The Expectation Value of a Random Variable:

The *expectation value* E[x] of a *random variable* x is the <u>mean</u> value of x, *i.e.* \hat{x} (aka μ).

For <u>discrete</u> x_i , E[x] is the sum of all x_i , weighted by their associated probabilities $P(x_i)$:

Discrete
$$x_i$$
: $E[x] \equiv \hat{x} \equiv \mu \equiv \sum_{i=1}^{N} x_i P(x_i)$

For <u>continuous</u> x, E[x] is the *integral* over all x, weighted by the **probability density function** of x, f(x):

Continuous x:
$$E[x] \equiv \hat{x} \equiv \mu \equiv \int x f(x) dx$$

<u>Note</u>: \hat{x} is <u>not</u> a *random variable* – it is a single, well-defined <u>number</u> that characterizes the *true mean* of the particular/specific distribution of a *random variable*.

It is clear that for *any* constant, *a*:

Discrete:
$$E[x+a] = \sum_{i=1}^{N} (x_i + a) P(x_i) = \sum_{i=1}^{N} x_i P(x_i) + a \sum_{i=1}^{N} P(x_i) = \hat{x} + a$$
Continuous: $E[x+a] = \int (x+a) f(x) dx = \int x f(x) dx + a \underbrace{\int f(x) dx}_{=1} = \hat{x} + a$

In fact, <u>any</u> function of x, H(x) has an expectation value:

Discrete:
$$E[H(x)] = \hat{H}(x) = \sum_{i=1}^{N} H(x_i) P(x_i)$$

Continuous:
$$E[H(x)] \equiv \hat{H}(x) \equiv \int H(x) f(x) dx$$

The *true mean*/expectation value $E[x] \equiv \hat{x} \equiv \mu$ is actually the *first moment* of the random variable x's probability distribution, *relative to/taken about the <u>origin</u>* x = 0.

The L^{th} moment of the random variable x taken about an <u>arbitrary</u> point x = c is defined as:

$$\mu_L \equiv E[(x-c)^L]$$

n.b. Moments taken about $x=0$ are known as algebraic moments.

Moments taken about the true mean $x=\mu$ are known as central moments.

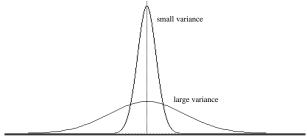
In particular, the 2^{nd} moment of x taken about/with respect to the **true mean** \hat{x} is:

$$\mu_2 \equiv E[(x - \hat{x})^2] \equiv \text{var}(x) \equiv \sigma_x^2 \equiv \text{the } variance \text{ of } x.$$

The quantity $\sigma_x = \sqrt{\mu_2} = \sqrt{\operatorname{var}(x)} = \sqrt{\sigma_x^2}$ is known as the <u>dispersion</u>, or <u>standard</u> <u>deviation</u> of x.

The *true mean* \hat{x} tells us where (in x) the distribution is "important" – i.e. \hat{x} tells us where the random variable x is most likely to be found).

The *dispersion/standard deviation* σ_x is a measure of the spread (or width) of the distribution over the space of x.

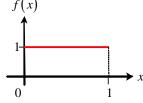


The two distributions in the above figure have the same *expectation value/true mean* $E[x] = \hat{x} = \mu$, but obviously have very different widths – and hence have different *variances*.

Example of the Uniform Distribution U(0,1) in the Range x=0 to x=1:

The Uniform Probability Density Function (P.D.F.) U(0,1) is:

$$f(x) = \begin{cases} 1 & \text{if } 0 \le x < 1 \\ 0 & \text{otherwise} \end{cases}$$



This is correctly "normalized", i.e. $\int_0^1 f(x) dx = \int_0^1 1 dx = \int_0^1 dx = 1$.

(i.e. the area under the curve of f(x) vs. x is = 1.)

The expectation value/*true mean* of U(0,1) is: $\hat{x} = \int_0^1 x \, f(x) \, dx = \int_0^1 x \, dx = \frac{1}{2}$

The *variance* associated with
$$U(0,1)$$
 is: $var(x) = \sigma_x^2 = \int_0^1 (x - \hat{x})^2 f(x) dx = \int_0^1 (x - \frac{1}{2})^2 f(x) dx$

We can certainly calculate the above *variance* σ^2 for U(0,1) directly, however, in general it is frequently often easier to calculate, if the *variance* is first transformed in the following way:

$$\operatorname{var}(x) = \sigma_{x}^{2} = \int (x - \hat{x})^{2} f(x) dx$$

$$= \int (x^{2} - 2x\hat{x} + \hat{x}^{2}) f(x) dx$$

$$= \int x^{2} f(x) dx - \int 2x\hat{x} f(x) dx + \int \hat{x}^{2} f(x) dx$$

$$= \int x^{2} f(x) dx - 2\hat{x} \int x f(x) dx + \hat{x}^{2} \int f(x) dx$$

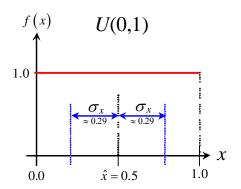
$$= E[x^{2}] - 2\hat{x} E[x] + \hat{x}^{2} = E[x^{2}] - 2\hat{x}^{2} + \hat{x}^{2} = E[x^{2}] - \hat{x}^{2}$$

$$\therefore \text{ for the } U(0,1) \text{ distribution: } \boxed{\operatorname{var}(x) = \sigma_x^2 = E[x^2] - E[x]^2 = \overline{x^2} - \overline{x}^2}$$

So for the U(0,1) distribution:

$$\operatorname{var}(x) = \sigma_x^2 = \int_0^1 x^2 dx - \left(\frac{1}{2}\right)^2 = \frac{1}{3} - \frac{1}{4} = \frac{1}{12}$$

$$\Rightarrow \sigma_x = \frac{1}{\sqrt{12}} \cong 0.29$$



This means that if we perform a measurement of a quantity x that results in the knowledge that it can be *anywhere* (with equal likelihood/equal/flat probability) within the range (0,1), our best estimate of its location is that is at the true mean, $x = \hat{x} = 0.5$ with a standard deviation of $\sigma_x = 1/\sqrt{12} \approx 0.29$.

More generally, for the flat/uniform distribution U(0,L) = 1/L:

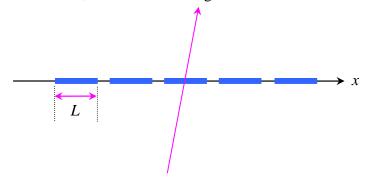
mean:
$$\hat{x} = \frac{L}{2}$$
 variance: $\sigma_x^2 = \frac{L^2}{12}$ \Rightarrow standard deviation: $\sigma_x = \frac{L}{\sqrt{12}}$

If the allowed range of x is $a \le x \le b$, it is also easy to show that U(a,b) = 1/(b-a), and that:

mean:
$$\hat{x} = \frac{b+a}{2}$$
 variance: $\sigma_x^2 = \frac{(b-a)^2}{12}$ \Rightarrow standard deviation: $\sigma_x = \frac{b-a}{\sqrt{12}}$

Example – The HEP Hodoscope:

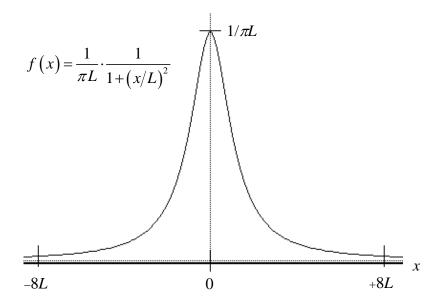
In particle physics experiments, a planar array of scintillation counters ("hodoscope") is often used to detect the passage of ionizing (charged) particles. If a particular counter is hit (and that is all that on one knows), then this information can be considered a measurement of (say) the x coordinate of the trajectory at the plane of the hodoscope. The result of the measurement is that \hat{x} is the x-coordinate of the center of the particular counter that was hit by the particle. The standard deviation of the measurement is $\sigma = L/\sqrt{12} \approx 0.29L$, where L is the lateral width (in the x-direction) of the counter, as shown in the figure below:



This example is also relevant for determining the uncertainty in ADC count data in the regime of noise-free data -e.g. a 12-bit ADC measures quasi-DC voltages over a ± 5.0 volt range. 2^{12} bits = 4096. Note that: 10 Volts/4096 bits = 2.44 mV/ADC count.

For noise-free {or nearly noise-free} ADC data, in the absence of using so-called ADC dithering techniques (adding and systematically {or randomly} varying a small voltage offset), then $\sigma_{ADC} = 1/\sqrt{12}$ ADC counts, corresponding to $\sigma_V = (1/\sqrt{12}) \cdot 2.44 mV = 0.70 mV$.

Example – The Cauchy Distribution:
$$f(x) = \frac{1}{\pi L} \cdot \frac{1}{1 + (x/L)^2}$$



A careful check shows that: $\int_{-\infty}^{+\infty} f(x) dx = 1 \iff \mathbf{OK}$

but:
$$\hat{x} = \int_{-\infty}^{+\infty} x f(x) dx = \frac{L}{2\pi} \ln(x^2 + L^2) \Big|_{-\infty}^{+\infty} \Leftarrow \text{Undefined !!!}$$

Obviously, the expectation value of x is/should be $\hat{x} = 0$. Hmmm... There are mathematical complications with this P.D.F. !!!

Using $\hat{x} = 0$, we can calculate the *variance* of the Cauchy Distribution (using a table of integrals):

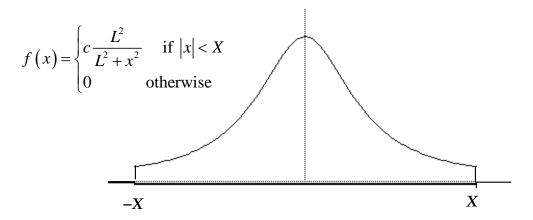
$$\operatorname{var}(x) = \sigma^{2} = \int_{-\infty}^{+\infty} x^{2} f(x) dx = \frac{L}{\pi} \left\{ x \Big|_{-\infty}^{+\infty} - L^{2} \int_{-\infty}^{+\infty} \frac{dx}{L^{2} + x^{2}} \right\} \quad \Leftarrow \quad \text{Infinite !!!}$$

$$\uparrow \qquad \qquad \uparrow \qquad \qquad \uparrow$$

$$\text{infinite} \qquad \text{finite}$$

More mathematical problems...!

In order to avoid these difficulties, in practice a truncated Cauchy distribution is used:



The requirement that: $\int_{-\infty}^{+\infty} f(x) dx = 1 \text{ requires: } c = \frac{1}{2L \tan^{-1}(X/L)}.$

(Note that for $X \to \infty$, $c \to 1/L\pi$ as we expect.)

Now:
$$\hat{x} = \int_{-X}^{+X} x f(x) dx = \frac{L}{2\pi} \ln(x^2 + L^2) \Big|_{-X}^{+X} = 0$$
 and $var(x) = \sigma^2$ is finite as well.

More About the Moments of a Probability Distribution Function / P.D.F.

The 1st central moment (i.e. taken about the true mean $\hat{x} = \mu$) is <u>not</u> interesting, since:

$$\mu_1 \equiv E[x - \hat{x}] = E[x] - E[\hat{x}] = \hat{x} - \hat{x} = 0$$

Higher moments of a probability distribution are *sometimes* useful.

For example – the *coefficients of <u>skewness</u>* and <u>kurtosis</u>, (Gr. "kurtos" = "bulging/swelling") respectively are the 3^{rd} and 4^{th} standardized *central* moments of a distribution (*i.e.* also taken about the *true mean* $\hat{x} = \mu$):

Skewness:
$$\gamma_1 = \frac{\mu_3}{\mu_2^{3/2}} = \frac{E[(x-\hat{x})^3]}{\{\underbrace{E[(x-\hat{x})^2]}\}^{3/2}} = \frac{E[(x-\hat{x})^3]}{\sigma^3}$$

Excess Kurtosis:
$$\gamma_2 = \frac{\mu_4}{\mu_2^{4/2}} - 3 = \frac{E[(x - \hat{x})^4]}{\{\underbrace{E[(x - \hat{x})^2]}\}^{4/2}} - 3 = \frac{E[(x - \hat{x})^4]}{\sigma^4} - 3$$

The -3 in the <u>excess</u> kurtosis definition is added so that $\gamma_2 = 0$ for a Gaussian (*aka* normal) probability distribution. Without the -3, "regular/normal" kurtosis is defined as $\mu_4/\mu_2^{4/2}$.

Skewness γ_1 tells us whether the P.D.F. f(x) is:

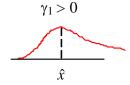
 $\gamma_1 < 0$: skewed toward $x < \hat{x}$

 $\gamma_1 = 0$: symmetric about \hat{x}

 $\gamma_1 > 0$: skewed toward $x > \hat{x}$





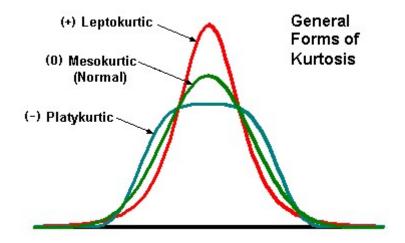


Obviously, the "long tail" dominates in $E[(x-\hat{x})^3]$.

A probability distribution (e.g. the Gaussian/normal distribution – see below...) that has **zero** kurtosis $(\gamma_2 = 0)$ is called **mesokurtic**, or mesokurotic (Gr. "**mesos**" = "middle").

Positive kurtosis ($\gamma_2 > 0$) indicates a peaked/slender/narrow probability distribution near the mean with abnormally **long** tails, and is called **leptokurtic**, or **leptokurotic** (Gr. "**leptos**" = "slender/thin").

Negative kurtosis ($\gamma_2 < 0$) indicates a flattened/wide/broad probability distribution near the mean with abnormally **shortened** tails, and is called **platykurtic**, or **platykurotic** (Gr. "**platys**" = "broad/flat").



Stock market investors are *very* interested in these particular higher moments (skewness & kurtosis) of various probability distribution functions...

Important Note:

<u>Each/every</u> measurement of a random variable x contains <u>information</u> about the PDF f(x) from which it originates – and its expectation value/true mean \hat{x} , variance $var(x) = \sigma_x^2$, and the PDF's higher order moments...

In the limit of an *infinite* # of measurements $N \to \infty$ of the random variable x, the PDF f(x) and all of its associated moments are *precisely* known...

<u>Any</u> random variable x can be "normalized" or "standardized" and can thus be reduced to a <u>dimensionless</u> quantity.

Suppose that the *random variable* x has P.D.F. f(x) and has a well-defined *expectation value* / *true mean* \hat{x} and *variance* σ_x^2 .

Consider a <u>new random variable</u> defined as: $u(x) \equiv \frac{(x-\hat{x})}{\sigma_x} \leftarrow n.b.$ dimensionless quantity

Aside: in general, if c is a **constant**, then: E[cg(x)] = cE[g(x)].

Thus: $E[u(x)] = \hat{u} = \int u(x) f(x) dx = \int \frac{x - \hat{x}}{\sigma_x} f(x) dx = \frac{1}{\sigma_x} \int (x - \hat{x}) f(x) dx = 0$, i.e. $\hat{u} = 0$ and:

$$\operatorname{var}(u) = \int (u - \hat{u})^{2} f(x) dx = \int u^{2}(x) f(x) dx - 2 \hat{u} \underbrace{\int u(x) f(x) dx}_{=\hat{u}} + \hat{u}^{2} \int f(x) dx = \int u^{2} f(x) dx$$

$$= \frac{1}{\sigma_{x}^{2}} \underbrace{\int (x - \hat{x})^{2} f(x) dx}_{=\operatorname{var}(x) = \sigma^{2}} = 1$$

Thus, the *random variable* $u(x) \equiv (x - \hat{x})/\sigma_x$ has:

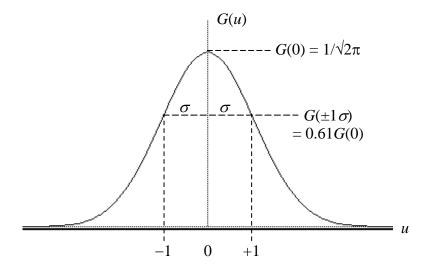
Expectation value/true mean $E[u(x)] = \hat{u} = 0$

Variance $var(u) = \sigma_u^2 = 1$, hence standard deviation $\sigma_u = 1$

The 1-D Gaussian (aka Normal) Distribution:

One of the most common P.D.F.'s that is encountered in the everyday world of doing experimental physics is the Gaussian (*aka* normal) probability distribution function (P.D.F.):

$$G(u) = \frac{1}{\sqrt{2\pi} \sigma} e^{-u^2/2\sigma^2} \quad \text{with: } \boxed{\sigma = 1}$$



It is easily verified that if u has the above P.D.F., then the *expectation value/true mean* of the Gaussian/normal distribution is E[u] = 0 (by symmetry) and its standard deviation, $\sigma_u = 1$.

In terms of probability, the probability that u will be found in an interval du is:

$$P = \begin{cases} \frac{1}{\sqrt{2\pi}} du & \text{at } u = 0\\ \frac{1}{\sqrt{2\pi}} e^{-1/2} du & \text{at } u = \pm \sigma \end{cases}$$
 The ratio $G(\pm \sigma)/G(0)$ is $e^{-1/2} = 0.6065... \approx 0.61$

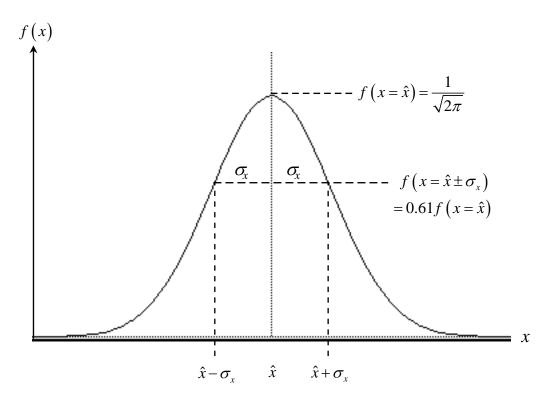
Later we shall show why the normal distribution is so universal in experimental physics...

In many situations in experimental physics, we need to describe a normally distributed **random variable** x with **expectation value** / **true mean** \hat{x} and standard deviation σ_x .

We can also *invert* the transformation $u = \frac{x - \hat{x}}{\sigma_x}$:

The P.D.F. of
$$x$$
 satisfies: $f(x)dx = G(u)du = \frac{1}{\sqrt{2\pi}}e^{-(x-\hat{x})^2/2\sigma_x^2}\frac{dx}{\sigma_x}$

i.e.
$$f(x) = \frac{1}{\sqrt{2\pi} \sigma_x} e^{-(x-\hat{x})^2/2\sigma_x^2} \iff \text{Shifted Gaussian distribution, centered at } x = \hat{x} \text{ with standard deviation } \sigma_x.$$



The *N*-D Probability Distributions Functions, *N*-D Cumulative Distribution Functions:

In physics experiments, we often measure more than one property of a system (e.g. the three spatial coordinates of a particle's position, its 3-momentum (or velocity), the energy and/or lifetime of an excited atomic state, etc.) An experiment may then result in k different random variables, which can be conveniently plotted on k mutually orthogonal axes (i.e. a mathematical hyper-space consisting of k orthogonal dimensions) and treated as if they were the components of a k-dimensional vector.

First, let us consider only \underline{two} such \underline{random} variables x and y.

The 2-D <u>Cumulative</u> Distribution Function: $F(X,Y) \equiv Prob(x < X \text{ and. } y < Y)$

The 2-D Probability Density Function:
$$f(x, y) = \frac{\partial}{\partial x} \frac{\partial}{\partial y} F(x, y)$$

If the *random variables* x and y are *independent*, then f(x, y) is <u>separable</u> in x & y, i.e. $f(x, y) = f_x(x) \cdot f_y(y) \iff x \& y$ have no <u>correlations</u> (i.e. are <u>uncorrelated</u>).

Then:
$$P(a \le x < b, c \le y < d) = \int_a^b dx \int_c^d dy f(x, y) \text{ with the requirement: } \int_{-\infty}^{+\infty} dx \int_{-\infty}^{+\infty} dy f(x, y) = 1$$

where the *x*, *y* integrals run over the allowed values of *x* and *y*: $-\infty \le a \le x \le b \le +\infty$ and: $-\infty \le c \le y \le d \le +\infty$.

In certain experimental situations, we may not care (or even know) what y is, hence we integrate y over its entire range and get: $g(x) = \int_{-\infty}^{+\infty} f(x, y) dy$.

In this situation:

$$P(a \le x < b, -\infty \le y < +\infty) = \int_a^b g(x) dx.$$

Thus we see that g(x) is <u>also</u> a P.D.F.; g(x) is known as the <u>marginal distribution</u> of x.

Change of *k* **Random Variables:**

Suppose we have k random variables $(x_1, x_2, ..., x_k)$, with corresponding P.D.F. $f(x_1, x_2, ..., x_k)$.

We may wish (or need) to make a change variables from $(x_1, x_2, ..., x_k)$ to $(y_1, y_2, ..., y_k)$, where <u>each</u> of the $y_i = y_i(x_1, x_2, ..., x_k)$.

For the case of a <u>single</u> random variable x, for a change of variables y = H(x), from P598AEM Lecture Notes 2, p. 8, the relation g(y)dy = f(x)dx leads to the relation (<u>here</u>):

$$g(y) = \frac{f(x)}{|H(x)|} \text{ where: } |H(x)| = |dy/dx| = |y\text{-slope}| \longleftarrow \boxed{n.b. \text{ absolute value}}$$

For *k* random variables, this generalizes to:

$$g(y_1, y_2, \dots, y_k) = \frac{f(x_1, x_2, \dots, x_k)}{|J|}$$
where J is the $\underline{Jacobian}$:
$$J\left(\frac{y_1y_2...y_k}{x_1x_2...x_k}\right) \equiv \begin{vmatrix} \frac{\partial y_1}{\partial x_1} & \frac{\partial y_1}{\partial x_2} & \dots & \frac{\partial y_1}{\partial x_k} \\ \frac{\partial y_2}{\partial x_1} & \frac{\partial y_2}{\partial x_2} & \dots & \frac{\partial y_2}{\partial x_k} \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\partial y_k}{\partial x_1} & \frac{\partial y_k}{\partial x_2} & \dots & \frac{\partial y_k}{\partial x_k} \end{vmatrix}$$

$$= \begin{bmatrix} n.b. \text{ The Jacobian } J \\ \text{is the } \underline{determinant} \text{ of the } k \times k \text{ matrix of derivatives } \partial y_i / \partial x_j \end{vmatrix}$$

i.e. compute the individual entries of the $k \times k$ matrix of partial derivatives $\partial y_i / \partial x_j$, then compute the determinant of this matrix, then take the absolute value of the determinant.