

Lecture 2: Bayesian inference in practice

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Outline

- 1 Model testing through posterior predictive checks
- 2 Why is exact Bayesian statistics hard?
- 3 Attempts to deal with the difficulty
- 4 Sampling

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Example: Modelling rainfall in Oxford

Example:

- Measure the average rainfall by month in Oxford.

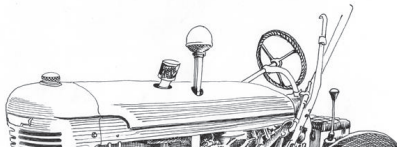


Modelling rainfall in Oxford

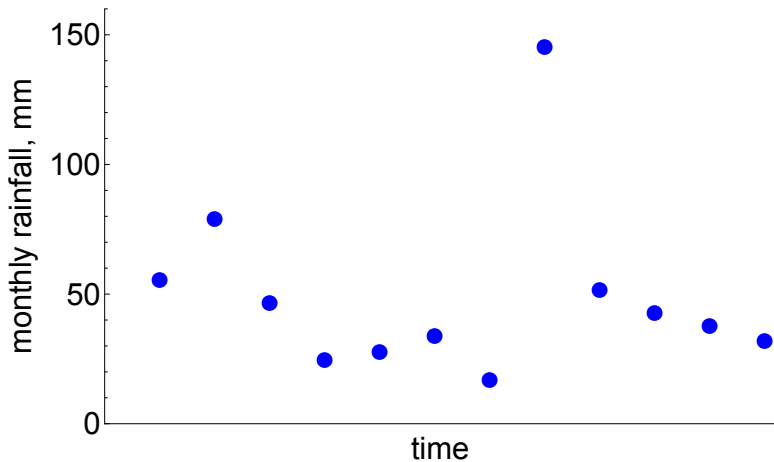
Scenario: modelling Oxford rainfall for farmers

- Government needs a model for rainfall to help plan the budget for farmers' subsidies over the next 5 years.
- Crop yields depend on rainfall following typical season patterns.
- If rainfall is persistently above normal for a number of months \implies yields \downarrow
- Assume crop more tolerant to drier spells.

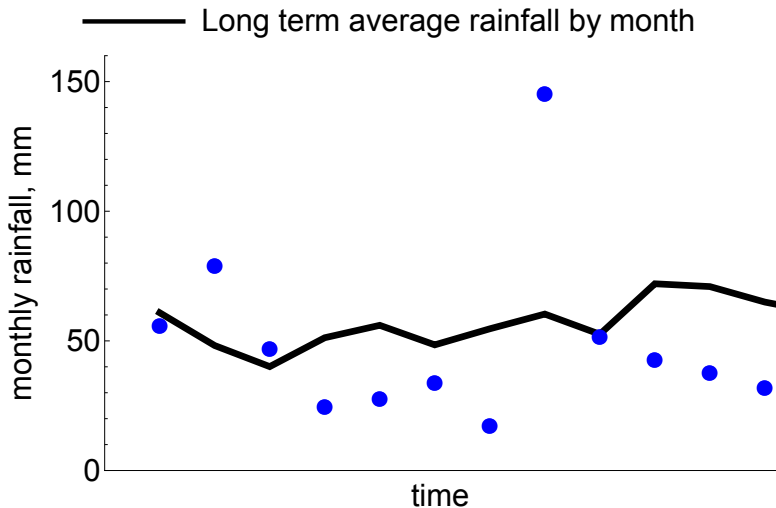
\implies create a binary variable equal to 1 if rainfall above average; 0 otherwise.



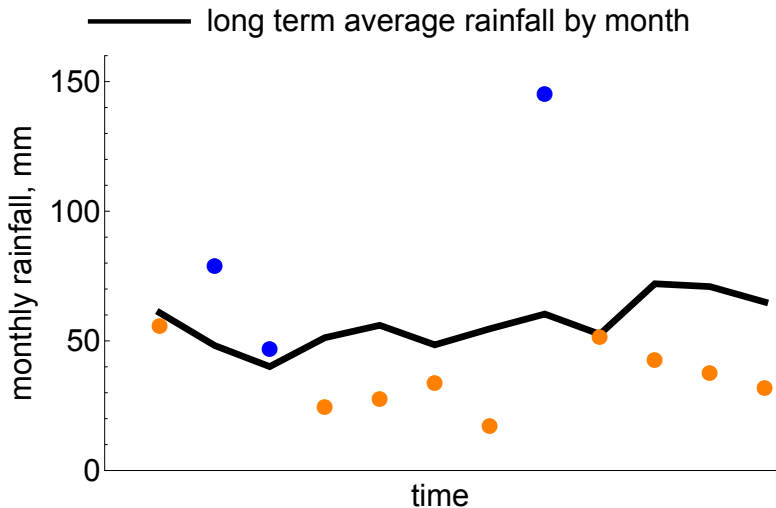
Scenario: modelling Oxford rainfall for farmers



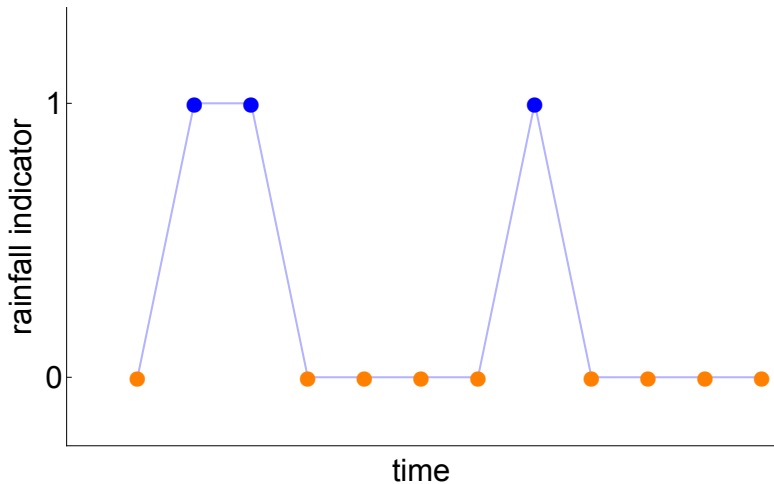
Scenario: modelling Oxford rainfall for farmers



Scenario: modelling Oxford rainfall for farmers



Scenario: modelling Oxford rainfall for farmers



Choosing a likelihood

Building a model to explain $X_t \in (0, 1)$; whether the rainfall in one month exceeds a long term monthly average.

- **Independence:** the value of X_t in month t is independent of that in the previous months.
- **Identical distribution:** all months in our sample have the same probability (θ) of rainfall exceeding long-term average.

Choosing a likelihood

Conditions:

- $X_t \in (0, 1)$ is a **discrete** random variable.
- Assume **independence** among X_t .
- Assume **identical distribution** for X_t ; probability of rainfall exceeding monthly average is θ .

\Rightarrow **Bernoulli** likelihood for each **individual** X_t .



The Bernoulli likelihood

X_t measures whether or not the rainfall in a month t is above a long term average. A Bernoulli likelihood for a single X_t has the form:

$$p(X_t|\theta) = \theta^{X_t}(1 - \theta)^{1-X_t} \quad (1)$$

But what does this mean? Work out the probabilities *given* θ :

- $p(X_t = 1|\theta) = \theta^1(1 - \theta)^0 = \theta$
- $p(X_t = 0|\theta) = \theta^0(1 - \theta)^1 = 1 - \theta$



Likelihood vs sampling distribution

Question: what is the difference between a likelihood and a sampling/probability distribution?

Answer: they are given by the same object, but under different conditions (“the equivalence relation”). Consider a single X_t :

$$L(\theta|X_t) = p(X_t|\theta) \quad (2)$$

- If hold θ constant \implies sampling distribution
 $X_t \sim p(X_t|\theta)$.
- If hold X_t constant \implies likelihood distribution
 $\theta \sim L(\theta|X_t)$.
- In Bayes' rule we vary $\theta \implies$ we use the **likelihood** interpretation.

Likelihood vs sampling distribution

Sampling distribution: hold **parameter** constant, for example $\theta = 0.75$:

$$Pr(X_t = 1 | \theta = 0.75) = 0.75^1 (1 - 0.75)^0 = 0.75$$

$$Pr(X_t = 0 | \theta = 0.75) = 0.75^0 (1 - 0.75)^1 = 0.25$$

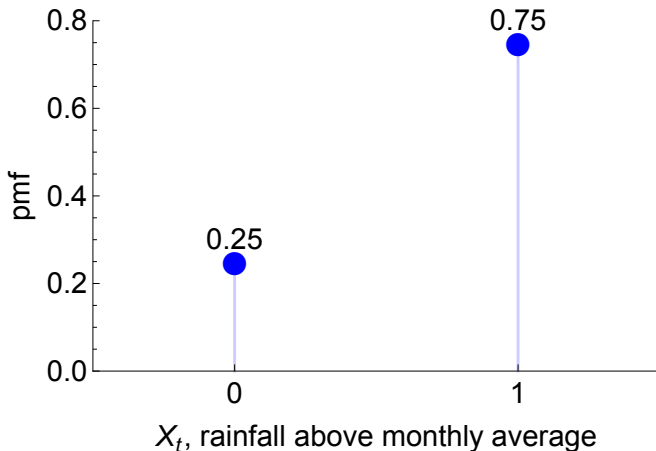
Likelihood distribution: hold **data** constant for example consider $X_t = 1$:

$$L(\theta | X_t = 1) = \theta^1 (1 - \theta)^0 = \theta \quad (3)$$

Therefore here the sampling distribution is **discrete** whereas the likelihood distribution is **continuous**.

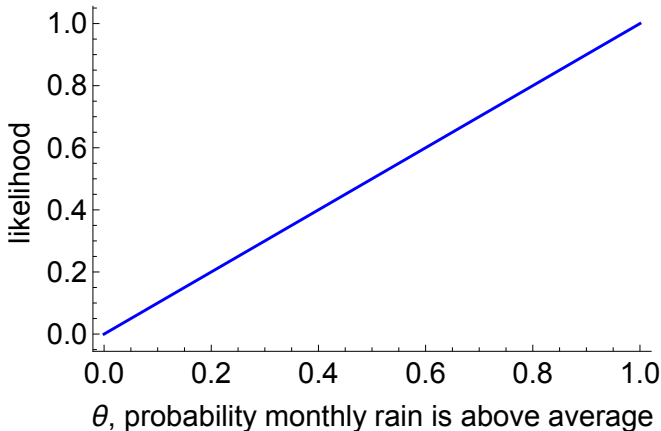
Likelihood vs sampling distribution

Sampling distribution: hold θ constant and vary the data X_t
 \Rightarrow valid probability distribution. For example for $\theta = 0.75$:



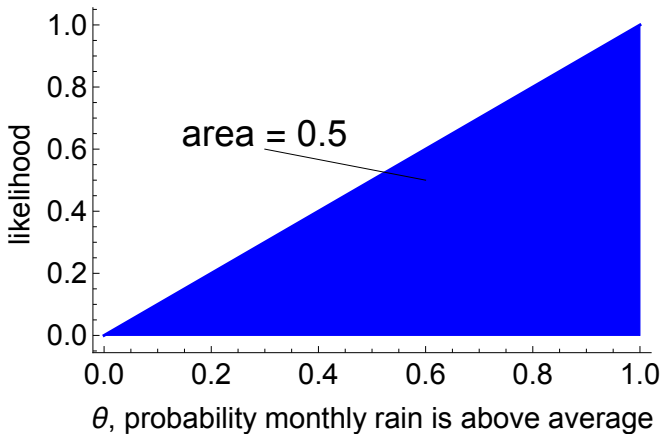
Likelihood vs sampling distribution

Likelihood: hold $X_t = 1$ and vary θ
 $\Rightarrow L(\theta|X_t = 1) = \theta^1(1 - \theta)^0 = \theta$:



Likelihood vs sampling distribution

Likelihood: hold $X_t = 1$ and vary θ . Not a valid probability distribution!



The overall likelihood

Now assuming that we have a series of $X = (X_1, X_2, \dots, X_T)$.

Question: How do we obtain the full likelihood? By **independence:**

$$\begin{aligned} p(X_1, X_2, \dots, X_T | \theta) &= \theta^{X_1} (1 - \theta)^{1-X_1} \times \theta^{X_2} (1 - \theta)^{1-X_2} \times \dots \\ &\quad \times \theta^{X_T} (1 - \theta)^{1-X_T} \\ &= \theta^{\sum X_t} (1 - \theta)^{T - \sum X_t} \end{aligned}$$

So if we suppose rain exceeded average in 4/12 months \implies

$$L(\theta | X) = \theta^4 (1 - \theta)^8 \quad (4)$$

Posterior predictive distribution

Defined:

“The probability distribution for a new data sample \tilde{X} given our current data X .”

We obtain this by the following recipe:

- 1 Sample a value of θ_i from posterior:

$$\theta_i \sim p(\theta|X) \quad (5)$$

where X is the current data.

- 2 Sample a value of \tilde{X}_i from the sampling distribution conditional on θ_i ;

$$\tilde{X}_i \sim p(\tilde{X}|\theta_i) \quad (6)$$

- 3 Graph histogram of \tilde{X}_i values \implies posterior predictive distribution.

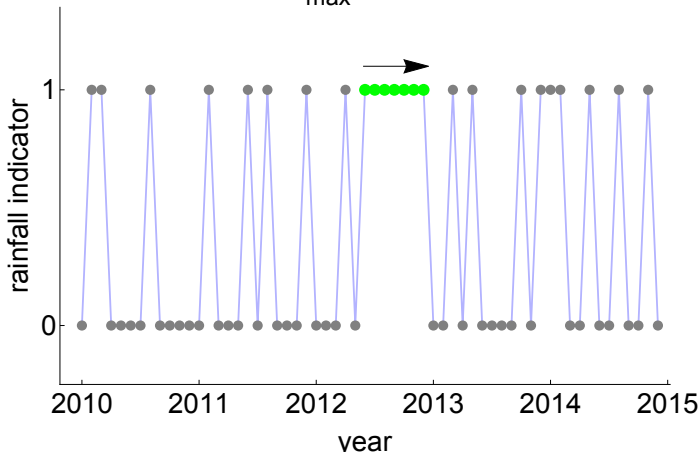
Scenario 1: key question

- Crop yields depend on whether rainfall is **persistently** above average.
- **Key question:** does the model allow for sufficient persistence in process?
- **Answer:** find the length of maximum run of consecutive $X_t = 1$ in real data. Then:
 - Draw a sample data series 60 months long from the posterior predictive distribution.
 - Find maximum run of consecutive $X_t = 1$ in simulated series.
- Repeat the above steps a number of times.
- **Compare** real maximum run length with distribution of simulated run lengths.

Scenario: maximum length run of wet months in real data

- Start with real data.
- Find maximum run of $X_t = 1$ (rainfall above average).

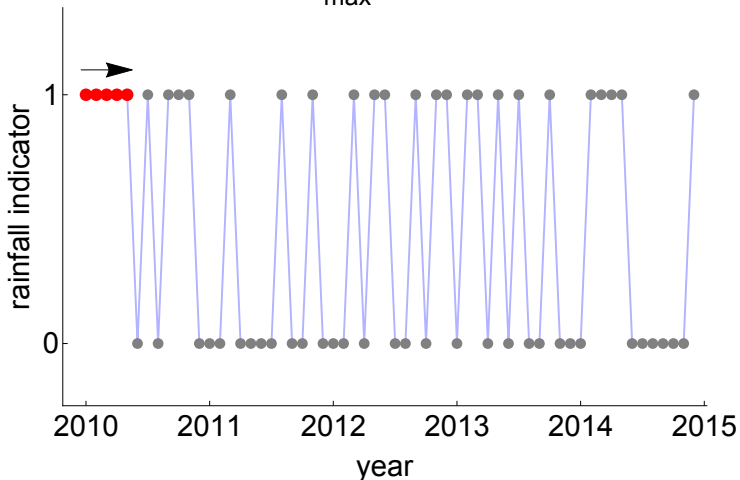
$$N_{\max}^{\text{real}} = 7$$



Scenario: posterior predictive checks

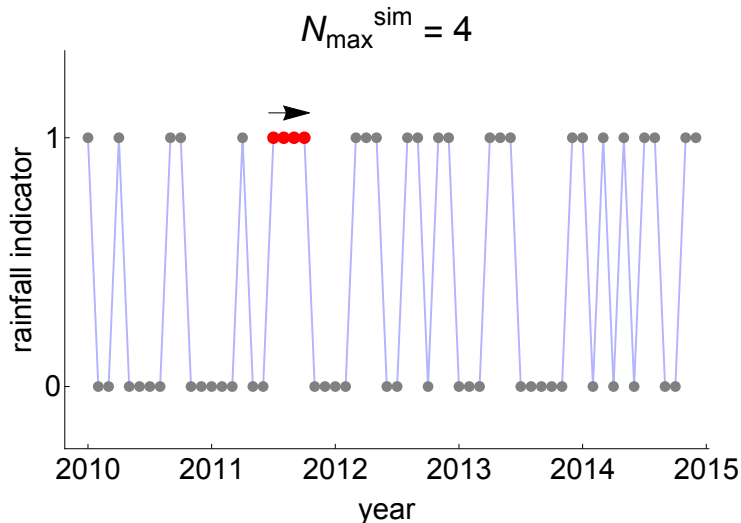
Repeating for data simulated from the posterior predictive.

$$N_{\max}^{\text{sim}} = 5$$



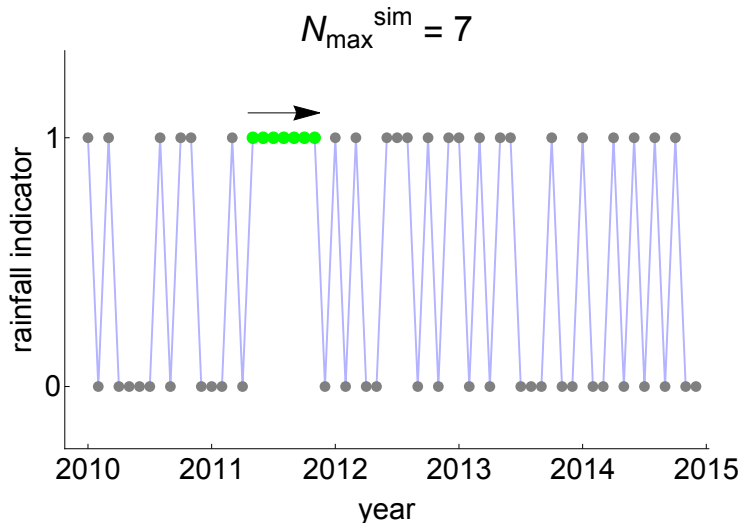
Scenario: posterior predictive checks

Another sample.



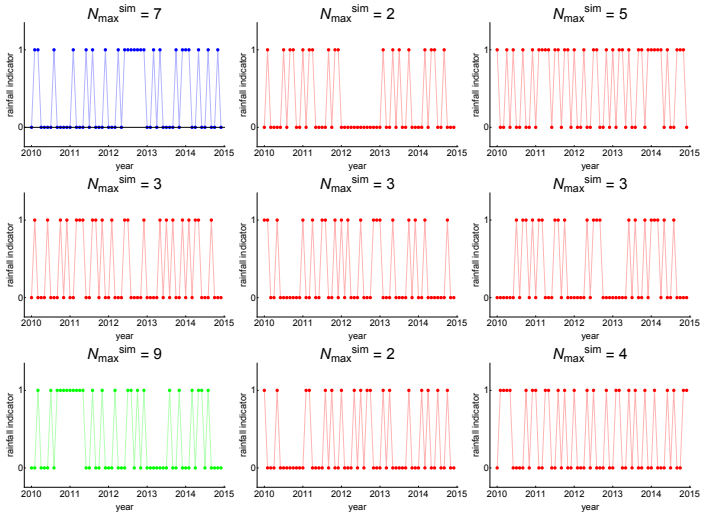
Scenario: posterior predictive checks

A further sample.



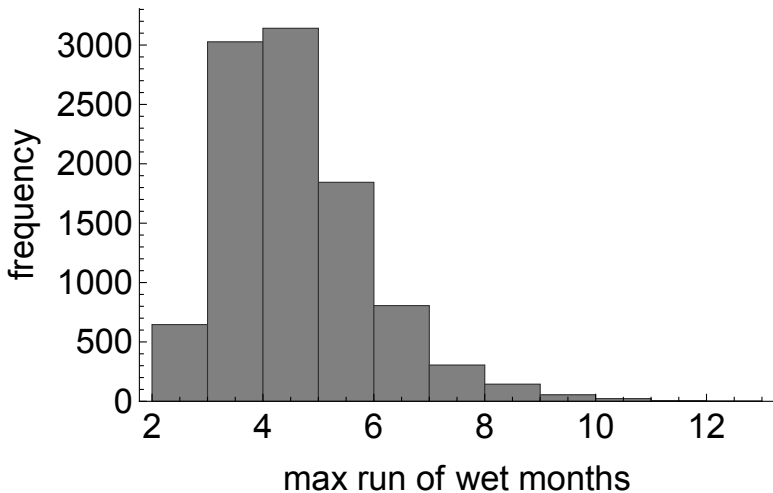
Scenario: posterior predictive checks

A number of samples.



Scenario: p value

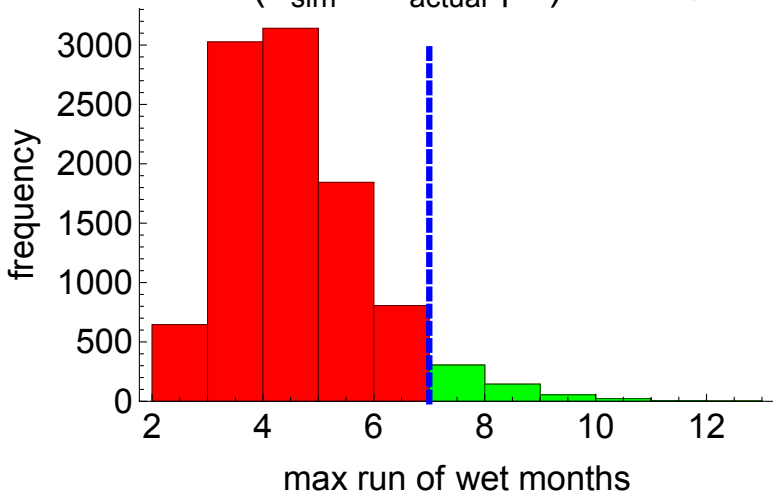
Repeat 10,000 times; each time recording maximum run length.



Scenario: p value

Find percentage of times where simulated exceeds real.

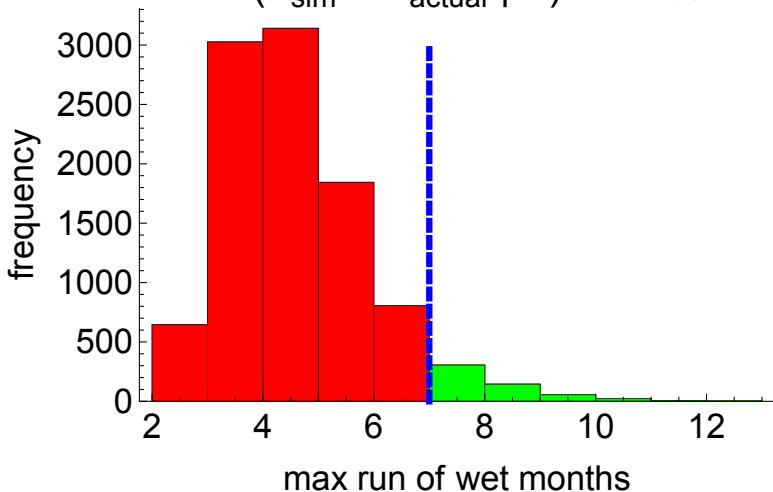
$$\Pr(T_{\text{sim}} \geq T_{\text{actual}} \mid X) = 5.0\%$$



Scenario: p value

Therefore conclude that model is not fit for purpose!

$$\Pr(T_{\text{sim}} \geq T_{\text{actual}} \mid X) = 5.0\%$$



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Example problem: paternal discrepancy

- **Paternal discrepancy** is the term given to a child who has a biological father different to their supposed biological father.
- **Question:** how common is it?
- **Answer:** a recent meta-analysis of studies of “paternal discrepancy” (PD) found a rate of $\sim 10\%^1$.
- Suppose we have data for a random sample of 10 children's presence/absence of PD.

Aim: infer the prevalence of PD in the population (θ).



Paternal discrepancy

Assume individual samples are:

- **Independent.**
- **Identically-distributed.**

Since sample size is fixed at 10 \implies binomial likelihood.

The denominator revisited

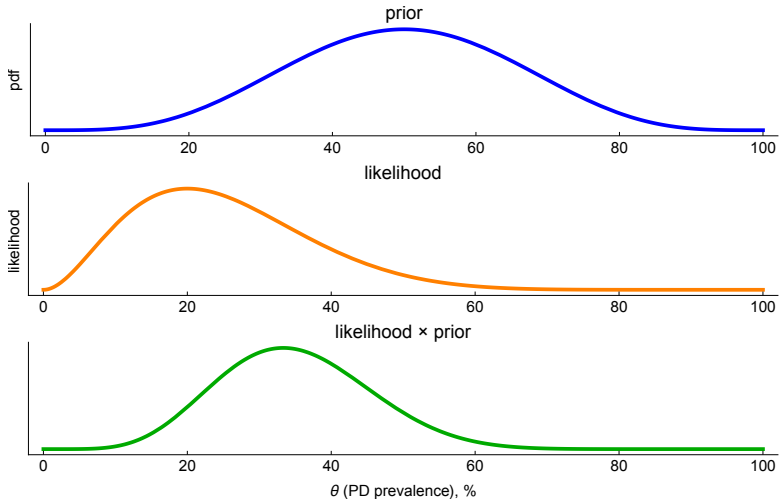
$$p(\theta|X=2) = \frac{p(X=2|\theta) \times p(\theta)}{p(X=2)} \quad (7)$$

Where we suppose we have data $X = 2$ out of a sample of 10 in our PD example. We obtain the denominator by averaging out all θ dependence. This is equivalent to integrating across all θ :

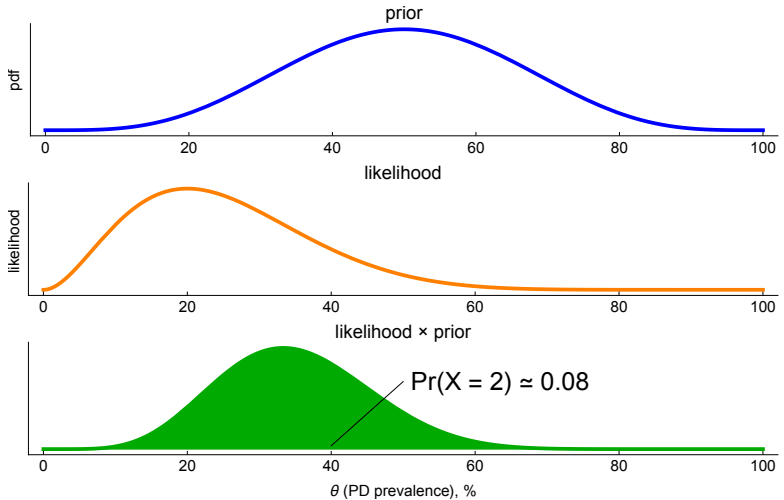
$$p(X=2) = \int_0^1 p(X=2|\theta) \times p(\theta) d\theta \quad (8)$$

(We approximately determined this using sampling previously.)

The denominator as an area



The denominator as an area

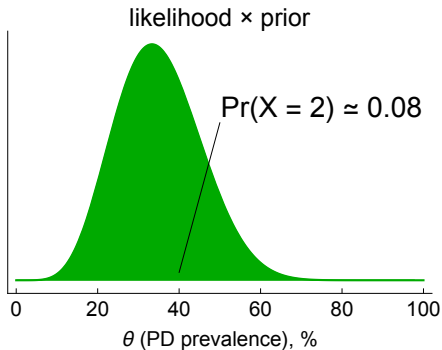


Calculating the denominator in 1 dimension

For our PD example there is a single parameter $\theta \implies$

$$p(X = 2) = \int_0^1 p(X = 2|\theta) \times p(\theta) d\theta \quad (9)$$

This is equivalent to working out an **area** under a curve.



Calculating the denominator in 2 dimensions

If we considered a different model where there were two parameters $\theta_1 \in (0, 1)$, $\theta_2 \in (0, 1) \implies$:

$$p(X = 2) = \int_0^1 \int_0^1 p(X = 2 | \theta_1, \theta_2) \times p(\theta_1, \theta_2) d\theta_1 d\theta_2 \quad (10)$$

This is equivalent to working out a **volume** contained within a surface.

Calculating the denominator in d dimensions

If we considered a different model where there were d parameters $(\theta_1, \dots, \theta_d)$ all defined to lie between 0 and 1 \implies :

$$p(X = 2) = \int_0^1 \dots \int_0^1 p(X = 2 | \theta_1, \dots, \theta_d) \times p(\theta_1, \dots, \theta_d) d\theta_1 \dots d\theta_d \quad (11)$$

This is equivalent to working out a $(d + 1)$ -dimensional **volume** contained within a d -dimensional (hyper-surface)!



The difficult denominator

- Calculating the denominator possible for $d < \sim 3$ using computers.
- Numerical quadrature and many other approximate schemes struggle for larger d .
- Many models have **thousands** of parameters.

Arrrrghhh!

Other difficult integrals

Assume we can calculate posterior:

$$p(\theta|X) = \frac{p(X|\theta) \times p(\theta)}{p(X)} \quad (12)$$

Typically we want summary measures of posterior, for example, the mean of θ_1 :

$$\begin{aligned} E(\theta_1|X) &= \int_{\Theta_1} \theta_1 \left[\int_{\Theta_2} \dots \int_{\Theta_d} p(\theta_1, \theta_2, \dots, \theta_d|X) d\theta_d \dots d\theta_2 \right] d\theta_1 \\ &= \int_{\Theta_1} \theta_1 p(\theta_1|X) d\theta_1 \end{aligned}$$

Nearly as difficult as denominator!

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What are conjugate priors?

Judicious choice of prior and likelihood can make posterior calculation trivial.

- Choose a likelihood L .
- Choose a prior $\theta \sim f \in F$, where:
 - F is a family of distributions.
 - f is a member of that **same** family.
- If posterior, $\theta|X \sim f' \in F \implies$ conjugate!
- In other words both the **prior** and **posterior** are members of the same distribution!

Conjugate priors: PD example revisited

Sample 10 children and count number (X) with PD:

- For likelihood (if independent and identically-distributed):

$$X \sim \text{Binomial}(10, \theta) \implies p(X|\theta) \propto \theta^X (1 - \theta)^{10-X} \quad (13)$$

- For prior assume a Beta distribution (a reasonable choice if $\theta \in (0, 1)$):

$$\theta \sim \text{Beta}(\alpha, \beta) \implies p(\theta) \propto \theta^{\alpha-1} (1 - \theta)^{\beta-1} \quad (14)$$

- Numerator of Bayes' rule for inference:

$$p(X|\theta) \times p(\theta) \propto \theta^X (1 - \theta)^{10-X} \times \theta^{\alpha-1} (1 - \theta)^{\beta-1} \quad (15)$$

Conjugate priors: PD example revisited

- Numerator of Bayes' rule for inference:

$$\begin{aligned} p(X|\theta) \times p(\theta) &\propto \theta^X (1 - \theta)^{10-X} \times \theta^{\alpha-1} (1 - \theta)^{\beta-1} \\ &= \theta^{X+\alpha-1} (1 - \theta)^{10-X+\beta-1} \end{aligned}$$

- This has same θ -dependence as a $Beta(X + \alpha, 10 - X + \beta)$ density \implies must be this distribution!
- \therefore a Beta prior is *conjugate* to a Binomial likelihood.

Table of common conjugate pairs of likelihoods and priors

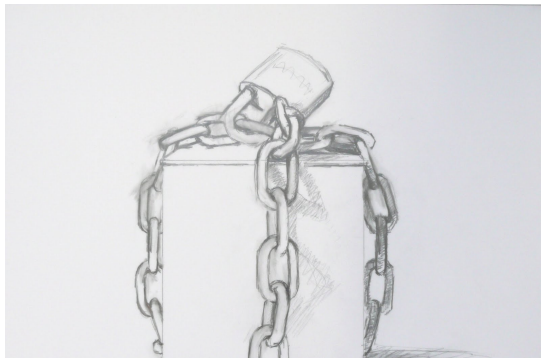
No need to do any integrals! Just lookup rules:

Likelihood	Prior	Posterior
Bernoulli	$\text{Beta}(\alpha, \beta)$	$\text{Beta}(\alpha + \sum_{i=1}^n X_i, \beta + n - \sum_{i=1}^n X_i)$
Binomial	$\text{Beta}(\alpha, \beta)$	$\text{Beta}(\alpha + \sum_{i=1}^n X_i, \beta + \sum_{i=1}^n N_i - \sum_{i=1}^n X_i)$
Poisson	$\text{Gamma}(\alpha, \beta)$	$\text{Gamma}(\alpha + \sum_{i=1}^n X_i, \beta + n)$
Multinomial	$\text{Dirichlet}(\alpha)$	$\text{Dirichlet}(\alpha + \sum_{i=1}^n \mathbf{X}_i)$
Normal	Normal-inv- Γ	Normal-inv- Γ

Limits of conjugate modelling

Using conjugate priors is limiting because:

- Often restricted to univariate problems.
 - \implies we could just use numerical quadrature instead.
- Required to use relevant conjugate prior for a given likelihood \iff may not be sufficient to capture pre-data beliefs of analyst.



Another solution: discrete Bayes' rule

- To calculate the denominator we need to do an integral, if parameters are continuous.
- If instead parameters are discrete \implies denominator is a sum over **finite** number of possible parameter values:

$$p(X) = \sum_{i=1}^p p(X|\theta_i) \times p(\theta_i) \quad (16)$$

- In general this sum is more tractable than an integral.
- **Question:** can we use this to help us with continuous parameter problems?



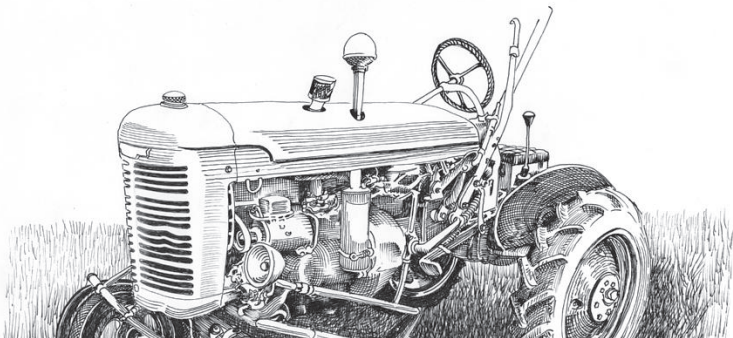
Discretised Bayesian inference

Method:

- Convert **continuous** parameter into k **discrete** values.
- Use discrete version of Bayes' rule.
- As $k \rightarrow \infty$ discrete posterior \rightarrow true posterior.

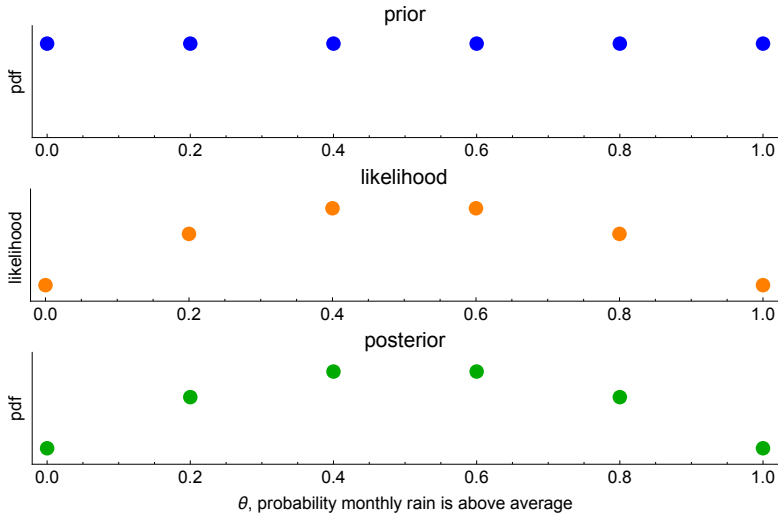
Scenario: discretised Bayesian inference

- X_t measures whether rainfall exceeds long term monthly average.
- Suppose $X_t = 1$ and $X_{t+1} = 0$.
- Assumed $p(X_t = 1, X_{t+1} = 0|\theta) = \theta(1 - \theta)$; i.e. likelihood.
- Also assume $p(\theta) = 1$; i.e. the prior.
- Discretise $\theta \in (0, 1) \rightarrow (0.0, 0.2, 0.4, 0.6, 0.8, 1.0)$.



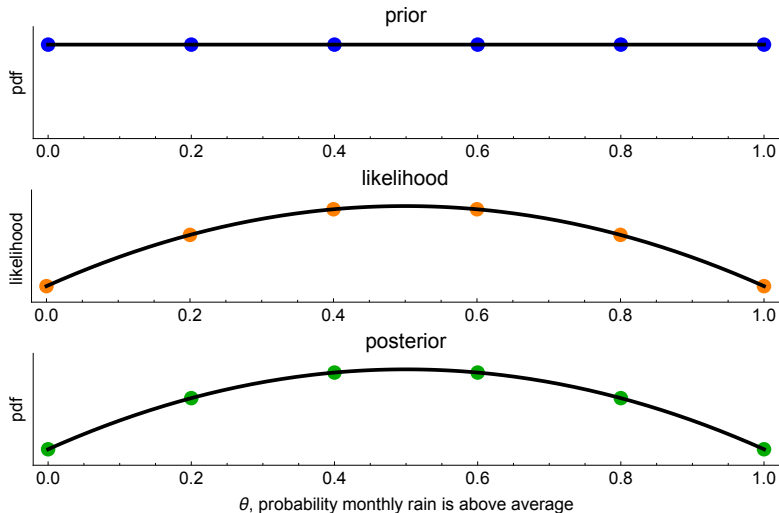
Scenario: discretised Bayesian inference

Discretise θ at intervals of 0.2.



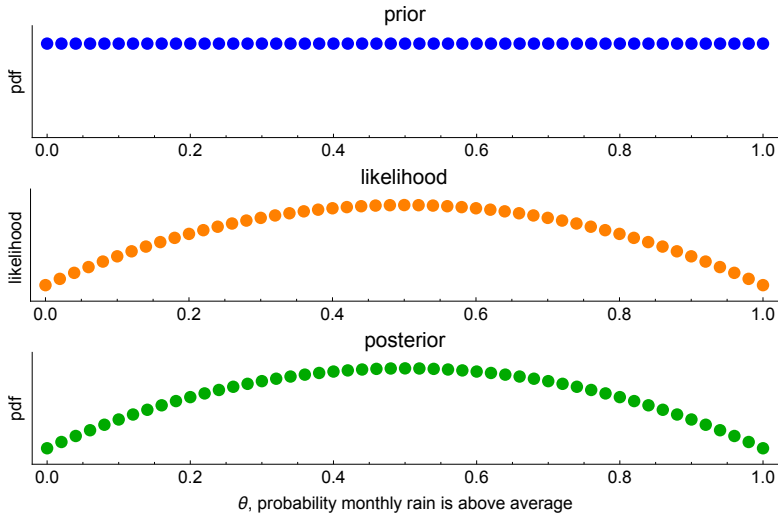
Scenario: discretised Bayesian inference

Discretise θ at intervals of 0.2.



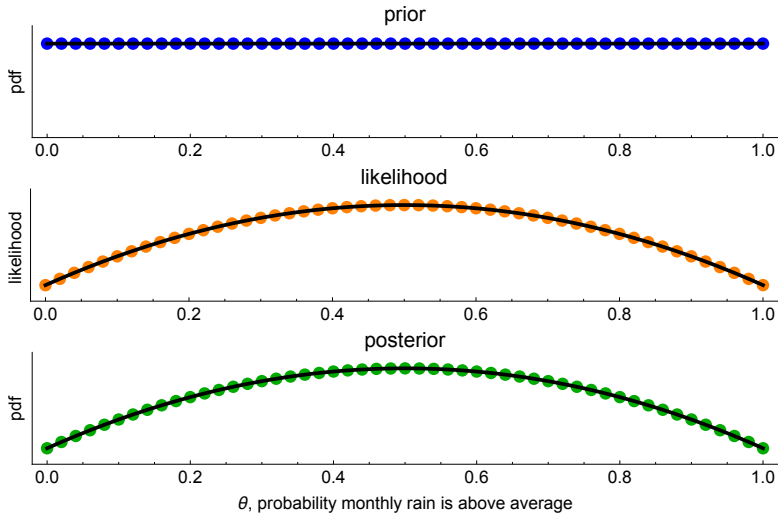
Scenario: discretised Bayesian inference

Discretise θ at intervals of 0.02.



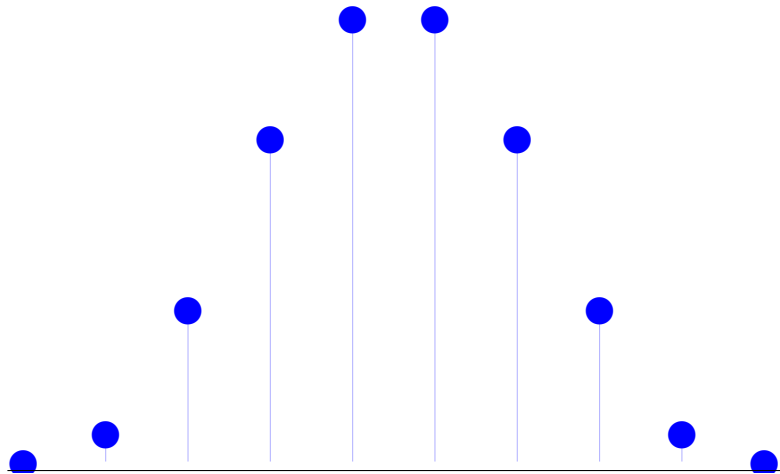
Scenario: discretised Bayesian inference

Discretise θ at intervals of 0.02.



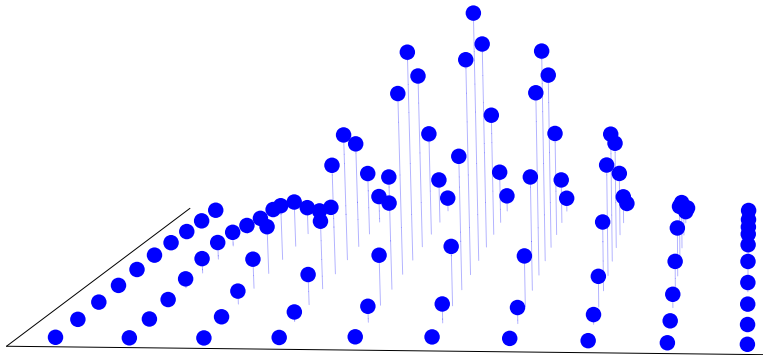
The problem with discretised Bayes

1 parameter \rightarrow 10 points



The problem with discretised Bayes

2 parameters $\rightarrow 10^2$ points



The problem with discretised Bayes and numerical quadrature

Question: how many grid points do we need for a 20-parameter model?

Answer: $10^{20} = 100,000,000,000,000,000,000,000$ grid points \therefore impossible!

Same goes for other methods that makes Bayesian inference discrete, for example **numerical quadrature**.



The problem of aforementioned methods: summary

- Bayesian inference requires us to difficult integrals; both for the denominator and posterior summaries.
- Conjugate priors are too simple for most real life examples.
- Another method is to approximate integrals by discretising them into sums.
- Method works ok for models with a few parameters.
- **But** doesn't scale well for models with more than about 3 parameters (curse of dimensionality).
- **Question:** can we find a method whose complexity is independent of the # of parameters?

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Black box die

- Black box containing a die with an **unknown** number of faces, and **weightings** towards sides.
- Shake the box and view the number that lands face up through a viewing window.
- Note: an individual shake represents one **sample** from the probability distribution of the die.



Black box die: estimating mean

- Question: How can we estimate the die's mean?
- Answer: shake it off! Then calculate the overall mean across all shakes.



Computational die in a box: results

Black box die: sampling to estimate a sum

- Mean of a **sample** of size n is:

$$\bar{X} = \frac{1}{n} \sum_{i=1}^n X_i \quad (17)$$

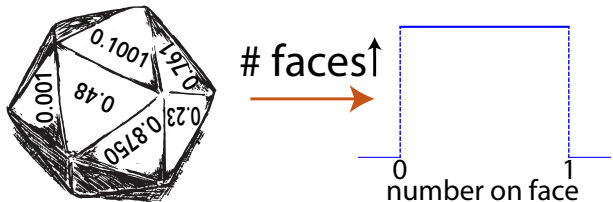
- Whereas the true mean of the die is given by:

$$E(X) = \sum_{j=1}^{\# \text{ faces}} Pr(X_j = x_j) \times x_j \quad (18)$$

- For a sample size of $< \sim 1000$ we were able to estimate:

$$\bar{X} \approx E(X) \quad (19)$$

An infinitely-sided die as a continuous distribution



- Imagine increasing the number of faces to infinity (a strange die indeed).
- Each face corresponds to one real number between 0 and 1.
- All possible numbers between 0 and 1 are covered.
- Basically like a **continuous uniform** distribution between 0 and 1.

An infinitely-sided die

- However its mean is now given by an **integral** rather than a **sum**.

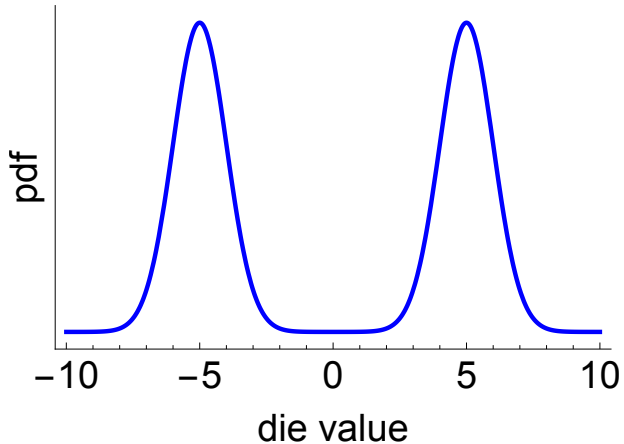
$$E(X) = \int_{\text{all faces}} p(X) \times X dX \quad (20)$$

- **Question:** can still estimate its true mean by the **sample** mean?
- If so this amounts to estimating the above integral!

Continuous distribution sampling

A stranger distribution

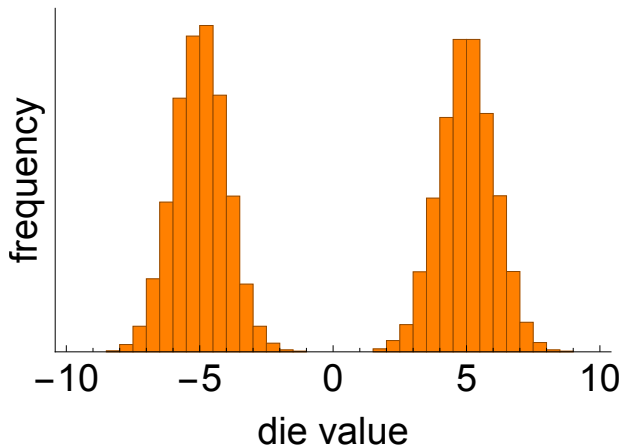
- Method seems to work for continuous uniform distribution.
- **Question:** does it work for other distributions?



A stranger distribution: sampling

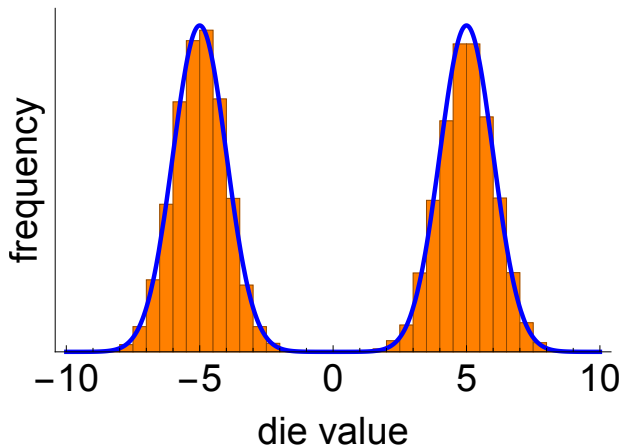
A stranger distribution: why does sampling work?

Compare samples...



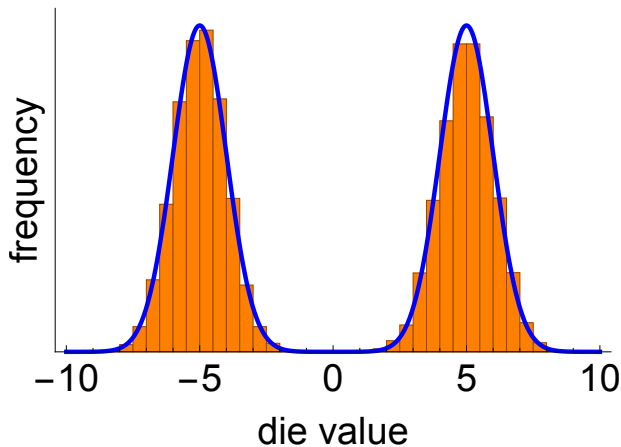
A stranger distribution: why does sampling work?

...with actual distribution \implies same shape!



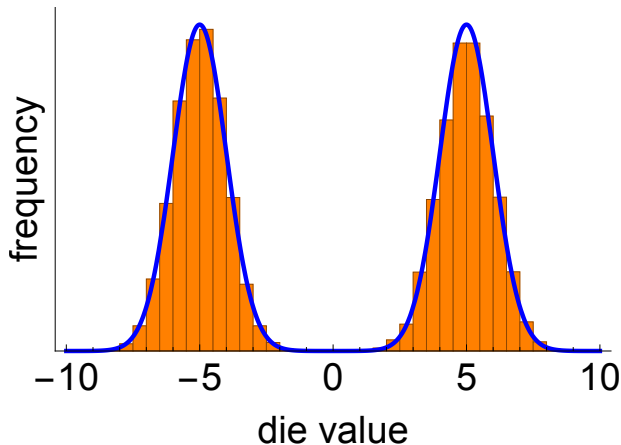
A stranger distribution: why does sampling work?

Therefore sample properties \rightarrow actual properties.



A stranger distribution: why does sampling work?

Note: nowhere have we explicitly mentioned the parameter dimension (complexity-free scaling?).



What is an independent sample?

- Aforementioned methods require us to generate **independent** samples from the distribution.
- **Question:** what *is* an independent sample?
- **Answer:** a value drawn from the distribution whose value is unconnected to other samples (apart from their joint reliance on the distribution.)

How to generate independent samples?

- By definition using independent sampling to estimate integrals requires us to be able to generate independent samples: $\theta_i \sim p(\theta)$.
- Not as simple as might first appear.
- Most statistical software has inbuilt ability to generate (pseudo-)independent samples for a few basic distributions: uniform, normal, poisson etc.
- However, for more complex distributions it is not trivial to create an independent sampler.

Summary

- Posterior is a weighted average of prior and likelihood, where weight of likelihood determined by amount of data.
- Posterior predictive distributions show implications of the posterior on the observable world.
- Exact Bayes is hard due to difficulty of calculating posterior, and other high dimensional integrals.
- Conjugate priors can make analysis simpler, although are highly restrictive.
- Discretisation can work for low dimensional problems but cannot cope with more complex models.
- Independent sampling can help to estimate integrals but can be hard to do in practice (see problem set).