

Suggestion:

Review the first two chapters of Cowan – Statistical Data Analysis

# 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 ( $0 \leq n \leq N$ ).

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  
variable

parameters

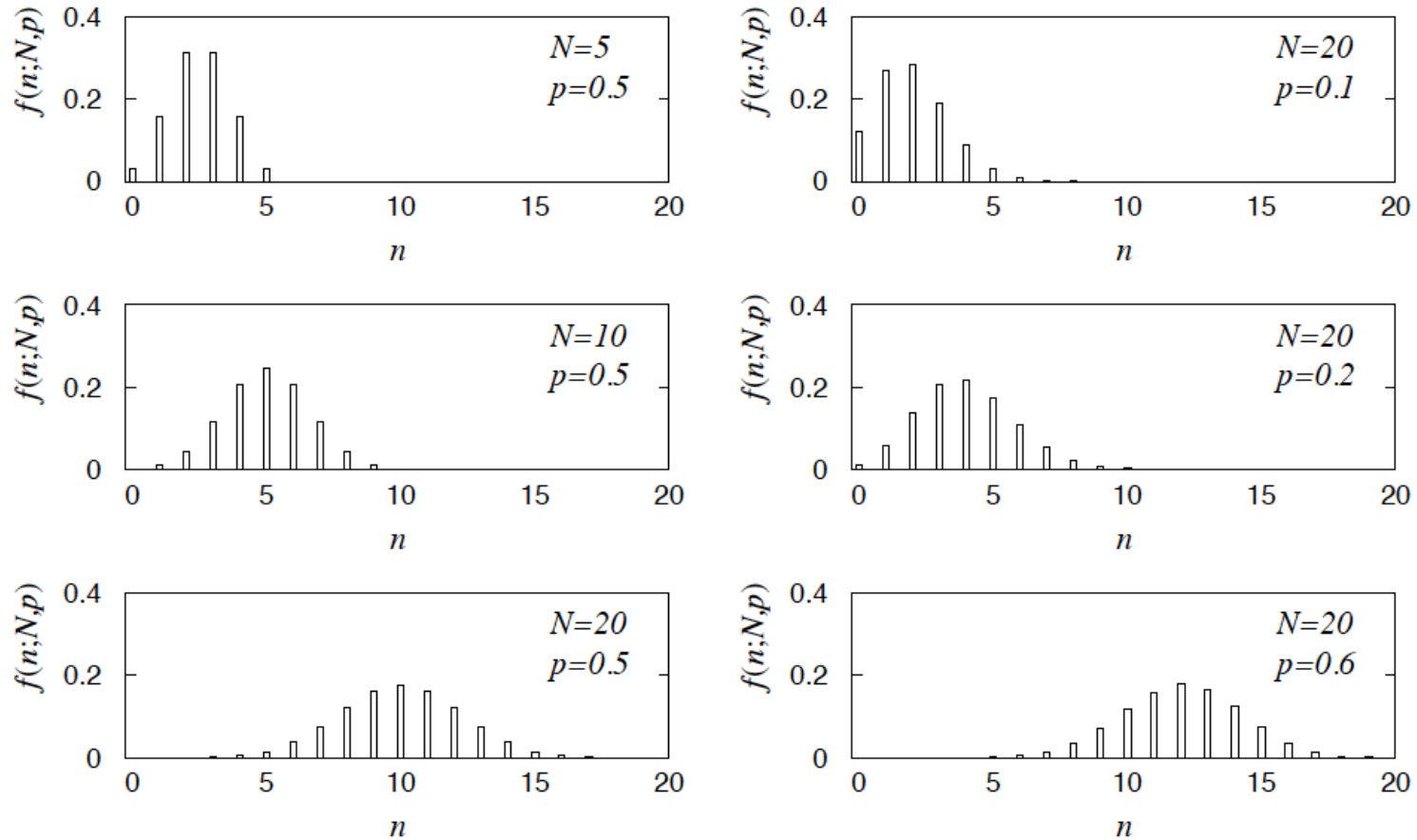
For the expectation value and variance we find:

$$E[n] = \sum_{n=0}^N n f(n; N, p) = Np$$

$$V[n] = E[n^2] - (E[n])^2 = Np(1-p)$$

# Binomial distribution (3)

Binomial distribution for several values of the parameters:



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



# Multinomial distribution

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

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

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  $\vec{n} = (n_1, \dots, n_m)$

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

## Multinomial distribution (2)

Now consider outcome  $i$  as ‘success’, all others as ‘failure’.

→ 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

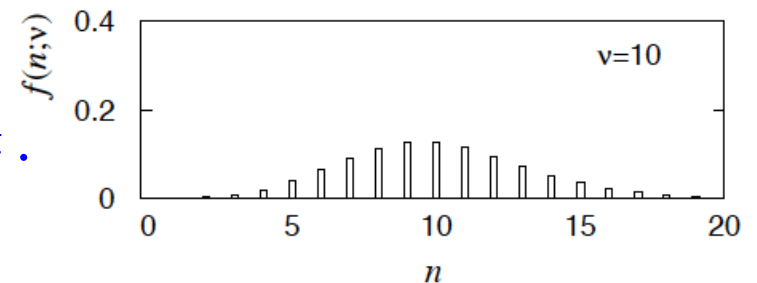
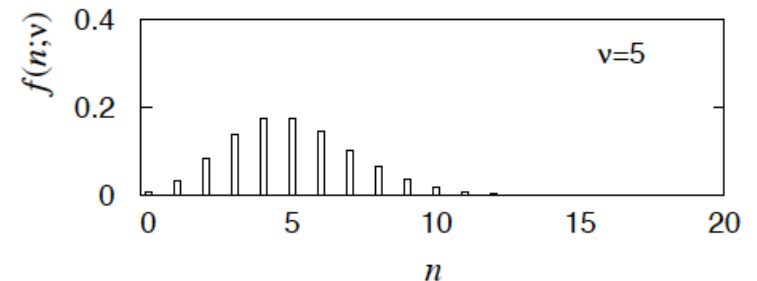
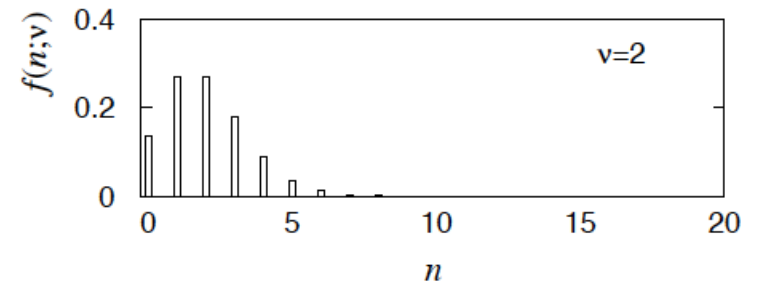
$$N \rightarrow \infty, \quad p \rightarrow 0, \quad E[n] = Np \rightarrow \nu .$$

→  $n$  follows the Poisson distribution:

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

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

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



# From Binomial to Poisson to Gaussian

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

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

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

$$k \rightarrow \infty \Rightarrow x = k$$

Using Stirling Formula

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

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

# Histograms

$N$  collisions

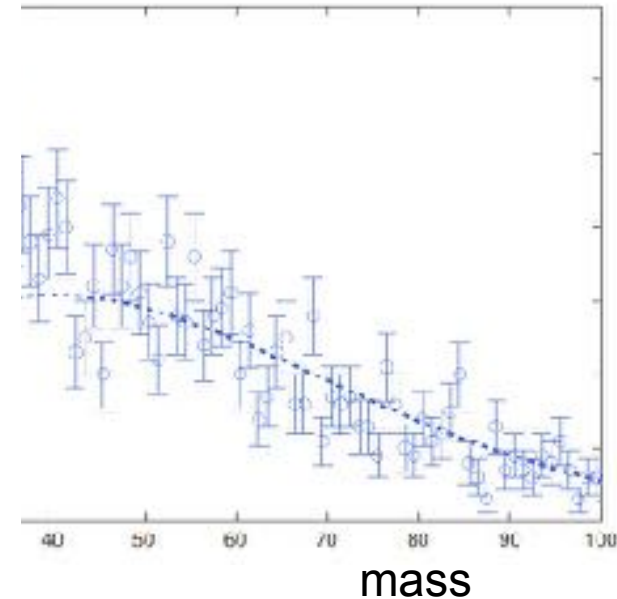
$$p(\text{Higgs event}) = \frac{\mathcal{L}\sigma(pp \rightarrow H) A\epsilon_{ff}}{\mathcal{L}\sigma(pp)}$$

Prob to see  $n_H^{obs}$  in  $N$  collisions is

$$P(n_H^{obs}) = \binom{N}{n_H^{obs}} p^{n_H^{obs}} (1-p)^{N-n_H^{obs}}$$

$$\lim_{N \rightarrow \infty} P(n_H^{obs}) = \text{Poiss}(n_H^{obs}, \lambda) = \frac{e^{-\lambda} \lambda^{n_H^{obs}}}{n_H^{obs}!}$$

$$\lambda = Np = \mathcal{L}\sigma(pp) \cdot \frac{\mathcal{L}\sigma(pp \rightarrow H) A\epsilon_{ff}}{\mathcal{L}\sigma(pp)} = n_H^{exp}$$



# Histograms

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

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

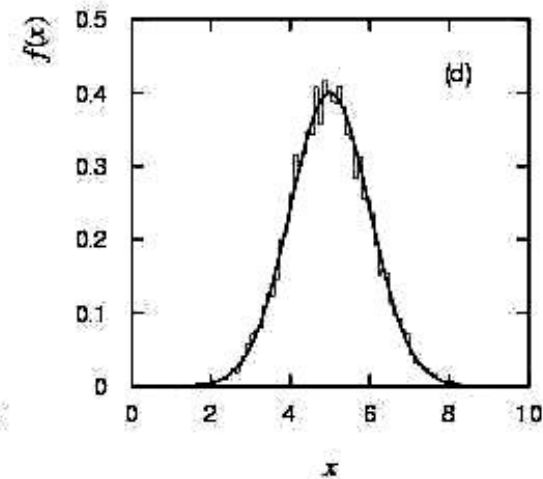
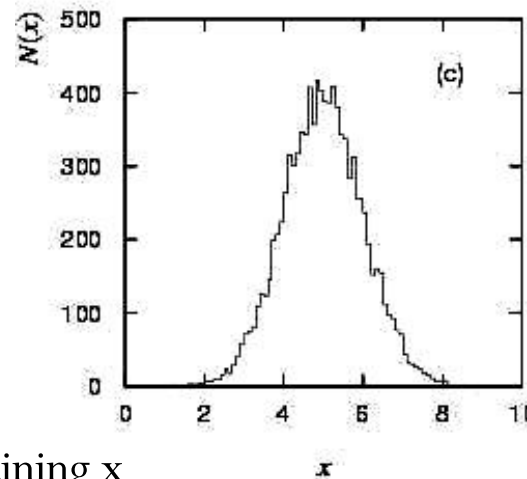
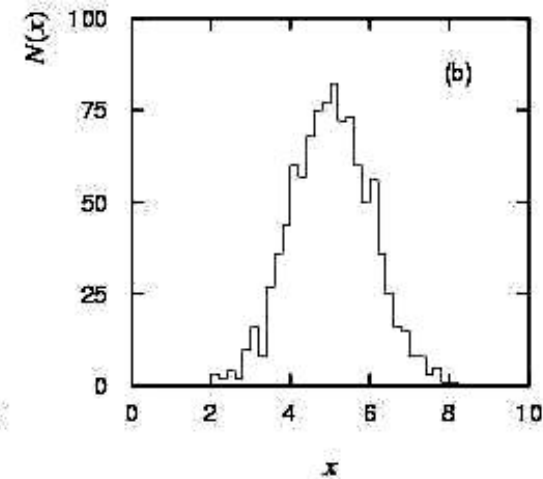
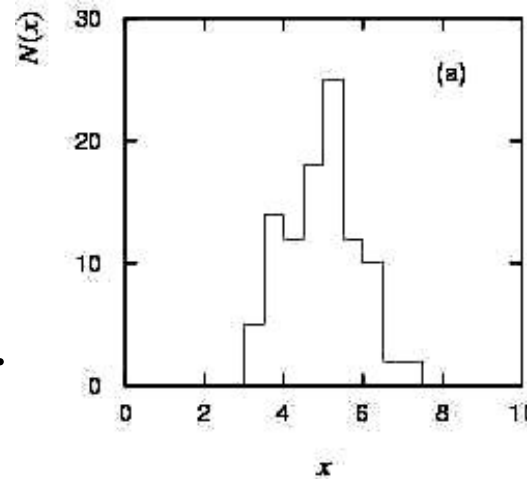
$n$  = number of entries

$\Delta x$  = bin width

$N(x)$  = number of entries in a bin containing  $x$

$f(x)$  = pdf

$$\int_{x_{\min}}^{x_{\max}} f(x) dx = 1$$



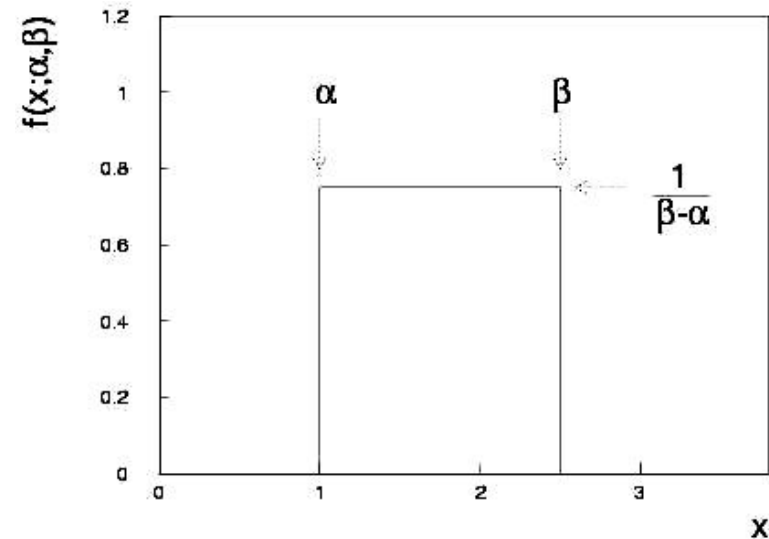
# Uniform distribution

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

$$f(x; \alpha, \beta) = \begin{cases} \frac{1}{\beta - \alpha} & \alpha \leq x \leq \beta \\ 0 & \text{otherwise} \end{cases}$$

$$E[x] = \frac{1}{2}(\alpha + \beta)$$

$$V[x] = \frac{1}{12}(\beta - \alpha)^2$$



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

Example: for  $\pi^0 \rightarrow \gamma\gamma$ ,  $E_\gamma$  is uniform in  $[E_{\min}, E_{\max}]$ , with

$$E_{\min} = \frac{1}{2}E_\pi(1 - \beta), \quad E_{\max} = \frac{1}{2}E_\pi(1 + \beta)$$

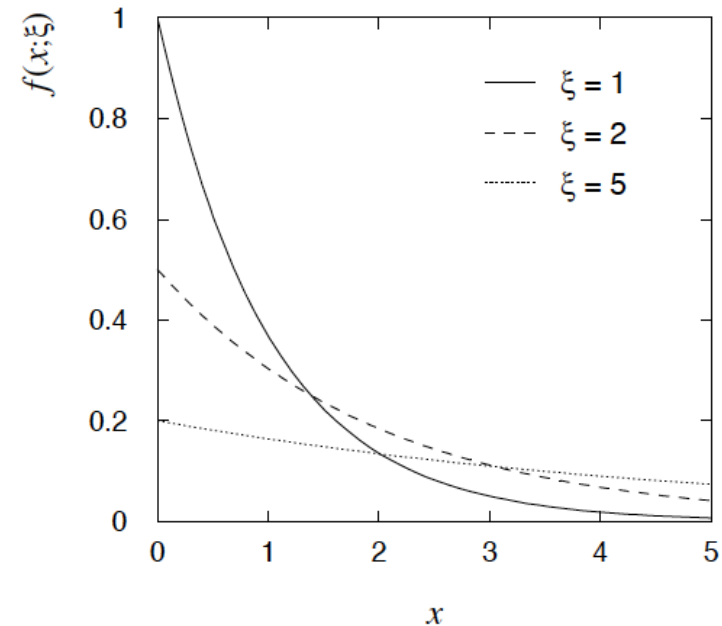
# Exponential distribution

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

$$f(x; \xi) = \begin{cases} \frac{1}{\xi} e^{-x/\xi} & x \geq 0 \\ 0 & \text{otherwise} \end{cases}$$

$$E[x] = \xi$$

$$V[x] = \xi^2$$



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

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

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



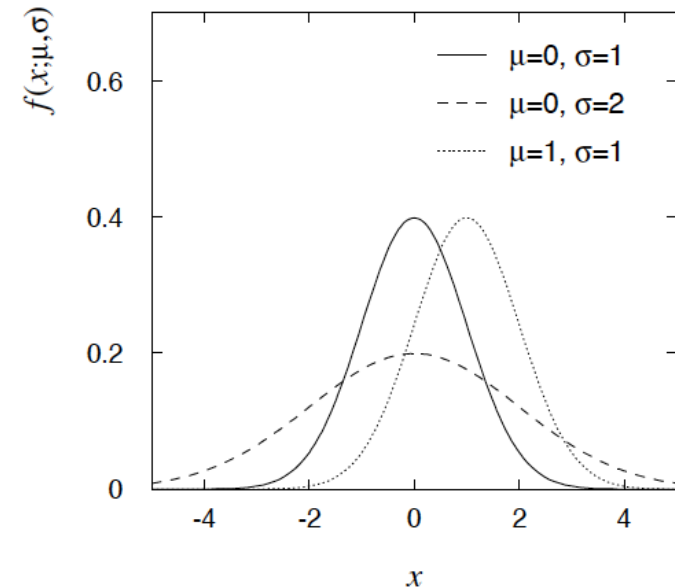
# Gaussian distribution

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

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

$$E[x] = \mu \quad (\text{N.B. often } \mu, \sigma^2 \text{ denote mean, variance of any}$$

$$V[x] = \sigma^2 \quad \text{r.v., not only Gaussian.})$$



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  $\varphi(x)$ .

# 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  $n \rightarrow \infty$ ,  $y$  is a Gaussian r.v. with

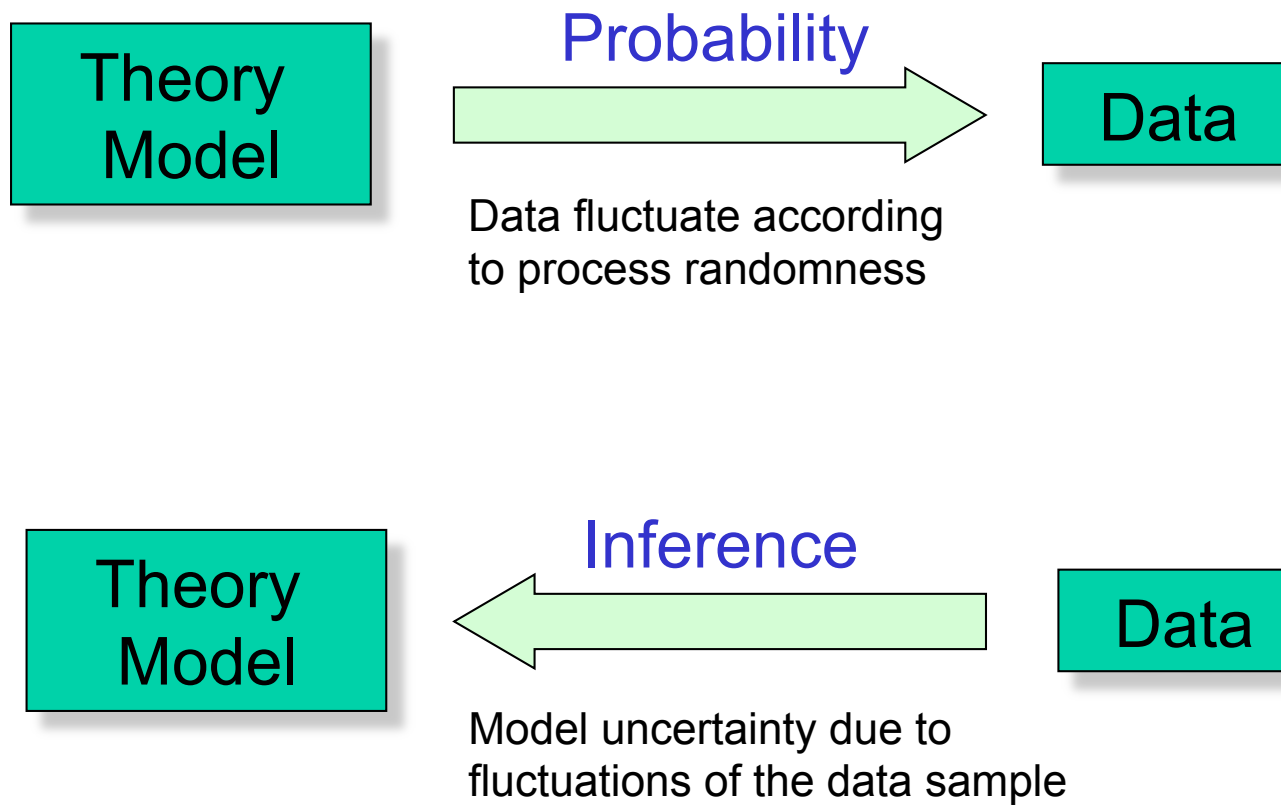
$$E[y] = \sum_{i=1}^n \mu_i \quad V[y] = \sum_{i=1}^n \sigma_i^2$$

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

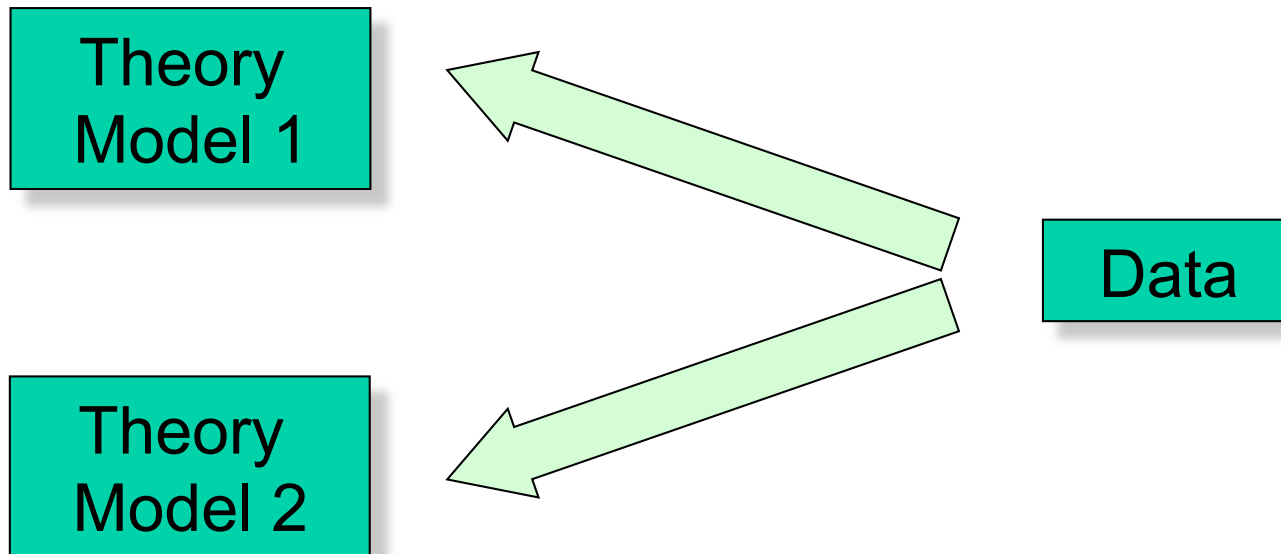
# Meaning of parameter estimate

- We are interested in some physical **unknown parameters**
- Experiments provide **samplings** of some PDF which has among its parameters the physical unknowns we are interested in
- Experiment's results are statistically "related" to the unknown PDF
  - PDF parameters can be **determined** from the sample within some **approximation** or **uncertainty**
- **Knowing** a parameter within some **error** may mean different things:
  - **Frequentist**: a large fraction (68% or 95%, usually) of the experiments will contain, in the limit of large number of experiments, the (fixed) unknown true value within the quoted confidence interval, usually  $[\mu - \sigma, \mu + \sigma]$  ('**coverage**')
  - **Bayesian**: we determine a **degree of belief** that the unknown parameter is contained in a specified interval can be quantified as 68% or 95%
- **We will see that there is still some more degree of arbitrariness in the definition of confidence intervals...**

# Statistical inference



# Hypothesis tests



Which hypothesis is the most consistent with the experimental data?

# Parameter estimators

- An **estimator** is a function of a given sample whose statistical properties are known and related to some PDF parameters
  - “Best fit”
- Simplest example:
  - Assume we have a Gaussian PDF with a *known*  $\sigma$  and an *unknown*  $\mu$
  - A single experiment will provide a measurement  $x$
  - We estimate  $\mu$  as  $\mu^{\text{est}} = x$
  - The distribution of  $\mu^{\text{est}}$  (repeating the experiment many times) is the original Gaussian
  - 68.27%, *on average*, of the experiments will provide an estimate within:  $\mu - \sigma < \mu^{\text{est}} < \mu + \sigma$
- We can determine:  $\mu = \mu^{\text{est}} \pm \sigma$

# Likelihood function

- Given a sample of  $N$  events each with variables  $(x_1, \dots, x_n)$ , the likelihood function expresses the probability density of the sample, as a function of the unknown parameters:

$$L = \prod_{i=1}^N f(x_1^i, \dots, x_n^i; \theta_1, \dots, \theta_m)$$

- Sometimes the used notation for parameters is the same as for conditional probability:

$$f(x_1, \dots, x_n | \theta_1, \dots, \theta_m)$$

- If the size  $N$  of the sample is also a random variable, the extended likelihood function is also used:

$$L = p(N; \theta_1, \dots, \theta_m) \prod_{i=1}^N f(x_1^i, \dots, x_n^i; \theta_1, \dots, \theta_m)$$

- Where  $p$  is most of the times a Poisson distribution whose average is a function of the unknown parameters

- In many cases it is convenient to use  $-\ln L$  or  $-2 \ln L$ :  $\prod_i \rightarrow \sum_i$

# Maximum likelihood estimates

- ML is the widest used parameter estimator
- The “best fit” parameters are the set that maximizes the likelihood function
  - “Very good” statistical properties
- The maximization can be performed analytically, for the simplest cases, and numerically for most of the cases
- **Minuit** is historically the most used minimization engine in High Energy Physics
  - F. James, 1970’s; rewritten in C++ recently



## CL & CI

measurement  $\hat{\mu} = 1.1 \pm 0.3$

$$L(\mu) = G(\mu; \hat{\mu}, \sigma_{\hat{\mu}})$$

$\Rightarrow$  CI of  $\mu = [0.8, 1.4]$  at 68% CL

- A confidence interval (CI) is a particular kind of interval estimate of a population parameter.
- Instead of estimating the parameter by a single value, an interval likely to include the parameter is given.
- How likely the interval is to contain the parameter is determined by the confidence level
- Increasing the desired confidence level will widen the confidence interval.

# Confidence Interval & Coverage

- Say you have a measurement  $\mu_{meas}$  of  $\mu$  with  $\mu_{true}$  being the unknown true value of  $\mu$
- Assume you know the probability distribution function  $p(\mu_{meas}|\mu)$
- based on your statistical method you deduce that there is a 95% Confidence interval  $[\mu_1, \mu_2]$ .  
(it is 95% likely that the  $\mu_{true}$  is in the quoted interval)

The correct statement:

- In an ensemble of experiments 95% of the obtained confidence intervals will contain the true value of  $\mu$ .

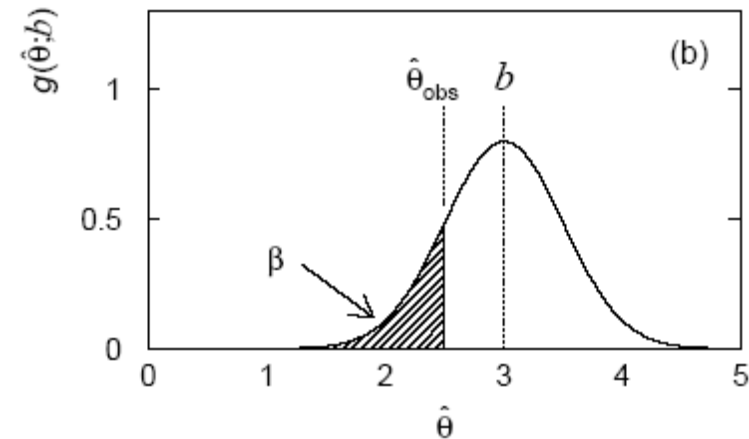
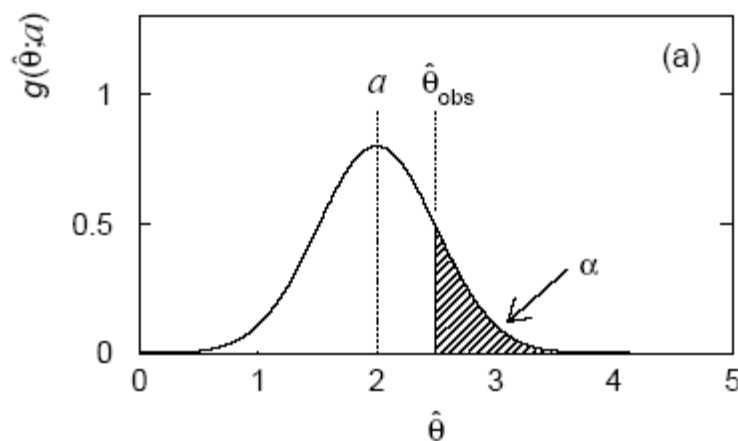


# Confidence intervals in practice

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

$$\alpha = \int_{u_\alpha(\theta)}^{\infty} g(\hat{\theta}; \theta) d\hat{\theta} = \int_{\hat{\theta}_{\text{obs}}}^{\infty} g(\hat{\theta}; a) d\hat{\theta},$$

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



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

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

## 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  $\hat{\theta} = 80.25_{-0.25}^{+0.31}$  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  $\theta$ .

# Confidence Interval & Coverage

- You claim,  $CI_{\mu}=[\mu_1, \mu_2]$  at the 95% CL  
i.e. In an ensemble of experiments CL (95%) of the obtained confidence intervals will contain the true value of  $\mu$ .
- If your statement is accurate, you have full coverage
- If the true CL is  $>95\%$ , your interval has an over coverage
- If the true CL is  $<95\%$ , your interval has an undercoverage

# How to deduce a CI

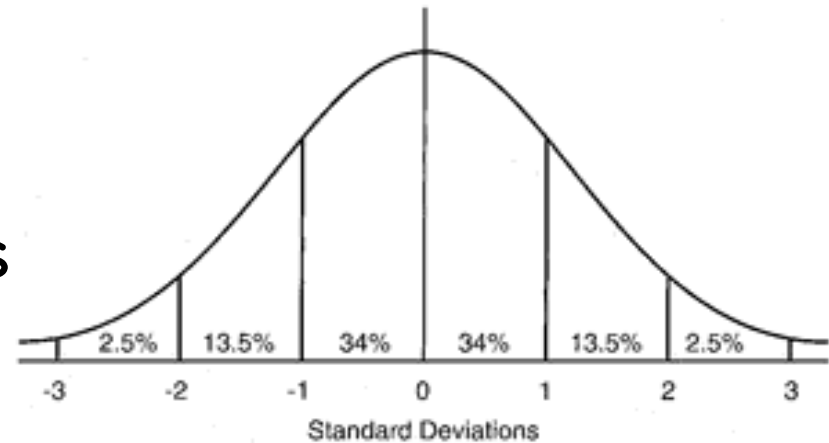
- One can show that if the data is distributed normal around the average i.e.  $P(\text{data}|\mu) = \text{normal}$

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

- then one can construct a 68% CI around the estimator of  $\mu$  to be

$$\hat{X} \pm \sigma \quad \text{i.e. } x_{\text{true}} \in [\hat{x} - \sigma_{\hat{x}}, \hat{x} + \sigma_{\hat{x}}] @ 68\% \text{ CL}$$

- However, not all distributions are normal, many distributions are even unknown and coverage might be a real issue



Side Note:

A CI is an interval in the true parameters phase-space

- One can guarantee a coverage with the Neyman Construction (1937)

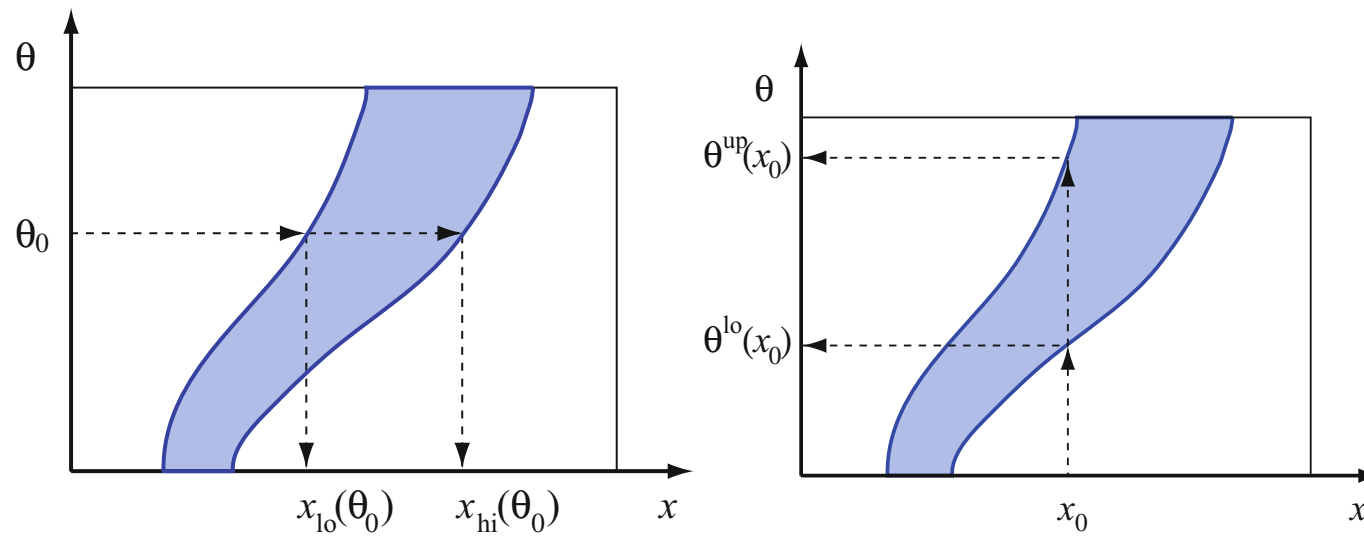
Neyman, J. (1937) "Outline of a Theory of Statistical Estimation Based on the Classical Theory of Probability" Philosophical Transactions of the Royal Society of London A, 236, 333-380.

# The Frequentist Game a 'la Neyman

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Or

How to ensure a Coverage with  
Neyman construction



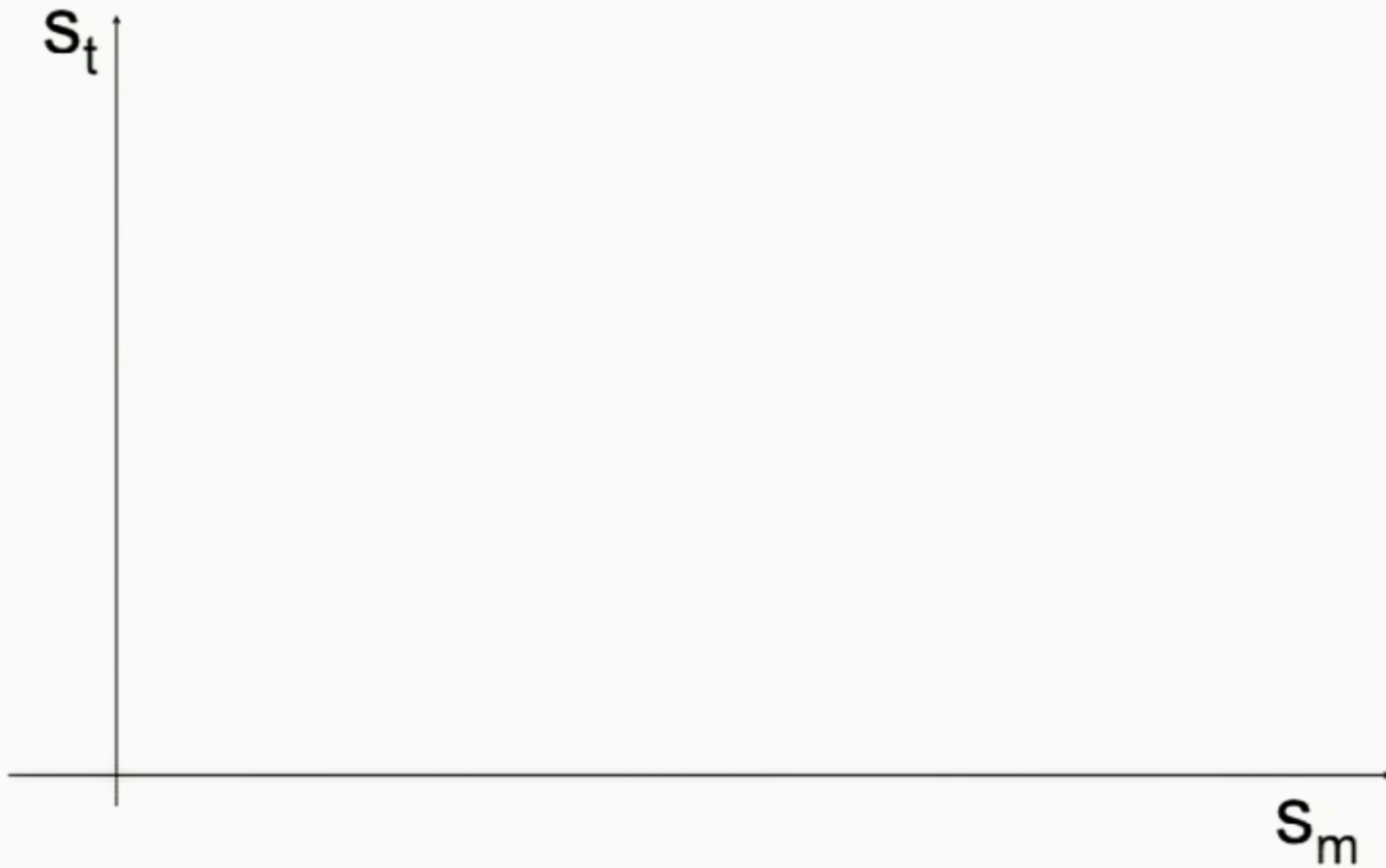
**Fig. 7.1** Graphical illustration of Neyman belt construction (*left*) and inversion (*right*)

$$1 - \alpha = \int_{x^{\text{lo}}(\theta_0)}^{x^{\text{up}}(\theta_0)} f(x | \theta_0) dx$$



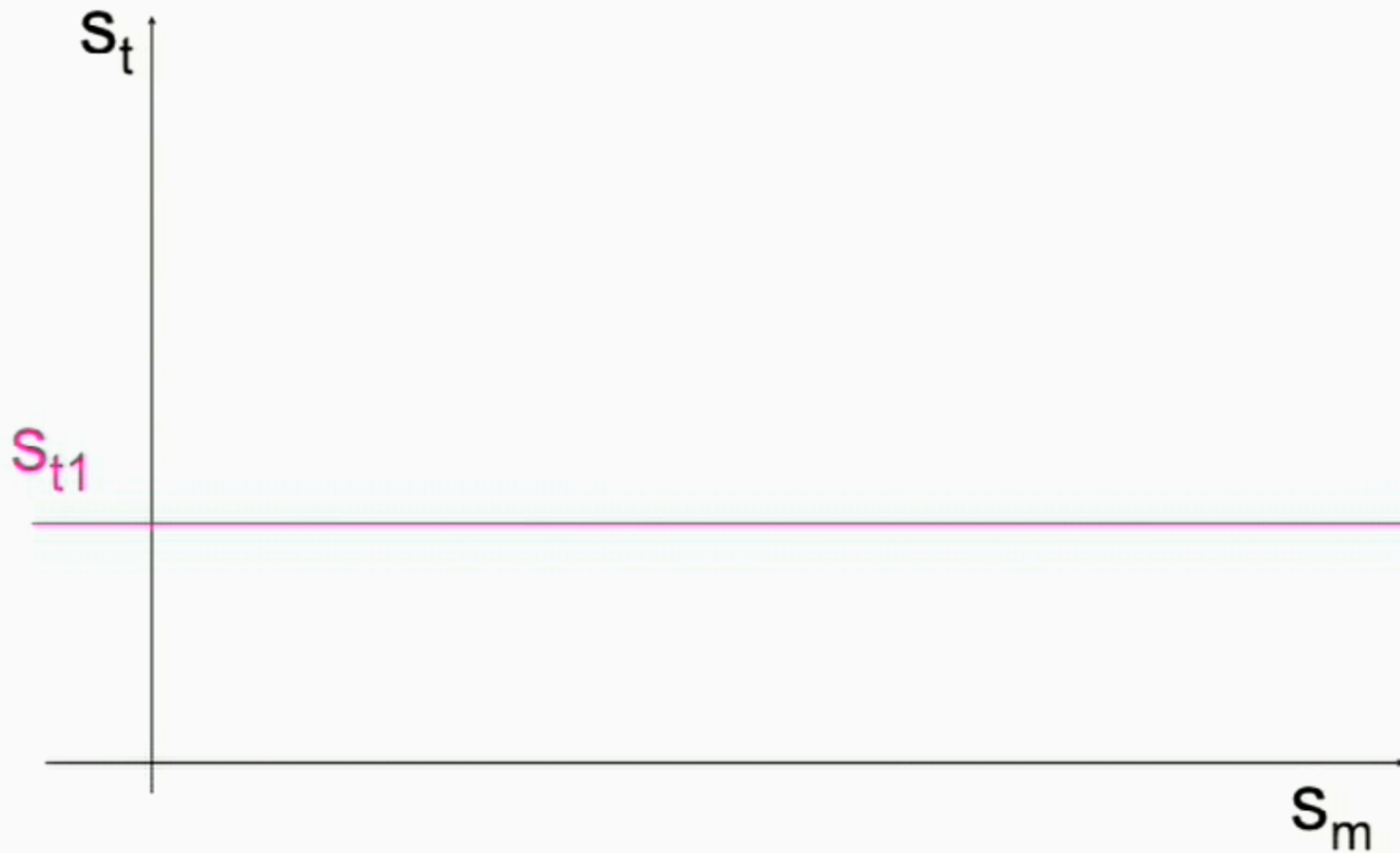
# Neyman Construction

$\text{Prob}(s_m | s_t)$  is known



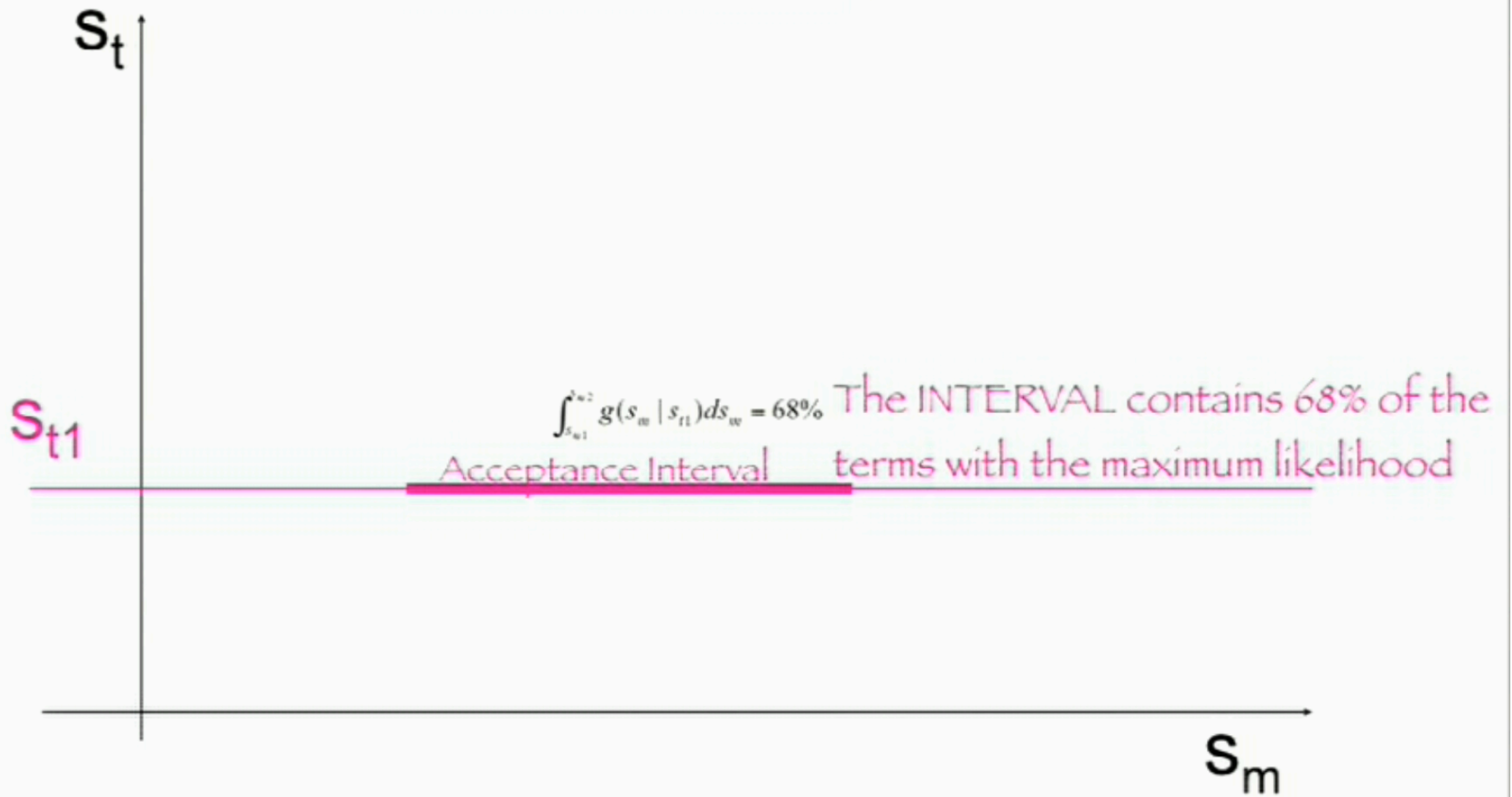
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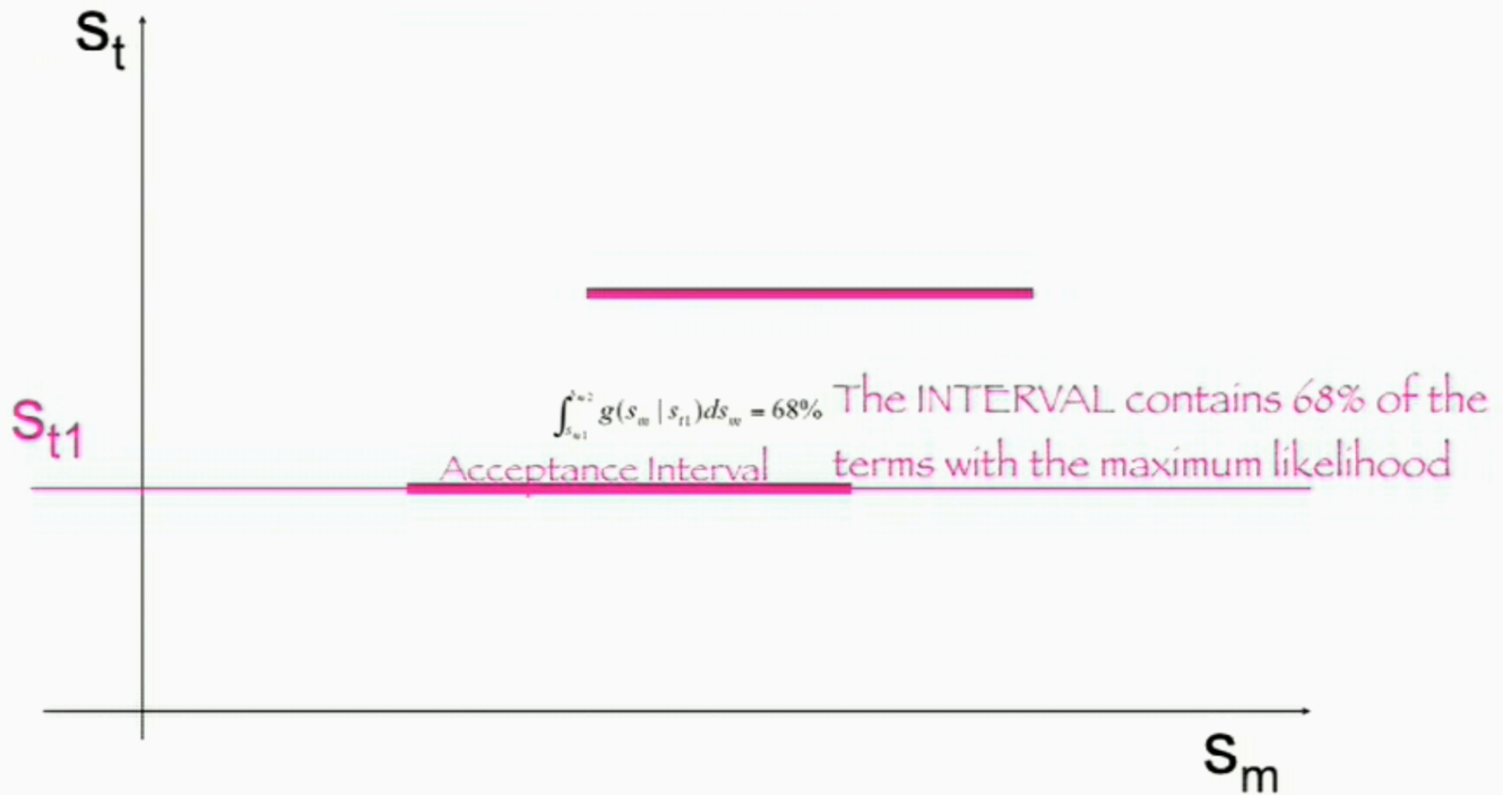
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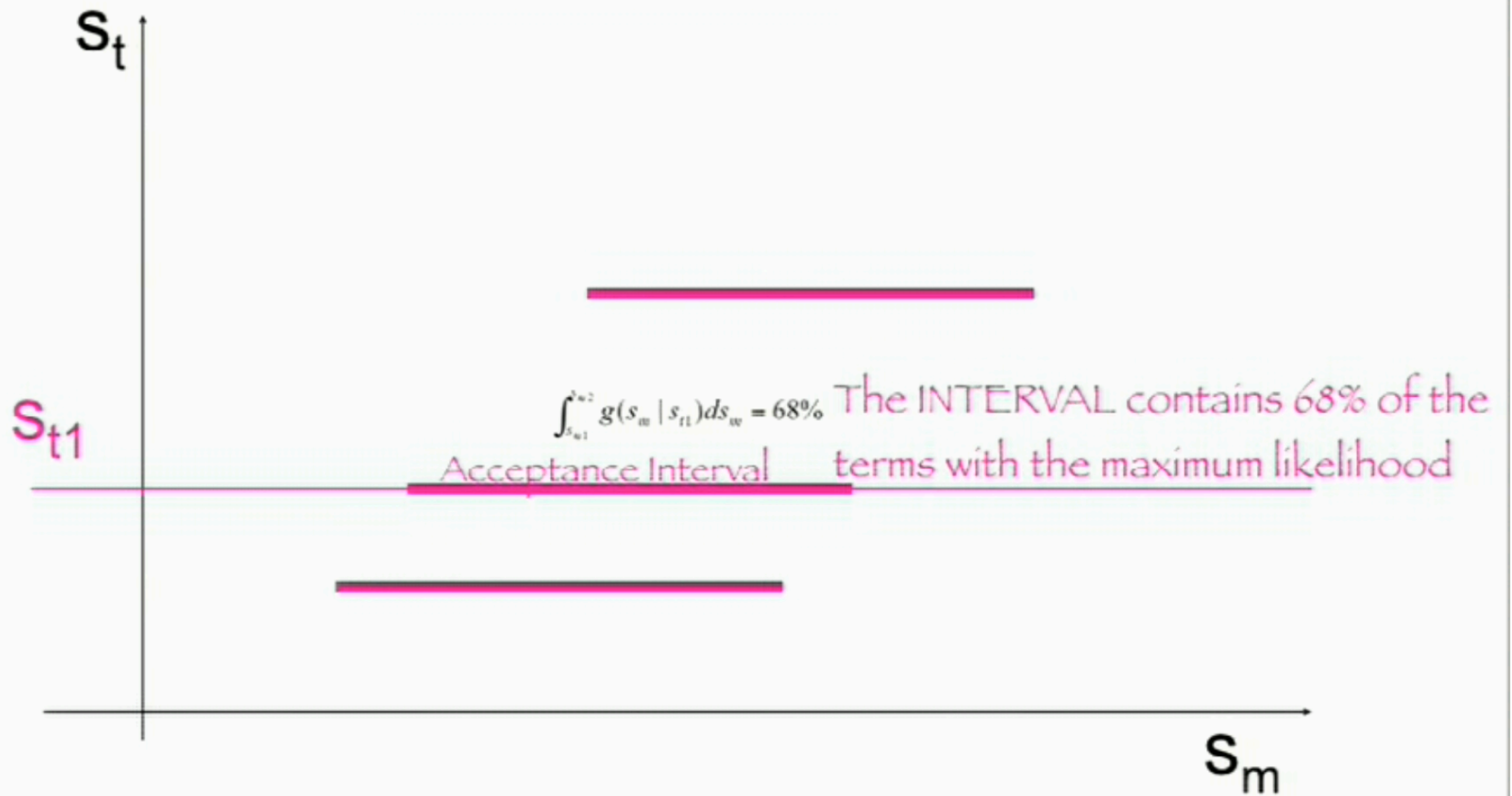
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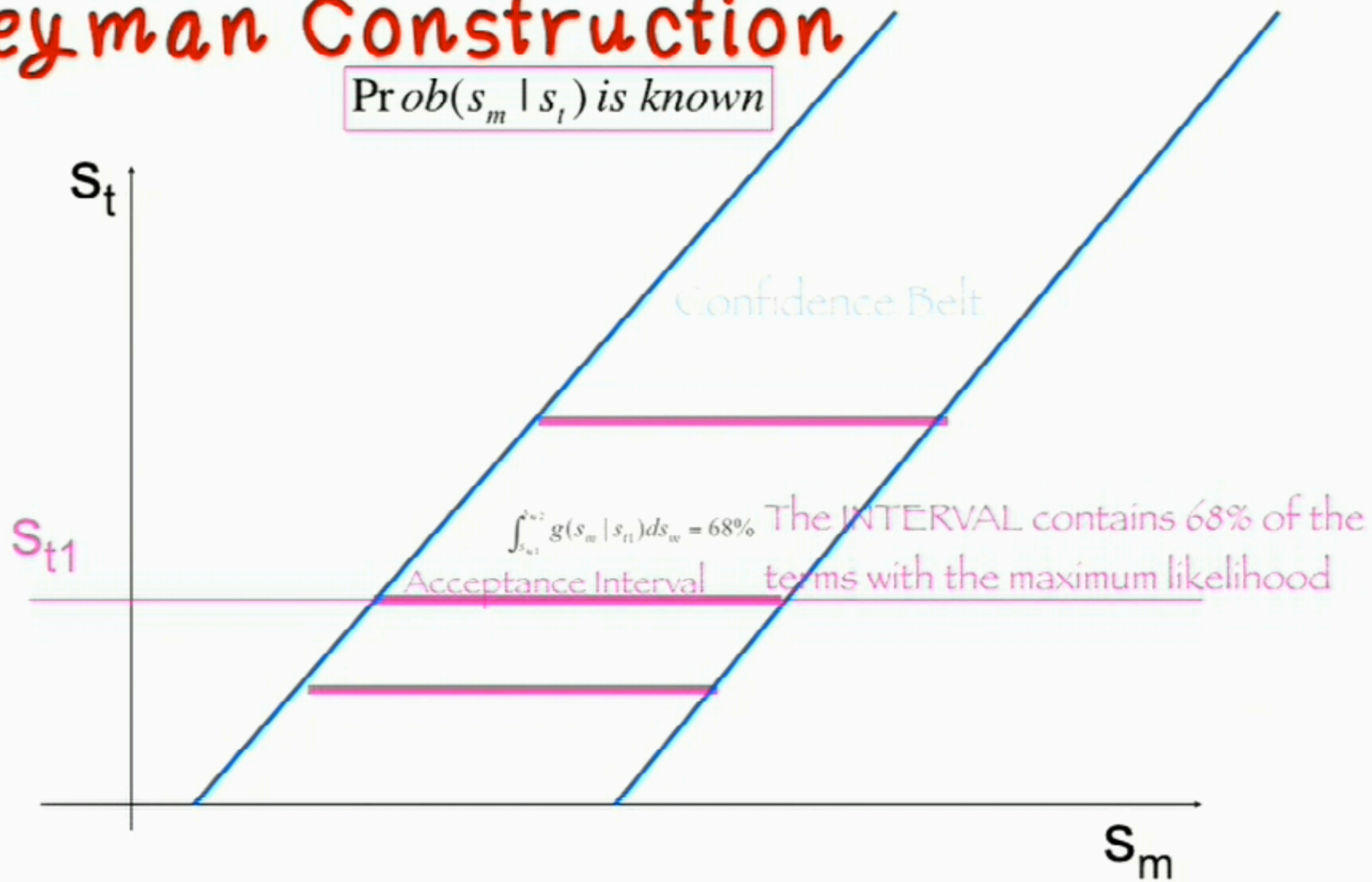
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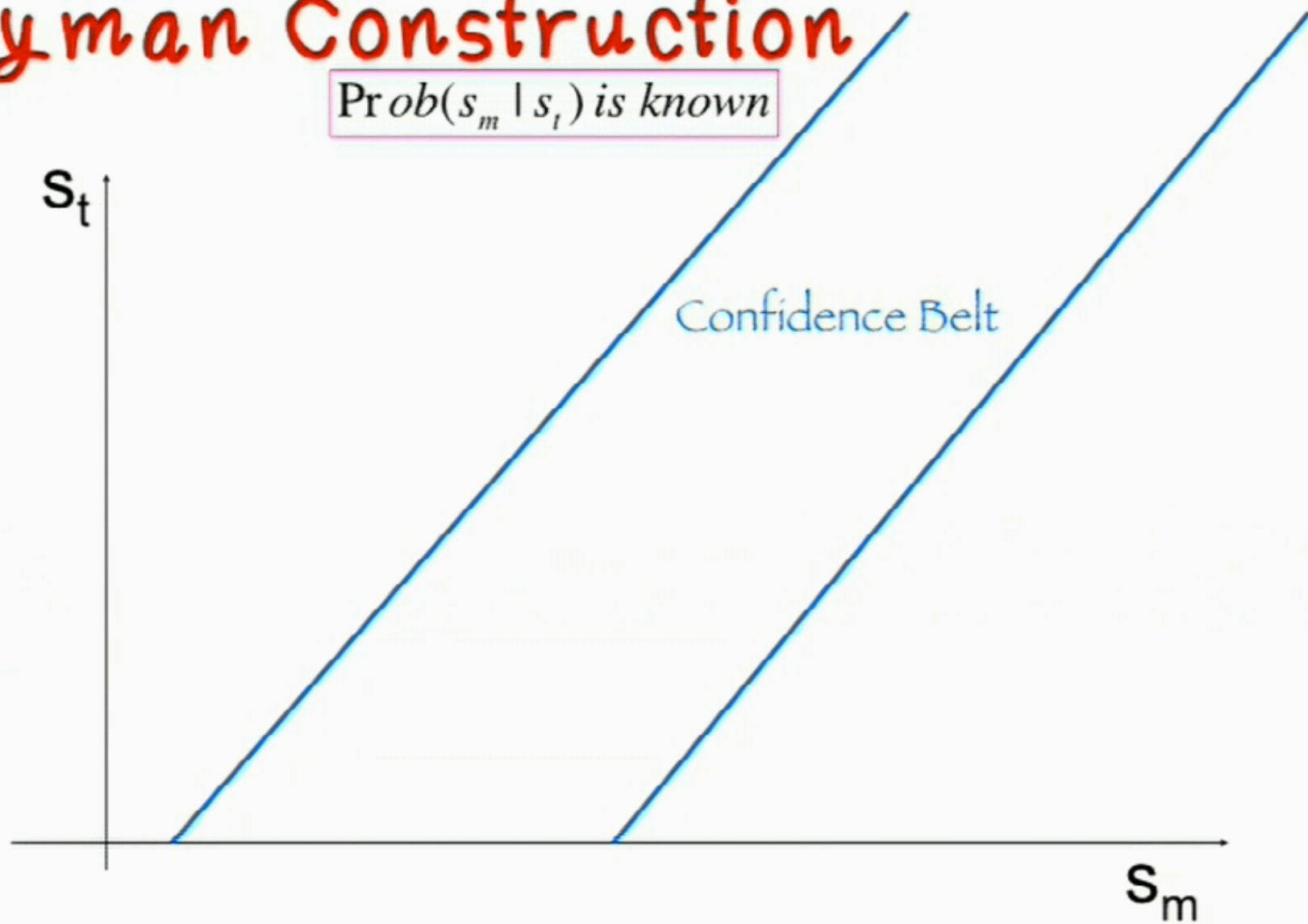
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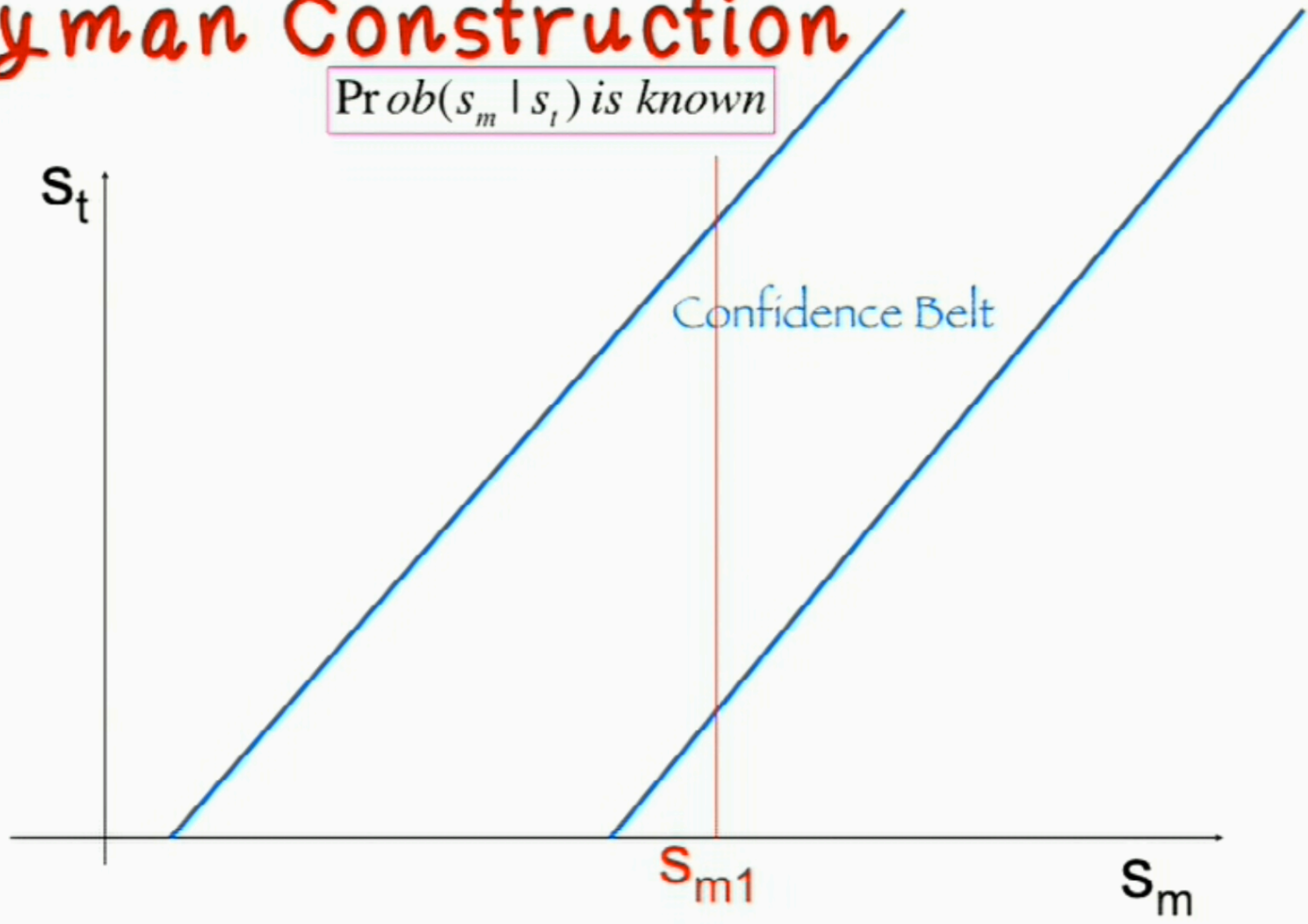
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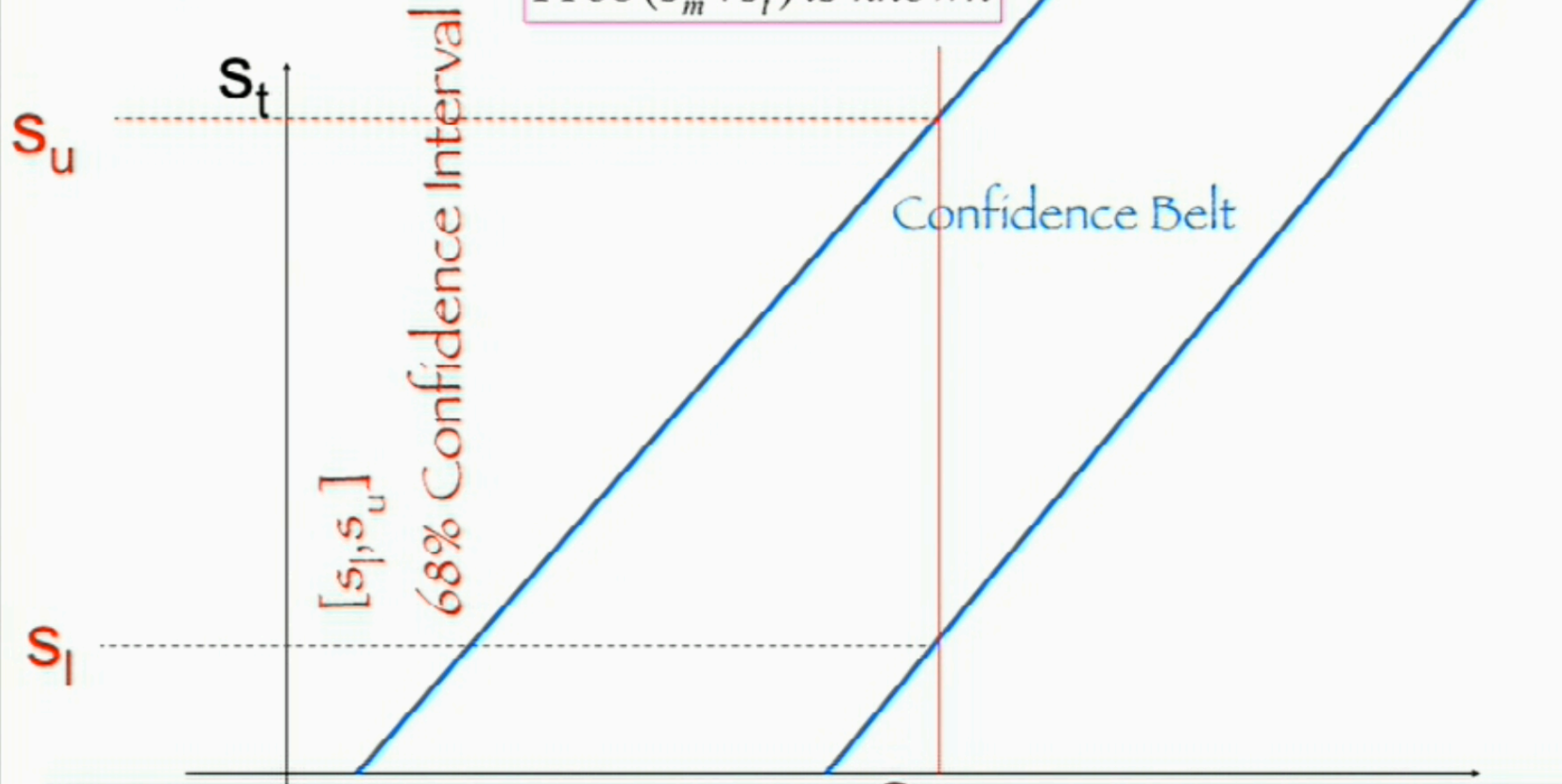
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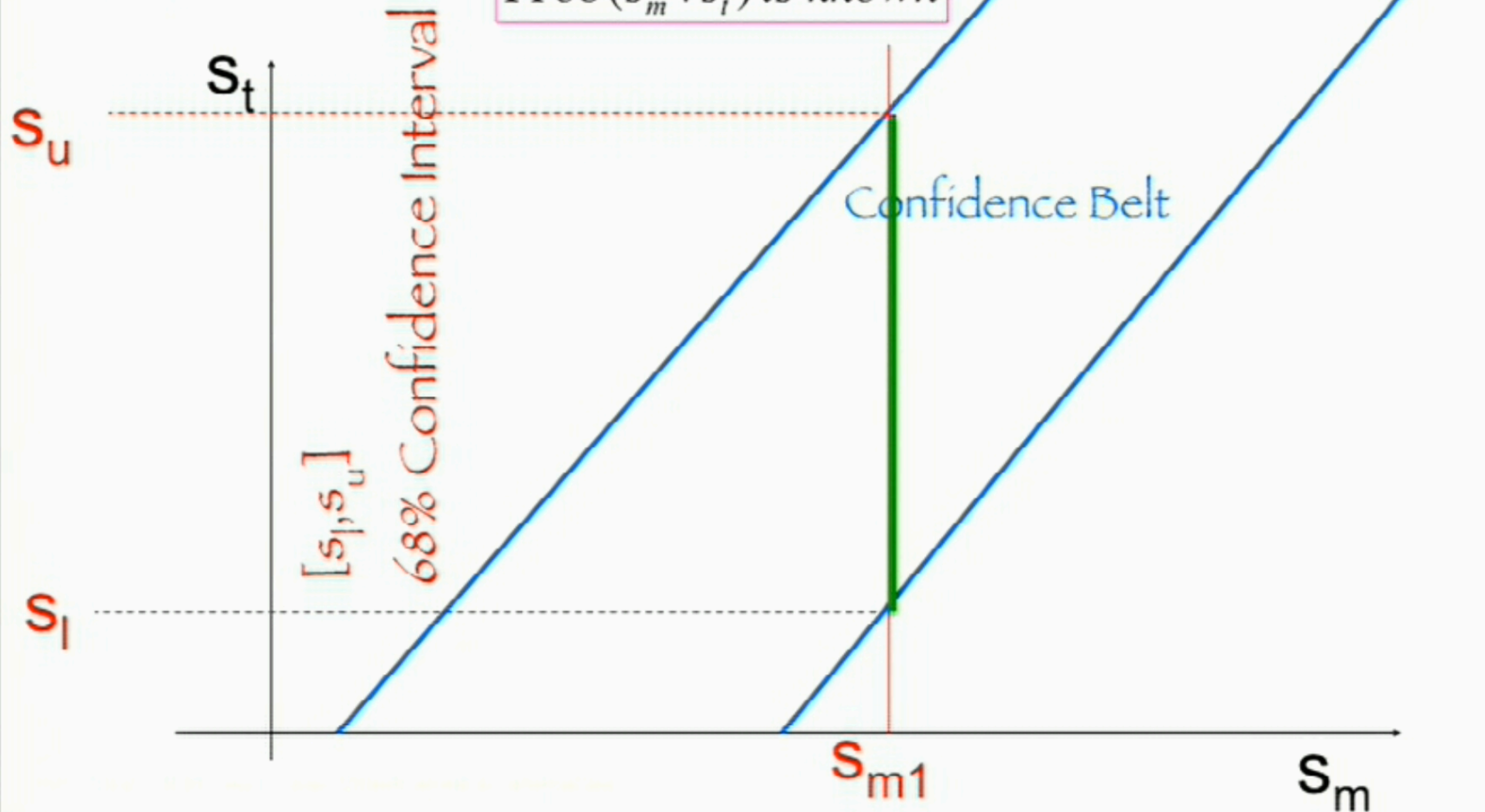


$[s_l, s_u]$  68% Confidence Interval

In 68% of the experiments the derived C.I. contains the unknown true value of  $s$

# Neyman Construction

$Pr ob(s_m | s_t)$  is known



- With Neyman Construction we guarantee a coverage via construction, i.e. for any value of the unknown true  $s$ , the Construction Confidence Interval will cover  $s$  with the correct rate.

# Neyman Construction

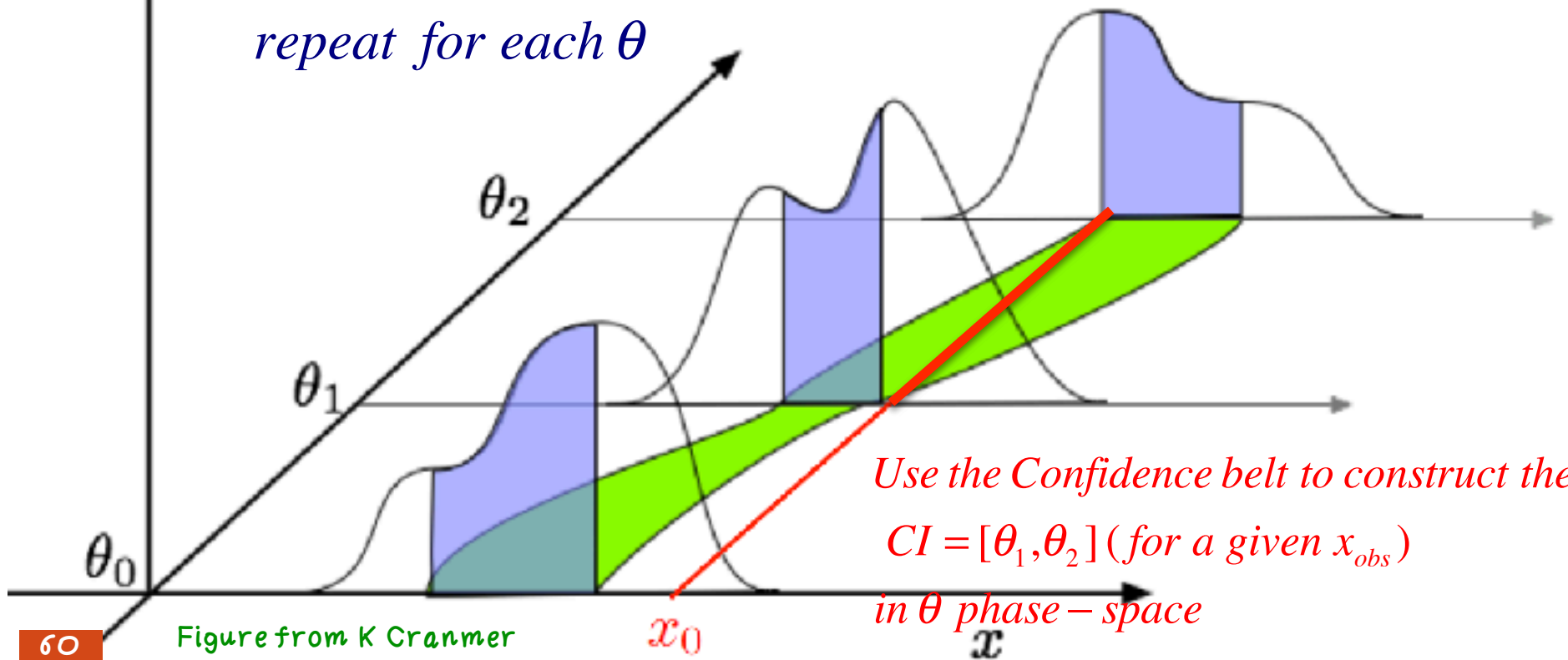
$\theta \equiv s_{true}$     $x \equiv s_{measured}$    pdf  $f(x|\theta)$  is known

for each prospective  $\theta$  generate  $x$

$f(x|\theta)$  construct an interval in DATA phase – space

$$\text{Interval} = \int_{x_l}^{x_h} f(x|\theta) dx = 68\%$$

repeat for each  $\theta$



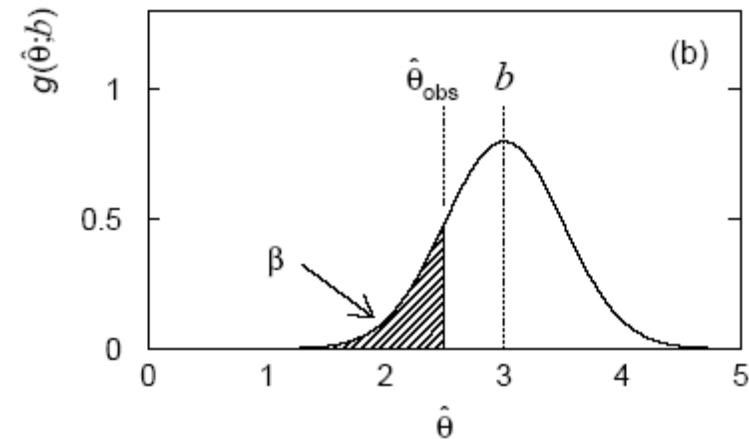
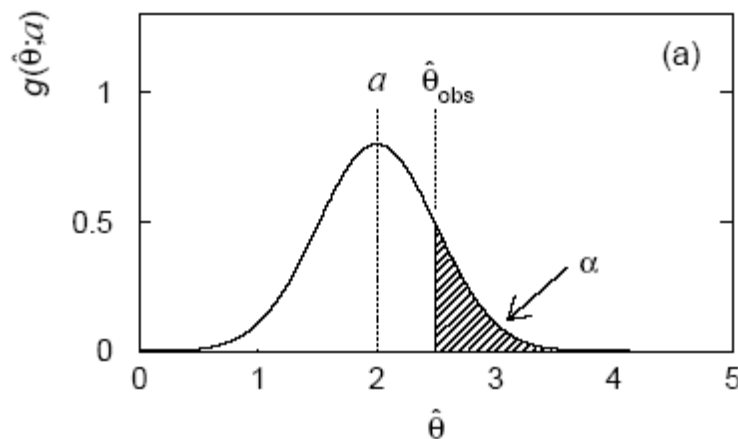
Use the Confidence belt to construct the  
 $CI = [\theta_1, \theta_2]$  (for a given  $x_{obs}$ )  
 in  $\theta$  phase – space

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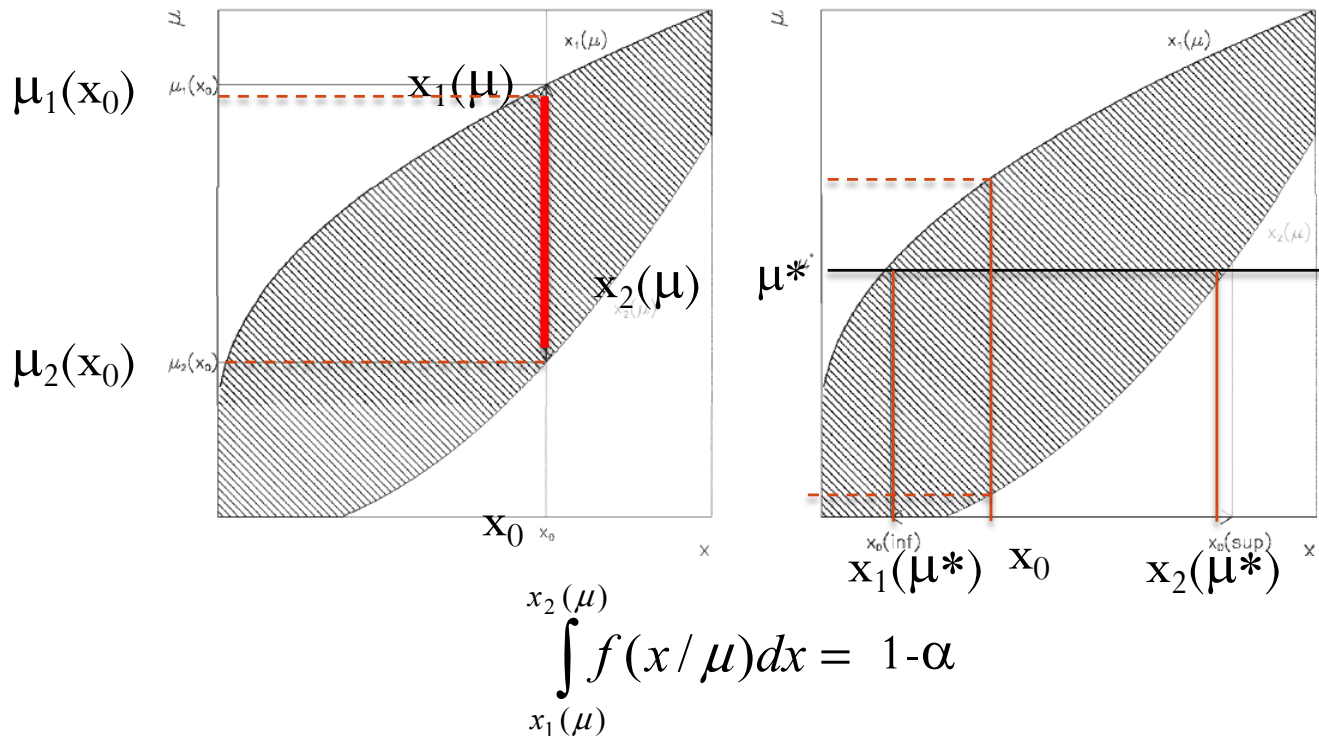
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repeat the experiment many times with same sample size,  
construct interval according to same prescription each time,  
in  $1 - \alpha - \beta$  of experiments, interval will cover  $\theta$ .



## Neyman's construction



By construction the probability to measure  $x_0' < x_0$  if the true value  $\mu = \mu_1(x_0)$  is  $\alpha/2$

$x_0' > x_0$  if the true value  $\mu = \mu_2(x_0)$  is  $\alpha/2$

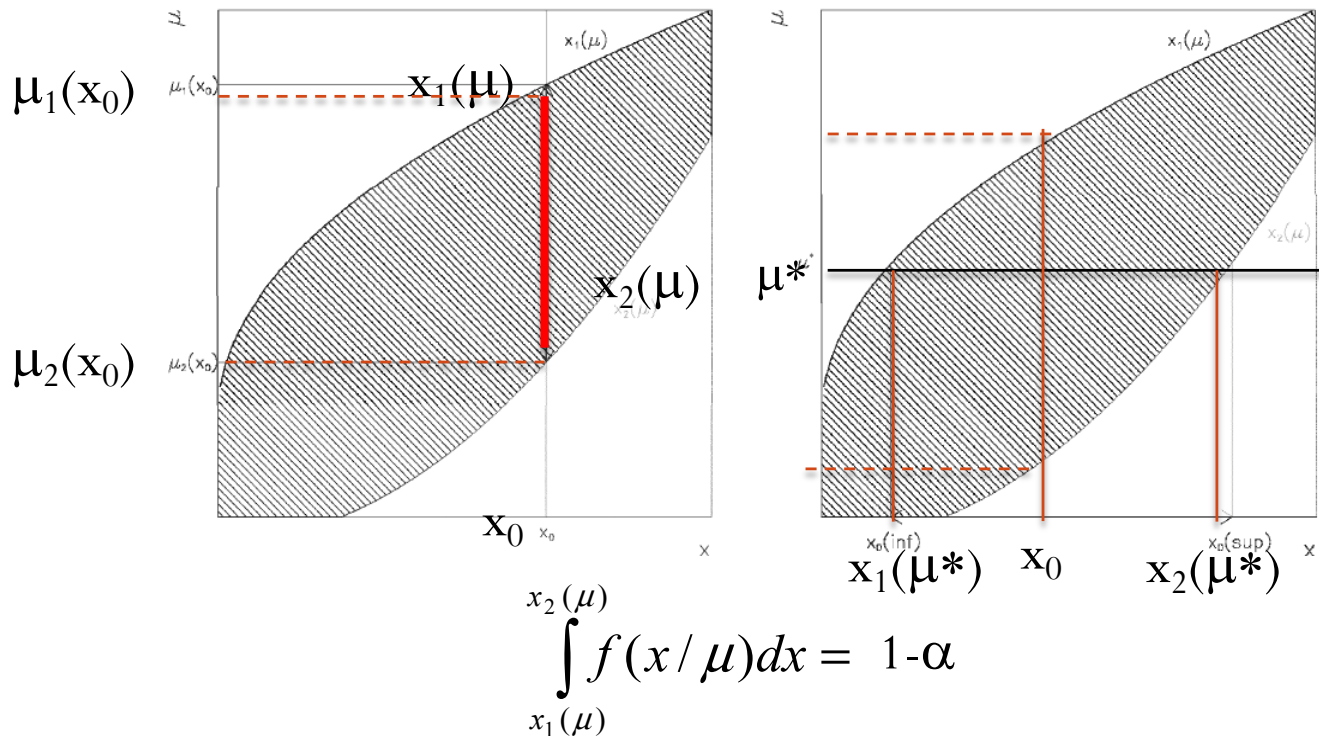
The determined C.I. is  $[\mu_2(x_0), \mu_1(x_0)]$ .

Check the correct coverage: suppose  $\mu^*$  is the true value. I repeat N times the measurement and determine each time the C.I. By construction in a fraction  $1 - \alpha$  of the cases  $x_0$  is within  $x_1(\mu^*)$  and  $x_2(\mu^*)$  and the corresponding C.I. provides coverage of  $\mu^*$ .

$$P(x_1(\mu^*) < x_0 < x_2(\mu^*)) = 1 - \alpha$$

In  $\alpha$  cases  $x_0$  lies outside the interval  $[x_1(\mu^*), x_2(\mu^*)]$  and the corresponding C.I. does not cover  $\mu^*$ .

## Neyman's construction



By construction the probability to measure  $x_0' < x_0$  if the true value  $\mu = \mu_1(x_0)$  is  $\alpha/2$

$x_0' > x_0$  if the true value  $\mu = \mu_2(x_0)$  is  $\alpha/2$

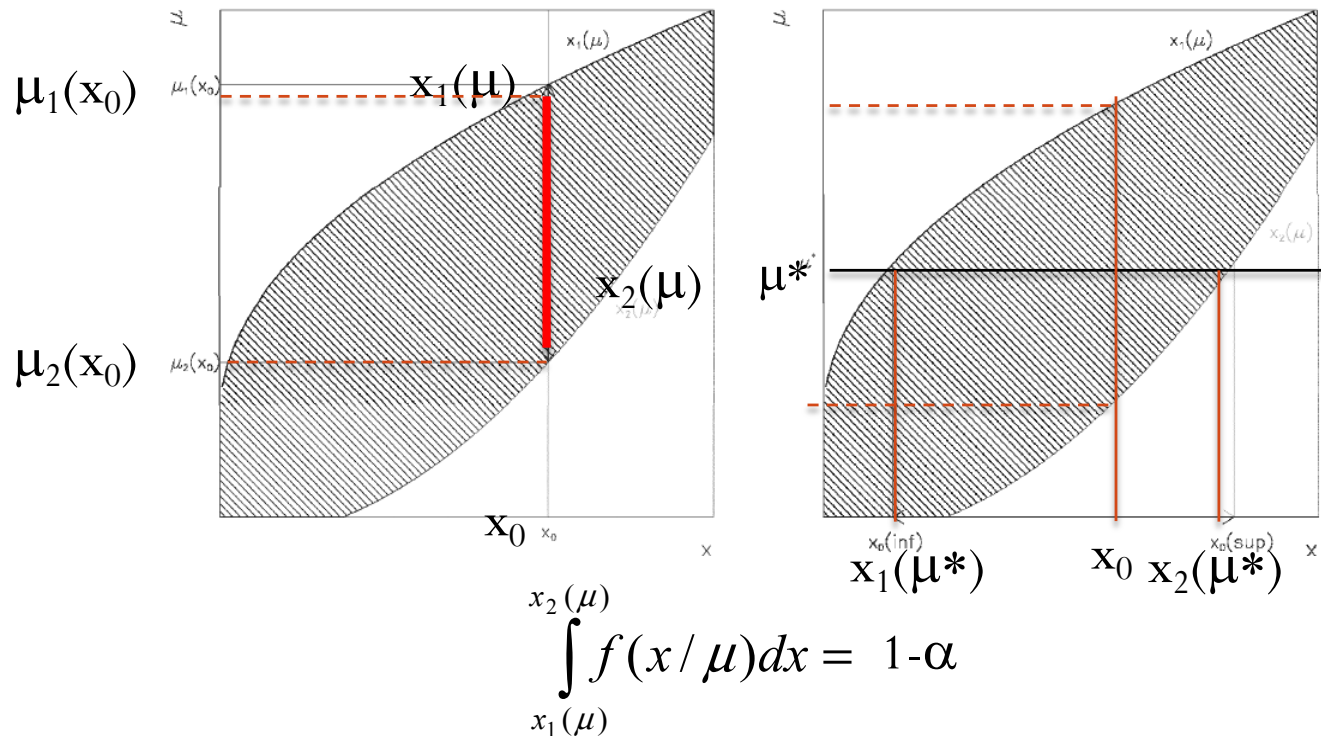
The determined C.I. is  $[\mu_2(x_0), \mu_1(x_0)]$ .

Check the correct coverage: suppose  $\mu^*$  is the true value. I repeat N times the measurement and determine each time the C.I. By construction in a fraction  $1-\alpha$  of the cases  $x_0$  is within  $x_1(\mu^*)$  and  $x_2(\mu^*)$  and the corresponding C.I. provides coverage of  $\mu^*$ .

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## Neyman's construction



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The determined C.I. is  $[\mu_2(x_0), \mu_1(x_0)]$ .

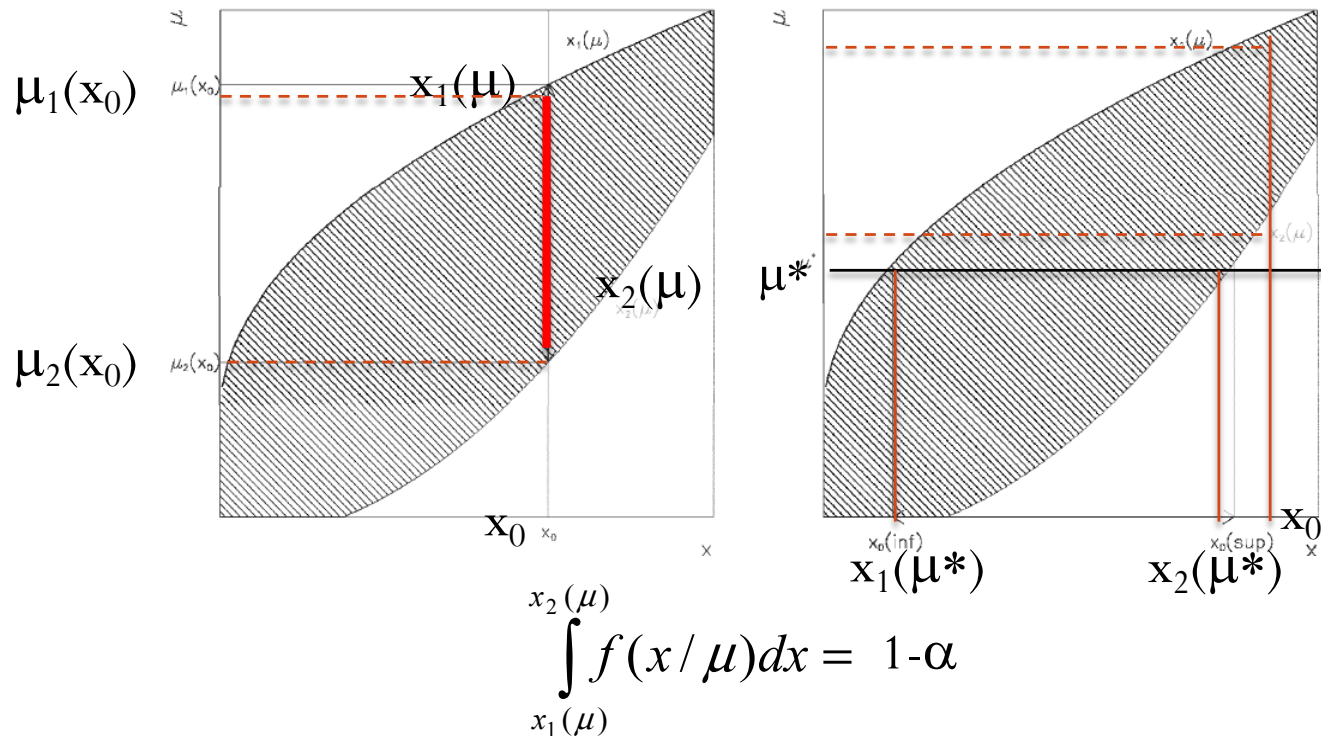
Check the correct coverage: suppose  $\mu^*$  is the true value. I repeat N times the measurement and determine each time the C.I. By construction in a fraction  $1-\alpha$  of the cases  $x_0$  is within  $x_1(\mu^*)$  and  $x_2(\mu^*)$  and the corresponding C.I. provides coverage of  $\mu^*$ .

$$P(x_1(\mu^*) < x_0 < x_2(\mu^*)) = 1-\alpha$$

In  $\alpha$  cases  $x_0$  lies outside the interval  $[x_1(\mu^*), x_2(\mu^*)]$  and the corresponding C.I. does not cover  $\mu^*$ .



## Neyman's construction



By construction the probability to measure  $x_0' < x_0$  if the true value  $\mu = \mu_1(x_0)$  is  $\alpha/2$

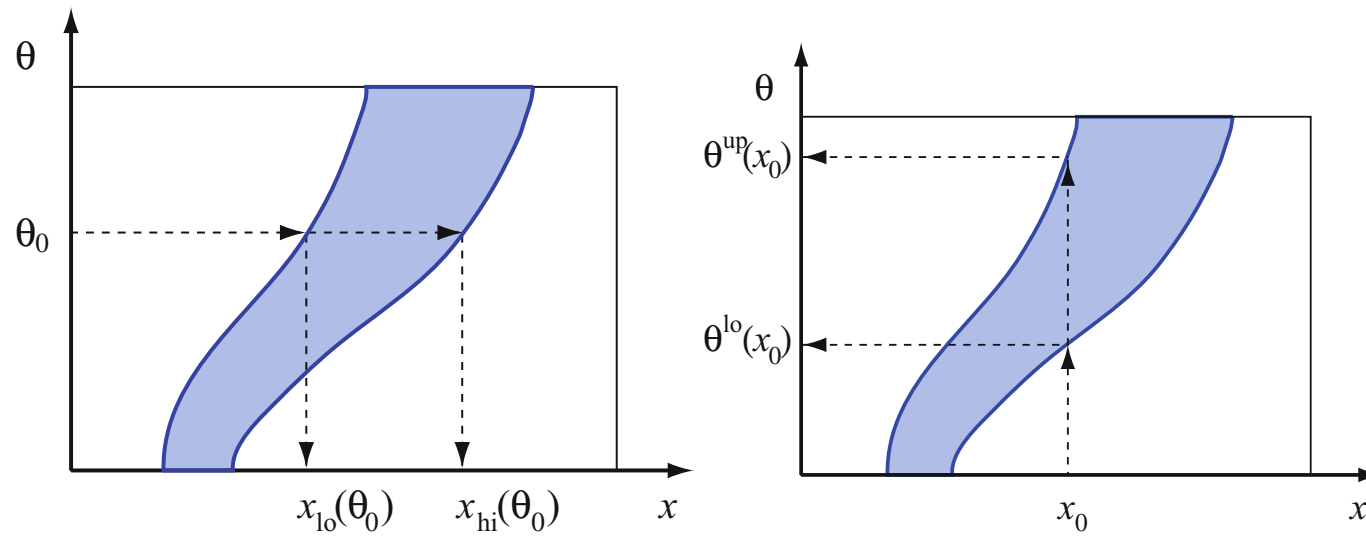
$x_0' > x_0$  if the true value  $\mu = \mu_2(x_0)$  is  $\alpha/2$

The determined C.I. is  $[\mu_2(x_0), \mu_1(x_0)]$ .

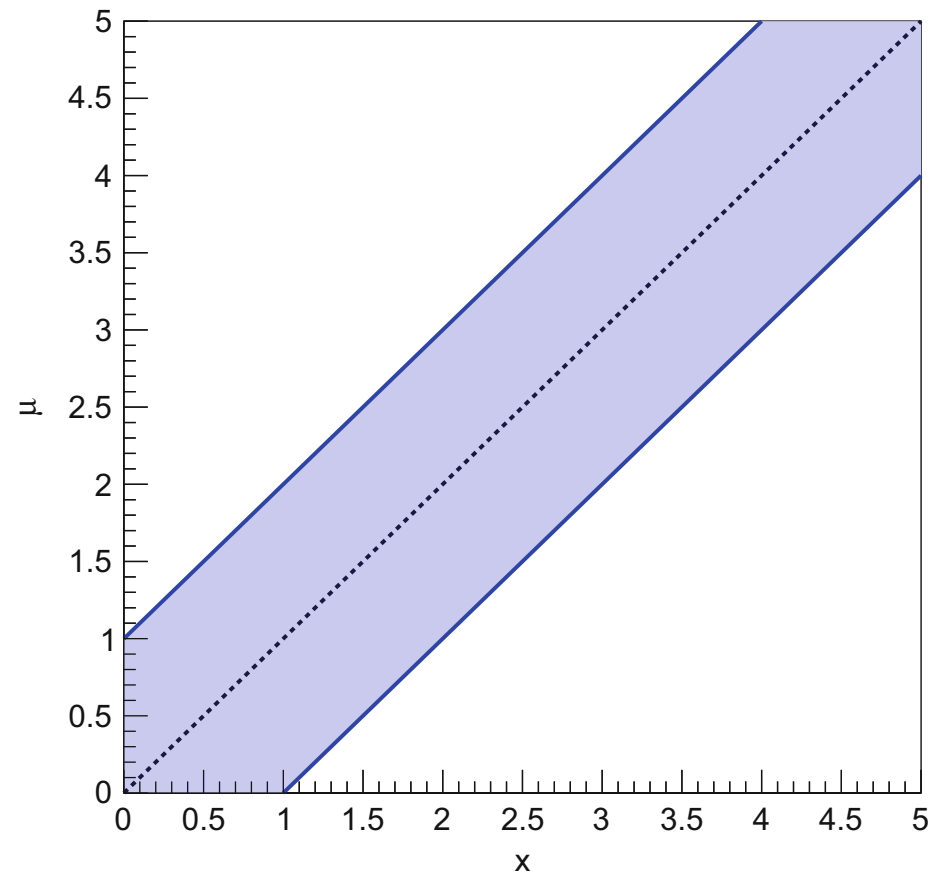
Check the correct coverage: suppose  $\mu^*$  is the true value. I repeat N times the measurement and determine each time the C.I. By construction in a fraction  $1-\alpha$  of the cases  $x_0$  is within  $x_1(\mu^*)$  and  $x_2(\mu^*)$  and the corresponding C.I. provides coverage of  $\mu^*$ .

$$P(x_1(\mu^*) < x_0 < x_2(\mu^*)) = 1-\alpha$$

In  $\alpha$  cases  $x_0$  lies outside the interval  $[x_1(\mu^*), x_2(\mu^*)]$  and the corresponding C.I. does not cover  $\mu^*$ .



**Fig. 7.1** Graphical illustration of Neyman belt construction (*left*) and inversion (*right*)



**Fig. 7.3** Neyman belt for the parameter  $\mu$  of a Gaussian with  $\sigma = 1$  at the 68.27% confidence level

Suppose Poisson variable and  $n=0$  is measured (no background) Upper limit (lower limit =0)  
 $\Rightarrow 0 \pm 0$  (freq) or  $1 \pm 1$  (Bayes) ?

By construction the probability to measure  $x_0' < x_0$  if the true value  $\mu = \mu_1(x_0)$  is  $(1-\alpha)$  (only one limit)  
 or the probability to measure  $x_0' > x_0$  if the true value  $\mu = \mu_1(x_0)$  is  $\alpha$

$$P(n > 0 / \lambda) = \sum_{n=1}^{\infty} \frac{\lambda^n e^{-\lambda}}{n!} = 1 - e^{-\lambda} = \alpha$$

frequentist

$$\bar{\lambda} = -\ln(1 - \alpha)$$

Note:

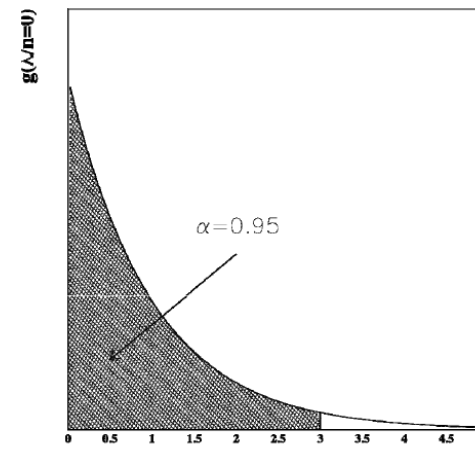
in this example  
 $\alpha$  has complementary  
 meaning than in  
 previous slides  
 $\alpha \Rightarrow 1 - \alpha$

$$g(\lambda / n = 0) = \frac{p(n = 0 / \lambda) f_0(\lambda)}{\int_0^{\infty} p(n = 0 / \lambda) f_0(\lambda) d\lambda} = \frac{e^{-\lambda}}{\int_0^{\infty} e^{-\lambda} d\lambda} = e^{-\lambda}$$

Bayesian  
 (uniform prior)

$$p(\lambda < \bar{\lambda}) = \int_0^{\bar{\lambda}} e^{-\lambda} d\lambda = 1 - e^{-\bar{\lambda}} = \alpha$$

	90%	95%	99%
$\bar{\lambda}$	2.3	3.0	4.6



## frequentist limits

By construction the probability to measure  $n_0' < n_0$  if the true value  $s = s_{up}(n_0)$  is  $(1 - \beta)$  (only one limit) or the probability to measure  $n_0' > n_0$  if the true value  $s = s_{up}(n_0)$  is  $\beta$

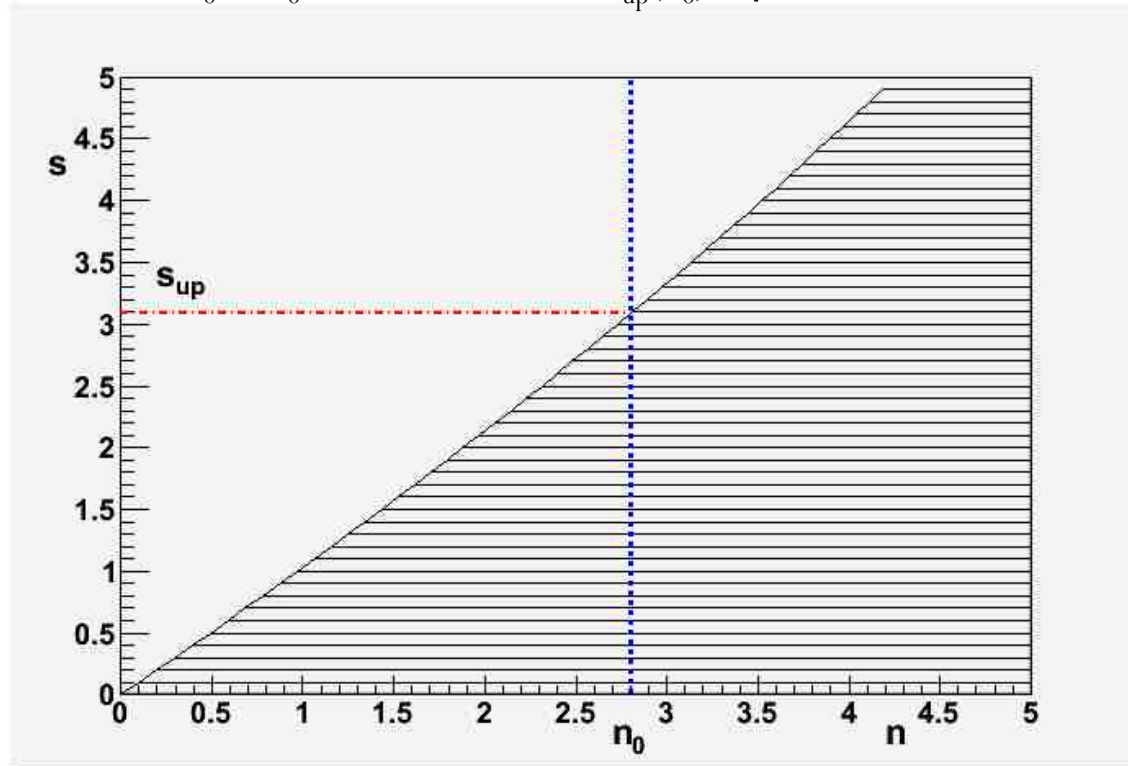
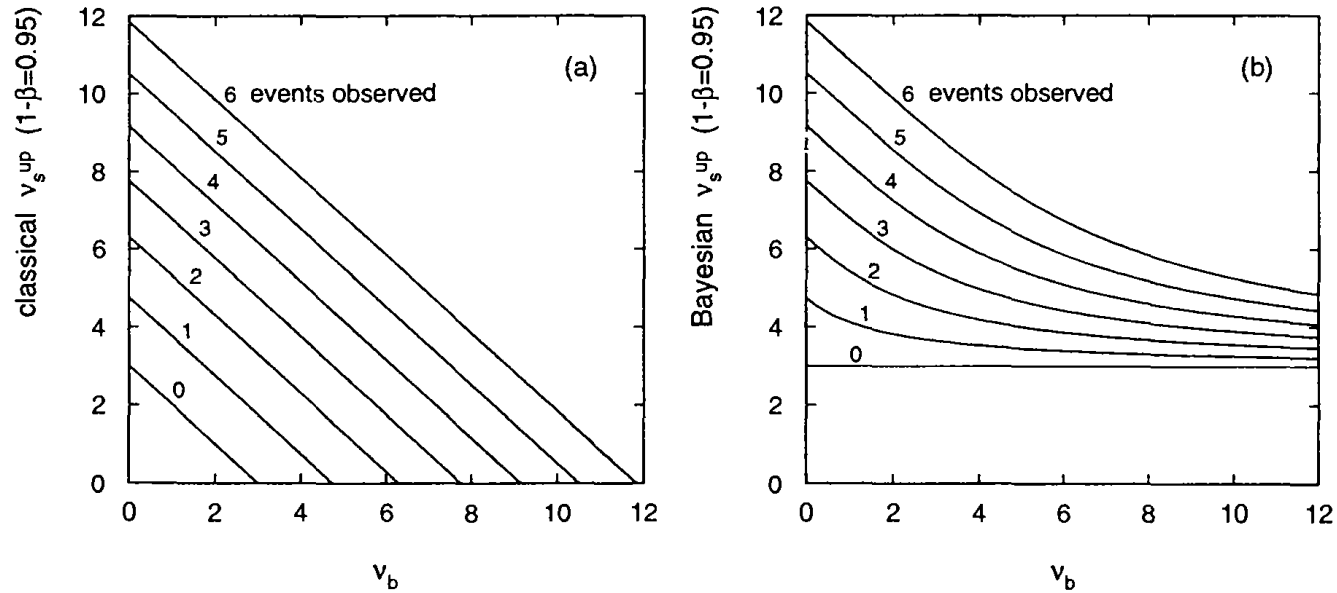


FIGURE 19. Neyman construction for the case of an upper limit. In this case a segment between  $n_1(\theta)$  and  $\infty$  is drawn for each value of the parameter  $\theta$ . The segments define the confidence region. Once a value of  $n$ ,  $n_0$  is obtained, the upper limit  $s_{up}$  is found. (For simplicity the discrete variable  $n$  is considered as a real number here).

Poisson



**Fig. 9.9** Upper limits  $\nu_s^{\text{up}}$  at a confidence level of  $1 - \beta = 0.95$  for different numbers of events observed  $n_{\text{obs}}$  and as a function of the expected number of background events  $\nu_b$ . (a) The classical limit. (b) The Bayesian limit based on a uniform prior density for  $\nu_s$ .