

# Bayesian, Laplace or De Finetti Statistics?

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Over the last twenty years, Oxford has been strongly associated with Bayesian Statistics:

- ▶ Chris Holmes, Yee Whye Teh, Arnaud Doucet, Tom Rainforth, Geoff Nichols, Christopher Yau, Yarin Gal, Steve Roberts, Mike Osborne, Francois Caron, Mark van der Wilk, Judith Rousseau, ...

... and many graduate students and postdocs.

A billiard ball is rolled onto a table at random, landing somewhere along a line between 0 and 1 (the “true probability” of success).

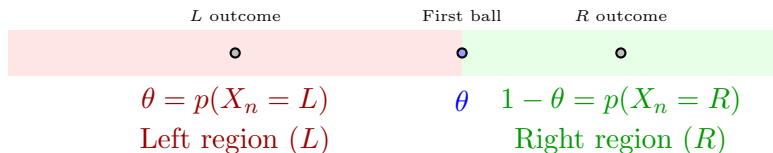
Then, another ball is rolled repeatedly, and each time we note whether it lands to the left or right of the first ball.

From these outcomes, Bayes asked:

*Given the number of successes and failures observed, what is the probability that the true chance of success lies within a certain interval?*

Let  $\theta \in (0, 1)$  denote the position of the first ball.

Let  $X_n \in \{L, R\}$  denote the left/right positioning of the  $n$ -th following balls.



The posterior distribution over  $\theta$  given  $X_1, \dots, X_N$ :

$$\begin{aligned} p(\theta|X_1, \dots, X_N) &= \frac{p(X_1, \dots, X_N|\theta)p(\theta)}{p(X_1, \dots, X_N)}, \\ &\propto \prod_{n=1}^N p(X_n|\theta)p(\theta), \\ &\propto \prod_{n=1}^N \underbrace{\theta^{\delta(X_n=L)}}_{p(X_n=L)} \underbrace{(1-\theta)^{\delta(X_n=R)}}_{p(X_n=R)} p(\theta), \\ &\propto \theta^{N_L} (1-\theta)^{N_R} p(\theta) \end{aligned}$$

where  $N_L = \sum_n \underbrace{\delta(X_n = L)}_{\text{Indicator function}}$  and  $N_R = N - N_L$ .

1700s: English Reverend Thomas Bayes wanted to know how to infer causes from effects.

Working problem:

*How could he learn the probability of a future event occurring if he only knew how many times it had occurred or not occurred in the past?*



Bayes considered problems similar to the motivating example and how you narrow down the area in which the first ball probably sits.

Each new piece of information constrains the area where the first ball probably is.

Bayes' system was:

*Initial Belief + New Data  $\Rightarrow$  Improved Belief.*

There were two enduring criticisms to Bayes' system:

- ▶ Mathematicians (even to this day) were horrified to see something as whimsical as a *guess* play a role in rigorous mathematics,
- ▶ Bayes said that if he didn't know what guess to make, he would just assign all possibilities equal probability to start. This concept of “prior belief” was too much for mathematicians.

Bayes never published his discovery, but his friend Richard Price found it among his notes after Bayes' death in 1761, re-edited it, and published it ( “*An Essay towards solving a Problem in the Doctrine of Chances*”)

Unfortunately, virtually no one seems to have read the paper, and Bayes' method lay unappreciated until Laplace.



In 1774, in “*Mémoire sur la probabilité des causes par les événements*”, Pierre Laplace independently identified the same idea but as the greater mathematician was able to systemise it using probability theory (**inductive reasoning**).

Laplace was the first true *Bayesian* and the work was integrated into his monumental work “*Théorie analytique des probabilités*” in 1812.

Strangely, Laplace also identified the **Central Limit Theorem** - a key result - underpinning what has become *frequentist* approaches.

1. Begin with **prior** belief over a parameter  $\theta$  encapsulated by a probability distribution  $p(\theta)$ .
2. Acquire data  $X$  according to some process/experiment which is controlled by  $\theta$ . The probability of observing  $X$  given  $\theta$  is known as the **likelihood**,  $p(X|\theta)$ .
3. Compute the **posterior** probability distribution  $p(\theta|X)$  of  $\theta$  given data  $X$  using **Bayes' Theorem**.

The Bayesian belief update is given by:

$$\underbrace{p(\theta|X)}_{\text{Posterior}} = \frac{\overbrace{p(X|\theta)}^{\text{Likelihood}} \overbrace{p(\theta)}^{\text{Prior}}}{\underbrace{p(X)}_{\text{Marginal Likelihood / Evidence}}}$$

where  $\theta$  denotes a parameter and  $X$  is data.

The potentially multi-dimensional integral

$$p(X) = \int p(X|\theta)dp(\theta)$$

is the computational barrier to the application of Bayesian belief updating.

## Definition

$$s(\theta) = \nabla_{\theta} \log p(X|\theta)$$

- ▶ Measures how sensitive the likelihood is to changes in  $\theta$ .
- ▶ Tells us the *direction* of steepest ascent of the likelihood.

## Connection to Bayes' Rule

Taking logs of Bayes' theorem:

$$\log p(\theta|X) = \log p(X|\theta) + \log p(\theta) - \log p(X)$$

Hence:

$$\nabla_{\theta} \log p(\theta|X) = \underbrace{\nabla_{\theta} \log p(X|\theta)}_{\text{Score Function}} + \nabla_{\theta} \log p(\theta) \text{ (+no terms involving } p(X)\text{)}$$

The update can be iterated:

$$\underbrace{p(\theta|X_{\text{new,old}})}_{\text{Updated Posterior}} = \frac{\overbrace{p(X_{\text{new}}|\theta)}^{\text{Likelihood}} \quad \overbrace{p(\theta|X_{\text{old}})}^{\text{Prior} = \text{Previous Posterior}}}{\underbrace{p(X_{\text{new}}|X_{\text{old}})}_{\text{Marginal Likelihood / Evidence}}}$$

where  $\theta$  denotes a parameter and  $X_{\text{old/new}}$  is old and new data respectively.

This allows *sequential updating* of knowledge.

Bayesian belief updating appeared on and off over the next couple of centuries but was never the mainstream approach to probability or statistics.

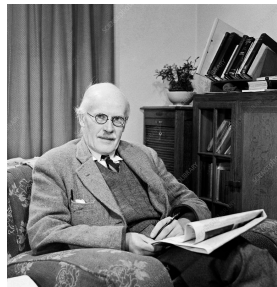
John Stuart Mill denounced probability as:

*“ignorance ... coined into science”*

Subjectivity and mathematics did not mix.

Bayesian theory continued to be developed sporadically (Emile Borel, Frank Ramsay, Bruno De Finetti):

- ▶ Harold Jeffreys, a geologist, made Bayesian Statistics slightly more fashionable in the 1930s.
- ▶  $p(\theta|X)$  vs  $p(X|\theta)$ .
- ▶ Jeffreys was a counterpoint to the great Ronald Fisher - who advocated sampling-based statistical thinking.
- ▶ Ironically famous for the *Jeffrey's prior* which is derived using *Fisher information* (variance of the score).



As the more robust personality, Fisher “won”.

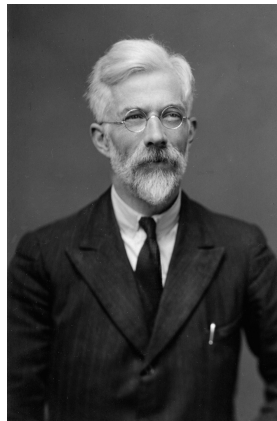
Ronald Fisher and Jerzi Neyman killed Bayesian thinking in the first half of the 1900s.

*“(Bayes) attempted to find, by observing a sample, the actual probability that the population value lay in any given range.”*

*“Such a problem is indeterminate without knowing the statistical mechanism under which different values of (a parameter) come into existence; it cannot be solved from the data supplied by a sample, or any number of samples, of the population.”*

*“We can know nothing of the probability of hypotheses or hypothetical quantities.”*

Fisher was never comfortable with  $p(\theta|X)$  and preferred  $p(X|\theta)$ .





Fisher’s problematic **prior** is often cited as one of the distinctive and advantageous features of Bayesian Statistics.

It allows you to:

1. Define *preferred* values or ranges for parameters.
2. Regularise and make models more *stable*.
3. Enforce bias or assumptions.

Together with **uncertainty quantification**, these are often reasons given for why it is good to be Bayesian.

- ▶ But how do we know the prior exists?
- ▶ How do we know if we keep updating then the true posterior is obtained?
- ▶ What is the posterior if we had an infinite number of observations?

### Definition

A  **$\sigma$ -algebra**  $\mathcal{F}$  on a sample space  $\Omega$  is a collection of subsets of  $\Omega$  such that:

1.  $\Omega \in \mathcal{F}$  (entire sample space is an event),
2. If  $A \in \mathcal{F}$ , then  $A^c \in \mathcal{F}$  (closed under complements),
3. If  $A_1, A_2, A_3, \dots \in \mathcal{F}$ , then  $\bigcup_{i=1}^{\infty} A_i \in \mathcal{F}$  (closed under countable unions).

### Intuition

$\mathcal{F}$  represents all events that are **knowable or observable**.

Each  $A \in \mathcal{F}$  corresponds to a “question” we can ask about the random world.

### Example: Coin flips

Let  $X_t$  = result of the  $t$ -th flip, and  $\mathcal{F}_t = \sigma(X_1, \dots, X_t)$ .

Then  $\mathcal{F}_t$  contains all events determined by the first  $t$  outcomes.

### Definition

A **filtration**  $(\mathcal{F}_t)_{t \geq 0}$  is an **increasing family of  $\sigma$ -algebras**:

$$\mathcal{F}_0 \subseteq \mathcal{F}_1 \subseteq \mathcal{F}_2 \subseteq \dots$$

### Meaning

As time passes, we only gain information:

More data  $\Rightarrow$  more events become measurable.

### Intuition

$\mathcal{F}_t$  encodes “everything known up to time  $t$ .” A filtration represents the process of knowledge accumulation.

### Definition

A stochastic process  $(X_t)$  is **adapted** to a filtration  $(\mathcal{F}_t)$  if

$X_t$  is  $\mathcal{F}_t$ -measurable for all  $t$ .

### Interpretation

- ▶ Being adapted means the process respects a causal flow of information.
- ▶  $X_t$  depends only on information available up to time  $t$  — not on the future.

## Definition

A process  $(X_t)$  is a **martingale** with respect to the filtration  $(\mathcal{F}_t)$  if:

$$\mathbb{E}[X_{t+1} \mid \mathcal{F}_t] = X_t,$$

and  $\mathbb{E}[|X_t|] < \infty$ .

## Intuition

- ▶ Given all current information, your expected future value is the same as the present value.
- ▶ Martingales model a “fair game” — there is no predictable gain over time.

*Note: A martingale is defined with respect to a filtration, but the condition at each step uses the  $\sigma$ -algebra  $\mathcal{F}_t$  for the corresponding time step in the filtration.*

### Definition

For any integrable random variable  $X$  and filtration  $(\mathcal{F}_t)$ :

$$M_t = \mathbb{E}[X \mid \mathcal{F}_t]$$

defines the **Doob martingale**.

### Martingale Property

By the Tower Property (Law of Iterated Conditional Expectations):

$$\mathbb{E}[M_{t+1} \mid \mathcal{F}_t] = \underbrace{\mathbb{E}[\mathbb{E}[X \mid \mathcal{F}_{t+1}] \mid \mathcal{F}_t]}_{\text{Tower Property}} = \mathbb{E}[X \mid \mathcal{F}_t] = M_t.$$

The next step conditional expectation (given no new information) is the same as the current conditional expectation.

## Intuition

- ▶ Imagine trying to predict a final outcome  $X$  (like tomorrow's stock price or total rainfall) as you receive information over time.
- ▶ Each  $M_t = \mathbb{E}[X|\mathcal{F}_t]$  is your best guess given what you currently know.
- ▶ The Doob martingale says that, although your estimates change as new data arrives, they do so fairly — there is no tendency to systematically over- or under-estimate the final outcome.
- ▶ In the absence of new information, things stay the same.



## Setup

Let  $\Theta$  be a parameter with prior  $\Pi_0$ , and let

$$\mathcal{F}_t = \sigma(X_1, \dots, X_t)$$

represent the information from the first  $t$  data points.

## Posterior probability as conditional expectation

For any measurable set  $A \subseteq \Theta$ :

$$M_t = \mathbb{E}[\mathbf{1}_{\{\Theta \in A\}} \mid \mathcal{F}_t] = \Pi_t(A) = P(\Theta \in A \mid X_1, \dots, X_t).$$

## Martingale property

$$\mathbb{E}[M_{t+1} \mid \mathcal{F}_t] = M_t.$$

Hence  $(\Pi_t(A))_{t \geq 0}$  is a bounded martingale.

## Doob's Martingale Convergence Theorem

If  $(M_t)$  is a bounded martingale, then

$$M_t \rightarrow M_\infty \quad \text{almost surely as } t \rightarrow \infty.$$

## Application to posterior probabilities

For every  $A \subseteq \Theta$ ,

$$\Pi_t(A) = \mathbb{E}[\mathbf{1}_{\{\Theta \in A\}} \mid \mathcal{F}_t] \quad \Rightarrow \quad \Pi_t(A) \rightarrow \Pi_\infty(A) \text{ a.s.}$$

Thus, each posterior probability converges almost surely.

As  $t$  increases,  $\mathcal{F}_t$  grows, so your conditional expectation “stabilises” — you have ve learned all there is to learn about  $X$  from the filtration.

Assume data come from the true parameter  $\theta_0$

That is,  $X_i \sim P_{\theta_0}$  i.i.d.

Likelihood behaviour

For any  $\theta \neq \theta_0$ ,

$$\frac{p_{\theta}(X_1, \dots, X_n)}{p_{\theta_0}(X_1, \dots, X_n)} \rightarrow 0 \quad \text{a.s. under } P_{\theta_0}.$$

Posterior collapse

Hence:

$$\Pi_t(A) \rightarrow \begin{cases} 1, & \text{if } \theta_0 \in A, \\ 0, & \text{if } \theta_0 \notin A, \end{cases} \quad P_{\theta_0}\text{-a.s.}$$

The limiting random measure is  $\Pi_{\infty} = \delta_{\theta_0}$ .

- ▶ Each posterior probability  $\Pi_t(A)$  is a martingale:

$$\mathbb{E}[\Pi_{t+1}(A) \mid \mathcal{F}_t] = \Pi_t(A).$$

- ▶ Bounded martingales converge (Doob, 1953).
- ▶ Under the true parameter, posterior mass outside neighborhoods of  $\theta_0$  vanishes.
- ▶  $\Rightarrow$  Posterior collapses:  $\Pi_t \Rightarrow \delta_{\theta_0}$ .

## Summary

- ▶ Bayesian updating = Doob martingale + Martingale convergence theorem  $\Rightarrow$  Almost sure posterior consistency.
- ▶ Remarkably this is a pure measure-theoretic result with few conditions.
- ▶ More detailed exposition: <https://arxiv.org/pdf/1801.03122>

The Bayesian belief update is given by:

$$\underbrace{p(\theta|X)}_{\text{Posterior}} = \frac{\overbrace{p(X|\theta)}^{\text{Likelihood}} \overbrace{p(\theta)}^{\text{Prior}}}{\underbrace{p(X)}_{\text{Marginal Likelihood / Evidence}}}$$

where  $\theta$  denotes a parameter and  $X$  is data.

**How do I now a prior exists?**

Suppose I flip 5 coins:

$$X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T$$

What is the probability of an  $H$  on the 6th flip?

$$P(X_6 = H | X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T)$$

If each flip is independent and identically distributed (iid) then:

$$\begin{aligned} &P(X_6 = H | X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T) \\ &= \frac{P(X_6 = H, X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T)}{P(X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T)} \end{aligned}$$

If each flip is independent and identically distributed (iid) then:

$$\begin{aligned} P(X_6 = H | X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T) &= \\ &= \frac{P(X_6 = H) \cancel{P(X_5 = T)} \cancel{P(X_4 = H)} \cancel{P(X_3 = H)} \cancel{P(X_2 = T)} \cancel{P(X_1 = H)}}{\cancel{P(X_5 = T)} \cancel{P(X_4 = H)} \cancel{P(X_3 = H)} \cancel{P(X_2 = T)} \cancel{P(X_1 = H)}}, \\ &= P(X_6 = H) \end{aligned}$$

The history of coin flips tells me nothing?



Ah, but there is a parameter - the probability of getting an  $H$  which we call  $\theta$ !

$$\begin{aligned} P(X_6 = H, \theta | X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T) = \\ = P(X_6 = H | \theta) p(\theta | X_1 = H, X_2 = T, X_3 = H, X_4 = H, X_5 = T) \end{aligned}$$

The history of flips tells us about  $\theta$  which then helps us to predict  $X_6$ .

The coin flips are only iid if I know  $\theta$  (conditionally iid).

In general, we can introduce **shared** parameters to make the joint distribution conditionally iid:

$$p(X_1 = x_1, \dots, X_n = x_n | \theta) = p(X_1 = x_1 | \theta) p(X_2 = x_2 | \theta) \cdots p(X_n = x_n | \theta)$$

The data  $X$  is **conditionally** iid **given** parameter  $\theta$  so historical information can feed into predicting the future:

$$\begin{aligned} p(X_{n+1}, \theta | X_1 = x_1, \dots, X_n = x_n) = \\ p(X_{n+1} | \theta) p(\theta | X_1 = x_1, X_2 = x_2, \dots, X_n = x_n) \end{aligned}$$

**Does this construction *always* exist? Must there be a suitable parameter?**

If a sequence of random variables is **exchangeable** then there is no information in their order, i.e.

$$p(X_1 = x_1, X_2 = x_2, \dots, X_n = x_n) = p(X_{\pi(1)} = x_1, X_{\pi(2)} = x_2, \dots, X_{\pi(n)} = x_n)$$

where  $\pi$  denotes a permutation operation (e.g.  $(1, 2, 3, 4, 5) \rightarrow (5, 3, 1, 4, 2)$  )

Exchangeability is a *weaker* assumption than independence.

Crucially, exchangeability assumptions apply to many real-world problems.

The remarkable finding by De Finetti is that for *any* infinite exchangeable random sequence, it is *always* possible to write:

$$p(X_1 = x_1, X_2 = x_2, \dots, X_n = x_n) = \int \prod_{i=1}^n p(X_i|\theta)p(\theta)d\theta$$

as  $n \rightarrow \infty$ .

The implications are profound, there is:

- ▶ a parameter  $\theta$  which makes  $X_1, \dots, X_n$  conditionally independent,
- ▶ this parameter must always have a distribution  $p(\theta)$  associated with it,
- ▶ and there is always a sampling model  $p(X|\theta)$ ,
- ▶ exchangeable sequences are mixtures of conditionally iid sequences under the prior measure.

We *should be* Bayesian when exchangeability applies!

We can also use the De Finetti representation theorem in reverse:

$$\int \prod_{i=1}^n p(X_i|\theta)p(\theta)d\theta \rightarrow p(X_1 = x_1, X_2 = x_2, \dots, X_n = x_n),$$

If we define a **latent variable model** then the marginal joint distribution of the data must be exchangeable.

This allows us to model complex exchangeable distributions in an *indirect* way.

- ▶ De Finetti proved for *binary* exchangeable sequences only.
- ▶ The Generalised Representation Theorem by Hewitt & Savage appeared later.
- ▶ Strictly speaking only applies to *infinite* sequences but Diaconis and Freedman have developed forms/bounds for finite sequences.
- ▶ Things become complex for partial exchangeability and more nuanced dependence structures.
- ▶ It is an *existence* theorem only.

Doob and De Finetti provide strong theoretical justification for Bayesian reasoning:

De Finetti shows that if data are exchangeable, a Bayesian model must exist.

Doob shows that if we update beliefs through Bayes' rule, they will converge consistently.

Exchangeability is not just a theoretical assumption — it is a practical modelling principle.

It gives us a situation when Bayesian inference *optimally* applies, and guides how to build our models.

## Prompt given to an LLM

Example 1: "I loved this movie!" → *Positive*

Example 2: "The plot was boring and predictable." → *Negative*

Example 3: "The soundtrack was beautiful." → *Positive*

Now classify: "The ending was disappointing." → ?

## Questions:

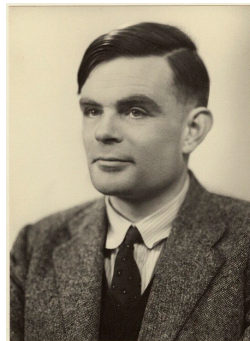
- ▶ What label do you expect the model to output?
- ▶ How can it infer that “disappointing” implies a *Negative* label?
- ▶ The model’s weights never changed — so where does this learning happen?



Throughout its history, Bayesian approaches have always been synonymous with military applications:

- ▶ In 1890, Joseph Bertrand helped French/Russia artillery officers to use Bayesian approaches to target enemy locations.
- ▶ Kolmogorov suggested the same idea in WW2.
- ▶ Turing used it for code cracking.

Modern military information fusion systems are all based on Bayesian frameworks.



Following the Second World War:

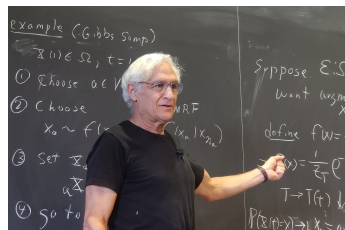
- ▶ I.J. Good began to revisit the system of Bayesian Inference.
- ▶ Lindley and Savage put mathematical rigour into the approach.
- ▶ Lindley began to put Bayesian academics into Professorships.



Bayesian approaches had a fundamental problem.

Outside of a few mathematically convenient examples, it was not tractable.

Seminal paper by Geman & Geman (1984) that linked Bayesian computation to Statistical Physics.



Suppose there are multiple parameters:

$$\theta = \{\theta_1, \theta_2\}.$$

The marginal likelihood:

$$p(X) = \int p(X|\theta)p(\theta)d\theta$$

requires a multi-dimensional integral over the entire parameter space.

Without the marginal likelihood we cannot directly compute,  $p(\theta|X)$ .

Geman & Geman overcame this with the Gibbs Sampler:

$$\theta_1 \sim p(\theta_1|\theta_2, X),$$

$$\theta_2 \sim p(\theta_2|\theta_1, X).$$

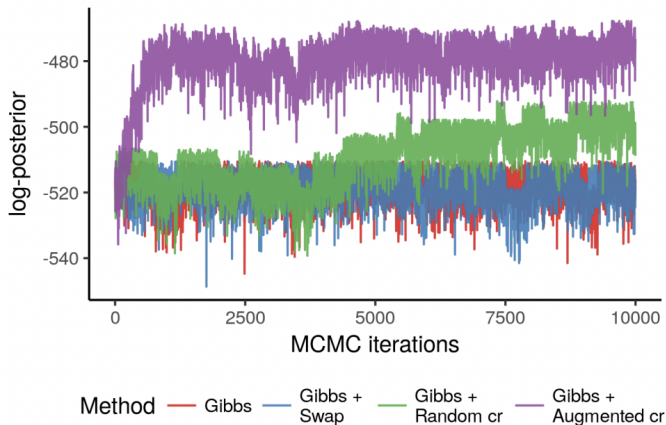
The Markov chain generated by this process has the posterior distribution as its stationary distribution. As the chain mixes, the samples converge in distribution to draws from this posterior.

The Gibbs Sampler was one of a family of techniques known now as **Markov Chain Monte Carlo** (MCMC) methods.

Physicists and mathematicians looking at statistical properties of molecular configurations had already worked it out (Metropolis (1953) and Hastings (1970)).

Suddenly through the 1990s and early 2000s, Bayesian computation now enabled widespread use of the technique.

MCMC algorithms only provide *asymptotic* guarantees of convergence.



<http://proceedings.mlr.press/v89/martens19a/martens19a.pdf>

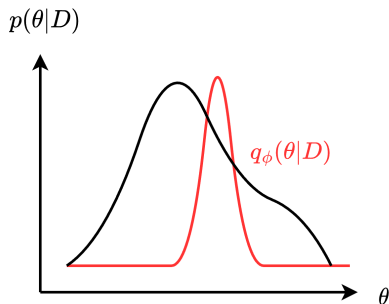
As datasets grew larger and models became higher-dimensional, traditional MCMC-based Bayesian inference began to struggle computationally.

- ▶ **Curse of dimensionality** - Sampling efficiency deteriorates exponentially with dimensionality.
- ▶ **Conditional sampling** - Sequential conditional updates make MCMC inherently difficult to parallelise.
- ▶ **Need to update all model parameters** - each MCMC iteration requires updating all parameters — in some models, the number of parameters grows with the data size.

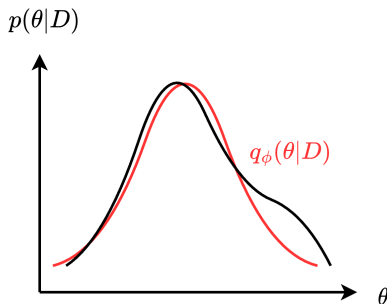
*Note - see Chris Nemeth's lecture on Monte Carlo Methods for recent work in this area.*



Physics also offered another approach. *What if I approximate the posterior distribution with a simpler form?*



Unoptimised variational approximation



Optimised variational approximation

Variational methods trade *asymptotic exactness* for *optimisation-based approximation*.

The **mean-field approximation** ( $O(NP)$ ):

$$p(\theta_1, \dots, \theta_P | X) \approx \prod_{p=1}^P q_{\phi_p}(\theta_p | X)$$

In mean-field approximation, a set of *variational parameters* are tied to each data point. The use of **amortisation** provides a second layer of approximation ( $O(N)$ ):

$$\phi_p = f_\nu(X_p)$$

where instead of optimising all variational parameters, you learn a function that maps relevant subsets of data to estimates of those parameters.

*Note - see Hanwen's lecture this PM*

In recent years, *explicit* Bayesian modelling has reduced in popularity or exposure.

But unlike previous centuries, it remains in widespread *implicit* use throughout science and engineering.

Traditional Bayesian analysis is *purist*, pragmatically we need to deal with the following:

- ▶ **Prior elicitation** - defining a “good” prior is hard (see George D lecture on Diffusion Models),
- ▶ **Uncertainty quantification** - MCMC and VB under-estimate posterior variance,
- ▶ **Causality** - Bayesian reasoning is concerned with probabilistic coherence not causal reasoning,
- ▶ **Model misspecification** - all the theorems only hold under the *true* model.

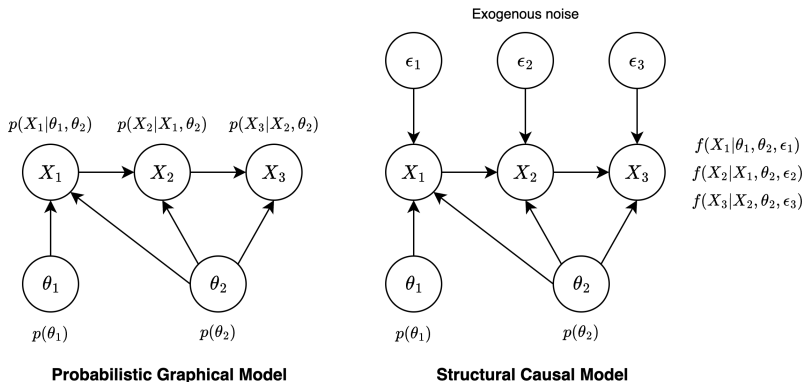
Modern Bayesian research has sought to address these issues.

The term originates from *likelihood-based inference* in Statistics but (can) mean much more in Bayesian Statistics:

$$\underbrace{p(\theta|X)}_{\text{Posterior}} = \frac{\overbrace{p(X|\theta)}^{\text{Likelihood}} \overbrace{p(\theta)}^{\text{Prior}}}{\underbrace{p(X)}_{\text{Marginal Likelihood / Evidence}}}$$

- ▶ The likelihood contains a **probabilistic graphical model** (PGM) for your problem.
- ▶ The PGM describes the **statistical dependencies** between all quantities in your model.
- ▶ Sometimes referred to as harbouring the **data generating process** (DGP) but not actually true ...

A **structural causal model** (SCM) allows **interventions** while a **probabilistic graphical model** (PGM) does not.



A PGM does not define a data generation model *under interventions*.

While causal models add structure to describe interventions, Bayesian reasoning doesn't require causality — it operates at the level of probabilistic coherence.

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## **Position: Probabilistic Modelling is Sufficient for Causal Inference**

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**Bruno Mlodozienec<sup>1,2</sup> David S. Krueger<sup>3</sup> Richard E. Turner<sup>1,4</sup>**

*“Always write down the probability of everything.”*

– Steve Gull (MacKay, 2003, p. 61)

A recent line of thought is to do away with causal inference approaches and to add specific mirror-like “counterfactual” variables to mirror an intervention.<sup>1</sup>

<sup>1</sup><https://openreview.net/forum?id=V1FP9WDKa7>

When discussing model specification, Bayesians refer to **M-closed** and **M-open**.

In the **M-closed world**:

- ▶ The true data-generating process lies within our model family  $\{p(X|\theta) : \theta \in \Theta\}$ .
- ▶ Classical Bayes is **well-calibrated and coherent**.
- ▶ Posterior concentrates around the true parameter  $\theta^*$ .

In the **M-open world**:

- ▶ The true data-generating process is **outside** our model family.
- ▶ The likelihood  $p(X|\theta)$  is **misspecified**.
- ▶ Classical Bayes can be **overconfident or misleading**.
- ▶ Traditionally, Bayesian tried to overcome this by using **Bayesian nonparametrics** to define infinite-dimensional function and random probability measures.

In M-open settings, we need a more flexible update rule —  
**Generalised Bayes** replaces the likelihood with a *loss*.



- ▶ Statistics can be viewed as a **decision problem under uncertainty**.
- ▶ Let  $\theta$  denote an unknown state of nature and  $a$  an action.
- ▶ Decisions are evaluated by a **loss function**  $L(a, \theta)$ .
- ▶ The optimal action minimises **expected loss**:

$$a^* = \arg \min_a \mathbb{E}_{p(\theta|X)}[L(a, \theta)].$$

- ▶ Contrast with classical decision theory where the expectation is wrt to sampling distribution of  $X$ .

For a Bayesian, rational decisions are those that *minimise* **posterior expected loss**.

- ▶ Bissiri, Holmes & Walker (2016): updating beliefs **is itself a decision problem**.
- ▶ Choose a posterior  $q(\theta)$  to minimise:

$$q^*(\theta) = \arg \inf_q \mathbb{E}_q[L(\theta; X)] + \frac{1}{\eta} \text{KL}(q||p).$$

- ▶ Here:
  - ▶  $L(\theta; X)$ : loss describing the data's informational content (replaces likelihood),
  - ▶  $\text{KL}(q||p)$ : penalty for deviating from the prior.
- ▶ Solution:

$$q^*(\theta) \propto p(\theta) \exp(-\eta L(\theta; X))$$

— the Generalised Bayes posterior.

*Updating = deciding which beliefs to hold, balancing data and prior.*

Also, known as a **Gibbs posterior**:

$$p(\theta|X) \propto p(\theta) \exp(-\eta L(\theta; X))$$

- ▶  $p(\theta)$ : prior belief
- ▶  $L(\theta; X)$ : data-dependent loss (not necessarily log-likelihood)
- ▶  $\eta > 0$ : learning-rate / temperature parameter

Retains the Bayesian **form**, but broadens what “evidence” means.

*Note: Standard Bayes is a special case of the Generalised Bayes update.*

$$L(\theta; X) = -\log p(X|\theta) \quad \Rightarrow \quad p(\theta|X) \propto p(\theta)p(X|\theta)$$

Generalised Bayes offers a way to:

- ▶ Retain **coherent belief updating** without a likelihood
- ▶ Incorporate **model misspecification and robustness**
- ▶ Align inference with **predictive or decision objectives**

*Generalised Bayes is Bayesian reasoning for an imperfect world.*

The update:

$$p(\theta|X) \propto p(\theta) \exp(-\eta L(\theta; X))$$

is only valid for **parameters that appear in the loss function**  $L(\theta; X)$ .

If a component  $\theta_j$  does not appear in  $L$ :

$$p(\theta_j|X) = p(\theta_j)$$

- ▶ Coherence requires belief revision only where the loss provides evidence.
- ▶ Prevents spurious updating of uninformed parameters.

*Generalised Bayes is local in scope: only beliefs connected to the loss are revised.*

- ▶ **PAC-Bayes / Learning Theory:** treat  $L$  as empirical risk  $\rightarrow$  connects to regularised risk minimisation.
- ▶ **Robust Bayes:** replace log-likelihood with  $\beta$ -divergence or Huber loss  $\rightarrow$  down-weight outliers.
- ▶ **Online learning:** adapt  $\eta$  dynamically  $\rightarrow$  information-theoretic updates.
- ▶ **Approximate inference:** variational or ensemble updates mimic generalised Bayes behaviour.

When our model isn't perfect, *Generalised Bayes keeps us Bayesian in spirit*:

- ▶ Still **belief updating**, just under broader notions of evidence
- ▶ **Retains coherence** preserves consistency – Doob still applies
- ▶ **Bridges with modern approaches in AI**, e.g. gives a framework for what were previously ad-hoc regularisation strategies

For further discussion, see:

“Position: Bayesian Deep Learning is Needed in the Age of Large-Scale AI”

<https://arxiv.org/abs/2402.00809>

BREAK HERE



A simple **latent variable model** has the form:

$$X|Z \sim N(\mu(Z), \sigma^2(Z)), \quad (1)$$

$$Z \sim N(0, 1), \quad (2)$$

and we want to learn  $p(Z|X)$  which is the posterior distribution over the unobserved latent variable  $Z$  given the observation  $X$ .

Frequently used where dimensionality of  $Z$  is much lower than  $X$ .

If I choose:

- ▶  $\mu(Z)$  is a function described by a neural network,
- ▶  $\sigma^2(Z)$  is a function described by a neural network,
- ▶ Approximate  $p(Z|X)$  via a Normal distribution,
- ▶ Apply amortisation to reduce the number of variational parameters.
- ▶ Perform optimisation via stochastic gradient descent.

The result is an algorithm called a **Variational Autoencoder** (VAE):

$$[\text{VAE}] = [\text{Latent Variable Model}] + [\text{Inference Choices}]$$

**Encoder** - maps data  $X$  to (distribution over) latent variables  $Z$

**Decoder** - maps latent variable  $Z$  to (distribution over) data  $X$

**Loss function**

$$l(\theta, \phi) = -\mathbb{E}_{Z \sim q_\theta(Z|X)} [\log p_\phi(X|Z)] + \mathbb{KL}(q_\theta(Z|X) || p(Z))$$

Black box building blocks have simplified and normalised the application of Bayesian techniques by unifying models and computation.

$$\underbrace{[\text{Structure}] - [\text{Distributions}] - [\text{Inference}]}_{\text{Model}} \\ \underbrace{\hspace{10em}}_{\text{Black Box Bayesian}}$$

This convergence underpins the utility of **probabilistic programming languages**, e.g. STAN, Pyro, PyMC.

But black box techniques intrinsically conflate the statistical and computing biases making diagnostic more difficult.

Bayesian thinking in 2025 goes beyond the mechanics of applying Bayes' rule and computational inference:

- ▶ Foundational motivations for Bayesian reasoning give strong theoretical justification for its optimality
- ▶ Existing and new AI approaches should be able to reason and compute in a Bayesian way
- ▶ Adds to the battery of expectations we might expect of the highest-performing models.