31. PROBABILITY

Revised September 2007 by G. Cowan (RHUL).

31.1. General [1–8]

An abstract definition of probability can be given by considering a set $S$, called the sample space, and possible subsets $A, B, \ldots$, the interpretation of which is left open. The probability $P$ is a real-valued function defined by the following axioms due to Kolmogorov [9]:

1. For every subset $A$ in $S$, $P(A) \geq 0$;
2. For disjoint subsets (i.e., $A \cap B = \emptyset$), $P(A \cup B) = P(A) + P(B)$;
3. $P(S) = 1$.

In addition, one defines the conditional probability $P(A|B)$ (read $P$ of $A$ given $B$) as

$$P(A|B) = \frac{P(A \cap B)}{P(B)}. \quad (31.1)$$

From this definition and using the fact that $A \cap B$ and $B \cap A$ are the same, one obtains Bayes’ theorem,

$$P(A|B) = \frac{P(B|A)P(A)}{P(B)}. \quad (31.2)$$

From the three axioms of probability and the definition of conditional probability, one obtains the law of total probability,

$$P(B) = \sum_i P(B|A_i)P(A_i), \quad (31.3)$$

for any subset $B$ and for disjoint $A_i$ with $\cup_i A_i = S$. This can be combined with Bayes’ theorem (Eq. (31.2)) to give

$$P(A|B) = \frac{P(B|A)P(A)}{\sum_i (B|A_i)P(A_i)}, \quad (31.4)$$

where the subset $A$ could, for example, be one of the $A_i$.

The most commonly used interpretation of the subsets of the sample space are outcomes of a repeatable experiment. The probability $P(A)$ is assigned a value equal to the limiting frequency of occurrence of $A$. This interpretation forms the basis of frequentist statistics.

The subsets of the sample space can also be interpreted as hypotheses, i.e., statements that are either true or false, such as ‘The mass of the $W$ boson lies between 80.3 and 80.5 GeV.’ In the frequency interpretation, such statements are either always or never true, i.e., the corresponding probabilities would be 0 or 1. Using subjective probability, however, $P(A)$ is interpreted as the degree of belief that the hypothesis $A$ is true. Subjective probability is used in Bayesian (as opposed to frequentist) statistics. Bayes’ theorem can be written
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\[ P(\text{theory}|\text{data}) \propto P(\text{data}|\text{theory})P(\text{theory}) , \]  

(31.5)

where ‘theory’ represents some hypothesis and ‘data’ is the outcome of the experiment. Here \( P(\text{theory}) \) is the prior probability for the theory, which reflects the experimenter’s degree of belief before carrying out the measurement, and \( P(\text{data}|\text{theory}) \) is the probability to have gotten the data actually obtained, given the theory, which is also called the likelihood.

Bayesian statistics provides no fundamental rule for obtaining the prior probability; this is necessarily subjective and may depend on previous measurements, theoretical prejudices, etc. Once this has been specified, however, Eq. (31.5) tells how the probability for the theory must be modified in the light of the new data to give the posterior probability, \( P(\text{theory}|\text{data}) \). As Eq. (31.5) is stated as a proportionality, the probability must be normalized by summing (or integrating) over all possible hypotheses.

31.2. Random variables

A random variable is a numerical characteristic assigned to an element of the sample space. In the frequency interpretation of probability, it corresponds to an outcome of a repeatable experiment. Let \( x \) be a possible outcome of an observation. If \( x \) can take on any value from a continuous range, we write \( f(x; \theta)dx \) as the probability that the measurement’s outcome lies between \( x \) and \( x + dx \). The function \( f(x; \theta) \) is called the probability density function (p.d.f.), which may depend on one or more parameters \( \theta \). If \( x \) can take on only discrete values (e.g., the non-negative integers), then \( f(x; \theta) \) is itself a probability.

The p.d.f. is always normalized to unit area (unit sum, if discrete). Both \( x \) and \( \theta \) may have multiple components and are then often written as vectors. If \( \theta \) is unknown, we may wish to estimate its value from a given set of measurements of \( x \); this is a central topic of statistics (see Sec. 32).

The cumulative distribution function \( F(a) \) is the probability that \( x \leq a \):

\[ F(a) = \int_{-\infty}^{a} f(x) \, dx . \]  

(31.6)

Here and below, if \( x \) is discrete-valued, the integral is replaced by a sum. The endpoint \( a \) is expressly included in the integral or sum. Then \( 0 \leq F(x) \leq 1 \), \( F(x) \) is nondecreasing, and \( P(a < x \leq b) = F(b) - F(a) \). If \( x \) is discrete, \( F(x) \) is flat except at allowed values of \( x \), where it has discontinuous jumps equal to \( f(x) \).

Any function of random variables is itself a random variable, with (in general) a different p.d.f. The expectation value of any function \( u(x) \) is

\[ E[u(x)] = \int_{-\infty}^{\infty} u(x) \, f(x) \, dx , \]  

(31.7)
assuming the integral is finite. For \( u(x) \) and \( v(x) \), any two functions of \( x \), \( E[u + v] = E[u] + E[v] \). For \( c \) and \( k \) constants, \( E[cu + k] = cE[u] + k \).

The \( n \)th moment of a random variable is

\[
\alpha_n \equiv E[x^n] = \int_{-\infty}^{\infty} x^n f(x) \, dx , \tag{31.8a}
\]

and the \( n \)th central moment of \( x \) (or moment about the mean, \( \alpha_1 \)) is

\[
m_n \equiv E[(x - \alpha_1)^n] = \int_{-\infty}^{\infty} (x - \alpha_1)^n f(x) \, dx . \tag{31.8b}
\]

The most commonly used moments are the mean \( \mu \) and variance \( \sigma^2 \):

\[
\mu \equiv \alpha_1 , \tag{31.9a}
\]
\[
\sigma^2 \equiv V[x] \equiv m_2 = \alpha_2 - \mu^2 . \tag{31.9b}
\]

The mean is the location of the “center of mass” of the p.d.f., and the variance is a measure of the square of its width. Note that \( V[cx + k] = c^2 V[x] \). It is often convenient to use the standard deviation of \( x \), \( \sigma \), defined as the square root of the variance.

Any odd moment about the mean is a measure of the skewness of the p.d.f. The simplest of these is the dimensionless coefficient of skewness \( \gamma_1 = m_3/\sigma^3 \).

The fourth central moment \( m_4 \) provides a convenient measure of the tails of a distribution. For the Gaussian distribution (see Sec. 31.4), one has \( m_4 = 3\sigma^4 \). The kurtosis is defined as \( \gamma_2 = m_4/\sigma^4 - 3 \), i.e., it is zero for a Gaussian, positive for a leptokurtic distribution with longer tails, and negative for a platykurtic distribution with tails that die off more quickly than those of a Gaussian.

Besides the mean, another useful indicator of the “middle” of the probability distribution is the median, \( x_{med} \), defined by \( F(x_{med}) = 1/2 \), i.e., half the probability lies above and half lies below \( x_{med} \). (More rigorously, \( x_{med} \) is a median if \( P(x \geq x_{med}) \geq 1/2 \) and \( P(x \leq x_{med}) \geq 1/2 \). If only one value exists, it is called ‘the median.’)

Let \( x \) and \( y \) be two random variables with a joint p.d.f. \( f(x, y) \). The marginal p.d.f. of \( x \) (the distribution of \( x \) with \( y \) unobserved) is

\[
f_1(x) = \int_{-\infty}^{\infty} f(x, y) \, dy , \tag{31.10}
\]

and similarly for the marginal p.d.f. \( f_2(y) \). The conditional p.d.f. of \( y \) given fixed \( x \) (with \( f_1(x) \neq 0 \)) is defined by \( f_3(y|x) = f(x, y)/f_1(x) \), and similarly \( f_4(x|y) = f(x, y)/f_2(y) \). From these, we immediately obtain Bayes’ theorem (see Eqs. (31.2) and (31.4)),

\[
f_4(x|y) = \frac{f_3(y|x)f_1(x)}{f_2(y)} = \frac{f_3(y|x)f_1(x)}{\int f_3(y|x')f_1(x') \, dx'} . \tag{31.11}
\]
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The mean of \( x \) is

\[
\mu_x = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} x f(x, y) \, dx \, dy = \int_{-\infty}^{\infty} x f_1(x) \, dx ,
\]

and similarly for \( y \). The covariance of \( x \) and \( y \) is

\[
\text{cov}[x, y] = E[(x - \mu_x)(y - \mu_y)] = E[xy] - \mu_x\mu_y .
\]

A dimensionless measure of the covariance of \( x \) and \( y \) is given by the correlation coefficient,

\[
\rho_{xy} = \frac{\text{cov}[x, y]}{\sigma_x \sigma_y} ,
\]

where \( \sigma_x \) and \( \sigma_y \) are the standard deviations of \( x \) and \( y \). It can be shown that \(-1 \leq \rho_{xy} \leq 1\).

Two random variables \( x \) and \( y \) are independent if and only if

\[
f(x, y) = f_1(x)f_2(y) .
\]

If \( x \) and \( y \) are independent, then \( \rho_{xy} = 0 \); the converse is not necessarily true. If \( x \) and \( y \) are independent, \( E[u(x)v(y)] = E[u(x)]E[v(y)] \), and \( V[x + y] = V[x] + V[y] \); otherwise, \( V[x + y] = V[x] + V[y] + 2\text{cov}[x, y] \), and \( E[uv] \) does not necessarily factorize.

Consider a set of \( n \) continuous random variables \( x = (x_1, \ldots, x_n) \) with joint p.d.f. \( f(x) \), and a set of \( n \) new variables \( y = (y_1, \ldots, y_n) \), related to \( x \) by means of a function \( y(x) \) that is one-to-one, i.e., the inverse \( x(y) \) exists. The joint p.d.f. for \( y \) is given by

\[
g(y) = f(x(y))|J| ,
\]

where \( |J| \) is the absolute value of the determinant of the square matrix \( J_{ij} = \partial x_i/\partial y_j \) (the Jacobian determinant). If the transformation from \( x \) to \( y \) is not one-to-one, the \( x \)-space must be broken in to regions where the function \( y(x) \) can be inverted, and the contributions to \( g(y) \) from each region summed.

Given a set of functions \( y = (y_1, \ldots, y_m) \) with \( m < n \), one can construct \( n - m \) additional independent functions, apply the procedure above, then integrate the resulting \( g(y) \) over the unwanted \( y_i \) to find the marginal distribution of those of interest.

To change variables for discrete random variables simply substitute; no Jacobian is necessary because now \( f \) is a probability rather than a probability density. If \( f \) depends on a set of parameters \( \theta \), a change to a different parameter set \( \eta(\theta) \) is made by simple substitution; no Jacobian is used.
31.3. Characteristic functions

The characteristic function $\phi(u)$ associated with the p.d.f. $f(x)$ is essentially its Fourier transform, or the expectation value of $e^{ixu}$:

$$
\phi(u) = E[e^{ixu}] = \int_{-\infty}^{\infty} e^{ixu} f(x) \, dx .
$$

(31.17)

Once $\phi(u)$ is specified, the p.d.f. $f(x)$ is uniquely determined and vice versa; knowing one is equivalent to the other. Characteristic functions are useful in deriving a number of important results about moments and sums of random variables.

It follows from Eqs. (31.8a) and (31.17) that the $n^{th}$ moment of a random variable $x$ that follows $f(x)$ is given by

$$
i^{-n} \frac{d^n \phi}{du^n} \bigg|_{u=0} = \int_{-\infty}^{\infty} x^n f(x) \, dx = \alpha_n .
$$

(31.18)

Thus it is often easy to calculate all the moments of a distribution defined by $\phi(u)$, even when $f(x)$ cannot be written down explicitly.

If the p.d.f.s $f_1(x)$ and $f_2(y)$ for independent random variables $x$ and $y$ have characteristic functions $\phi_1(u)$ and $\phi_2(u)$, then the characteristic function of the weighted sum $ax + by$ is $\phi_1(ax)\phi_2(by)$. The additional rules for several important distributions (e.g., that the sum of two Gaussian distributed variables also follows a Gaussian distribution) easily follow from this observation.

Let the (partial) characteristic function corresponding to the conditional p.d.f. $f_2(x|z)$ be $\phi_2(u|z)$, and the p.d.f. of $z$ be $f_1(z)$. The characteristic function after integration over the conditional value is

$$
\phi(u) = \int \phi_2(u|z)f_1(z) \, dz .
$$

(31.19)

Suppose we can write $\phi_2$ in the form

$$
\phi_2(u|z) = A(u)e^{ig(u)z} .
$$

(31.20)

Then

$$
\phi(u) = A(u)\phi_1(g(u)) .
$$

(31.21)

The cumulants (semi-invariants) $\kappa_n$ are defined by

$$
\phi(u) = \exp \left[ \sum_{n=1}^{\infty} \frac{\kappa_n}{n!} (iu)^n \right] = \exp \left( i\kappa_1 u - \frac{1}{2} \kappa_2 u^2 + \ldots \right) .
$$

(31.22)
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The values $\kappa_n$ are related to the moments $\alpha_n$ and $m_n$. The first few relations are

\[
\begin{align*}
\kappa_1 &= \alpha_1 \ (= \mu, \text{the mean}) \\
\kappa_2 &= m_2 = \alpha_2 - \alpha_1^2 \ (= \sigma^2, \text{the variance}) \\
\kappa_3 &= m_3 = \alpha_3 - 3\alpha_1\alpha_2 + 2\alpha_1^2.
\end{align*}
\] (31.23)

31.4. Some probability distributions

Table 31.1 gives a number of common probability density functions and corresponding characteristic functions, means, and variances. Further information may be found in Refs. [1–8], [10], and [11], which has particularly detailed tables. Monte Carlo techniques for generating each of them may be found in our Sec. 33.4 and in Ref. 10. We comment below on all except the trivial uniform distribution.

31.4.1. Binomial distribution:

A random process with exactly two possible outcomes which occur with fixed probabilities is called a Bernoulli process. If the probability of obtaining a certain outcome (a “success”) in an individual trial is $p$, then the probability of obtaining exactly $r$ successes ($r = 0, 1, 2, \ldots, N$) in $N$ independent trials, without regard to the order of the successes and failures, is given by the binomial distribution $f(r; N, p)$ in Table 31.1. If $r$ and $s$ are binomially distributed with parameters ($N, r, p$) and ($N, s, p$), then $t = r + s$ follows a binomial distribution with parameters ($N, r + s, p$).

31.4.2. Poisson distribution:

The Poisson distribution $f(n; \nu)$ gives the probability of finding exactly $n$ events in a given interval of $x$ (e.g., space and time) when the events occur independently of one another and at an average rate of $\nu$ per the given interval. The variance $\sigma^2$ equals $\nu$. It is the limiting case $p \to 0, N \to \infty, Np = \nu$ of the binomial distribution. The Poisson distribution approaches the Gaussian distribution for large $\nu$.

In an accelerator experiment, for example, an opportunity for any pair of particles to come into collision and produce an event of a given type can be viewed as an independent Bernoulli trial. The total number of trials $N$ may be extremely large, but the number of such events that occur represents in general a tiny fraction, $p$, of this. Therefore, the number of events is well modeled as a Poisson variable.

31.4.3. Normal or Gaussian distribution:

The normal (or Gaussian) probability density function $f(x; \mu, \sigma^2)$ given in Table 31.1 has mean $E[x] = \mu$ and variance $V[x] = \sigma^2$. Comparison of the characteristic function $\phi(u)$ given in Table 31.1 with Eq. (31.22) shows that all cumulants $\kappa_n$ beyond $\kappa_2$ vanish; this is a unique property of the Gaussian distribution. Some other properties are:

\[
\begin{align*}
P(x \text{ in range } \mu \pm \sigma) &= 0.6827, \\
P(x \text{ in range } \mu \pm 0.6745\sigma) &= 0.5, \\
E[|x - \mu|] &= \sqrt{2/\pi}\sigma = 0.7979\sigma.
\end{align*}
\]
half-width at half maximum = $\sqrt{2 \ln 2} \sigma = 1.177 \sigma$.

For a Gaussian with $\mu = 0$ and $\sigma^2 = 1$ (the standard Gaussian), the cumulative distribution, Eq. (31.6), is related to the error function erf($y$) by

$$F(x; 0, 1) = \frac{1}{2} \left[ 1 + \text{erf}(x/\sqrt{2}) \right]. \quad (31.24)$$

The error function and standard Gaussian are tabulated in many references (e.g., Ref. [11]) and are available in software packages such as ROOT [12] and CERNLIB [13]. For a mean $\mu$ and variance $\sigma^2$, replace $x$ by $(x - \mu)/\sigma$. The probability of $x$ in a given range can be calculated with Eq. (32.45).

For $x$ and $y$ independent and normally distributed, $z = ax + by$ follows $f(z; a\mu_x + b\mu_y, a^2\sigma_x^2 + b^2\sigma_y^2)$; that is, the weighted means and variances add.

The Gaussian derives its importance in large part from the central limit theorem: If independent random variables $x_1, \ldots, x_n$ are distributed according to any p.d.f.s with finite means and variances, then the sum $y = \sum_{i=1}^{n} x_i$ will have a p.d.f. that approaches a Gaussian for large $n$. The mean and variance are given by the sums of corresponding terms from the individual $x_i$. Therefore, the sum of a large number of fluctuations $x_i$ will be distributed as a Gaussian, even if the $x_i$ themselves are not.

(Note that the product of a large number of random variables is not Gaussian, but its logarithm is. The p.d.f. of the product is log-normal. See Ref. [8] for details.)

For a set of $n$ Gaussian random variables $\mathbf{x}$ with means $\mu$ and corresponding Fourier variables $\mathbf{u}$, the characteristic function for a one-dimensional Gaussian is generalized to

$$\phi(\mathbf{u}; \mu, V) = \exp \left[ i \mu \cdot \mathbf{u} - \frac{1}{2} \mathbf{u}^T V \mathbf{u} \right]. \quad (31.25)$$

From Eq. (31.18), the covariance of $x_i$ and $x_j$ is

$$E [(x_i - \mu_i)(x_j - \mu_j)] = V_{ij}. \quad (31.26)$$

If the components of $\mathbf{x}$ are independent, then $V_{ij} = \delta_{ij}\sigma_i^2$, and Eq. (31.25) is the product of the c.f.s of $n$ Gaussians.

The characteristic function may be inverted to find the corresponding p.d.f.,

$$f(\mathbf{x}; \mu, V) = \frac{1}{(2\pi)^{n/2} \sqrt{|V|}} \exp \left[ -\frac{1}{2} (\mathbf{x} - \mu)^T V^{-1} (\mathbf{x} - \mu) \right], \quad (31.27)$$

where the determinant $|V|$ must be greater than 0. For diagonal $V$ (independent variables), $f(\mathbf{x}; \mu, V)$ is the product of the p.d.f.s of $n$ Gaussian distributions.

For $n = 2$, $f(\mathbf{x}; \mu, V)$ is

$$f(x_1, x_2; \mu_1, \mu_2, \sigma_1, \sigma_2, \rho) = \frac{1}{2\pi \sigma_1 \sigma_2 \sqrt{1 - \rho^2}}.$$
\begin{align*}
\times \exp \left\{ -\frac{1}{2(1-\rho^2)} \left[ \frac{(x_1 - \mu_1)^2}{\sigma_1^2} - \frac{2\rho(x_1 - \mu_1)(x_2 - \mu_2)}{\sigma_1 \sigma_2} + \frac{(x_2 - \mu_2)^2}{\sigma_2^2} \right] \right\}.
\end{align*}

(31.28)

The marginal distribution of any \( x_i \) is a Gaussian with mean \( \mu_i \) and variance \( V_{ii} \). \( V \) is \( n \times n \), symmetric, and positive definite. Therefore, for any vector \( \mathbf{X} \), the quadratic form \( \mathbf{X}^T \mathbf{V}^{-1} \mathbf{X} = C \), where \( C \) is any positive number, traces an \( n \)-dimensional ellipsoid as \( \mathbf{X} \) varies. If \( X_i = x_i - \mu_i \), then \( C \) is a random variable obeying the \( \chi^2 \) distribution with \( n \) degrees of freedom, discussed in the following section. The probability that \( \mathbf{X} \) corresponding to a set of Gaussian random variables \( x_i \) lies outside the ellipsoid characterized by a given value of \( C \) (= \( \chi^2 \)) is given by \( 1 - F_{\chi^2}(C; n) \), where \( F_{\chi^2} \) is the cumulative \( \chi^2 \) distribution. This may be read from Fig. 32.1. For example, the “s-standard-deviation ellipsoid” occurs at \( C = s^2 \). For the two-variable case \( (n = 2) \), the point \( \mathbf{X} \) lies outside the one-standard-deviation ellipsoid with 61% probability. The use of these ellipsoids as indicators of probable error is described in Sec. 32.3.2.4; the validity of those indicators assumes that \( \mu \) and \( V \) are correct.

31.4.4. \( \chi^2 \) distribution:

If \( x_1, \ldots, x_n \) are independent Gaussian random variables, the sum \( z = \sum_{i=1}^n (x_i - \mu_i)^2 / \sigma_i^2 \) follows the \( \chi^2 \) p.d.f. with \( n \) degrees of freedom, which we denote by \( \chi^2(n) \). More generally, for \( n \) correlated Gaussian variables as components of a vector \( \mathbf{X} \) with covariance matrix \( \mathbf{V} \), \( z = \mathbf{X}^T \mathbf{V}^{-1} \mathbf{X} \) follows \( \chi^2(n) \) as in the previous section. For a set of \( z_i \), each of which follows \( \chi^2(n_i) \), \( \sum z_i \) follows \( \chi^2(\sum n_i) \). For large \( n \), the \( \chi^2 \) p.d.f. approaches a Gaussian with mean \( \mu = n \) and variance \( \sigma^2 = 2n \).

The \( \chi^2 \) p.d.f. is often used in evaluating the level of compatibility between observed data and a hypothesis for the p.d.f. that the data might follow. This is discussed further in Sec. 32.2.2 on tests of goodness-of-fit.

31.4.5. Student’s \( t \) distribution:

Suppose that \( x \) and \( x_1, \ldots, x_n \) are independent and Gaussian distributed with mean 0 and variance 1. We then define

\[
z = \sum_{i=1}^n x_i^2 \quad \text{and} \quad t = \frac{x}{\sqrt{z/n}}.
\]

(31.29)

The variable \( z \) thus follows a \( \chi^2(n) \) distribution. Then \( t \) is distributed according to Student’s \( t \) distribution with \( n \) degrees of freedom, \( f(t; n) \), given in Table 31.1.

The Student’s \( t \) distribution resembles a Gaussian with wide tails. As \( n \to \infty \), the distribution approaches a Gaussian. If \( n = 1 \), it is a Cauchy or Breit–Wigner distribution. The mean is finite only for \( n > 1 \) and the variance is finite only for \( n > 2 \), so the central limit theorem is not applicable to sums of random variables following the \( t \) distribution for \( n = 1 \) or 2.

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As an example, consider the sample mean $\bar{x} = \frac{\sum x_i}{n}$ and the sample variance $s^2 = \frac{\sum (x_i - \bar{x})^2}{n-1}$ for normally distributed $x_i$ with unknown mean $\mu$ and variance $\sigma^2$. The sample mean has a Gaussian distribution with a variance $\sigma^2/n$, so the variable $(\bar{x} - \mu)/\sqrt{\sigma^2/n}$ is normal with mean 0 and variance 1. The quantity $(n-1)s^2/\sigma^2$ is independent of this and follows $\chi^2(n-1)$. The ratio

$$t = \frac{(\bar{x} - \mu)/\sqrt{\sigma^2/n}}{\sqrt{(n-1)s^2/\sigma^2(n-1)}} = \frac{\bar{x} - \mu}{\sqrt{s^2/n}}$$

(31.30)

is distributed as $f(t; n-1)$. The unknown variance $\sigma^2$ cancels, and $t$ can be used to test the probability that the true mean is some particular value $\mu$.

In Table 31.1, $n$ in $f(t; n)$ is not required to be an integer. A Student’s $t$ distribution with non-integral $n > 0$ is useful in certain applications.

31.4.6. Gamma distribution:

For a process that generates events as a function of $x$ (e.g., space or time) according to a Poisson distribution, the distance in $x$ from an arbitrary starting point (which may be some particular event) to the $k^{th}$ event follows a gamma distribution, $f(x; \lambda, k)$. The Poisson parameter $\mu$ is $\lambda$ per unit $x$. The special case $k = 1$ (i.e., $f(x; \lambda, 1) = \lambda e^{-\lambda x}$) is called the exponential distribution. A sum of $k'$ exponential random variables $x_i$ is distributed as $f(\sum x_i; \lambda, k')$.

The parameter $k$ is not required to be an integer. For $\lambda = 1/2$ and $k = n/2$, the gamma distribution reduces to the $\chi^2(n)$ distribution.
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Table 31.1. Some common probability density functions, with corresponding characteristic functions and means and variances. In the Table, $\Gamma(k)$ is the gamma function, equal to $(k-1)!$ when $k$ is an integer.

<table>
<thead>
<tr>
<th>Distribution</th>
<th>Probability density function $f$ (variable; parameters)</th>
<th>Characteristic function $\phi(u)$</th>
<th>Mean</th>
<th>Variance $\sigma^2$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Uniform</td>
<td>$f(x; a, b) = \begin{cases} 1/(b-a) &amp; a \leq x \leq b \ 0 &amp; \text{otherwise} \end{cases}$</td>
<td>$e^{i\mu u} - e^{i\mu u}/(b-a)iu$</td>
<td>$a + b$</td>
<td>$(b-a)^2/12$</td>
</tr>
<tr>
<td>Binomial</td>
<td>$f(r; N, p) = \frac{N!}{r!(N-r)!} p^r q^{N-r}$</td>
<td>$(q + pe^{iu})^N$</td>
<td>$Np$</td>
<td>$Npq$</td>
</tr>
<tr>
<td>Poisson</td>
<td>$f(n; \nu) = \frac{\nu^ne^{-\nu}}{n!}$; $n = 0, 1, 2, \ldots$; $\nu &gt; 0$</td>
<td>$\exp[\nu(e^{iu} - 1)]$</td>
<td>$\nu$</td>
<td>$\nu$</td>
</tr>
<tr>
<td>Normal (Gaussian)</td>
<td>$f(x; \mu, \sigma^2) = \frac{1}{\sigma\sqrt{2\pi}}\exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right)$</td>
<td>$\exp(i\mu u - \frac{1}{2}\sigma^2 u^2)$</td>
<td>$\mu$</td>
<td>$\sigma^2$</td>
</tr>
<tr>
<td>Multivariate Gaussian</td>
<td>$f(x; \mu, V) = \frac{1}{(2\pi)^{n/2}\sqrt{</td>
<td>V</td>
<td>}}\exp\left[i\mu \cdot u - \frac{1}{2}u^TVu\right]$</td>
<td>$\mu$</td>
</tr>
<tr>
<td>$\chi^2$</td>
<td>$f(z; n) = \frac{z^{n/2-1}e^{-z/2}}{2^{n/2}\Gamma(n/2)}$; $z \geq 0$</td>
<td>$(1 - 2iu)^{-n/2}$</td>
<td>$n$</td>
<td>$2n$</td>
</tr>
<tr>
<td>Student’s t</td>
<td>$f(t; n) = \frac{1}{\sqrt{n\pi}}\frac{\Gamma((n+1)/2)}{\Gamma(n/2)}\left(1 + \frac{t^2}{n}\right)^{-(n+1)/2}$</td>
<td>$-$</td>
<td>$0$ for $n \geq 2$</td>
<td>$n/(n-2)$ for $n \geq 3$</td>
</tr>
<tr>
<td>Gamma</td>
<td>$f(x; \lambda, k) = \frac{x^{k-1}\lambda^k e^{-\lambda x}}{\Gamma(k)}$; $0 &lt; x &lt; \infty$</td>
<td>$(1 - iu/\lambda)^{-k}$</td>
<td>$k/\lambda$</td>
<td>$k/\lambda^2$</td>
</tr>
</tbody>
</table>

**References:**


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31. Probability

13. The CERN Program Library (CERNLIB); see cernlib.web.cern.ch/cernlib.