

The evidence:

$$\mathcal{E} = \int L(d\vec{\theta}(\mu)) \cdot p(\vec{\theta}|\mu) d\vec{\theta}$$

→ completely integrates out all parameters : also called „marginal likelihood“
 = the likelihood of obtaining the data at all

\mathcal{E} is a numerical beast.

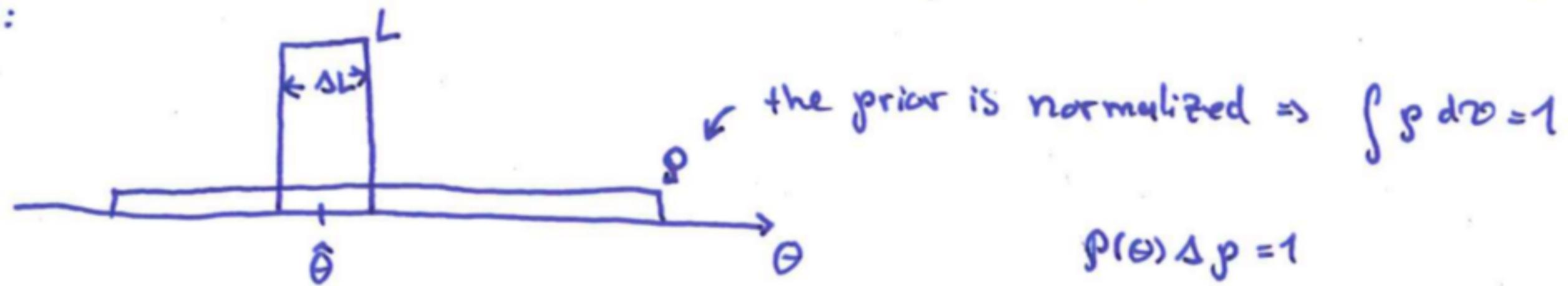
But it does the following / is needed for the following:

- Normalization of the Posterior
- penalize too complicated models (Def. "too": too complicated in the light of the available data)
 - ↪ „Occam's Razor“
- Compare the likelihoods of different models
- test how many parameters the data want

Occam's razor:

The evidence will sink with wide prior ranges → predictivity of a model is rewarded

Toy model:



$$\Rightarrow \mathcal{E} = \int L(d\vec{\theta}) p(\vec{\theta}) d\vec{\theta} = L(\vec{\theta}) p(\vec{\theta}) \Delta L = L(\vec{\theta}) \frac{\Delta L}{\Delta p}$$

will boost
the evidence
of a very good fit.

wide prior ranges (i.e. an
"unpredictive model")
bring the evidence down

→ we'll apply the priors and the evidence for
Model Comparisons

Model comparisons

E.g.: "standard model or, beyond standard model?"

- Model with 4 or 40 parameters?
- Statistical fluke or new physics?

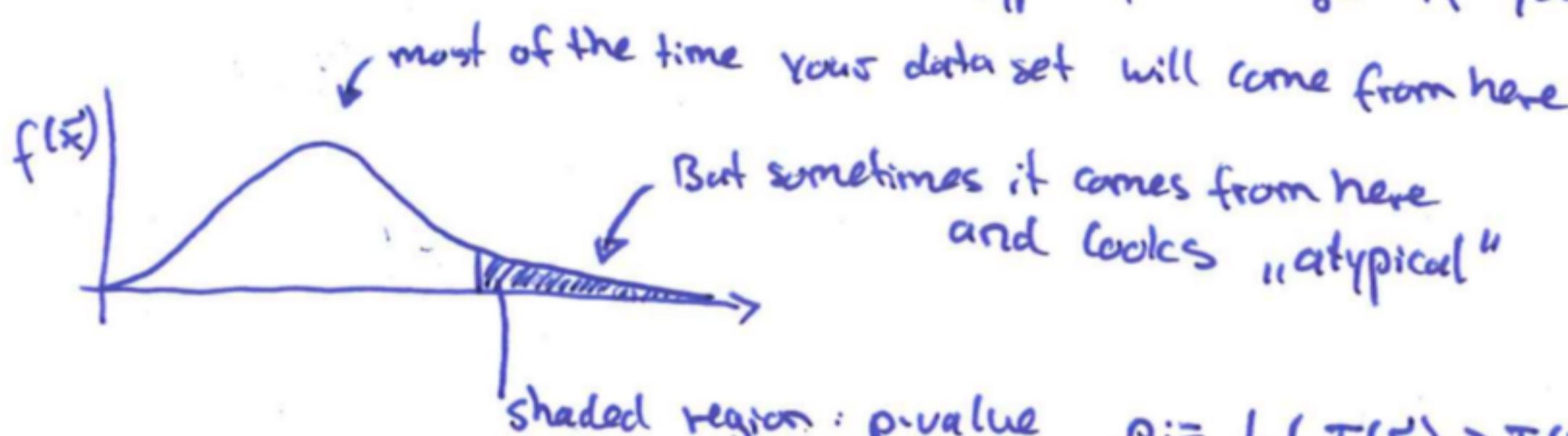
Model comparisons are rather easy to deal with in the Bayesian context, because the data are taken as fixed, and the credibility of the fitted model is investigated.

In contrast: Frequentist has $L(x|H)$, which is the likelihood that Hypothesis H has produced the data x.

But: $L(x|H) \neq L(H)$! From $L(x|H)$ we can't tell the probability of the Hypothesis.
 ↓
 in the sense of "true" with a certain ~~prob~~ probability.

Let's investigate a Frequentist's p-value:

the measured data set \tilde{x} stems from a probability distribution $f(\tilde{x})$, which tells you how your measured data set would typically change if you measure another time



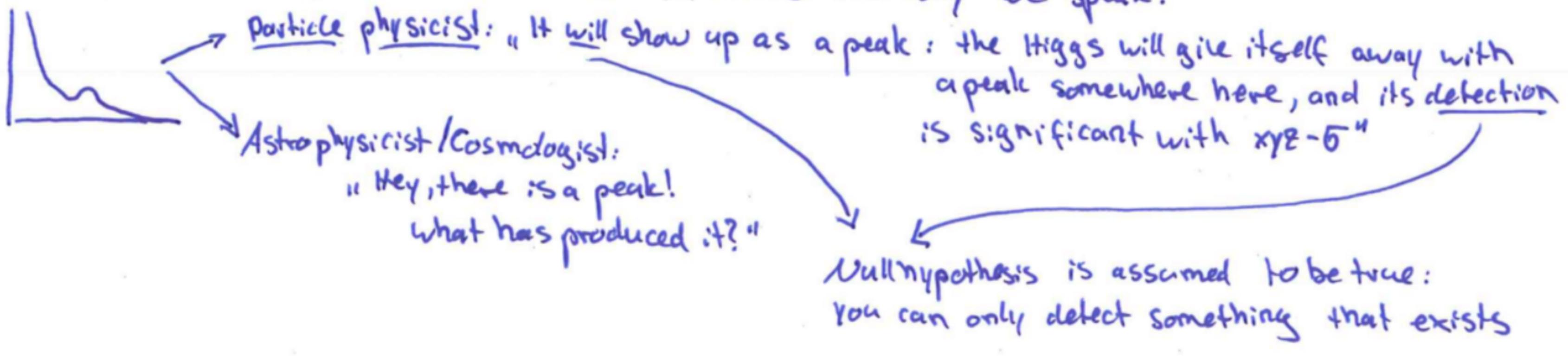
shaded region: p-value $p := L(T(\tilde{x}) \geq T(x_{\text{obs}}))$ if T is some statistics on your data set
 ↓
 "the likelihood that the statistics can turn out as extreme, or more extreme, than in the data set which you analyse."

E.g. $p\text{-value} = 0.05$: If the Hypothesis were true (which we don't know), then only in 5% of the times that you measure, it would produce you a data set as extreme, or more extreme, than the one you just calculated your p-value from.

{→ already while calculating the p-value, you assume the hypothesis to be true}

⇒ $L(x|H) \neq L(H)$ the likelihood that Hypothesis H has produced the data x is not the likelihood of H to be true.

(Not) Assuming the null hypothesis to be true influences the way we speak:



Why can't you use a p-value for model comparison?

- Because while calculating it, you already assume the Hypothesis to be true.
- Because it answers the Question „If the Hypothesis is true, then x% of the times...“

→ Example: one data set \vec{x} , two competing Hypothesis H_1 and H_2 , only one is true.

Calculate $p_1 = L(T(\vec{x}|H_1) \geq T(\vec{x}_{\text{obs}}|H_1)) \rightsquigarrow$ assumes H_1 is true } one of these calculations must base on wrong assumptions

Calculate $p_2 = L(T(\vec{x}|H_2) \geq T(\vec{x}_{\text{obs}}|H_2)) \rightsquigarrow$ assumes H_2 is true }

calculating p-values doesn't allow to compare Models to each other

Bayesian Model comparison

We had the evidence

$$E = p(d|M) = \int \underbrace{L(d|\theta(M))}_{\substack{\text{likelihood} \\ (\text{data go in here})}} \underbrace{p(\theta(M))}_{\substack{\text{prior} \\ (\text{theory goes in here})}} d\theta$$

⇒ marginalizes out all parameters → evidence is the likelihood that the model M can produce the data at all.

Question of model comparison; e.g. „How likely is my model, in light of the available data?“

↓
Likelihoods are normalized, so this question would only have an answer if you know „all possible models“.

→ But you can still compare multiple models with each other:

Evidence: $E = L(d|M)$ \rightsquigarrow $L(M|d)$ is the likelihood of your model, given the data

↓
Bayes theorem:

$$L(M|d) = \frac{L(d|M)}{p(d)} \cdot \frac{p(M)}{p(M)}$$

↑ prior on model ?? \rightarrow take a ratio :)

↑ prior on data ??

2nd model: $L(M_x|d) = \frac{L(d|M_x)}{p(d)} \cdot \frac{p(M_x)}{p(M)}$

$$\Rightarrow \left\{ \frac{L(M|d)}{L(M_x|d)} = \frac{p(M)}{p(M_x)} \cdot \frac{L(d|M)}{L(d|M_x)} = \frac{p(M)}{p(M_x)} \cdot \frac{E}{E_x} \right\}$$

→ prior on data drops out,

since in both cases, the data are the same

→ $\frac{p(M)}{p(M_x)}$ is usually set to 1.

Evidence ratio:

$$\frac{L(M|d)}{L(M_x|d)} = \frac{p(d|M)}{p(d|M_x)} = \frac{\epsilon}{\epsilon_x}$$

The evidence ratio quantifies whether the data prefer M or M_x .

"In the light of data", is M or M_x more likely?

Your beforehand belief of whether M or M_x are more likely.

ϵ is what you care about.
 ϵ_x The larger the evidence, the more likely your model.

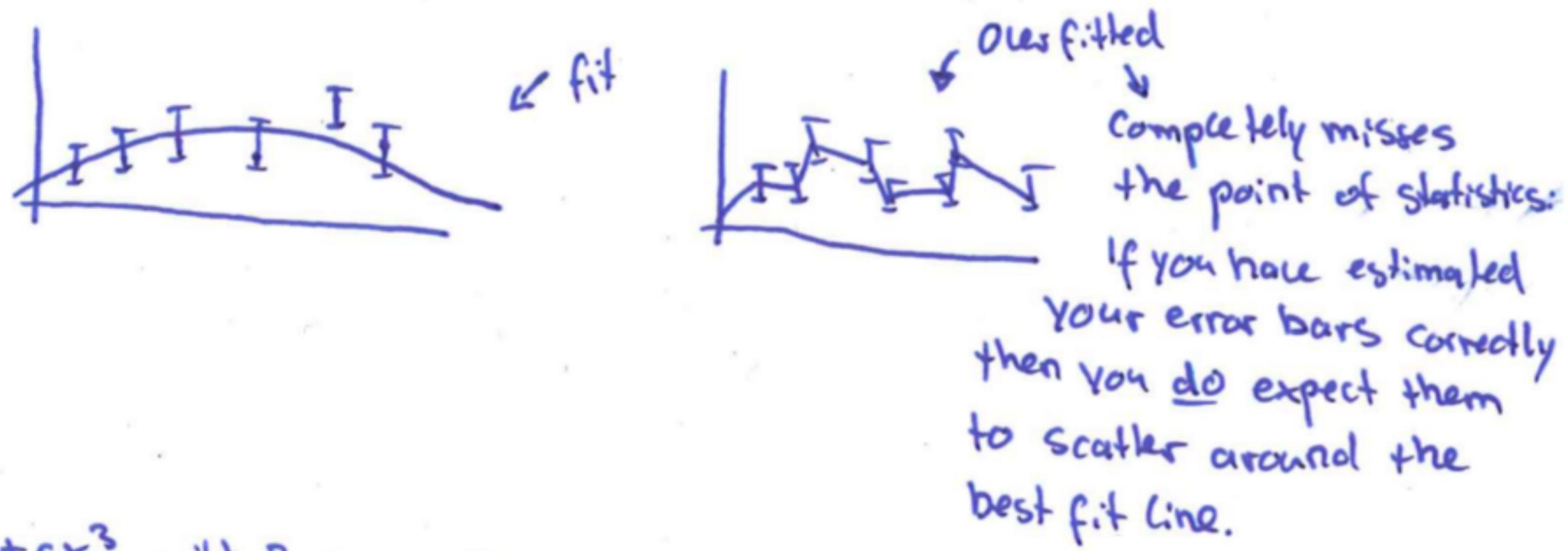
Example:

"With how many parameters should I fit my data?"

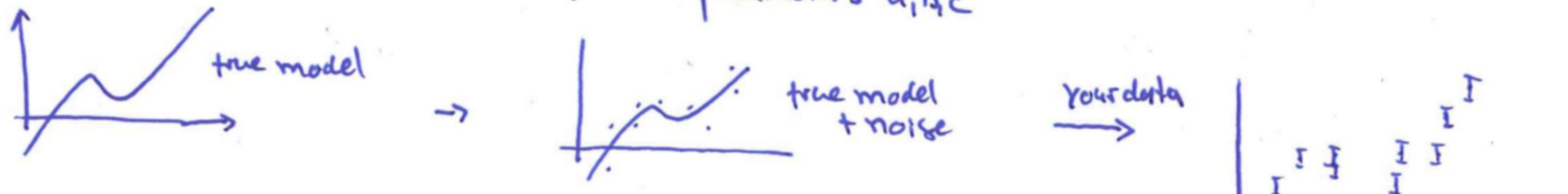
Problem: Increasing the number of parameters will always lead to an improvement χ^2 of the fit, with $I \leq 0$. \rightarrow When do you stop?

Classical problem: Overfitting / „Maßen nach Zahlen“.

usually done with χ^2



E.g. true model: $ax + bx^2 + cx^3$, with 3 parameters a, b, c

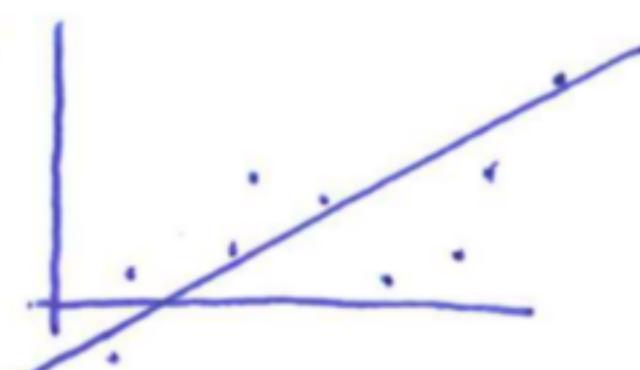


Let's find out with the evidence how many parameters the model has:
Reminder:

$$L(\hat{\theta}) \cdot \frac{\Delta L}{\Delta p} \Rightarrow \epsilon = \text{goodness of fit} \cdot \text{Occams Razor}$$

Start with 1 parameter:

fit ax :



\Rightarrow bad fit, bad $\chi^2 \Rightarrow$ low $L(\hat{\theta}) \Rightarrow$ low ϵ

fit with 2 parameters:

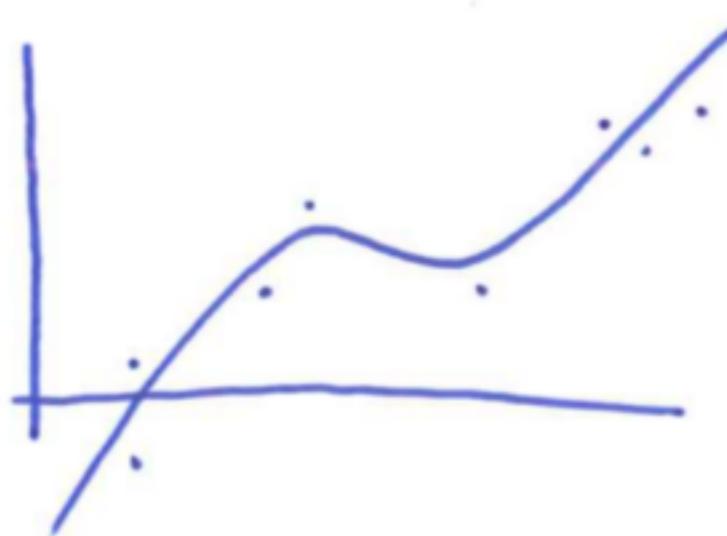
$ax + bx^2$:



\Rightarrow better fit, better $\chi^2 \Rightarrow$ larger $L(\hat{\theta}) \Rightarrow$ larger ϵ

fit with 3 parameters:

$$ax^3 + bx^2 + cx^3 \quad (\text{the actual number of parameters of the true model})$$



$$\Rightarrow \frac{\chi^2}{\deg F} \approx 1 \Rightarrow \text{large } L(\hat{\theta}) \approx \text{large } \epsilon$$

(over)fit with 4 parameters

$$ax + bx^2 + cx^3 + dx^4$$



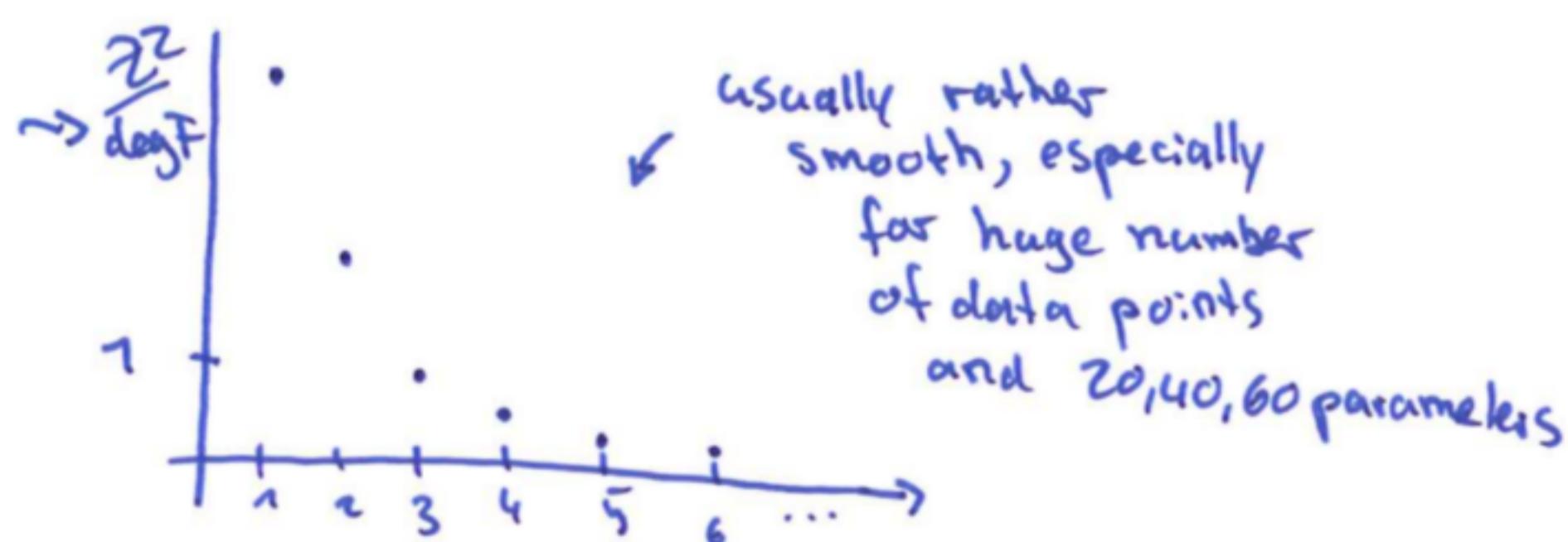
$$\frac{\chi^2}{\deg F}$$

keeps sinking. But the introduction of d does not make the fit so much better that the Occams razor factor for the new dimension is outweighed.

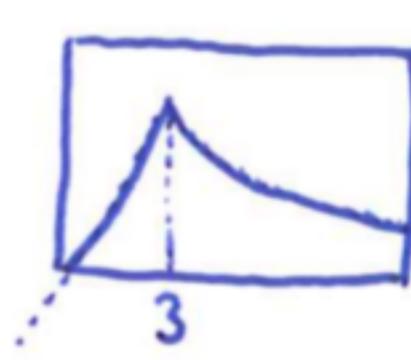
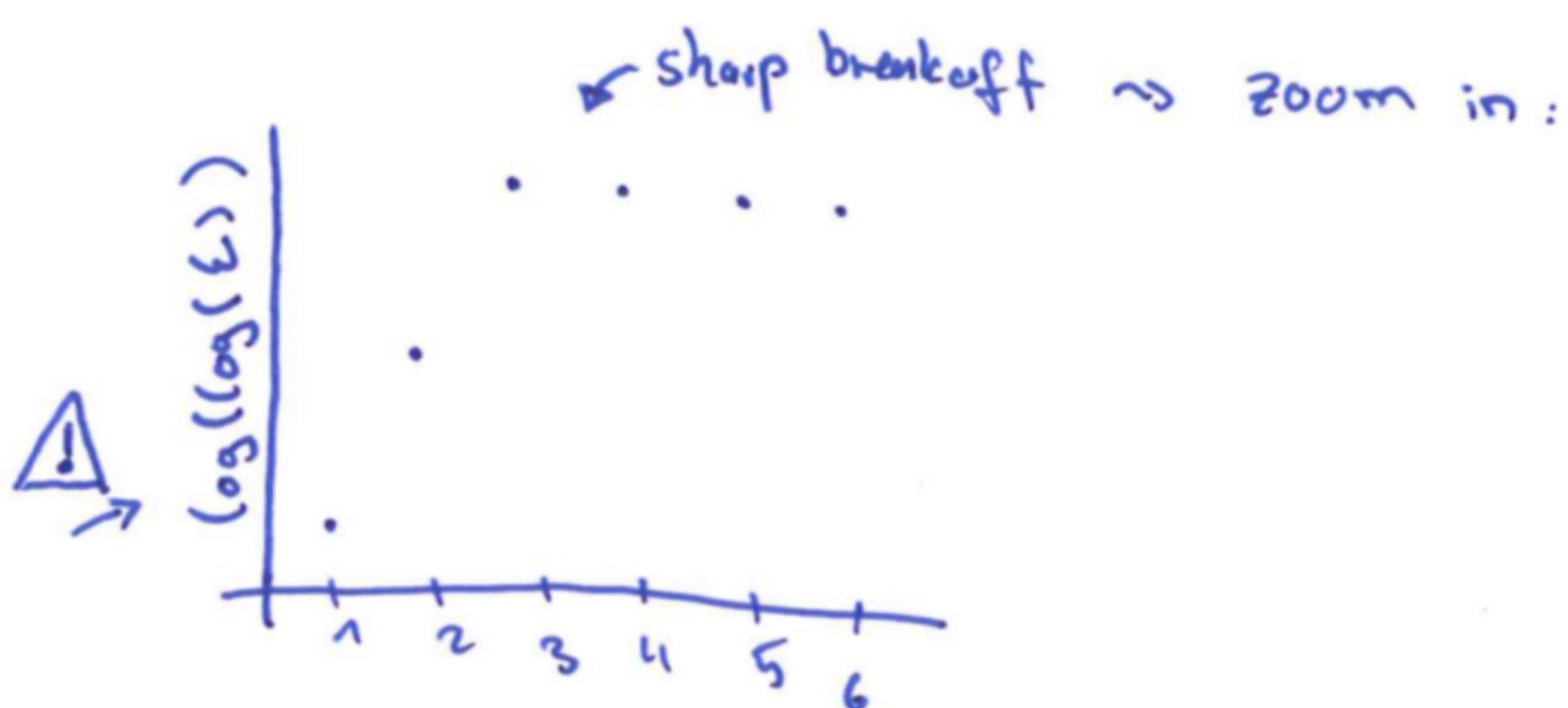
$$\frac{\Delta L(d\text{-direction})}{\Delta p(d\text{-parameter})} \text{ brings Evidence down} \leftarrow \text{large}$$

and so on, with ever more params

Results:

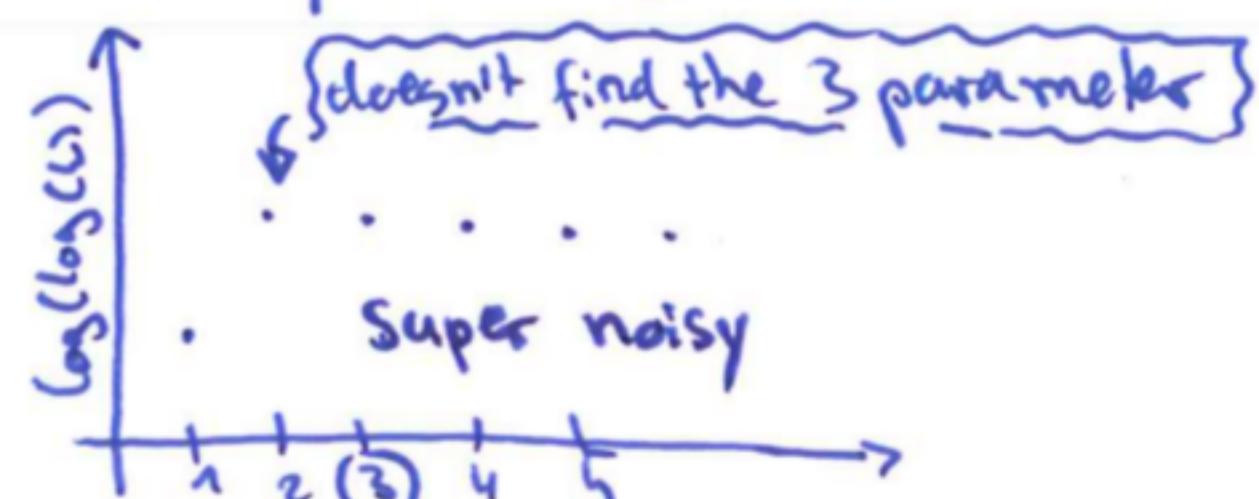
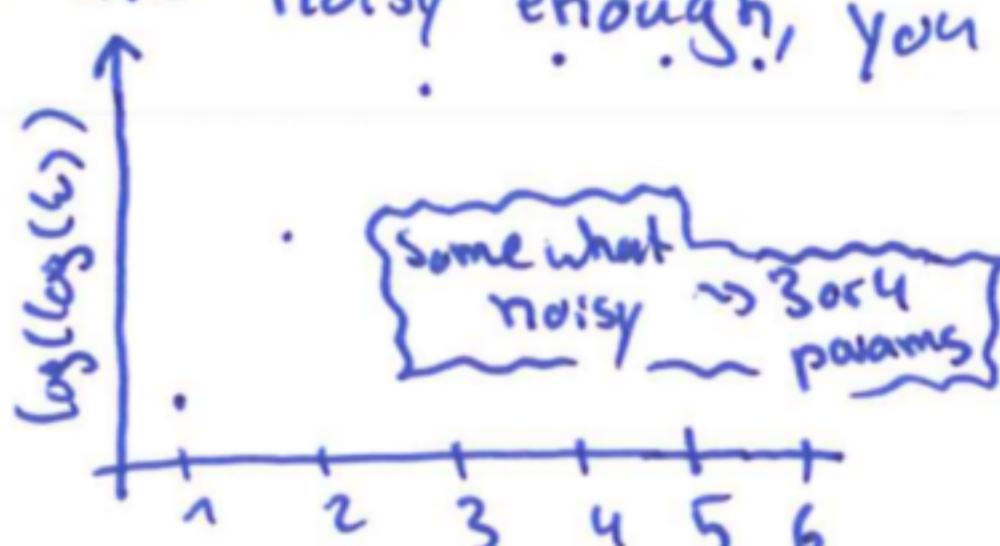


→ it can be hard to judge how many parameters are needed



↑ dropping with Occams razor is invisible on the $\log(\log(L))$ scale

- Of course, if the data are noisy enough, you can miss parameters:



2nd Example: Standard model or Beyond Standard model?

$$\frac{L(SM|d)}{L(BSM|d)} = \underbrace{\frac{p(SM)}{p(BSM)}}_{\text{Set to unity}} \cdot \frac{E(SM)}{E(BSM)}$$

without specifying here what exactly your BSM does. (If you want to code it, of course you have to specify)

$\Rightarrow \ln\left(\frac{E(SM)}{E(BSM)}\right) > 0 \Rightarrow$ Standardmodel is favoured

$< 0 \Rightarrow$ Beyond standardmodel is favoured



But with which significance? \Rightarrow Run tests and calibrate:

Jeffreys Scale:

$ \ln\left(\frac{E(SM)}{E(BSM)}\right) $	odds	"conventionally called"	Probability of the favoured model
≤ 1.0	3:1	"better data is needed"	≤ 0.75
≤ 2.5	12:1	weak evidence	0.923
≤ 5.0	$\leq 150:1$	moderate evidence	0.993
≥ 5.0	$> 150:1$	strong evidence	> 0.993

Jeffreys scale

\Rightarrow measuring more precisely (= "improving parameter constraints") will ~~bring~~ push the evidence for good fitting model up

\Rightarrow but only the evidence ratio tells whether this also happens with another model

if nature is indeed a BSM, then with the improvement of data we will observe

how $\ln\left(\frac{E(SM)}{E(BSM)}\right)$ tips from ~~>0~~ to < 0 .

— Approximations of Likelihoods & Gaussianity —

We'll find Gaussians everywhere in statistics. Why?

- Because:
- they can be handled analytically \rightarrow fast!
 - because many data are naturally Gauss-distributed
 - Physicists like to count and Poisson-statistics has a Gaussian limit
 - and if all else fails: The Central Limit Theorem

CTL: For large n , the sum of n random numbers that were drawn from a probability distribution with a finite variance σ^2 , and higher moments, tends to be gaussian, with mean of the sums expectation value and variance $n\sigma^2$ (and no higher moments, because CTL is exactly that the higher moments decrease and asymptote to zero.) So only the first two moments survive \Rightarrow Gaussian.

This works in the data space.

Does it mean that also a likelihood is usually Gaussian? \rightsquigarrow Gruplot script

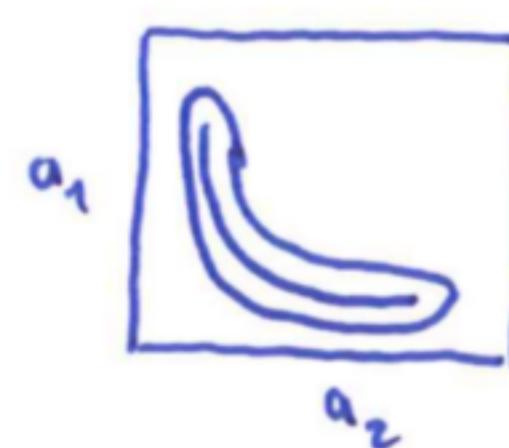
Other examples of non-Gaussian likelihoods:

- Parameter degeneracies: Imagine your data measure a function $f(A, x_1, x_2, x_3, \dots)$, where the x_i are independent variables, and A a model parameter. Let's use $f = Ax_1$



Gaussian likelihood in A , if data are Gaussian

\rightsquigarrow But maybe you know from theory that $A = a_1/a_2$, and you are interested in a_1 and $a_2 \Rightarrow$ degeneracy:



\rightsquigarrow non-Gaussian likelihood although data are Gaussian

\rightsquigarrow So one can't always approximate a likelihood with a Gaussian

\rightsquigarrow no analytical solutions

\rightsquigarrow need with numerical methods to keep computation times down (MCUC, Nested Sampling)

\rightsquigarrow or need a well-working non-Gaussian approximation