

#### Confidence intervals in practice

The recipe to find the interval [a, b] boils down to solving

$$\alpha = \int_{u_{\alpha}(\theta)}^{\infty} g(\hat{\theta}; \theta) \, d\hat{\theta} = \int_{\hat{\theta}_{obs}}^{\infty} g(\hat{\theta}; a) \, d\hat{\theta} \,,$$
  
$$\beta = \int_{-\infty}^{v_{\beta}(\theta)} g(\hat{\theta}; \theta) \, d\hat{\theta} = \int_{-\infty}^{\hat{\theta}_{obs}} g(\hat{\theta}; b) \, d\hat{\theta} \,.$$



→ *a* is hypothetical value of  $\theta$  such that  $P(\hat{\theta} > \hat{\theta}_{obs}) = \alpha$ . → *b* is hypothetical value of  $\theta$  such that  $P(\hat{\theta} < \hat{\theta}_{obs}) = \beta$ .

#### Meaning of a confidence interval

N.B. the interval is random, the true  $\theta$  is an unknown constant. Often report interval [a, b] as  $\hat{\theta}_{-c}^{+d}$ , i.e.  $c = \hat{\theta} - a, d = b - \hat{\theta}$ . So what does  $\hat{\theta} = 80.25^{+0.31}_{-0.25}$  mean? It does not mean:  $P(80.00 < \theta < 80.56) = 1 - \alpha - \beta$ , but rather: repeat the experiment many times with same sample size, construct interval according to same prescription each time, in  $1 - \alpha - \beta$  of experiments, interval will cover  $\theta$ .





Coverage: suppose  $\mu$ \* the true value

 $P(x_1(\mu^*) < x_0 < x_2(\mu^*)) = \alpha$ 





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## "Logic" of an EPP experiment - VI

- End of the selection: CANDIDATES sample  $N_{cand}$
- Which relation is there between  $N_{cand}$  and  $N_X$ ?
  - *Efficiency*: not all searched final states are selected and go to the candidates sample.(Trigger efficiencies are particularly delicate to treat.) Efficiency includes also the **acceptance**.
  - **Background**: few other final states are faking good ones and go in the candidates sample.

$$\varepsilon N_X = N_{cand} - N_b$$

- where:
  - $\varepsilon = \text{efficiency} \ (0 < \varepsilon < 1); \ \varepsilon = A \times \varepsilon_d$
  - $N_b$  = number of background events
- Estimate  $\varepsilon$  and  $N_b$  is a crucial work for the experimentalist and can be done either using simulation (this is tipically done before the experiment and updated later) or using data themselves.



#### **Binomial distribution**

Consider *N* independent experiments (Bernoulli trials): outcome of each is 'success' or 'failure', probability of success on any given trial is *p*.

Define discrete r.v. n = number of successes ( $0 \le n \le N$ ).

Probability of a specific outcome (in order), e.g. 'ssfsf' is  $pp(1-p)p(1-p) = p^n(1-p)^{N-n}$ 

But order not important; there are

 $\frac{N!}{n!(N-n)!}$ 

ways (permutations) to get *n* successes in *N* trials, total probability for *n* is sum of probabilities for each permutation.

#### Binomial distribution (2)

The binomial distribution is therefore

$$f(n; N, p) = \frac{N!}{n!(N-n)!}p^n(1-p)^{N-n}$$
random parameters
variable

For the expectation value and variance we find:

$$E[n] = \sum_{n=0}^{N} nf(n; N, p) = Np$$
$$V[n] = E[n^{2}] - (E[n])^{2} = Np(1-p)$$



#### Binomial distribution (3)

Binomial distribution for several values of the parameters:



Example: observe *N* decays of  $W^{\pm}$ , the number *n* of which are  $W \rightarrow \mu v$  is a binomial r.v., *p* = branching ratio.

Methods in Experimental Particle Physics

#### Multinomial distribution

Like binomial but now m outcomes instead of two, probabilities are

$$\vec{p} = (p_1, \dots, p_m)$$
, with  $\sum_{i=1}^m p_i = 1$ .

For N trials we want the probability to obtain:

 $n_1$  of outcome 1,  $n_2$  of outcome 2,  $\vdots$  $n_m$  of outcome *m*.

This is the multinomial distribution for  $\vec{n} = (n_1, \ldots, n_m)$ 

$$f(\vec{n}; N, \vec{p}) = \frac{N!}{n_1! n_2! \cdots n_m!} p_1^{n_1} p_2^{n_2} \cdots p_m^{n_m}$$



#### Multinomial distribution (2)

Now consider outcome *i* as 'success', all others as 'failure'.

 $\rightarrow$  all  $n_i$  individually binomial with parameters  $N, p_i$ 

$$E[n_i] = Np_i, \quad V[n_i] = Np_i(1-p_i) \quad \text{for all } i$$

One can also find the covariance to be

$$V_{ij} = Np_i(\delta_{ij} - p_j)$$

Example:  $\vec{n} = (n_1, \dots, n_m)$  represents a histogram with *m* bins, *N* total entries, all entries independent.



#### Poisson distribution

Consider binomial n in the limit

$$N \to \infty, \qquad p \to 0, \qquad E[n] = Np \to \nu.$$

 $\rightarrow$  *n* follows the Poisson distribution:

$$f(n;\nu) = \frac{\nu^n}{n!}e^{-\nu} \quad (n \ge 0)$$

$$E[n] = \nu, \quad V[n] = \nu.$$

Example: number of scattering events *n* with cross section  $\sigma$  found for a fixed integrated luminosity, with  $\nu = \sigma \int L dt$ .









The binomial distribution with parameters n and p is

the discrete probability distribution of the number of successes in a sequence of n independent experiments. (Wikipedia)

$$P(k:n,p) = \binom{n}{k} p^{k} (1-p)^{n-k}$$

$$[f X \sim B(n, p)]$$

$$E[X] = np$$





$$\mathbf{P}(k:n,p) = \begin{pmatrix} n \\ k \end{pmatrix} p^k (1-p)^{n-k}$$

The Poisson distribution with parameter  $\lambda = np$  can be used as an approximation to B(n, p) of the binomial distribution if n is sufficiently large and p is sufficiently small.

$$P(k:n,p) \xrightarrow{n \to \infty, np = \lambda} Poiss(k;\lambda) = \frac{\lambda^k e^{-k}}{k!}$$
  
If  $X \sim Poiss(k;\lambda)$   
 $E[X] = Var[X] = \lambda$ 



# **From Binomial to Poisson to Gaussian** $P(k:n,p) = \binom{n}{k} p^{k} (1-p)^{n-k}$

$$P(k:n,p) \xrightarrow{n \to \infty, np = \lambda} Poiss(k;\lambda) = \frac{\lambda^k e^{-k}}{k!}$$

$$\langle k \rangle = \lambda, \ \sigma_k = \sqrt{\lambda}$$

$$k \to \infty \Longrightarrow x = k$$

Using Stirling Formula

prob(x)=G(x, 
$$\sigma = \sqrt{\lambda}$$
) =  $\frac{1}{\sqrt{2\pi\sigma}}e^{-(x-\lambda)^2/2\sigma^2}$ 

This is a Gaussian, or Normal distribution with mean and variance of  $\lambda$ 



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Histograms N collisions  $p(Higgs \; event) = \frac{\mathcal{L}\sigma(pp \to H) A\epsilon_{ff}}{\mathcal{L}\sigma(nn)}$ Prob to see  $n_{H}^{obs}$  in N collisions is  $P(n_{H}^{obs}) = \begin{pmatrix} N \\ n_{H}^{obs} \end{pmatrix} p^{n_{H}^{obs}} (1-p)^{N-n_{H}^{obs}}$  $\ell im_{N \to \infty} P(n_H^{obs}) = Poiss(n_H^{obs}, \lambda) = \frac{e^{-\lambda} \lambda^{n_H}}{n_H^{obs}!}$ 80 100 mass  $\lambda = Np = \mathcal{L}\sigma(pp) \cdot \frac{\mathcal{L}\sigma(pp \to H) A\epsilon_{ff}}{\mathcal{L}\sigma(pp)} = n_H^{exp}$ 



#### Histograms

pdf = histogram with
 infinite data sample,
 zero bin width,
 normalized to unit area.

$$f(x) = \frac{N(x)}{n\Delta x}$$

$$n =$$
 number of entries

 $\Delta x = \text{bin width}$ 





#### Uniform distribution

Consider a continuous r.v. *x* with  $-\infty < x < \infty$ . Uniform pdf is:



N.B. For any r.v. *x* with cumulative distribution F(x), y = F(x) is uniform in [0,1].

Example: for  $\pi^0 \to \gamma \gamma$ ,  $E_{\gamma}$  is uniform in  $[E_{\min}, E_{\max}]$ , with  $E_{\min} = \frac{1}{2} E_{\pi} (1 - \beta)$ ,  $E_{\max} = \frac{1}{2} E_{\pi} (1 + \beta)$ 

#### Exponential distribution

The exponential pdf for the continuous r.v. *x* is defined by:



Example: proper decay time *t* of an unstable particle

 $f(t;\tau) = \frac{1}{\tau}e^{-t/\tau}$  ( $\tau$  = mean lifetime)

Lack of memory (unique to exponential):  $f(t - t_0 | t \ge t_0) = f(t)$ 



#### Gaussian distribution

The Gaussian (normal) pdf for a continuous r.v. *x* is defined by:

$$f(x; \mu, \sigma) = \frac{1}{\sqrt{2\pi\sigma}} e^{-(x-\mu)^2/2\sigma^2}$$

$$E[x] = \mu$$

$$K[x] = \mu$$

$$F[x] = \mu$$

$$F[x] = \mu$$

$$F[x] = \mu$$

$$F[x] = \sigma^2$$

$$F[x] =$$

Special case:  $\mu = 0$ ,  $\sigma^2 = 1$  ('standard Gaussian'):

$$\varphi(x) = \frac{1}{\sqrt{2\pi}} e^{-x^2/2} , \quad \Phi(x) = \int_{-\infty}^x \varphi(x') \, dx'$$

If  $y \sim$  Gaussian with  $\mu$ ,  $\sigma^2$ , then  $x = (y - \mu) / \sigma$  follows  $\varphi(x)$ .



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#### Gaussian pdf and the Central Limit Theorem

The Gaussian pdf is so useful because almost any random variable that is a sum of a large number of small contributions follows it. This follows from the Central Limit Theorem:

For *n* independent r.v.s  $x_i$  with finite variances  $\sigma_i^2$ , otherwise arbitrary pdfs, consider the sum

$$y = \sum_{i=1}^{n} x_i$$

In the limit  $n \to \infty$ , y is a Gaussian r.v. with

$$E[y] = \sum_{i=1}^{n} \mu_i$$
  $V[y] = \sum_{i=1}^{n} \sigma_i^2$ 

Measurement errors are often the sum of many contributions, so frequently measured values can be treated as Gaussian r.v.s.



## Quantities to measure

- In order to estimate  $N_X$  we need to measure:
  - N<sub>cand</sub>
  - E
  - N<sub>b</sub>
- We already know that each of these variables have a fluctuation model:
  - $N_{cand}$  is described by a Poisson process
  - $\boldsymbol{\varepsilon}$  is described by a Bernoulli process
  - N<sub>b</sub>



## N<sub>cand</sub>: a Poisson variable

- If events come in a random way (without any time structure) the event count *N* is a Poisson variable.
- $\rightarrow$  if I count *N*, the best estimate of  $\lambda$  is *N* itself and the uncertainty is  $\sqrt{N}$
- If N is large enough (N>20) Poisson  $\rightarrow$  Gaussian.  $\rightarrow N \pm \sqrt{N}$  is a 68% probability interval for N.
- If *N* is small (close to 0) the Gaussian limit is not ok, a specific treatment is required (see later in the course).



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$$P(N,\lambda) = \lambda^{N} e^{-\lambda} / N! \Longrightarrow P(\lambda \mid N) = \lambda^{N} e^{-\lambda} / N!$$
$$E[\lambda] = N + 1$$
$$var[\lambda] = N + 1$$

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## Efficiency: a binomial variable - I

• Bernoulli process: success/failure N proofs, 0 < n < N, p = success probability.  $p == \varepsilon$ 

$$P(n / N, p) = \binom{N}{n} p^n (1 - p)^{N - n}$$
$$E[n] = Np$$
$$var[n] = Np(1 - p)$$

• Inference: given n and N which is the best estimate of p? And its uncertainty ? *(see previous lectures)* 

$$\begin{split} \varepsilon &= \hat{p} = \frac{n+1}{N+2} \\ \sigma \left( \varepsilon \right) = \frac{\sigma \left( n \right)}{N} = \frac{1}{\sqrt{N+2}} \sqrt{\hat{p}(1-\hat{p})} \end{split}$$



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## Efficiency: a binomial variable - II

- How measure it ?
  - From data: Sample of *N* true particles and I measure how many, out of these give rise to a signal in my detector
  - From MC: I generate  $N_{gen}$  "signal" events. If I select  $N_{sel}$  of these events out of  $N_{gen}$ , the efficiency is (assume  $N_{gen}$  and  $N_{sel}$  large numbers):

$$\varepsilon = \frac{N_{sel}}{N_{gen}}$$

$$\sigma(\varepsilon) = \frac{\sigma(N_{sel})}{N_{gen}} = \frac{1}{\sqrt{N_{gen}}} \sqrt{\frac{N_{sel}}{N_{gen}}} \left(1 - \frac{N_{sel}}{N_{gen}}\right)$$



## Background N<sub>b</sub>

- Simulation of  $N_{gen}$  "bad final states";  $N_{sel}$  are selected. What about  $N_b$  ?
- We define the "rejection factor"  $R = N_{gen} / N_{sel} > 1$
- We also need a correct normalization in this case: we need to know  $N_{exp}$  = total number of expected "bad final states" in our sample ( $N_{exp}$  related to luminosity and cross-section).

$$N_{b} = N_{sel} \frac{N_{exp}}{N_{gen}} = \frac{N_{exp}}{R}$$

$$\sigma(N_{b}) = \sigma(N_{sel}) \frac{N_{exp}}{N_{gen}} = \sqrt{N_{sel}} \frac{N_{exp}}{N_{gen}} = \frac{N_{exp}}{\sqrt{RN_{gen}}}$$

### **Statistical Errors**

- In alla cases there is an unreducible error on  $N_X$  given by limited statistics. It is a random error, coming from the procedure of "sampling" that is intrinsic in our experiments.
- In all cases increasing the statistics, the error decreases

$$\frac{\sigma(N_{cand})}{N_{cand}} = \frac{1}{\sqrt{N_{cand}}}$$
$$\sigma(\varepsilon) \approx \frac{1}{\sqrt{N_{gen}}}$$
$$\sigma(N_b) \approx \frac{1}{\sqrt{N_{gen}}}$$

