Analysis of Experimental Data

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1 Statistical Distributions

1.1 Statistics of Experiments

A primary purpose of science is to predict the results of experiments. Consider a simple experiment with five possible outcomes which we repeat ten times. If our theoretical prediction is that each of these five outcomes is equally probable, than our prediction for a typical series of ten experiments would be for each outcome to occur two times. Now suppose we perform the experiment ten times and present the results like this:

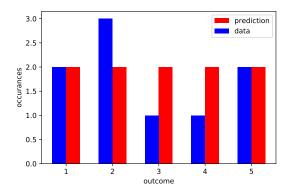


Figure 1: Comparison of experimental results with a prediction.

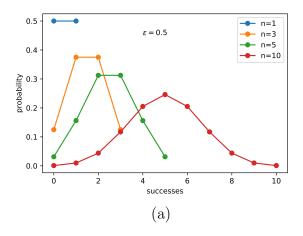
Scientist almost never display experimental data this way (as a bar graph) because it is nearly impossible to answer the crucial question is this data consistent with this prediction? Even if every outcome has an equal probability, the results of individual experiments experience statistical fluctuations. So even if the theory is correct, we well seldom reproduce exactly the theory prediction.

To interpret scientific experiments, it isn't enough to have a single prediction for the outcome of an experiment, instead, you need a prediction for the statistical distribution of outcomes: a probability distribution function. We'll start this discussion, therefore, by deriving three of the most frequently encountered probability distributions: the Binomial Distribution, the Poisson Distribution, and the Gaussian Distribution.

1.2 The Binomial Distribution

The Binomial Distribution is the most general of the distributions we'll consider, but it is a bit cumbersome to use in practice. The more familiar Poisson and Gaussian distributions are limiting

cases of this distribution.



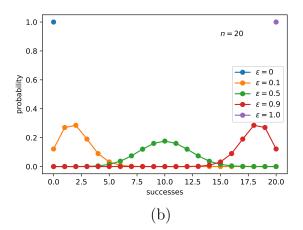


Figure 2: The binomial distribution for several different values of the parameters (a) n and (b) ϵ .

Suppose you repeat a particular process n times, and each time you have the same probability ϵ of a particular outcome, which, without losing generality, we'll call "success". The probability of having exactly m successes after n trials is simply given by:

$$P = \sum_{i} p_{i}$$

where i runs over all specific outcomes with m successes and p_i is the probability of each specific outcome. However, as these specific outcomes all contain exactly m successes, they share the same probability, namely:

$$p_i = \epsilon^m (1 - \epsilon)^{n - m}$$

and so we are left to consider simply the total number of specific outcomes containing m successes. The quantity we need is provided by the binomial theorem from mathematics, which states that:

$$(p+q)^n = \sum_{m=0}^n \binom{n}{m} p^m q^{n-m}$$
 (1)

where the binomial coefficients are defined by

$$\binom{n}{m} = \frac{n!}{m! (n-m)!} \tag{2}$$

and are also often referred to in other contexts as n-choose-m. The binomial coefficient simply tells us how many times we can choose m instances of p instead of q, from n factors, and so it is precisely the combinatoric factor that we need.

The probability of obtaining m successes after n trials with probability ϵ is therefore given by:

$$P(m; n, \epsilon) = \binom{n}{m} \epsilon^m (1 - \epsilon)^{n-m}$$
(3)

which is called the Binomial Distribution.

1.3 Mean and Variance

Given a probability distribution, the most urgent questions are generally "what is the mean value we can expect from this distribution?" and "how close to the mean value are most of the outcomes?" The first answer localizes the distribution while the second answer describes its width.

To calculate the mean value in answer to the first question, we simply calculate a weighted average:

$$\langle m \rangle \equiv \bar{m} \equiv \sum_{m} m P(m)$$
 (4)

For a continuous probability distribution, we would integrate instead:

$$\langle x \rangle \equiv \bar{x} \equiv \int x P(x) dx$$
 (5)

We usually answer the second question in terms of the variance, σ^2 , of the distribution:

$$\sigma^2 \equiv \langle (x - \bar{x})^2 \rangle$$

Other answers have problems, e.g. $\langle x - \bar{x} \rangle$ can be zero or nearly so, even for wide distributions, as long as it is symmetric. You could fix this by calculating $\langle |x - \bar{x}| \rangle$ but this is generally much harder to calculate, and less useful, than the variance. For instance, it is left as an exercise to show that:

$$\langle (x - \bar{x})^2 \rangle = \langle x^2 \rangle - \bar{x}^2 \tag{6}$$

using the fact that \bar{x} is simply a number, and so $\langle \bar{x} \rangle = \bar{x}$. We often write this result equivalently as:

$$\sigma^2 \equiv \langle (x - \langle x \rangle)^2 \rangle = \langle x^2 \rangle - \langle x \rangle^2 \tag{7}$$

Which shows explicitly that we need only calculate $\langle x \rangle$ and $\langle x^2 \rangle$ in order to determine the variance of a distribution.

1.4 Mean and Variance of the Binomial Distribution

The mean value of Binomial Distribution is given by:

$$\bar{m} = \sum_{m=0}^{n} m P(m)$$
$$= \sum_{m=0}^{n} m \binom{n}{m} \epsilon^{m} (1 - \epsilon)^{n-m}$$

which looks rather daunting! The trick is to use the Binomial Theorem (1) and define a function of two independent variables p and q given by:

$$f(p,q) = (p+q)^n = \sum_{m=0}^n \binom{n}{m} p^m q^{n-m}$$

We then calculate:

$$\frac{\partial f}{\partial p} = n(p+q)^{n-1} = \sum_{m=0}^{n} m \binom{n}{m} p^{m-1} q^{n-m}$$

and multiplying by p we have:

$$np(p+q)^{n+1} = \sum_{m=0}^{n} m \binom{n}{m} p^m q^{n-m}$$

which is true for any p and q. We now substitute the particular values $p = \epsilon$ and $q = 1 - \epsilon$ and find that:

$$n\epsilon = \sum_{m=0}^{n} m \binom{n}{m} \epsilon^m (1 - \epsilon)^{n-m} \equiv \sum_{m=0}^{n} m P(m) = \bar{m}$$

So the mean value is given by:

$$\bar{m} = n\epsilon$$
 (8)

or the total number of trials times the probability of success for each trial, a wholly plausible answer. For the variance, we use a variation of the same trick, this time using the second partial derivative:

$$p^{2} \cdot \frac{\partial^{2} f}{\partial p^{2}} = n(n-1)p^{2}(p+q)^{n-2} = \sum_{m=0}^{n} m(m-1) \binom{n}{m} p^{m} q^{n-m}$$

and again putting $p = \epsilon$ and $q = 1 - \epsilon$ to find that:

$$n(n-1)\epsilon^{2} = \sum_{m=0}^{n} (m^{2} - m) \binom{n}{m} p^{m} q^{n-m}$$
$$= \sum_{m=0}^{n} (m^{2} - m) P(m)$$
$$= \langle m^{2} - m \rangle = \langle m^{2} \rangle - \langle m \rangle$$

and as $\langle m \rangle = n\epsilon$ we have:

$$\langle m^2 \rangle = n(n-1)\epsilon^2 + n\epsilon$$

And so:

$$\sigma^2 = \langle m^2 \rangle - \langle m \rangle^2 = n(n-1)\epsilon^2 + n\epsilon - n^2\epsilon^2$$

or simply:

$$\sigma^2 = n \,\epsilon \,(1 - \epsilon) \tag{9}$$

Note that if $\epsilon = 0$ or $\epsilon = 1$, there is only one outcome (all failures or all success) and so the variation is zero.

1.5 The Poisson Distribution

Suppose we have some time interval over which we expect to observe a mean number of events λ . The events must be independent of one another: an event occurring at a particular time cannot affect the time at which the next event occurs. We divide the time interval over which the λ events are expected to occur into into n sub-intervals, each with an equal probability to contain an event. These intervals will be all the same size if the events are uniformly distributed in time, but if the events are not uniformly distributed, the sub-intervals are simply chosen to ensure the probability

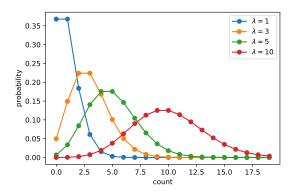


Figure 3: The Poisson distribution for several values of parameter λ .

is the same in each interval. Once cast this way, we can interpret this as a binomial distribution, with probability to contain an event, by construction, given by $\epsilon = \lambda/n$:

$$P(m) = \binom{n}{m} \epsilon^m (1 - \epsilon)^{n-m}$$

$$= \frac{n!}{m! (n - m)!} \left(\frac{\lambda}{n}\right)^m \left(1 - \frac{\lambda}{n}\right)^{n-m}$$

$$= \left(\frac{\lambda^m}{m!}\right) \left(1 - \frac{\lambda}{n}\right)^n \left[\frac{n!}{(n - m)!} \cdot \frac{1}{n^m}\right]_1 \left[\left(1 - \frac{\lambda}{n}\right)^{-m}\right]_2$$

We obtain the Poisson distribution by considering the limit that $n \to \infty$. It is left as an exercise to show that both $[\ldots]_1 \to 1$ and $[\ldots]_2 \to 1$ as $n \to \infty$. Recalling that

$$\lim_{n \to \infty} \left(1 - \frac{\lambda}{n} \right)^n = e^{-\lambda}$$

we obtain the Poisson distribution, the probability for observing m events for a mean of λ :

$$P(m; \lambda) = \frac{\lambda^m}{m!} e^{-\lambda} \tag{10}$$

Notice that there is no longer a parameter n, since we took $n \to \infty$, and so m now ranges from 0 to ∞ . Note also that λ is a real number, not necessarily an integer, even though it is the mean of integer values. For example, the mean of the integers 1 and 2 is the real value 1.5.

1.6 Mean and Variance of The Poisson Distribution

The mean of the Poisson distribution is given by:

$$\bar{m} = \sum_{m \ge 0} m P(m)$$
$$= \sum_{m \ge 0} m \frac{\lambda^m}{m!} e^{-\lambda}$$

Since the first term (m = 0) is zero, we have:

$$\bar{m} = e^{-\lambda} \sum_{m \ge 1} \frac{\lambda^m}{(m-1)!}$$

$$= \lambda e^{-\lambda} \sum_{m \ge 1} \frac{\lambda^{m-1}}{(m-1)!}$$

$$= \lambda e^{-\lambda} \sum_{n \ge 0} \frac{\lambda^n}{n!}$$

$$= \lambda e^{-\lambda} e^{\lambda}$$

$$\bar{m} = \lambda$$
(11)

which should come as no surprise, as the assumption in the derivation was the that mean number of events was λ .

For the variance, we use a similar manipulation to calculate:

$$\langle m^2 \rangle = \sum_{m \ge 0} m^2 P(m)$$

$$= \sum_{m \ge 0} m^2 \frac{\lambda^m}{m!} e^{-\lambda}$$

$$= \lambda \sum_{m \ge 1} m \frac{\lambda^{m-1}}{(m-1)!} e^{-\lambda}$$

$$= \lambda \sum_{n \ge 0} (n+1) \frac{\lambda^n}{(n)!} e^{-\lambda}$$

$$= \lambda \langle m+1 \rangle = \lambda (\lambda+1)$$

And so:

$$\sigma^{2} = \langle m^{2} \rangle - \langle m \rangle^{2}$$

$$= \lambda (\lambda + 1) - \lambda^{2}$$

$$\sigma^{2} = \lambda.$$
(12)

That is, the variance of a Poisson distribution is simply the mean.

1.7 The Gaussian Distribution

The last and most important probability distribution we will consider is the Gaussian distribution:

$$P(x) = \frac{1}{\sqrt{2\pi}\sigma} \cdot \exp\left(-\frac{(x-\lambda)^2}{2\sigma^2}\right) \tag{13}$$

which is plotted in Fig. 4. The Gaussian distribution (also sometimes called the normal distribution) appears frequently as a consequence of the central limit theorem, which states that the sum of independent random variables, when appropriately average (such as by taking the mean value) approaches a Gaussian distribution. As one example of the CLT, we'll show in the next section that the Gaussian distribution emerges from the Poisson distribution in the limit that $\lambda \to \infty$.

The Binomial and Poisson distributions are discrete functions: they describe the probability of outcomes which are integer quantities. The Poisson probability for having 2.31 events is either

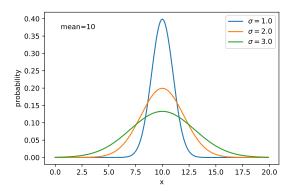


Figure 4: The Gaussian distribution for a mean of 10 and several values of parameter σ .

undefined or taken to be zero. The value of the Binomial and Poisson distribution function at a particular integer value is simply the probability of that particular outcome. To determine the probability that an outcome is within a range of integers, say m_1 to m_2 the probability distribution function is simply added:

$$P = \sum_{m=m_1}^{m_2} P(m).$$

The Gaussian distribution is a continuous function that describes the probability density at each position x. The probability of any particular outcome, say x = 1.24323, is vanishing small. But the probability that the value lies within a range of values is non-zero, and is determined by integrating:

$$P = \int_{x_{\min}}^{x_{\max}} P(x; \lambda, \sigma) \, dx$$

The Gaussian distribution is called a probability density function (PDF). A discrete probability distribution is sometimes referred to as probability mass function (PMF), to draw upon an analogy with physical mass and density: a PDF is probability per unit volume while a PMF is simply a probability.

The Gaussian distribution is normalized so that:

$$\int_{-\infty}^{+\infty} P(x) \, dx = 1$$

To determine the mean value of a function f(x) we likewise integrate:

$$\langle f(x) \rangle \equiv \int_{-\infty}^{+\infty} f(x)P(x) dx = 1$$

In particular the mean value of x is:

$$\langle x \rangle \equiv \int_{-\infty}^{+\infty} x P(x) \, dx = 1$$

and:

$$\langle x^2 \rangle \equiv \int_{-\infty}^{+\infty} x^2 P(x) \, dx = 1$$

We have defined the Gaussian distribution with the suggestive parameters λ and σ . It is left as an exercise to show that indeed, the mean is λ and the variance is σ^2 .

1.8 The Gaussian Distribution as a Limiting Case

As expected from the Central Limit Theorem, the Poisson Distribution in the limit $\lambda \to \infty$ approaches the Gaussian distribution. In this case, we can apply the Stirling Approximation:

$$\lim_{n \to \infty} n! = \sqrt{2\pi n} \ e^{-n} \ n^n$$

to the Poisson distribution as follows:

$$P(m) = \frac{\lambda^m}{m!} e^{-\lambda}$$

$$\rightarrow \frac{\lambda^m e^{-\lambda}}{\sqrt{2\pi m}} e^{-m} m^m$$

$$= \frac{e^{m-\lambda}}{\sqrt{2\pi \lambda}} \left(\frac{\lambda}{m}\right)^{m+\frac{1}{2}}$$

Now we consider a new variable δ , defined by

$$\delta \equiv \frac{m - \lambda}{\lambda}$$

which measures the difference between the observed number of events m and the mean of the distribution, as a fraction of the mean. Intuitively, the function is getting very narrow, and so we expect this to be a small quantity, but let's check this. First we have:

$$\langle \delta \rangle = \frac{\langle m \rangle - \lambda}{\lambda} = \frac{\lambda - \lambda}{\lambda} = 0$$

but also:

$$\langle \delta^2 \rangle = \frac{\langle (m-\lambda)^2 \rangle}{\lambda^2} = \frac{\lambda}{\lambda^2} = \frac{1}{\lambda}$$

where we have used the fact that the variance is given by $\langle (m-\lambda)^2 \rangle = \lambda$, and so as $\lambda \to \infty$ we have

$$\langle \delta^2 \rangle \to 0$$

So we can write:

$$m = \lambda(1+\delta) \tag{14}$$

where we expect the approximation $\delta \to 0$ to hold as long as we require $\lambda \to \infty$. So now we can write the distribution in terms of the small quantity δ and the large quantity λ as:

$$P(\delta) = \frac{e^{\lambda \delta}}{\sqrt{2\pi \lambda}} \left(\frac{\lambda}{\lambda(1+\delta)}\right)^{\lambda(1+\delta)+\frac{1}{2}}$$
$$= \frac{e^{\lambda \delta}}{\sqrt{2\pi \lambda}} \cdot \frac{1}{X}$$
(15)

where we define the quantity:

$$X = (1+\delta)^{\lambda(1+\delta) + \frac{1}{2}}$$

which can be approximated as follows:

$$\ln X = \left(\lambda(1+\delta) + \frac{1}{2}\right) \cdot \ln(1+\delta)$$

$$= \left(\lambda(1+\delta) + \frac{1}{2}\right) \cdot \left(\delta - \frac{\delta^2}{2} + \mathcal{O}(\delta^3)\right)$$

$$\frac{\ln X}{\lambda} = \left(1 + \delta + \frac{1}{2\lambda}\right) \cdot \left(\delta - \frac{\delta^2}{2} + \mathcal{O}(\delta^3)\right)$$

$$= \left(1 + \delta + \mathcal{O}\left(\delta^2\right)\right) \cdot \left(\delta - \frac{\delta^2}{2} + \mathcal{O}(\delta^3)\right)$$

$$= \delta + \frac{\delta^2}{2} + \mathcal{O}(\delta^3),$$

where in the second to last step we used $\mathcal{O}\left(\frac{1}{\lambda}\right) \sim \mathcal{O}\left(\delta^2\right)$. Neglecting the small quantities, we can approximate

$$X = \exp\left(\lambda\delta + \lambda \frac{\delta^2}{2}\right)$$

which, when plugged backed into Equation 15 yields:

$$P(\delta) = \frac{e^{\lambda \delta}}{\sqrt{2\pi \lambda}} \cdot \frac{1}{\exp\left(\lambda \delta + \lambda \frac{\delta^2}{2}\right)}$$
$$= \frac{1}{\sqrt{2\pi \lambda}} \cdot \exp\left(-\lambda \frac{\delta^2}{2}\right)$$
(16)

Recalling that:

$$\delta \equiv \frac{m - \lambda}{\lambda}$$

$$P(x) = \frac{1}{\sqrt{2\pi\lambda}} \cdot \exp\left(-\frac{(m - \lambda)^2}{2\lambda}\right)$$
(17)

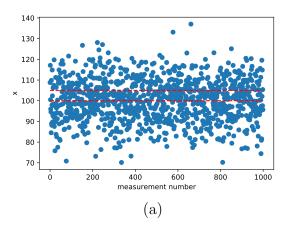
Which we can identify as a special case of the Gaussian distribution of Equation 13 with $\sigma^2 = \lambda$ and x = m:

$$P(x) = \frac{1}{\sqrt{2\pi}\sigma} \cdot \exp\left(-\frac{(x-\lambda)^2}{2\sigma^2}\right)$$

It may trouble you that x is continuous while m here was an integer. But note that in the limit $\lambda \to \infty$, the difference between m/λ and $(m+1)/\lambda$ is infinitesimal, and so this is effectively a continuous function, consistent with the Gaussian distribution.

1.9 Histograms in Scientific Python

Suppose a particular variable x is measured 1000 times. One way to visualize the collected data is shown in Fig. 5a, which simply plots each measurement value above the measurement number (from 0 to 1000). In this example, the number of measurements that occur within the range from x = 100 to x = 105 is 181. This count is plotted as the red data point in Fig. 5b, 181 entries located above x = 102.5, the center of the range. If we repeat this exercise across a number or ranges,



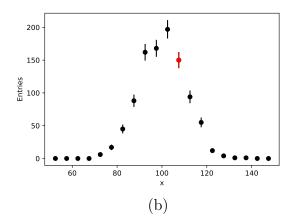


Figure 5: The 1000 measurements of variable x in (a) are used to produce the histogram in (b). The red data point in (b) is the count of the number of entries in range indicated by the red dashed lines in (a).

the resulting plot in Fig. 5b is called a histogram. A histogram reports the number of entries that occur within each of a sequence of consecutive ranges of x-values. Each range considered is called a histogram bin, and the choice of which bins to use is at the discretion of the analyzer.

The content of each bin is a single number, a count, and is therefore subject to the Poisson distribution. We can estimate the mean of the Poisson distribution by the measured value, so that $\lambda = N$. For the Poisson distribution, the variance $\sigma^2 = \lambda$, and so $\sigma = \sqrt{N}$. It is customary to draw a line of size \sqrt{N} when plotting a histogram value N. This is an example of an error bar, which indicates how well our measurement has determined a particular value.

An example producing a histogram in Scientific Python is shown in Fig. 6. The data to plot is simply a sequence of 1000 values randomly and uniformly chosen in the range [0, 20]:

```
x = np.random.uniform(high=20.0,size=1000)
```

To create a histogram from these 1000 values, we use the np.histogram function:

```
counts, edges = np.histogram(x,bins=10,range=(0,20))
```

where we have specified 10 bins, uniformly covering the range from 0 to 20. The function returns to arrays, which we save as counts and edges. The counts array contains the bin contents, the count of the number of values in each bin:

```
counts: [ 92 82 123 96 85 106 105 99 99 113]
```

The edges array contains the edges of the bins:

```
edges: [ 0. 2. 4. 6. 8. 10. 12. 14. 16. 18. 20.]
```

You'll notice that 10 consecutive bins have 11 edges. For plotting continuous data, one choice is to plot the contents at the center of each bin:

```
cbins = (edges[:-1] + edges[1:])/2.0
```

the two slices edges[:-1] and edges[1:] are all but the last and all but the first. The average of the two is the center of each bin:

```
x = np.random.uniform(high=20.0,size=1000)
counts,edges = np.histogram(x,bins=10,range=(0,20))
print("counts: ", counts)
print("edges: ", edges)
cbins = (edges[:-1] + edges[1:])/2.0
err = counts**0.5
print("cbins: ", cbins)
plt.errorbar(cbins,counts,yerr=err,fmt="ko")
plt.xlabel("x")
plt.ylabel("Entries")
```

```
99 102 110
                                                 98
counts:
           [ 81
                 96 104 108 104
                                                     98]
edges:
                      4.
                          6.
                               8. 10. 12. 14. 16. 18. 20.]
           [ 0.
cbins:
           [ 1.
                 3.
                      5.
                          7.
                               9. 11. 13. 15. 17. 19.]
```

Text(0,0.5,'Entries')

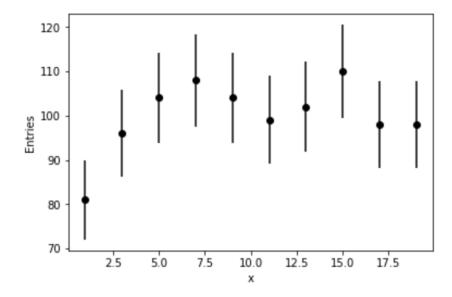


Figure 6: Example producing a histogram in Scientific Python.

cbins: [1. 3. 5. 7. 9. 11. 13. 15. 17. 19.]

The error bar values are chosen as the square root of the bin values:

err = counts**0.5

The histogram is plotted using the plt.errorbar function:

plt.errorbar(cbins,counts,yerr=err,fmt="ko")

which plots the bin contents counts at the bin center values cbins using the square root error bars in the array err, using the format "ko" for black circles.

1.10 Comparing a Histogram to a Probability Distribution Function

Our theoretical models often predict a PDF for some observable variable x. As experimentalists, we are often therefore concerned with the question as to whether our collected data for an observable x is consistent with the theoretical PDF. A visual approach to answering this question is to plot the data in a histogram, and to draw the PDF as a curve normalized to the histogram.

To predict the number of events in a bin with edges x_{lower} and x_{upper} , in principle we need to integrate the PDF and normalize to the number of experiments:

$$N_{\text{pred}} = N_{\text{meas}} \int_{x_{\text{lower}}}^{x_{\text{upper}}} p(x) \, dx$$

In practice, we generally choose the bin sizes small enough that the PDF is approximately constant during the entire bin, and in this case, the prediction can be taken as:

$$N_{\text{pred}} = N_{\text{meas}} \, \Delta x \, p(x)$$

where Δx is the width of each bin. This scale factor $N_{\text{meas}} \Delta x$ allows us to compare a continuous function to data collected in discrete bins, as shown in Fig. 7.

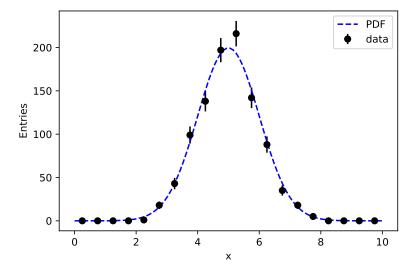


Figure 7: The Gaussian PDF scaled to compare to data from a Gaussian distribution. In this case, there are 1000 total entries in the histogram and the bin size is 0.5, for a scale factor of 500.

1.11 Homework Exercises for Distributions

Problem 1: Show that the Binomial distribution, P(m), in Equation 3 is properly normalized:

$$\sum_{m=0}^{n} P(m) = 1$$

as a consequence of the Binomial Theorem (Equation 1).

Problem 2: Show that Equation 6 is correct.

Problem 3: Show that the Poisson distribution, P(m), in Equation 10 is properly normalized:

$$\sum_{m \ge 0} P(m) = 1.$$

Hint: recall the Taylor series expansion for e^{λ} .

Problem 4: Show that the Gaussian distribution, P(x), in Equation 13 is properly normalized:

$$\int_{-\infty}^{\infty} P(x)dx = 1.$$

Problem 5: Show that the mean of the Gaussian distribution has been correctly identified in Equation 13. That is, show explicitly that:

$$\int_{-\infty}^{\infty} x P(x) dx = \lambda$$

Problem 6: Show that the variance of the Gaussian distribution has been correctly identified in Equation 13. That is, show explicitly that:

$$\int_{-\infty}^{\infty} x^2 P(x) dx = \sigma^2$$

when we take $\lambda=0$ (which is equivalent to simply changing variables $y=x-\lambda.$)

2 Statistical Uncertainties

2.1 Reporting Experimental Uncertainties

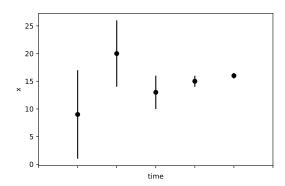


Figure 8: The evolution of a measurement of a quantity x over time.

A scientific measurement is meaningless without an associated uncertainty. For instance, we might report a measured distance as:

$$x = 1.2 \pm 0.3 \text{ m}$$

and a time as:

$$t = 4.76 \pm 0.13 \ \mu s.$$

When reporting a number with an associated uncertainty, we round both the value and the uncertainty to one significant digit in the uncertainty, so:

$$x = 1.245 \pm 0.313 \text{ m} \rightarrow x = 1.2 \pm 0.3 \text{ m}$$

If the first nonzero digit in the uncertainty is a zero or one, we keep two digits instead:

$$x=1.245\pm0.113~\mathrm{m} \rightarrow x=1.25\pm0.11~\mathrm{m}.$$

We'll define this uncertainty quite precisely, but let's start with a conceptual definition. Each particular measurement of a quantity x makes a claim about where the true value of x is most likely to lie. If we assume that measurements improve with time and better reproduce the true value, we expect a series of measurements to proceed like in Fig. 8, with future measurements shown to be consistent with the uncertainties reported by previous experiments.

We can therefore think of a measurement and it's uncertainty as describing a PDF for the outcome of experiments in the distant future with uncertainties so small as to effectively measure the true value of the quantity. Because most measured quantities have many independent sources of uncertainties, the Central Limit Theorem implies that the true value of the measured quantity is best described by a Gaussian Distribution. We define the measured value x as the mean of this distribution, and the uncertainty as the σ of the Gaussian.

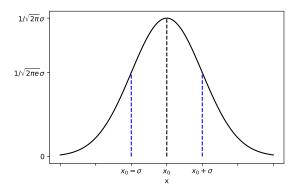


Figure 9: The interpretation of experimental uncertainties as the parameters of a Gaussian distribution for the true value.

2.2 The Error Function

Suppose we measure the speed of light in vacuum to be:

$$3.3 \pm 0.2 \times 10^8 \text{ m/s}^2$$

How consistent is this with the generally accepted value:

$$3.0 \times 10^8 \text{ m/s}^2$$

where we have omitted the uncertainty here because it is much smaller than our uncertainty! The interpretation of the uncertainty as the σ of a Gaussian distribution allows us to give quite precise answers. To begin we might say that it is consistent within:

$$\frac{3.3 - 3.0}{0.2} = 1.5 \ \sigma.$$

but we can also ask, what is the probability enclosed in 1.5σ . To answer this question, we use the error function:

$$\operatorname{erf}(x) = \frac{1}{\sqrt{\pi}} \int_{-x}^{x} \exp(-t^{2}) dt$$

which is simply the integral from -x to x of a Gaussian with $\sigma = 1/sqrt2$. So if we want to calculate the probability contained in n sigma we would calculate:

$$\operatorname{erf}(n \cdot \frac{1}{\sqrt{2}})$$

which is tabulated in Table 1.

2.3 Mean and Variance of a Sum

So far we've been calculating expectation values of functions of one random variable:

$$\langle f(x) \rangle = \int f(x) P(x) dx$$

interval	error function	integrated probability
$\pm 1\sigma$	$\operatorname{erf}\left(\frac{1}{\sqrt{2}}\right)$	68.3%
$\pm 2\sigma$	$\operatorname{erf}\left(\frac{2}{\sqrt{2}}\right)$	95.4%
$\pm 3\sigma$	$\operatorname{erf}\left(\frac{3}{\sqrt{2}}\right)$	98.5%
$\pm 4\sigma$	$\operatorname{erf}\left(\frac{4}{\sqrt{2}}\right)$	99.5%
$\pm 5\sigma$	$\operatorname{erf}\left(\frac{5}{\sqrt{2}}\right)$	99.8%

Table 1: The integrated probability for a Gaussian distribution within the stated bounds.

but this easily generalizes to two or more random variables:

$$\langle f(x,y)\rangle = \int \int f(x,y) P(x,y) dx dy$$

If the random variables are independent, then we have P(x,y) = P(x) P(y) and so

$$\langle f(x,y)\rangle = \int f(x,y) P(x) P(y) dx dy$$

And also we have

$$\langle f(x) + g(y) \rangle = \int \int dx \, dy \, P(x) \, P(y) \, (f(x) + g(y))$$

$$= \left(\int P(y) \, dy \right) \cdot \left(\int P(x) \, f(x) \, dx \right) + \left(\int P(x) \, dx \right) \cdot \left(\int P(y) \, f(y) \, dy \right)$$

$$= 1 \cdot \int P(x) \, f(x) \, dx + 1 \cdot \int P(y) \, f(y) \, dy$$

which we can write much more simply as:

$$\langle f(x) + g(y) \rangle = \langle f(x) \rangle + \langle g(y) \rangle$$

revealing the power of the expectation value notation for tackling independent random variables. In the homework you will show that similarly:

$$\langle f(x) \cdot g(y) \rangle = \langle f(x) \rangle \cdot \langle g(y) \rangle$$
 (18)

We can therefore conclude that the mean of a sum of two random variables:

$$\langle x + y \rangle = \langle x \rangle + \langle y \rangle$$

is just the sum of the mean.

For simplicity let's assume that $\langle x \rangle = \langle y \rangle = 0$, and so:

$$\langle (x+y)^2 \rangle = \langle x^2 + 2xy + y^2 \rangle$$

$$= \langle x^2 \rangle + 2 \langle xy \rangle + \langle y^2 \rangle$$

$$= \langle x^2 \rangle + 2 \langle x \rangle \langle y \rangle + \langle y^2 \rangle$$

$$= \langle x^2 \rangle + \langle y^2 \rangle$$

from which it follows that:

$$\sigma^{2}(x+y) = \langle (x+y)^{2} \rangle - (\langle x+y \rangle)^{2}$$

$$= \langle (x+y)^{2} \rangle$$

$$= \langle x^{2} \rangle + \langle y^{2} \rangle$$

$$= \sigma_{x}^{2} + \sigma_{y}^{2}$$

In the exercises you will show that even if $\langle x \rangle$ and $\langle y \rangle$ are non-zero, we still obtain:

$$\sigma^2(x+y) = \sigma_x^2 + \sigma_y^2$$

that is, the uncertainties in a sum add in quadrature.

2.4 Uncertainties Add in Quadrature

In the previous section, we discussed that the variance of the sum of two random variables drawn from any distribution is just the sum of their variances:

$$\sigma^2(x+y) = \sigma_x^2 + \sigma_y^2$$

This is a quite general result. In the context of experimental uncertainties, we consider experimental measurements to be drawn from a Gaussian distribution and the square-root of the variance (σ) is called the uncertainty.

If we calculate the sum s of two independent measurements x and y with uncertainties σ_x and σ_y , it's clear that the variance of the sum will be simply $\sigma_x^2 + \sigma_y^2$. If the resulting distribution of x + y is a Gaussian distribution, as seems probable due to the Central Limit Theorem, then the uncertainties simply add in quadrature:

$$\sigma_{x+y} = \sqrt{\sigma_x^2 + \sigma_y^2}.$$

Let's show explicitly that this is the case. If we wish to know the probability that two random variables x and y add to some particular value u = x + y, we simply integrate the total probability of x and y subject to the requirement u = x + y:

$$P(u) = \int dx \int dy P_x(x) P_y(y) \delta(u - (x+y))$$
$$= \int dx P_x(x) P_y(u-x)$$

If we make the (very often valid) assumption that x and y are Gaussian distributed, and, for simplicity, assume that the mean values are zero (or simply change coordinates), so that we have:

$$P_x(x) = \frac{1}{\sqrt{2\pi a}} \exp\left(-\frac{x^2}{2a^2}\right)$$

$$P_y(y) = \frac{1}{\sqrt{2\pi a}} \exp\left(-\frac{y^2}{2b^2}\right)$$

And so the mean value probability distribution function for u is now:

$$P(u) = \frac{1}{2\pi ab} \int dx \exp\left(-\frac{x^2}{2a^2}\right) \exp\left(-\frac{(u-x)^2}{2b^2}\right)$$
$$= \frac{1}{2\pi ab} \int dx \exp\left(-\frac{a^2+b^2}{2a^2b^2} \left\{x^2 - \frac{2a^2}{a^2+b^2}ux + \frac{a^2}{a^2+b^2}u^2\right\}_1\right)$$

We deal with the term in brackets $(\{\}_1)$ by completing the square. Simply note that

$$\left(x - \frac{a^2}{a^2 + b^2}u\right)^2 = x^2 - \frac{2a^2}{a^2 + b^2}ux + \frac{a^4}{(a^2 + b^2)^2}u^2 \tag{19}$$

reproduces the first and second terms, so we can replace:

$$x^{2} - \frac{2a^{2}}{a^{2} + b^{2}}ux = \left(x - \frac{a^{2}}{a^{2} + b^{2}}u\right)^{2} - \frac{a^{4}}{(a^{2} + b^{2})^{2}}u^{2}$$
(20)

to obtain:

$$\{\}_1 = \left(x - \frac{a^2}{a^2 + b^2}u\right)^2 - \frac{a^4}{(a^2 + b^2)^2}u^2 + \frac{a^2}{a^2 + b^2}u^2$$
$$= \left(x - \frac{a^2}{a^2 + b^2}u\right)^2 + \frac{a^2b^2}{(a^2 + b^2)^2}u^2$$

and substituting back into the original expression we obtain:

$$P(u) = \frac{1}{2\pi ab} \left\{ \int_{-\infty}^{+\infty} dx \, \exp\left(-\frac{a^2 + b^2}{2a^2b^2} \left(x - \frac{a^2}{a^2 + b^2}u\right)^2\right) \right\}_2 \exp\left(-\frac{1}{2}\frac{u^2}{a^2 + b^2}\right)$$

making a substitution of variables (u is constant during the integration):

$$y = x - \frac{a^2}{a^2 + b^2}u$$

the definite integral in brackets yields:

$$\{\}_2 = \sqrt{(2\pi)} \frac{ab}{\sqrt{a^2 + b^2}} \tag{21}$$

and we have at last:

$$P(u) = \frac{1}{\sqrt{2\pi}\sqrt{a^2 + b^2}} \exp\left(-\frac{1}{2}\frac{u^2}{a^2 + b^2}\right)$$

which is a Gaussian distribution with variance:

$$\sigma^2 = a^2 + b^2,$$

that is, uncertainties add in quadrature.

2.5 Handling Constants

When adding a known constant to a measured quantity:

$$y = x + C$$

we can consider the constant to have uncertainty $\sigma = 0$. And therefore, from the addition in quadrature rule:

$$\sigma_y = \sigma_x$$

If we multiply a measured quantity by a known constant

$$y = kx$$

we consider that:

$$dx \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right) = d(kx) \frac{1}{\sqrt{2\pi}k\sigma} \exp\left(-\frac{(kx-k\mu)^2}{2(k\sigma)^2}\right)$$

and conclude that:

$$\sigma_y = k\sigma_x$$

2.6 General Propagation of Uncertainties

We are now ready to handle propagation of uncertainties for a general function. Suppose we measure $x = x_0 \pm \sigma_x$ and $y = y_0 \pm \sigma_y$ and we wish to know the resulting uncertainty on the calculated quantity f(x, y).

Now we Taylor expand the function about the measured values x_0 and y_0 :

$$f(x,y) = f(x_0 + \Delta x, y_0 + \Delta y) \sim f(x_0, y_0) + \left. \frac{df}{dx} \right|_{x_0, y_0} \Delta x + \left. \frac{df}{dy} \right|_{x_0, y_0} \Delta y$$

We note that

$$\Delta x \equiv x - x_0$$
$$\Delta y \equiv y - y_0$$

where we can consider x_0 and y_0 to be precisely known constants that happen to coincide with our best measured value of x and y. In this case, our measurement amounts to:

$$\Delta x = 0 \pm \sigma_x$$
$$\Delta y = 0 \pm \sigma_y$$

and our uncertainties are associated with Δx and Δy , not the "constants" x_0 and y_0 .

Turning back to the Taylor expansion, we see now that:

$$f(x,y) \sim A + B\Delta x + C\Delta y$$

where the constants are:

$$A = f(x_0, y_0)$$

$$B = \frac{df}{dx}\Big|_{x_0, y_0}$$

$$C = \frac{df}{dy}\Big|_{x_0, y_0}$$

We know how to propagate uncertainties in this case, which involves scaling, adding in quadrature, and adding a constant:

$$\sigma_f^2 = B^2 \sigma_r^2 + C^2 \sigma_r^2$$

And plugging in the constants:

$$\sigma_f^2 = \left(\frac{df}{dx} \Big|_{x_0 y_0} \right)^2 \sigma_x^2 + \left(\frac{df}{dy} \Big|_{x_0 y_0} \right)^2 \sigma_y^2$$

2.7 Variance of the Mean

Suppose we have an apparatus that can measure a quantity x with uncertainty σ , and we make a series of N measurements $x_1, x_2, x_3, \ldots, x_N$.

We can calculate the mean value:

$$X = \frac{x_1 + x_2 + x_3 + \ldots + x_N}{N}$$

The uncertainty on the mean value can be calculated by propagating the uncertainties in the usual way:

$$\sigma_X = \sqrt{\left(\frac{dX}{dx_1}\sigma\right)^2 + \left(\frac{dX}{dx_2}\sigma\right)^2 + \dots + \left(\frac{dX}{dx_N}\sigma\right)^2}$$

$$= \sqrt{N\left(\frac{\sigma}{N}\right)^2}$$

$$= \frac{\sigma}{\sqrt{N}}$$

If we assume that the mean value is approaching the true value of the quantity x, then repeating the measurement N times reduces the uncertainty by a factor \sqrt{N} .

It would seem, therefore, that by simply repeating the same experiment many times, we could reach very high precision even with very lousy equipment. But this is, unfortunately, not generally the case. The statistical uncertainties which we have considered here do indeed diminish with more data. But there are also systematic uncertainties, which are the same for each measurement. The mean of many measurements tends toward the true value plus some constant (and generally unknown) offset known as the systematic uncertainty.

Generally in the early stages, experiments are statistics limited, and taking more data helps reduce the overall uncertainty. But at a later stage, the experiment becomes systematics limited. More data does not help unless the apparatus is improved. Often, in complicated experiments, increased statistics also gives the experimenter a better handle on the systematic uncertainties, allowing for improvements to both the statistical and systematic uncertainties.