LTAT.02.004 Machine Learning II

Basics of probabilistic modelling

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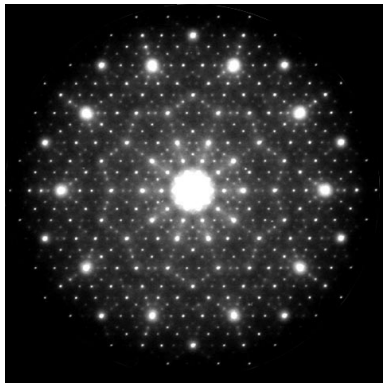
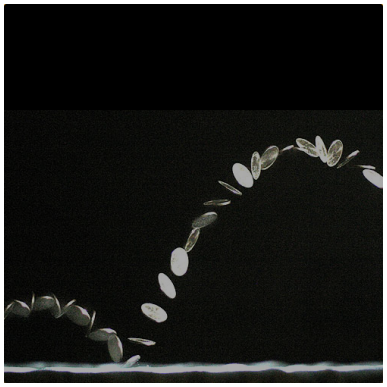
What is probability?





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.,

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

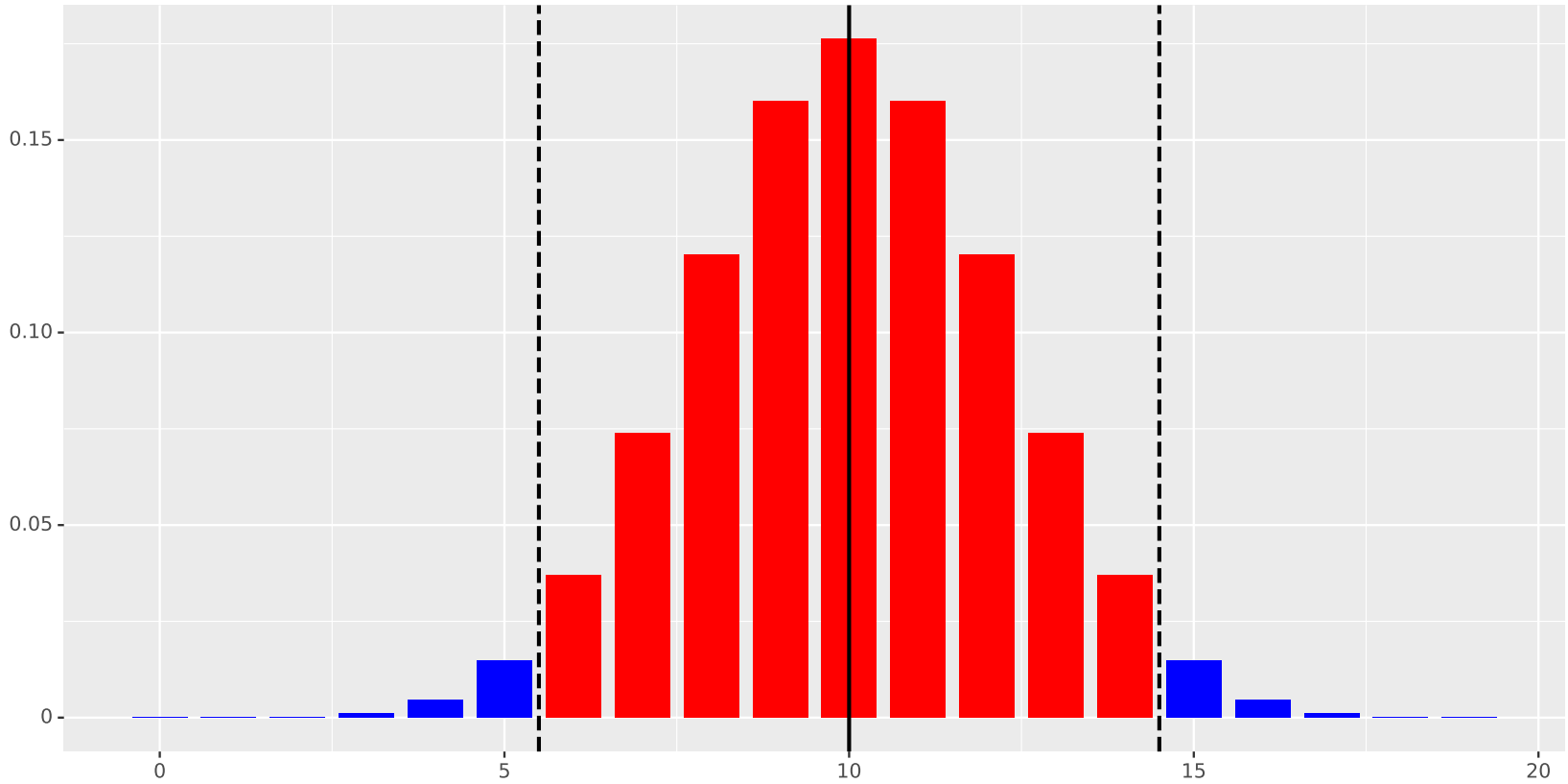
Then the probability is to get k ones in n independent trials is

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

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

$$B_1 + \cdots + B_n \sim \mathsf{Bin}(n, p)$$

Illustration



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

How to build a statistical test

I. Null hypothesis:

▷ The probability of a single one is $\Pr[B_i = 1] = p$.

II. Choose value to compute aka test statistic:

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

III. Consequences on the observations:

▷ The observed sum $B_1 + \ldots + B_n \sim \mathrm{Bin}(n = 20, p = 0.5)$.

▷ Limit on the tail probability $\Pr[|B_1 + \ldots + B_n - 10| \geq 6] \leq 5\%$

IV. Test procedure

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

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

- ▷ There might exists 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.

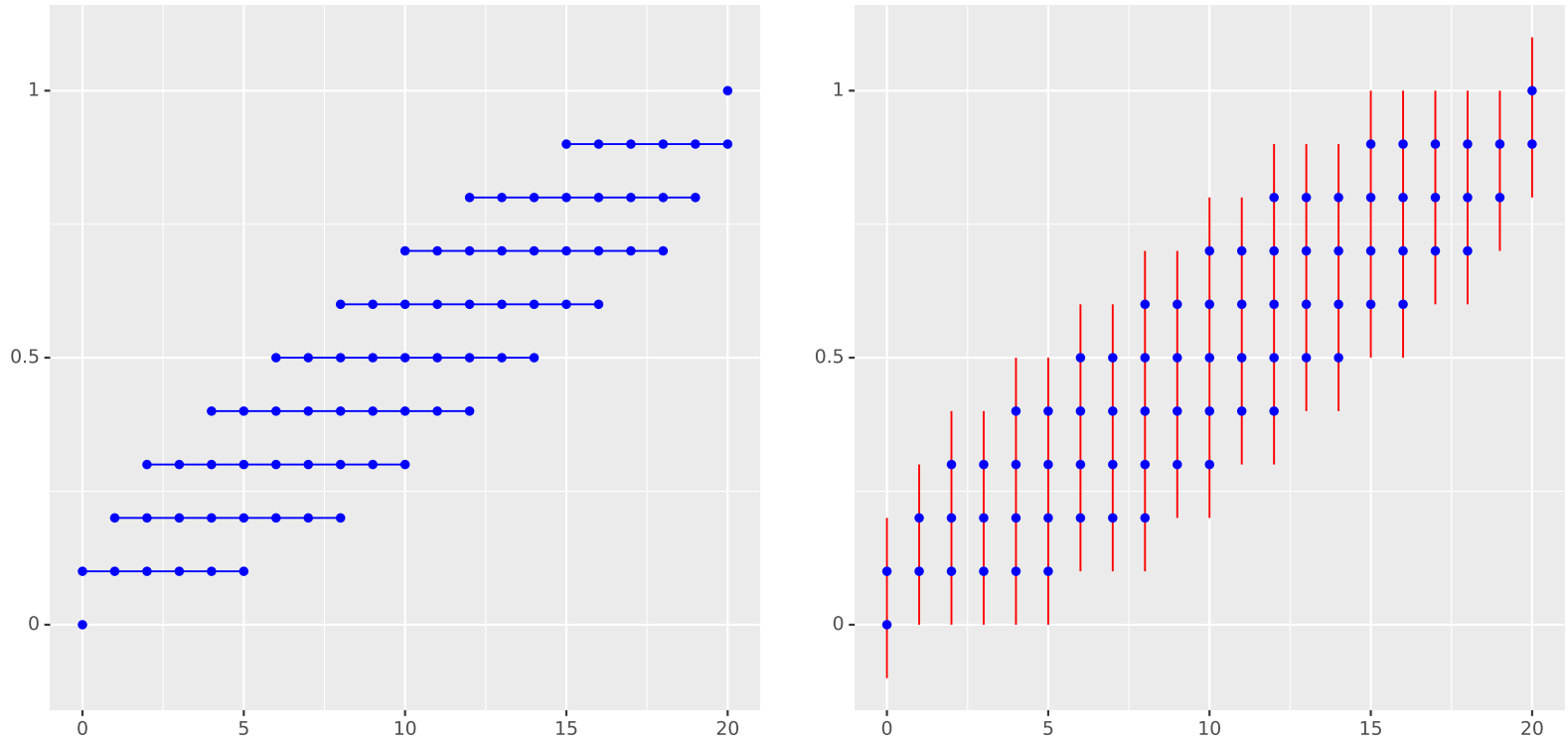
II. Perform multiple hypotesis testing for all parameter values:

- ▷ Accept all parameters values for which pvalue is greather than α.
- ▷ Output a minimal interval that covers all accepted parameter values.

Rationale

- ▷ The true parameter value is rejected on α-fraction of possible observations.
- ▷ Otherwise, 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 out of this range on α fraction of cases.

Paradoxical inapplicability

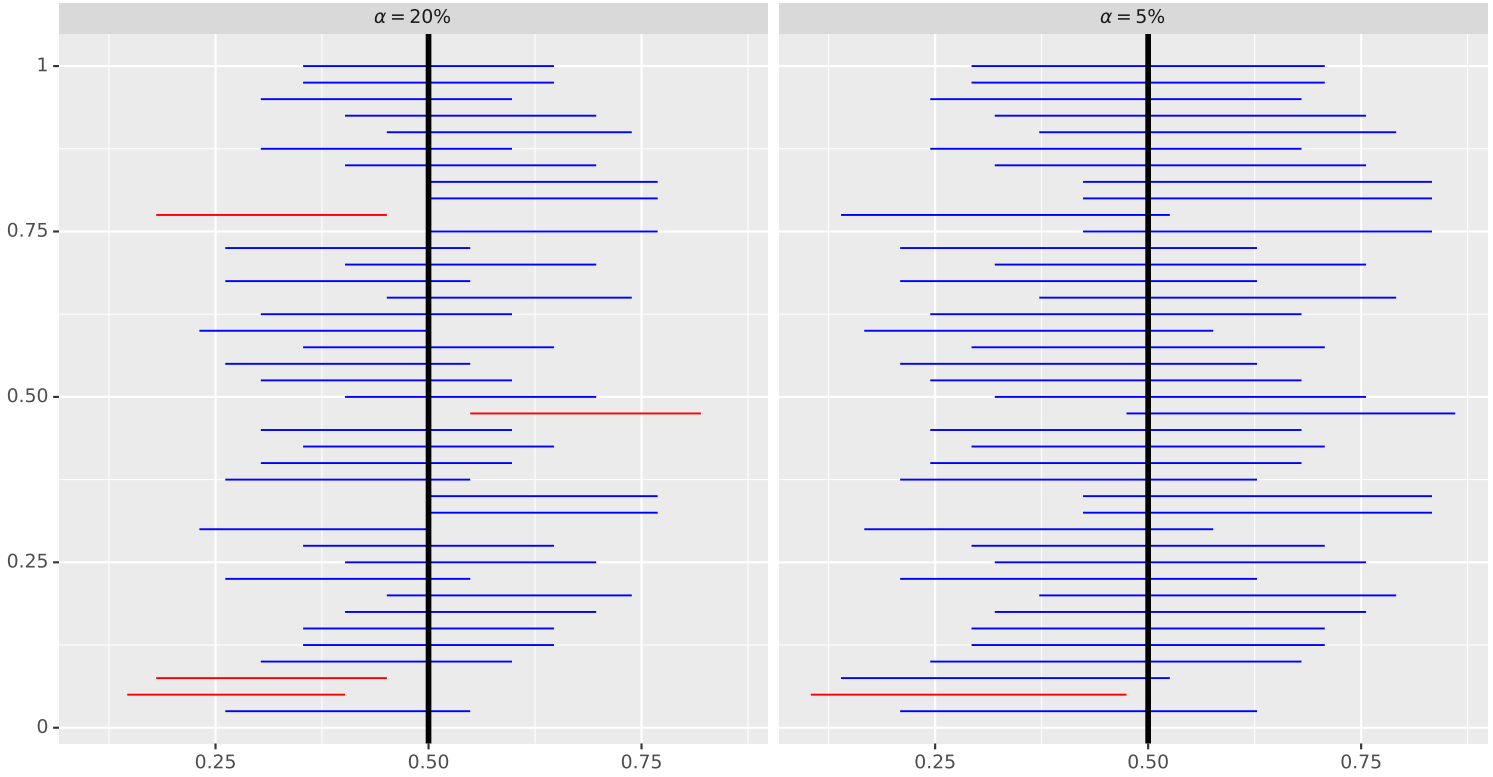
The definition does not state that the probability $p \in [\hat{p}-\varepsilon, \hat{p}+\varepsilon]$ is α!

- ▷ 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 α 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 $B_1 = \ldots = B_N = 1$?
- ▷ 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 $y = f(\boldsymbol{x})$ we cannot predict the observation $y_i = f(\boldsymbol{x}_i) + \varepsilon_i$, as the noise term ε_i is not known ahead.

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

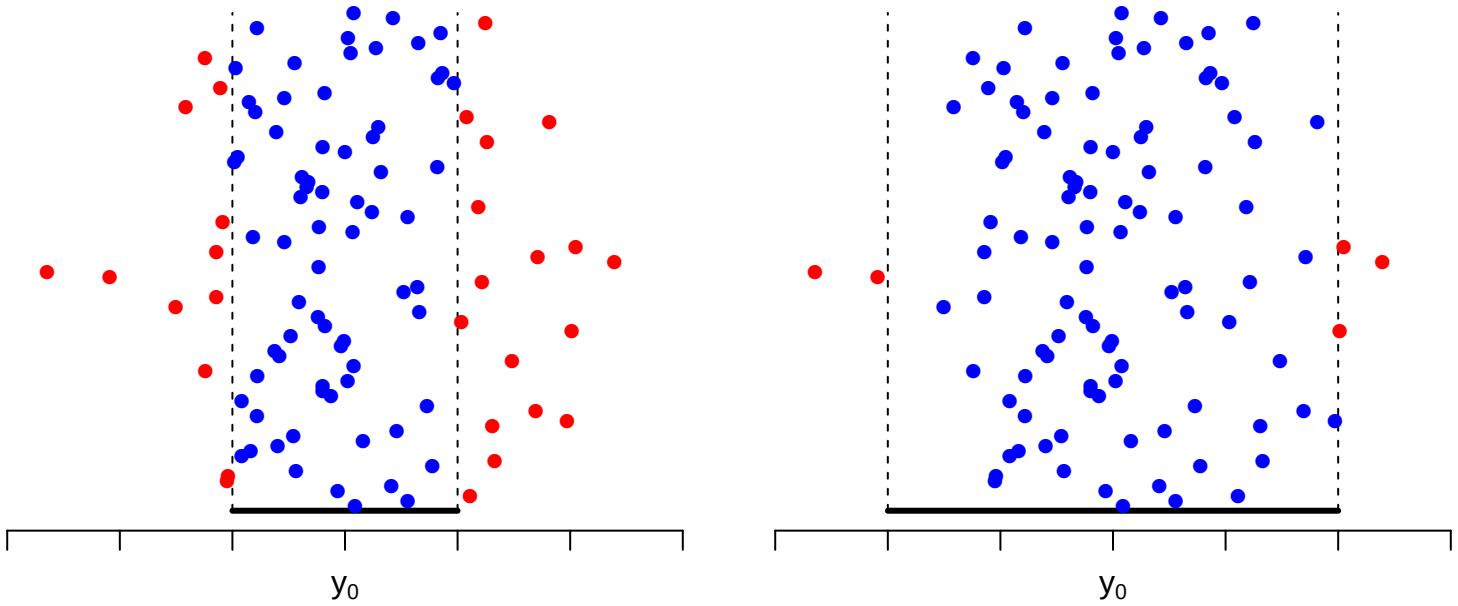
Instead, we can specify a prediction interval $[y_* - \varepsilon, y_* + \varepsilon]$ so that with probability 95% the resulting measurement y_i is in the range.

▷ 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

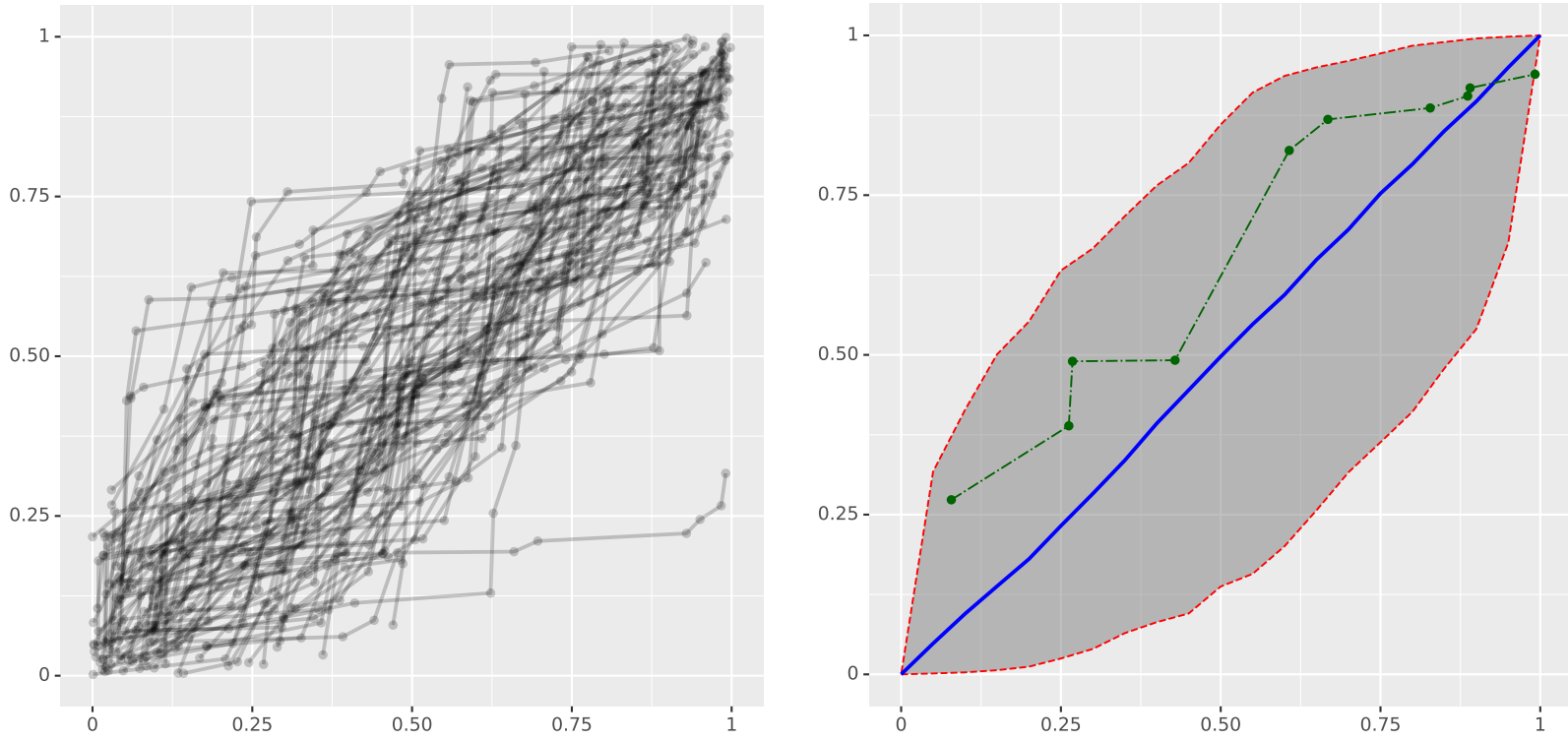
Confidence intervals is a good way to visualise uncertainty of a particular paramater. However, we are sometimes interested in the uncertainty many parameters or in the uncertainty of a function:

- ▷ How a predictor $f : [0, 1] \to \mathbb{R}$ depends on the training set
- ▷ How a ROC curve $\mathrm{ROC} : [0, 1] \to [0, 1]$ 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 that encloses α-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.

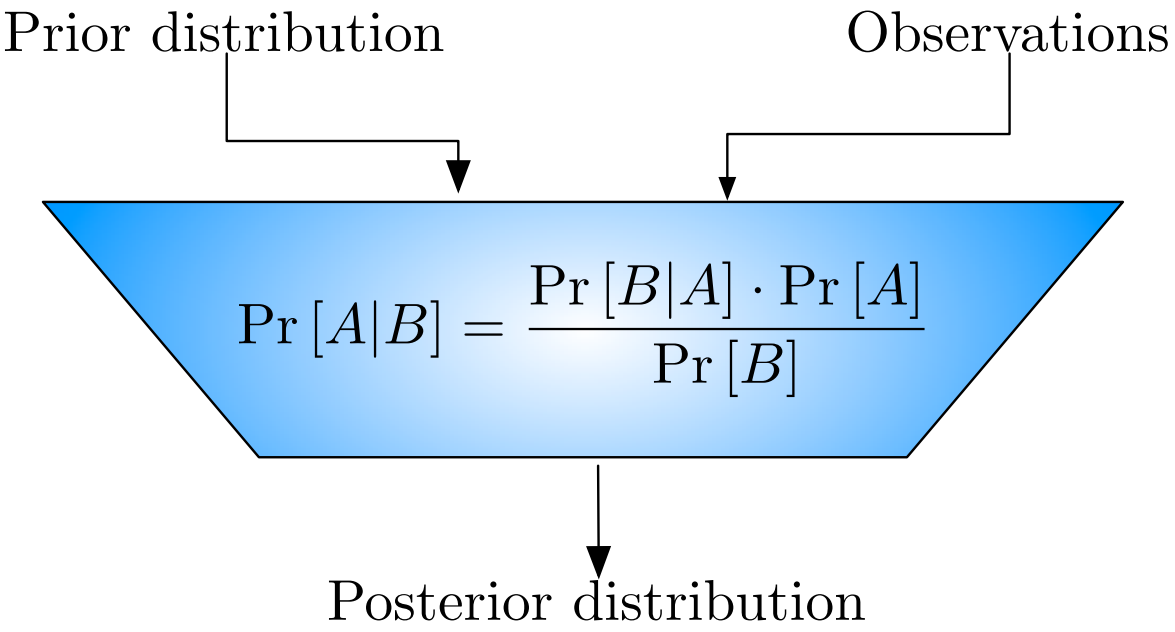
Label permutation. A random permutation π of outputs y_i destroys correlations between input-output pairs $(\boldsymbol{x}_i, \boldsymbol{y}_{\pi(i)})$ but preserves marginal distribution of inputs and outputs.

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!

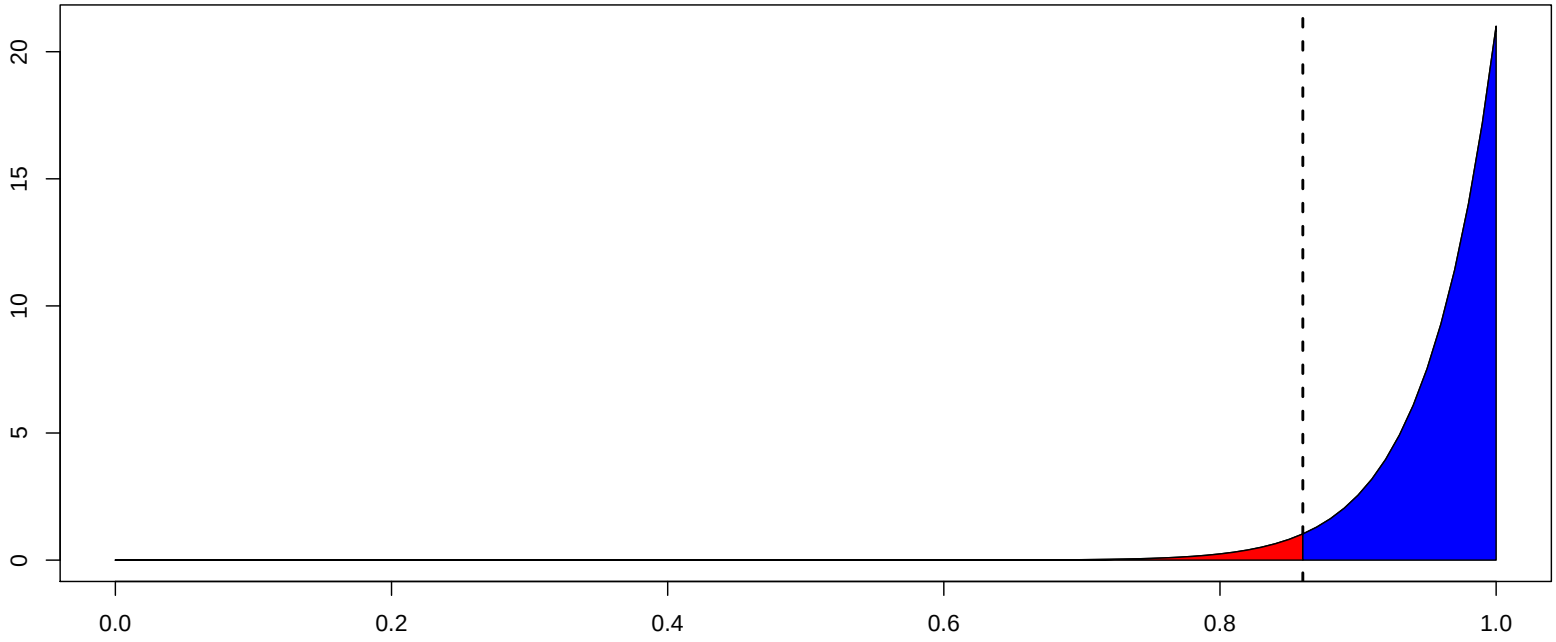
Bayesian methods

Bayesian inference procedure



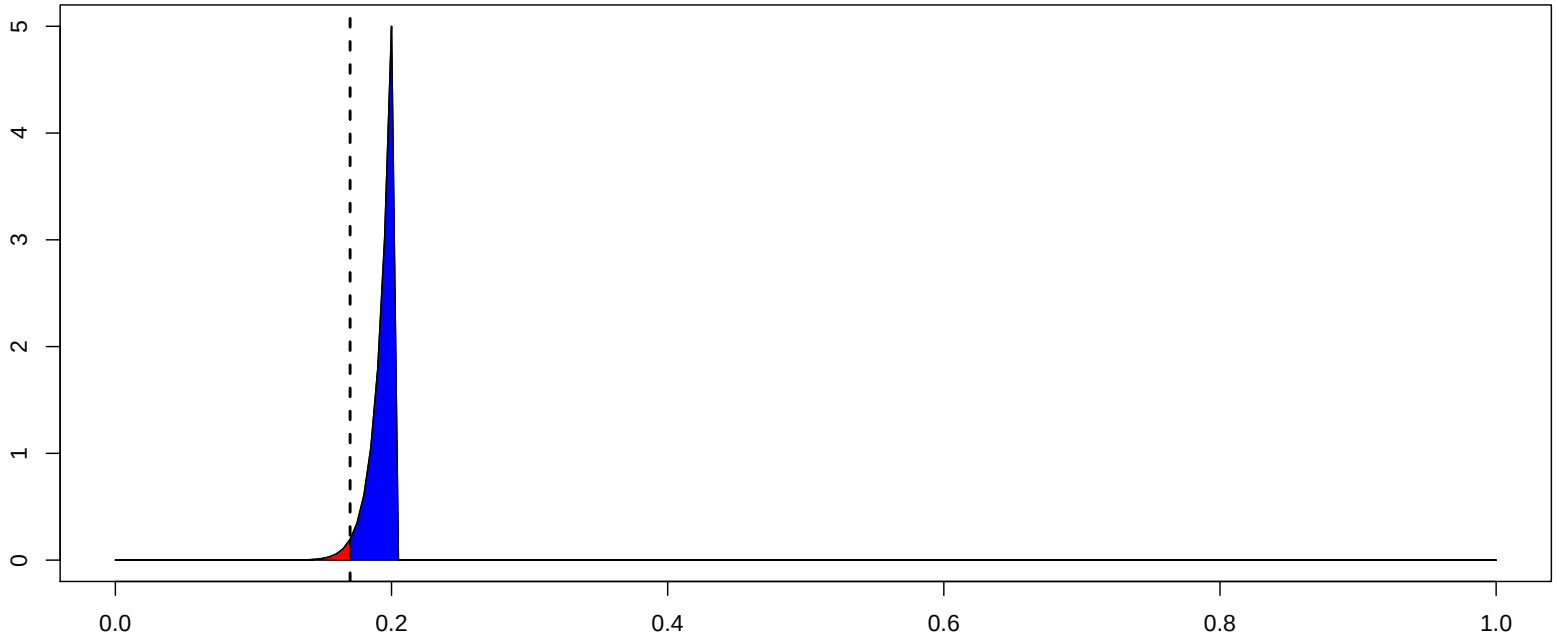
- ▷ Prior distribution $\Pr[A]$ encodes the background knowledge
- ▷ The model $\Pr[B|A]$ determines how the posterior $\Pr[A|B]$ is updated

Posterior of an uninformed person



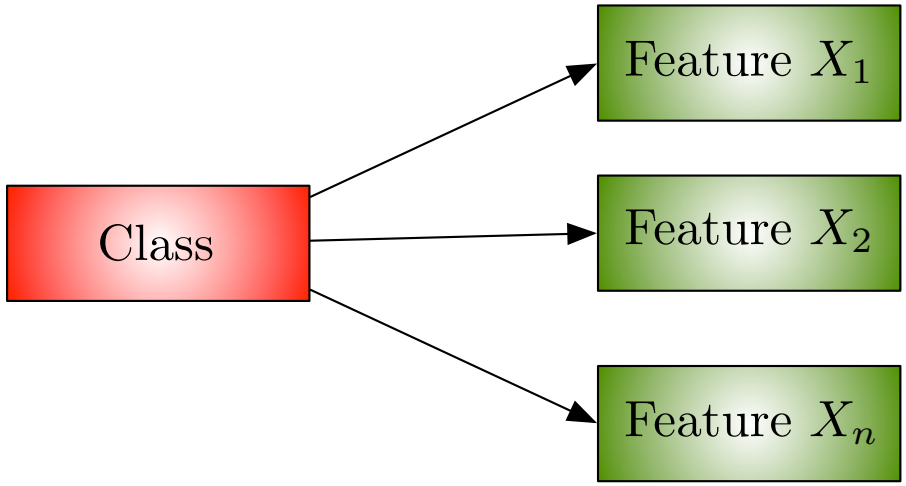
With no preferences on the value of p the posterior is strongly skewed towards one and the range $p \in [0.86, 1]$ contains 95% of posterior probability.

Posterior of an informed person



With the knowledge $p \in [0.1, 0.2]$ the posterior is strongly skewed towards 0.2 and the range $p \in [0.17, 0.2]$ contains 95% of posterior probability.

Model behind naive Bayes classifier



Underlying class value determines observed attributes

- ▷ Each attribute X_i is binary
- ▷ All variables are independent if class is fixed
- ▷ Sometimes we just ignore dependancies for easier modelling

Likelihood of the data

Let us assume that we know the probabilities

$$p_i = \Pr\left[X_i = 1 | Class = 0\right]$$
$$q_i = \Pr\left[X_i = 1 | Class = 1\right]$$

Then using the independence assumption we get

$$\Pr\left[X_1 = a_1, \ldots, X_n = a_n | Class = 0\right] = \prod_{i=1}^{n} p_i^{a_i}(1-p_i)^{1-a_i}$$

$$\Pr\left[X_1 = a_1, \ldots, X_n = a_n | Class = 1\right] = \prod_{i=1}^{n} q_i^{a_i}(1-q_i)^{1-a_i}$$

Prior and posterior for the class labels

Now it is straightforward to derive

$$\Pr\left[Class = 0 | \boldsymbol{X} = \boldsymbol{a}\right] = \frac{\prod_{i=1}^{n} p_i^{a_i}(1-p_i)^{1-a_i} \cdot \Pr\left[Class = 0\right]}{\Pr\left[\boldsymbol{X} = \boldsymbol{a}\right]}$$

$$\Pr\left[Class = 1 | \boldsymbol{X} = \boldsymbol{a}\right] = \frac{\prod_{i=1}^{n} q_i^{a_i}(1-q_i)^{1-a_i} \cdot \Pr\left[Class = 1\right]}{\Pr\left[\boldsymbol{X} = \boldsymbol{a}\right]}$$

which gives an *odd ratio*

$$\frac{\Pr\left[Class = 0 | \boldsymbol{X} = \boldsymbol{a}\right]}{\Pr\left[Class = 1 | \boldsymbol{X} = \boldsymbol{a}\right]} = \frac{\Pr\left[Class = 0\right]}{\Pr\left[Class = 1\right]} \cdot \frac{\prod_{i=1}^{n} p_i^{a_i}(1-p_i)^{1-a_i}}{\prod_{i=1}^{n} q_i^{a_i}(1-q_i)^{1-a_i}}$$

The resulting classifier is a linear classifer

By taking logarithm form the odd ratio we get

$$\log\left(\frac{\Pr\left[Class=0|\boldsymbol{X}=\boldsymbol{a}\right]}{\Pr\left[Class=1|\boldsymbol{X}=\boldsymbol{a}\right]}\right)=w_0+\sum_{i=1}^{n}w_i a_i$$

where

$$w_0=\log\left(\frac{\Pr\left[Class=0\right]}{\Pr\left[Class=1\right]}\right)+\sum_{i=1}^{n}\log\left(\frac{1-p_i}{1-q_i}\right)$$

$$w_i=\log\left(\frac{p_i}{1-p_i}\cdot\frac{1-q_i}{q_i}\right)$$

How to train the classifier?

A frequentistic approach is to fix probabilities from the training sample

$$p_i = \frac{\#\{\text{data points form class 0 with } X_i = 1\}}{\#\{\text{data points form class 0}\}}$$

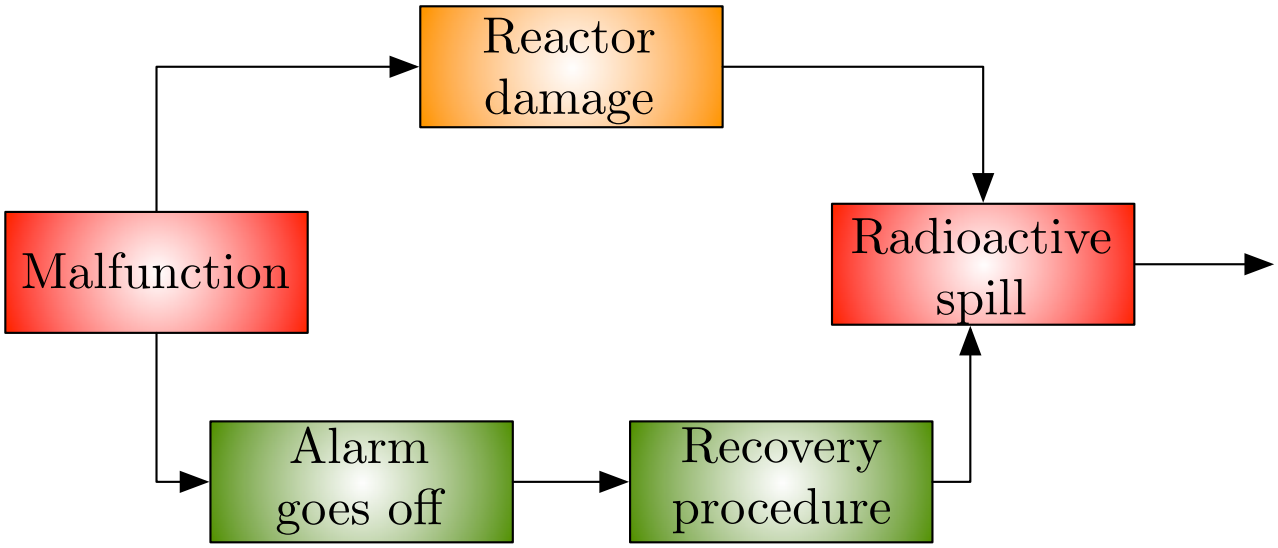
$$q_i = \frac{\#\{\text{data points form class 1 with } X_i = 1\}}{\#\{\text{data points form class 1}\}}$$

However if some value does not occur for X_i in the training sample we get overly confident results. Thus, Bayesian mean estimate is better alternative

$$p_i = \frac{\#\{\text{data points form class 0 with } X_i = 1\} + 1}{\#\{\text{data points form class 0}\} + 2}$$

$$q_i = \frac{\#\{\text{data points form class 1 with } X_i = 1\} + 1}{\#\{\text{data points form class 1}\} + 2}$$

Going beyond naive Bayesian models



Complex causal models are often defined through Bayesian networks

- ▷ A complex processes is first split into sub-events
- ▷ Direct causal dependencies between sub-events are detected
- ▷ Causation mechanisms are characterised with probability tables

Strength and weaknesses of Bayesian networks

Strengths

- ▷ Bayesian networks are easy to interpret
- ▷ Bayesian networks are good for formalising fuzzy background knowledge
- ▷ Estimation of individual probability tables is tractable
- ▷ There are tools for doing inference with Bayesian networks

Weaknesses

- ▷ You must know the causal structure of sub-events
- ▷ Identification of causal structure form data alone is very difficult
- ▷ It is notoriously difficult to model non-trivial causal dependencies
- ▷ Standard inference procedures often do not have closed solutions