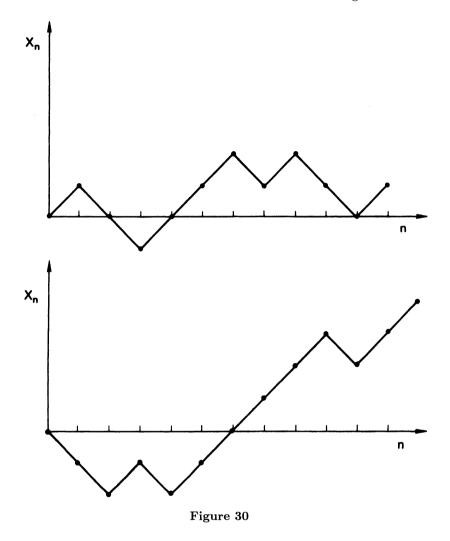
From Random Walks to Markov Chains

8.1. Problems of the wanderer or gambler

The simplest $random\ walk$ may be described as follows. A particle moves along a line by steps; each step takes it one unit to the right or to the left with probabilities p and q=1-p, respectively, where 0< p<1. For verbal convenience we suppose that each step is taken in a unit of time so that the nth step is made instantaneously at time n; furthermore we suppose that the possible positions of the particle are the set of all integers on the coordinate axis. This set is often referred to as the "integer lattice" on $R^1=(-\infty,\infty)$ and will be denoted by I. Thus the particle executes a walk on the lattice, back and forth, and continues ad infinitum. If we plot its position X_n as a function of the time n, its path is a zigzag line of which some samples are shown below in Figure 30.

A more picturesque language turns the particle into a wanderer or drunkard and the line into an endless street divided into blocks. In each unit of time, say 5 minutes, he walks one block from street corner to corner, and at each corner he may choose to go ahead or turn back with probabilities p or q. He is then taking a random walk and his track may be traced on the street with a lot of doubling and redoubling. This language suggests an immediate extension to a more realistic model where there are vertical as well as horizontal streets, regularly spaced as in parts of New York City. In this case each step may take one of the four possible directions as in Figure 31. This scheme corresponds to a random walk on the integer lattice of the plane R^2 . We shall occasionally return to this below, but for the most

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part we confine our discussion to the simplest situation of one dimension.

A mathematical formulation is near at hand. Let ξ_n be the *n*th step taken or *displacement*, so that

$$\xi_n = \begin{cases} +1 & \text{with probability } p, \\ -1 & \text{with probability } q; \end{cases}$$
 (8.1.1)

and the ξ_n 's are independent random variables. If we denote the initial position by X_0 , then the position at time n (or after n steps) is just

$$X_n = X_0 + \xi_1 + \dots + \xi_n. \tag{8.1.2}$$

Thus the random walk is represented by the sequence of random variables $\{X_n, n \geq 0\}$ which is a stochastic process in discrete time. In fact, $X_n - X_0$

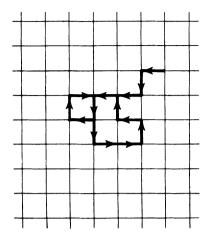


Figure 31

is a sum of independent Bernoullian random variables much studied in Chapters 5, 6, and 7. We have changed our previous notation [see, e.g., (6.3.11)] to the present one in (8.1.2) to conform with later usage in §8.3. But apart from this what is new here?

The answer is that our point of view will be new. We are going to study the entire walk, or process, as it proceeds, or develops, in the course of time. In other words, each path of the particle or wanderer will be envisioned as a possible development of the process subject to the probability laws imposed on the motion. Previously we have been interested mostly in certain quantitative characteristics of X_n (formerly S_n) such as its mean, variance, and distribution. Although the subscript n there is arbitrary and varies when $n \to \infty$, a probability like $P(a \le X_n \le b)$ concerns only the variable X_n taken one at a time, so to speak. Now we are going to probe deeper into the structure of the sequence $\{X_n, n \geq 0\}$ by asking questions that involve many of them all at once. Here are some examples. Will the moving particle ever "hit" a given point? If so, how long will this take, and will it happen before or after the particle hits some other point? One may also ask how frequently the particle hits a point or a set; how long it stays within a set, etc. Some of these questions will be made precise below and answered. In the meantime you should let your fancy go free and think up a few more such questions and perhaps relate them to concrete models of practical significance.

Let us begin with the following problem.

Problem 1. Consider the interval [0,c], where c=a+b and $a \ge 1, b \ge 1$. If the particle starts at the point "a" what is the probability that it will hit one endpoint of the interval before the other?

This is a famous problem in another setting, discussed by Fermat and Pascal and solved in general by Montmart. Two gamblers Peter and Paul play a series of games in which Peter wins with probability p and Paul wins with probability q, and the outcomes of the successive games are assumed to be independent. For instance, they may toss a coin repeatedly or play ping-pong or chess in which their skills are rated as p to q. The loser pays a dollar each time to the winner. Now if Peter has a and Paul has a at the outset and they continue to play until one of them is ruined (bankrupt), what is the probability that Peter will be ruined?

In this formulation the position of the particle at time n becomes the number of dollars Peter has after n games. Each step to the right is \$1 won by him, each step to the left is \$1 lost. If the particle reaches 0 before c, then Peter has lost all his initial capital and is ruined; on the other hand, if the particle reaches c before 0, then Paul has lost all his capital and is ruined. The game terminates when one of these eventualities occurs. Hence the historical name of "gambler's ruin problem."

We are now going to solve Problem 1. The solution depends on the following smart "put," for $1 \le j \le c - 1$:

$$u_j$$
 = the probability that the particle will reach 0 (8.1.3) before c , when it starts from j .

The problem is to find u_a , but since "a" is arbitrary we really need all the u_j 's. Indeed the idea is to exploit the relations between them and trap them together. These relations are given by the following set of difference equations:

$$u_j = pu_{j+1} + qu_{j-1}, \quad 1 \le j \le c - 1,$$
 (8.1.4)

together with the boundary conditions:

$$u_0 = 1, \quad u_c = 0.$$
 (8.1.5)

To argue (8.1.4), think of the particle as being at j and consider what will happen after taking one step. With probability p it will then be at j+1, under which hypothesis the (conditional) probability of reaching 0 before c will be u_{j+1} ; similarly with probability q it will be at j-1, under which hypothesis the said probability will be u_{j-1} . Hence the total probability u_j is equal to the sum of the two terms on the right side of (8.1.4), by an application of Proposition 2 of §5.2. This argument spelled out in the extremal cases j=1 and j=c-1 entails the values of u_0 and u_c given in (8.1.5). These are not included in (8.1.3) and strictly speaking are not well defined by the verbal description given there, although it makes sense by a kind of extrapolation.

The rest of our work is purely algebraic. Since p + q = 1 we may write the left member of (8.1.4) as $pu_j + qu_j$; after a transposition the equation

becomes

$$q(u_j - u_{j-1}) = p(u_{j+1} - u_j).$$

Using the abbreviations

$$r = \frac{q}{p}, \quad d_j = u_j - u_{j+1},$$

we obtain the basic recursion between successive differences below:

$$d_j = rd_{j-1}. (8.1.6)$$

Iterating we get $d_j = r^j d_0$; then summing by telescoping:

$$1 = u_0 - u_c = \sum_{j=0}^{c-1} (u_j - u_{j+1})$$

$$= \sum_{j=0}^{c-1} d_j = \sum_{j=0}^{c-1} r^j d_0 = \frac{1 - r^c}{1 - r} d_0$$
(8.1.7)

provided that $r \neq 1$. Next we have similarly

$$u_{j} = u_{j} - u_{c} = \sum_{i=j}^{c-1} (u_{i} - u_{i+1})$$

$$= \sum_{i=j}^{c-1} d_{j} = \sum_{i=j}^{c-1} r^{i} d_{0} = \frac{r^{j} - r^{c}}{1 - r} d_{0}.$$
(8.1.8)

It follows that

$$u_j = \frac{r^j - r^c}{1 - r^c}, \quad 0 \le j \le c.$$
 (8.1.9)

In case r = 1 we get from the penultimate terms in (8.1.7) and (8.1.8) that

$$1 = cd_0,$$

$$u_j = (c - j) d_0 = \frac{c - j}{c},$$

$$u_a = \frac{b}{c}.$$
(8.1.10)

One half of Problem 1 has been completely solved; it remains to find

 v_j = the probability that the particle will reach c before 0, when it starts from j.

Exactly the same argument shows that the set of equations (8.1.4) will be valid when the u's are replaced by v's, while the boundary conditions in (8.1.5) are merely interchanged: $v_0 = 0, v_c = 1$. Hence we can find all v_j by a similar method, which you may wish to carry out as an excellent exercise. However, there are quicker ways without this effort.

One way is perhaps easier to understand by thinking in terms of the gamblers. If we change p into q (namely r into 1/r), and at the same time j into c-j (because when Peter has j, Paul has (c-j), and vice versa), then their roles are interchanged and so u_j will go over into v_j (not v_{c-j} , why?). Making these changes in (8.1.9) and (8.1.10), we obtain

$$v_j = \frac{1 - r^j}{1 - r^c}$$
 if $p \neq q$, $v_j = \frac{j}{c}$ if $p = q$.

Now it is a real pleasure to see that in both cases we have

$$u_j + v_j = 1, \quad 0 \le j \le c.$$
 (8.1.11)

Thus as a by-product, we have solved the next problem that may have occurred to you in the course of the preceding discussion.

Problem 2. If the particle starts inside the interval [0, c], what is the probability that it will ever reach the boundary?

Since the boundary consists of the two endpoints 0 and c, the answer is given by (8.1.11) and is equal to 1. In terms of the gamblers, this means that one of them is bound to be ruined sooner or later if the game is continued without a time limit; in other words, it cannot go on forever. Now you can object that surely it is conceivable for Peter and Paul to seesaw endlessly as, e.g., indicated by the sequence +1-1+1-1+1-1... The explanation is that while this eventually is a *logical possibility* its probability is equal to zero as just shown. Namely, it will *almost never* happen in the sense discussed at the end of §7.6, and this is all we can assert.

Next, let us mention that Problem 2 can be solved without the intervention of Problem 1. Indeed, it is clear the question raised in Problem 2 is a more broad "qualitative" one that should not depend on the specific numerical answers demanded by Problem 1. It is not hard to show that even if the ξ 's in (8.1.2) are replaced by independent random variables with an arbitrary common distribution, which are not identically zero, so that we have a generalized random walk with all kinds of possible steps, the answer to Problem 2 is still the same in the broader sense that the particle will sooner or later get out of any finite interval (see, e.g., [Chung 1, Theorem 9.2.3]). Specializing to the present case where the steps are ± 1 , we see that

the particle must go through one of the endpoints before it can leave the interval [0, c]. If this conclusion tantamount to (8.1.11) is accepted with or without a proof, then of course we get $v_j = 1 - u_j$ without further calculation.

Let us state the answer to Problem 2 as follows.

Theorem 1. For any random walk (with arbitrary p), the particle will almost surely* not remain in any finite interval forever.

As a consequence, we can define a random variable that denotes the waiting time until the particle reaches the boundary. This is sometimes referred to as "absorption time" if the boundary points are regarded as "absorbing barriers," namely the particle is supposed to be stuck there as soon as it hits them. In terms of the gamblers, it is also known as the "duration of play." Let us put for $1 \le j \le c-1$:

$$S_j$$
 = the first time when the particle reaches 0 or c (8.1.12) starting from j ;

and denote its expectation $E(S_j)$ by e_j . The answer to Problem 2 asserts that S_j is almost surely finite, hence it is a random variable taking positive integer values. [Were it possible for S_j to be infinite it would not be a random variable as defined in §4.2, since " $+\infty$ " is not a number. However, we shall not elaborate on the sample space on which S_j is defined; it is not countable!] The various e_j 's satisfy a set of relations like the u_j 's, as follows:

$$e_j = pe_{j+1} + qe_{j-1} + 1, \quad 1 \le j \le c - 1,$$

 $e_0 = 0, \quad e_c = 0.$ (8.1.13)

The argument is similar to that for (8.1.4) and (8.1.5), provided we explain the additional constant "1" on the right side of the first equation above. This is the unit of time spent in taking the one step involved in the argument from j to $j \pm 1$. In the above we have tacitly assumed that all e_j are finite; for a not-so-easy proof see Exercise 48 at the end of the chapter.

The complete solution of (8.1.13) may be carried out directly as before, or more expeditiously by falling back on a standard method in solving difference equations detailed in Exercise 13. Since the general solution is not enlightening we will indicate the direct solution only in the case p = q = 1/2, which is needed in later discussion. Let $f_j = e_j - e_{j+1}$; then

$$f_j = f_{j-1} + 2, \quad f_j = f_0 + 2j,$$

$$0 = \sum_{j=0}^{c-1} f_j = c(f_0 + c - 1).$$

^{*}In general, "almost surely" means "with probability 1."

Hence $f_0 = 1 - c$, and after a little computation,

$$e_j = \sum_{i=j}^{c-1} f_i = \sum_{i=j}^{c-1} (1 - c + 2i) = j(c - j).$$
 (8.1.14)

Since the random walk is symmetric, the expected absorption time should be the same when the particle is at distance j from 0, or from c (thus at distance c-j from 0), hence it is a priori clear that $e_j = e_{c-j}$, which checks out with (8.1.14).

8.2. Limiting schemes

We are now ready to draw important conclusions from the preceding formulas. First of all, we will convert the interval [0, c] into the half-line $[0, \infty)$ by letting $c \to +\infty$. It follows from (8.1.9) and (8.1.10) that

$$\lim_{c \to \infty} u_j = \begin{cases} r^j & \text{if } r < 1, \\ 1 & \text{if } r \ge 1. \end{cases}$$
 (8.2.1)

Intuitively, this limit should mean the probability that the particle will reach 0 before "it reaches $+\infty$," starting from j; or else the probability that Peter will be ruined where he plays against an "infinitely rich" Paul, who cannot be ruined. Thus it simply represents the probability that the particle will ever reach 0 from j, or that of Peter's eventual ruin when his capital is j. This interpretation is correct and furnishes the answer to the following problem, which is a sharpening of Problem 2.

Problem 3. If the particle starts from $a (\geq 1)$, what is the probability that it will ever hit 0?

The answer is 1 if $p \leq q$; and $(q/p)^a$ if p > q. Observe that when $p \leq q$ the particle is at least as likely to go left as to go right, so the first conclusion is most plausible. Indeed, in case p < q we can say more by invoking the law of large numbers in its strong form given in §7.6. Remembering our new notation in (8.1.2) and that $E(\xi_n) = p - q$, we see that in the present context (7.6.17) becomes the assertion that almost surely we have

$$\lim_{n \to \infty} \frac{(X_n - X_0)}{n} = p - q < 0.$$

This is a much stronger assertion than that $\lim_{n\to\infty} X_n = -\infty$. Now our particle moves only one unit at a time, hence it can go to $-\infty$ only by passing through all the points to the left of the starting point. In particular, it will almost surely hit 0 from a.

In case p > q the implication for gambling is curious. If Peter has a definite advantage, then even if he has only \$1 and is playing against an unruinable Paul, he still has a chance 1-q/p to escape ruin forever. Indeed, it can be shown that in this happy event Peter will win big in the following precise sense, where X_n denotes his fortune after n games:

$$P\{X_n \to +\infty \mid X_n \neq 0 \text{ for all } n\} = 1.$$

[This is a conditional probability given the event $\{X_n \neq 0 \text{ for all } n\}$.] Is this intuitively obvious? Theorem 1 helps the argument here but does not clinch it.

When p = q = 1/2, the argument above does not apply, and since in this case there is symmetry between left and right, our conclusion may be stated more forcefully as follows.

Theorem 2. Starting from any point in a symmetric random walk, the particle will almost surely hit any point any number of times.

Proof: Let us write $i \Rightarrow j$ to mean that starting from i the particle will almost surely hit j, where $i \in I$, $j \in I$. We have already proved that if $i \neq j$, then $i \Rightarrow j$. Hence also $j \Rightarrow i$. But this implies $j \Rightarrow j$ by the obvious diagram $j \Rightarrow i \Rightarrow j$. Hence also $i \Rightarrow j \Rightarrow j \Rightarrow j \Rightarrow j \dots$, which means that starting from i the particle will hit j as many times as we desire, and note that j = i is permitted here.

We shall say briefly that the particle will hit any point in its range I infinitely often and that the random walk is recurrent (or persistent). These notions will be extended to Markov chains in §8.4.

In terms of gambling, Theorem 2 has the following implication. If the game is fair, then Peter is almost sure to win any amount set in advance as his goal, provided he can afford to go into debt for an arbitrarily large amount. For Theorem 2 only guarantees that he will eventually win say \$1000000 without any assurance as to how much he may have lost before he gains this goal. Not a very useful piece of information this—but strictly fair from Paul's viewpoint! A more realistic prediction is given in (8.1.10), which may be rewritten as

$$u_a = \frac{b}{a+b}, \quad v_a = \frac{a}{a+b};$$
 (8.2.2)

which says that the chance of Peter winning his goal b before he loses his entire capital a is in the exact inverse proportion of a to b. Thus if he has \$100, his chance of winning \$1000000 is equal to 100/1000100, or about 1 in 10000. This is about the state of affairs when he plays in a casino, even if the house does not reserve an advantage over him.

Another wrinkle is added when we let $c \to +\infty$ in the definition of e_j . The limit then represents the expected time that the particle starting at $j \geq 1$ will first reach 0 (without any constraint as to how far it can go to the right of j). Now this limit is infinite according to (8.1.14). This means even if Peter has exactly \$1 and is playing against an infinitely rich casino, he can "expect" to play a long, long time provided the game is fair. This assertion sounds fantastic as stated in terms of a single gambler, whereas the notion of mathematical expectation takes on practical meaning only through the law of large numbers applied to "ensembles." It is common knowledge that on any given day many small gamblers walk away from the casino with pocketed gains—they have happily escaped ruin because the casino did not have sufficient time to ruin them in spite of its substantial profit margin!

Let us mention another method to derive (8.2.2) which is stunning. In the case p = q we have $E(\xi_n) = 0$ for every n, and consequently we have from (8.1.2) that

$$E(X_n) = E(X_0) + E(\xi_1) + \dots + E(\xi_n) = a.$$
 (8.2.3)

In terms of the gamblers this means that Peter's expected capital remains constant throughout the play since the game is fair. Now consider the duration of play S_a in (8.1.12). It is a random variable that takes positive integer values. Since (8.2.3) is true for every such value, might it not remain so when we substitute S_a for n there? This is in general risky business but it happens to be valid here by the special nature of S_a as well as that of the process $\{X_n\}$. We cannot justify it here (see Appendix 3) but will draw the conclusion. Clearly X_{S_a} takes only the two values 0 and c by its definition; let

$$P(X_{S_a} = 0) = \rho, \quad P(X_{S_a} = c) = 1 - \rho.$$
 (8.2.4)

Then

$$E(X_{S_a}) = \rho \cdot 0 + (1 - \rho) \cdot c = (1 - \rho)c.$$

Hence $E(X_{S_a}) = a$ means

$$\rho = 1 - \frac{a}{c} = \frac{b}{a+b},$$

in agreement with (8.2.2). Briefly stated, the argument above says that the game remains fair up to and including the time of its termination. Is this intuitively obvious?

We now proceed to describe a limiting procedure that will lead from the symmetric random walk to *Brownian motion*. The English botanist Brown observed (1826) that microscopic particles suspended in a liquid are subject to continual molecular impacts and execute zigzag movements. Einstein and Smoluchovski found that in spite of their apparent irregularity these movements can be analyzed by laws of probability; in fact, the displacement over a period of time follows a normal distribution. Einstein's result (1906) amounted to a derivation of the central limit theorem (see §7.4) by the method of differential equations. The study of Brownian motion as a stochastic process was undertaken by Wiener* in 1923, preceded by Bachelier's heuristic work, and soon was developed into its modern edifice by Paul Lévy and his followers. Together with the Poisson process (§7.2) it constitutes one of the two fundamental species of stochastic processes, in both theory and application. Although the mathematical equipment allowed in this book is not adequate to treat the subject properly, it is possible to give an idea of how the Brownian motion process can be arrived at through random walk and to describe some of its basic properties.

The particle in motion observed by Brown moved of course in three-dimensional space, but we can think of its projection on a coordinate axis. Since numerous impacts are received per second, we will shorten the unit of time; but we must also shorten the unit of length in such a way as to lead to the correct model. Let δ be the new time unit – in other words, the time between two successive impacts. Thus in our previous language t/δ steps are taken by the particle in old time t. Each step is still a symmetrical Bernoullian random variable, but we now suppose that the step is of magnitude $\sqrt{\delta}$, namely for all k:

$$P(\xi_k = \sqrt{\delta}) = P(\xi_k = -\sqrt{\delta}) = \frac{1}{2}.$$

We then have

$$E(\xi_k) = 0, \quad \sigma^2(\xi_k) = \frac{1}{2}(\sqrt{\delta})^2 + \frac{1}{2}(-\sqrt{\delta})^2 = \delta.$$

Let $X_0 = 0$ so that by (8.1.2)

$$X_t = \sum_{k=1}^{t/\delta} \xi_k. (8.2.5)$$

If δ is much smaller than t, t/δ is large and may be thought of as an integer. Hence we have by Theorem 4 of §6.3:

$$E(X_t) = 0, \quad \sigma^2(X_t) = \frac{t}{\delta} \cdot \delta = t. \tag{8.2.6}$$

Furthermore if t is fixed and $\delta \to 0$, then by the De Moivre–Laplace central limit theorem (Theorem 6 of §7.3), X_t will have the normal distribution

^{*}Norbert Wiener (1894-1964), renowned U.S. mathematician, father of cybernetics.

N(0,t). This means we are letting our approximate scheme, in which the particle moves a distance of $\pm\sqrt{\delta}$ with equal probability in old time δ , go to the limit as $\delta \to 0$. This limiting scheme is the Brownian motion, also called *Wiener process*, and here is its formal definition.

Definition of Brownian Motion. A family of random variables $\{X(t)\}$, indexed by the continuous variable t ranging over $[0, \infty)$, is called the *Brownian motion* iff it satisfies the following conditions:

- (i) X(0) = 0;
- (ii) the increments $X(s_i + t_i) X(s_i)$, over an arbitrary finite set of disjoint intervals $(s_i, s_i + t_i)$, are independent random variables;
- (iii) for each $s \ge 0, t \ge 0, X(s+t) X(s)$ has the normal distribution N(0,t).

For each constant a, the process $\{X(t) + a\}$, where X(t) is just defined, is called the *Brownian motion starting at a*.

We have seen that the process constructed above by a limiting passage from symmetric random walks has property (iii). Property (ii) comes from the fact that increments over disjoint intervals are obtained by summing the displacements ξ_k in disjoint blocks; hence the sums are independent by a remark made after Proposition 6 of §5.5.

The definition above should be compared with that of a Poisson process given in §7.2, the only difference being in (iii). However, by the manner in which a Poisson process is constructed there, we know the general appearance of its paths as described under Figure 29. The situation is far from obvious for Brownian motion. It is one of Wiener's major discoveries that almost all its paths are continuous; namely, for almost all ω , the function $t \to X(t,\omega)$ is a continuous function of t in $[0,\infty)$. In practice, we can discard the null set of ω 's which yield discontinuous functions from the sample space Ω , and simply stipulate that all Brownian paths are continuous. This is a tremendously useful property that may well be added to the definition above. On the other hand, Wiener also proved that almost every path is nowhere differentiable, i.e., the curve does not have a tangent anywhere—which only goes to show that one cannot rely on intuition any more in these matters.

However, it is not hard to guess the answers to our previous questions restated for Brownian motion. In fact, the analogue in Theorem 1 holds: starting at any point, the path will go through any other point infinitely many times. Note that because of the continuity of the path this will follow from the "intermediate value theorem" in calculus once we show that it will reach out as far as we wish. Since each approximating random walk has this property, it is obvious that the Brownian motion does too. Finally, let us show that formula (8.2.2) holds also for Brownian motion, where u_a

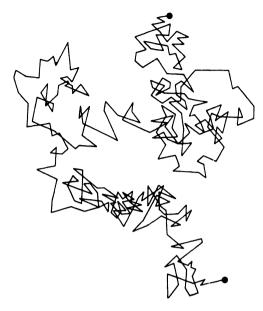


Figure 32

and v_a retain the same meanings as before but now a and c are arbitrary numbers such that 0 < a < c. Consider the Brownian motion starting at a; then it follows from property (i) that $E(X_t) = a$ for all $t \geq 0$, which is just the continuous analogue of (8.2.3). Now we substitute T_a^* for t to get $E(X_{T_a}) = a$ as before. This time, the continuity of paths assures us that at the instant T_a , the position of the particle must be exactly at 0 or at c. In fact, the word "reach" used in the definition of u_a , v_a , and T_a would have to be explained more carefully if the path could jump over the boundary. Thus we can again write (8.2.4) and get the same answer as for the symmetric random walk.

8.3. Transition probabilities

The model of random walks can be greatly generalized to that of *Markov chains*, named after A.A. Markov (see §7.6). As the saying goes, one may fail to see the forest on account of the trees. By doing away with some cumbersome and incidental features of special cases, a general theory emerges that is clearer and simpler and covers a wider range of applications. The remainder of this chapter is devoted to the elements of such a theory.

We continue to use the language of a moving particle as in the random walk scheme, and denote its range by I. This may now be a finite or infinite set of integers, and it will soon be apparent that in general no geometric

^{*}See (8.4.2) for definition of T_a with j = a there.

or algebraic structure (such as right or left, addition and subtraction) is required of I. Thus it may be an arbitrary countable set of elements, provided that we extend our definition of random variables to take values in such a set. [In §4.2 we have defined a random variable to be numerically valued.] We shall call I the state space and an element of it a state. For example, in physical chemistry a state may be a certain level of energy for an atom; in public opinion polls it may be one of the voter's possible states of mind, etc. The particle moves from state to state, and the probability law governing its change of states or transition will be prescribed as follows. There is a set of transition probabilities p_{ij} , where $i \in I, j \in I$, such that if the particle is in the state i at any time, regardless of what state i has been in before then, the probability that it will be in the state j after one step is given by p_{ij} . In symbols, if X_n denotes the state of the particle at time n, then we have

$$P\{X_{n+1} = j \mid X_n = i; A\} = P\{X_{n+1} = j \mid X_n = i\} = p_{ij}, \tag{8.3.1}$$

for an arbitrary event A determined by $\{X_0,\ldots,X_{n-1}\}$ alone. For instance, A may be a completely specified "past" of the form " $X_0=i_0,X_1=i_1,\ldots,X_{n-1}=i_{n-1}$," or a more general past event where the states i_0,\ldots,i_{n-1} are replaced by sets of states: " $X_0\in J_0,X_1\in J_1,\ldots,X_{n-1}\in J_{n-1}$." In the latter case some of these sets may be taken to be the whole space I, so that the corresponding random variables are in effect omitted from the conditioning: thus " $X_0\in J_0,X_1\in I,X_2\in J_2$ " is really just " $X_0\in J_0,X_2\in J_2$." The first equation in (8.3.1) renders the precise meaning of the phrase "regardless of prior history" and is known as the Markov property. The second equation says that the conditional probability there does not depend on the value of n; this is referred to as the stationarity (or $temporal\ homogeneity$) of the transition probabilities. Together they yield the following definition.

Definition of Markov Chain. A stochastic process $\{X_n, n \in \mathbb{N}^0\}^*$ taking values in a countable set I is called a *homogeneous Markov chain*, or *Markov chain with stationary transition probabilities*, iff (8.3.1) holds.

If the first equation in (8.3.1) holds without the second, then the Markov chain is referred to as being "nonhomogeneous," in which case the probability there depends also on n and must be denoted by $p_{ij}(n)$, say. Since we shall treat only a homogeneous chain, we mean this case when we say "Markov chain" or "chain" below without qualification.

As a consequence of the definition, we can write down the probabilities of successive transitions. Whenever the particle is in the state i_0 , and regardless of its prior history, the *conditional* probability that it will be in

^{*}It may be more convenient in some verbal descriptions to begin with n=1 rather than n=0.

the states i_1, i_2, \ldots, i_n , in the order given, during the next n steps may be suggestively denoted by the left member below and evaluated by the right member:

$$p\{\ldots i_0 \to i_1 \to i_2 \to \cdots \to i_n\} = p_{i_0 i_1} p_{i_1 i_2} \cdots p_{i_{n-1} i_n},$$
 (8.3.2)

where the five dots at the beginning serve to indicate the irrelevant and forgotten past. This follows by using (8.3.1) in the general formula (5.2.2) for joint probabilities; for instance,

$$P\{X_4 = j, X_5 = k, X_6 = l \mid X_3 = i\} = P\{X_4 = j \mid X_3 = i\}$$

$$\cdot P\{X_5 = k \mid X_3 = i, X_4 = j\} P\{X_6 = l \mid X_3 = i, X_4 = j, X_5 = k\}$$

$$= P\{X_4 = j \mid X_3 = i\} P\{X_5 = k \mid X_4 = j\} P\{X_6 = l \mid X_5 = k\}$$

$$= p_{ij} p_{jk} p_{kl}.$$

Moreover, we may adjoin any event A determined by $\{X_0, X_1, X_2\}$ alone behind the bars in the first two members above without affecting the result. This kind of calculation shows that given the state of the particle at any time, its prior history is not only irrelevant to the next transition as postulated in (8.3.1), but equally so to any future transitions. Symbolically, for any event B determined by $\{X_{n+1}, X_{n+2}, \ldots\}$, we have

$$P\{B \mid X_n = i; A\} = P\{B \mid X_n = i\}$$
(8.3.3)

as an extension of the Markov property. But that is not yet the whole story; there is a further and more sophisticated extension revolving around the three little words "at any time" italicized above, which will be needed and explained later.

It is clear from (8.3.2) that all probabilities concerning the chain are determined by the transition probabilities, provided that it starts from a fixed state, e.g., $X_0 = i$. More generally we may randomize the initial state by putting

$$P\{X_0 = i\} = p_i, \quad i \in I.$$

Then $\{p_i, i \in I\}$ is called the *initial distribution* of the chain and we have for arbitrary states i_0, i_1, \ldots, i_n :

$$P\{X_0 = i_0, X_1 = i_1, \dots, X_n = i_n\} = p_{i_0} p_{i_0 i_1} \cdots p_{i_{n-1} i_n}$$
 (8.3.4)

as the joint distribution of random variables of the process. Let us pause to take note of the special case where for every $i \in I$ and $j \in I$ we have

The right member of (8.3.4) then reduces to $p_{i_0}p_{i_1} \dots p_{i_n}$, and we see that the random variables X_0, X_1, \dots, X_n are independent with the common distribution given by $\{p_j, j \in I\}$. Thus, a sequence of independent, identically distributed, and countably valued random variables is a special case of a Markov chain, which has a much wider scope. The basic concept of such a scheme is due to Markov, who introduced it around 1907.

It is clear from the definition of p_{ij} that we have

(a)
$$p_{ij} \ge 0$$
 for every i and j ,
(b) $\sum_{j \in I} p_{ij} = 1$ for every i . (8.3.5)

Indeed, it can be shown that these are the only conditions that must be satisfied by the p_{ij} 's in order that they be the transition probabilities of a homogeneous Markov chain. In other words, such a chain can be constructed to have a given matrix satisfying those conditions as its transition matrix. Examples are collected at the end of the section.

Let us denote by $p_{ij}^{(n)}$ the probability of transition from i to j in exactly n steps, namely:

$$p_{ij}^{(n)} = P\{X_n = j \mid X_0 = i\}.$$
(8.3.6)

Thus $p_{ij}^{(1)}$ is our previous p_{ij} and we may add

$$p_{ij}^{(0)} = \delta_{ij} = \begin{cases} 0 & \text{if } i \neq j, \\ 1 & \text{if } i = j, \end{cases}$$

for convenience. The δ_{ij} above is known as *Kronecker's symbol*, which you may have seen in linear algebra. We proceed to show that for $n \geq 1$, $i \in I$, $k \in I$, we have

$$p_{ik}^{(n)} = \sum_{j} p_{ij} p_{jk}^{(n-1)} = \sum_{j} p_{ij}^{(n-1)} p_{jk}, \tag{8.3.7}$$

where the sum is over I, an abbreviation that will be frequently used below. To argue this, let the particle start from i, and consider the outcome after taking one step. It will then be in the state j with probability p_{ij} ; and conditioned on this hypothesis, it will go to the state k in n-1 more steps with probability $p_{jk}^{(n-1)}$, regardless of what i is. Hence the first equation in (8.3.7) is obtained by summing over all j according to the general formula for total probabilities; see (5.2.3) or (5.2.4). The second equation in (8.3.7) is proved in a similar way by considering first the transition in n-1 steps, followed by one more step.

For n = 2, (8.3.7) becomes

$$p_{ik}^{(2)} = \sum_{j} p_{ij} p_{jk}, \tag{8.3.8}$$

which suggests the use of matrices. Let us arrange the p_{ij} 's in the form of a matrix

$$\Pi = [p_{ij}], \tag{8.3.9}$$

so that p_{ij} is the element at the *i*th row and *j*th column. Recall that the elements of Π satisfy the conditions in (8.3.5). Such a matrix is called *stochastic*. Now the product of two square matrices $\Pi_1 \times \Pi_2$ is another such matrix whose element at the *i*th row and *j*th column is obtained by multiplying the corresponding elements of the *i*th row of Π_1 with those of the *j*th column of Π_2 , and then adding all such products. In case both Π_1 and Π_2 are the same Π , this yields precisely the right member of (8.3.8). Therefore we have

$$\Pi^2 = \Pi \times \Pi = [p_{ij}^{(2)}],$$

and it follows by induction on n and (8.3.7) that

$$\Pi^n = \Pi \times \Pi^{n-1} = \Pi^{n-1} \times \Pi = [p_{ij}^{(n)}].$$

In other words, the *n*-step transition probabilities $p_{ij}^{(n)}$ are just the elements in the *n*th power of Π . If I is the finite set $\{1,2,\ldots,r\}$, then the rule of multiplication described above is of course the same as the usual one for square matrices (or determinants) of order r. When I is an infinite set, the same rule applies but we must make sure that the resulting infinite series such as the one in (8.3.8) are all convergent. This is indeed so, by virtue of (8.3.7). We can now extend the latter as follows. For $n \in N^0$, $m \in N^0$, and $i \in I, k \in I$, we have

$$p_{ik}^{(n+m)} = \sum_{j} p_{ij}^{(n)} p_{jk}^{(m)}.$$
(8.3.10)

This set of equations is known as the *Chapman–Kolmogorov equations* [Sydney Chapman, 1888–1970, English applied mathematician]. It is simply an expression of the law of exponentiation for powers of Π :

$$\Pi^{n+m} = \Pi^n \times \Pi^m$$
.

and can be proved, either by induction of m from (8.3.7), purely algebraically, or by a probabilistic argument along the same line as that for

(8.3.7). Finally, let us record the trivial equation, valid for each $n \in \mathbb{N}^0$ and $i \in \mathbb{I}$:

$$\sum_{i} p_{ij}^{(n)} = 1. {(8.3.11)}$$

The matrix Π^n may be called the *n*-step transition matrix. Using $p_{ij}^{(n)}$ we can express joint probabilities when some intermediate states are not specified. An example will make this clear:

$$P\{X_4 = j, X_6 = k, X_9 = l \mid X_2 = i\} = p_{ij}^{(2)} p_{jk}^{(2)} p_{kl}^{(3)}.$$

We are now going to give some illustrative examples of homogeneous Markov chains, and one that is nonhomogeneous.

Example 1. $I = \{..., -2, -1, 0, 1, 2, ...\}$ is the set of all integers.

$$p_{ij} = \begin{cases} p & \text{if } j = i+1, \\ q & \text{if } j = i-1, \\ 0 & \text{otherwise;} \end{cases}$$

$$\Pi = \begin{bmatrix} \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & q & 0 & p & 0 & 0 & \cdot & \cdot \\ \cdot & \cdot & 0 & q & 0 & p & 0 & \cdot & \cdot \\ \cdot & \cdot & 0 & 0 & q & 0 & p & \cdot & \cdot & \cdot \\ \cdot & \cdot & 0 & 0 & q & 0 & p & \cdot & \cdot & \cdot \end{cases},$$

$$(8.3.12)$$

where $p+q=1, p\geq 0, q\geq 0$. This is the free random walk discussed in §8.1. In the extreme cases p=0 or q=0, it is of course deterministic (almost surely).

Example 2. $I = \{0, 1, 2, ...\}$ is the set of nonnegative integers; p_{ij} is the same as in Example 1 for $i \neq 0$; but $p_{00} = 1$, which entails $p_{0j} = 0$ for all $j \neq 0$. This is the random walk with one *absorbing state* 0. It is the model appropriate for Problem 3 in §8.2. The absorbing state corresponds to the ruin (state of bankruptcy) of Peter, whereas Paul is infinitely rich so that I is unlimited to the right.

Example 3. $I = \{0, 1, \dots, c\}, c \ge 2.$

$$\Pi = \begin{bmatrix} 1 & 0 & 0 & . & . & . & . & . \\ q & 0 & p & 0 & . & . & . & . & . \\ 0 & q & 0 & p & 0 & . & . & . & . \\ . & . & . & . & . & 0 & q & 0 & p \\ . & . & . & . & . & . & 0 & 0 & 0 & 1 \end{bmatrix}.$$

For $1 \le i \le c - 1$, the p_{ij} 's are the same as in Example 1, but

$$p_{00} = 1, \quad p_{cc} = 1. \tag{8.3.13}$$

This is the random walk with two absorbing barriers 0 and c and is appropriate for Problem 1 of §8.1. Π is a square matrix of order c+1.

Example 4. In Example 3 replace (8.3.13) by

This represents a random walk with two reflecting barriers such that after the particle reaches either endpoint of the interval [0, c], it is bound to turn back at the next step. In other words, either gambler will be given a \$1 reprieve whenever he becomes bankrupt, so that the game can go on forever—for fun! We may also eliminate the two states 0 and c, and let $I = \{1, 2, \ldots, c-1\}$,

$$\Pi = \begin{bmatrix}
q & p & 0 & 0 & \dots & \dots \\
q & 0 & p & 0 & \dots & \dots \\
\vdots & \vdots & \vdots & \ddots & \vdots & \dots & \dots \\
\vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\
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Example 5. Let $p \ge 0$, $q \ge 0$, $r \ge 0$, and p+q+r=1. In Examples 1 to 4 replace each row of the form $(\ldots q0p\ldots)$ by $(\ldots qrp\ldots)$. This means that at each step the particle may stay put, or that the game may be a draw, with probability r. When r=0, this reduces to the preceding examples.

Example 6. Let $\{\xi_n, n \geq 0\}$ be a sequence of independent integer-valued random variables such that all except possibly ξ_0 have the same distribution given by $\{a_k, k \in I\}$, where I is the set of all integers. Define X_n as in (8.1.2): $X_n = \sum_{k=0}^n \xi_k, n \geq 0$. Since $X_{n+1} = X_n + \xi_{n+1}$, and ξ_{n+1} is independent of X_0, X_1, \ldots, X_n , we have for any event A determined by X_0, \ldots, X_{n-1} alone:

$$P\{X_{n+1} = j \mid A; X_n = i\} = P\{\xi_{n+1} = j - i \mid A; X_n = i\}$$
$$= P\{\xi_{n+1} = j - i\} = a_{j-i}.$$

Hence $\{X_n, n \geq 0\}$ constitutes a homogeneous Markov chain with the transition matrix $[p_{ij}]$, where

$$p_{ij} = a_{j-i}. (8.3.14)$$

The initial distribution is the distribution of ξ_0 , which need not be the same as $\{a_k\}$. Such a chain is said to be *spatially homogeneous* as p_{ij} depends only on the difference j-i. Conversely, suppose $\{X_n, n \geq 0\}$ is a chain with the transition matrix given in (8.3.14); then we have

$$P\{X_{n+1} - X_n = k \mid X_n = i\} = P\{X_{n+1} = i + k \mid X_n = i\} = p_{i,i+k} = a_k.$$

It follows that if we put $\xi_{n+1} = X_{n+1} - X_n$, then the random variables $\{\xi_n, n \geq 1\}$ are independent (why?) and have the common distribution $\{a_k\}$. Thus a spatially as well as temporally homogeneous Markov chain is identical with the successive partial sums of independent and identically distributed integer-valued random variables. The study of the latter has been one of our main concerns in previous chapters.

In particular, Example 1 is the particular case of Example 6 with $a_1 = p$, $a_{-1} = q$; we may add $a_0 = r$ as in Example 5.

Example 7. For each $i \in I$, let p_i and q_i be two nonnegative numbers satisfying $p_i + q_i = 1$. Take I to be the set of all integers and put

$$p_{ij} = \begin{cases} p_i & \text{if } j = i+1, \\ q_i & \text{if } j = 1-1, \\ 0 & \text{otherwise.} \end{cases}$$
 (8.3.15)

In this model the particle can move only to neighboring states as in Example 1, but the probabilities may now vary with the position. The model can be generalized as in Example 5 by also allowing the particle to stay put at each position i with probability r_i , with $p_i + q_i + r_i = 1$. Observe that this example contains also Examples 2, 3, and 4 above. The resulting chain is no longer representable as sums of independent steps as in Example 6. For a full discussion of the example, see [Chung 2].

Example 8. (Ehrenfest model). This may be regarded as a particular case of Example 7 in which we have $I = \{0, 1, ..., c\}$ and

$$p_{i,i+1} = \frac{c-i}{c}, \quad p_{i,i-1} = \frac{i}{c}.$$
 (8.3.16)

It can be realized by an urn scheme as follows. An urn contains c balls, each of which may be red or black; a ball is drawn at random from it and

replaced by one of the other color. The state of the urn is the number of black balls in it. It is easy to see that the transition probabilities are as given above and the interchange can go on forever. P. and T. Ehrenfest used the model to study the transfer of heat between gas molecules. Their original urn scheme is slightly more complicated (see Exercise 14).

Example 9. Let $I = \{0, 1, 2, ...\}$ and

$$p_{i,0} = p_i, p_{i,i+1} = 1 - p_i$$
 for $i \in I$.

The p_i 's are arbitrary numbers satisfying $0 < p_i < 1$. This model is used to study a recurring phenomenon represented by the state 0. Each transition may signal an occurrence of the phenomenon, or else prolong the waiting time by one time unit. It is easy to see that the event " $X_n = k$ " means that the last time $\le n$ when the phenomenon occurred is at time n - k, where $0 \le k \le n$; in other words, there has been a waiting period equal to k units since that occurrence. In the particular case where all p_i are equal to p, we have

$$P\{X_v \neq 0 \text{ for } 1 \le v \le n-1; X_n = 0 \mid X_0 = 0\} = (1-p)^{n-1}p.$$

This gives the geometric waiting time discussed in Example 8 of §4.4.

Example 10. Let I be the integer lattice in \mathbb{R}^d , the Euclidean space of d dimensions. This is a countable set. We assume that starting at any lattice point, the particle can go only to one of the 2d neighboring points in one step, with various (not necessarily equal) probabilities. For d=1 this is just Example 1; for d=2 this is the street wanderer mentioned in §8.1. In the latter case we may represent the states by (i,i'), where i and i' are integers; then we have

$$p_{(i,i')(j,j')} = \begin{cases} p_1 & \text{if } j = i+1, j' = i', \\ p_2 & \text{if } j = i-1, j' = i', \\ p_3 & \text{if } j = i, j' = i'+1, \\ p_4 & \text{if } j = i, j' = i'-1, \end{cases}$$

where $p_1 + p_2 + p_3 + p_4 = 1$. If all these four probabilities are equal to 1/4, the chain is a symmetric two-dimensional random walk. Will the particle still hit every lattice point with probability 1? Will it do the same in three dimensions? These questions will be answered in the next section.

Example 11. (Nonhomogeneous Markov chain). Consider the Pólya urn scheme described in §5.4 with $c \ge 1$. The number of black balls in the urn is called its state so that " $X_n = i$ " means that after n drawings and insertions

there are i black balls in the urn. Clearly each transition either increases this number by c or leaves it unchanged, and we have

$$P\{X_{n+1} = j \mid X_n = i; A\} = \begin{cases} \frac{i}{b+r+nc} & \text{if } j = i+c, \\ 1 - \frac{i}{b+r+nc} & \text{if } j = i, \\ 0 & \text{otherwise;} \end{cases}$$
(8.3.17)

where A is any event determined by the outcomes of the first n-1 drawings. The probability above depends on n as well as i and j, hence the process is a nonhomogeneous Markov chain. We may also allow c=-1, which is the case of sampling without replacement and yields a finite sequence of $\{X_n: 0 \le n \le b+r\}$.

Example 12. It is trivial to define a process that is not Markovian. For instance, in Example 8 or 11, let $X_n = 0$ or 1 according to whether the nth ball drawn is red or black. Then it is clear that the probability of " $X_{n+1} = 1$ " given the values of X_1, \ldots, X_n will not in general be the same as given the value of X_n alone. Indeed, the latter probability is not very useful.

8.4. Basic structure of Markov chains

We begin a general study of the structure of homogeneous Markov chains by defining a binary relation between the states. We say "i leads to j" and write " $i \leadsto j$ " iff there exists $n \ge 1$ such that $p_{ij}^{(n)} > 0$; we say "i communicates with j" and write " $i \leadsto j$ " iff we have both $i \leadsto j$ and $j \leadsto i$. The relation " \leadsto " is transitive, namely if $i \leadsto j$ and $j \leadsto k$ then $i \leadsto k$. This follows from the inequality

$$p_{ik}^{(n+m)} \ge p_{ij}^{(n)} p_{jk}^{(m)}, \tag{8.4.1}$$

which is an algebraic consequence of (8.3.10), but perhaps even more obvious from its probabilistic meaning. For if it is possible to go from i to j in n steps, and also possible to go from j to k in m steps, then it is possible by combining these steps to go from i to k in n+m steps. Here and henceforth we shall use such expressions as "it is possible" or "one can" to mean with positive probability; but observe that even in the trivial argument just given the Markov property has been used and cannot be done without. The relation " \iff " is clearly both symmetric and transitive and may be used to divide the states into disjoint classes as follows.

Definition of Class. A *class of states* is a subset of the state space such that any two states (distinct or not) in the class communicate with each other.

This kind of classification may be familiar to you under the name of "equivalence classes." But here the relation " \iff " is not necessarily reflexive; in other words, there may be a state that does not lead to itself, hence it does not communicate with any state. Such states are simply unclassified! On the other hand, a class may consist of a single state i: this is the case when $p_{ii} = 1$. Such a state is called an absorbing state. Two classes that are not identical must be disjoint, because if they have a common element they must merge into one class via that element.

For instance, in Examples 1, 4, 5, 8 and 9 all states form a single class provided p > 0 and q > 0; as also in Example 7 provided $p_i > 0$ and $q_i > 0$ for all i. In Example 2 there are two classes: the absorbing state 0 as a singleton and all the rest as another class. Similarly in Example 3 there are three classes. In Example 6 the situation is more complicated. Suppose, for instance, the a_k 's are such that $a_k > 0$ if k is divisible by 5, and $a_k = 0$ otherwise. Then the state space I can be decomposed into five classes. Two states i and j belong to the same class if and only if i - j is divisible by 5. In other words, these classes coincide with the residue classes modulo 5. It is clear that in such a situation it would be more natural to take one of these classes as the reduced state space, because if the particle starts from any class it will (almost surely) remain in that class forever, so why bother dragging in those other states it will never get to?

In probability theory, particularly in Markov chains, the first instance of occurrence of a sequence of events is an important notion. Let j be an arbitrary state and consider the first time that the particle enters it, namely:

$$T_j(\omega) = \min\{n \ge 1 \mid X_n(\omega) = j\},$$
 (8.4.2)

where the right member reads as follows: the minimum positive value of n such that $X_n = j$. For some sample point ω , $X_n(\omega)$ may never be j, so that no value of n exists in the above and T_j is not really defined for that ω . In such a case we shall define it by the decree: $T_j(\omega) = \infty$. In common language, "it will never happen" may be rendered into "one can wait until eternity (or 'hell freezes over')." With this convention T_j is a random variable that may take the value ∞ . Let us denote the set $\{1, 2, \ldots, \infty\}$ by N_{∞} . Then T_j takes values in N_{∞} ; this is a slight extension of our general definition in §4.2.

We proceed to write down the probability distribution of T_j . For simplicity of notation we shall write $P_i\{\cdots\}$ for probability relations associated

with a Markov chain starting from the state i. We then put, for $n \in N_{\infty}$,

$$f_{ij}^{(n)} = P_i \{ T_j = n \}, \tag{8.4.3}$$

and

$$f_{ij}^* = \sum_{n=1}^{\infty} f_{ij}^{(n)} = P_i \{ T_j < \infty \}.$$
 (8.4.4)

Remember that $\sum_{n=1}^{\infty}$ really means $\sum_{1 \leq n < \infty}$; since we wish to stress the fact that the value ∞ for the superscript is not included in the summation. It follows that

$$P_i\{T_j = \infty\} = f_{ij}^{(\infty)} = 1 - f_{ij}^*.$$
 (8.4.5)

Thus $\{f_{ij}^{(n)}, n \in N_{\infty}\}$ is the probability distribution of T_j for the chain starting from i.

We can give another more explicit expression for $f_{ij}^{(n)}$, etc. as follows:

$$f_{ij}^{(1)} = p_{ij} = P_i \{ X_1 = j \},$$

$$f_{ij}^{(n)} = P_i \{ X_v \neq j \text{ for } 1 \leq v \leq n - 1; X_n = j \}, n \geq 2,$$

$$f_{ij}^{(\infty)} = P_i \{ X_v \neq j \text{ for all } v \geq 1 \},$$

$$f_{ij}^* = P_i \{ X_v = j \text{ for some } v \geq 1 \}.$$
(8.4.6)

Note that we may have i=j in the above, and "for some v" means "for at least one value of v."

The random variable T_j is called the first entrance time into the state j; the terms "first passage time" and "first hitting time" are also used. It is noteworthy that by virtue of homogeneity we have

$$f_{ij}^{(n)} = P\{X_{m+v} \neq j \text{ for } 1 \le v \le n-1; X_{m+n} = j \mid X_m = i\}$$
 (8.4.7)

for any m for which the conditional probability is defined. This kind of interpretation will be consistently used without specific mention.

The key formula connecting the $f_{ij}^{(n)}$ and $p_{ij}^{(n)}$ will now be given.

Theorem 3. For any i and j, and $1 \le n < \infty$, we have

$$p_{ij}^{(n)} = \sum_{v=1}^{n} f_{ij}^{(v)} p_{ij}^{(n-v)}.$$
 (8.4.8)

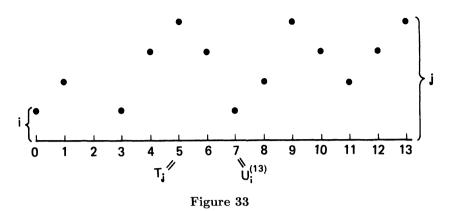
Proof: This result is worthy of a formal treatment in order to bring out the basic structure of a homogeneous Markov chain. Everything can be set down in a string of symbols:

$$\begin{aligned} p_{ij}^{(n)} &= P_i \{ X_n = j \} = P_i \{ T_j \le n; X_n = j \} = \sum_{v=1}^n P_i \{ T_j = v; X_n = j \} \\ &= \sum_{v=1}^n P_i \{ T_j = v \} P_i \{ X_n = j \mid T_j = v \} \\ &= \sum_{v=1}^n P_i \{ T_j = v \} P_i \{ X_n = j \mid X_1 \ne j, \dots, X_{v-1} \ne j, X_v = j \} \\ &= \sum_{v=1}^n P_i \{ T_j = v \} P_i \{ X_n = j \mid X_v = j \} \\ &= \sum_{v=1}^n P_i \{ T_j = v \} P_j \{ X_{n-v} = j \} \\ &= \sum_{v=1}^n f_{ij}^{(v)} p_{jj}^{(n-v)}. \end{aligned}$$

Let us explain each equation above. The first is the definition of $p_{ij}^{(n)}$; the second because $\{X_n = j\}$ implies $\{T_j \leq n\}$; the third because the events $\{T_j = v\}$ for $1 \leq v \leq n$ are disjoint; the fourth is by definition of conditional probability; the fifth is by the meaning of $\{T_j = v\}$ as given in (8.4.6); the sixth by the Markov property in (8.3.1) since $\{X_1 \neq j, \ldots, X_{v-1} \neq j\}$, as well as $\{X_0 = i\}$ implicit in the notation P_i , constitutes an event prior to the time v; the seventh is by the temporal homogeneity of a transition from j to j in n-v steps; the eighth is just notation. The proof of Theorem 3 is therefore completed.

True, a quicker verbal account can be and is usually given for (8.4.8), but if you spell out the details and pause to ask "why" at each stage, it will come essentially to a rough translation of the derivation above. This is a pattern of argument much used in a general context in the advanced theory of Markov processes, so a thorough understanding of the simplest case as this one is well worth the pains.

For $i \neq j$, formula (8.4.8) relates the transition matrix elements at (i,j) to the diagonal element at (j,j). There is a dual formula that relates them to the diagonal element at (i,i). This is obtained by an argument involving the last exit from i as the dual of first entrance into j. It is slightly more tricky in its conception and apparently known only to a few specialists. We will present it here for the sake of symmetry—and mathematical beauty. Actually the formula is a powerful tool in the theory of Markov chains, although it is not necessary for our discussions here.



Define for $n \geq 1$:

$$U_i^{(n)}(\omega) = \max\{0 \le v \le n \mid X_v(\omega) = i\};$$
 (8.4.9)

namely $U_i^{(n)}$ is the last exit time from the state i before or at the given time n. This is the dual of T_j but complicated by its dependence on n. Next we introduce the counterpart of $f_{ij}^{(n)}$, as follows:

$$g_{ij}^{(1)} = p_{ij};$$

$$g_{ij}^{(n)} = P_i \{ X_v \neq i \text{ for } 1 \le v \le n - 1; X_n = j \}, \quad 2 \le n < \infty.$$
(8.4.10)

Thus $g_{ij}^{(n)}$ is the probability of going from i to j in n steps without going through i again (for n=1 the restriction is automatically satisfied since $i \neq j$). Contrast this with $f_{ij}^{(n)}$, which may now be restated as the probability of going from i to j in n steps without going through j before. Both kinds of probability impose a taboo on certain passages and are known as taboo probabilities (see [Chung 2, §1.9] for a fuller discussion). We can now state the following result.

Theorem 4. For $i \neq j$, and $n \geq 1$, we have

$$p_{ij}^{(n)} = \sum_{v=0}^{n-1} p_{ii}^{(v)} g_{ij}^{(n-v)}.$$
 (8.4.11)

Proof: We shall imitate the steps in the proof of Theorem 3 as far as possible; thus

$$\begin{split} p_{ij}^{(n)} &= P_i \{ X_n = j \} = P_i \{ 0 \le U_i^{(n)} \le n - 1, X_n = j \} \\ &= \sum_{v=0}^{n-1} P_i \{ U_i^{(n)} = v; X_n = j \} \\ &= \sum_{v=0}^{n-1} P_i \{ X_v = i, X_u \ne i \text{ for } v + 1 \le u \le n - 1; X_n = j \} \\ &= \sum_{v=0}^{n-1} P_i \{ X_v = i \} P_i \{ X_u \ne i \text{ for } 1 \le u \le n - v - 1; X_{n-v} = j \} \\ &= \sum_{v=0}^{n-1} p_{ii}^{(v)} g_{ij}^{(n-v)}. \end{split}$$

The major difference lies in the fourth equation above, but this is obvious from the meaning of $U_i^{(n)}$. We leave the rest to the reader.

We also put

$$g_{ij}^* = \sum_{n=1}^{\infty} g_{ij}^{(n)}.$$
 (8.4.12)

However, while each term in the series above is a probability, it is not clear whether the series converges (it does, provided $i \rightsquigarrow j$; see Exercise 33). In fact, g_{ij}^* may be seen to represent the expected number of entrances in j between two successive entrances in i.

Theorems 3 and 4 may be called the first entrance and last exit decomposition formulas, respectively. Used together they work like the two hands of a human being, though one can do many things with one hand tied behind one's back, as we shall see later. Here as a preliminary ambidextrous application let us state the following little proposition as a lemma.

Lemma. $i \rightsquigarrow j$ is equivalent to $f_{ij}^* > 0$ and to $g_{ij}^* > 0$.

Proof: If $f_{ij}^* = 0$, then $f_{ij}^{(n)} = 0$ for every n and it follows from (8.4.8) that $p_{ij}^{(n)}=0$ for every n. Hence it is false that $i\leadsto j$. Conversely, if $f_{ij}^*>0$, then $f_{ij}^{(n)} > 0$ for some n; since $p_{ij}^{(n)} \ge f_{ij}^{(n)}$ from the meaning of these two proabilities, we get $p_{ij}^{(n)} > 0$ and so $i \leadsto j$. Now the argument for g_{ij}^* is exactly the same when we use (8.4.11) in

lieu of (8.4.8), demonstrating the beauty of dual thinking.

Let us admit that the preceding proof is unduly hard in the case of f_{ij}^* , since a little reflection should convince us that " $i \rightsquigarrow j$ " and " $f_{ij}^* > 0$ " both mean: "it is possible to go from i to j in some steps" (see also Exercise 31). However, is it equally obvious that " $g_{ij}^* > 0$ " means the same thing? The latter says that it is possible to go from i to j in some steps without going through i again. Hence the asserted equivalence will imply this: if it is possible to go from i to j, then it is also possible to do so without first returning to i. For example, since one can drive from New York to San Francisco, does it follow that one can do that without coming back for repairs, forgotten items, or a temporary postponement? Is this so obvious that no proof is needed?

An efficient way to exploit the decomposition formulas is to introduce generating functions associated with the sequences $\{p_{ij}^{(n)}, n \geq 0\}$ (see §6.5)

$$P_{ij}(z) = \sum_{n=0}^{\infty} p_{ij}^{(n)} z^n,$$
 $F_{ij}(z) = \sum_{n=1}^{\infty} f_{ij}^{(n)} z^n,$ $G_{ij}(z) = \sum_{n=1}^{\infty} g_{ij}^{(n)} z^n,$

where |z| < 1. We have then by substitution from (8.4.8) and inverting the order of summation:

$$P_{ij}(z) = \delta_{ij} + \sum_{n=1}^{\infty} \left(\sum_{v=1}^{n} f_{ij}^{(v)} p_{jj}^{(n-v)} \right) z^{v} z^{n-v}$$

$$= \delta_{ij} + \sum_{v=1}^{\infty} f_{ij}^{(v)} z^{v} \sum_{n=0}^{\infty} p_{jj}^{(n-v)} z^{n-v}$$

$$= \delta_{ij} + F_{ij}(z) P_{ij}(z).$$
(8.4.13)

The inversion is justified because both series are absolutely convergent for |z| < 1. In exactly the same way we obtain for $i \neq j$:

$$P_{ij}(z) = P_{ii}(z)G_{ij}(z). (8.4.14)$$

The first application is to the case i = j.

Theorem 5. For any state i we have $f_{ii}^* = 1$ if and only if

$$\sum_{n=0}^{\infty} p_{ii}^{(n)} = \infty; \tag{8.4.15}$$

if $f_{ii}^* < 1$, then we have

$$\sum_{n=0}^{\infty} p_{ii}^{(n)} = \frac{1}{1 - f_{ii}^*}.$$
(8.4.16)

Proof: From (8.4.13) with i = j and solving for $P_{ii}(z)$ we obtain

$$P_{ii}(z) = \frac{1}{1 - F_{ii}(z)}. (8.4.17)$$

If we put z = 1 above and observe that

$$P_{ii}(1) = \sum_{n=0}^{\infty} p_{ii}^{(n)}, \quad F_{ii}(1) = f_{ii}^*;$$

both assertions of the theorem follow. Let us point out that strictly speaking we must let $z \uparrow 1$ in (8.4.17) (why?), and use the following theorem from calculus. If $c_n \geq 0$ and the power series $C(z) = \sum_{n=0}^{\infty} c_n z^n$ converges for |z| < 1, then $\lim_{z \uparrow 1} C(z) = \sum_{n=0}^{\infty} c_n$, finite or infinite. This important result is called an *Abelian theorem* (after the great Norwegian mathematician Abel) and will be used again later.

The dichotomy in Theorem 5 yields a fundamental property of a state.

Definition of recurrent and nonrecurrent state. A state i is called recurrent iff $f_{ii}^* = 1$, and nonrecurrent iff $f_{ii}^* < 1$.

The adjectives "persistent" and "transient" are used by some authors for "recurrent" and "nonrecurrent." For later use let us insert a corollary to Theorem 5 here.

Corollary to Theorem 5. If j is nonrecurrent, then $\sum_{n=0}^{\infty} p_{ij}^{(n)} < \infty$ for every i. In particular, $\lim_{n\to\infty} p_{ij}^{(n)} = 0$ for every i.

Proof: For i = j, this is just (8.4.16). If $i \neq j$, this follows from (8.4.13) since

$$P_{ij}(1) = F_{ij}(1)P_{jj}(1) \le P_{jj}(1) < \infty.$$

It is easy to show that two communicating states are either both recurrent or both nonrecurrent. Thus either property pertains to a class and may be called a class property. To see this let $i \iff j$; then there exist $m \ge 1$ and $m' \ge 1$ such that $p_{ij}^{(m)} > 0$ and $p_{ji}^{(m')} > 0$. Now the same argument for (8.4.1) leads to the inequality

$$p_{ii}^{(m'+n+m)} \ge p_{ii}^{(m')} p_{ii}^{(n)} p_{ij}^{(m)}$$
.

Summing this over $n \geq 0$, we have

$$\sum_{n=0}^{\infty} p_{jj}^{(n)} \ge \sum_{n=0}^{\infty} p_{jj}^{(m'+n+m)} \ge p_{ji}^{(m')} \left(\sum_{n=0}^{\infty} p_{ii}^{(n)}\right) p_{ij}^{(m)}. \tag{8.4.18}$$

If i is recurrent, then by (8.4.15) the last term above is infinite, hence so is the first term, and this means j is recurrent by Theorem 5. Since i and j are interchangeable we have proved our assertion regarding recurrent states. The assertion regarding nonrecurrent states then follows because nonrecurrence is just the negation of recurrence.

The preceding result is nice and useful, but we need a companion that says that it is impossible to go from a recurrent to a nonrecurrent state. [The reverse passage is possible as shown by Example 3 of §8.3.] This result lies deeper and will be proved twice below by different methods. The first relies on the dual Theorems 3 and 4.

Theorem 6. If i is recurrent and $i \rightsquigarrow j$, then j is also recurrent.

Proof: There is nothing to provide if i = j, hence we may suppose $i \neq j$. We have by (8.4.13) and (8.4.14)

$$P_{ij}(z) = F_{ij}(z)P_{jj}(z), \quad P_{ij}(z) = P_{ii}(z)G_{ij}(z),$$

from which we infer

$$F_{ij}(z)P_{jj}(z) = P_{ii}(z)G_{ij}(z). (8.4.19)$$

If we let $z \uparrow 1$ as at the end of proof of Theorem 5, we obtain

$$F_{ij}(1)P_{jj}(1) = P_{ii}(1)G_{ij}(1) = \infty$$

since $G_{ij}(1) > 0$ by the lemma and $P_{ii}(1) = \infty$ by Theorem 5. Since $F_{ij}(1) > 0$ by the lemma, we conclude that $P_{jj}(1) = \infty$, hence j is recurrent by Theorem 5. This completes the proof of Theorem 6, but let us note that the formula (8.4.19) written in the form

$$\frac{P_{ii}(z)}{P_{jj}(z)} = \frac{F_{ij}(z)}{G_{ij}(z)}$$

leads to other interesting results when $z \uparrow 1$, called "ratio limit theorems" (see [Chung 2, §1.9]).

8.5. Further developments

To probe the depth of the notion of recurrence we now introduce a new "transfinite" probability, that of entering a given state *infinitely often*:

$$q_{ij} = P_i \{ X_n = j \text{ for an infinite number of values of } n \}.$$
 (8.5.1)

We have already encountered this notion in Theorem 2 of §8.2; in fact, the latter asserts in our new notation that $q_{ij} = 1$ for every i and j in a symmetric random walk. Now what exactly does "infinitely often" mean? It means "again and again, without end," or more precisely: "given any large number, say m, it will happen more than m times." This need not strike you as anything hard to grasp, but it may surprise you that if we want to express q_{ij} in symbols, it looks like this (cf. the end of §1.3):

$$q_{ij} = P_i \left\{ \bigcap_{m=1}^{\infty} \bigcup_{n=m}^{\infty} [X_n = j] \right\}.$$

For comparison let us write also

$$f_{ij}^* = P_i \left\{ \bigcup_{n=1}^{\infty} [X_n = j] \right\}.$$

However, we will circumvent such formidable formalities in our discussion below.

To begin with, it is trivial from the meaning of the probabilities that

$$q_{ij} \le f_{ij}^* \tag{8.5.2}$$

because "infinitely often" certainly entails "at least once." The next result is crucial.

Theorem 7. For any state i, we have

$$q_{ii} = \begin{cases} 1 & \text{if i is recurrent,} \\ 0 & \text{if i is nonrecurrent.} \end{cases}$$

Proof: Put $X_0 = i$, and $\alpha = f_{ii}^*$. Then α is the probability of at least one return to i. At the moment of the first return, the particle is in i and its prior history is irrelevant; hence from that moment on it will move as if making a fresh start from i ("like a newborn baby"). If we denote by R_m the event of "at least m returns," then this implies that the conditional probability $P(R_2 \mid R_1)$ is the same as $P(R_1)$ and consequently

$$P(R_2) = P(R_1R_2) = P(R_1)P(R_2 \mid R_1) = \alpha \cdot \alpha = \alpha^2.$$

Repeating this argument, we have by induction from $m \geq 1$

$$P(R_{m+1}) = P(R_m R_{m+1}) = P(R_m) P(R_{m+1} \mid R_m) = \alpha^m \cdot \alpha = \alpha^{m+1}.$$

Therefore the probability of infinitely many returns is equal to

$$\lim_{m \to \infty} P(R_m) = \lim_{m \to \infty} \alpha^m = \begin{cases} 1 & \text{if } \alpha = 1, \\ 0 & \text{if } \alpha < 1, \end{cases}$$
 (8.5.3)

proving the theorem.

Now is a good stopping time to examine the key point in the preceding proof:

$$P(R_{m+1} \mid R_m) = \alpha,$$

which is explained by considering the moment of the mth return to the initial state i and starting anew from that moment on. The argument works because whatever has happened prior to the moment is irrelevant to future happenings. Otherwise one can easily imagine a situation in which previous returns tend to inhibit a new one, such as visiting the same old tourist attraction.] This seems to be justified by the Markov property except for one essential caveat. Take m=1 for definiteness; then the moment of the first return is precisely the T_i defined in (8.4.2), and the argument above is based on applying the Markovian assumption (8.3.3) at the moment T_i . But T_i is a random variable, its value depends on the sample point ω ; can we substitute it for the constant time n in those formulas? You might think that since the latter holds true for any n, and $T_i(\omega)$ is equal to some n whatever ω may be, such a substitution must be "OK." (Indeed, we have made a similar substitution in §8.2 without justification.) The fallacy in this thinking is easily exposed,* but here we will describe the type of random variables for which the substitution is legitimate.

Given the homogeneous Markov chain $\{X_n, n \in N^0\}$, a random variable T is said to be *optional* [or a *stopping time*] iff for each n, the event $\{T=n\}$ is determined by $\{X_0, X_1, \ldots, X_n\}$ alone. An event is *prior to* T iff it is determined by $\{X_0, X_1, \ldots, X_{T-1}\}$, and *posterior to* T iff it is determined by $\{X_{T+1}, X_{T+2}, \ldots\}$. (When T=0 there is no prior event to speak of.) The state of the particle at the moment T is of course given by X_T [note: this is the random variable $\omega \to X_{T(\omega)}(\omega)$]. In case T is a constant n, these notions agree with our usual interpretation of "past" and "future" relative to the "present" moment n. In the general case they may depend on the sample point. There is nothing far-fetched in this; for instance, phrases such as "prenatal care," "postwar construction," or "the day after the locusts"

^{*}E.g., suppose $X_0=i_0\neq k$, and take $T=T_k-1$ in (8.5.5) below. Since $X_{T+1}=k$ the equation cannot hold in general.

contain an uncertain and therefore random date. When a gambler decides that he will bet on red "after black has appeared three times in a row," he is dealing with X_{T+1} , where the value of T is a matter of chance. However, it is essential to observe that these relative notions make sense by virtue of the way an optional T is defined. Otherwise if the determination of T involves the future as well as the past and present, then "pre-T" and "post-T" will be mixed up and serve no useful purpose. If the gambler can foresee the future, he would not need probability theory! In this sense an optional time has also been described as being "independent of the future"; it must have been decided upon as an "option" without the advantage of clairvoyance.

We can now formulate the following extension of (8.3.3). For any optional T, any event A prior to T, and any event B posterior to T, we have

$$P\{B \mid X_T = i; A\} = P\{B \mid X_T = i\}; \tag{8.5.4}$$

and in particular for any state i and j:

$$P\{X_{T+1} = j \mid X_T = i; A\} = p_{ij}. \tag{8.5.5}$$

This is known as the *strong Markov property*. It is actually implied by the apparently weaker form given in (8.3.3), hence also in the original definition (8.3.1). Probabilists used to announce the weak form and use the strong one without mentioning the difference. Having flushed the latter out in the open we will accept it as the definition for a homogeneous Markov chain. For a formal proof see [Chung 2, §I.13]. Let us observe that the strong Markov property was needed as early as in the proof of Theorem 2 of §8.2, where it was deliberately concealed in order not to sound a premature alarm. Now it is time to look back with understanding.

To return to Theorem 7 we must now verify that the T_i used in the proof there is indeed optional. This has been effectively shown in (8.4.6), for the event

$${T_i = n} = {X_v \neq i \text{ for } 1 \le v \le n - 1; X_n = i}$$

is clearly determined by $\{X_1, \ldots, X_n\}$ only. This completes the rigorous proof of Theorem 7, to which we add a corollary.

Corollary to Theorem 7. For any i and j,

$$q_{ij} = \begin{cases} f_{ij}^* & \text{if } j \text{ is recurrent,} \\ 0 & \text{if } f \text{ is nonrecurrent.} \end{cases}$$

Proof: This follows at once from the theorem and the relation:

$$q_{ij} = f_{ij}^* q_{jj}. (8.5.6)$$

For to enter j infinitely many times means to enter it at least once and then return to it infinitely many times. As in the proof of Theorem 8, the reasoning involved here is based on the strong Markov property.

The next result shows the power of "thinking infinite."

Theorem 8. If i is recurrent and $i \rightsquigarrow j$, then

$$q_{ij} = q_{ji} = 1. (8.5.7)$$

Proof: The conclusion implies that $i \leftrightarrow j$ and that j is recurrent by the corollary above. Thus the following proof contains a new proof of Theorem 6.

Let us note that for any two events A and B, we have $A \subset AB \cup B^c$, and consequently

$$P(A) \le P(B^c) + P(AB).$$
 (8.5.8)

Now consider

$$A = \{ \text{enter } i \text{ infinitely often} \},$$

 $B = \{ \text{enter } j \text{ at least once} \}.$

Then $P_i(A) = q_{ii} = 1$ by Theorem 7 and $P_i(B^c) = 1 - f_{ij}^*$. As for $P_i(AB)$ this means the probability that the particle will enter j at some finite time and thereafter enter i infinitely many times, because "infinite minus finite is still infinite." Hence if we apply the strong Markov property at the first entrance time into j, we have

$$P(AB) = f_{ij}^* q_{ji}.$$

Substituting into the inequality (8.5.8), we obtain

$$1 = q_{ii} \le 1 - f_{ij}^* + f_{ij}^* q_{ji}$$

and so

$$f_{ij}^* \le f_{ij}^* q_{ji}.$$

Since $f_{ij}^* > 0$ this implies $1 \le q_{ji}$, hence $q_{ji} = 1$. Since $q_{ji} \le f_{ji}^*$ it follows that $f_{ji}^* = 1$, and so $j \leadsto i$. Thus i and j communicate and therefore j is recurrent by (8.4.18). Knowing this we may interchange the roles of i and j in the preceding argument to infer $q_{ij} = 1$.

Corollary. In a recurrent class (8.5.7) holds for any two states i and j.

When the state space of a chain forms a single recurrent class, we shall call the chain recurrent; similarly for "nonrecurrent." The state of affairs for a recurrent chain described in the preceding corollary is precisely that for a symmetric random walk in Theorem 2 of §8.2. In fact, the latter is a particular case as we now proceed to show.

We shall apply the general methods developed above to the case of random walk discussed in §8.1, namely Example 1 of §8.3. We begin by evaluating $p_{ii}^{(n)}$. This is the probability that the particle returns to its initial position i in exactly n steps. Hence $p_{ii}^{(2n-1)} = 0$ for $n \ge 1$; and in the notation of (8.1.2)

$$p_{ii}^{(2n)} = P\{\xi_1 + \dots + \xi_{2n} = 0\} = {2n \choose n} p^n q^n$$
 (8.5.9)

by Bernoulli's formula (7.3.1), since there must be n steps to the right and n steps to the left, in some order. Thus we obtain the generating function

$$P_{ii}(z) = \sum_{n=0}^{\infty} {2n \choose n} (pqz^2)^n.$$
 (8.5.10)

Recalling the general binomial coefficients from (5.4.4), we record the pretty identity:

$$\binom{-\frac{1}{2}}{n} = \frac{(-1)^n 1 \cdot 3 \cdots (2n-1)}{2^n \cdot n!} = \frac{(-1)^n}{2^{2n}} \binom{2n}{n},\tag{8.5.11}$$

where the second equation is obtained by multiplying both the denominator and numerator of its left member by $2^n \cdot n! = 2 \cdot 4 \cdots (2n)$. Substituting into (8.5.10), we arrive at the explicit analytic formula:

$$P_{ii}(z) = \sum_{n=0}^{\infty} {\binom{-\frac{1}{2}}{n}} (-4pqz^2)^n = (1 - 4pqz^2)^{-1/2},$$
 (8.5.12)

where the second member is the binomial (Taylor's) series of the third member.

It follows that

$$\sum_{n=0}^{\infty} p_{ii}^{(n)} = P_{ii}(1) = \lim_{z \uparrow 1} P_{ii}(z) = \lim_{z \uparrow 1} (1 - 4pqz^2)^{-1/2}.$$
 (8.5.13)

Now $4pq = 4p(1-p) \le 1$ for $0 \le p \le 1$; and = 1 if and only if p = 1/2 (why?). Hence the series above diverges if p = 1/2 and converges if $p \ne 1/2$. By Theorem 5, i is recurrent if and only if p = 1/2. The calculations above

do not depend on the integer i because of spatial homogeneity. Thus for p = 1/2 the chain is recurrent; otherwise it is nonrecurrent. In other words, the random walk is recurrent if and only if it is symmetric.

There is another method of showing this directly from (8.5.9), without the use of generating functions. For when p = 1/2 we have

$$p_{ii}^{(2n)} = {2n \choose n} \frac{1}{2^{2n}} \sim \frac{1}{\sqrt{\pi n}},$$
 (8.5.14)

by (7.3.6) as an application of Stirling's formula. Hence by the comparison test for positive series, the series in (8.5.13) diverges because $\sum_{n} 1/\sqrt{n}$ does. This method has the merit of being applicable to random walks in higher dimensions. Consider the symmetric random walk in R^2 (Example 10 of §8.3 with all four probabilities equal to 1/4). To return from any state (i,j) to (i,j) in 2n steps means that: for some k, $0 \le k \le n$, the particle takes, in some order, k steps each to the east and west, and n-k steps each to the north and south. The probability for this, by the multinomial formula (6.4.6), is equal to

$$p_{(i,j)(i,j)}^{(2n)} = \frac{1}{4^{2n}} \sum_{k=0}^{n} \frac{(2n)!}{k! \, k! \, (n-k)! \, (n-k)!}$$

$$= \frac{(2n)!}{4^{2n} n! \, n!} \sum_{k=0}^{n} \binom{n}{k}^{2} = \frac{(2n)!}{4^{2n} n! \, n!} \binom{2n}{n} = \left[\frac{1}{2^{2n}} \binom{2n}{n}\right]^{2},$$
(8.5.15)

where in the penultimate equation we have used a formula given in Exercise 28 of Chapter 3. The fact that this probability turns out to be the exact square of the one in (8.5.14) is a pleasant coincidence. [It is not due to any apparent independence between the two components of the walk along the two coordinate axes.] It follows by comparison with (8.5.14) that

$$\sum_{n} p_{(i,j)(i,j)}^{(2n)} \sim \sum_{n} \frac{1}{\pi n} = \infty.$$

Hence another application of Theorem 5 shows that the symmetric random walk in the plane as well as on the line is a recurrent Markov chain. A similar but more complicated argument shows that it is nonrecurrent in R^d for $d \geq 3$, because the probability analogous to that in (8.5.15) is bounded by $c/n^{d/2}$ (where c is a constant), and $\sum_n 1/n^{d/2}$ converges for $d \geq 3$. These results were first discovered by Pólya in 1921. The nonsymmetric case can be treated by using the normal approximation given in (7.3.13), but there the nonrecurrence is already implied by the strong law of large numbers as in R^1 ; see §8.2.

As another illustration, we will derive an explicit formula for $f_{ii}^{(n)}$ in case p = 1/2. By (8.4.17) and (8.5.12), we have

$$F_{ii}(z) = 1 - \frac{1}{P_{ii}(z)} = 1 - (1 - z^2)^{1/2}.$$

Hence another expansion by means of binomial series gives

$$\begin{split} F_{ii}(z) &= 1 - \sum_{n=0}^{\infty} \binom{\frac{1}{2}}{n} (-z^2)^n \\ &= \frac{1}{2} z^2 + \sum_{n=2}^{\infty} \frac{1 \cdot 3 \cdots (2n-3)}{2^n \cdot n!} z^{2n}. \end{split}$$

Thus $f_{ii}^{(2n-1)} = 0$; and

$$f_{ii}^{(2n)} = \frac{1}{2^{2n}} {2n \choose n} \frac{1}{2n-1}, \quad n \ge 1;$$
 (8.5.16)

by a calculation similar to (8.5.11). In particular, we have

Comparison with (8.5.14) shows that

$$f_{ii}^{(2n)} \sim rac{1}{2\sqrt{\pi} \, n^{3/2}}$$

and so $\sum_{n=1}^{\infty} n f_{ii}^{(n)} = \infty$. This can also be gotten by calculating $F'_{ii}(1)$. Thus, although return is almost certain, the expected time before return is infinite. This result will be seen in a moment to be equivalent to the remark made in §8.2 that $e_1 = \infty$.

We can calculate $f_{ii}^{(n)}$ for any i and j in a random walk by a similar method. However, sometimes a combinatorial argument is quicker and more revealing. For instance, we have

$$f_{00}^{(2n)} = \frac{1}{2}f_{10}^{(2n-1)} + \frac{1}{2}f_{-1,0}^{(2n-1)} = f_{10}^{(2n-1)} = f_{01}^{(2n-1)}.$$
 (8.5.17)

To argue this let the particle start from 0 and consider the outcome of its first step as in the derivation of (8.1.4); then use the symmetry and spatial homogeneity to get the rest. The details are left to the reader.

8.6. Steady state

In this section we consider a recurrent Markov chain, namely we suppose that the state space forms a single recurrent class.

After the particle in such a chain has been in motion for a long time, it will be found in various states with various probabilities. Do these settle down to limiting values? This is what the physicists and engineers call a "steady state" (distribution).* They are accustomed to thinking in terms of an "ensemble" or large number of particles moving according to the same probability laws and independently of one another, such as in the study of gaseous molecules. In the present case the laws are those pertaining to a homogeneous Markov chain as discussed in the preceding sections. After a long time, the proportion (percentage) of particles to be found in each state gives approximately the steady-state probability of that state. Note the double usage of the word "state" in the last sentence; we shall use "stationary" for the adjective "steady-state." In effect, this is the frequency interpretation of probability mentioned in Example 3 of §2.1, in which the limiting proportions are taken to determine the corresponding probabilities. In our language, if the particle starts from the state i, then the probability of the set of paths in which it moves to state j at time n, namely $\{\omega \mid$ $X_n(\omega) = j$, is given by $P_i\{X_n = j\} = p_{ij}^{(n)}$. We are therefore interested in the asymptotic behavior of $p_{ij}^{(n)}$ as $n \to \infty$. It turns out that a somewhat more amenable quantity is its average value over a long period of time, namely

$$\frac{1}{n+1} \sum_{v=0}^{n} p_{ij}^{(v)} \quad \text{or} \quad \frac{1}{n} \sum_{v=1}^{n} p_{ij}^{(v)}. \tag{8.6.1}$$

The difference between these two averages is negligible for large n but we shall use the former. This quantity has a convenient interpretation as follows. Fix our attention on a particular state j and imagine that a counting device records the number of time units the particle spends in j. This is done by introducing the random variables below that count 1 for the state j but 0 for any other state:

$$\xi_v(j) = \begin{cases} 1 & \text{if } X_v = j, \\ 0 & \text{if } X_v \neq j. \end{cases}$$

We have used such indicators, e.g., in (6.4.11). Next we put

$$N_n(j) = \sum_{v=0}^n \xi_v(j),$$

^{*}Strange to relate, they call a "distribution" a "state"!

which represents the total occupation time of the state j in n steps. Now if E_i denotes the mathematical expectation associated with the chain starting from i [this is a conditional expectation; see end of §5.2], we have

$$E_i(\xi_v(j)) = p_{ij}^{(v)},$$

and so by Theorem 1 of $\S6.1$:

$$E_i(N_n(j)) = \sum_{v=0}^n E_i(\xi_v(j)) = \sum_{v=0}^n p_{ij}^{(v)}.$$
 (8.6.2)

Thus the quantity in (8.6.1) turns out to be the average expected occupation time.

In order to study this we consider first the case i = j and introduce the expected return time from j to j as follows:

$$m_{jj} = E_j(T_j) \sum_{v=1}^{\infty} v f_{jj}^{(v)},$$
 (8.6.3)

where T_j is defined in (8.4.2). Since j is a recurrent state we know that T_j is almost surely finite, but its expectation may be finite or infinite. We shall see that the distinction between these two cases is essential.

Here is the heuristic argument linking (8.6.2) and (8.6.3). Since the time required for a return is m_{jj} units on the basis of expectation, there should be about n/m_{jj} such returns in a span of n time units on the same basis. In other words the particle spends about n/m_{jj} units of time in the state j during the first n steps, namely $E_j(N_n(j)) \approx n/m_{jj}$. The same argument shows that it makes no difference whether the particle starts from j or any other state i, because after the first entrance into j the initial i may be forgotten and we are concerned only with the subsequent returns from j to j. Thus we are led to the following limit theorem.

Theorem 9. For any i and j we have

$$\lim_{n \to \infty} \frac{1}{n+1} \sum_{v=0}^{n} p_{ij}^{(v)} = \frac{1}{m_{jj}}.$$
 (8.6.4)

The argument indicated above can be made rigorous by invoking a general form of the strong law of large numbers (see §7.5), applied to the successive return times that form a sequence of independent and identically distributed random variables. Unfortunately the technical details are above the level of this book. There is another approach, which relies on a powerful analytical result due to Hardy and Littlewood. [This is the same Hardy as in the Hardy-Weinberg theorem of §5.6.] It is known as a *Tauberian theorem*

(after Tauber, who first found a result of the kind) and may be stated as follows.

Theorem 10. If $A(z) = \sum_{n=0}^{\infty} a_n z^n$, where $a_n \ge 0$ for all n and the series converges for $0 \le z < 1$, then we have

$$\lim_{n \to \infty} \frac{1}{n+1} \sum_{v=0}^{n} a_v = \lim_{z \to 1} (1-z) A(z). \tag{8.6.5}$$

To get a feeling for this theorem, suppose all $a_n = c > 0$. Then

$$A(z) = c \sum_{n=0}^{\infty} z^n = \frac{c}{1-z}$$

and the relation in (8.6.5) reduces to the trivial identity

$$\frac{1}{n+1} \sum_{n=0}^{n} c = c = (1-z) \frac{c}{1-z}.$$

Now take A(z) to be

$$P_{ij}(z) = \sum_{n=0}^{\infty} p_{ij}^{(n)} z^n = F_{ij}(z) P_{ij}(z) = \frac{F_{ij}(z)}{1 - F_{jj}(z)},$$

where the last two equations come from (8.4.13) and (8.4.17). Then we have

$$\lim_{z \to 1} (1 - z) P_{ij}(z) = F_{ij}(1) \lim_{z \to 1} \frac{1 - z}{1 - F_{ij}(z)} = \lim_{z \to 1} \frac{1 - z}{1 - F_{ij}(z)}$$

since $F_{ij}(1) = f_{ij}^* = 1$ by Theorem 8 of §8.5. The last-written limit may be evaluated by l'Hospital rule:

$$\lim_{z \to 1} \frac{(1-z)'}{(1-F_{jj}(z))'} = \lim_{z \to 1} -\frac{-1}{F'_{jj}(z)} = \frac{1}{F'_{jj}(1)},$$

where "t" stands for differentiation with respect to z. Since $F'_{jj}(z) = \sum_{v=1}^{\infty} v f_{jj}^{(v)} z^{v-1}$ we have $F'_{jj}(1) = \sum_{v=1}^{\infty} f_{jj}^{(v)} = m_{jj}$, and so (8.6.4) is a special case of (8.6.5).

We now consider a finite state space I in order not to strain our mathematical equipment. The finiteness of I has an immediate consequence.

Theorem 11. If I is finite and forms a single class (namely if there are only a finite number of states and they all communicate with each other), then the chain is necessarily recurrent.

Proof: Suppose the contrary; then each state is transient and so almost surely the particle can spend only a finite number of time units in it by Theorem 7. Since the number of states is finite, the particle can spend altogether only a finite number of time units in the whole space I. But time increases ad infinitum and the particle has nowhere else to go. This absurdity proves that the chain must be recurrent. (What then can the particle do?)

Let $I = \{1, 2, ..., l\}$, and put

$$x=(x_1,\ldots,x_l),$$

which is a row vector of l components. Consider the "steady-state equation"

$$x = x\Pi \quad \text{or} \quad x(\Delta - \Pi) = 0, \tag{8.6.6}$$

where Δ is the identity matrix with l rows and l columns: $\Delta = (\delta_{ij})$, and Π is the transition matrix in (8.3.9). This is a system of l linear homogeneous equations in l unknowns. Now the determinant of the matrix $\Delta - \Pi$

$$\begin{vmatrix} 1 - p_{11} & -p_{12} & \cdots & -p_{1l} \\ -p_{21} & 1 - p_{22} & \cdots & -p_{2l} \\ \vdots & \vdots & \vdots & \vdots \\ -p_{l1} & -p_{l2} & \cdots & 1 - p_{ll} \end{vmatrix}$$

is equal to zero because the sum of all elements in each row is $1 - \sum_j p_{ij} = 0$. Hence we know from linear algebra that the system has a nontrivial solution, namely one that is not the zero vector. Clearly if x is a solution, then so is $cx = (cx_1, cx_2, \ldots, cx_l)$ for any constant c. The following theorem identifies all solutions where I is a single finite class.

We shall write

$$w_j = \frac{1}{m_{jj}}, \quad j \in I,$$

 $w = (w_1, w_2, \dots, w_l),$ (8.6.7)

and \sum_{i} for $\sum_{i \in I}$ below.

Theorem 12. If I is finite and forms a single class, then

- (i) w is a solution of (8.6.6),
- (ii) $\sum_{i} w_j = 1$,
- (iii) $w_j > 0$ for all j,
- (iv) any solution of (8.6.6) is a constant multiple of w.

Proof: We have from (8.3.7), for every $v \ge 0$,

$$p_{ik}^{(v+1)} = \sum_{j} p_{ij}^{(v)} p_{jk}.$$

Taking an average over v, we get

$$\frac{1}{n+1} \sum_{v=0}^{n} p_{ik}^{(v+1)} = \sum_{j} \left(\frac{1}{n+1} \sum_{v=0}^{n} p_{ij}^{(v)} \right) p_{jk}.$$

The left member differs from $1/(n+1)\sum_{v=0}^{n}p_{ik}^{(v)}$ by $1/(n+1)(p_{ik}^{(n+1)}-p_{ik}^{(0)})$, which tends to zero as $n\to\infty$; hence its limit is equal to w_k by Theorem 9. Since I is finite, we may let $n\to\infty$ term by term in the right member. This yields

$$w_k = \sum_j w_j p_{jk},$$

which is $w = w\Pi$; hence (i) is proved. We can now iterate:

$$w = w\Pi = (w\Pi)\Pi = w\Pi^2 = (w\Pi)\Pi^2 = w\Pi^3 = \cdots,$$
 (8.6.8)

to obtain $w = w\Pi^n$, or explicitly for $n \ge 1$:

$$w_k = \sum_{j} w_j p_{jk}^{(n)}. (8.6.9)$$

Next we have $\sum_{j} p_{ij}^{(v)} = 1$ for every i and $v \ge 1$. Taking an average over v, we obtain

$$\frac{1}{n+1} \sum_{v=0}^{n} \sum_{j} p_{ij}^{(v)} = 1.$$

It follows that

$$\sum_{j} w_{j} = \sum_{j} \lim_{n \to \infty} \left(\frac{1}{n+1} \sum_{v=0}^{n} p_{ij}^{(v)} \right) = \lim_{n \to \infty} \frac{1}{n+1} \sum_{v=0}^{n} \sum_{j} p_{ij}^{(v)} = 1,$$

where the second equation holds because I is finite. This establishes (ii) from which we deduce that at least one of the w_j 's, say w_i , is positive. For any k we have $i \leadsto k$ and so there exists n such that $p_{ik}^{(n)} > 0$. Using this value of n in (8.6.9), we see that w_k is also positive. Hence (iii) is true.

Finally suppose x is any solution of (8.6.6). Then $x = x\Pi^v$ for every $v \ge 1$ by iteration as before, and

$$x = \frac{1}{n+1} \sum_{v=0}^{n} x \Pi^v$$

by averaging. In explicit notation this is

$$x_k = \sum_{j} x_j \left(\frac{1}{n+1} \sum_{v=0}^{n} p_{jk}^{(v)} \right).$$

Letting $n \to \infty$ and using Theorem 9, we obtain

$$x_k = \left(\sum_j x_j\right) w_k.$$

Hence (iv) is true with $c = \sum_{j} x_{j}$. Theorem 12 is completely proved.

We call $\{w_j, j \in I\}$ the stationary (steady-state) distribution of the Markov chain. It is indeed a probability distribution by (ii). The next result explains the meaning of the word "stationary."

Theorem 13. Suppose that we have, for every j,

$$P\{X_0 = j\} = w_j; (8.6.10)$$

then the same is true when X_0 is replaced by any $X_n, n \geq 1$. Furthermore the joint probability

$$P\{X_{n+v} = j_v, 0 \le v \le l\}$$
(8.6.11)

for arbitrary j_v is the same for all $n \geq 0$.

Proof: We have by (8.6.9)

$$P\{X_n = j\} = \sum_{i} P\{X_0 = i\} P_i\{X_n = j\} = \sum_{i} w_i p_{ij}^{(n)} = w_j.$$

Similarly the probability in (8.6.11) is equal to

$$P\{X_n = j_0\}p_{j_0j_1}\cdots p_{j_{l-1}j_l} = w_{j_0}p_{j_0j_1}\cdots p_{j_{l-1}j_l},$$

which is the same for all n.

Thus, with the stationary distribution as its initial distribution, the chain becomes a stationary process as defined in §5.4. Intuitively, this means

that if a system is in its steady state, it will hold steady indefinitely there, that is, so far as distributions are concerned. Of course, changes go on in the system, but they tend to balance out to maintain an overall equilibrium. For instance, many ecological systems have gone through millions of years of transitions and may be considered to have reached their stationary phase—until human intervention abruptly altered the course of evolution. However, if the new process is again a homogeneous Markov chain as supposed here, then it too will settle down to a steady state in due time according to our theorems.

The practical significance of Theorem 12 is that it guarantees a solution of (8.6.6) that satisfies conditions (ii) and (iii). In order to obtain this solution, we may proceed as follows. Discard one of the l equations and solve the remaining equations for w_2, \ldots, w_l in terms of w_1 . These are of the form $w_j = c_j w_1, 1 \le j \le l$, where $c_1 = 1$. The desired solution is then given by

$$w_j = \frac{c_j}{\sum\limits_{j=1}^{l} c_j}, \quad 1 \le j \le l.$$

Example 13. A switch may be on or off; call these two positions states 1 and 2. After each unit of time the state may hold or change, but the respective probabilities depend only on the present position. Thus we have a homogeneous Markov chain with $I = \{1, 2\}$ and

$$\Pi = egin{bmatrix} p_{11} & p_{12} \ p_{21} & p_{22} \end{bmatrix},$$

where all elements are supposed to be positive. The steady-state equations are

$$(1 - p_{11})x_1 - p_{21}x_2 = 0,$$

$$-p_{12}x_1 + (1 - p_{22})x_2 = 0.$$

Clearly the second equation is just the negative of the first and may be discarded. Solving the first equation we get

$$x_2 = \frac{1 - p_{11}}{p_{21}} x_1 = \frac{p_{12}}{p_{21}} x_1.$$

Thus

$$w_1 = \frac{p_{21}}{p_{12} + p_{21}}, \quad w_2 = \frac{p_{12}}{p_{12} + p_{21}}.$$

In view of Theorem 9, this means: in the long run the switch will be on or off for a total amount of time in the ratio of $p_{21}: p_{12}$.

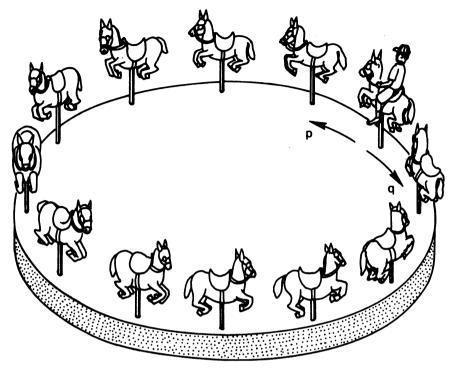


Figure 34

Example 14. At a carnival Daniel won a prize for free rides on the merrygo-round. He therefore took "infinitely many" rides, but each time when the bell rang he moved onto the next hobby-horse forward or backward, with probability p or q = 1 - p. What proportion of time was he on each of these horses?

This may be described as "random walk on a circle." The transition matrix looks like this:

$\begin{bmatrix} 0 \\ q \\ 0 \end{bmatrix}$	$egin{array}{c} p \ 0 \ q \end{array}$	$egin{matrix} 0 \\ p \\ 0 \end{matrix}$	$egin{matrix} 0 \ 0 \ p \end{matrix}$	•••	0 0 0	0 0 0	$egin{pmatrix} q \\ 0 \\ 0 \\ \end{bmatrix}$
$\begin{vmatrix} 0 \\ 0 \end{vmatrix}$	$0 \\ 0$	$0 \\ 0$	$0 \\ 0$		$\frac{0}{q}$	$p \\ 0$	p = 0

The essential feature of this matrix is that the elements in each column (as well as in each row) add up to 1. In general notation, this means that we

have for every $j \in I$:

$$\sum_{i \in I} p_{ij} = 1. \tag{8.6.12}$$

Such a matrix is called *doubly stochastic*. Now it is trivial that under the condition (8.6.12), x = (1, 1, ..., 1) where all components are equal to 1, is a solution of the equation (8.6.6). Since the stationary distribution w must be a multiple of this by (iv) of Theorem 12, and also satisfy (iii), it must be

$$w = \left(\frac{1}{l}, \frac{1}{l}, \dots, \frac{1}{l}\right),\,$$

where as before l is the number of states in I. This means if Daniel spent 4 hours on the merry-go-round and there are 12 horses, his occupation time of each horse is about 20 minutes, provided that he changed horses sufficiently many times to make the limiting relation in (8.6.4) operative.

For a recurrent Markov chain in an infinite state space, Theorem 12 must be replaced by a drastic dichotomy as follows:

- (a) either all $w_j > 0$, then (ii) and (iii) hold as before, and Theorem 13 is also true;
- (b) or all $w_j = 0$.

The chain is said to be positive-recurrent (or strongly ergodic) in case (a) and null-recurrent (or weakly ergodic) in case (b). The symmetric random walk discussed in §§8.1–8.2 is an example of the latter (see Exercise 38). It can be shown (see [Chung 2, §I.7]) that if the equation (8.6.6) has a solution $x=(x_1,x_2,\ldots)$ satisfying the condition $0<\sum_j|x_j|<\infty$, then in fact all $x_j>0$ and the stationary distribution is given by

$$w_j = \frac{x_j}{\sum_j x_j}, \quad j \in I.$$

The following example illustrates this result.

Example 15. Let $I = \{0, 1, 2, ...\}$, and $p_{ij} = 0$ for |i - j| > 1, whereas the other p_{ij} 's are arbitrary positive numbers. These must then satisfy the equation

$$p_{j,j-1} + p_{jj} + p_{j,j+1} = 1 (8.6.13)$$

for every j. This may be regarded as a special case of Example 7 in §8.3 with $p_{0,-1}=0$ and a consequent reduction of state space. It may be called

a simple birth-and-death process (in discrete time) in which j is the population size and $j \to j+1$ or $j \to j-1$ corresponds to a single birth or death. The equation (8.6.6) becomes

$$x_0 = x_0 p_{00} + x_1 p_{10},$$

$$x_j = x_{j-1} p_{j-1,j} + x_j p_{jj} + x_{j+1} p_{j+1,j}, \quad j \ge 1.$$
(8.6.14)

This is an infinite system of linear homogeneous equations, but it is clear that all possible solutions can be obtained by assigning an arbitrary value to x_0 , and then solve for x_1, x_2, \ldots successively from the equations. Thus we get

$$x_1 = \frac{p_{01}}{p_{10}} x_0,$$

$$x_2 = \frac{1}{p_{21}} \{ x_1 (1 - p_{11}) - x_0 p_{01} \} = \frac{p_{01} (1 - p_{11} - p_{10})}{p_{21} p_{10}} x_0 = \frac{p_{01} p_{12}}{p_{10} p_{21}} x_0.$$

It is easy to guess (perhaps after a couple more steps) that we have in general

$$x_j = c_j x_0$$
 where $c_0 = 1$, $c_j = \frac{p_{01} p_{12} \cdots p_{j-1,j}}{p_{10} p_{21} \cdots p_{j,j-1}}$, $j \ge 1$. (8.6.15)

To verify this by induction, let us assume that $p_{j,j-1}x_j = p_{j-1,j}x_{j-1}$; then we have by (8.6.14) and (8.6.13)

$$p_{j+1,j}x_{j+1} = (1 - p_{jj})x_j - p_{j-1,j}x_{j-1}$$
$$= (1 - p_{jj} - p_{j,j-1})x_j = p_{j,j+1}x_j.$$

Hence this relation holds for all j and (8.6.15) follows. We therefore have

$$\sum_{j=0}^{\infty} x_j = \left(\sum_{j=0}^{\infty} c_j\right) x_0, \tag{8.6.16}$$

and the dichotomy cited above is as follows, provided that the chain is recurrent. It is easy to see that this is true in case (a).

Case (a). If $\sum_{j=0}^{\infty} c_j < \infty$, then we may take $x_0 = 1$ to obtain a solution satisfying $\sum_j x_j < \infty$. Hence the chain is positive-recurrent and the stationary distribution is given by

$$w_j = \frac{c_j}{\sum_{j=0}^{\infty} c_j}, \quad j \ge 0.$$

Case (b). If $\sum_{j=0}^{\infty} c_j = \infty$, then for any choice of x_0 , either $\sum_j |x_j| = \infty$ or $\sum_j |x_j| = 0$ by (8.6.16). Hence $w_j = 0$ for all $j \geq 0$, and the chain is null-recurrent or transient.

The preceding example may be modified to reduce the state space to a finite set by letting $p_{c,c+1} = 0$ for some $c \ge 1$. A specific case of this is Example 8 of §8.3, which will now be examined.

Example 16. Let us find the stationary distribution for the Ehrenfest model. We can proceed exactly as in Example 15, leading to the formula (8.6.15), but this time it stops at j = c. Substituting the numerical values from (8.3.16), we obtain

$$c_j = \frac{c(c-1)\cdots(c-j+1)}{1\cdot 2\cdots j} = {c \choose j}, \quad 0 \le j \le c.$$

We have $\sum_{j=0}^{c} c_j = 2^c$ from (3.3.7); hence

$$w_j = \frac{1}{2^c} {c \choose j}, \quad 0 \le j \le c.$$

This is just the binomial distribution B(c, 1/2).

Thus the steady state in Ehrenfest's urn may be simulated by coloring the c balls red or black with probability 1/2 each, and independently of one another; or again by picking them at random from an infinite reservoir of red and black balls in equal proportions.

Next, recalling (8.6.3) and (8.6.7), we see that the mean recurrence times are given by

$$m_{jj} = 2^c {c \choose j}^{-1}, \quad 0 \le j \le c.$$

For the extreme cases j=0 (no black ball) and j=c (no red ball) this is equal to 2^c , which is enormous even for c=100. It follows (see Exercise 42) that the expected time for a complete reversal of the composition of the urn is very long indeed. On the other hand, the chain is recurrent; hence starting, e.g., from an urn containing all black balls, it is almost certain that they will eventually be all replaced by red balls at some time in the Ehrenfest process, and vice versa. Since the number of black balls can change only one at a time the composition of the urn must go through all intermediate "phases" again and again. The model was originally conceived to demonstrate the reversibility of physical processes, but with enormously long cycles for reversal. "If we wait long enough, we shall grow younger again!"

Finally, let us describe without proof a further possible decomposition of a recurrent class. The simplest illustration is that of the classical random walk. In this case the state space of all integers may be divided into two subclasses: the even integers and the odd integers. At one step the particle must move from one subclass to the other, so that the alternation of the two subclasses is a deterministic part of the transition. In general, for each recurrent class C containing at least two states, there exists a unique positive integer d, called the period of the class, with the following properties:

- (a) for every $i \in C$, $p_{ii}^{(n)} = 0$ if $d \nmid n$;* on the other hand, $p_{ii}^{(nd)} > 0$ for all sufficiently large n (how large depending on i);
- (b) for every $i \in C$ and $j \in C$, there exists an integer $r, 1 \le r \le d$, such that $p_{ij}^{(n)} = 0$ if $d \nmid n r$; on the other hand, $p_{ij}^{(nd+r)} > 0$ for all sufficiently large n (how large depending on i and j).

Fixing the state i, we denote by C_r the set of all states j associated with the same number r in (b), for $1 \le r \le d$. These are disjoint sets whose union is C. Then we have the deterministic cyclic transition:

$$C_1 \to C_2 \to \cdots \to C_d \to C_1$$
.

Here is the diagram of such an example with d = 4:

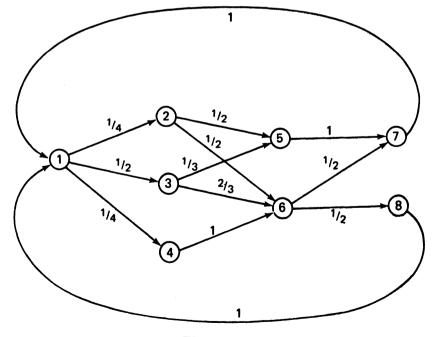


Figure 35

^{* &}quot; $d \nmid n$ " reads "d does not divide n."

where the transition probabilities between the states are indicated by the numbers attached to the directed lines joining them.

The period d of C can be found as follows. Take any $i \in C$ and consider the set of all $n \geq 1$ such that $p_{ii}^{(n)} > 0$. Among the common divisors of this set there is a greatest one: this is equal to d. The fact that this number is the same for all choices of i is part of the property of the period. Incidentally, the decomposition described above holds for any class that is stochastically closed (see §8.7 for definition); thus the free random walk has period 2 whether it is recurrent or transient.

When d=1, the class is said to be aperiodic. A sufficient condition for this is: there exists an integer m such that all elements in Π^m are positive. For then it follows from the Chapman–Kolmogorov equations that the same is true of Π^n for all $n \geq m$, and so property (a) above implies that d=1. In this case the fundamental limit theorem given in (8.6.4) can be sharpened as follows:

$$\lim_{n \to \infty} p_{ij}^{(n)} = \frac{1}{m_{ij}}; \tag{8.6.17}$$

namely the limit of averages may be replaced by a strict individual limit. In general if the period is d, and i and j are as in (b) above, then

$$\lim_{n \to \infty} p_{ij}^{(nd+r)} = \frac{d}{m_{jj}}.$$
 (8.6.18)

We leave it to the reader to show: granted that the limit above exists, its value must be that shown there as a consequence of (8.6.4). Actually (8.6.18) follows easily from the particular case (8.6.17) if we consider d steps at a time in the transition of the chain, so that it stays in a fixed subclass. The sharp result above was first proved by Markov, who considered only a finite state space, and was extended by Kolmogorov in 1936 to the infinite case. Several different proofs are now known; see [Chung 2, $\S I.6$] for one of them.

8.7. Winding up (or down?)

In this section we shall give some idea of the general behavior of a homogeneous Markov chain when there are both recurrent and transient states. Let R denote the set of all recurrent states, T the set of all transient states, so that $I = R \cup T$. We being with a useful definition: a set of states will be called [stochastically] closed iff starting from any state in the set the particle will remain forever in the set. Here and hereafter we shall omit the tedious repetition of the phrase "almost surely" when it is clearly indicated. The salient features of the global motion of the particle may be summarized as follows.

- (i) A recurrent class is closed. Hence, once the particle enters such a class it will stay there forever.
- (ii) A finite set of transient states is not closed. In fact, starting from such a set the particle will eventually move out and stay out of it.
- (iii) If T is finite, then the particle will eventually enter into one of the various recurrent classes.
- (iv) In general the particle will be absorbed into the recurrent classes with total probability α , and remain forever in T with probability $1 - \alpha$, where $0 \le \alpha \le 1$.

Let us prove assertion (i). The particle cannot go from a recurrent state to any transient state by Theorem 6; and it cannot go to any recurrent state in a different class because two states from different classes do not communicate by definition, hence one does not lead to the other by Theorem 8 if these states are recurrent. Therefore from a recurrent class the particle can only move within the class. Next, the truth of assertion (ii) is contained in the proof of Theorem 11, according to which the particle can only spend a finite number of time units in a finite set of transient states. Hence from a certain instant on it will be out of the set. Assertion (iii) is a consequence of (ii) and is illustrated by Example 3 of §8.3 (gambler's ruin problem). Assertion (iv) states an obvious alternative on account of (i), and is illustrated by Example 1 of §8.3 with p > 1/2, in which case $\alpha = 0$; or by Example 9. In the latter case it is clear that starting from $i \geq 1$, either the particle will be absorbed in the state 0 with probability f_{i0}^* in the notation of (8.4.6); or it will move steadily through the infinite set of transient states $\{i+1, i+2, \dots\}$ with probability $1-f_{i0}^*$.

Let us further illustrate some of the possibilities by a simple numerical example.

Example 17. Let the transition matrix be as follows:

The state space may be finite or infinite according to the specification of R_2 , which may be the transition matrix of any recurrent Markov chain such as Example 4 or 8 of §8.3, or Example 1 there with p = 1/2.

Here $T = \{1, 2, 3\}, R_1 = \{4, 5\}$ and R_2 are two distinct recurrent classes. The theory of communication between states implies that the four blocks of 0's in the matrix will be preserved when it is raised to any power. Try to confirm this fact by a few actual schematic multiplications. On the other hand, some of the single 0's will turn positive in the process of multiplication. There are actually two distinct transient classes: $\{1,3\}$ and $\{2\}$; it is possible to go from the first to the second but not vice versa. [This is not important; in fact, a transient class that is not closed is not a very useful entity. It was defined to be a class in $\S 8.4$ only by the force of circumstance!] All three transient states lead to both R_1 and R_2 , but it would be easy to add another that leads to only one of them. The problem of finding the various absorption probabilities can be solved by the general procedure below.

Let $i \in T$ and C be a recurrent class. Put for $n \ge 1$:

$$y_i^{(n)} = \sum_{j \in C} p_{ij}^{(n)} = P_i \{ X_n \in C \}.$$
 (8.7.2)

This is probability that the particle will be in C at time n, given that it starts from i. Since C is closed it will then also be in C at time n+1; thus $y_i^{(n)} \leq y_i^{(n+1)}$ and so by the monotone sequence theorem in calculus the

limit exists as $n \to \infty$:

$$y_i = \lim_{n \to \infty} y_i^{(n)} = P_i \{ X_n \in C \text{ for some } n \ge 1 \}$$

(why the second equation?) and gives the probability of absorption.

Theorem 14. The $\{y_i\}$ above satisfies the system of equations

$$x_i = \sum_{j \in T} p_{ij} x_j + \sum_{j \in C} p_{ij}, \quad i \in T.$$
 (8.7.3)

If T is finite, it is the unique solution of this system. Hence it can be computed by standard method of linear algebra.

Proof: Let the particle start from i, and consider its state j after one step. If $j \in T$, then the Markov property shows that the conditional probability of absorption becomes y_j ; if $j \in C$, then it is already absorbed; if $j \in (I-T)-C$, then it can never be absorbed in C. Taking into account these possibilities, we get

$$y_i = \sum_{j \in T} p_{ij} y_j + \sum_{j \in C} p_{ij} \cdot 1 + \sum_{j \in (I-T)-C} p_{ij} \cdot 0.$$

This proves the first assertion of the theorem. Suppose now T is the finite set $\{1, 2, ..., t\}$. The system (8.7.3) may be written in matrix form as follows:

$$(\Delta_T - \Pi_T)x = y^{(1)}, \tag{8.7.4}$$

where Δ_T is the identity matrix indexed by $T \times T$; Π_T is the restriction of Π on $T \times T$, and $y^{(1)}$ is given in (8.7.2). According to a standard result in linear algebra, the equation above has a unique solution if and only if the matrix $\Delta_T - \Pi_T$ is nonsingular, namely it has an inverse $(\Delta_T - \Pi_T)^{-1}$, and then the solution is given by

$$x = (\Delta_T - \Pi_T)^{-1} y^{(1)}. (8.7.5)$$

Suppose the contrary; then the same result asserts that there is a nonzero solution to the associated homogeneous equation. Namely there is a column vector $v = (v_1, \ldots, v_t) \neq (0, \ldots, 0)$ satisfying

$$(\Delta_T - \Pi_T)v = 0$$
, or $v = \Pi_T v$.

It follows by iteration that

$$v = \Pi_T(\Pi_T v) = \Pi_T^2 v = \Pi_T^2(\Pi_T v) = \Pi_T^3 v = \cdots$$

and so for every $n \geq 1$:

$$v = \Pi_T^n v$$

[cf. (8.6.8) but observe the difference between right-hand and left-hand multiplications]. This means

$$v_i = \sum_{j \in T} p_{ij}^{(n)} v_j, \quad i \in T.$$

Letting $n \to \infty$ and using the corollary to Theorem 5 we see that every term in the sum converges to zero and so $v_i = 0$ for all $i \in T$, contrary to the hypothesis. This contradiction establishes the nonsingularity of $\Delta_T - \Pi_T$ and consequently the existence of a unique solution given by (8.7.5). Since $\{y_i, i \in T\}$ is a solution, the theorem is proved.

For Example 17 above, the equations in (8.7.3) for absorption probabilities into R_1 are

$$x_1 = \frac{1}{8}x_1 + \frac{3}{8}x_2 + \frac{1}{4}x_3 + \frac{1}{4},$$

$$x_2 = \frac{1}{2}x_2 + \frac{1}{3},$$

$$x_3 = \frac{1}{5}x_1 + \frac{3}{10}x_2 + \frac{2}{5}.$$

We get x_2 at once from the second equation, and then x_1 , x_3 from the others:

$$x_1 = \frac{26}{33}, \quad x_2 = \frac{2}{3}, \quad x_3 = \frac{25}{33}.$$

For each i, the absorption probabilities into R_1 and R_2 add up to 1, hence those for R_2 are just $1 - x_1, 1 - x_2, 1 - x_3$. This is the unique solution to another system of equations in which the constant terms above are replaced by 0, 1/6, 1/10. You may wish to verify this as it is a good habit to double-check these things, at least once in a while.

It is instructive to remark that the problem of absorption into recurrent classes can always be reduced to that of absorbing states. For each recurrent class may be merged into a single absorbing state since we are not interested in the transitions within the class; no state in the class leads outside, whereas the probability of entering the class at one step from any transient state i is precisely the $y_i^{(1)}$ used above. Thus, the matrix in (8.7.1)

may be converted to the following one:

$$\begin{bmatrix} \frac{1}{8} & \frac{3}{8} & \frac{1}{4} & \frac{1}{4} & 0 \\ 0 & \frac{1}{2} & 0 & \frac{1}{3} & \frac{1}{6} \\ \frac{1}{5} & \frac{3}{10} & 0 & \frac{2}{5} & \frac{1}{10} \\ 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix}$$

in which the last two states $\{4\}$ and $\{5\}$ take the place of R_1 and R_2 . The absorption probabilities become just f_{i4}^* and f_{i5}^* in the notation of (8.4.6). The two systems of equations remain of course the same.

When T is finite and there are exactly two absorbing states, there is another interesting method. As before let $T=\{1,2,\ldots,t\}$ and let the absorbing states be denoted by 0 and t+1, so that $I=\{0,1,\ldots,t+1\}$. The method depends on the discovery of a positive nonconstant solution of the equation $(\Delta-\Pi)x=0$, namely some such $v=(v_0,v_1,\ldots,v_{t+1})$ satisfying

$$v_i = \sum_{j=0}^{t+1} p_{ij} v_j, \quad i = 0, 1, \dots, t+1.$$
 (8.7.6)

Observe that the two equations for i=0 and i=t+1 are automatically true for any v, because $p_{0j}=\delta_{0j}$ and $p_{t+1,j}=\delta_{t+1,j}$; also that $v_i=1$ is always a solution of the system, but it is constant. Now iteration yields

$$v_i = \sum_{j=0}^{t+1} p_{ij}^{(n)} v_j$$

for all $n \geq 1$; letting $n \to \infty$ and observing that

$$\lim_{n \to \infty} p_{ij}^{(n)} = 0 \quad \text{for } 1 \le j \le t,$$

$$\lim_{n \to \infty} p_{ij}^{(n)} = f_{ij}^* \quad \text{for } j = 0 \quad \text{and} \quad j = t+1,$$

we obtain

$$v_i = f_{i0}^* v_0 + f_{i,t+1}^* v_{t+1}. (8.7.7)$$

Recall also that

$$1 = f_{i0}^* + f_{i,t+1}^*. (8.7.8)$$

We claim that $v_0 \neq v_{t+1}$; otherwise it would follow from the last two equations that $v_i = v_0$ for all i, contrary to the hypothesis that v is nonconstant. Hence we can solve these equations as follows:

$$f_{i0}^* = \frac{v_i - v_{t+1}}{v_0 - v_{t+1}}; \qquad f_{i,t+1}^* = \frac{v_0 - v_i}{v_0 - v_{t+1}}.$$
 (8.7.9)

Example 18. Let us return to Problem 1 of §8.1, where t = c-1. If $p \neq q$, then $v_i = (q/p)^i$ is a nonconstant solution of (8.7.6). This is trivial to verify but you may well demand to know how on earth did we discover such a solution? The answer in this case is easy (but motivated by knowledge of difference equations used in §8.1): try a solution of the form λ^i and see what λ must be. Now if we substitute this v_i into (8.7.9) we get f_{i0}^* equal to the u_i in (8.1.9).

If p = q = 1/2, then $v_i = i$ is a nonconstant solution of (8.7.6) since

$$i = \frac{1}{2}(i+1) + \frac{1}{2}(i-1).$$
 (8.7.10)

This leads to the same answer as given in (8.1.10). The new solution has to do with the idea of a martingale (see Appendix 3). Here is another similar example.

Example 19. The following model of random reproduction was introduced by S. Wright in his genetical studies (see, e.g., [Karlin] for further details). In a *haploid* organism the genes occur singly rather than in pairs as in the diploid case considered in §5.6. Suppose 2N genes of types A and a (the alleles) are selected from each generation. The number of A genes is the state of the Markov chain and the transition probabilities are given below: $I = \{0, 1, \ldots, 2N\}$, and

$$p_{ij} = {2N \choose j} \left(\frac{i}{2N}\right)^i \left(1 - \frac{i}{2N}\right)^{2N-j} \tag{8.7.11}$$

Thus if the number of A genes in any generation is equal to i, then we may suppose that there is an infinite pool of both types of genes in which the proportion of A to a is as i:2N-i, and 2N independent drawings are made from it to give the genes of the next generation. We are therefore dealing with 2N independent Bernoullian trials with success probability i/2N, which results in the binomial distribution B(2N;i/2N) in (8.7.11). It follows that [see (4.4.16) or (6.3.6)] the expected number of A genes is equal to

$$\sum_{j=0}^{2N} p_{ij}j = 2N \frac{i}{2N} = i. (8.7.12)$$

This means that the expected number of A genes in the next generation is equal to the actual (but random) number of these genes in the present generation. In particular, this expected number remains constant through the successive generations. The situation is the same as in the case of a fair game discussed in §8.2 after (8.2.3). The uncertified trick used there is again applicable and in fact leads to exactly the same conclusion except for notation. However, now we can also apply the proven formula (8.7.9), which gives at once

$$f_{i0}^* = \frac{2N - i}{2N}, \qquad f_{i,2N}^* = \frac{i}{2N}.$$

These are the respective probabilities that the population will wind up being pure a-type or A-type.

Our final example deals with a special but important kind of homogeneous Markov chain. Another specific example, queuing process, is outlined with copious hints in Exercises 29–31.

Example 20. A subatomic particle may split into several particles after a nuclear reaction; a male child bearing the family name may have a number of male children or none. These processes may be repeated many times unless extinction occurs. These are examples of a *branching process* defined below.

There is no loss of generality to assume that at the beginning there is exactly one particle: $X_0 = 1$. It gives rise to X_1 descendants of the first generation, where

$$P(X_1 = j) = a_j, \quad j = 0, 1, 2, \dots$$
 (8.7.13)

Unless $X_1 = 0$, each of the particles in the first generation will give rise to descendants of the second generation, whose number follows the same probability distribution given in (8.7.13), and the actions of the various particles are assumed to be stochastically independent. What is the distribution of the number of particles of the second generation? Let the generating function of X_1 be g:

$$g(z) = \sum_{j=0}^{\infty} a_j z^j.$$

Suppose the number of particles in the first generation is equal to j, and we denote the numbers of their descendants by Z_1, \ldots, Z_j , respectively. Then by hypothesis these are independent random variables each having g as its generating function. The total number of particles in the second generation is $X_2 = Z_1 + \cdots + Z_j$ and this has the generating function g^j

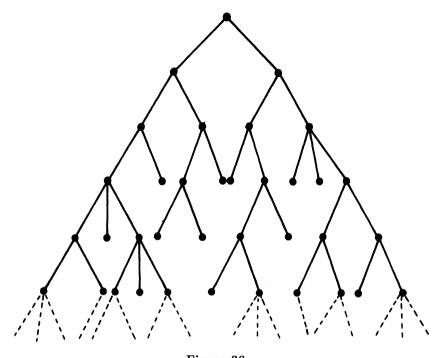


Figure 36

by Theorem 6 of $\S6.5$. Recalling (6.5.12) and the definition of conditional expectation in (5.2.11), this may be written as follows:

$$E(z^{X_2} \mid X_1 = j) = g(z)^j,$$
 (8.7.14)

and consequently by (5.2.12):

$$E(z^{X_2}) = \sum_{j=0}^{\infty} P(X_1 = j) E(z^{X_2} \mid X_1 = j) = \sum_{j=0}^{\infty} a_j g(z)^j = g(g(z)).$$

Let g_n be the generating function of X_n so that $g_1 = g$; then the above says $g_2 = g(g_1)$. Exactly the same argument gives $g_n = g(g_{n-1}) = g \circ g \circ \cdots \circ g$ (there are n appearances of g), where " \circ " denotes the composition of functions. In other words g_n is just the n-fold composition with g with itself. Using this new definition of g_n , we record this as follows:

$$g_n(z) = E(z^{X_n}) = \sum_{k=0}^{\infty} P(X_n = k) z^k.$$
 (8.7.15)

Since the distribution of the number of descendants in each succeeding generation is determined solely by the number in the existing generation, regardless of past evolution, it is clear that the sequence $\{X_n, n \geq 0\}$ has

the Markov property. It is a homogeneous Markov chain because the law of reproduction is the same from generation to generation. In fact, it follows from (8.7.14) that the transition probabilities are given below:

$$p_{jk}$$
 = coefficient of z^k in the power series for $g(z)^j$. (8.7.16)

To exclude trivial cases, let us now suppose that

$$0 < a_0 < a_0 + a_1 < 1. (8.7.17)$$

The state space is then (why?) the set of all nonnegative integers. The preceding hypothesis implies that all states lead to 0 (why?), which is an absorbing state. Hence all states except 0 are transient but there are infinitely many of them. The general behavior under (iii) at the beginning of the section does not apply and only (iv) is our guide. [Observe the term "particle" was used in a different context there.] Indeed, we will now determine the value of α which is called the *probability of extinction* in the present model.

Putting z = 0 in (8.7.15) we see that $g_n(0) = p_{10}^{(n)}$; on the other hand, our general discussion about absorption tells us that

$$\alpha = \lim_{n \to \infty} p_{10}^{(n)} = \lim_{n \to \infty} g_n(0). \tag{8.7.18}$$

Since $g_n(0) = g(g_{n-1}(0))$, by letting $n \to \infty$ we obtain

$$\alpha = g(\alpha). \tag{8.7.19}$$

Thus the desired probability is a root of the equation $\varphi(z) = 0$ where $\varphi(z) = g(z) - z$; we shall call it simply a root of φ . Since g(1) = 1, one root is z = 1. Next we have

$$\varphi''(z) = g''(z) = \sum_{j=2}^{\infty} j(j-1)a_j z^{j-2} > 0$$

for z>0, on account of (8.7.17). Hence the derivative φ' is an increasing function. Now recall *Rolle's theorem* from calculus: between two roots of a differentiable function there is at least one root of its derivative. It follows that φ cannot have more than two roots in [0,1], for then φ' would have more than one root, which is impossible because φ' increases. Thus φ can have at most one root different from 1 in [0,1], and we have two cases to consider.*

^{*}It is customary to draw two pictures to show the two cases below. The reader is invited to do this and see if he or she is more readily convinced than the author.

Case 1. φ has no root in [0,1). Then since $\varphi(0)=a_0>0$, we must have $\varphi(z)>0$ for all z in [0,1), for a continuous function cannot take both positive and negative values in an interval without vanishing somewhere. Thus we have

$$\varphi(1) - \varphi(z) < \varphi(1) = 0, \quad 0 \le z < 1;$$

and it follows that

$$\varphi'(1) = \lim_{z \uparrow 1} \frac{\varphi(1) - \varphi(z)}{1 - z} \le 0;$$

hence $g'(1) \leq 1$.

Case 2. φ has a unique root r in [0,1). Then by Rolle's theorem φ' must have a root s in [r,1), i.e., $\varphi'(s)=g'(s)-1=0$, and since g' is an increasing function we have

$$g'(1) > g'(s) = 1.$$

To sum up: the equation g(z) = z has a positive root less than 1 if and only if g'(1) > 1.

In Case 1, we must have $\alpha = 1$ since $0 \le \alpha \le 1$ and α is a root by (8.7.19). Thus the population is almost certain to become extinct.

In Case 2, we will show that α is the root r < 1. For g(0) < g(r) = r, and supposing for the sake of inducting $g_{n-1}(0) < r$, then $g_n(0) = g(g_{n-1}(0)) < g(r) = r$ because g is an increasing function. Thus $g_n(0) < r$ for all n and so $\alpha \le r$ by (8.7.18). But then α must be equal to r because both of them are roots of the equation in [0, 1).

What will happen in Case 2 if the population escapes extinction? According to the general behavior under (iv) it must then remain forever in the transient states $\{1,2,\ldots\}$ with probability $1-\alpha$. Can its size sway back and forth from small to big and vice versa indefinitely? This question is answered by the general behavior under (ii), according to which it must stay out of every finite set $\{1,2,\ldots,\ell\}$ eventually, no matter how large ℓ is. Therefore it must in fact become infinite (not necessarily monotonically, but as a limit), namely:

$$P\{\lim_{n\to\infty} X_n = +\infty \mid X_n \neq 0 \text{ for all } n\} = 1.$$

The conclusion is thus a "boom or bust" syndrome. The same is true of the gambler who has an advantage over an infinitely rich opponent (see §8.2): if he is not ruined, he will also become infinitely rich. Probability theory contains a lot of such extreme results, some of which are known as zero-or-one ("all or nothing") laws.

In the present case there is some easy evidence for the conclusions reached above. Let us compute the expectation of the population of the nth generation. Let $\mu = E(X_1)$ be the expected number of descendants of each particle. Observe that $\mu = g'(1)$ so that we have $\mu \leq 1$ in Case 1 and $\mu > 1$ in Case 2. Suppose $\mu < \infty$; then if the number of particles in the n-1st generation is j, the expected number of particles in the nth generation will be $j\mu$ (why?). Using conditional expectation, this may be written as

$$E\{X_n \mid X_{n-1} = j\} = \mu j.$$

It follows from (5.2.12) that

$$E(X_n) = \sum_{j=0}^{\infty} \mu j P(X_{n-1} = j) = \mu E(X_{n-1}),$$

and consequently by iteration

$$E(X_n) = \mu^n E(X_0) = \mu^n.$$

Therefore we have

$$\lim_{n \to \infty} E(X_n) = \lim_{n \to \infty} \mu^n = \begin{cases} 0 & \text{if } \mu < 1, \\ 1 & \text{if } \mu = 1, \\ \infty & \text{if } \mu > 1. \end{cases}$$

This tends to support our conclusion in Case 1 for certain extinction; in fact it is intuitively obvious that if $\mu < 1$ the population fails to be self-replacing on the basis of averages. The case $\mu = 1$ may be disposed of with a bit more insight, but let us observe that here we have the strange situation that $E(X_n) = 1$ for all n, but $P(\lim_{n \to \infty} X_n = 0) = 1$ by Case 1. In case $\mu > 1$ the crude interpretation would be that the population will certainly become infinite. But we have proved under Case 2 that there is a definite probability that it will die out as a dire contrast. This too is interesting in relating simple calculations to more sophisticated theory. These comments are offered as an invitation to the reader for further wonderment about probability and its meaning.

Exercises

- 1. Let X_n be as in (8.1.2) with $X_0 = 0$. Find the following probabilities:
 - (a) $P\{X_n \ge 0 \text{ for } n = 1, 2, 3, 4\};$
 - (b) $P\{X_n \neq 0 \text{ for } n = 1, 2, 3, 4\};$
 - (c) $P\{X_n \le 2 \text{ for } n = 1, 2, 3, 4\};$

- (d) $P\{|X_n| \le 2 \text{ for } n = 1, 2, 3, 4\}.$
- 2. Let $Y_n = X_{2n}$, where X_n is as in No. 1. Show that $\{Y_n, n \geq 0\}$ is a Markov chain and find its transition matrix. Similarly for $\{Z_n, n \geq 0\}$ where $Z_n = X_{2n+1}$; what is its initial distribution?
- 3. Let a coin be tossed indefinitely; let H_n and T_n respectively denote the numbers of heads and tails obtained in the first n tosses. Put $X_n = H_n, Y_n = H_n T_n$. Are these Markov chains? If so, find the transition matrix.
- *4. As in No. 3 let $Z_n = |H_n T_n|$. Is this a Markov chain? [Hint: compute, e.g., $P\{Y_{2n} = 2i \mid Z_{2n} = 2i\}$ by Bernoulli's formula, then $P\{Z_{2n+1} = 2i \pm 1 \mid Z_{2n} = 2i, Y_{2n} > 0\}$.]
- 5. Let the transition matrix be given below:

(a)
$$\begin{pmatrix} \frac{1}{2} & \frac{1}{2} \\ \frac{1}{3} & \frac{2}{3} \end{pmatrix}$$
 (b) $\begin{pmatrix} p_1 & q_1 & 0 \\ 0 & p_2 & q_2 \\ q_3 & 0 & p_3 \end{pmatrix}$.

Find $f_{11}^{(n)}$, $f_{12}^{(n)}$, $g_{12}^{(n)}$ for n = 1, 2, 3 [for notation see (8.4.6) and (8.4.10)].

- 6. In a model for the *learning process* of a rat devised by Estes, the rodent is said to be in state 1 if he has learned a certain trick (to get a peanut or avoid an electric shock), and to be in state 2 if he has not yet learned it. Suppose that once it becomes learned it will remain so, while if it is not yet learned it has a probability α of becoming so after each trial run. Write down the transition matrix and compute $p_{21}^{(n)}, f_{21}^{(n)}$ for all $n \geq 1$; and m_{21} [see (8.6.3) for notation].
- 7. Convince yourself that it is a trivial matter to construct a transition matrix in which there are any given number of transient and recurrent classes, each containing a given number of states, provided that either (a) I is infinite, or (b) I is finite but not all states are transient.
- 8. Given any transition matrix Π , show that it is trivial to enlarge it by adding new states that lead to old ones, but it is impossible to add any new state that communicates with any old one.
- 9. In the "double or nothing" game, you bet all you have and you have a fifty-fifty chance to double it or lose it. Suppose you begin with \$1 and decide to play this game up to n times (you may have to quit sooner because you are broke). Describe the Markov chain involved with its transition matrix.
- 10. Leo is talked into playing heads in a coin-tossing game in which the probability of heads is only 0.48. He decides that he will quit as soon as he is one ahead. What is the probability that he may never quit?
- 11. A man has two girlfriends, one uptown and one downtown. When he wants to visit one of them for a weekend he chooses the uptown girl

with probability p. Between two visits he stays home for a weekend. Describe the Markov chain with three states for his weekend whereabouts: "uptown," "home," and "downtown." Find the long-run frequencies of each. This is the simplest case of Example 4 of §8.3, but here is a nice puzzle related to the scheme. Suppose that the man decides to let chance make his choice by going to the bus stop where buses go both uptown and downtown and jumping aboard the first bus that comes. Since he knows that buses run in both directions every 15 minutes, he figures that these equal frequencies must imply p = 1/2 above. But after a while he realizes that he has been visiting uptown twice as frequently as downtown. How can this happen? This example carries an important lesson to the practicing statistician, namely that the relevant datum may not be what appears at first right. Assume that the man arrives at the bus stop at random between 6 p.m. and 8 p.m. Figure out the precise bus schedules that will make him board the uptown buses with probability p = 2/3.

- 12. Solve Problem 1 of $\S 8.1$ when there is a positive probability r of the particle remaining in its position at each step.
- *13. Solve (8.1.13) when $p \neq q$ as follows. First determine the two values λ_1 and λ_2 such that $x_j = \lambda^j$ is a solution of $x_j = px_{j+1} + qx_{j-1}$. The general solution of this system is then given by $A\lambda_j^1 + B\lambda_j^2$, where A and B are constants. Next find a particular solution of $x_j = px_{j+1} + qx_{j-1} + 1$ by trying $x_j = Cj$ and determine the constant C. The general solution of the latter system is then given by $A\lambda_j^1 + B\lambda_j^2 + Cj$. Finally, determine A and B from the boundary conditions in (8.1.13).
 - 14. The original Ehrenfest model is as follows. There are N balls in each of two urns. A ball is chosen at random from the 2N balls from either urn and put into the other urn. Let X_n denote the number of balls in a fixed urn after n drawings. Show that this is a Markov chain having the transition probabilities given in (8.3.16) with c = 2N.
 - 15. A scheme similar to that in No. 14 was used by Daniel Bernoulli [son of Johann, who was younger brother of Jakob] and Laplace to study the flow of incompressible liquids between two containers. There are N red and N black balls in two urns containing N balls each. A ball is chosen at random from each urn and put into the other. Find the transition probabilities for the number of red balls in a specified urn.
 - 16. In certain ventures such as doing homework problems one success tends to reinforce the chance for another by imparting experience and confidence; in other ventures the opposite may be true. Anyway let us assume that the aftereffect is carried over only two consecutive trials so that the resulting sequence of successes and failures constitutes a Markov chain on two states $\{s, f\}$. Let

$$p_{ss} = \alpha, \quad p_{ff} = \beta,$$

where α and β are two arbitrary members between 0 and 1. Find the long-run frequency of successes.

- 17. The following model has been used for the study of contagion. Suppose that there are N persons, some of whom are sick with influenza. The following assumptions are made:
 - (a) when a sick person meets a healthy one, the chance is α that the latter will be infected;
 - (b) all encounters are between two persons;
 - (c) all possible encounters in pairs are equally likely;
 - (d) one such encounter occurs in every (chosen) unit of time.

Define a Markov chain for the spread of the disease and write down its transition matrix. [Are you overwhelmed by all these oversimplifying assumptions? Applied mathematics is built upon the shrewd selection and exploitation of such simplified models.]

- 18. The age of a light bulb is measured in days, and fractions of a day do not count. If a bulb is burned out during the day, then it is replaced by a new one at the beginning of the next day. Assume that a bulb that is alive at the beginning of the day, possibly one that has just been installed, has probability p of surviving at least one day so that its age will be increased by 1. Assume also that the successive bulbs used lead independent lives. Let $X_0 = 0$ and X_n denote the age of the bulb that is being used at the beginning of the n + 1st day. (We begin with the first day, thus $X_1 = 1$ or 0 depending on whether the initial bulb is still in place or not at the beginning of the second day.) The process $\{X_n, n \geq 0\}$ is an example of a renewal process. Show that it is a recurrent Markov chain, find its transition probabilities and stationary distribution. [Note: the life span of a bulb being essentially a continuous variable, a lot of words are needed to describe the scheme accurately in discrete time, and certain ambiguities must be resolved by common sense. It would be simpler and clearer to formulate the problem in terms of heads and tails in coin tossing (how?), but then it would have lost the flavor of application!
- 19. Find the stationary distribution for the random walk with two reflecting barriers (Example 4 of §8.3).
- 20. In a sociological study of "conformity" by B. Cohen, the following Markov chain model was used. There are four states: $S_1 = \text{consistently}$ nonconforming; $S_2 = \text{indecisively nonconforming}$; $S_3 = \text{indecisively}$ conforming; $S_4 = \text{consistently conforming}$. In a group experiment subjects were found to switch states after each session according to the

following transition matrix:

	S_1	S_2	S_3	S_4
$\overline{S_1}$	1	0	0	0
S_2	.06	.76	.18	0
S_3	0	.27	.69	.04
S_4	0	0	0	1

Find the probabilities of ultimate conversion from the "conflict" states S_2 and S_3 into the "resolution" states S_1 and S_4 .

21. In a genetical model similar to Example 19 of §8.7, we have $I = \{0, 1, \dots, 2N\}$ and

$$p_{ij} = \binom{2i}{j} \binom{2N-2i}{N-j} / \binom{2N}{N}.$$

How would you describe the change of genotypes from one generation to another by some urn scheme? Find the absorption probabilities. [Hint: compute $\sum_{j=0}^{2N} j p_{ij}$ by simplifying the binomial coefficients, or by Theorem 1 of §6.1.]

- 22. For the branching process in Example 20 of §8.7, if a_0 , a_1 , and a_2 are positive but the other a_j 's are all zero, find the probability of extinction.
- 23. Suppose that the particles in the first generation of a branching process follow a probability law of splitting given by $\{b_j, j \geq 0\}$ that may be different from that of initial particle given by (8.7.13). What then is the distribution of the number of particles in the second generation?
- 24. A sequence of electric impulses is measured by a meter that records the highest voltage that has passed through it up to any given time. Suppose that the impulses are uniformly distributed over the range $\{1, 2, ..., \ell\}$. Define the associated Markov chain and find its transition matrix. What is the expected time until the meter records the maximum value ℓ ? [Hint: argue as in (8.1.13) for the expected absorption time into the state ℓ ; use induction after computing e_{l-1} and e_{l-2} .]
- 25. In proofreading a manuscript each reader finds at least one error. But if there are j errors when she begins, she will leave it with any number of errors between 0 and j-1 with equal probabilities. Find the expected number of readers needed to discover all the errors. [Hint: $e_j = j^{-1}(e_1 + \cdots + e_{j-1}) + 1$; now simplify $e_j e_{j-1}$.]
- 26. A deck of m cards may be shuffled in various ways. Let the state space be the m! different orderings of the cards. Each particular mode of shuffling sends any state (ordering) into another. If the various modes are randomized, this results in various transition probabilities between the states. Following my tip (a) in §3.4 for combinatorial problems, let us begin with m=3 and the following two modes of shuffling:

- (i) move the top card to the bottom, with probability p;
- (ii) interchange the top and middle cards, with probability 1 p.

Write down the transition matrix. Show that it is doubly stochastic and all states communicate. Show that if either mode alone is used the states will not all communicate.

- 27. Change the point of view in No. 26 by fixing our attention on a particular card, say the queen of spades if the three cards are the king, queen, and knight of spades. Let X_n denote its position after n shufflings. Show that this also constitutes a Markov chain with a doubly stochastic transition matrix.
- *28. Now generalize Nos. 26 and 27: for any m and any randomized shuffling, the transition matrices in both formulations are doubly stochastic. [Hint: each mode of shuffling as a permutation on m cards has an inverse. Thus if it sends the ordering j into k then it sends some ordering i into j. For fixed j the correspondence i = i(k) is one-to-one and $p_{ij} = p_{jk}$. This proves the result for the general case of No. 26. Next consider two orderings j_1 and j_2 with the fixed card in the topmost position, say. Each mode of shuffling that sends j_1 into an ordering with the given card second from the top does the same to j_2 . Hence the sum of probabilities of such modes is the same for j_1 or j_2 , and gives the transition probability $1 \to 2$ for the displacement of the card in question.]
- *29. Customers arrive singly at a counter and enter into a queue if it is occupied. As soon as one customer finishes, the service for the next customer begins if there is anyone in the queue, or upon the arrival of the next customer if there is no queue. Assume that the service time is constant (e.g., a taped recording or automatic hand-dryer), then this constant may be taken as the unit of time. Assume that the arrivals follow a Poisson process with parameter α in this unit. For $n \geq 1$ let X_n denote the number of customers in the queue at the instant when the nth customer finishes his service. Let $\{Y_n, n \geq 1\}$ be independent random variables with the Poisson distribution $\pi(\alpha)$; see §7.1. Show that

$$X_{n+1} = (X_n - 1)^+ + Y_n, \quad n \ge 1,$$

where $x^+ = x$ if x > 0 and $x^+ = 0$ if $x \le 0$. Hence conclude that $\{X_n, n \ge 1\}$ is a Markov chain on $\{0, 1, 2, ...\}$ with the following transition matrix:

$$\begin{bmatrix} c_0 & c_1 & c_2 & c_3 & \cdots \\ c_0 & c_1 & c_2 & c_3 & \cdots \\ 0 & c_0 & c_1 & c_2 & \cdots \\ 0 & 0 & c_0 & c_1 & \cdots \\ \cdots & \cdots & \cdots & \cdots \end{bmatrix}$$

where $c_j = \pi_j(\alpha)$. [Hint: this is called a queuing process and $\{X_n, n \geq 1\}$ is an imbedded Markov chain. At the time when the nth customer finishes there are two possibilities. (i) The queue is not empty; then the n+1st customer begins his service at once and during his (unit) service time Y_n customers arrive. Hence when he finishes the number in the queue is equal to X_n-1+Y_n . (ii) The queue is empty; then the counter is free and the queue remains empty until the arrival of the n+1st customer. He begins service at once and during his service time Y_n customers arrive. Hence when he finishes the number in the queue is equal to Y_n . The Y_n 's are independent and have $\pi(\alpha)$ as distribution, by Theorems 1 and 2 of §7.2.]

*30. Generalize the scheme in No. 29 as follows. The service time is a random variable S such that $P\{S=k\}=b_k, k\geq 1$. Successive service times are independent and identically distributed. Show that the conclusions of No. 29 hold with

$$c_j = \sum_{k=1}^{\infty} b_k \pi_j(k\alpha).$$

*31. In No. 29 or 30, let $\mu = \sum_{j=0}^{\infty} jc_j$. Prove that the Markov chain is transient, null-recurrent, or positive-recurrent depending on whether $\mu < 1$, $\mu = 1$ or $\mu > 1$. [This result is due to Lindley; here are the steps for a proof within the scope of Chapter 8. In the notation of §8.4 let

$$F_{10}(z) = f(z), \quad g(z) = \sum_{j=0}^{\infty} c_j z^j.$$

- (a) $F_{j,j-1}(z) = f(z)$ for all $j \ge 1$; because, e.g., $f_{j,j-1}^{(4)} = P\{Y_n \ge 1, Y_n + Y_{n+1} \ge 2, Y_n + Y_{n+1} + Y_{n+2} \ge 3, Y_n + Y_{n+1} + Y_{n+2} + Y_{n+3} = 3 \mid X_n = j\}$.
- (b) $F_{j0}(z) = f(z)^i$ for $j \ge 1$, because the queue size can decrease only by one at a step.
- (c) $f_{10}^{(1)} = c_0, f_{10}^{(v)} = \sum_{j=1}^{\infty} c_j f_{j0}^{(v-1)}$ for $v \ge 2$; hence

$$f(z) = c_0 z + \sum_{j=1}^{\infty} c_j z F_{j0}(z) = zg(f(z)).$$

- (d) $F_{00}(z) = zg(f(z))$ by the same token.
- (e) If $f(1) = \rho$, then ρ is the smallest root of the equation $\rho = g(\rho)$ in [0,1]; hence $F_{00}(1) = f(1) < 1$ or = 1 according as g'(1) > 1 or ≤ 1 by Example 4 of §8.7.

(f) f'(1) = f'(1)g'(1) + g(1); hence if $g'(1) \le 1$ then in the notation of (8.6.3), $m_{00} = F'_{00}(1) = f'(1) = \infty$ or $< \infty$ depending on whether g'(1) = 1 or < 1. Q.E.D.

For more complicated queuing models see e.g., [Karlin].]

*32. A company desires to operate s identical machines. These machines are subject to failure according to a given probability law. To replace these failed machines the company orders new machines at the beginning of each week to make up the total s. It takes one week for each new order to be delivered. Let X_n be the number of machines in working order at the beginning of the nth week and let Y_n denote the number of machines that fail during the nth week. Establish the recursive formula

$$X_{n+1} = s - Y_n$$

and show that $\{X_n, n \geq 1\}$ constitutes a Markov chain. Suppose that the failure law is uniform, i.e.:

$$P{Y_n = j \mid X_n = i} = \frac{1}{i+1}, \quad j = 0, 1, \dots, i.$$

Find the transition matrix of the chain, its stationary distribution, and the expected number of machines in operation in the steady state.

*33. In No. 32 suppose the failure law is binomial:

$$P{Y_n = j \mid X_n = i} = {i \choose j} p^j (1-p)^{i-j}, \quad j = 0, 1, \dots, i,$$

with some probability p. Answer the same questions as before. [These two problems about $machine\ replacement$ are due to D. Iglehart.]

- 34. The matrix $[p_{ij}], i \in I, j \in I$ is called *substochastic* iff for every i we have $\sum_{j \in I} p_{ij} \leq 1$. Show that every power of such a matrix is also substochastic.
- 35. Show that the set of states C is stochastically closed if and only if for every $i \in C$ we have $\sum_{j \in C} p_{ij} = 1$.
- 36. Show that

$$\max_{0 \le n < \infty} P_i \{ X_n = j \} \le P_i \left\{ \bigcup_{n=0}^{\infty} [X_n = f] \right\} \le \sum_{n=0}^{\infty} P_i \{ X_n = j \}.$$

Hence deduce that $i \rightsquigarrow j$ if and only if $f_{ij}^* > 0$.

- 37. Prove that if $q_{ij} > 0$, then $\sum_{n=0}^{\infty} p_{ij}^{(n)} = \infty$.
- *38. Prove that if $j \rightsquigarrow i$, then $g_{ij}^* < \infty$. Give an example where $g_{ij}^* = \infty$. [Hint: show that $g_{ij}^{(n)} f_{ji}^{(v)} \leq f_{ii}^{(n+v)}$ and choose v so that $f_{ji}^{(v)} > 0$.]

- 39. Prove that if there exists j such that $i \rightsquigarrow j$ but not $j \rightsquigarrow i$, then i is transient. [Hint: use Theorem 9; or argue as in the proof of Theorem 9 to get $q_{ii} \leq p_{ij}^{(n)} \cdot 0 + (1 p_{ij}^{(n)}) \cdot 1$ for every n.]
- 40. Define for arbitrary i, j and k in I and $n \ge 1$:

$$_{k}p_{ij}^{(n)} = P_{i}\{X_{v} \neq k \text{ for } 1 \leq v \leq n-1; X_{n} = j\}.$$

Show that if k = j this reduces to $f_{ij}^{(n)}$, while if k = i it reduces to $g_{ij}^{(n)}$. In general, prove that

$$\sum_{\ell \neq k} {}_{k} p_{i\ell}^{(n)} {}_{k} p_{\ell j}^{(m)} = {}_{k} p_{ij}^{(n+m)}.$$

These are called $taboo\ probabilities$ because the passage through k during the transition is taboo.

- 41. If the total number of states is r, and $i \leadsto j$, then there exists n such that $1 \le n \le r$ and $p_{ij}^{(n)} > 0$. [Hint: any sequence of states leading from i to j in which some k occurs twice can be shortened.]
- 42. Generalize the definition in (8.6.3) as follows:

$$m_{ij} = E_i(T_j) = \sum_{v=1}^{\infty} v f_{ij}^{(v)}.$$

Prove that $m_{ij} + m_{ji} \ge m_{ii}$ for any two states i and j. In particular, in Example 16 of §8.6, we have $m_{0c} \ge 2^{c-1}$.

- 43. Prove that the symmetric random walk is null-recurrent. [Hint: $P_{ij}^{(n)} = P\{\xi_1 + \dots + \xi_n = j i\}$; use (7.3.7) and the estimate following it.]
- 44. For any state *i* define the *holding time* in *i* as follows: $S = \max\{n \ge 1 \mid X_v = i, \text{ for all } v = 1, 2, \dots, n\}$. Find the distribution of *S*.
- *45. Given the Markov chain $\{X_n, n \geq 1\}$ in which there is no absorbing state, define a new process as follows. Let n_1 be the smallest value of n such that $X_n \neq X_1$, n_2 the smallest value $> n_1$ such that $X_n \neq X_{n_1}$, n_3 the smallest value $> n_2$ such that $X_n \neq X_{n_2}$ and so on. Now put $Y_v = X_{nv}$; show that $\{Y_v, v \geq 1\}$ is also a Markov chain and derive its transition matrix from that of $\{X_n, n \geq 1\}$. Prove that if a state is recurrent in one of them, then it is also recurrent in the other.
- 46. In the notation of No. 3, put $H_n^{(2)} = \sum_{v=1}^n H_v$. Show that $\{X_n\}$ does not form a Markov chain but if we define a process whose value Y_n at time n is given by the ordered pair of states (X_{n-1}, X_n) , then $\{Y_n, n \geq 1\}$ is a Markov chain. What are its state space and transition matrix? The process $\{H_n^{(2)}, n \geq 0\}$ is sometimes called a *Markov chain of order* 2. How would you generalize this notion to a higher order?

47. There is a companion to the Markov property that shows it in reverse time. Let $\{X_n\}$ be a homogeneous Markov chain. For $n \geq 1$ let B be any event determined by X_{n+1}, X_{n+2}, \ldots Show that we have for any two states i and j:

$$P\{X_{n-1} = j \mid X_n = i; B\} = P\{X_{n-1} = j \mid X_n = i\};$$

but this probability may depend on n. However, if $\{X_n\}$ is stationary as in Theorem 13, show that the probability above is equal to

$$\tilde{p}_{ij} = \frac{w_j p_{ji}}{w_i}$$

and so does not depend on n. Verify that $[\tilde{p}_{ij}]$ is a transition matrix. A homogeneous Markov chain with this transition matrix is said to be a *reverse chain* relative to the original one.

48. Prove that $E(S_j) < \infty$ for each j, thus strengthening Theorem 1. [Hint: starting from any j in [1, c-1], at most c-1 steps get us out of the interval. Let W denote the waiting time, then $P(W \ge c-1) \le 1 - p^{c-1} = \delta$ say. Repeating the argument (how?) we obtain for any $n \ge 1$, $P(W \ge n(c-1)) \le \delta^n$. Put $V = W(c-1)^{-1}$, then $P(V = n) \le \delta^n$ and so $E(V) \le \sum_{n=1}^{\infty} n\delta^n < \infty$. Tricky? Mathematics can be.]

Appendix 3

Martingale

Let each X_n be a random variable having a finite expectation, and for simplicity we will suppose it to take integer values. Recall the definition of conditional expectation from the end of §5.2. Suppose that for every event A determined by X_0, \ldots, X_{n-1} alone, and for each possible value i of X_n , we have

$$E\{X_{n+1} \mid A; X_n = i\} = i; (A.3.1)$$

then the process $\{X_n, n \geq 0\}$ is called a *martingale*. This definition resembles that of a Markov chain given in (8.3.1) in the form of the conditioning, but the equation is a new kind of hypothesis. It is more suggestively exhibited in the symbolic form below:

$$E\{X_{n+1} \mid X_0, X_1, \dots, X_n\} = X_n. \tag{A.3.2}$$

This means: for arbitrary given values of X_0, X_1, \ldots, X_n , the conditional expectation of X_{n+1} is equal to the value of X_n , regardless of the other values. The situation is illustrated by the symmetric random walk or the genetical model in Example 19 of §8.7. In the former case, if the present position of the particle is X_n , then its position after one step will be X_n+1 or X_n-1 with probability 1/2 each. Hence we have, whatever the value of X_n ,

$$E\{X_{n+1} \mid X_n\} = \frac{1}{2}(X_n + 1) + \frac{1}{2}(X_n - 1) = X_n;$$

furthermore this relation remains true when we add to the conditioning the previous positions of the particle represented by $X_0, X_1, \ldots, X_{n-1}$. Thus

the defining condition (A.3.1) for a martingale is satisfied. In terms of the gambler, it means that if the game is fair then at each stage his expected gain or loss cancels out so that his expected future worth is exactly equal to his present assets. A similar assertion holds true of the number of A genes in the genetical model. More generally, when the condition (8.7.6) is satisfied, then the process $\{v(X_n) \ n \geq 0\}$ constitutes a martingale, where v is the function $i \to v(i), i \in I$. Finally in Example 20 of §8.7, it is easy to verify that the normalized population size $\{X_n/\mu^n, n \geq 0\}$ is a martingale.

If we take A to be an event with probability 1 in (A.3.1), and use (5.2.12), we obtain

$$E(X_{n+1}) = \sum_{i} P(X_n = i) E(X_{n+1} \mid X_n = i)$$

$$= \sum_{i} P(X_n = i) i = E(X_n).$$
(A.3.3)

Hence in a martingale all the random variables have the same expectation. This is observed in (8.2.3), but the fact by itself is not significant. The following result from the theory of martingales covers the applications mentioned there and in §8.7. Recall the definition of an optional random variable from §8.5.

Theorem. If the martingale is bounded, namely if there exists a constant M such that $|X_n| \leq M$ for all n, then for any optional T we have

$$E(X_T) = E(X_0).$$
 (A.3.4)

For any martingale, this equation holds if T is bounded.

In the case of Problem 1 of §8.1 with p = 1/2, we have $|X_n| \le c$; in the case of Example 3 of §8.3 we have $|X_n| \le 2N$. Hence the theorem is applicable and the absorption probabilities fall out from it as shown in §8.2.

The extension of (A.3.3) to (A.3.4) may be false for a martingale and an optional T, without some supplementary condition such as boundedness. In this respect, the theorem above differs from the strong Markov property discussed in §8.5. Here is a trivial but telling example for the failure of (A.3.4). Let the particle start from 0 and let T be the first entrance time into 1. Then T is finite by Theorem 2 of §8.2; hence X_T is well defined and must equal 1 by its definition. Thus $E(X_T) = 1$ but $E(X_0) = 0$.

Martingale theory was largely developed by J.L. Doob (1910–) and has become an important chapter of modern probability theory; for an introduction see [Chung 1, Chapter 9].

We conclude this brief introduction with a wonderful exhibit.

Borel's St. Petersburg Martingale. Let $\{y_n, n \in N\}$ be independent random variables with $P\{y_n = +1\} = P\{y_n = -1\} = 1/2$, namely the fair

coin-tossing sequence. Define a sequence of numbers as follows:

$$b_1 = 2;$$
 for $n \ge 2$, $b_n = \sum_{j=1}^{n-1} b_j + 2^n$.

Actually we can verify that $b_n = (n+1).2^{n-1}$. Now define

$$X_n = \sum_{j=1}^n b_j y_j.$$

Then $\{X_n, n \in N\}$ is a martingale. Indeed, successive sums of independent random variables with mean zero form a martingale (Exercise). Next define the random variable

$$T(w) = \min\{n \in N : y_n(w) = +1\},\$$

namely T is the first time the coin comes up "heads." Then T is optional with $P\{T=n\}=2^{-n}$ and E(T)=2. It follows that T is almost surely finite. When $T<\infty$ we have $X_T=2^T$ because when T=n, we have

$$X_n = -\sum_{j=1}^{n-1} b_j + b_n = 2^n.$$

Thus $E(X_T) = \sum_{n=1}^{\infty} 2^{-n} \cdot 2^n = \infty$; whereas $E(X_n) = 0$ for all n. In sum, the game is fair and yet the gambler has a sure win (almost sure to be exact!) of infinite expectation. This is the St. Petersburg Paradox discussed in Exercise 28 of Chapter 4, over which numerous ancient and modern mathematicians have wracked their brains.

Mathematically, the "fair" equation (A.3.4) (read X_1 for X_0) is false. The theorem above does not apply because neither the martingale nor T is bounded. Now for any positive integer t let $T \wedge t$ denote the minimum of T and t. Borel showed

$$E(X_{T \wedge t}) = 0 \qquad \text{for any } t \ge 1 \tag{A.3.5}$$

by a direct, explicit computation. This is not hard and is an excellent exercise, but it is a particular case of the second assertion in the Theorem above.

Equation (A.3.5) restores the fairness of Borel's reincarnation of the St. Petersburg game. In the real world, of course, "Time must have a stop." Historically the idea of limiting the duration of the game, or equivalently the liability of the gamblers, had been variously suggested as realistic caveats. But not until Borel's martingaling and bounded stopping (1938 to 1949) was the denouement of the paradox presented in a simple mathematical proposition, which turns out to be valid for all martingales.

To prove the Theorem we begin with the notation $E(A; X) = E(I_A.X)$ for a set A and random variable X. Then the martingale property (A.3.3) (omitting X_0) means: for any set determined by X_1, \ldots, X_n alone we have

$$E(A; X_{n+1}) = E(A; X_n).$$
 (A.3.6)

Since T is optional, $\{T \leq n\}$ is such a set, so is its complement $\{T > n\}$. Now

$$E(X_{T \wedge (n+1)}) = E(T \leq n; X_T) + E(T > n; X_{n+1}),$$

$$E(X_{T \wedge n}) = E(T \leq n; X_T) + E(T > n; X_n);$$

hence by using (A.3.6) with $A = \{T > n\}$ we see the equality of the two expectations displayed. Thus all $E(X_{T \wedge n})$ are equal for all $n \in N$, so equal to $E(X_1)$ since $T \wedge 1 = 1$. The second assertion in the theorem is proved because if T is bounded then it is identical with $T \wedge n$ for some n. To prove the first assertion, let n increase to infinity, then $T \wedge n$ becomes T in the limit if T is finite by assumption, and so $X_{T \wedge n}$ becomes X_T . Since $X_{T \wedge n}$ for all n is bounded by assumption, $E(X_{T \wedge n})$ converges to $E(X_T)$ by Lebesgue's bounded convergence theorem [Chung 1, §3.2]. This is a bit of advanced analysis appropriate for the occasion. Here is the instructive direct proof:

$$E(X_{T \wedge t}) = \sum_{n=1}^{t} E(T = n; X_T) + E(T > t; X_t).$$

The last-written expectation is bounded by E(T > t; M) = P(T > t).M, which converges to zero as t goes to infinity since T is finite. Hence

$$\lim_{t \to \infty} E(X_{T \wedge t}) = \sum_{n=1}^{\infty} E(T = n; X_T) = E(X_T). \text{ Q.E.D.}$$

Mean-Variance Pricing Model

In this and the next chapter we will illustrate the applications of probability concepts to the field of mathematical finance, starting with the concept of sample space, all the way to stochastic processes including martingales. The present chapter examplifies "one-period" models, where the analysis is based on random variables evaluated at a specific time. The next chapter will address time-dependent ("multiperiod") models, with analysis based on stochastic processes. The two chapters differ in their financial perspectives: the first is basic to the concept of equilibrium pricing, the next describes the application of the arbitrage-free pricing argument.

9.1. An investments primer

The field of finance is replete with vocabulary reflecting both its rich conceptual and practical aspects. Numerous introductory books have been written on the subject. We list here only two that appear to give comprehensive and understandable descriptions of the practice of finance: Investments by Sharpe et al., and Investments by Bodie et al. The book Capital Ideas by Bernstein is a lively account of the paths followed by individuals who shaped the field of finance as a quantitative discipline and as a distinct branch of economics. In this section we give a short and basic introduction to the financial concepts we will be using in this book.

When you invest you either buy, sell, borrow, or lend financial instruments such as stock and cash. The theory and practice of finance involve several concepts that sometimes differ only in a subtle way. To simplify matters, though, we will use certain words interchangeably. In particular,

we refer to financial instruments also as assets or securities. We will use either wealth or fortune when referring to the total cash value of an investor's assets.

At their most fundamental level the financial instruments we deal with here are of either equity-type or debt-type. Equity-type instruments represent ownership of companies. In this chapter, and for simplicity, we will consider publicly traded stocks, on stock markets such as the New York Stock Exchange, as the only equity-type securities. Many of the very large and well-established companies pay dividends to their stock holders. These are payments made, generally quarterly, sometimes less frequently, by the companies. Debt-type instruments represent cash loans. Investors lend money to corporations or to the government (both federal and local) by purchasing bonds. As with any loan, an interest payment schedule (so-called coupon payment) is specified as well as a date (so-called maturity date) by which the loan (also called *principal* or *par-value*) is to be repaid. For example, an investor who purchases a 30-year U.S. government bond at \$10,000 for which he or she receives a coupon payment schedule of \$800 a year is lending the U.S. government \$10,000, to be repaid in 30 years, with an interest rate of 8% per year. For certain types of bonds the investor does not receive any coupon payment. These are called "zero-coupon" or "discount" bonds. In effect, with these bonds, investors lend the money that they receive back with interest at the maturity date. Another example is where you buy a certificate of deposit ("CD") at your local bank. You can buy a CD for any amount you wish, subject to certain minimum imposed by your bank, say \$5000. You will earn interest on this amount, which you technically loaned to your bank. A CD comes with a term or duration, which is the amount of time you have to wait before your bank pays you back, usually 6 or 12 months. For simplicity, we will consider only one type of debt-type asset: a money market instrument, which represents a pool of several loans with possibly different (short) maturities and interest rates. If we invest in it, we get paid a fixed interest rate.

The basic financial instruments of stocks and bonds are traded and priced according to the laws of supply and demand, as well as expectations about certain economic factors such as inflation. They are the focus of this chapter. The other category of financial instruments we consider in this book are the so-called financial derivatives. Their values depend directly on those of the basic instruments of stocks and bonds. An American put option on a stock is an example of an equity-type derivative. It is a contract between two parties giving one the right to sell to the other, by an agreed-upon date, a specific stock at an agreed upon price. Derivatives are the subject of the next chapter.

9.2. Asset return and risk

Suppose today you hold an amount x_0 dollars in an asset, say a company stock. A year from now the value of this asset, i.e., the price of the company stock, is likely to be different from its value today. You need only look at the financial pages of the newspapers, or the Internet, to see this variability in action. A number of factors can affect the value of an asset: the general economic and political climates, the past performance, and the perceived ability of the leaders of the company in the case of company stock, the weather for an agriculture-related company stock, technological changes, natural disasters, etc. It may be futile to try to find all the factors that affect the value of a particular asset. Thus it is more efficient to think of the value of your asset holdings as a random variable X. As defined in §4.2, X is a numerical function defined on a sample space $\Omega: \omega \to X(\omega), \omega \in \Omega$, where all the unknown factors that can cause the fluctuations in the observable values of X are encapsulated in that sample (point) ω .

We may interpret a sample point in Ω as a "state of the economic world." For example, Ω may consist of two elements: a "good" state of the economic world, ω_1 , and a "bad" state of the economic world, ω_2 . Then we may assume that the value X of your stock holdings a year from now is of the form $X(\omega_1) = ux_0$ and $X(\omega_2) = dx_0$, where u > 1 and $0 \le d < 1$. This example shows your fortune X as rising if the state of the economic world is ω_1 , and dropping if ω_2 is sampled instead (stock prices cannot be negative; the condition d > 0 is the mathematical expression of what financial analysts call "limited liability"). This example, however, is a particular instance. There can be other stocks that behave differently. Their prices drop if the "good" state of the economic world ω_1 is sampled and increase if the "bad" state of the economic world is sampled. This example highlights the non-necessity of the literal interpretation of the elements of the sample space (see also the discussion on sample spaces on pp. 78–80). In fact, we might as well have chosen our sample space Ω to be that of a not-necessarily fair coin toss, so often discussed in this book (see its introduction as Example 8 in §2.4), where $\omega_1 = T$ (tails) and $\omega_2 = H$ (heads). This two-point sample space can in fact generalize to a countable sample space. It is also common to consider Ω as an uncountable sample space representing all shades of possible states of the economic world, such as the set of all real (or positive) numbers. As a result, X can take real values as well.

Consider now the random variable Δ defined by

$$\omega \in \Omega : \omega \to \Delta(\omega) = X(\omega) - x_0.$$

 Δ represents the change in the value of your asset holdings. It is called a

gain if positive, a loss if negative. Next define the random variable

$$R = \frac{\Delta}{x_0}$$
.

This is the one-year return on your intial investment x_0 . Notice that we have omitted the ω in the notation as usual. In our simple example, we can also write

$$X = (1+R)x_0.$$

The dollar amounts x_0 and X can be expressed as

$$x_0 = np_0$$
 and $X = nP$,

where n is the number of shares of the asset you hold, and p_0 and P are, respectively, the price of the asset today (a constant) and the price of the asset a year from now (a not-necessarily constant random variable). We can therefore also represent R as

$$R = (P - p_0)/p_0.$$

In other words, we do not need to know the number of shares you hold in an asset to determine the return of your investment in this asset. We can determine it by simply using the price information of this asset.

One-Period and Multiperiod Models. For the situation considered so far we were concerned with the value of a single random variable denoted by R above. To define R we have two dates: today (date t_0), where the value of the asset is known with certainty, namely it is a constant x_0 ; and a year from now (date t_1), where the value of the asset is a random variable that is not necessarily constant.

Suppose now that we are interested in the prices of an asset in each of the next 3 years, or 12 months, or 4 quarters. Then we say we are fixing a horizon, to be denoted T, and dates t_0, t_1, \ldots, t_N , where N = 3, 12, 4 for the examples above. The interval [0, T] is thus subdivided into N periods $(t_{i-1}, t_i]$, $i = 1, \ldots, N$, where $t_i - t_{i-1} = T/N$. If we let P_i be the asset price at date t_i , then we can define returns

$$r_i = \frac{P_i - P_{i-1}}{P_{i-1}} = \frac{P_i}{P_{i-1}} - 1 \text{ for period } i = 1, 2, \dots, N.$$
 (9.2.1)

This is the context of a *multiperiod* model, of which the one-period model is a particular case with N=1. The latter will be the focus of most of this chapter.

An asset return expresses a relative price change over a specific period, e.g., 1% over one month, 5% over a quarter, 17% over a two–year period. As

we have defined it, a return is in effect a rate of return, i.e., a change relative to time, but we shall use both the word and expression interchangeably. It is customary to express return rates in a single unit, % per year, to allow for rapid comparison and to use standard units for numerical values in formulas. This procedure is called annualization. It is easier to estimate the annualized rate of return by working with the gross return, 1+r, instead of r. For period i we have $1+r_i=P_i/P_{i-1}$. So now if we compute the gross return over N periods, P_N/P_0 , and relate it to that of every period between 1 and N, we have

$$\frac{P_N}{P_0} = \frac{P_N}{P_{N-1}} \times \frac{P_{N-1}}{P_{N-2}} \times \dots \times \frac{P_1}{P_0}$$
$$= \prod_{i=1}^{N} (1 + r_i).$$

If each period is less than a year, and a year consists of N of these periods, then the annualized rate of return $AR^{(N)}$ satisfies

$$\frac{P_N}{P_0} = 1 + AR^{(N)} = \prod_{i=1}^{N} (1 + r_i).$$

For small values of r_i we can write $AR^{(N)} \sim \sum_{i=1}^{N} r_i$ (check it as exercise, use Taylor expansion). So our examples of 1% per month and 5% per quarter become approximately annualized to 12% and 20%, respectively.

If each period is exactly a year, then the annualization consists of finding a fixed (annual) rate $AR^{(N)}$ such that $P_N/P_0 = \left(1 + AR^{(N)}\right)^N$. Thus

$$AR^{(N)} = (P_N/P_0)^{1/N} - 1.$$

Our example of 17% over a two-year period converts to $\sqrt{1.17}-1$, or 8.17%, per year. Note that in the last example, the return of 17% is for the *entire* two-year period, and its conversion to 8.17% per year is for each of the two years. Often we observe returns r_1, r_2, \ldots, r_N for a number N of successive years and would like to find an average of r_1, r_2, \ldots, r_N . Because of the effect of compounding, which is often illustrated via the expression "earning interest on interest" in the case of savings accounts, the precise way to find the average annual rate $AR^{(N)}$ is to start with $P_N/P_0 = (1 + AR^{(N)})^N$, which yields

$$\frac{1}{N}\log\left(P_N/P_0\right) = \log\left(1 + AR^{(N)}\right).$$

Thus

$$\log\left(1 + AR^{(N)}\right) = \frac{1}{N}\log\left[\prod_{i=1}^{N}(1+r_i)\right]$$
$$= \left[\frac{1}{N}\sum_{i=1}^{N}\log(1+r_i)\right].$$

For small values of r_i we have the approximation $AR^{(N)} \sim (1/N) \sum_{i=1}^{N} r_i$ (check it using Taylor expansion).

Asset Risk. As we just saw in the example above, and as you may know from experience, investments can be risky: you could end up with less than what you started out with, or even lose everything! But how do we express (model) this risk mathematically? We have already encountered a similar situation where a gambler risks losing everything (cf. Problem 1 in §8.1). So is it reasonable to think of your investment risk as the probability of losing all of your investment, i.e., $P\{X \leq 0\}$? Or half your initial investment, i.e., $P\{X < x_0/2\}$? Note in passing that in the former case we use $X \leq 0$ instead of X = 0 as, in general, X may become negative when borrowing is allowed as we shall see.

There are, in fact, several ways in which financial theorists model risk. Obviously we want to model risk so that we can assess whether some financial instruments such as stocks are too risky for our tastes, and so that we can decide how much, if at all, we should invest in them. Among financial economists, a commonly accepted way to capture an investor's attitude towards risky investments is through the use of the so-called utility functions that can be theoretically identified for each individual. We do not pursue this approach in this book. In this chapter we use another risk measure that is popular with both financial theorists and practitioners.

Definition 1. The risk of an asset with return R is its variance $\sigma^2(R)$ (see §6.3 for the definitions of the variance and standard deviation of a random variable).

Remark. We could have used the standard deviation $\sigma(R)$ instead of its square $\sigma^2(R)$. The advantage of this alternative is that its unit of account (% per unit of time) is the same as that of R. But we prefer to use the variance for mathematical convenience.

Riskless Security. A particular type of security (asset) plays an important role in finance theory: it is one where we know with certainty the rate of return if we invest in it. Namely, if X is the value of this asset a year from now, then $X(\omega) = (1+r_f)x_0$, for all $\omega \in \Omega$, where the annual rate of return

 r_f is a constant. For simplicity, we will consider this particular security in the form of a money market instrument introduced in §9.1. Its return is $R(\omega) = (X(\omega) - x_0)/x_0 = 1 + r_f, \omega \in \Omega$ (for simplicity, we assume $r_f \geq 0$). Its return variance is $E\left\{(R - E(R))^2\right\} = E\left\{((1 + r_f) - (1 + r_f))^2\right\} = E\left\{0\right\} = 0$. Because its return has zero variance, such a security is called riskless.

9.3. Portfolio allocation

Suppose that you currently own $n_1 = 100$ shares of a company stock, say AT&T, and $n_2 = 80$ shares of another company, say Ford. In addition, assume that the current price of the first stock is $P_1 = \$60$ and that of the second is $P_2 = \$50$. Then your current wealth is

$$x_0 = n_1 P_1 + n_2 P_2$$

= 6,000 + 4,000
= 10,000 dollars.

The quantities $\alpha_1 = n_1 P_1/x_0 = 6,000/10,000 = 0.60$ and $\alpha_2 = n_2 P_2/x_0 = 4,000/10,000 = 0.40$ represent the proportions of your wealth invested in AT&T and Ford, respectively. Alternatively, α_1 and α_2 are expressed in percentages so that in this particular example you are holding $\alpha_1 = 60\%$ of your wealth in the first stock and $\alpha_2 = 40\%$ of your wealth in the second. Notice that we have

$$\alpha_1 + \alpha_2 = 1$$
.

Definition 2. Given M investment opportunities, an investor's portfolio allocation, or investment strategy, consists of an M-tuple $(\alpha_1, \alpha_2, \ldots, \alpha_M)$ of reals such that

$$\sum_{i=1}^{M} \alpha_i = 1.$$

For $i \in \{1, ..., M\}$, α_i represents the proportion of the investor's total wealth invested in asset i and is also called the investor's portfolio weight for asset i.

Definition 3. Given a portfolio allocation $(\alpha_1, \alpha_2, \dots, \alpha_M)$, the corresponding portfolio return is defined as the random variable

$$\sum_{i=1}^{M} \alpha_i R_i,$$

where R_i is the one-period return (a random variable) for asset i.

9.4. Diversification

Ideally, we make investment decisions with an objective in mind. One such objective is that of reducing *portfolio risk*, defined as the variance of the portfolio return, similarly to individual assets (see Definition 1). One way to reduce portfolio risk is through *diversification*, which is a portfolio allocation defined as follows:

Definition 4. Let $(\alpha_1, \alpha_2, \ldots, \alpha_M)$ be a portfolio allocation with corresponding risk $\sigma^2\left(\sum_{i=1}^M \alpha_i R_i\right)$, where R_i is the one-period return for asset i. A portfolio $(\alpha_1^{'}, \alpha_2^{'}, \ldots, \alpha_M^{'})$ is a diversification of the original portfolio if $\sigma^2\left(\sum_{i=1}^M \alpha_i^{'} R_i\right) < \sigma^2\left(\sum_{i=1}^M \alpha_i R_i\right)$.

Example 1. Consider M=2 assets. Assume that $\sigma^2(R_1)=\sigma^2(R_2)>0$, i.e., the two assets have the same risk. Assume further that $-1 \le \rho_{12} < 1$, where ρ_{12} is the correlation coefficient between R_1 and R_2 (see §6.3 for the definitions of correlation and covariance). If the current allocation is $(\alpha_1, \alpha_2) = (1, 0)$, then diversification occurs with a new portfolio $(\alpha'_1, \alpha'_2) = (\alpha, 1 - \alpha)$ such that $0 < \alpha < 1$. Indeed,

$$\sigma^{2}(\alpha R_{1} + (1 - \alpha)R_{2}) = \sigma^{2}(R_{1}) + 2\alpha(1 - \alpha)\left[\operatorname{Cov}(R_{1}, R_{2}) - \sigma(R_{1})\sigma(R_{2})\right]$$

$$< \sigma^{2}(R_{1}) = \sigma^{2}(\alpha_{1}R_{1} + \alpha_{2}R_{2}), \qquad (9.4.1)$$

where the first equality results from $\sigma^2(R_1) = \sigma^2(R_2) > 0$, and the inequality is a consequence of $0 < \alpha < 1$ and $\operatorname{Cov}(R_1, R_2) / (\sigma(R_1)\sigma(R_2)) = \rho_{12} < 1$. In the particular case where $\operatorname{Cov}(R_1, R_2) \leq 0$, we have the following interpretation: when offered an opportunity to invest in two assets with equal risk and with returns that are either uncorrelated $(\operatorname{Cov}(R_1, R_2) = 0)$ or negatively correlated $(\operatorname{Cov}(R_1, R_2) < 0)$, e.g., when their returns tend to be of opposite signs, investing in both is less risky than investing in one only.

Example 2. In the setting of Example 1, is there an allocation $(\alpha, 1 - \alpha)$, with $0 < \alpha < 1$, that is least risky? In other words, is there $\alpha \in (0, 1)$ that minimizes the function $V(\alpha)$ defined by

$$V(\alpha) = \sigma^2 (\alpha R_1 + (1 - \alpha)R_2)?$$

From (9.4.1) we see that V is a quadratic function in α , with second derivative

$$V^{"}(\alpha) = 2\sigma^{2}(R_{1}) + 2\sigma^{2}(R_{2}) - 4\rho_{12}\sigma(R_{1})\sigma(R_{2}).$$

Under the assumptions $-1 \le \rho_{12} < 1$ and $\sigma(R_1) = \sigma(R_2) > 0$, we have

$$V''(\alpha) > 2\sigma^{2}(R_{1}) + 2\sigma^{2}(R_{2}) - 4\sigma(R_{1})\sigma(R_{2}) = 2(\sigma(R_{1}) - \sigma(R_{2}))^{2} = 0.$$

Since $V''(\alpha) > 0$, to get the minimizing α , say α^* , we solve:

$$V'(\alpha^*) = 2\alpha^*\sigma^2(R_1) - 2(1 - \alpha^*)\sigma^2(R_2) + 2\rho_{12}\sigma(R_1)\sigma(R_2)(1 - 2\alpha^*) = 0.$$

The second equation above simplifies, after canceling $2\sigma^2(R_1)$, to

$$\alpha^* - (1 - \alpha^*) + \rho_{12} (1 - 2\alpha^*) = 0,$$

from which we deduce

$$\alpha^* = 1/2.$$

To summarize, we have two assets with the same risk (same return variance) and the most diversified portfolio, i.e., with the smallest return variance, is the one where we invest our fortune equally in each of these assets. This result is interesting because this allocation is independent of the strength of the return correlation ($\rho_{12} \neq 1$) between the two assets. The corresponding risk is

$$V(\alpha^*) = \frac{1}{2} (1 + \rho_{12}) \sigma^2(R_1) < \sigma^2(R_1)$$

because $\rho_{12} < 1$. Note that if the two assets are such that their returns are perfectly negatively correlated, i.e., $\rho_{12} = -1$, then $V(\alpha^*) = 0$. The expected rate of return of this portfolio is $(E(R_1) + E(R_2))/2$. So when $\rho_{12} = -1$, we are in fact almost surely assured of the expected return of the portfolio.

Example 3. Consider again two assets such that $\sigma(R_1) > 0$ and $\sigma(R_2) = 0$. The second asset is a riskless security as we saw previously. Then for $0 \le \alpha \le 1$, $V(\alpha) = \sigma^2 (\alpha R_1 + (1 - \alpha) R_2) = \alpha^2 \sigma^2 (R_1) \ge 0$. Thus the minimand α^* of the return variance of the portfolio $(\alpha, 1 - \alpha)$ is $\alpha^* = 0$ and $V(\alpha^*) = 0$ is the smallest risk.

9.5. Mean-variance optimization

As just seen in the examples above, one can diversify with a judicious choice of portfolio allocation. However, can such risk reduction be too severe so as to result in corresponding returns that are significantly smaller than the original, less diversified portfolio allocation? It is generally assumed that investors are willing to take on additional risk for a chance of getting higher

returns, but can this trade-off be done systematically? Harry Markowitz, an economist credited with the first formalization of this trade-off for which he won the Nobel prize, proposes the following approach: for a given level of expected returns, find the portfolio allocation with smallest risk. A dual problem is that of fixing the level of portfolio risk and then looking for the corresponding allocation that maximizes the expected return of this portfolio. Mathematically, it is easier to deal with the former problem expressed as follows. Find $(\alpha_1, \alpha_2, \ldots, \alpha_M)$ that

minimize
$$\frac{1}{2}\sigma^2\left(\sum_{i=1}^M \alpha_i R_i\right)$$
, (9.5.1)

where

$$\sum_{i=1}^{M} \alpha_i = 1, \tag{9.5.2}$$

$$E\left(\sum_{i=1}^{M} \alpha_i R_i\right) = \mu,\tag{9.5.3}$$

and μ is given (desired expected portfolio return). In practice, additional conditions are imposed. They are generally of the form

$$l_i \le \alpha_i \le u_i, \quad 1 \le i \le M, \tag{9.5.4}$$

where u_i is generally nonnegative and l_i can be negative.

Short and Long Positions. An investor with $\alpha_i \neq 0$ is said to hold a position in asset i. If $\alpha_i < 0$, then the investor is said to short (or hold short) asset i. If $\alpha_i > 0$ the investor is said to go long (or hold long) asset i. In the above problem, if $l_i < 0$ (typically $l_i = -1$ or -.5), then the investor is allowed to short asset i. An investor shorts an asset by borrowing this asset. When investors short the riskless asset, they borrow money to invest for example in other assets as part of their optimal strategies. When you hold your money in a savings account, you long the riskless security. When investors short stocks, they borrow shares from their brokers and sell them on the market. Later they return these shares to their brokers by repurchasing them on the market.

A Three-Asset Example (1). In its general form, problem (9.5.1) - (9.5.4) does not necessarily lead to closed-form solutions for $\alpha_1, \ldots, \alpha_M$. It is generally solved using the tools of quadratic programming (see, for example, [Huang and Litzenberger] or [Luenberger]). When condition (9.5.4) is omitted, one can solve (9.5.1) - (9.5.3) in closed form, if a solution exists.

We illustrate here the case M=3, using techniques appropriate for the general level of this book. For i=1,2,3 let $\mu_i=E(R_i)$, $\sigma_i=\sigma(R_i)$ and let ρ_{ij} be the correlation coefficient between R_i and R_j , j=1,2,3. Given μ , (9.5.2) and (9.5.3) can be restated as

$$\alpha_3 = 1 - \alpha_2 - \alpha_1, \tag{9.5.5}$$

$$\alpha_1(\mu_1 - \mu_3) + \alpha_2(\mu_2 - \mu_3) = \mu - \mu_3.$$
 (9.5.6)

Assume for now that $\mu_2 \neq \mu_3$. Then from (9.5.5) and (9.5.6) we have

$$\alpha_2 = \frac{\mu - \mu_3}{\mu_2 - \mu_3} - \frac{\mu_1 - \mu_3}{\mu_2 - \mu_3} \alpha_1, \tag{9.5.7}$$

$$\alpha_3 = \frac{\mu - \mu_2}{\mu_3 - \mu_2} - \frac{\mu_1 - \mu_2}{\mu_3 - \mu_2} \alpha_1. \tag{9.5.8}$$

Thus (9.5.1) can be expressed as

minimize
$$V(\alpha_1) = \frac{1}{2} \sum_{i,j=1}^{3} \alpha_i \alpha_j \rho_{ij} \sigma_i \sigma_j,$$
 (9.5.9)

with α_2 and α_3 expressed in terms of α_1 . We can rewrite (9.5.7) and (9.5.8) as

$$\alpha_i = a_i \alpha_1 + b_i \mu + c_i$$
, for $i = 2, 3$,

where

$$a_2 = -\frac{\mu_1 - \mu_3}{\mu_2 - \mu_3}, \quad b_2 = \frac{1}{\mu_2 - \mu_3}, \quad c_2 = -\frac{\mu_3}{\mu_2 - \mu_3},$$

$$a_3 = -\frac{\mu_1 - \mu_2}{\mu_3 - \mu_2}, \quad b_3 = \frac{1}{\mu_3 - \mu_2}, \quad c_3 = -\frac{\mu_2}{\mu_3 - \mu_2}.$$

Therefore we can write the derivative of V as

$$V^{'}(\alpha_1) = A\alpha_1 + B\mu + C,$$

where

$$\begin{split} A &= \sigma_1^2 + a_2^2 \sigma_2^2 + a_3^2 \sigma_3^2 + 2a_2 \rho_{12} \sigma_1 \sigma_2 + 2a_3 \rho_{13} \sigma_1 \sigma_3 + 2a_3 \rho_{23} \sigma_3 \sigma_2, \\ B &= a_2 b_2 \sigma_2^2 + a_3 b_3 \sigma_3^2 + b_2 \rho_{12} \sigma_1 \sigma_2 + b_3 \rho_{13} \sigma_1 \sigma_3 + (a_3 b_2 + a_2 b_3) \rho_{23} \sigma_2 \sigma_3, \\ C &= a_2 c_2 \sigma_2^2 + a_3 c_3 \sigma_3^2 + c_2 \rho_{12} \sigma_1 \sigma_2 + c_3 \rho_{13} \sigma_1 \sigma_3 + (a_3 c_2 + a_2 c_3) \rho_{23} \sigma_2 \sigma_3. \end{split}$$

With α_i thus expressed, V in (9.5.9) is a quadratic function of α_1 . Therefore a solution to $V'(\alpha_1) = 0$ corresponds to a global minimand for the problem

(9.5.1) – (9.5.3) with M=3 if $V^{''}(\alpha_1)=A>0$. Assume then that the parameters μ_i , σ_i , and ρ_{ij} , for $1 \leq i, j \leq 3$, are such that the latter condition is satisfied. The solution of $V^{'}(\alpha_1)=0$ yields

$$\alpha_i = A_i \mu + B_i, \quad \text{for } 1 \le i \le 3,$$
(9.5.10)

where $A_1 = -B/A$, $B_1 = -C/A$, and for i = 2, 3, $A_i = a_i (B/A) + b_i$ and $B_i = -a_i (C/A) + c_i$.

The return variance σ^2 of the portfolio with allocations (or weights) α_i , i = 1, 2, 3, as given in (9.5.10) is a function of μ and can be written as

$$\sigma^2 = a\mu^2 + b\mu + c, (9.5.11)$$

where

$$a=\sum_{i,j}^3 A_i A_j
ho_{ij} \sigma_i \sigma_j, \quad b=\sum_{i,j}^3 (A_i B_j + A_j B_i)
ho_{ij} \sigma_i \sigma_j,$$
 and $c=\sum_{i,j}^3 B_i B_j
ho_{ij} \sigma_i \sigma_j.$

Recall that we arrived at (9.5.11) by fixing μ and looking for portfolio weights (α_i) that minimize the return variance of the portfolios with expected returns equal to μ . Recall also that the problem just solved can be viewed as first fixing the return variance σ^2 and then looking for the weights (α_i) maximizing the expected return of the corresponding portfolio. Equation (9.5.11) shows that in the variance-expected return space the graph of σ^2 as a function of μ is a parabola. Note that the units of μ (in percent per unit of time; e.g., 2% per year) are not directly comparable to those of σ^2 (in square percent per square unit of time). A more meaningful and in fact more common representation of the relation between μ and σ^2 is in the standard deviation-expected rate of return space.

Suppose the following holds:

$$a > 0$$
, $b^2 < 4ac$, and $b/2a < 0$.

Then (9.5.11) can be restated as

$$\frac{\sigma^2}{(4ac-b^2)/4a} - \frac{(\mu - (-b/2a))^2}{(4ac-b^2)/4a^2} = 1,$$
 (9.5.12)

thus expressing the relation between σ , the standard deviation of the return of a portfolio with allocations (α_i) as in (9.5.10), and μ , the corresponding expected return, as a hyperbola with asymptotes $\mu = -b/2a \pm \sigma/\sqrt{a}$ (see

Fig. 37). The pair $(\bar{\sigma}, \bar{\mu})$ corresponds to the portfolio that achieves the smallest variance $(\bar{\sigma} = \sqrt{(4ac - b^2)/4a})$ among all possible portfolios, with resulting expected return $\bar{\mu} = -(b/2a)$.

Efficient Frontier. Notice that when we solved the optimization problem (9.5.1)–(9.5.3) we did not impose any restrictions on the desired level μ . As a result, μ can conceivably span the entire real line. However, as a quick look at Figure 37 would indicate, if we fix a risk level at a certain value σ_0 then there correspond two values, μ^* and μ^{**} , of μ for which this level of risk σ_0 is minimal. At this point, economic theory about the behavior of investors is invoked to distinguish between the two points $A^* = (\sigma_0, \mu^*)$ and $A^{**} = (\sigma_0, \mu^{**})$ on the curve. In particular, it is postulated that investors always prefer more, even if only in expected returns, when comparing two alternatives that bear the same risk. For this reason, at risk level σ_0 , a typical investor is going to select the portfolio that yields the expected return μ^* instead of that which yields the expected return μ^{**} . In this manner, we only need focus on the part of the graph containing points (σ, μ) such that $\mu > \bar{\mu}$. This portion of the graph, which is separately plotted in Figure 38, is called the efficient frontier. This label comes from the interpretation that any point (σ^*, μ^*) in Figure 38 corresponds to a portfolio allocation α^* such that for any other portfolio allocation α' yielding expected portfolio return μ' and standard deviation return σ' we have the following: if $\sigma' < \sigma^*$, then $\mu' < \mu^*$, and if $\mu' > \mu^*$, then $\sigma' > \sigma^*$. In other words, no other portfolio allocation than α^* can achieve a higher expected return than μ^* with smaller risk than σ^* .

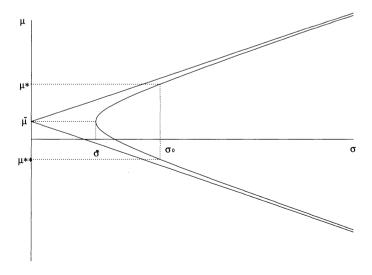


Figure 37 Portfolio frontier in the standard deviation (σ) - mean (μ) space

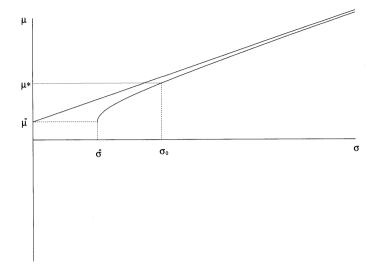


Figure 38 Efficient frontier in the standard deviation (σ) - mean (μ) space

The Case of M Securities. The result just obtained for three assets generalizes in exactly the same manner to that of M securities (i.e., the relation between the maximum expected return for a given level of risk is a hyperbola in the standard deviation-expected rate of return space.) Recall that in the derivation of this result we made some assumptions, particularly about the second derivative of the portfolio variance as a function of the weights. These assumptions are in fact easier to express using linear algebra in vector-matrix form. In the general setting of M assets, the problem is to find allocations $(\alpha_1, \alpha_2, \ldots, \alpha_M)$, denoted by the $M \times 1$ vector α , that

$$\text{minimize} \quad \frac{1}{2}\alpha^{\mathrm{T}}W\alpha$$

such that

$$\alpha^{\mathrm{T}} \epsilon = \mu,$$
 $\alpha^{\mathrm{T}} \mathbf{1} = 1,$

where W is the return covariance matrix for the M assets, i.e., $W_{ij} = \text{Cov}(R_i, R_j)$, $1 \leq i, j \leq M$, ϵ is an $M \times 1$ vector such that $\epsilon_i = \mu_i$, $i = 1, 2, \ldots, M$, and $\mathbf{1}$ is the M vector whose elements are each equal to 1.

At this point the derivation of the actual minimizing portfolios requires some knowledge that is slightly beyond the level of this book. The interested reader can consult Chapter 3 in [Huang and Litzenberger] where full details are displayed in the spirit of the presentation of this chapter. The matrix

formulation helps us identify more readily some of the conditions used in the three-asset example worked out earlier. In particular, we required that the second derivative of the portfolio variance as a function of the portfolio weights be positive. From the general matrix formulation, this condition is satisfied if and only if the covariance matrix W is positive definite, which by definition means that for every nonzero vector x we have $x^{\mathrm{T}}Wx > 0$. Suppose now that a riskless security is among those considered in the optimal allocation problem. Without loss of generality, we can assume that it is asset 1. Then, since this asset has a deterministic value at the end of the time period, the standard deviation σ_1 of the one-period return of asset 1 is zero and the correlation coefficient ρ_{1i} of the riskless asset with any other asset i is also zero. In other words, this means that all the elements in the first column and in the first row of W are equal to zero. Therefore, for the matrix W to be positive definite it is necessary that the riskless asset be excluded from consideration in the optimization problem. We will soon see that the inclusion of the riskless asset in the optimal asset allocation problem can still be handled, leading to a different characterization of the trade-off between optimal expected return and risk level.

Effect of Riskless Security. In addition to M risky assets, i.e., assets for which $\sigma^2(R_i) > 0$, $1 \le i \le M$, we now include asset 0, which is riskless, i.e., $\sigma^2(R_0) = 0$. We are now looking for M+1 portfolio weights $(\alpha_0, \alpha_1, \alpha_2, \ldots, \alpha_M)$ that again minimize the corresponding portfolio risk

$$\sigma^2 \left(\sum_{i=0}^M \alpha_i R_i \right) \tag{9.5.13}$$

under the constraints

$$\sum_{i=0}^{M} \alpha_i = 1, \tag{9.5.14}$$

$$E\left(\sum_{i=0}^{M} \alpha_i R_i\right) = \mu, \tag{9.5.15}$$

where μ is the desired expected return, with the possible inclusion of additional constraints of the form (9.5.4). Since $\sigma(R_0) = 0$, it follows that $\text{Cov}(R_i, R_0) = \rho_{i0}\sigma(R_i)\sigma(R_0) = 0$, and since $\alpha_0 = 1 - \sum_{i=1}^{M} \alpha_i$ by (9.5.15), the problem becomes that of finding $(\alpha_1, \alpha_2, \dots, \alpha_M)$ that minimize (9.5.13) such that

$$E\left(\sum_{i=1}^{M} \alpha_i (R_i - R_0)\right) = \mu - R_0. \tag{9.5.16}$$

A Three-Asset Example (2). This case is similar to the three-asset example (1) we saw earlier, now with M=2 risky assets and one riskless asset. Here too, conditions of the type (9.5.4) are not imposed. Following the same notation we also let $\mu_0 = R_0$. From (9.5.16) we have (assuming $\mu_2 \neq \mu_0$)

$$\alpha_2 = a + b\alpha_1, \tag{9.5.17}$$

where $a = (\mu - \mu_0)/(\mu_2 - \mu_0)$ and $b = -(\mu_1 - \mu_0)/(\mu_2 - \mu_0)$. Therefore the problem reduces to that of finding α_1 that minimizes the function

$$V(\alpha_1) = \alpha_1^2 \sigma_1^2 + (a + b\alpha_1)^2 \sigma_2^2 + 2\alpha_1 (a + b\alpha_1)\rho_{12}\sigma_1\sigma_2$$
 (9.5.18)

from which we derive

$$V'(\alpha_1) = 2\alpha_1 \left(\sigma_1^2 + b^2 \sigma_2^2 + 2b\rho_{12}\sigma_1\sigma_2\right) + 2a\rho_{12}\sigma_1\sigma_2,$$

$$V''(\alpha_1) = 2\left(\sigma_1^2 + b^2 \sigma_2^2 + 2b\rho_{12}\sigma_1\sigma_2\right)$$

$$= 2\sigma^2 \left(R_1 + bR_2\right),$$
(9.5.19)

where we recognize in (9.5.19) the variance of the random variable $R_1 + bR_2$ thanks to the properties of the variance (see §6.3), which also includes $\sigma^2(Z) \geq 0$ for any random variable Z. Notice again that (9.5.18) expresses V as a quadratic function of α_1 . Therefore, as long as $V^{"}(\alpha_1) > 0$, in order to find the minimizing α_1 we need only solve $V^{'}(\alpha_1) = 0$. The solution of the latter is

$$\alpha_1^* = -a \frac{\rho_{12}\sigma_1\sigma_2}{\sigma_1^2 + b^2\sigma_2^2 + 2b\rho_{12}\sigma_1\sigma_2}$$
(9.5.20)

under the condition that

$$\sigma_1^2 + b^2 \sigma_2^2 + 2b\rho_{12}\sigma_1\sigma_2 \neq 0. (9.5.21)$$

Note that this condition will make $V^{"}(\alpha_1) > 0$.

Assuming condition (9.5.21), with α_1^* given by (9.5.20), the other optimal portfolio weights α_0^* and α_2^* are obtained directly through (9.5.17), where $\alpha_1 = \alpha_1^*$, and through

$$\alpha_0^* = 1 - (\alpha_1^* + \alpha_2^*).$$

In a similar manner to that of the first three-asset example, we now derive a relationship between the desired expected level μ and the optimal portfolio variance σ^2 resulting from the application of the portfolio $(\alpha_0^*, \alpha_1^*, \alpha_2^*)$ in order to achieve μ . Let

$$C = -\frac{\rho_{12}\sigma_{1}\sigma_{2}}{\sigma_{1}^{2} + b^{2}\sigma_{2}^{2} + 2b\rho_{12}\sigma_{1}\sigma_{2}}$$

and

$$K = \frac{\left[C^2 \sigma_1^2 + (1 + bC)^2 \sigma_2^2 + 2C(1 + bc)\rho_{12}\sigma_1\sigma_2\right]}{(\mu_2 - \mu_0)^2}.$$

Then from (9.5.18) and (9.5.20) we have

$$\sigma^2 = V(\alpha_1^*) = (\mu - \mu_0)^2 K, \tag{9.5.22}$$

where we remark that

$$K = \frac{\sigma^2 \left(CR_1 + (1 + bC)R_2 \right)}{(\mu_2 - \mu_0)^2} \ge 0$$

since the variance of a random variable is always nonnegative and since $(\mu_2 - \mu_0)^2 > 0$ by the assumption $\mu_2 \neq \mu_0$. We can also write (9.5.22) as

$$\sigma = |\mu - \mu_0|\sqrt{K},\tag{9.5.23}$$

therefore showing that the graph of the minimum standard deviation-expected portfolio return is the union of two half-lines emanating from the point $(0, \mu_0)$, more precisely of the form:

$$\sigma = \left\{ \begin{array}{ll} \sqrt{K}(\mu - \mu_0) & \text{if } \mu \geq \mu_0, \\ -\sqrt{K}(\mu - \mu_0) & \text{if } \mu < \mu_0. \end{array} \right.$$

What happens to security prices when all investors adopt the same mean-variance optimization, thus allocating their money in the same proportions across all the stocks? In simple terms, the prices will be determined by the available supply of the stock shares and by the demand for these shares as dictated by the optimal portfolio allocation α^* for all stocks. This equilibrium argument forms the basis of what is referred to as the Capital Asset Pricing Model, fully explored by William Sharpe, and for which he shared the Nobel Prize with Harry Markowitz, mentioned earlier. The derivation of these prices can be made mathematically complete (cf., for example [Duffie] or [Huang and Litzenberger]). This theory is an approach that predicts prices one period at a time. Prior to this approach, a number of economists had attempted to find the right price at which stocks should trade. Often these economists were academically oriented and their views clashed with those who participated actively in the stock markets. Regularly, economists were scorned for their rigid and "unprofitable" views there were indeed very few economists who used their theories to become "rich and famous"! One exception though, is that of Keynes, who acquired a large fortune by investing and who held what he called gamblers, who buy stocks only for short-term gains, in contempt.

The main point to be noted about this section is that it illustrates the importance of diversification. There are several ways in which this message is put into practice. For example, mutual funds* are investment assets where small investors, such as teachers or even students, can pool their money together to buy stocks in a diversified manner. Since the early 1980s, mutual funds have become a favorite vehicle for workers to save and invest for retirement. Unfortunately, some employees have most of their retirement money invested only in the stock of the company they work for. As it lacks diversification, this approach is at the mercy of the fortunes of a single company. A prominent example where this strategy failed miserably and affected thousands of employees is that of Enron, which in the year 2001 was one of the largest corporations in the world. To conclude, we mention the opinion of the renowned American economist Frank H. Knight who in his book [Knight] argued philosophically that not only is risk a fact of life, but it is necessary to entrepreneurship and the realization of profits from investments. Incidentally, his son Frank B. is the author of a treatise on Brownian motion (see Chapter 8, Section 2), the theory of which is now widely used as a model for stock market movements.

9.6. Asset return distributions

For the results we have derived so far we made use almost exclusively of means, variances, and covariances. Little was said about the distributions of these random variables. Clearly, two random variables can have the same mean and variance and yet have completely different distribution functions. An obvious example is that of one random variable that is integer-valued, e.g. Bernoulli, and the other has a density, e.g., a normally distributed random variable. If you need to convince yourself with two random variables with densities, do Exercise 7.

With asset returns being at the heart of most financial investment analyses, it is not surprising that their distribution functions have received a lot of attention. On the one hand, actual data (and there is a lot of it) suggest certain types of distributions, and on the other, the desire to obtain mathematically tractable results imposes additional assumptions. In anticipation of the next chapter we discuss the distribution of asset returns in the multi-period context. Consider a situation with N periods $(t_0,t_1],(t_1,t_2],\ldots,(t_{N-1},t_N]$. Let S_n be the price of a security at time n and define the one-period returns $r_n=(S_n-S_{n-1})/S_{n-1},\,n=1,2,\ldots,N$. Define also the N-period return $R_N=(S_N-S_0)/S_0$ and the random variables

^{*}There are thousands of mutual funds generally grouped in families, but we mention one in particular, the Alger Group, where the former second grader (at the time of the first edition) appearing at the beginning of §1.1, who rode the merry-go-round horses in Example 14 of Chapter 8 (see Fig. 34), is currently Chief Investment Officer.

$$\xi_n = \ln(1 + r_n)$$
, for $1 \le n \le N$, and $\Psi_N = \ln(1 + R_N)$.

It is commonly assumed that the random variables $\xi_1, \xi_2, \ldots, \xi_n, \ldots, \xi_N$ are identically and independently distributed (equivalently, r_1, r_2, \ldots, r_N are identically and independently distributed — this is a direct application of Proposition 6 in §5.5). There is some evidence from actual data to support this assumption. In addition, it is commonly assumed that the random variables ξ_n have normal distributions. Below we discuss whether this assumption is reasonable.

If n refers to a small period index (e.g., r_n represents a daily or hourly return), actual data suggest that the distribution of ξ_n has "fatter" tails than the normal distribution. Formally, a random variable Y is said to have a fatter right tail than the random variable Z if $P\{Y > \xi\} > P\{Z > \xi\}$ for all ξ sufficiently large. Similarly, Y is said to have a fatter left tail than Z if $P\{Y < \xi\} > P\{Z < \xi\}$ for all $\xi < 0$ with $|\xi|$ sufficiently large. So for a small period index n, if we assume ξ_n to have a normal distribution we are likely to underestimate the probability of large gains or losses (corresponding to large absolute values of ξ_n).

Cauchy Distribution. A Cauchy distribution with location parameter μ and scale parameter $\sigma > 0$ is defined as having the density function

$$f(x) = \left[\pi\sigma\left(1 + \left(\frac{x-\mu}{\sigma}\right)^2\right)\right]^{-1}$$
.

Logarithmic Scale. As we have seen, the one-period model considered in this chapter requires finiteness of return variances. Therefore the Cauchy distribution is ruled out for this type of model. In the next chapter we shall consider multiperiod models, where the distribution of the one-period returns r_n or their log transforms ξ_n will be assessed on the basis of the random variable $\sum_{i=1}^{N} \xi_n$. To understand the significance of the latter, recall that

$$1 + r_n = \frac{S_n}{S_{n-1}},\tag{9.6.1}$$

and therefore

$$\xi_n = \log(1 + r_n) = \log S_n - \log S_{n-1},$$
(9.6.2)

thus showing that $\{\xi_n\}$ represents the successive price changes on a logarithmic scale. ξ_n is sometimes called the continuously compounded return or log return for period n (see also the examples of §9.2 for multi-period

models). Therefore

$$\sum_{i=1}^{N} \xi_n = \Psi_N \tag{9.6.3}$$

represents the cumulative log return between t_0 and t_N . If we assume, as we will, that the random variables $\xi_1, \xi_2, \ldots, \xi_N$ are independently and identically distributed, then we may be able to use readily available tools such as the random walk properties of Chapter 8 and the central limit theorem of §7.5.

Recall that because all prices S_n are required to remain positive, r_n is equivalently required to remain above -1, thus potentially limiting our choices for a distribution for r_n . On the other hand, ξ_n may take any value on the real line and thus allows us more flexibility in the choice of its distribution. More importantly though is the fact that the log transform allows us to work with more tractable models, i.e., where we can obtain results relatively quickly.

9.7. Stable probability distributions

Let N be the number of business days in a given month. Suppose t_1, t_2, \dots, t_N represent the times at which the stock market closes on the successive business days in that month. Let t_0 correspond to the last closing time in the preceding month. Then for $1 \leq n \leq N$, S_n , r_n , and ξ_n represent, respectively, the closing price, the return, and the log return for day (period) n. Let S_0 be the stock price at closing on the last business day of the preceding month. We now have two ways of estimating the distribution of the monthly log return Ψ_N : one is to estimate the distribution of the daily log returns ξ_n , using daily returns over 5 years, say, and then use (9.6.3); the other is to estimate this distribution directly by using monthly returns, through closing prices at the end of each month of the same 5 years. For simplicity, assume there are N (say N=22) business days in each month over the five-year span. Then we observe $5 \times 12 \times N$ daily returns over the five years. These observations help us decide on the distribution of the daily returns r_n , thus that of their logarithms ξ_n , and finally that of the monthly returns via their logarithm Ψ_N through $\Psi_N = \sum_{n=1}^N \xi_n$. The second approach is to observe 5×12 monthly returns over the five years and infer the distribution of monthly returns (or equivalently their logarithm Ψ_N). Notice the difference between the two approaches: with the first, we use actual data to infer the distribution of the daily returns. The distribution of the monthly returns is then obtained as a result of the addition of N random variables (the daily returns). With the second approach, the distribution of the monthly returns is inferred directly from the data, not as a sum of the N random variables (daily returns). Ideally,

the latter distribution should be close to that which results from the sum of the daily returns. This property, known as stability, is our next topic.

From Theorem 7 in §7.4, if X_1 and X_2 are two independent random variables with normal distributions $N\left(m_1,\sigma_1^2\right)$ and $N\left(m_2,\sigma_2^2\right)$, respectively, then the random variable X_1+X_2 also has a normal distribution (with mean m_1+m_2 and variance $\sigma_1^2+\sigma_2^2$). In other words, the specific distributions of X_1 , X_2 and X differ only through their means and variances (these are called the parameters of the normal distribution — once known, they specify the distribution precisely). Observe that $Z_i=(X_i-m_i)/\sigma_i$, i=1,2, are independent and follow the unit (or standard) normal distribution. Let $Z=[X_1+X_2-(m_1+m_2)]/\sigma$, where $\sigma^2=\sigma_1^2+\sigma_2^2$. Then we can write:

$$\sigma Z = \sigma_1 \frac{X_1 - m_1}{\sigma_1} + \sigma_2 \frac{X_2 - m_2}{\sigma_2}$$

= $\sigma_1 Z_1 + \sigma_2 Z_2$.

Notice that each of Z_1 , Z_2 and Z follows the unit normal distribution.

Definition. A distribution function is said to be stable if for any two independent random variables X_1 and X_2 with this distribution, and any two positive constants c_1 and c_2 , we have

$$c_1 X_1 + c_2 X_2 = cX + d, (9.7.1)$$

where c is a positive constant, d a constant, both depending on c_1 and c_2 , and X is a random variable with the same distribution as X_1 and X_2 .

Example. As observed above, if X_1 and X_2 are independent, both following a unit normal distribution, then the random variable $Y = X_1 + X_2$ is normally distributed with mean 0 and variance 2. If we want Y to be such that Y = cX + d, where c > 0, d is real and X follows a unit normal distribution, then we must have 0 = E(Y) = cE(X) + d = d and $2 = Var(Y) = c^2Var(X) = c^2$. In other words, $c = \sqrt{2}$ and d = 0 will do.

Notice that if $c_1 = c_2 \neq 0$, then we can write (9.7.1) as $X_1 + X_2 = c'X + d'$, where $c' = c/c_1$ and $d' = d/c_1$. This way, stability refers to obtaining, as a result of the addition of two independent and identically distributed random variables X_1 and X_2 , a similarly distributed random variable modulo a possible scale change (captured by c') and/or a shift from the origin (captured by d'). We have already encountered this form of stability, not only for a single distribution as above, but for an entire family of distributions: in Theorem 3 of §7.2 for the Poisson family and in Theorem 7 of §7.4 for the normal family. In fact, stability is a property that is most useful at the distribution level. Accordingly, we shall refer to both distributions and distribution families (or types*) when considering the stability property.

^{*}As in most texts, we use family and type interchangeably.

Lévy Characterization of Stable Laws. In Theorem 7 of §7.4 we used the fact that, when it is well defined, the moment-generating function (also known as the Laplace transform) of a nonnegative random variable uniquely determines its probability distribution function (cf. §6.5). As mentioned also in §6.5, the moment-generating function of a random variable may not always exist, but its characteristic function (also known as its Fourier transform) always does. Similarly to the moment-generating function, the characteristic function also uniquely determines the probability distribution function of a random variable (cf. the proofs of Theorems 8 and 9 in §7.5).

Let F be a stable law (this is another way of calling a distribution function or a distribution family) and let $\varphi(\theta)$ be its characteristic function [see (6.5.17)]. If X_1 and X_2 are independent with distribution F, then we have, for given positive constants c_1 and c_2 ,

$$E\left[e^{i\theta(c_1X_1+c_2X_2)}\right] = E\left[e^{i\theta c_1X_1}\right] \cdot E\left[e^{i\theta c_2X_2}\right]$$
$$= \varphi(\theta c_1) \cdot \varphi(\theta c_2),$$

where the first equality is justified by the independence assumption, and the second results from the identity of distributions of X_1 and X_2 [cf. Theorem 7 in §6.5]. For a random variable X following the same distribution, and in order to have $c_1X_1 + c_2X_2 = cX + d$ for some positive constant c and real constant d, we must have

$$E\left[e^{i\theta(c_1X_1+c_2X_2)}\right] = E\left[e^{i\theta(cX+d)}\right]$$
$$= E\left[e^{i\theta cX}\right] \cdot e^{i\theta d},$$

or equivalently

$$\varphi(\theta c_1) \cdot \varphi(\theta c_2) = \varphi(\theta c) \cdot e^{i\theta d}. \tag{9.7.2}$$

Example (revisited). For the unit normal distribution, the characteristic function of which is $\varphi(\theta) = e^{-\theta^2/2}$ (see 7.5.7), (9.7.2) becomes

$$e^{-c_1^2\theta^2/2} \cdot e^{-c_2^2\theta^2/2} = e^{-c^2\theta^2/2} \cdot e^{i\theta d},$$
 or

$$e^{-(c_1^2+c_2^2)\theta^2/2} = e^{-c^2\theta^2/2} \cdot e^{i\theta d}$$
.

Thus with $c = \sqrt{c_1^2 + c_2^2}$ and d = 0 we verify that the unit normal distribution is stable.

Paul Lévy has shown that a random variable X with stable distribution has a characteristic function of the form

$$\varphi(\theta) = E\left(e^{i\theta X}\right) = e^{-\gamma_{\alpha}|\theta|^{\alpha} + id\theta},\tag{9.7.3}$$

where $0 < \alpha \le 2$, γ_{α} is a complex constant and d is a real constant (cf. (5) on p. 95 in [Lévy] and the Theorem of §34 on p. 164 in [Gnedenko and Kolmogorov]).

For example, the characteristic function corresponding to a normal distribution with mean m and variance v is

$$\varphi(\theta) = e^{-\frac{v}{2}\theta^2 + im\theta}.$$

Notice that the exponent α is 2 no matter what m or v is. In this sense, $\alpha = 2$ is associated with the normal family.

As another example, the characteristic function for a Cauchy distribution with parameters μ and σ (see §9.6) is

$$\varphi(\theta) = e^{-\sigma|\theta| + i\mu\theta}.$$

Here $\alpha = 1$ is associated with the Cauchy type of distribution.

Remarkably, the exponent α of a stable law uniquely determines whether it has finite mean or variance. If $\alpha=2$, both mean and variance are finite; if $1<\alpha<2$, only the mean is finite; and if $0<\alpha\leq1$, neither mean nor variance is finite (see [Gnedenko and Kolmogorov], p. 182).

As we shall see in the next chapter, we are interested in the distribution of $\sum_{n=1}^N \xi_n$. Then one may ask, when N is large, if the distribution of $\sum_{n=1}^N \xi_n$ could become, as in the central limit theorem (cf. §7.5), independent of the actual distribution of $\{\xi_n\}$. As we alluded to at the end of §7.5, there exist limit theorems for stable laws (see also Appendix 4). However, only the tail distributions of $\sum_{n=1}^N \xi_n$ can be characterized usefully when the tail distributions of ξ_n are Pareto (i.e., for large |x|, $P\{\xi_n > x\}$, if x > 0, or $P\{\xi_n < x\}$, if x < 0, has the form $A/|x|^\beta$ for some constants A and β .) With the exception of the normal distribution, one has to resort to numerical procedures to determine the "middle" (i.e. when |x| is not large) of the limit distribution. For these reasons, and with some supporting empirical evidence, the normal distribution assumption for ξ_n has become standard for practical mathematical models in finance. In this case, one also says that r_n has the lognormal distribution.

Exercises

- 1. For the example given at the beginning of §9.2, express R explicitly in terms of u and d. To familiarize yourself with some common values, consider also the numerical example of u = 1.10 and d = .75 and interpret the corresponding values of R in percentages.
- 2. Consider now the case where R maps to the set of reals. Assume that R is normally distributed with mean μ and variance σ^2 ; see §7.4 and (7.4.6).

- (a) If $\mu = 0$, are you more likely to lose money than to make money?
- (b) If $\mu < 0$, what is the probability of your losing money?
- (c) If $\mu > 0$, what is this probability?

This exercise shows that if we accept the return variance as a risk gauge, then different assets with different probabilities of loss may be considered equally risky.

- 3. Show that for any asset with positive variance, there exists a diversifying portfolio that involves an asset with smaller variance (assuming the second asset exists).
- 4. For the three-asset example (1), exhibit a set of conditions on μ_i , σ_i , and ρ_{ij} for which the function $V^{"}(\alpha_1) \leq 0$ for all α_1 .
- 5. Recall that σ_1 and σ_2 are positive. Are there values of b and ρ_{12} for which condition (9.5.21) is violated?
- 6. Check what happens when $\mu_1 = \mu_0$ or $\mu_2 = \mu_0$. Are there any solutions for $V'(\alpha) = 0$?
- 7. Let X be an exponentially distributed random variable with mean 1 and variance 1, and let Y be a normally distributed random variable with mean 1 and variance 1. Given x > 0, compare $P\{X < x\}$ and $P\{0 < Y < x\}$. [Note: recall that an exponentially distributed random variable X is positive and so $P\{0 < X < x\} = P\{X < x\}$; cf. Example 12 in §4.5.]
- 8. Show that the Pareto distribution (defined in Appendix 4) has a fatter right tail than the standard normal distribution [its tail distribution is given right after (7.4.2)]. [You need only look at the case $x \to \infty$.]
- 9. (a) Show that the mean of a Cauchy distribution (see §9.6) is undefined. What can you say about its variance?
 - (b) Show that if X_1 and X_2 are independent random variables with Cauchy distributions, then $X_1 + X_2$ also has a Cauchy distribution.
 - (c) Plot the density functions of both a Cauchy distribution with $\mu=0$ and $\sigma=1$, and a standard normal distribution. Notice that both curves are symmetric and that the Cauchy plot is above that of the normal for |x|>2. This is a property of fatter tails.
- 10. (a) Show that the sum of two independent random variables with gamma distributions [see Exercise 37(c) in §6.5 for a definition] also has a gamma distribution. [Hint: this is straightforward if you look at exercise 42 in §6.5.]
 - (b) Show that the Pareto distribution is not stable. [Hint: show that for $\alpha = 1$ and A = 1 we have $P\{X+Y > z\} = 2/z + 2\log(z-1)/z^2$].
- 11. The density function of a random variable X with a lognormal distribution is

$$f(x) = \frac{1}{x\sigma\sqrt{2\pi}} \exp^{-(\log x - \mu)^2/(2\sigma^2)}, \text{ for } 0 < x,$$

where μ and $\sigma > 0$ are given parameters. Show that

- (a) the random variable $Y = \log X$ has a normal distribution with mean μ and variance σ^2 ,
- (b) all the moments of X exist and are finite, and
- (c) X does not have a moment-generating function.

Appendix 4

Pareto and Stable Laws

We list here some basic facts about stable laws and the Pareto distribution, which received a lot of attention in the 1960s and 1970s to model stock returns. This distribution is named after Vilfredo Pareto, an engineer turned economist, who studied the distribution of wealth among people at the end of the 19th century. A random variable X is said to have the (standard) Pareto distribution with parameter $\alpha>0$ if

$$P\{X > x\} = \begin{cases} \frac{A}{x^{\alpha}} & \text{if } x \ge 1, \\ 0 & \text{otherwise,} \end{cases}$$

for some constant A > 0. If $0 < \alpha \le 1$, then the mean of X is infinite. In probability theory, the Pareto distribution has been useful to motivate the generalization of the central limit theorem in §8.5.

Recall that if we let $S_n = \sum_{i=1}^n X_i$, where X_i , i = 1, ..., n are random variables that are independent and identically distributed, and if $\sigma^2 \equiv \sigma^2(X_i) < \infty$ and $\mu \equiv E(X_i) < \infty$, then

$$Y_n = \frac{S_n - n\mu}{\sigma n^{1/2}}$$

has approximately a unit (or standard) normal distribution when n is large enough. Namely, for n large enough we have

$$P\{Y_n > x\} \sim \int_x^\infty \frac{e^{-y^2/2}}{\sqrt{2\pi}} \, dy$$

But what happens if μ or σ^2 is not defined or infinite? A different way of "normalizing" the sums S_n is possible, so that, in certain cases, when n

is large, their distributions are found to be asymptotically (i.e., for large |x|) of Pareto type.

Let us introduce a bit of notation first. Define F(x) as the distribution function of X_i and let

$$Y_n = \frac{1}{B_n} S_n - A_n \tag{A.4.1}$$

so that, when $\mu < \infty$ and $\sigma < \infty$, $B_n = \sigma n^{1/2}$ and $A_n = (\mu/\sigma) n^{1/2}$ for the central limit theorem. Define also $F_n(x)$ to be the distribution function of Y_n . A generalization to the central limit theorem is to find conditions on F, B_n and A_n under which $\lim_{n\to\infty} F_n(x)$ exists. For example, for the central limit theorem, F must be such that μ and σ are finite, and A_n and B_n as above. The proof of the solution to the general problem is beyond the level of this book, but we summarize below the main results.

Theorem 1. (cf. theorem of §33, p. 162 in [Gnedenko-Kolmogorov]) When it exists, the limiting distribution of Y_n , defined by $\Phi(x) = \lim_{n \to \infty} F_n(x)$ must be a stable law.

Recall that by Lévy's theorem cited in §9.7, a stable distribution family is identified by the exponent α of its characteristic function, where $0 < \alpha \le 2$ (see (9.7.3)). This property is exploited in the following two theorems.

Theorem 2. (cf. theorem 5, p. 181 in [Gnedenko-Kolmogorov]) If the distribution function F(x) of X_i satisfies for large |x|:

$$F(x) \sim \frac{A}{|x|^{\alpha}}$$
 if $x < 0$,

$$1 - F(x) \sim \frac{B}{|x|^{\alpha}} \text{ if } x > 0,$$

for some positive constants A, B and $\alpha \in (0,2)$, then $F_n(x)$ converges to a stable law with characteristic function exponent α .

Theorem 3. (cf. Theorem 1, p. 172 in [Gnedenko-Kolmogorov]). In order for $F_n(x)$ to converge to the unit normal distribution function (i.e., with $\alpha = 2$), it is sufficient that $Var(X_i)$ be finite.

The above sufficient conditions are in fact also necessary if we want to impose that B_n in (A.4.1) be of the form $an^{1/\alpha}$, where a is a positive constant. The reference theorems above in [Gnedenko-Kolmogorov] actually address these necessary and sufficient conditions. In the central limit

theorem ($\alpha=2$ and finite variance) we have $a=\sqrt{Var(X_i)}$, so that the limiting distribution is the *unit* normal distribution. More generally, one way to justify the need for B_n to be of the form $an^{1/\alpha}$ is through the following theorem.

Theorem 4. Let $\{X_i\}$ be independent, following the same stable distribution with characteristic function $\varphi(\theta) = e^{-\gamma_{\alpha}|\theta|^{\alpha}+id\theta}$. Then, for a given positive a, $Y_n = (1/a)n^{-1/\alpha}\sum_{k=1}^n (X_k - d)$ has a stable distribution with characteristic exponent α .

Proof: For $1 \le k \le n$, let $X'_k = (X_k - d)/(an^{1/\alpha})$. Its characteristic function is

$$\varphi'(\theta) = E\left[e^{i\theta X_k'}\right]$$

$$= E\left[e^{i\theta \frac{X_k - d}{an^{1/\alpha}}}\right]$$

$$= e^{-i\frac{\theta d}{an^{1/\alpha}}}E\left[e^{i\frac{\theta}{an^{1/\alpha}}X_k}\right]$$

$$= e^{-i\frac{\theta d}{an^{1/\alpha}}}\varphi\left(\frac{\theta}{an^{1/\alpha}}\right).$$

Given that the random variables $\{X_k\}$, and thus $\{X'_k\}$, are independent and identically distributed, Theorem 7 in §6.5 enables us to write the characteristic function of Y_n as:

$$\begin{split} E\left[e^{i\theta Y_n}\right] &= E\left[e^{i\theta \sum_{k=1}^n X_k'}\right] \\ &= \prod_{k=1}^n \varphi'(\theta) \\ &= e^{-in\frac{\theta d}{an^{1/\alpha}}} \left[\varphi\left(\frac{\theta}{an^{1/\alpha}}\right)\right]^n \\ &= e^{-\frac{\gamma_\alpha}{a}|\theta|^\alpha} \end{split}$$

Through this theorem we ensure that the sum of (properly scaled and shifted) independent random variables that follow an identical stable law converges to the same type.

We conclude by observing that when $0 < \alpha < 2$, the Pareto-type behavior that is manifested under the conditions of Theorem 2 above is also present in the limiting stable distribution. This property was determined by Lévy (cf. (37) on p. 201 of [Lévy] and §36 in [Gnedenko-Kolmogorov]). It is hardly surprising as Theorem 4 above shows that if F is a stable type with characteristic exponent α , then Φ must be of the same type.

When $\alpha=2$ (i.e., in the central limit theorem) the property above cannot hold. For a random variable Z with unit normal distribution we have, for $z \geq 1$:

$$\begin{split} P\left\{|Z|>z\right\} &= \int_z^\infty \frac{e^{-\zeta^2/2}}{\sqrt{2\pi}} \, d\zeta \\ &\leq \int_z^\infty \zeta \frac{e^{-\zeta^2/2}}{\sqrt{2\pi}} \, d\zeta \\ &= \frac{1}{\sqrt{2\pi}} e^{-z^2/2}. \end{split}$$

This, by the way, means that the tail probability is at least "square–exponential". It decreases to zero as $x \to \infty$ infinitely faster than the exponential e^{-z} , and even more so than the power $x^{-\alpha}$ ($\alpha > 0$) as in Pareto's law. This demonstrates that the upper bound on $P\{|Z| > z\}$ given by the Chebyshev inequality (see (7.6.5)) is quite loose in this case.

Final note: we have drawn the material on stable laws mostly from [Lévy] and [Gnedenko–Kolmogorov]. However, similar results with alternative proofs are also given in [Feller 2].

10

Option Pricing Theory

The previous chapter treats one-period models that form the foundation of the equilibrium pricing of primary securities such as stocks. In this chapter, we focus on the so called derivative securities or contingent claims. These are assets whose prices depend on those of primary securities. In this chapter we will illustrate the application of some of the most advanced material on stochastic processes presented in this book. The ideas presented here form the basis of many developments in the field of mathematical finance which have had a profound impact on both theory and practice.

10.1. Options basics

An option is a contract giving its holder the right, without the obligation, to either buy or sell a security such as a stock (the underlying security), at a predetermined price (the exercise or strike price). If the right granted is to buy, then the option is labeled a call. If the right granted is to sell, the option is labeled a put. This type of contract comes with an expiration or maturity date, which may be infinite. Options that are exercisable only at expiration are called European, and those that can be exercised any time up to expiration are called American. Both American and European option types are traded worldwide, but American options are more common. Options are the most representative type of contingent claims. For standard comprehensive texts that include actual contract transactions and market descriptions, we refer to [Cox and Rubinstein] and [Hull].

Example 1. Suppose today you hold a contract giving you the right to buy one share of General Electric stock (the underlying security) at \$30 any time during the next three months. Then you are holding an American call option on General Electric stock with a strike price of \$30 and a maturity of 3 months.

Example 2. If you hold a contract giving you the right to sell one Chinese Renminbi at 16 Japanese Yen exactly six months from today then you are holding a European currency put on the exchange rate between Chinese Renminbi and Japanese Yen with a strike price of 16 Japanese Yen and a maturity of 6 months.

In both of these examples the price at which the underlying security will be bought or sold is fixed and known in advance. These types of options are called standard options, in contrast to so called "exotic" options, where more generally the rule determining the exercise price is fixed but not known in advance. For example, you may hold a contract giving you the right to buy one share of General Electric stock at the average price it will be selling for between now and the moment you exercise the option within the next three months. In the remainder of this chapter our focus will be on standard options of either European or American type. To simplify the exposition, we will deal only with options where the underlying security is a stock that pays no dividend, unless noted otherwise.

Option contracts can be used for several purposes. Prices of standard options are significantly lower than those of their underlying securities. For speculators, these lower prices give them an opportunity to potentially benefit from favorable changes in the stock prices without committing as much as the full stock price initially. Option contracts can also be viewed as insurance contracts as they offer their holders protection against adverse movement in the underlying security. For example, an investor who currently owns stock of a company and who is concerned about a possible significant drop in the value of the stock in the next three months may want to purchase a put on the stock in order to guarantee a minimum price at which he or she will sell this stock over the next three months.

In order to get the right to exercise an option, its holder must first pay a premium at the time the contract is established. This price, also known as the value of the option at inception, is the focus of option pricing theory. In the case of American options, this theory allows us to also determine the optimal exercise time, i.e., the best time to exercise the option.

Option Payoff. A standard option is characterized by its exercise price K (\$30 in the case of example 1 above), its expiration date T, expressed in years (so that it is .25 year in the case of example 1), and whether it is a call or a put. The price S_T of the underlying security T units of time from now is a random variable; i.e., there exists a sample space Ω such that

the value of the underlying security at T is $S_T(\omega)$ if $\omega \in \Omega$ is drawn. Fast forward to time T and assume that a particular $\tilde{\omega}$ is sampled. Then, in the case of a call option for example, the option is exercised only if $S_T(\tilde{\omega}) > K$. This makes sense since there is no reason to buy a security at price K if it is selling at a price $S_T(\tilde{\omega}) \leq K$ [by convention, we assume that when $S_T(\tilde{\omega}) = K$ we are not exercising the option]. So, assuming $S_T(\tilde{\omega}) > K$, the holder of the option with strike K can exercise this option, i.e., buy at K dollars per share, and then sell the shares back on the market at $S_T(\tilde{\omega})$ to realize a profit of $S_T(\tilde{\omega}) - K$ per share. If $S_T(\tilde{\omega}) \leq K$, the call holder does not exercise the option, thus realizing no profit. In summary we have the following.

Definition 1. For a given T, let S_T be the price of a stock at T. Consider a standard European option on this stock with exercise price K and expiration date T. Then the payoff of this option upon exercise is the random variable $g = \max(S_T - K, 0)$, in the case of a call, and $g = \max(K - S_T, 0)$, in the case of a put.

From a speculator's or gambler's point of view, the option price is the amount that she or he is willing to pay at time 0 in order to collect the option payoff T units of time later.

Generally, options are considered a less expensive way to benefit from the price changes of an asset without having to commit to the full price of the asset up front. For example, on March 8, 2002, IBM stock was trading around \$105 per share, while it cost only about \$11 to buy a call option on IBM stock with strike at \$100 and which expired in July 2002. On the other hand, on the same day, it cost approximately \$6 to buy an IBM put with strike at \$100, with the same expiration month.

We should also remember that options involve two parties to a contract: the option buyer (or holder) and the option writer (or seller), with opposing views about the future performance of the underlying security. Often the writer of a call option does not own the underlying security, hoping that the buyer of the option would never exercise the option. Throughout history, there have been some phenomenal episodes where speculation about prices of the underlying went wild and major crashes ensued as a result. A famous example is that of "tulipmania" in the Netherlands in the 1630s. Tulips were cultivated originally in Turkey and were brought to Holland in the 16th century. By the 17th century, many new breeds had been created through horticultural experiments. These exotic and beautiful flowers became expensive, and thus a status symbol. As they became more desirable, people realized they could make significant profits by speculating on their prices. People with limited means were buying and selling tulip bulbs. To protect themselves against endlessly rising prices, many speculators bought call options (with high strikes) from other speculators who did not believe that they would be called on to deliver bulbs at these high prices. Eventually, option writers failed to honor their contracts. Many could no longer afford the high prices anymore and the market crashed.

In recent history, similar phenomena involving speculation and the use of derivatives occured. In the 1990s particularly, aggressive bets on currency or interest rates, made by investors exposing their institutions, led to the collapse of major organizations. Barings, one of the oldest banks, which played a major role in the expansion of the British Empire; Orange County, one of the largest in California; and Long-Term Capital Management, an asset management company which used the services of two Nobel laureates in economics and which invested for very wealthy individuals, large financial institutions, and universities*; all defaulted on their contracts and went bankrupt or had to be rescued in a dramatic fashion.

Discounting. In the course of pricing options we will need to account for cash values at different times. A common theme in finance is that of the "time value of money"; e.g., in the case of price inflation, where an object that can be purchased at one dollar today will have to be purchased at more than a dollar say a year from now. In other words, the "value" of a dollar today may be different than that of a dollar a year from now. To allow for a fair comparison between cash values at different times, financial economists have devised the concept of present-value discounting using a discount rate, which in our case will simply be the rate of return r of the riskless security introduced in §9.2. For our purposes, there exists only one riskless security in our economy and we shall discount relative to time 0 over discrete dates. If we hold an asset with cash value C_n at time t_n , its time 0 discounted value is $C_n/(1+r)^n$.

Submartingales and Supermartingales. As we shall see shortly, option pricing is closely related to taking expectations of random variables. The martingales introduced in Appendix 3 illustrate the concept of a fair game, i.e., the odds are neither in favor nor against the gambler. There are, however, situations where the odds, defined by the probability under which the conditional expectation such as (A.3.1) is taken, are either in favor or against the gambler. Using the same notation as in Appendix 3, the first case is expressed as

$$E(X_{n+1}|X_0, X_1, \dots, X_n) \ge X_n, \tag{10.1.1}$$

and the second as

$$E(X_{n+1}|X_0, X_1, \dots, X_n) \le X_n. \tag{10.1.2}$$

^{*}In February 2000, the PBS television network broadcast a documentary with the principal actors involved and many industry observers. As of the writing of this book, a transcript of this program ("Trillion Dollar Bet") is available online at www.pbs.org/wgbh/nova/stockmarket.

A process $\{X_n\}$ is called a *submartingale* if it satisfies (10.1.1), and a *supermartingale* if (10.1.2) is true. There is an unfortunate verbal discordance between these names and their meanings. In general, a word qualified with *super* is associated with an outcome that is desirable. Here "supermartingale" means a losing game.

Theorem 1. Under the same conditions of the theorem in Appendix 3, we have for a finite optional time T

$$E\left(X_{T}\right) \geq E\left(X_{0}\right)$$
 for a submartingale and $E\left(X_{T}\right) \leq E\left(X_{0}\right)$ for a supermartingale.

For any supermartingale (respectively, submartingale) the same result holds as long as T is bounded.

Proof: The proof is exactly as in Appendix 3 with equalities replaced with the corresponding inequalities (left as an exercise).

We now illustrate the reach of this theorem with the following.

An Example of Greed. Consider two persons with same initial wealth X_0 . The first one is timidly greedy and says that she will quit investing as soon as she is ahead by one unit (you can pick your unit of account as \$1 or \$10000). The second person is more aggressive. He would like to quit as soon as his wealth is 10 times his initial amount. For the first person define $T = \min\{n : X_n > X_0 + 1\}$, where X_n is this investment value in period n. For the second individual, define $T = \min\{n : X_n > 10X_0\}$. T is not bounded. By our own nature, we all have a limited number of years to live (some like the Queen mother of the United Kingdom can get as many as 101 years). Let's call the upper bound on the number of years we can possibly live \bar{t} . Define $\bar{T} = T \wedge \bar{t}$; then \bar{T} is bounded. If the wealth process X_n of these investors is a supermartingale, then we have $E(X_{\bar{T}}) \leq E(X_0)$. In either case he/she is expected to lose! All we can say in probability is expectation (hope). Now people will retort: "But I know Mike bought Enron at 8.52 and sold at 88.87!" Lord Keynes had the answer long ago, cited on page 135. He wrote that in his doctorate in philosophy before delving in "Money" (title of his famous treatise in economics). His philosophical treatise is listed in the references.

We can even tell these investors their chances of making their goals. For a given b positive, by Chebyshev's inequality – see Exercise 21 of §7.6, analogue of (7.6.5) – we have

$$P\left(X_{\bar{T}} > b\right) \le \frac{E\left(X_{\bar{T}}\right)}{b}.$$

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From Theorem 1 above, we have $E(X_{\bar{T}}) \leq E(X_0)$. Therefore

$$P\{X_{\bar{T}} > b\} \le E(X_0)/b.$$

Take, for example, X_0 to be a constant C and $b = a \times C$, for some a > 0. Then

$$P\left\{X_{\bar{T}} > aC\right\} \le 1/a.$$

With a = 10, for example, our aggressive investor has less than 10% chance of reaching his goal.

Stock Price Evolution and Expectations. If you follow the evolution of a particular stock price through media such as newspapers, radio, television, or the Internet, you learn that it is "up" or "down," or that it is likely to go up or down. This description may remind you of a random walk (cf. §8.1 for a definition). In fact we shall rely upon the framework introduced in §9.2 to formalize this observation.

Recalling our discussion about discounting above, if we needed to compare two stock values S_n , at time t_n , and S_m at time t_m , we would need to look at their discounted values. So let

$$Y_n = \frac{1}{(1+r)^n} S_n$$
 for $n = 0, 1, 2, \dots, N$.

Similarly, any expectation about the prospects of a stock would have to be done on the basis of the discounted values $\{Y_n\}$.

Recall that one of the main reasons investors are interested in options is because of their expectations about the underlying securities. They may view an option as insurance, for example, should they be interested in the future purchase of a stock. If they expect a stock price to go up, then they would lock in the purchase price (exercise price) today by acquiring an option. For this coverage, they pay an insurance premium (option price) for the right to use the locked-in price later.

In the spirit of the remainder of this chapter, assume that the oneperiod returns $\{r_n\}$, $n=1,2,\ldots,N$, are identically and independently distributed. Then following the notation introduced in Appendix 3, we write the conditional expectation of Y_n given all the history up to time n-1 as

$$E[Y_n|Y_0, Y_1, \dots, Y_{n-1}] = E[Y_n|Y_{n-1}]$$

$$= E\left[\frac{1}{(1+r)^n} S_n | S_{n-1}\right]$$

$$= \frac{1}{(1+r)^n} E[S_n | S_{n-1}]$$

$$= \frac{1}{(1+r)^n} E[(1+r_n) S_{n-1} | S_{n-1}]$$

$$= \frac{1}{(1+r)^n} [1+E(r_n)] S_{n-1}$$

$$= \frac{1+E(r_n)}{1+r} \frac{S_{n-1}}{(1+r)^{n-1}}.$$

Therefore

$$E[Y_n|Y_0, Y_1, \dots, Y_{n-1}] = \frac{1 + E(r_n)}{1 + r} Y_{n-1},$$

indicating that $\{Y_n\}$, the discounted stock price process, is a martingale, submartingale, or supermartingale whether s $E(r_n) = r$, $E(r_n) \ge r$, or $E(r_n) \le r$, respectively.

Discrete and Continuous Models. To price an option we need to know the evolution of the underlying security as a stochastic process. So far in this book we have seen two kinds of stochastic processes: those that move only at discrete times, e.g., a random walk as in §8.1, and those that evolve continuously, e.g., a Brownian motion as in §8.2. Historically the major breakthrough in pricing options happened in the context of the harder case of continuous stochastic processes. It resulted in a neat formula (the famous Black-Scholes formula) that we will present shortly. As an indication of its impact on the field of economics, both Myron Scholes and Robert Merton were awarded the Nobel Prize for this formula and its fundamental extension (Fisher Black had passed away at the time of the award decision in 1997). On the other hand, the discrete-time model, which we will study in detail, is easier to understand but lacks the elegance of a closedform formula. In this chapter we will show an approach to approximate with the discrete-time model the closed-form expressions obtained via the continuous-time model.

Black–Scholes Formula. Consider a stock with instantaneous mean return rate μ and instantaneous return rate variance σ^2 , a market where the

instantaneous return rate of the riskless asset is r, and a call with expiration date T given. We assume that the stock does not pay a dividend. [Black and Scholes] have derived prices C_0 and P_0 at time 0 for, respectively, a European call and a European put written on this stock as follows:

$$C_0 = S_0 \Phi(d_1) - K e^{-rT} \Phi(d_2)$$

and

$$P_0 = Ke^{-rT}\Phi(-d_2) - S_0\Phi(-d_1),$$

where $d_1 = \left[\log(S_0/K) + (r + \sigma^2/2)T\right]/\sigma\sqrt{T}$, $d_2 = d_1 - \sigma\sqrt{T}$, and Φ is the standard normal cumulative distribution function. Notice that the mean instantaneous rate of return appears in neither formula. We will show in the discrete-time setting that this situation is related to the concept of arbitrage-free pricing, to be defined soon, and the associated concept of equivalent pricing probability.

10.2. Arbitrage-free pricing: 1-period model

The one-period model is the building block for the multi-period model in the discrete-time context of option pricing. To determine the option premium, we need to know the nature of the stochastic behavior of the underlying stock. In the simple one-period pricing setting, our sample space Ω consists of only two elements H and T, such that $P\{H\} = p$ and $P\{T\} = 1 - p$, $p \in (0,1)$. Let S_0 be the initial price of the underlying stock of the option $[S_0$ is a constant random variable: $S_0(H) = S_0(T)$. If $\omega = H$ is sampled, then the price $S_1(H)$ of this stock at date 1 is uS_0 , and if $\omega = T$ is sampled, then $S_1(T) = dS_0$, where u and d are given such that

$$0 < d < 1 + r < u, (10.2.1)$$

where $r \geq 0$ is the one-period rate of return of the riskless security. Recall that the one-period rate of return of the stock is $(S_1 - S_0)/S_0$. Thus

$$\frac{S_1(\omega) - S_0}{S_0} = \left\{ \begin{array}{ll} u - 1 & \text{if } \omega = H, \\ d - 1 & \text{if } \omega = T. \end{array} \right.$$

We can rewrite the last two inequalities in (10.2.1) as

$$d-1 < r < u-1$$
.

With the interpretation that u-1 and d-1 are, respectively, the favorable (i.e., we have a gain) and unfavorable rates of returns on the stock, we see that the rate of return r of the riskless asset must satisfy (10.2.1).

Otherwise, if $d \geq 1 + r$, then $(S_1(\omega) - S_0)/S_0 \geq r \ \forall \omega \in \Omega$, indicating that the return of the stock is always as good as that of the riskless asset. In this situation we have no reason to invest in the riskless asset; i.e., there is no need for this asset to exist at all. Similarly, if $1 + r \geq u$, then $(S_1(\omega) - S_0)/S_0 \leq r \ \forall \omega \in \Omega$, making the investment in the stock unattractive.

Let V_0 and V_1 be the values of a European option at times 0 and 1, respectively. As seen before, V_1 is known to be the payoff g at the expiration date 1. To determine the option price V_0 at time 0, we consider an investment environment where we can invest in a market that consists of only the following assets: the option, its underlying stock, and the riskless security, which we label bond as is generally the case in standard presentations on the subject. A portfolio will therefore consist of a triple (α, β, γ) , where, respectively, α is the number of options, β the number of stock shares, and γ the number of bond shares held. Define also B_n and S_n to be, respectively, the bond and stock prices at n = 0, 1.

Arbitrage-free markets. It is clear that we cannot have a trading strategy (α, β, γ) such that, starting with an initial wealth $W_0 = \alpha V_0 + \beta S_0 + \gamma B_0 = 0$, we end up at time 1 with wealth $W_1 = \alpha V_1 + \beta S_0 + \gamma B_1$ such that $W_1(\omega) > 0$ for all $\omega \in \Omega$; i.e., starting with zero wealth we are guaranteed positive wealth at time 1. This strategy is economically untenable, for if it existed it would soon be discovered by many traders and, through their desire to hold exactly (α, β, γ) , would bid up or down the time-0 prices of either the option, the stock, or the bond so that $W_0 = 0$ would soon be violated. A more rigorous argument using so-called separation theorems can be articulated (cf. [Duffie]). In fact, we define an arbitrage opportunity or portfolio to be any strategy (α, β, γ) such that

- (i) $\alpha V_0 + \beta S_0 + \gamma B_0 = 0$,
- (ii) $\alpha V_1(\omega) + \beta S_1(\omega) + \gamma B_1(\omega) \ge 0$ for all $\omega \in \Omega$, and
- (iii) $E\{\alpha V_1 + \beta S_1 + \gamma B_1\} > 0.$

As we shall see shortly, there can be arbitrage opportunities such that (ii) is satisfied strongly; i.e., $\alpha V_1(\omega) + \beta S_1(\omega) + \gamma B_1(\omega) > 0$ for all $\omega \in \Omega$. This means that the strategy (α, β, γ) is a "sure win": starting with no wealth, you are guaranteed a positive wealth no matter which ω is sampled.

What is arbitrage? Most finance books describe immediately the *mechanics* of arbitrage without dwelling too much on the meaning of the word itself. If you look it up in the dictionary, you might be left with a sense of incomplete definition. We have found one source [Weisweiller], particularly its Chapter 1, where the author traces the subtle evolution of the meaning of this word through several editions of dictionaries. The concept of arbitrage as defined here is in fact an extension of its original meaning in the context of finance, where it refers to a situation such that a person buys an asset in one market (e.g., the New York Stock Exchange) and sells it immediately in another market (e.g. the Paris Bourse) where it is trading

at a higher price. The Oxford English dictionary follows this interpretation while noticing its original French root, namely the verb "arbitrer," which means to arbiter or to judge. In this sense, arbitrage describes a person (a speculator really) who is evaluating an asset, or judging its value, to identify mispricing, i.e., the difference between what the asset is selling for and what it should sell for. An arbitrage-free market is then defined as one where no arbitrage opportunities exist. It is this type of market that we shall assume henceforth.

Lemma. Let $\Omega = \{H, T\}$. There exists a portfolio $(\alpha_0, \beta_0, \gamma_0)$ such that $\alpha_0 = 0$ and

$$\beta_0 S_1(\omega) + \gamma_0 B_1(\omega) = g(\omega)$$
 for all $\omega \in \Omega$.

The proof is obvious, as it simply requires us to find β_0 and γ_0 as solutions to a system of two linear equations. Since $\Omega = \{H, T\}$, β_0 and γ_0 must satisfy

$$\beta_0 S_1(H) + \gamma_0 B_1(H) = g(H)$$
, and $\beta_0 S_1(T) + \gamma_0 B_1(T) = g(T)$,

or equivalently

$$\beta_0 u S_0 + \gamma_0 (1+r) B_0 = g(H)$$
, and $\beta_0 d S_0 + \gamma_0 (1+r) B_0 = g(T)$,

from which we deduce

$$\beta_0 = \frac{g(H) - g(T)}{(u - d)S_0},$$

$$\gamma_0 = \frac{1}{(1 + r)B_0} \left[\frac{ug(T) - dg(H)}{u - d} \right].$$
(10.2.2)

This lemma shows that starting with a portfolio consisting of only the stock and the bond (no option) we can replicate the value of the option at maturity. This strategy is labeled a replicating strategy. The next proposition shows that the price of the option V_0 has to be the value of this portfolio $(\beta_0 S_0 + \gamma_0 B_0)$ at time 0.

Proposition 1. The time-0 price of a European option on a stock with price S and payoff g at time 1 is

$$V_0 = \frac{1}{1+r} \left[\tilde{p}g(H) + (1-\tilde{p})g(T) \right], \qquad (10.2.3)$$

where

$$\tilde{p} = \frac{1 + r - d}{u - d}. (10.2.4)$$

Proof: We first show that $V_0 = \beta_0 S_0 + \gamma_0 B_0$, where β_0 and γ_0 are defined by (10.2.2). Suppose $\epsilon = V_0 - (\beta_0 S_0 + \gamma_0 B_0) > 0$. Consider the following portfolio $(-1, \beta_0, \gamma_0 + \epsilon/B_0)$. At time 0 the value of this portfolio is

$$W_0 = -V_0 + \beta_0 S_0 + \left(\gamma_0 + \frac{\epsilon}{B_0}\right) B_0 = 0.$$

At time 1 its value is

$$W_{1} = -V_{1} + \beta_{0}S_{1} + \left(\gamma_{0} + \frac{\epsilon}{B_{0}}\right)B_{1}$$
$$= -V_{1} + \beta_{0}S_{1} + \gamma_{0}B_{1} + \epsilon \frac{B_{1}}{B_{0}}.$$

Recalling that $V_1 = g$ and using the lemma above, we have

$$-V_1(\omega) + \beta_0 S_1(\omega) + \gamma_0 B_1(\omega) = 0$$
 for all $\omega \in \Omega$.

Since B_0 and B_1 are positive, we have $W_1(\omega) = \epsilon(B_1/B_0) > 0$ for all $\omega \in \Omega$, thus showing that $(-1, \beta_0, \gamma_0 + \epsilon/B_0)$ is an arbitrage portfolio. A similar proof applies to the case $\epsilon = V_0 - (\beta_0 S_0 + \gamma_0 B_0) < 0$ with portfolio $(1, -\beta_0, -\gamma_0 - \epsilon/B_0)$.

Using (10.2.2) we now write the value of the option as

$$\begin{split} V_0 &= \beta_0 S_0 + \gamma_0 B_0 \\ &= \frac{g(H) - g(T)}{u - d} + \frac{1}{1 + r} \left(\frac{ug(T) - dg(H)}{u - d} \right) \\ &= \frac{1}{1 + r} \left[\frac{1 + r - d}{u - d} g(H) + \frac{u - (1 + r)}{u - d} g(T) \right], \end{split}$$

where we recall

$$g(H) = \begin{cases} \max(uS_0 - K, 0) & \text{for a call,} \\ \max(K - uS_0, 0) & \text{for a put, and} \end{cases}$$

$$g(T) = \begin{cases} \max(dS_0 - K, 0) & \text{for a call,} \\ \max(K - dS_0, 0) & \text{for a put.} \end{cases}$$

Remarks.

- 1. With condition (10.2.1) we have $0 < \tilde{p} < 1$, and thus $\tilde{p}g(H) + (1 \tilde{p})g(T)$ can be interpreted as the expected value of the random variable g under the probability \tilde{P} defined on $\Omega = \{H, T\}$ such that $\tilde{P}\{H\} = \tilde{p}$ and $\tilde{P}\{T\} = 1 \tilde{p}$ (since this argument will show up again when $\Omega = \{0,1\} \times \{0,1\} \times \cdots \times \{0,1\}$, you may want to consult §2.4 now to check how we can construct probabilities on countable spaces). The probability \tilde{P} is often referred to as the risk-neutral probability; however, we simply call it the pricing probability. Thus, in the one-period model the price at time 0 of a European option with payoff g at time 1 is the discounted conditional expected value of this payoff under the pricing probability. It is a conditional expectation because both g(T) and g(H) depend on information at time 0, namely S_0 .
- 2. The pricing probability \tilde{P} is said to be *equivalent* to P. This property is formally defined as: $P(\omega) > 0$ if and only if $\tilde{P}(\omega) > 0$, for all $\omega \in \Omega$. It means that P and \tilde{P} agree on the sample points with nonzero probability. In the current setting of $\Omega = \{H, T\}$, this property seems trivial, as $0 and <math>0 < \tilde{p} < 1$ imply $P\{H\} = p > 0$, $P\{T\} = 1 p > 0$, $\tilde{P}\{H\} = \tilde{p} > 0$, and $\tilde{P}\{T\} = 1 \tilde{p} > 0$. But for more abstract Ω you may not be able to see it as easily.
- 3. The price V_0 is independent of the original probability defined on Ω . In fact, under \tilde{P} we have the following conditional expectation:

$$\tilde{E}[S_1|S_0] = \tilde{p}uS_0 + (1-\tilde{p})dS_0
= \frac{1+r-d}{u-d}uS_0 + \frac{u-(1+r)}{u-d}dS_0
= (1+r)S_0,$$

which we can rewrite as

$$\tilde{E}[Y_1|Y_0] = Y_0, \tag{10.2.5}$$

where we define

$$Y_k = \frac{1}{(1+r)^k} S_k, \quad k = 0, 1.$$

Relation (10.2.5) states that under the pricing probability \tilde{P} , the discounted stock price $\{Y_k\}_{k=0}^1$ is a martingale in this rather trivial case of one-period; in contrast to the original probability, under which the discounted stock price process could be a submartingale or a supermartingale, depending on the expected one-period returns. In fact, under the pricing probability, all the discounted price processes, including the bond and the option, are martingales. For the

bond, this result is obvious since its price process is nonrandom, and for the option we notice that we can rewrite (10.2.3) as

$$\frac{1}{(1+r)^0}V_0 = \frac{1}{1+r}\tilde{E}[V_1|V_0].$$

Indeed, by definition, $V_1(\omega) = g(\omega)$ for all $\omega \in \Omega$, and $g(\omega)$ is a function of S_0 and ω , which by the relation $V_0 = \beta_0 S_0 + \gamma_0 B_0$ implies that $\tilde{E}[V_1|S_0] = \tilde{E}[V_1|V_0]$.

4. Numéraire Invariance Principle

In our arbitrage pricing we presented an argument where prices are evaluated in (units of) dollars. This way, cash is what is called the unit of account or *numéraire*. In other words, we treated cash as a security whose price is always 1. In fact, it turns out that we could have used any other security with prices that are always positive. For example, with the stock as a numéraire, we define new prices for the bond as

$$\bar{B}_0 = B_0/S_0$$
 and $\bar{B}_n = B_n/S_n$,

with the assumption $S_0 > 0$ and $S_n(\omega) > 0$ for all n and $\omega \in \Omega$. Of course, $\bar{S}_n \equiv 1$.

Using stock as numéraire, we can rewrite (10.2.3) as

$$\frac{V_0}{S_0} = \frac{1}{(1+r)S_0} \left[\tilde{p}S_1(H) \frac{g(H)}{S_1(H)} + (1-\tilde{p})S_1(T) \frac{g(T)}{S_1(T)} \right]$$

or

$$\bar{V}_0 = \left[\tilde{p} \frac{S_1(H)}{(1+r)S_0} \bar{g}(H) + (1-\tilde{p}) \frac{S_1(T)}{(1+r)S_0} \bar{g}(T) \right],$$

where $\bar{V}_0 = V_0/S_0$ and $\bar{g}(\omega) = g(\omega)/S_1(\omega)$, for $\omega \in \{H, T\}$. Thus, with the stock price as numéraire, we can define yet another probability \bar{P} , equivalent to P, by:

$$\bar{P}(H) = \tilde{p} \frac{S_1(H)}{(1+r)S_0}, \quad \text{and}$$

$$\bar{P}(T) = (1 - \tilde{p}) \frac{S_1(T)}{(1+r)S_0}$$

so that $\bar{V}_0 = [\bar{P}(H)\bar{g}(H) + \bar{P}(T)\bar{g}(T)]$, which expresses the price of the option in this unit of account as yet another expectation (as an exercise, verify that $\bar{P}(H) + \bar{P}(T) = 1$). In fact, one can show that

there is no arbitrage in this new market (with prices \bar{S}_n and \bar{B}_n) if and only if there is no arbitrage in the original market (with prices S_n and B_n). This one-period result generalizes to multiple periods (try it as an exercise and see [Duffie], for example, for the continuous case).

10.3. Arbitrage-free pricing: N-period model

For the general case of $N \geq 1$ periods: the arbitrage-free argument used for one period will be applied to each of the N periods, and the properties listed in the remarks at the end of the last section will hold as well. However, because the previous analysis was limited to one period only, it didn't allow us to illustrate two other concepts that are important in the general case and that connect adjacent periods: (i) dynamic replication, and (ii) self-financing strategies, which are described below.

The holder of an option does not have to keep the option until exercise. He or she can trade it on an exchange (e.g., the Chicago Board Options Exchange) and therefore the next holder of this option will have to face the same problem regarding its value: i.e., evaluating its fair price at the time of purchase. As a result, in option pricing theory we are in fact interested in the value of the option at all times. For multiple periods we shall exhibit a strategy that replicates the value of the option at every discrete date. which in the one-period case is reduced to the origin and the exercise date. This process is called *dynamic replication* in the multi-period situation. It is predicated upon the constraint that this strategy be self-financing: at every date it will neither generate nor require external cash. For example, suppose you are in period n (i.e., in the interval $(t_{n-1}, t_n]$) and that you hold one share of stock and two of bond. If after observing the prices of these two assets at time t_n you decide to hold .8 share of stock and 3 of bond for ensuing period $(t_n, t_{n+1}]$ it must be the case that the purchase of the additional share of bond uses exactly the cash amount generated by the sale of .2 share of stock (we are of course ignoring the pesky problems associated with transaction costs, taxes, etc.). In other words, the amount that is generated by the sale of one asset finances exactly the additional purchase of the other security.

For N periods the stock returns for two distinct periods are assumed to be independent and identically distributed. In the simplest case, which generalizes the one-period Bernoulli model we have just studied, the N-period model is identical to the coin-tossing scheme that we have studied in this book (cf. Example 8 in §2.4, §3.2 and Example 9 in §4.4). For the one-period model we could have equally used the sample space $\Omega = \{0,1\}$ instead of $\{H,T\}$. To make for a neat presentation we shall, as in the coin tossing example, adopt the sample space $\Omega = \{0,1\}^N$ for the N-period model. Thus a sample point (also called sample path) $\omega \in \Omega$ is of the form

 $(\omega_1, \omega_2, \dots, \omega_N)$. In period $n \leq N$, price S_n is observed at time t_n and depends only on the outcome values of $\hat{\omega}_n = (\omega_1, \omega_2, \dots, \omega_n)$. Therefore we can write $S_n(\omega) \equiv S_n(\hat{\omega}_n)$ such that

$$S_n(\hat{\omega}_n) = \begin{cases} uS_{n-1}(\hat{w}_{n-1}, 1) & \text{with probability } p, \\ dS_{n-1}(\hat{w}_{n-1}, 0) & \text{with probability } 1 - p, \end{cases}$$

where u, d, and p are defined in (10.2.1). The probability P on Ω is then defined by

$$P\{(\omega_1, \omega_2, \dots, \omega_N)\} = p^{\sum_{i=1}^N \omega_i} (1-p)^{N-\sum_{i=1}^N \omega_i},$$

which as we saw in the binomial case is the distribution of the total number of 1s.

At discrete times $0, 1, 2, \ldots, N$ (equivalently, actual times t_0, t_1, \ldots, t_N), let V_0, V_1, \ldots, V_N be the values of a European option with underlying stock prices S_0, S_1, \ldots, S_N and exercise price K. As in the one-period model, the value V_N of the option at expiration is a random variable that has an explicit form, namely:

$$V_N = \left\{ egin{array}{ll} \max(S_N - K, 0) & ext{ for a call, and} \\ \max(K - S_N, 0) & ext{ for a put.} \end{array}
ight.$$

We can even express V_N as a function of the sample path $\omega \in \Omega$:

$$V_N = \begin{cases} \max(u^{\sum_{i=1}^{N} \omega_i} d^{N - \sum_{i=1}^{N} \omega_i} S_0 - K, 0) & \text{for a call,} \\ \max(K - u^{\sum_{i=1}^{N} \omega_i} d^{N - \sum_{i=1}^{N} \omega_i} S_0, 0) & \text{for a put.} \end{cases}$$

When we have N periods, we determine the option values recursively by moving backward in time: first, given V_N we determine V_{N-1} , then V_{N-2} , etc. until we reach V_0 . Note that when we say "we determine V_n , $1 \le n \le N$," we do not necessarily mean that the value V_n is fixed; it will actually be in the form of a random variable, as is V_N (recall the case N=1 in the previous section). Therefore at any time $1 \le n \le N$, if the value V_n is known explicitly, then V_{n-1} is to be determined as in the one-period context. At (discrete) time n-1 (which is in period n) prices $S_0, S_1, \ldots, S_{n-1}$ and $S_0, S_1, \ldots, S_{n-1}$ have been observed. Presumably, $V_0, V_1, \ldots, V_{n-1}$ have also been observed, but the point of the pricing theory is to ensure that there are no arbitrage opportunities, in a manner identical to the one-period case.

N-Period Portfolio Strategies. The concept of no-arbitrage introduced for the one-period model extends to the N-period context, one period at a time.

Definition. An N-period portfolio strategy is a collection of triples $\{(\alpha_n, \beta_n, \gamma_n), n = 0, 1, \dots, N\}$ where for $0 \le n \le N$, α_n , β_n , and γ_n are random variables defined on $\Omega = \{0, 1\}^N$, e.g., for $\omega = (\omega_1, \omega_2, \dots, \omega_N) \in \Omega$, $\alpha_n(\omega) \in (-\infty, \infty)$.

For $0 \le n \le N$, we also require that a portfolio $(\alpha_n, \beta_n, \gamma_n)$ be determined only upon the information available up to period n. For example, we rule out "insider trading": in period n you are privy to some information that will be made public in period n+2, say. We have not formally defined information, but for our purposes, it suffices to say that all the information we need to determine a portfolio is the stock price history. As remarked earlier, a price S_n observed at time t_n depends only on the outcome values of $\hat{\omega}_n = (\omega_1, \omega_2, \dots, \omega_n)$. In a similar manner, a portfolio $(\alpha_n, \beta_n, \gamma_n)$ depends only on $\hat{\omega}_n$ at time t_n , e.g., we write $\alpha_n(\omega) \equiv \alpha_n(\hat{\omega}_n)$.

Lemma. In period n there exists a portfolio $(\alpha_{n-1}(\hat{\omega}_{n-1}), \beta_{n-1}(\hat{\omega}_{n-1}), \gamma_{n-1}(\hat{\omega}_{n-1}))$ such that $\alpha_{n-1}(\hat{\omega}_{n-1}) = 0$ and

$$\beta_{n-1}(\hat{\omega}_{n-1})S_n(\hat{\omega}_{n-1}, \omega_n) + \gamma_{n-1}(\hat{\omega}_{n-1})B_n(\hat{\omega}_{n-1}, \omega_n) = V_n(\hat{\omega}_{n-1}, \omega_n)$$
for $\omega_n \in \{0, 1\}$.

Proof: For ease of notation, we let $\alpha_{n-1} \equiv \alpha_{n-1}(\hat{\omega}_{n-1})$, $\beta_{n-1} \equiv \beta_{n-1}(\hat{\omega}_{n-1})$, $\gamma_{n-1} \equiv \gamma_{n-1}(\hat{\omega}_{n-1})$. As in Lemma 10.2, we now seek β_{n-1} and γ_{n-1} such that (recall that in the N-period model we substitute 0 for H and 1 for T):

$$\beta_{n-1}S_n(\hat{\omega}_{n-1},0) + \gamma_{n-1}B_n(\hat{\omega}_{n-1},0) = V_n(\hat{\omega}_{n-1},0), \text{ and}$$

$$\beta_{n-1}S_n(\hat{\omega}_{n-1},1) + \gamma_{n-1}B_n(\hat{\omega}_{n-1},1) = V_n(\hat{\omega}_{n-1},1).$$

Equivalently, this linear system of equations can be written as

$$\beta_{n-1}uS_{n-1}(\hat{\omega}_{n-1}) + \gamma_{n-1}(1+r)B_{n-1}(\hat{\omega}_{n-1}) = V_n(\hat{\omega}_{n-1}, 0), \text{ and}$$

$$\beta_{n-1}dS_{n-1}(\hat{\omega}_{n-1}) + \gamma_{n-1}(1+r)B_{n-1}(\hat{\omega}_{n-1}) = V_n(\hat{\omega}_{n-1}, 1),$$

therefore leading to the solution

$$\beta_{n-1} = \frac{V_n(\hat{\omega}_{n-1}, 0) - V_n(\hat{\omega}_{n-1}, 1)}{(u - d)S_{n-1}(\hat{\omega}_{n-1})},$$

$$\gamma_{n-1} = \frac{1}{(1+r)B_{n-1}(\hat{\omega}_{n-1})} \left[\frac{uV_n(\hat{\omega}_{n-1}, 1) - dV_n(\hat{\omega}_{n-1}, 0)}{u - d} \right] (10.3.1)$$

Proposition 2. The values V_0, V_1, \ldots, V_N of a European option expiring at time t_N satisfy the recursive relations

$$V_{n-1}(\hat{\omega}_{n-1}) = \frac{1}{1+r} \left[\tilde{p} V_n(\hat{\omega}_{n-1}, 0) + (1-\tilde{p}) V_n(\hat{\omega}_{n-1}, 1) \right],$$

for $1 \le n \le N$, where \tilde{p} s defined as in (10.2.4).

Proof: As in the proof of Proposition 1, we need to first show that $V_{n-1} = \beta_{n-1}S_{n-1} + \gamma_{n-1}B_{n-1}$, where β_{n-1} and γ_{n-1} are defined in (10.3.1). This step is straightforward and is left as an exercise.

Remark. As we just saw, for any period $1 \le n \le N$, an arbitrage-free pricing strategy satisfies

$$\beta_{n-1}S_n + \gamma_{n-1}B_n = V_n$$

and

$$\beta_{n-1}S_n + \gamma_{n-1}B_n = \beta_n S_n + \gamma_n B_n.$$

The first equation expresses dynamic replication while the second illustrates the concept of self-financing strategies. Its left-hand side corresponds to the portfolio value before transaction at time t_n , and the right-hand side to the portfolio value after transaction at time t_n .

As an extension to the one-period case, the resulting N-period pricing probability \tilde{P} is defined as

$$\tilde{P}\{(\omega_1,\omega_2,\ldots,\omega_N)\} = \tilde{p}^{\sum_1^N \omega_i} (1-\tilde{p})^{N-\sum_1^N \omega_i}.$$

A Bit More on Conditional Expectations. The notion of conditional expectation relative to a single event was introduced at the end of §5.2. In fact, we can define the conditional expectation of a random variable relative to a *collection* of events. We have indirectly seen this aspect already. Namely, in the case of martingales (cf. Appendix 3) and sub- and supermartingales [cf. (10.1) and (10.2)], we implicitly referred to (historic) events when we used the history of the process X, X_0, X_1, \ldots, X_n , up to time n.

We already know that an event is formally defined as a set of sample points. A collection of events will therefore be defined as a collection of sets (of sample points). We have already encountered a particular collection of sets, the Borel field introduced in Appendix 1, which satisfies certain properties [cf. (a) and (b) in Appendix 1]. It is precisely these properties that we require of a collection of events relative to which conditional expectations are defined.

In what follows, we shall need the following property: let \mathcal{F}_1 and \mathcal{F}_2 be collections of events satisfying (a) and (b) in Appendix 1. Assume $\mathcal{F}_1 \subset \mathcal{F}_2$ and that Y is a random variable with well-defined expectation. Then

$$E\{\{Y|\mathcal{F}_2\}|\mathcal{F}_1\} = E\{Y|\mathcal{F}_1\}.$$
 (10.3.2)

If you stare at this expression, you will notice on the left-hand side two conditional expectations: the first one is $Z = E\{Y|\mathcal{F}_2\}$, and the second is

 $X = E\{Z|\mathcal{F}_1\}$. We do not prove this result here (called the tower property of conditional expectations), but the interested reader can consult [Chung] or [Williams]. Despite its forbidding look, this relation expresses the fact that when conditioning over two sets of information, the overriding information will be the minimal (smallest) one. In more advanced books, a collection of sets $\{\mathcal{F}_n\}$ such that $\mathcal{F}_n \subset \mathcal{F}_{n+1}$ is sometimes called a filtration.

Proposition 3. The time-0 price of a European call option on a stock with price process S defined above and payoff q at time N is

$$V_0 = \frac{1}{(1+r)^N} \tilde{E} \{g \mid (S_0, B_0)\},$$

where \tilde{E} refers to the expectation under the pricing probability \tilde{P} . For an option with strike K, we have more specifically

$$V_0 = \begin{cases} \frac{1}{(1+r)^N} \sum_{n=0}^{N} {N \choose n} \tilde{p}^n (1-\tilde{p})^{N-n} \max \left(u^n d^{N-n} S_0 - K, 0 \right) \text{ for a call,} \\ \frac{1}{(1+r)^N} \sum_{n=0}^{N} {N \choose n} \tilde{p}^n (1-\tilde{p})^{N-n} \max \left(K - u^n d^{N-n} S_0, 0 \right) \text{ for a put.} \end{cases}$$

Proof: As in the one-period case (cf. Proposition 1) we can write, for $1 \le i \le N$,

$$V_{i-1} = E\left\{\frac{1}{1+r}V_i|(S_{i-1}, B_{i-1})\right\}.$$

Then

$$V_{i-2} = E\left\{\frac{1}{1+r}V_{i-1}|(S_{i-2}, B_{i-2})\right\}$$

$$= E\left\{\frac{1}{(1+r)^2}E\left\{\frac{1}{1+r}V_{i}|(S_{i-1}, B_{i-1})\right\}|(S_{i-2}, B_{i-2})\right\}$$

$$= E\left\{\frac{1}{(1+r)^2}V_{i}|(S_{i-2}, B_{i-2})\right\},$$

where we used the tower property in the last equality. The rest of the proof follows easily and is left as an exercise. [Hint: use the binomial distribution as in (4.4.15).]

10.4. Fundamental asset pricing theorems

In the binomial model above we were able to obtain a unique price for each of the European put and European call. This price was derived using the assumption of no-arbitrage, which in turn led to the construction of a probability under which all prices are martingales. These results are in fact illustrations of the following asset pricing theorems:

- A finite market model is arbitrage-free if and only if there exists an equivalent probability under which all asset price processes are martingales;
- 2. An arbitrage-free market model is complete (i.e., where every European contingent claim can be priced) if and only if the equivalent probability is unique.

Note that in both (1) and (2) above we mention "market model" to refer to the particular probability model (i.e., sample space, probability, etc.) to represent the market. This aspect is crucial. For example, if you replace in the one-period case Ω with a set of three sample points, then there is no pricing probability (check it as an exercise). To fully see the extent of these theorems in the context of finite probability spaces with finite time horizon, we refer to [Jacod and Shiryaev]. Despite its appearance, the last reference is accessible even with knowledge at the level of this book. New notions are defined and the point of the authors is to show that some deep results can be proved with elementary methods. These results pertain to so-called finite market models; i.e., Ω is finite, and such that $P(\omega) > 0$ for all $\omega \in \Omega$; there are finitely many assets (in our case three), the horizon is finite and trading occurs at a finite number of dates.

On the other hand, the continuous analogue (continuous time and/or state space) is more complicated. Interestingly, the continuous model of Black and Scholes and the binomial model we have considered here are connected. When the length of each of the intervals in the N-period model tends to zero while N increases to infinity, then the value V_0 obtained in Proposition 3 tends to that of the Black and Scholes (for readable details at the level of this book, see, for example, Chapter 5 in [Cox and Rubinstein]).

Exercises

- 1. Convince yourself that $\max(K-S_T,0)$ is the right payoff for a European put by repeating the same argument we used for the call.
- 2. Prove that the discounted stock price process $\{Y_n\}_{n=0}^N$ is also a martingale under the pricing probability \tilde{P} . (This should be very straightforward from the one-period case we saw previously and from the proof of Proposition 3.)
- 3. Using the Black-Scholes formulas show that the put price P_0 is a decreasing function of the initial stock price S_0 , while that of the call C_0 is increasing. In other words, the put is less valuable than the call for large stock prices (think if it makes sense by looking at the payoff functions).

- 4. Show that both P_0 and C_0 are increasing functions of σ (volatility). This means that both the put and call are more valuable when the prospects of a stock are increasingly uncertain.
- 5. Show that C_0 is an increasing function of the time to maturity T, i.e., a European call option is more valuable with more distant exercise dates. What can you say about the European put in this regard?
- 6. Put-Call Parity (1)

Verify that $C_0 = P_0 + S_0 - Ke^{-rT}$. This means that for the same exercise date and strike, the call price can be deducted from that of a put. This is an important property because in some cases it is easier, for example, to derive certain mathematical results with puts rather than calls. The latter is true because of the boundedness of the payoff of the put (check this fact as an exercise).

7. Put-Call Parity (2)

To prove the parity result, consider a portfolio consisting (at conception, i.e., now) of one call held long, a put held short, a share of the stock held short, and a borrowed amount Ke^{-rT} . Here, both the call and put have strike K and exercise date T.

Show that the value of this portfolio on date T is 0. As a result, under the no-arbitrage requirement, it must be that the value of the portfolio at conception is 0, which results in the put—call parity relation given in the previous exercise.

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Answers to Problems

Chapter 1

- 7. $(A \cup B)(B \cup C) = ABC + ABC^c + A^cBC^c + A^cBC^c + AB^cC$; $A \setminus B = AB^cC + AB^cC^c$; {the set of ω which belongs to exactly one of the sets $A, B, C \in AB^cC^c + A^cBC^c + A^cB^cC$.
- 10. The dual is true.
- 14. Define $A \# B = A^c \cup B^c$, or $A^c \cap B^c$.
- 19. $I_{A \setminus B} = I_A I_A I_B; I_{A-B} = I_A I_B.$
- 20. $I_{A \cup B \cup C} = I_A + I_B + I_C I_{AB} I_{AC} I_{BC} + I_{ABC}$.

- 4. $P(A + B) \le P(A) + P(B)$.
- 5. $P(S_1 + S_2 + S_3 + S_4) \ge P(S_1) + P(S_2) + P(S_3) + P(S_4)$.
- 11. Take $AB = \emptyset, P(A) > 0, P(B) > 0$.
- 13. 17.
- 14. 126.
- 15. $|A \cup B \cup C| = |A| + |B| + |C| |AB| |AC| |BC| + |ABC|$.
- 16. $P(A \Delta B) = P(A) + P(B) 2P(AB) = 2P(A \cup B) P(A) P(B)$.
- 17. Equality holds when m and n are relatively prime.
- 20. $p_n = 1/2^n, n \ge 1; p_n = 1/(n(n+1)), n \ge 1.$
- 22. 14/60.

- 24. If A is independent of itself, then P(A) = 0 or P(A) = 1; if A and B are disjoint and independent, then P(A)P(B) = 0.
- 28. $p_1p_2q_3p_4q_5$, where p_k = probability that the kth coin falls heads, $q_k = 1 p_k$. The probability of exactly 3 heads for 5 coins is equal to $\sum p_{k_1}p_{k_2}p_{k_3}q_{k_4}q_{k_5}$ where the sum ranges over the 10 unordered triples (k_1, k_2, k_3) of (1, 2, 3, 4, 5) and (k_4, k_5) denotes the remaining unordered pair.

- 1. 3+2; $3+2+(3\times 2)$.
- 2. 3^2 , $\binom{3+2-1}{2}$.
- 3. Three shirts are delivered in two different packages each of which may contain 0 to 3. If the shirts are distinguishable: 2^3 ; if not: $\binom{2+3-1}{3}$.
- 4. $3 \times 4 \times 3 \times 5 \times 3$; $3 \times 4 \times (3+1) \times (2+1) \times 3$.
- 5. $26^2 + 26^3$; 100.
- $6.9^7.$
- 7. $\binom{12}{6}$.
- 8. 4!; $2 \times 4!$ 4!.
- 9. $\binom{20}{3}$; $(20)_3$.
- 10. 35 (0 sum being excluded); 23.
- 11. 1/2 if the missing ones are as likely to be of the same size as of different sizes; 2/3 if each missing one is equally likely to be of any of the sizes.
- 12. 2/3; 4!/6! or $(2 \times 4!)/6!$ depending on whether the two keys are tried in one or both orders (how is the lost key counted?).
- 13. 20/216 (by enumeration); some interpreted "steadily increasing" to mean "forming an arithmetical progression," if you know what that means.
- 14. (a) $1/6^3$; (b) $\{6 \times 1 + 90 \times 3 + 120 \times 6\}/6^6$.
- 15. $\binom{6}{4}4!$; $\binom{6}{3}\binom{4}{3}3!$.
- 16. $1 \left\{ \binom{5}{0} \binom{5}{4} + \binom{5}{1} \binom{4}{3} + \binom{5}{2} \binom{3}{2} + \binom{5}{3} \binom{2}{1} + \binom{5}{4} \binom{1}{0} \right\} / \binom{10}{4}$.
- 17. From an outside lane: 3/8; from an inside lane: 11/16.
- $18. \left(\frac{m-1}{m}\right)^n; \frac{(m-1)_n}{(m)_n}.$
- 19. 1/6, 4/6, 1/6.

20. (a)
$$4 / {18 \choose 15}$$
; (b) ${14 \choose 11} / {18 \choose 15}$.

- 21. Assuming that neither pile is empty: (a) both distinguishable: $2^{10} 2$; (b) books distinguishable but piles not: $(2^{10} 2)/2$; (c) piles distinguishable but books not: 9; (d) both indistinguishable: 5.
- $22. \ \frac{10!}{3! \ 3! \ 2! \ 2!}; \frac{10!}{3! \ 3! \ 2! \ 2!} \times \frac{4!}{2! \ 2!}; \frac{10!}{3! \ 3! \ 2! \ 2!} \times \frac{6!}{2! \ 2! \ 2!}.$
- 23. (a) $\binom{31}{15}^7 \binom{30}{15}^4 \binom{29}{15} (180)!/(366)_{180}$,
- (b) $(305)_{30}/(366)_{30}$.
- $24. \binom{29}{10} / \binom{49}{30}.$
- $25. \ \binom{n-100}{93} \binom{100}{7} / \binom{n}{100}.$
- 27. Divide the following numbers by $\binom{52}{5}$:
 - (a) $4 \times {13 \choose 5}$; (b) 9×4^5 ; (c) 4×9 ;
 - (d) 13×48 ; (e) $13 \times 12 \times 4 \times 6$.
- 29. Divide the following numbers by 6^6 :

$$\begin{aligned} &6; 6 \times 5 \times \frac{6!}{5! \, 1!}; 6 \times 5 \times \frac{6!}{4! \, 2!}; 6 \times \binom{5}{2} \times \frac{6!}{4!} \\ &\binom{6}{2} \times \frac{6!}{3! \, 3!}; (6)_3 \times \frac{6!}{3! \, 2!}; 6 \times \binom{5}{3} \times \frac{6!}{3!}; \\ &\binom{6}{3} \times \frac{6!}{2! \, 2! \, 2!}; \binom{6}{2} \binom{4}{2} \times \frac{6!}{2! \, 2!}; \binom{6}{1} \binom{5}{4} \times \frac{6!}{2!}; 6!. \end{aligned}$$

Add these up for a check; use your calculator if you have one.

30. Do the problem first for n=2 by enumeration to see the situation. In general, suppose that the right pocket is found empty and there are k matches remaining in the left pocket. For $0 \le k \le n$ the probability of this event is equal to $\frac{1}{2^{2n-k}} \binom{2n-k}{n} \frac{1}{2}$. This must be multiplied by 2 because right and left may be interchanged. A cute corollary to the solution is the formula below:

$$\sum_{k=0}^{n} \frac{1}{2^{2n-k}} \binom{2n-k}{n} = 1.$$

- 2. $P\{X + Y = k\} = 1/3 \text{ for } k = 3, 4, 5; \text{ same for } Y + Z \text{ and } Z + X.$
- 3. $P\{X+Y-Z=k\}=1/3 \text{ for } k=0,2,4;$ $P\{\sqrt{(X^2+Y^2)Z}=x\}=1/3 \text{ for } x=\sqrt{13},\sqrt{15},\sqrt{20};$ $P\{Z/|X-Y|=3\}=1/3,P\{Z/|X-Y|=1\}=2/3.$
- 4. Let $P(\omega_j) = 1/10$ for j = 1, 2; = 1/5 for j = 3, 4; = 2/5 for j = 5; $X(\omega_j) = j$ for $1 \le j \le 5$; $Y(\omega_j) = \sqrt{3}$ for j = 1, 4; $= \pi$ for j = 2, 5; $= \sqrt{2}$ for j = 3.
- 5. Let $P(\omega_i) = p_i, X(\omega_i) = v_i, 1 \le i \le n$.
- 6. $\{X + Y = 7\} = \{(1,6), (2,5), (3,4), (4,3), (5,2), (6,1)\}.$
- 8. $P{Y = 14000 + 4n} = 1/5000 \text{ for } 1 \le n \le 5000; E(Y) = 24002.$
- 9. $P\{Y = 11000 + 3n\} = 1/10000$ for $1 \le n \le 1000$; $P\{Y = 10000 + 4n\} = 1/10000$ for $1001 \le n \le 10000$; E(Y) = 41052.05.
- 10. $E(Y) = 29000 + 7000 \cdot e^{-2/7}$.
- 11. $\lambda e^{-\lambda x}, x > 0$.
- 12. $2xf(x^2), x > 0; 2x/(b-a)$ for $\sqrt{a} \le x \le \sqrt{b}$.
- 13. (i) f((x-b)/a)/|a| if $a \neq 0$. (ii) $\frac{1}{2\sqrt{x}} \{f(\sqrt{x}) + f(-\sqrt{x})\}, x > 0$.
- 15. $c = 1/(1 q^m)$.
- 16. $P(Y=j) = \left(\frac{n}{n+j}\right) \frac{1}{2^n}$, for $-n \le j \le n$ such that n+j is even. E(Y) = 0.
- 17. $P(X=j) = {11 \choose j} {539 \choose 25-j} / {550 \choose 25}, 0 \le j \le 25.$
- 18. If there are r rotten apples in a bushel of n apples and k are picked at random, the expected number of rotten ones among those picked is equal to kr/n.
- 19. $P(X \ge m) = \frac{1}{m}, E(X) = +\infty.$
- 20. 1.
- 21. Choose $v_n = (-1)^n 2^n / n$, $p_n = 1/2^n$.
- 23. According to the three hypotheses of Example 11 in §4.5: (1) $\sqrt{3}/2$; (2) 3/4; (3) 2/3.
- 24. 2.
- 26. $F_R(r) = r^2/100, f_R(r) = r/50$ for $0 \le r \le 100$; E(R) = 20/3.
- 27. $Y = d \tan \theta$, where d is the distance from the muzzle to the wall and θ is the angle the pistol makes with the horizontal direction. $P(Y \le y) = \arctan(y/d); E(Y) = +\infty.$
- 28. $E(2^X) = +\infty$.

29. If at most m tosses are allowed, then his expectation is m cents.

31.
$$P((X,Y) = (m,m')) = \binom{n}{2}^{-1}$$
 for $1 \le m < m' \le n$; $P(X = m) = (n-m)\binom{n}{2}^{-1}$; $P(Y = m') = (m'-1)\binom{n}{2}^{-1}$; $P(Y - X = k) = (n-k)\binom{n}{2}^{-1}$, $1 \le k \le n-1$.

32. Joint density of
$$(X,Y)$$
 is $f(u,v) = \begin{cases} 2 & \text{if } 0 \le u < v \le 1, \\ 0 & \text{otherwise.} \end{cases}$

- 1. 1050/6145, 95/1095.
- $2. \ \frac{18826}{19400}.$
- $3.\ 5/9.$
- 4. (a) 1/2; (b) 1/10.
- 5. 1/4; 1/4.
- 6. 1/4.
- $7. \ 1/2.$
- 8. $2\beta(1-\alpha+\beta)^{-1}$.
- 9. 6/11, 3/11, 2/11.
- 10. 400/568.
- $17. \ 1/2.$
- 18. $p^3 + (3/2) \times p^3(1-p)$.
- 19. 379/400.
- 20. $P(\text{no umbrella} \mid \text{rain}) = 2/9$; $P(\text{no rain} \mid \text{umbrella}) = 5/12$.
- 21. 27/43.
- 22. $[p^2 + (1-p)^2]/[3p^2 + (1-p)^2]$.
- 23. (a) 3/8; (b) 3/4; (c) 1/3.
- 25. (a) $\frac{1}{6} \sum_{n=1}^{6} \binom{n}{k} \frac{1}{2^n}$; (b) $\binom{n}{3} \frac{1}{2^n} \left\{ \sum_{n=3}^{6} \binom{n}{3} \frac{1}{2^n} \right\}^{-1}$ for $3 \le n \le 6$.
- 26. The probabilities that the number is equal to 1, 2, 3, 4, 5, 6 are equal, respectively, to:

 - (1) p_1^2 ; (2) $p_1p_2 + p_1^2p_2$; (3) $p_1p_3 + 2p_1p_2^2 + p_1^3p_3$; (4) $2p_1p_2p_3 + p_2^3 + 3p_1^2p_2p_3$; (5) $2p_2^2p_3 + 3p_1p_2^2p_3 + 3p_1^2p_3^2$; (6) $p_2p_3^2 + p_2^3p_3 + 6p_1p_2p_3^2$; (7) $3p_1p_3^3 + 3p_2^2p_3^2$; (8) $3p_2p_3^3$; (9) p_3^4 . Tedious work? See Example 20 and Exercise No. 23 of Chapter 8 for the general method.

27.
$$\left(\frac{4}{6}\right)^n - 2\left(\frac{3}{6}\right)^n + \left(\frac{2}{6}\right)^n$$
.

28.
$$\sum_{n=0}^{\infty} p_n \left(\sum_{k=0}^{n} p_k \right) ; \sum_{n=0}^{\infty} p_n^2.$$

- $29. \ 2/7.$
- 30. $P(\text{maximum} < y \mid \text{minimum} < x) = y^2/(2x x^2) \text{ if } y \le x; = (2xy x^2)/(2x x^2) \text{ if } y > x.$
- 31. 1/4.
- 33. (a) (r+c)/(b+r+2c); (b) (r+2c)/(b+r+2c); (c), (e), (f): (r+c)/(b+r+c); (d) same as (b).
- 34. $\{b_1(b_2+1)r_1+b_1r_2(r_1+1)+r_1b_2(r_1-1)+r_1(r_2+1)r_1\}/(b_1+r_1)^2(b_2+r_2+1).$

35.
$$\left(\sum_{k=1}^{N} k^{n+1}\right) / N\left(\sum_{k=1}^{N} k^{n}\right)$$
.

39.
$$(1+p)^2/4$$
; $(1+pq)/2$.

- 1. \$.1175; \$.5875.
- 2. \$94000; \$306000.

3.
$$2\left(\frac{3}{13} + \frac{2}{12} + \frac{4}{13} + \frac{3}{14} + \frac{4}{14}\right)$$
.

- 4. 21; 35/2.
- 5. .5; 2.5.

6.
$$13/4$$
; $4\left\{1-\binom{39}{13}\right\}\binom{52}{13}$.

7.
$$(6/7)^{25}$$
; $7\left\{1-\left(\frac{6}{7}\right)^{25}\right\}$.

- 8. (a) $1 (364/365)^{500} 500(364)^{499}/(365)^{500}$.
 - (b) 500/365.

(c)
$$365 \left\{ 1 - \left(\frac{364}{365} \right)^{500} \right\}$$
.

- (d) 365p, where p is the number in (a).
- 9. Expected number of boxes getting k tokens is equal to $m\binom{n}{k}\frac{(m-1)^{n-k}}{m^n}$; expected number of tokens alone in a box is equal to $n\left(\frac{m-1}{m}\right)^{n-1}$.

10.
$$P(n_j \text{ tokens in } j \text{th box for } 1 \leq j \leq m) = \frac{n!}{n_1! \cdots n_m!} \frac{1}{m^n}$$
, where $n_1 + \cdots + n_m = n$.

- 11. 49.
- 12. 7/2.
- 13. 100p; $10\sqrt{p(1-p)}$.
- $14. \ 46/5.$

15. (a)
$$N+1$$
; (b) $\sum_{n=1}^{N+1} \frac{(N)_{n-1}}{N^{n-1}}$.

16. Let M denote the maximum. With replacement:

$$P(M = k) = \frac{k^n - (k-1)^n}{N^n}, 1 \le k \le N;$$

$$E(M) = \sum_{k=1}^{N} \left\{ 1 - \left(\frac{k-1}{N}\right)^n \right\};$$

without replacement

without replacement:
$$P(M=k) = \binom{k-1}{n-1} / \binom{N}{n}, n \le k \le N;$$

$$E(M) = n(N+1)/(n+1).$$

- 17. (a) nr/(b+r); (b) $(r^2+br+cnr)/(b+r)$.
- 19. 1/p.
- 22. $E(X) = 1/\lambda$

23.
$$E(T) = \frac{a}{\lambda} + \frac{1-a}{\mu}; \ \sigma^2(T) = \frac{2a}{\lambda^2} + \frac{2(1-a)}{\mu^2} - \left(\frac{a}{\lambda} + \frac{1-a}{\mu}\right)^2.$$

- 24. $E(T | T > n) = 1/\lambda$.
- 25. (a) $1/5\lambda$; (b) $137/60\lambda$.
- 26. .4%.

27.
$$E(aX + b) = aE(X) + b, \sigma^{2}(aX + b) = a^{2}\sigma^{2}(X).$$

- 28. Probability that he quits winning is 127/128, having won \$1; probability that he quits because he does not have enough to double his last bet is 1/128, having lost \$127. Expectation is zero. So is it worth it? In the second case he has probability 1/256 of losing \$150, same probability of losing \$104, and probability 127/128 of winning \$1. Expectation is still zero.
- 29. E(maximum) = n/(n+1); E(minimum) = 1/(n+1); E(range) =(n-1)/(n+1).
- 30. $g(z) = \prod_{i=1}^{n} (q_i + p_i z); g'(1) = \sum_{i=1}^{n} p_i$
- 31. $u_k = P\{S_n \le k\}; g(z) = (q + pz)^n/(1-z).$
- 32. $g(z) = (1 z^{2N+1})/(2N+1)z^N(1-z); g'(1) = 0.$
- 33. $\binom{2n}{n} \frac{1}{4^n}$.

34.
$$g(z) = z^N \prod_{j=0}^{N-1} \frac{N-j}{N-jz}; g'(1) = N \sum_{j=0}^{N-1} \frac{1}{N-j}.$$

35.
$$m_1 = g'(1); m_2 = g''(1) + g'(1); m_3 = g'''(1) + 3g''(1) + g'(1); m_4 = g^{(iv)}(1) + 6g'''(1) + 7g''(1) + g'(1).$$

36.
$$(-1)^n L^{(n)}(0)$$
.

37. (a)
$$(1-e^{-c\lambda})/c\lambda$$
, $\lambda > 0$; (b) $2(1-e^{-c\lambda}-c\lambda e^{-c\lambda})/c^2\lambda^2$, $\lambda > 0$; (c) $L(\mu) = \lambda^n/(\lambda+\mu)^n$.

38. Laplace transform of S_n is equal to $\lambda^n/(\lambda+\mu)^n$;

$$P(a < S_n < b) = \frac{\lambda^n}{(n-1)!} \int_a^b u^{n-1} e^{-\lambda u} \, du.$$

1.
$$1 - \frac{5}{2}e^{-2/3}$$
.

2.
$$\left(1 - \frac{4}{100}\right)^{25} \approx e^{-1}$$
.

- 3. $e^{-\alpha} \alpha^k / k!$, where $\alpha = 1000/324$.
- 4. $e^{-20} \sum_{k=20}^{30} (20)^k / k!$.
- 5. Let $\alpha_1 = 4/3$, $\alpha_2 = 2$. $P\{X_1 = j \mid X_1 + X_2 = 2\} = \frac{2!}{i!(2-i)!} \frac{\alpha_1^j \alpha_2^{2-j}}{(\alpha_1 + \alpha_2)^2}$ for i = 0, 1, 2.
- 6. If (n+1)p is not an integer, the maximum term of $B_k(n;p)$ occurs at $k = \lceil (n+1)p \rceil$ where $\lceil x \rceil$ denotes the greatest integer not exceeding x; if (n+1)p is an integer, there are two equal maximum terms for k = (n+1)p - 1 and (n+1)p.
- 7. If α is not an integer, the maximum term of $\pi_k(\alpha)$ occurs at $k = [\alpha]$; if α is an integer, at $k = \alpha - 1$ and $k = \alpha$.
- 8. $\exp[-\lambda c + \alpha(e^{-\lambda h} 1)]$.
- 9. $\pi_k(\alpha+\beta)$.

11.
$$e^{-50} \sum_{k=50}^{60} (50)^k / k!$$
.
12. $\frac{1}{(n-1)! \, 2^n} \int_{N}^{\infty} u^{n-1} e^{-u/2} \, du$.

13.
$$\Phi\left(3\sqrt{\frac{12}{35}}\right) - \Phi\left(-2\sqrt{\frac{12}{35}}\right).$$

- 14. Find n such that $2\Phi(\sqrt{n}/10) 1 \ge .95$. We may suppose p > 1/2 (for the tack I used).
- 15. 537.
- 16. 475.
- 24. $(1/(2\pi x))^{-1/2} e^{-x/2}$.
- 27. $P\{\delta'(t) > u\} = e^{-\alpha u}; P\{\delta(t) > u\} = e^{-\alpha u} \text{ for } u < t; = 0 \text{ for } u \ge t.$
- 28. No!

Chapter 8

1. (a)
$$p^4 + 3p^3q + 2p^2q^2$$
; (b) $p^4 + 2p^3q + 2pq^3 + q^4$; (c) $1 - (p^4 + p^3q)$; (d) $1 - (p^4 + p^3q + pq^3 + q^4)$.

2. For
$$Y_n$$
: $I =$ the set of even integers; $p_{2i,2i+2} = p^2, p_{2i,2i} = 2pq, p_{2i,2i-2} = q^2;$

For Z_n : I = the set of odd integers; $p_{2i-1,2i+1} = p^2, p_{2i-1,2i-1} = 2pq, p_{2i-1,2i-3} = q^2; P\{Z_0 = 1\} = p, P\{Z_0 = -1\} = q.$

3. For
$$X_n: I =$$
 the set of nonnegative integers; $p_{i,i} = q, p_{i,i+1} = p$. For $Y_n: I =$ the set of all integers, $p_{i,i-1} = q, p_{i,i+1} = p$.

4.
$$P\{|Y_{2n+1}| = 2i+1 \mid \mid Y_{2n}| = 2i\} = (p^{2i+1}+q^{2i+1})/(p^{2i}+q^{2i}), P\{|Y_{2n+1}| = 2i-1 \mid \mid Y_{2n}| = 2i\} = (p^{2i}q+pq^{2i})/(p^{2i}+q^{2i}).$$

6.
$$\begin{bmatrix} 1 & 0 \\ \alpha & 1 - \alpha \end{bmatrix}; f_{21}^{(n)} = (1 - \alpha)^{n-1}\alpha; p_{21}^{(n)} = 1 - (1 - \alpha)^n; m_{21} = 1/\alpha.$$

9.
$$I = \{0; 2^i, 0 \le i \le n-1\}$$

 $p_{2^i, 2^{i+1}} = 1/2, p_{2^i, 0} = 12 \text{ for } 0 \le i \le n-1.$

10. 1/13.

10. 1/13.
11.
$$\begin{array}{c|cccc} U & H & D \\ \hline U & 0 & 1 & 0 \\ H & p & 0 & q \\ D & 0 & 1 & 0 \\ \end{array}$$
 $w_U = p/2, w_H = 1/2, w_D = q/2.$

12. Same as given in (8.1.9) and (8.1.10).

13.
$$e_j = \frac{l(1-r^j)}{(p-q)(1-r^l)} - \frac{j}{p-q}$$
 where $r = q/p$.

15.
$$p_{i,i-1} = (i/N)^2, p_{i,i} = 2i(N-i)/N^2, p_{i,i+1} = ((N-i)/N)^2;$$

$$w_i = \binom{N}{i}^2 / \binom{2N}{N}; 0 \le i \le N.$$

16.
$$w_s = (1 - \beta)/(2 - \alpha - \beta); w_f = (1 - \alpha)/(2 - \alpha - \beta).$$

18.
$$p_{j,j+1} = p, p_{j,0} = 1 - p; w_j = p^j q, 0 \le j \le \infty.$$

19. Let
$$r=pq, A^{-1}=1+p^{-1}\sum_{k=1}^{c-1}r^k+r^{c-1};$$
 then $w_0=A; w_k=p^{-1}r^kA, 1\leq k\leq c-1; w_c=r^{c-1}A.$

20.
$$f_{21}^* = .721; f_{31}^* = .628.$$

21.
$$f_{i,2N}^* = i/2N, f_{i,0}^* = 1 - (i/2N), 0 \le i \le 2N.$$

22.
$$(1-a_2+\sqrt{a_1^2-2a_1+1-4a_0a_2})/2a_2$$
.

23. Coefficient of
$$z^j$$
 in $g(h(z))$, where $h(z) = \sum_{j=0}^{\infty} b_j z^j$.

24. Let e_j denote the expected number of further impulses received until the meter registers the value l, when the meter reads j; then $e_j = l$ for $1 \le j \le l-1$, $e_l = 0$.

25.
$$e_m = \sum_{j=1}^m \frac{1}{j}$$
.

32.
$$p_{ij} = 1/(i+1)$$
 for $s - i \le j \le s$; = 0 otherwise;

$$w_j = 2(j+1)/(s+1)(s+2)$$
 for $0 \le j \le s$; $\sum_{j=0}^{s} jw_j = 2s/3$.

33.
$$p_{ij} = {i \choose s-j} p^{s-j} (1-p)^{i-s+j}$$
 for $s-i \le j \le s$; = 0 otherwise; $w_j = {s \choose j} \left(\frac{1}{1+p}\right)^j \left(\frac{p}{1+p}\right)^{s-j}$ for $0 \le j \le s$; $\sum_{j=0}^s j w_j = s/(1+p)$.

44.
$$P\{S=k \mid X_0=i\} = p_{ii}^k (1-p_{ii}^k), k \ge 1.$$

45.
$$\tilde{p}_{ij} = p_{ij}/(1 - p_{ii})$$
 for $i \neq j$; $\tilde{p}_{ii} = 0$.

46.
$$P\{(X_n, X_{n+1}) = (k, 2k - j + 1) \mid (X_{n-1}, X_n) = (j, k)\} = p,$$

 $P\{(X_n, X_{n+1}) = (k, 2k - j) \mid (X_{n-1}, X_n) = (j, k)\} = q.$
Let $H_n^{(3)} = \sum_{v=1}^n H_v^{(2)}$, then $\{H_n^{(3)}\}$ is a Markov chain of order 3, etc.

Chapter 9

- 1. $R(\omega_1) = u 1$, $R(\omega_2) = d 1$. With u = 1.10, $R(\omega_1) = .10$, i.e., a gain of 10% should ω_1 be drawn. With d = .75, $R(\omega_2) = -.25$, or a drop of 25% if ω_2 is drawn.
- 2. We are interested in comparing $P\{R \leq 0\}$ versus $P\{R > 0\}$. $P\{R \leq 0\} = P\{(R \mu)/\sigma \leq -\mu/\sigma\}$. Recall that $(R \mu)/\sigma$ has a standard normal distribution. So when $\mu = 0$, $P\{R \leq 0\} = 1/2 = P\{R > 0\}$; when $\mu < 0$, $-\mu/\sigma > 0$ and therefore $P\{R \leq 0\} > 1/2 > P\{R > 0\}$; and $\mu > 0$ then $-\mu/\sigma < 0$, leading to $P\{R \leq 0\} < 1/2 < P\{R > 0\}$.

3. Let R_1 and R_2 be the returns of the first and second assets, respectively. Our hypothesis is $\sigma(R_2) < \sigma(R_1)$. For ease of notation, we let $\sigma_i \equiv \sigma(R_i)$, i = 1, 2. Then, for $\alpha \in (0, 1)$, we have

$$\begin{split} \sigma^2 \left(\alpha R_1 + (1 - \alpha) R_2 \right) &= \alpha^2 \sigma_1^2 + (1 - \alpha)^2 \sigma_2^2 + 2\alpha (1 - \alpha) \rho_{12} \sigma_1 \sigma_2 \\ &< \alpha^2 \sigma_1^2 + (1 - \alpha)^2 \sigma_1^2 + 2\alpha (1 - \alpha) \rho_{12} \sigma_1 \sigma_2 \\ &< \alpha^2 \sigma_1^2 + (1 - \alpha)^2 \sigma_1^2 + 2\alpha (1 - \alpha) \sigma_1^2 \\ &= \left(\alpha \sigma_1 + (1 - \alpha) \sigma_1 \right)^2 \\ &= \sigma_1^2, \end{split}$$

where the first inequality results from $\sigma_2 < \sigma_1$, and the second from $|\rho_{12}| < 1$, $\sigma_2 < \sigma_1$, and $0 < \alpha < 1$.

- 4. $\sigma_1 = 0$ and $\mu_1 = \mu_2 = \mu_3$. In this case, there is no point trying to diversify across assets since they all yield the same expected return and one of them is riskless.
- 5. If such b and ρ₁₂ exist then b is a solution of the quadratic in b expressed through (9.5.21). For its discriminant to be nonnegative, we must have |ρ₁₂| ≥ 1. Since, by definition, |ρ₁₂| ≤ 1, only two cases are possible: ρ₁₂ = 1 or ρ₁₂ = −1. In either case, we must have b = −ρ₁₂σ₁/σ₂. If ρ₁₂ = 1, then (μ₁ − μ₀)/σ₁ = (μ₂ − μ₀)/σ₂. We have an equality between two values of a type of ratio that is referred to as the Sharpe ratio. It is used to compare risky assets when a riskless asset is available. It measures the expected excess return of an asset over the riskless asset, i.e., μ_i − μ₀, relative to the risk of the asset, σ_i. Another way to interpret this ratio is to view it as measuring the expected excess return over the riskless rate as the number of standard deviations of this asset's return. When ρ₁₂ = 1 the returns of assets 1 and 2 are perfectly correlated. In addition, they have the same Sharpe ratio. We thus have an indication that no diversification across these two assets would be meaningful.

If $\rho_{12} = -1$, then $(\sigma_1 + \sigma_2)\mu_0 = -(\sigma_1\mu_2 + \sigma_2\mu_1)$, which shows that μ_0 , μ_1 , and μ_2 cannot all be positive. Thus there cannot be a diversification involving all three assets.

Chapter 10

1. Since the put gives you the right to sell at time T, you do so only if $K > S_T$. In this case, you can first purchase the stock at price S_T and sell it at price K to obtain a net gain (payoff) of $K - S_T$. If $K \leq S_T$, there is no point trying to sell at K when the market price is S_T .

2.

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$$\tilde{E}[Y_{k+1}|S_1,\dots,S_k] = \frac{1}{(1+r)^{k+1}} \tilde{E}[S_{k+1}|S_1,\dots,S_k]$$

$$= \frac{1}{(1+r)^{k+1}} \tilde{E}[S_{k+1}|S_k]$$

$$= \frac{1}{(1+r)^{k+1}} (1+r)S_k$$

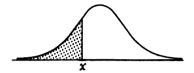
$$= Y_k.$$

- 3. For the call, $\partial C_0/\partial S_0 = \Phi(d_1) > 0$, where $\Phi(x)$ is the cumulative standard normal distribution. For the put, $\partial P_0/\partial S_0 = -\Phi(-d_1) < 0$.
- 4. $\partial C_0/\partial \sigma = S_0\sqrt{T}\phi(d_1) > 0$ and $\partial P_0/\partial \sigma = S_0\sqrt{T}\phi(-d_1) > 0$, where $\phi(x)$ is the density of the standard normal distribution.
- 5. For a call, $\partial C_0/\partial T = -S_0\sigma\phi(d_1)/\left(2\sqrt{T}\right) rKe^{-rT}\Phi(d_2) < 0$. For a put, $\partial C_0/\partial T = -S_0\sigma\phi(-d_1)/\left(2\sqrt{T}\right) + rKe^{-rT}\Phi(-d_2)$, which can be of any sign.
- 7. The value of this portfolio at time 0 is $V_0 = C_0 P_0 S_0 + Ke^{-rT}$. At time T, its value is $V_T = C_T P_T S_T + K$, where $C_T = \max(S_T K, 0)$ and $P_T = \max(K S_T, 0)$.
 - Case 1: $K > S_T$. Then $C_T = 0$, $P_T = K S_T$, and $V_T = -K + S_T S_T + K = 0$.
 - Case 2: $K < S_T$. Then $C_T = S_T K$, $P_T = 0$, and $V_T = S_T K S_T + K = 0$.
 - Case 3: $K = S_T$. Then $C_T = P_T = 0$ and $V_T = 0$.

Values of the Standard Normal Distribution Function

TABLE 1 Values of the standard normal distribution function

$$\Phi(x) = \int_{-\infty}^{x} \frac{1}{\sqrt{2\pi}} e^{-u^{2}/2} du = P(X \le x)$$



x	0	1	2	3	4	5	6	7	8	9
-3.	.0013	.0010	.0007	.0005	.0003	.0002	.0002	.0001	.0001	.0000
-2.9	.0019	.0018	.0017	.0017	.0016	.0016	.0015	.0015	.0014	.0014
-2.8	.0026	.0025	.0024	.0023	.0023	.0022	.0021	.0020	.0020	.0019
-2.7	.0035	.0034	.0033	.0032	.0031	.0030	.0029	.0028	.0027	.0026
-2.6	.0047	.0045	.0044	.0043	.0041	.0040	.0039	.0038	.0037	.0036
-2.5	.0062	.0060	.0059	.0057	.0055	.0054	.0052	.0051	.0049	.0048
-2.4	.0082	.0080	.0078	.0075	.0073	.0071	.0069	.0068	.0066	.0064
-2.3	.0107	.0104	.0102	.0099	.0096	.0094	.0091	.0089	.0087	.0084
-2.2	.0139	.0136	.0132	.0129	.0126	.0122	.0119	.0116	.0113	.0110
-2.1	.0179	.0174	.0170	.0166	.0162	.0158	.0154	.0150	.0146	.0143
-2.0	.0228	.0222	.0217	.0212	.0207	.0202	.0197	.0192	.0188	.0183
-1.9	.0287	.0281	.0274	.0268	.0262	.0256	.0250	.0244	.0238	.0233
-1.8	.0359	.0352	.0344	.0336	.0329	.0322	.0314	.0307	.0300	.0294
-1.7	.0446	.0436	.0427	.0418	.0409	.0401	.0392	.0384	.0375	.0367
-1.6	.0548	.0537	.0526	.0516	.0505	.0495	.0485	.0475	.0465	.0455
-1.5	.0668	.0655	.0643	.0630	.0618	.0606	.0594	.0582	.0570	.0559
-1.4	.0808	.0793	.0778	.0764	.0749	.0735	.0722	.0708	.0694	.0681
-1.3	.0968	.0951	.0934	.0918	.0901	.0885	.0869	.0853	.0838	.0823
-1.2	.1151	.1131	.1112	.1093	.1075	.1056	.1038	.1020	.1003	.0985
-1.1	.1357	.1335	.1314	.1292	.1271	.1251	.1230	.1210	.1190	.1170
-1.0	.1587	.1562	.1539	.1515	.1492	.1469	.1446	.1423	.1401	.1379
9	.1841	.1814	.1788	.1762	.1736	.1711	.1685	.1660	.1635	.1611
8	.2119	.2090	.2061	.2033	.2005	.1977	.1949	.1922	.1894	.1867
7	.2420	.2389	.2358	.2327	.2297	.2266	.2236	.2206	.2177	.2148
6	.2743	.2709	.2676	.2643	.2611	.2578	.2546	.2514	.2483	.2451
- .5	.3085	.3050	.3015	.2981	.2946	.2912	.2877	.2843	.2810	.2776
4	.3446	.3409	.3372	.3336	.3300	.3264	.3228	.3192	.3156	.3121
3	.3821	.3783	.3745	.3707	.3669	.3632	.3594	.3557	.3520	.3483
2	.4207	.4168	.4129	.4090	.4052	.4013	.3974	.3936	.3897	.3859
1	.4602	.4562	.4522	.4483	.4443	.4404	.4364	.4325	.4286	.4247
0	.5000	.4960	.4920	.4880	.4840	.4801	.4761	.4721	.4681	.4641

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TABLE 1 Values of the standard normal distribution function

x	0	1	2	3	4	5	6	7	8	9
.0	.5000	.5040	.5080	.5120	.5160	.5199	.5239	.5279	.5319	.5359
.1	.5398	.5438	.5478	.5517	.5557	.5596	.5363	.5675	.5714	.5753
.2	.5793	.5832	.5871	.5910	.5948	.5987	.6026	.6064	.6103	.6141
.3	.6179	.6217	.6255	.6293	.6331	.6368	.6406	.6443	.6480	.6517
.4	.6554	.6591	.6628	.6664	.6700	.6736	.6772	.6808	.6844	.6879
.5	.6915	.6950	.6985	.7019	.7054	.7088	.7123	.7157	.7190	.7224
.6	.7257	.7291	.7324	.7357	.7389	.7422	.7454	.7486	.7517	.7549
.7	.7580	.7611	.7642	.7673	.7703	.7734	.7764	.7794	.7823	.7852
.8	.7881	.7910	.7939	.7967	.7995	.8023	.8051	.8078	.8106	.8133
.9	.8159	.8186	.8212	.8238	.8264	.8289	.8315	.8340	.8365	.8389
1.0	.8413	.8438	.8461	.8485	.8508	.8531	.8554	.8577	.8599	.8621
1.1	.8643	.8665	.8686	.8708	.8729	.8749	.8770	.8790	.8810	.8830
1.2	.8849	.8869	.8888	.8907	.8925	.8944	.8962	.8980	.8997	.9015
1.3	.9032	.9049	.9066	.9082	.9099	.9115	.9131	.9147	.9162	.9177
1.4	.9192	.9207	.9222	.9236	.9251	.9265	.9278	.9292	.9306	.9319
1.5	.9332	.9345	.9357	.9370	.9382	.9394	.9406	.9418	.9430	.9441
1.6	.9452	.9463	.9474	.9484	.9495	.9505	.9515	.9525	.9535	.9545
1.7	.9554	.9564	.9573	.9582	.9591	.9599	.9608	.9616	.9625	.9633
1.8	.9641	.9648	.9656	.9664	.9671	.9678	.9686	.9693	.9700	.9706
1.9	.9713	.9719	.9726	.9732	.9738	.9744	.9750	.9756	.9762	.9767
2.0	.9772	.9778	.9783	.9788	.9793	.9798	.9803	.9808	.9812	.9817
2.1	.9821	.9826	.9830	.9834	.9838	.9842	.9846	.9850	.9854	.9857
2.2	.9861	.9864	.9868	.9871	.9874	.9878	.9881	.9884	.9887	.9890
2.3	.9893	.9896	.9898	.9901	.9904	.9906	.9909	.9911	.9913	.9916
2.4	.9918	.9920	.9922	.9925	.9927	.9929	.9931	.9932	.9934	.9936
2.5	.9938	.9940	.9941	.9943	.9945	.9946	.9948	.9949	.9951	.9952
2.6	.9953	.9955	.9956	.9957	.9959	.9960	.9961	.9962	.9963	.9964
2.7	.9965	.9966	.9967	.9968	.9969	.9970	.9971	.9972	.9973	.9974
2.8	.9974	.9975	.9976	.9977	.9977	.9978	.9979	.9979	.9980	.9981
2.9	.9981	.9982	.9982	.9983	.9984	.9984	.9985	.9985	.9986	.9986
3.	.9987	.9990	.9993	.9995	.9997	.9998	.9998	.9999	.9999	1.0000

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