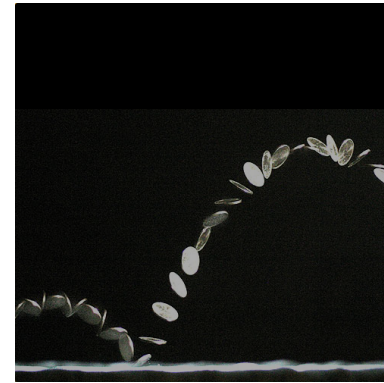
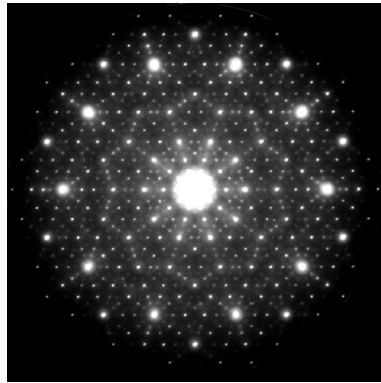


LTAT.02.004 MACHINE LEARNING II

## **Basics of probabilistic modelling**

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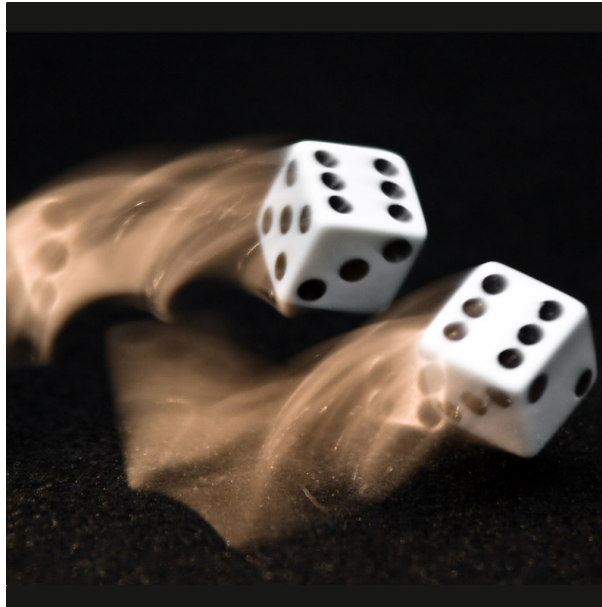
# What is probability?



Probability is a measure of uncertainty which can rise in several ways

- ▷ Intrinsic uncertainty in the system
- ▷ Uncertainty caused by inherent instability of the system
- ▷ Uncertainty caused by lack of knowledge or control over the system

# Frequentistic interpretation of probability



Probability is an average occurrence rate in long series of experiments.

- ▷ Law of large numbers
- ▷ Probability is a collective property
- ▷ Probabilities can be assigned only to future events

# Bayesian interpretation of probability



Probability reflects persons individual beliefs on future or unknown events.

- ▷ Belief updates through the Bayes rule
- ▷ Probability is an inherently subjective property
- ▷ Probabilities can be assigned to past, present and future events

# Ultra-frequentistic interpretation of probability



Events with small enough probability do not occur

- ▷ The main tool in classical statistics
- ▷ Errors in judgement does not matter if a gamma ray pulse kills us.
- ▷ One must avoid the lottery paradox in the reasoning

# The goal of statistical inference

## Frequentist goal

- ▷ The aim of statistics is to design algorithms that work well on average.
- ▷ For that one needs to specify probabilistic model for data sources.
- ▷ Confidence is the fraction of cases the algorithm works as specified.

## Bayesian goal

- ▷ The aim of statistics is to design algorithms that allow *rational individuals* to reliably update their beliefs through Bayes formula
- ▷ Besides the data source model one has to provide model for initial beliefs.
- ▷ Correctness of an algorithm does not make sense.

# Frequentistic methods

## Causation between zero-one events

Assume that condition  $A$  causes the event  $B = 1$  with probability  $p$ , i.e.,

$$\Pr[B = 1|A] = p$$

Then the probability is to get  $k$  ones in  $n$  independent trials is

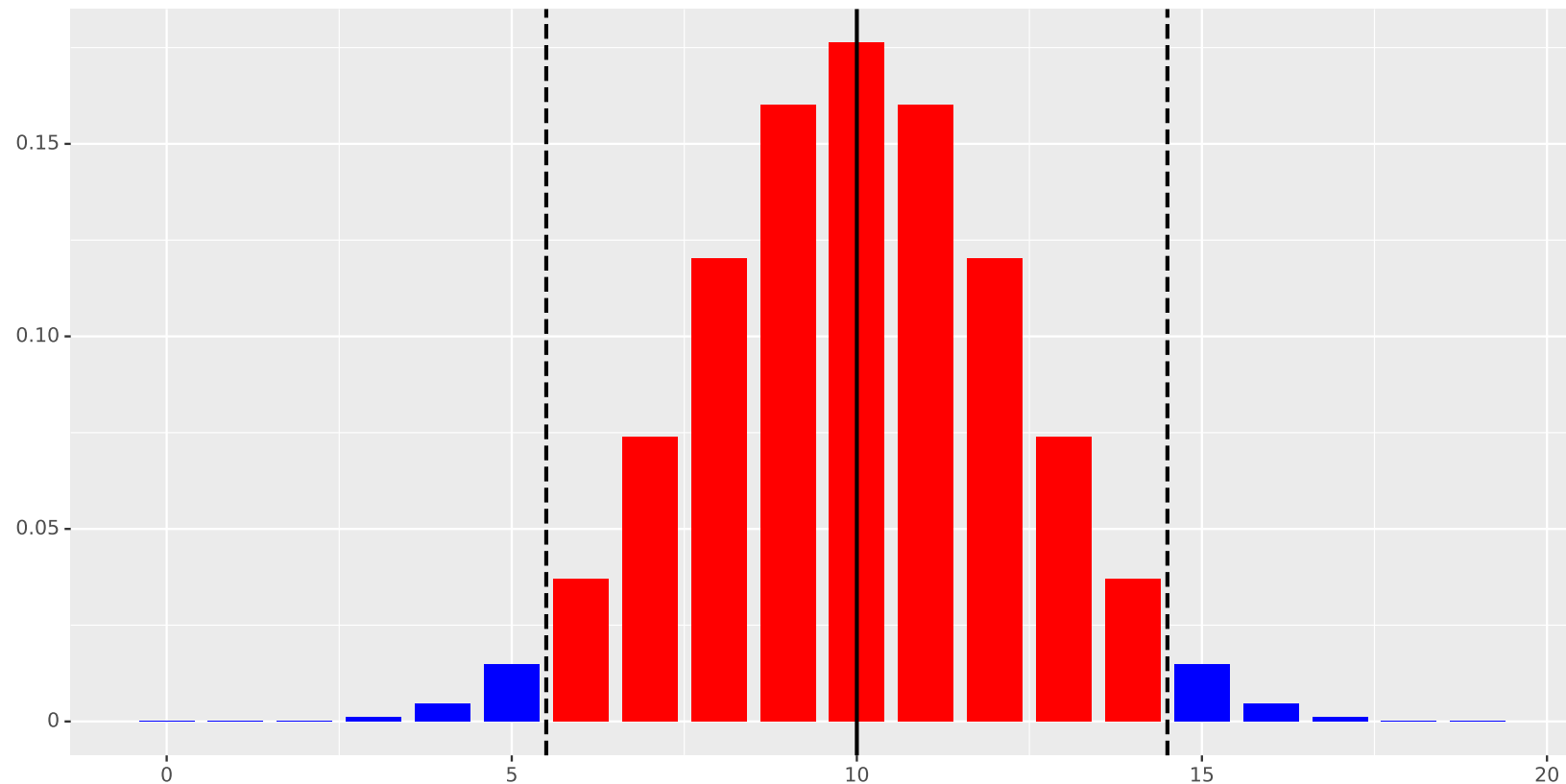
$$\Pr[B_1 + \dots + B_n = k|A] = \binom{n}{k} p^k (1 - p)^{n-k}$$

The number of ones is known to have a *binomial distribution*

$$B_1 + \dots + B_n \sim \text{Bin}(n, p)$$



## Illustration



The distribution of  $B_1 + \dots + B_n$  depends solely on the number of trials  $n$  and the probability  $p$ . Some values of  $B_1 + \dots + B_n$  are very unlikely.

# How to build a statistical test

## I. Null hypothesis:

- ▷ The probability of heads in a coinflip is  $\Pr[B_i = 1] = p$ .

## II. Choose value to compute aka test statistic:

- ▷ Our test statistic will be  $B_1 + \dots + B_n$ .

## III. Consequences on the observations:

- ▷ The observed sum  $B_1 + \dots + B_n \sim \text{Bin}(n = 20, p = 0.5)$ .
- ▷ Limit on the tail probability  $\Pr[|B_1 + \dots + B_n - 10| \geq 6] \leq 5\%$

## IV. Test procedure

- ▷ Reject null hypothesis at *significance level* 5% if  $|B_1 + \dots + B_n - 10| \geq 6$ .

## Properties of statistical tests

Statistical test is a classification algorithm designed to distinguish a fixed distribution of negative examples specified by a null hypothesis.

Any *fixed* classification *rule* can be converted to a statistical test by finding out the percentage of false positives aka *p-value*:

- ▷ There might exist a closed form solution.
- ▷ We can always estimate p-values using simulations.
- ▷ Observations must be compressed into a single decision value.

Testing several hypothesis in parallel increases the number of false positives. Several p-value adjustment methods are used to correct the issue:

- ▷ Bonferroni correction is almost optimal
- ▷ FDR correction controls the expected number false positives

# How to build confidence intervals

## I. Construct a family of statistical tests:

- ▷ Define a statistical test  $T_p$  for all possible parameter values  $p$ .
- ▷ All tests should share the same test statistic.

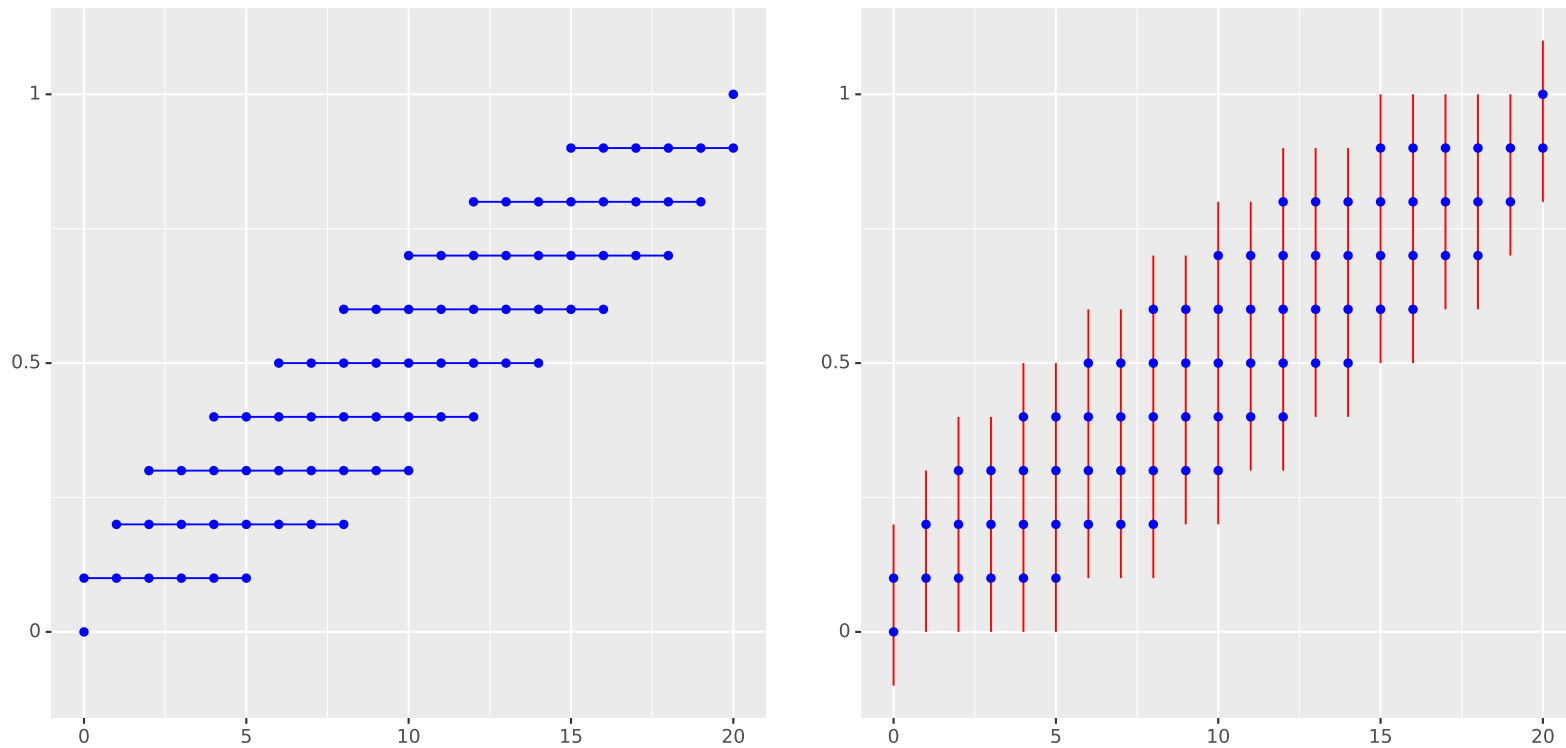
## II. Perform multiple hypothesis testing for all parameter values:

- ▷ Accept all parameters values for which p-value is greater than  $1 - \alpha$ .
- ▷ Output a minimal interval that covers all accepted parameter values.

## Rationale

- ▷ The true parameter value is rejected on  $\alpha$ -fraction of possible observations.
- ▷ For the remaining cases the true value is inside the predicted interval.

# Illustration



- ▷ Acceptance ranges for different parameter values on the left.
- ▷ Extended parameter ranges covering all accepted parameters on the right.
- ▷ These ranges are the desired confidence intervals.

## Interpretation of confidence intervals

**Definition.** Confidence interval for a parameter  $p$  is an outcome of an approximation algorithm. The algorithm must output an interval  $[\hat{p} - \varepsilon, \hat{p} + \varepsilon]$  such that the true estimate is in the range on  $\alpha$ -fraction of cases.

### Paradoxical inapplicability

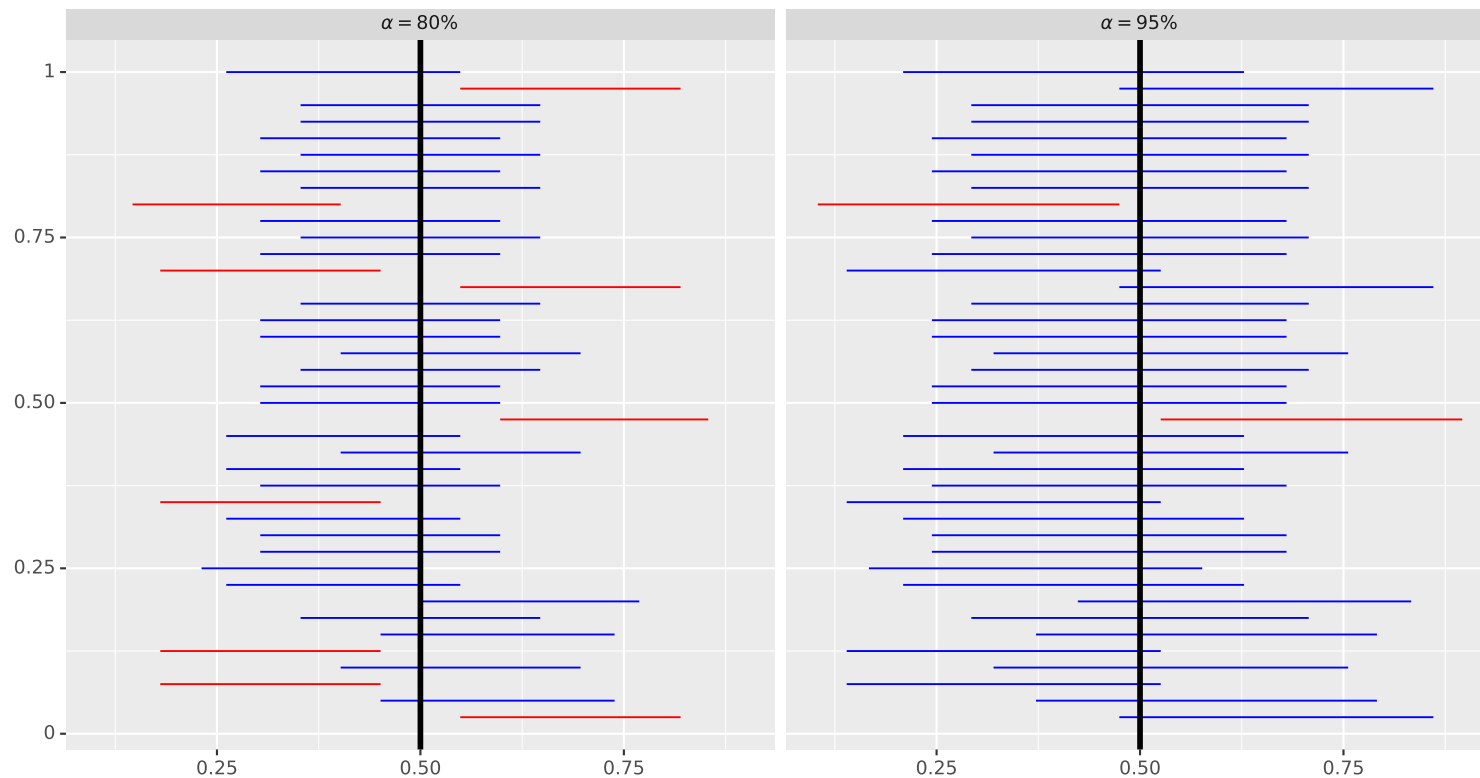
The definition does not state that the probability  $p \in [\hat{p} - \varepsilon, \hat{p} + \varepsilon]$  is  $\alpha$ !

- ▷ The statement  $p \in [\hat{p} - \varepsilon, \hat{p} + \varepsilon]$  is either true or false.
- ▷ There is no probability left. We just *do not know* the answer!

### Ultra-frequentistic resolution

- ▷ If  $1 - \alpha$  is small enough say 5% then the algorithm is always correct.

## Illustrative example



By increasing the length of the interval we increase the fraction of runs for which the true value of  $p$  lies in the interval.

# Problems with confidence intervals

## Inability to capture background knowledge

- ▷ What if I know that  $p \in [0.1, 0.2]$  and observe  $B_1 = \dots = B_N = 1$ ?
- ▷ Then the estimate  $[\hat{p} - \varepsilon, \hat{p} + \varepsilon]$  is clearly wrong although on average this confidence interval is reasonable.

## Multiple hypothesis testing

- ▷ Using several confidence intervals in parallel increases the fraction of cases where some true estimate is out of the predicted range.
- ▷ We can use p-value adjustment methods are used to correct the issue.



## Prediction intervals

Even if we know the true relation  $y = f(x)$  we cannot predict the observation  $y_i = f(x_i) + \varepsilon_i$ , as the noise term  $\varepsilon_i$  is not known ahead.

- ▷ We cannot give upper and lower bounds for  $y_i$  which always hold.

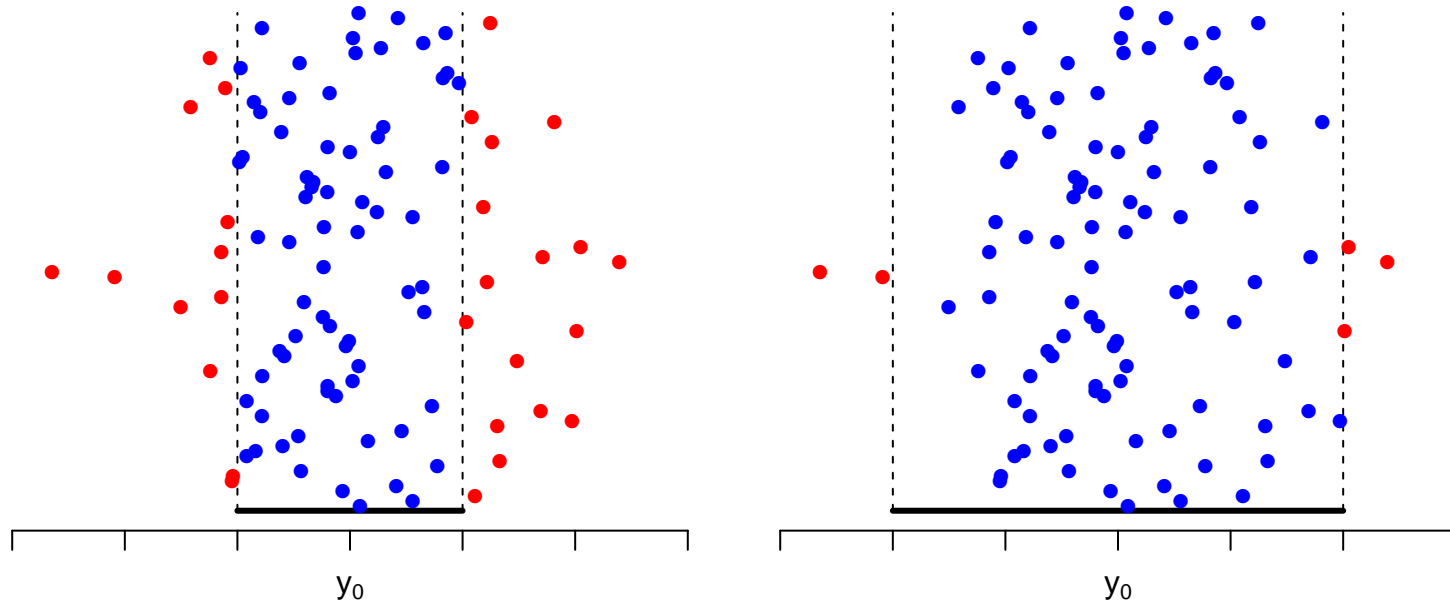
Instead, we can specify a prediction interval  $[y_* - \varepsilon, y_* + \varepsilon]$  so that with probability 95% the resulting measurement  $y_i$  is in the range.

- ▷ Usually, the analysis is similar to confidence interval derivation.

Interpretation of prediction intervals is different from confidence intervals.

- ▷ The probability estimate holds for the particular interval.

## Illustrative example



By increasing the length of the prediction interval we increase the fraction of future measurements which fall into interval.

# Confidence envelopes

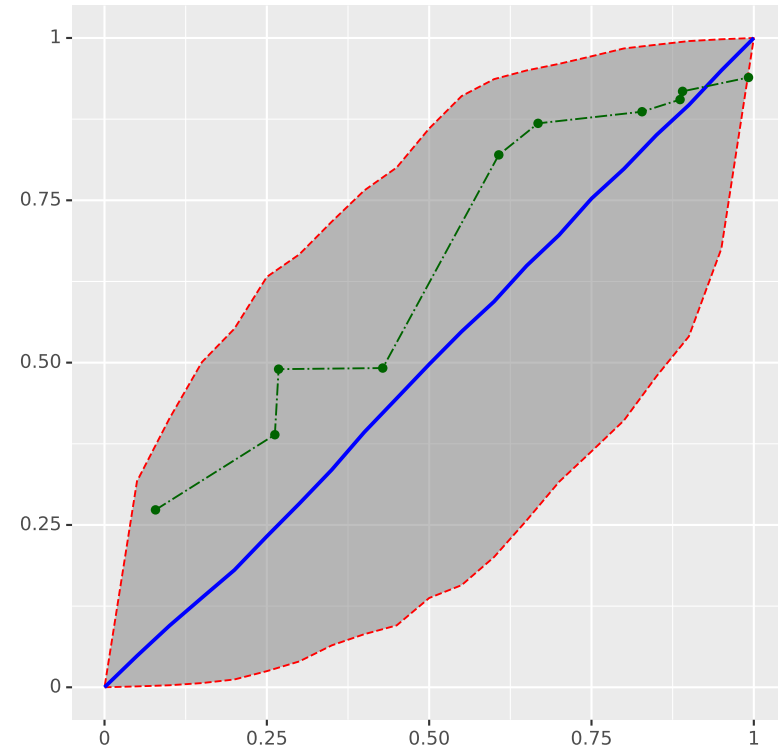
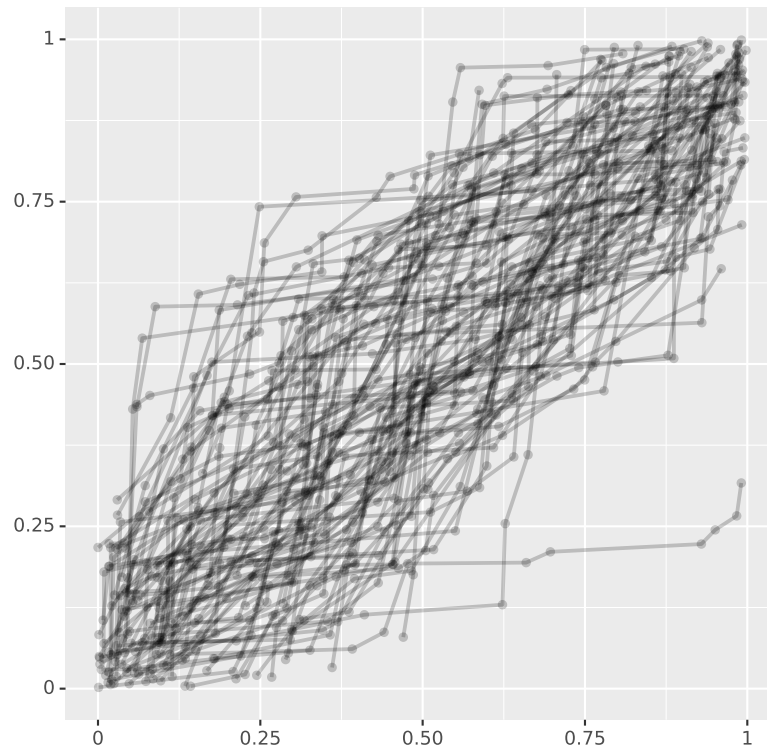
Confidence intervals is a good way to visualise uncertainty of a particular parameter. However, we are sometimes interested in the uncertainty many parameters or in the uncertainty of a function:

- ▷ How a predictor  $f : [0, 1] \rightarrow \mathbb{R}$  depends on the training set
- ▷ How a ROC curve  $\text{ROC} : [0, 1] \rightarrow [0, 1]$  depends on the test set
- ▷ How should a quantile-quantile plot be distributed.

Confidence bands are generalisations of confidence intervals

- ▷ Pointwise confidence band is a collection of confidence intervals
- ▷ Simultaneous confidence band must enclose  $\alpha$ -fraction of functions.
- ▷ Simultaneous confidence bands are much wider than pointwise bands.

## Illustrative example



- ▷ Distribution of qq-lines visualised through a sample on the left.
- ▷ A simulation based pointwise 95% confidence envelope on the right.
- ▷ The significance level that qq-line is inside the envelope is ca 50%.

# Permutation tests

## Baseline problem:

- ▷ Achievable accuracy depends on the data distribution.
- ▷ Artefacts in the dataset may bias performance measures.

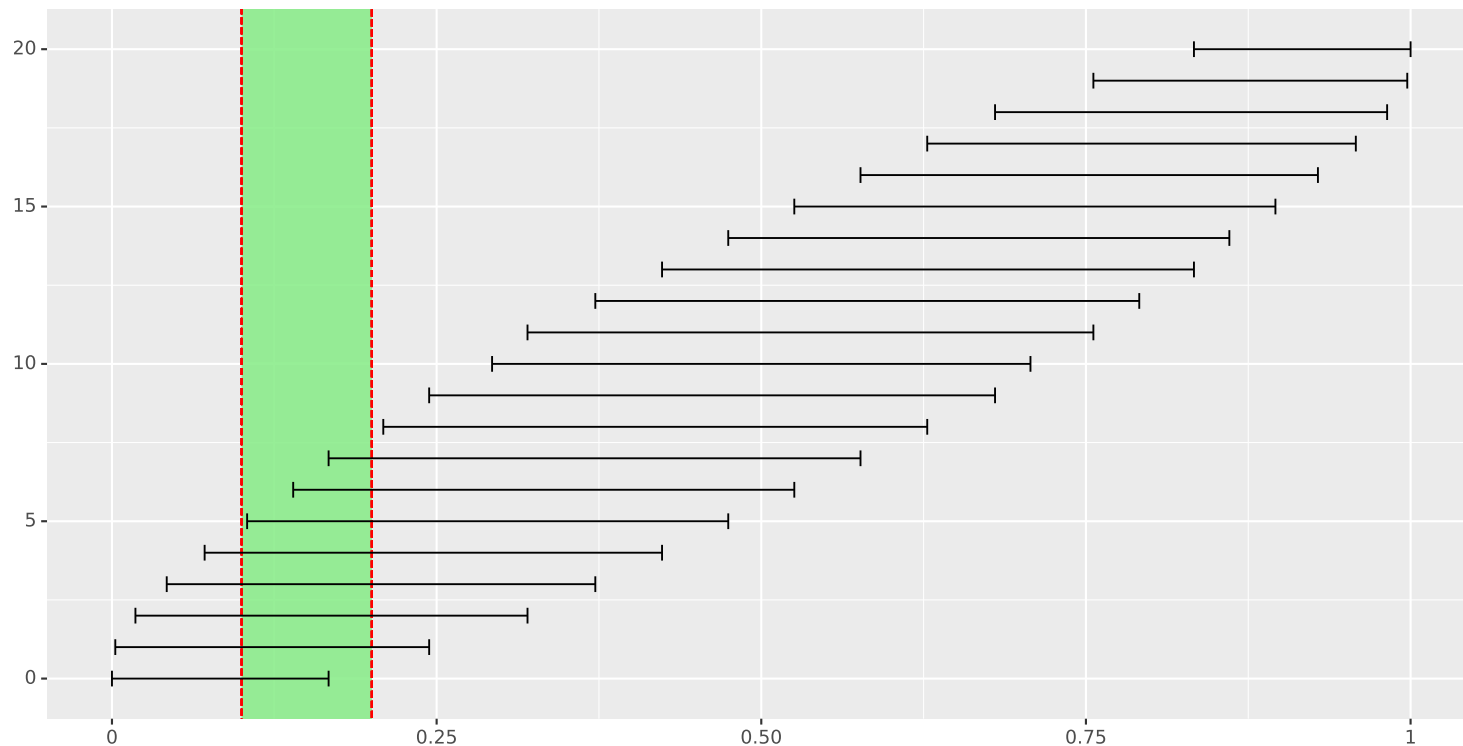
**Label permutation.** A random permutation  $\pi$  on outputs  $y_i$  destroys correlations between input-output pairs  $(x_i, y_{\pi(i)})$  but preserves marginal distribution of inputs and outputs.

**Permutation test.** Estimate how probable is to achieve equal or higher accuracy than was observed on the real data.

- ▷ If this probability is small then there must be signal in the data.
- ▷ The test completely neglect the effect size, i.e., how much results differ.
- ▷ Statistical significance does not imply utility!

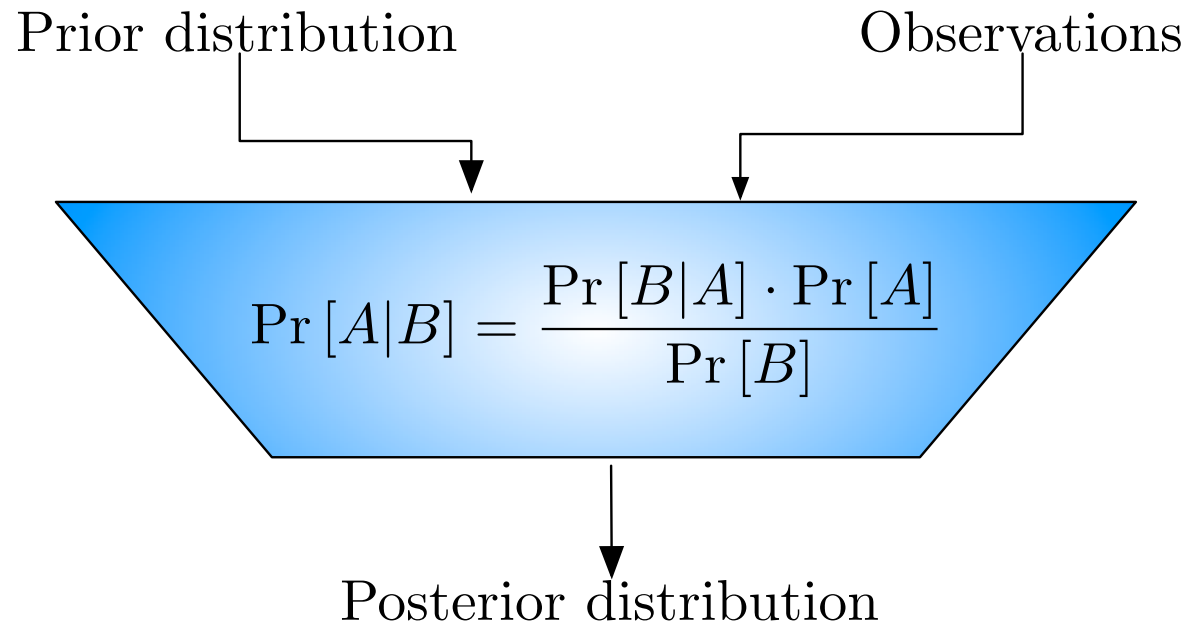
# Bayesian methods

# Confidence intervals vs background knowledge



- ▷ Confidence intervals do not capture background knowledge  $p \in [0.1, 0.2]$ .
- ▷ Thus we must accept absurd or suboptimal parameter estimations.

# Bayesian inference procedure



- ▷ Prior distribution  $\Pr[A]$  encodes the background knowledge
- ▷ The model  $\Pr[B|A]$  determines how the posterior  $\Pr[A|B]$  is updated



## Prior and likelihood

Likelihood  $\mathcal{L}(\mathcal{D}|\mathcal{M})$  is a probability of observations  $\mathcal{D}$  when the data generation model  $\mathcal{M}$  is fixed. The model is fixed by the set of parameters.

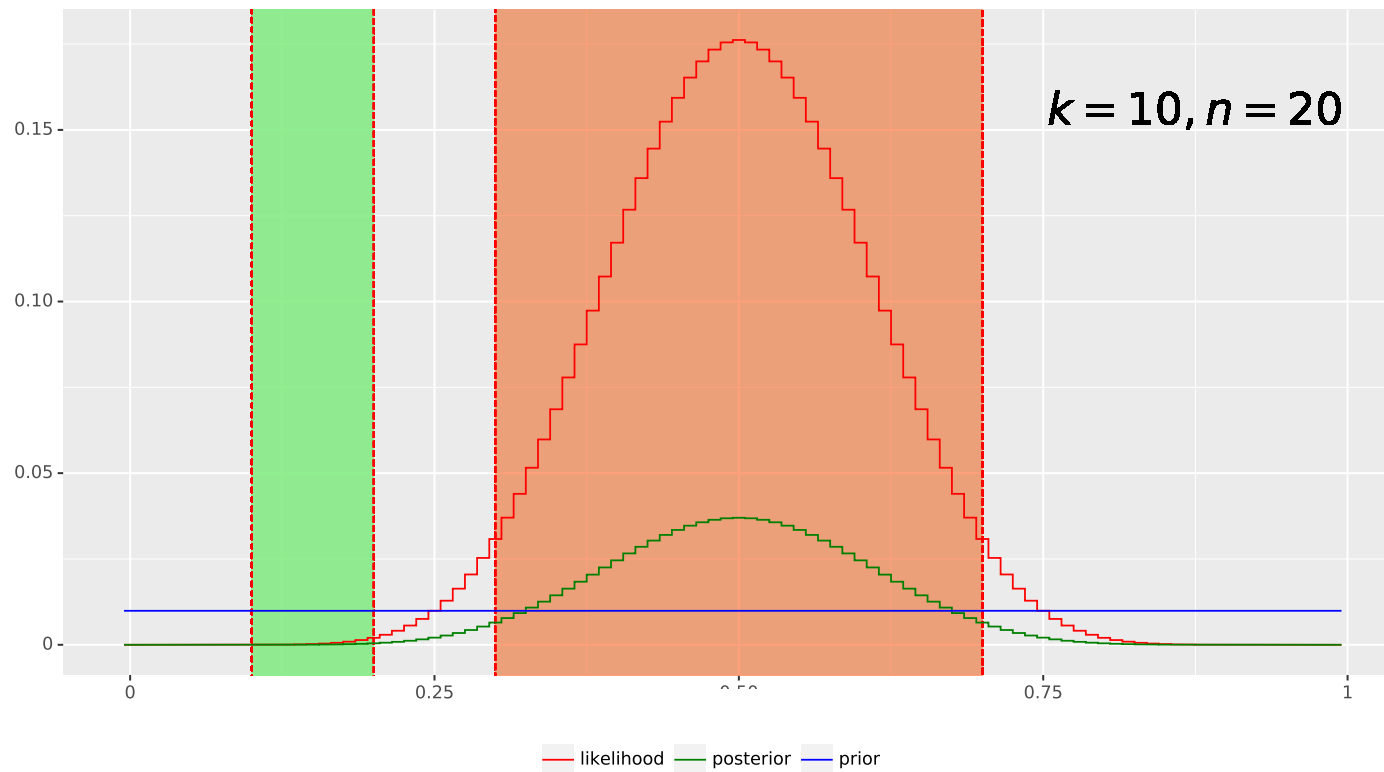
For coin flipping experiment the number of ones  $k$  is the observation and the coin bias  $p$  is the model parameter and thus

$$\mathcal{L}[k|p] = \binom{n}{k} p^k (1 - p)^{n-k}$$

Prior is a distribution over models that encodes our preferences of models before we observe any data.

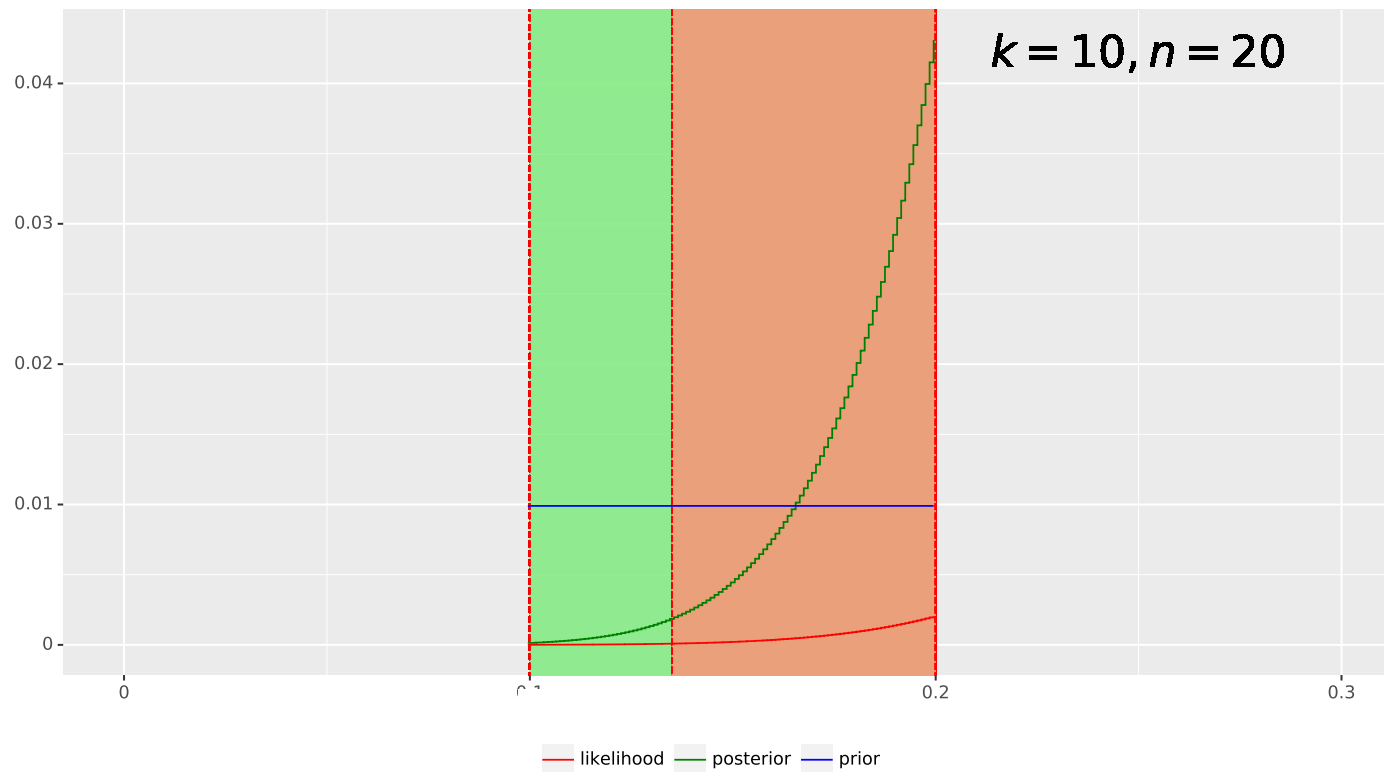
- ▷ Uninformative prior assigns uniform probability to all models.
- ▷ Uninformative prior is not well-defined for continuous parameters.

## Posterior of an uninformed person



- ▷ With no preferences the posterior is concentrated around 0.5.
- ▷ Credibility interval  $p \in [0.3, 0.7]$  contains 95% of posterior probability.

## Posterior of an informed person



- ▷ With preferences the posterior is concentrated to the left of 0.2.
- ▷ Credibility interval  $p \in [0.135, 0.2]$  contains 95% of posterior probability.

## Beta distribution as a posterior

By increasing the number of grid points in the non-informative prior we reach a continuous distribution with a density function

$$p[p|k] = \frac{\Gamma(n+2)}{\Gamma(k+1)\Gamma(n-k+1)} \cdot p^k(1-p)^{n-k} .$$

This distribution is known as *beta distribution*  $\text{Beta}(\alpha = k+1, \beta = n-k+1)$ . The parameter value that maximises the posterior is

$$p_* = \frac{\alpha - 1}{\beta - \alpha} = \frac{k}{n} .$$

## Dice throwing vs coin flipping

A behaviour of a dice with faces  $\{1, \dots, m\}$  is determined by probabilities

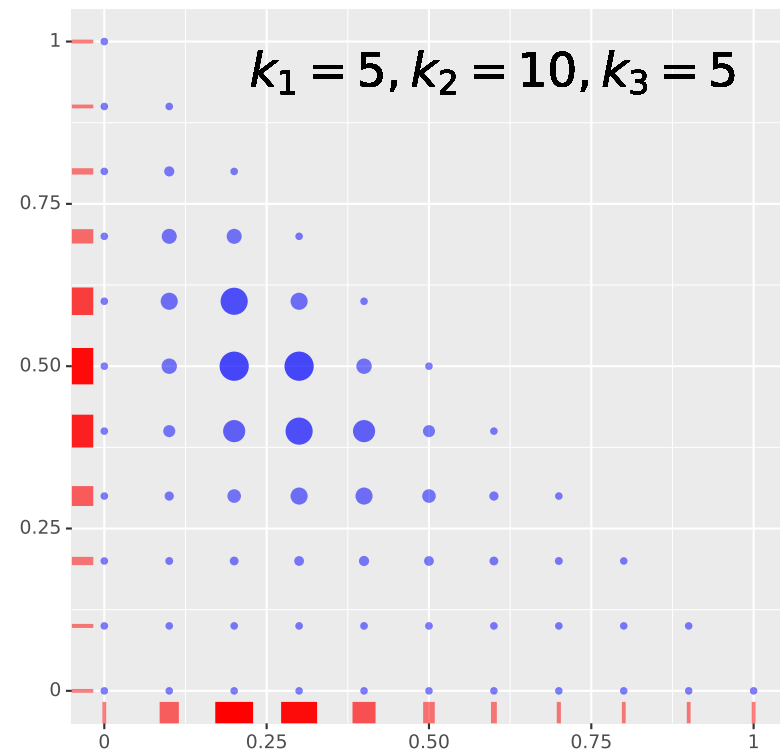
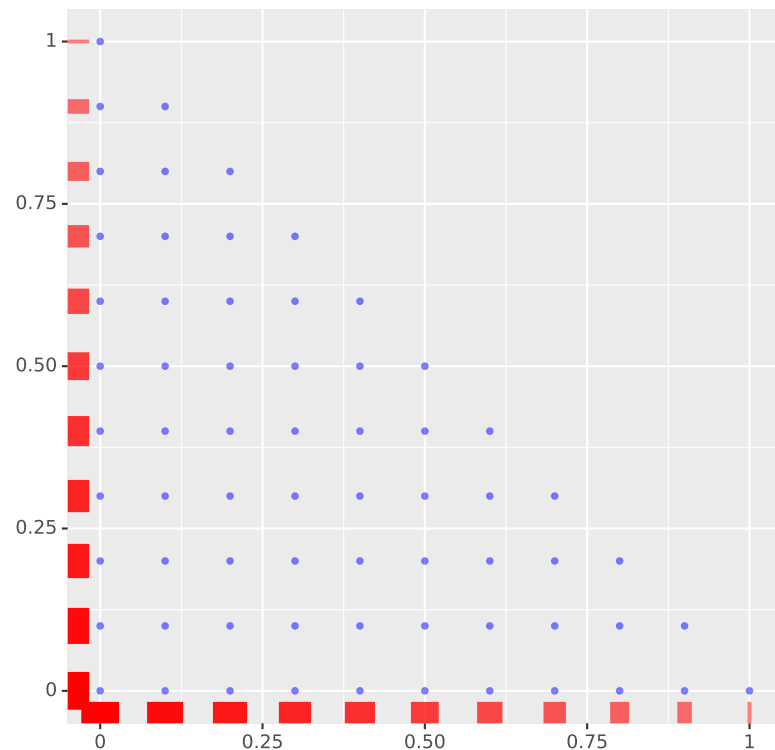
$$p_1 = \Pr[D_i = 1], \quad \dots, \quad p_m = \Pr[D_i = m]$$

### Reduction to coin flipping

- ▷ Let  $B_i$  denote the event that  $D_i = j$ .
- ▷ Then  $B_1, \dots, B_n$  is a coinflipping sequence with bias  $\Pr[B_i = 1] = p_j$ .
- ▷ Non-informative prior for dice throwing goes to the non-informative prior.
- ▷ Informative priors can be marginalised to the right format.
- ▷ The same reduction can be done for all faces of the dice.

**Caution:** Marginal posteriors do not determine the full posterior in general.

# Illustration



- ▷ Uniform prior over parameter pairs yields non-uniform marginal priors.
- ▷ The joint MAP estimate coincides with the marginal MAP estimates.

## Dirichlet distribution as a posterior

By increasing the number of grid points in the non-informative prior over simplex we reach a continuous distribution with a density function

$$p[p_1, \dots, p_m | k_1, \dots, k_m] = \frac{\Gamma(n + m)}{\Gamma(k_1 + 1) \cdots \Gamma(k_m + 1)} \cdot p_1^{k_1} \cdots p_m^{k_m} .$$

This distribution is known as *Dirichlet distribution*

$$\text{Dirichlet}(\alpha_1 = k_1 + 1, \dots, \alpha_m = k_m + 1) .$$

The parameter value that maximises the posterior is

$$p_i^* = \frac{\alpha_i - 1}{\alpha_1 + \dots + \alpha_m - m} = \frac{k_i}{n} .$$

## Laplace smoothing

Assume that we throw a dice with  $m$  faces and  $B_i$  encodes the event that the dice lands on a specific face. Then it is natural to assign the maximum prior probability to the parameter value  $p_* = \frac{1}{m}$ .

Such prior can be defined through a following though experiment:

- ▷ We start with non-informative prior.
- ▷ We observe all possible outcomes of the dice  $\alpha$  times.
- ▷ We use the resulting posterior as a prior for real observations.

Thus the posterior can be obtained by starting with non-informative prior and observing  $k + \alpha$  ones among  $n + m\alpha$  throws.

- ▷ The ratio  $p = \frac{k+\alpha}{n+m\alpha}$  is the maximal a posteriori estimate for  $p$ .