## Quiz

Suppose we are testing $H_{0}$ vs $H_{1}$ and find $p<0.05$ and reject $H_{0}$ at the $5 \%$ level. Which, if any, of these statements is true:

- The probability $H_{0}$ is true is 0.05
- If $H_{0}$ is false, then we will correctly reject $H_{0} 95 \%$ of the time.
- If we observe $p=0.05$, the probability we falsely reject $H_{0}$ given that it is true is 0.05 .
- If $H_{0}$ is true, we will falsely reject it $5 \%$ of the time


## Quiz

Suppose we are testing $H_{0}$ vs $H_{1}$ and find $p<0.05$ and reject $H_{0}$ at the $5 \%$ level. Which, if any, of these statements is true:

- The probability $\mathrm{H}_{0}$ is true is 0.05 False
- If $H_{0}$ is false, then we will correctly reject $H_{0} 95 \%$ of the time. False
- If we observe $p=0.05$, the probability we falsely reject $H_{0}$ given that it is true is 0.05 .False
- If $H_{0}$ is true, we will falsely reject it $5 \%$ of the time True

Personal view: p-values are not the problem - it is the belief that every dataset should give a yes/no answer.

# Approximate Bayesian Computation (ABC): inference for intractable computer models 

Richard Wilkinson<br>Newton Institute 2019: The Fickle Heart<br>School of Mathematics and Statistics<br>University of Sheffield

May 14, 2019

## Why be Bayesian?

a man's attitude toward inference, like his attitude towards religion, is determined by his emotional make-up, not by reason or mathematics.

M Kendall

## Why be Bayesian do Bayesian analyses？

Why be Bayesian do Bayesian analyses?

- Coherence: under various sets of axioms, Bayes is the only sensible choice.... cf Jaynes, de Finetti, Jeffreys etc.


## Why be Bayesian do Bayesian analyses?

- Coherence: under various sets of axioms, Bayes is the only sensible choice.... of Jaynes, de Finetti, Jeffreys etc.
- Build in expert belief / rule out things we know are unlikely / regularise the solution


## Why be Bayesian do Bayesian analyses?

- Coherence: under various sets of axioms, Bayes is the only sensible choice.... cf Jaynes, de Finetti, Jeffreys etc.
- Build in expert belief / rule out things we know are unlikely / regularise the solution
- It is all just probability
- makes combining different uncertainties easy/possible e.g. calibrated prediction

$$
p(y \mid x)=\int p(y \mid \theta) p(\theta \mid x) \mathrm{d} \theta
$$

- deals with equifinality (ie multiple feasible values, under-specified systems)
- Simpler! Q:What's the difference between probability, significance, coverage, confidence, $p$-values etc


## Why be Bayesian do Bayesian analyses?

- Coherence: under various sets of axioms, Bayes is the only sensible choice.... cf Jaynes, de Finetti, Jeffreys etc.
- Build in expert belief / rule out things we know are unlikely / regularise the solution
- It is all just probability
- makes combining different uncertainties easy/possible e.g. calibrated prediction

$$
p(y \mid x)=\int p(y \mid \theta) p(\theta \mid x) \mathrm{d} \theta
$$

- deals with equifinality (ie multiple feasible values, under-specified systems)
- Simpler! Q:What's the difference between probability, significance, coverage, confidence, p -values etc
- It is possible / increasingly easy to do Bayesian inference
- Frequentist procedures are 'mathematically challenging' to derive for complex models


## Why be Bayesian do Bayesian analyses?

- Coherence: under various sets of axioms, Bayes is the only sensible choice.... cf Jaynes, de Finetti, Jeffreys etc.
- Build in expert belief / rule out things we know are unlikely / regularise the solution
- It is all just probability
- makes combining different uncertainties easy/possible e.g. calibrated prediction

$$
p(y \mid x)=\int p(y \mid \theta) p(\theta \mid x) \mathrm{d} \theta
$$

- deals with equifinality (ie multiple feasible values, under-specified systems)
- Simpler! Q:What's the difference between probability, significance, coverage, confidence, p -values etc
- It is possible / increasingly easy to do Bayesian inference
- Frequentist procedures are 'mathematically challenging' to derive for complex models

The downsides:

- We have to choose a prior.
- In practice, 'priors of convenience' are often used. You need to really care about the answer to bother with expert elicitation.
- You can check robustness wrt your choices.


## Why be Bayesian do Bayesian analyses?

- Coherence: under various sets of axioms, Bayes is the only sensible choice.... cf Jaynes, de Finetti, Jeffreys etc.
- Build in expert belief / rule out things we know are unlikely / regularise the solution
- It is all just probability
- makes combining different uncertainties easy/possible e.g. calibrated prediction

$$
p(y \mid x)=\int p(y \mid \theta) p(\theta \mid x) \mathrm{d} \theta
$$

- deals with equifinality (ie multiple feasible values, under-specified systems)
- Simpler! Q:What's the difference between probability, significance, coverage, confidence, p-values etc
- It is possible / increasingly easy to do Bayesian inference
- Frequentist procedures are 'mathematically challenging' to derive for complex models

The downsides:

- We have to choose a prior.
- In practice, 'priors of convenience' are often used. You need to really care about the answer to bother with expert elicitation.
- You can check robustness wrt your choices.
- There is no need for your posterior to relate to the world
- Post-hoc checks (calibration etc) can help, but there are no frequency guarantees:


## Calibration

- For most simulators we specify parameters $\theta$ and i.c.s and the simulator, $f(\theta)$, generates output $X$.
- The inverse-problem: observe data $D$, estimate parameter values $\theta$ which explain the data.

The Bayesian approach is to find the posterior distribution

$$
\begin{aligned}
& \pi(\theta \mid D) \propto \pi(\theta) \pi(D \mid \theta) \\
& \text { posterior } \propto \\
& \quad \text { prior } \times \text { likelihood }
\end{aligned}
$$



## Intractability

$$
\pi(\theta \mid D)=\frac{\pi(D \mid \theta) \pi(\theta)}{\pi(D)}
$$

- usual intractability in Bayesian inference is not knowing $\pi(D)$.


## Intractability

$$
\pi(\theta \mid D)=\frac{\pi(D \mid \theta) \pi(\theta)}{\pi(D)}
$$

- usual intractability in Bayesian inference is not knowing $\pi(D)$.
- a problem is doubly intractable if $\pi(D \mid \theta)=c_{\theta} p(D \mid \theta)$ with $c_{\theta}$ unknown (cf Murray, Ghahramani and MacKay 2006)
- a problem is completely intractable if $\pi(D \mid \theta)$ is unknown and can't be evaluated (unknown is subjective). I.e., if the analytic distribution of the simulator, $f(\theta)$, run at $\theta$ is unknown.

Completely intractable models are where we need to resort to ABC methods

## Approximate Bayesian Computation (ABC)

If the likelihood function is intractable, then ABC (approximate Bayesian computation) is one of the few approaches we can use to do inference.

## Approximate Bayesian Computation (ABC)

If the likelihood function is intractable, then ABC (approximate Bayesian computation) is one of the few approaches we can use to do inference.

ABC algorithms are a collection of Monte Carlo methods used for calibrating simulators

- they do not require explicit knowledge of the likelihood function
- inference is done using simulation from the model (they are 'likelihood-free').


## Approximate Bayesian Computation (ABC)

If the likelihood function is intractable, then ABC (approximate Bayesian computation) is one of the few approaches we can use to do inference.

ABC algorithms are a collection of Monte Carlo methods used for calibrating simulators

- they do not require explicit knowledge of the likelihood function
- inference is done using simulation from the model (they are 'likelihood-free').

ABC methods are widely used primarily because they are

- Simple to implement
- Intuitive
- Embarrassingly parallelizable
- Can usually be applied


## Plan

i. Basics
ii. Efficient sampling algorithms
iii. Regression adjustments/post-hoc corrections
iv. Summary statistics
v. Inference for misspecified models

Basics

## ‘Likelihood-Free' Inference

## Rejection Algorithm

- Draw $\theta$ from prior $\pi(\cdot)$
- Accept $\theta$ with probability $\pi(D \mid \theta)$

Accepted $\theta$ are independent draws from the posterior distribution, $\pi(\theta \mid D)$.

## ‘Likelihood-Free' Inference

## Rejection Algorithm

- Draw $\theta$ from prior $\pi(\cdot)$
- Accept $\theta$ with probability $\pi(D \mid \theta)$

Accepted $\theta$ are independent draws from the posterior distribution, $\pi(\theta \mid D)$. If the likelihood, $\pi(D \mid \theta)$, is unknown:

## 'Mechanical' Rejection Algorithm

- Draw $\theta$ from $\pi(\cdot)$
- Simulate $X \sim f(\theta)$ from the computer model
- Accept $\theta$ if $D=X$, i.e., if computer output equals observation

The acceptance rate is $\int \mathbb{P}(D \mid \theta) \pi(\theta) \mathrm{d} \theta=\mathbb{P}(D)$.

## Rejection ABC

If $\mathbb{P}(D)$ is small (or $D$ continuous), we will rarely accept any $\theta$. Instead, there is an approximate version:

## Uniform Rejection Algorithm

- Draw $\theta$ from $\pi(\theta)$
- Simulate $X \sim f(\theta)$
- Accept $\theta$ if $\rho(D, X) \leq \epsilon$


## Rejection ABC

If $\mathbb{P}(D)$ is small（or $D$ continuous），we will rarely accept any $\theta$ ．Instead，there is an approximate version：

## Uniform Rejection Algorithm

－Draw $\theta$ from $\pi(\theta)$
－Simulate $X \sim f(\theta)$
－Accept $\theta$ if $\rho(D, X) \leq \epsilon$
$\epsilon$ reflects the tension between computability and accuracy．
－As $\epsilon \rightarrow \infty$ ，we get observations from the prior，$\pi(\theta)$ ．
－If $\epsilon=0$ ，we generate observations from $\pi(\theta \mid D)$ ．

## $\epsilon=10$

theta vs D


Density


$$
\begin{gathered}
\theta \sim U[-10,10], \quad X \sim N\left(2(\theta+2) \theta(\theta-2), 0.1+\theta^{2}\right) \\
\rho(D, X)=|D-X|, \quad D=2
\end{gathered}
$$

$$
\epsilon=7.5
$$

theta vs D


Density


$$
\begin{gathered}
\theta \sim U[-10,10], \quad X \sim N\left(2(\theta+2) \theta(\theta-2), 0.1+\theta^{2}\right) \\
\rho(D, X)=|D-X|, \quad D=2
\end{gathered}
$$

$$
\epsilon=5
$$

theta vs D


Density


$$
\begin{gathered}
\theta \sim U[-10,10], \quad X \sim N\left(2(\theta+2) \theta(\theta-2), 0.1+\theta^{2}\right) \\
\rho(D, X)=|D-X|, \quad D=2
\end{gathered}
$$

$$
\epsilon=2.5
$$

theta vs D


Density


$$
\begin{gathered}
\theta \sim U[-10,10], \quad X \sim N\left(2(\theta+2) \theta(\theta-2), 0.1+\theta^{2}\right) \\
\rho(D, X)=|D-X|, \quad D=2
\end{gathered}
$$

## $\epsilon=1$

theta vs D


Density


$$
\begin{gathered}
\theta \sim U[-10,10], \quad X \sim N\left(2(\theta+2) \theta(\theta-2), 0.1+\theta^{2}\right) \\
\rho(D, X)=|D-X|, \quad D=2
\end{gathered}
$$

## Rejection ABC

If the data are too high dimensional we never observe simulations that are 'close' to the field data - curse of dimensionality

Reduce the dimension using summary statistics, $S(D)$.

## Approximate Rejection Algorithm With Summaries

- Draw $\theta$ from $\pi(\theta)$
- Simulate $X \sim f(\theta)$
- Accept $\theta$ if $\rho(S(D), S(X))<\epsilon$

If $S$ is sufficient this is equivalent to the previous algorithm.

## Rejection ABC

If the data are too high dimensional we never observe simulations that are 'close' to the field data - curse of dimensionality

Reduce the dimension using summary statistics, $S(D)$.

## Approximate Rejection Algorithm With Summaries

- Draw $\theta$ from $\pi(\theta)$
- Simulate $X \sim f(\theta)$
- Accept $\theta$ if $\rho(S(D), S(X))<\epsilon$

If $S$ is sufficient this is equivalent to the previous algorithm.

## ABC as a probability model

## W. 2008/13

We wanted to solve the inverse problem

$$
D=f(\theta)
$$

but instead $A B C$ solves

$$
D=f(\theta)+e
$$

## ABC as a probability model

W. 2008/13

We wanted to solve the inverse problem

$$
D=f(\theta)
$$

but instead $A B C$ solves

$$
D=f(\theta)+e
$$

## $A B C$ gives 'exact' inference under a different model!

We can show that

## Proposition

If $\rho(D, X)=|D-X|$, then ABC samples from the posterior distribution of $\theta$ given $D$ where we assume $D=f(\theta)+e$ and that

$$
e \sim U[-\epsilon, \epsilon]
$$

We can generalise $A B C$ to assume non-uniform errors

## Key challenges for ABC (or perhaps for all inference)

## Scoring $\theta$

- The tolerance $\epsilon$, distance $\rho$, summary $S(D)$ (or variations thereof) determine the theoretical 'accuracy' of the approximation


## Key challenges for ABC (or perhaps for all inference)

Scoring $\theta$

- The tolerance $\epsilon$, distance $\rho$, summary $S(D)$ (or variations thereof) determine the theoretical 'accuracy' of the approximation
Computing acceptable $\theta$
- Computing the approximate posterior for any given score is usually hard.
- There is a trade-off between accuracy achievable in the approximation (size of $\epsilon$ ), and the information loss incurred when summarizing


## Key challenges for ABC (or perhaps for all inference)

Scoring $\theta$

- The tolerance $\epsilon$, distance $\rho$, summary $S(D)$ (or variations thereof) determine the theoretical 'accuracy' of the approximation
Computing acceptable $\theta$
- Computing the approximate posterior for any given score is usually hard.
- There is a trade-off between accuracy achievable in the approximation (size of $\epsilon$ ), and the information loss incurred when summarizing

Likelihood-free methods should generally be avoided when possible ${ }^{1}$
There are fewer and fewer likelihood-free problems.

[^0]
## Efficient Algorithms

## ABCifying Monte Carlo methods

Rejection $A B C$ is the basic $A B C$ algorithm

- Inefficient as it repeatedly samples from prior

More efficient sampling algorithms allow us to make better use of the available computational resource: spend more time in regions of parameter space likely to lead to accepted values.

- allows us to use smaller values of $\epsilon$

Most Monte Carlo algorithms now have ABC versions for when we don't know the likelihood: IS, MCMC, SMC $(\times n)$, EM, EP etc

## MCMC-ABC

Marjoram et al. 2003, Sisson and Fan 2011, Lee 2012
We are targeting the joint distribution

$$
\pi_{A B C}(\theta, x \mid D) \propto \mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta)
$$

## МСМС-АВС

Marjoram et al. 2003, Sisson and Fan 2011, Lee 2012
We are targeting the joint distribution

$$
\pi_{A B C}(\theta, x \mid D) \propto \mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta)
$$

To explore the $(\theta, x)$ space, proposals of the form

$$
Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)=q\left(\theta, \theta^{\prime}\right) \pi\left(x^{\prime} \mid \theta^{\prime}\right)
$$

seem to be inevitable. The Metropolis-Hastings (MH) acceptance probability is then

$$
r=\frac{\pi_{A B C}\left(\theta^{\prime}, x^{\prime} \mid D\right) Q\left(\left(\theta^{\prime}, x^{\prime}\right),(\theta, x)\right)}{\pi_{A B C}(\theta, x \mid D) Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)}
$$

## MCMC-ABC

## Marjoram et al. 2003, Sisson and Fan 2011, Lee 2012

We are targeting the joint distribution

$$
\pi_{A B C}(\theta, x \mid D) \propto \mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta)
$$

To explore the $(\theta, x)$ space, proposals of the form

$$
Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)=q\left(\theta, \theta^{\prime}\right) \pi\left(x^{\prime} \mid \theta^{\prime}\right)
$$

seem to be inevitable. The Metropolis-Hastings (MH) acceptance probability is then

$$
\begin{aligned}
r & =\frac{\pi_{A B C}\left(\theta^{\prime}, x^{\prime} \mid D\right) Q\left(\left(\theta^{\prime}, x^{\prime}\right),(\theta, x)\right)}{\pi_{A B C}(\theta, x \mid D) Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)} \\
& =\frac{\mathbb{I}_{\rho\left(D, x^{\prime}\right) \leq \epsilon} \pi\left(x^{\prime} \mid \theta^{\prime}\right) \pi\left(\theta^{\prime}\right) q\left(\theta^{\prime}, \theta\right) \pi(x \mid \theta)}{\mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta) q\left(\theta, \theta^{\prime}\right) \pi\left(x^{\prime} \mid \theta^{\prime}\right)}
\end{aligned}
$$

## MCMC-ABC

## Marjoram et al. 2003, Sisson and Fan 2011, Lee 2012

We are targeting the joint distribution

$$
\pi_{A B C}(\theta, x \mid D) \propto \mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta)
$$

To explore the $(\theta, x)$ space, proposals of the form

$$
Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)=q\left(\theta, \theta^{\prime}\right) \pi\left(x^{\prime} \mid \theta^{\prime}\right)
$$

seem to be inevitable. The Metropolis-Hastings (MH) acceptance probability is then

$$
\begin{aligned}
r & =\frac{\pi_{A B C}\left(\theta^{\prime}, x^{\prime} \mid D\right) Q\left(\left(\theta^{\prime}, x^{\prime}\right),(\theta, x)\right)}{\pi_{A B C}(\theta, x \mid D) Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)} \\
& =\frac{\mathbb{I}_{\rho\left(D, x^{\prime}\right) \leq \epsilon} \pi\left(x^{\prime} \mid \theta^{\prime}\right) \pi\left(\theta^{\prime}\right) q\left(\theta^{\prime}, \theta\right) \pi(x \mid \theta)}{\mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta) q\left(\theta, \theta^{\prime}\right) \pi\left(x^{\prime} \mid \theta^{\prime}\right)} \\
& =\frac{\mathbb{I}_{\rho\left(D, x^{\prime}\right) \leq \epsilon} q\left(\theta^{\prime}, \theta\right) \pi\left(\theta^{\prime}\right)}{\mathbb{I}_{\rho(D, x) \leq \epsilon} q\left(\theta, \theta^{\prime}\right) \pi(\theta)}
\end{aligned}
$$

## MCMC-ABC

## Marjoram et al. 2003, Sisson and Fan 2011, Lee 2012

We are targeting the joint distribution

$$
\pi_{A B C}(\theta, x \mid D) \propto \mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta)
$$

To explore the $(\theta, x)$ space, proposals of the form

$$
Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)=q\left(\theta, \theta^{\prime}\right) \pi\left(x^{\prime} \mid \theta^{\prime}\right)
$$

seem to be inevitable. The Metropolis-Hastings (MH) acceptance probability is then

$$
\begin{aligned}
r & =\frac{\pi_{A B C}\left(\theta^{\prime}, x^{\prime} \mid D\right) Q\left(\left(\theta^{\prime}, x^{\prime}\right),(\theta, x)\right)}{\pi_{A B C}(\theta, x \mid D) Q\left((\theta, x),\left(\theta^{\prime}, x^{\prime}\right)\right)} \\
& =\frac{\mathbb{I}_{\rho\left(D, x^{\prime}\right) \leq \epsilon} \pi\left(x^{\prime} \mid \theta^{\prime}\right) \pi\left(\theta^{\prime}\right) q\left(\theta^{\prime}, \theta\right) \pi(x \mid \theta)}{\mathbb{I}_{\rho(D, x) \leq \epsilon} \pi(x \mid \theta) \pi(\theta) q\left(\theta, \theta^{\prime}\right) \pi\left(x^{\prime} \mid \theta^{\prime}\right)} \\
& =\frac{\mathbb{I}_{\rho\left(D, x^{\prime}\right) \leq \epsilon} q\left(\theta^{\prime}, \theta\right) \pi\left(\theta^{\prime}\right)}{\mathbb{I}_{\rho(D, x) \leq \epsilon} q\left(\theta, \theta^{\prime}\right) \pi(\theta)}
\end{aligned}
$$

NB: HMC is not possible (w/o a surrogate)

## Sequential $A B C$ algorithms

Sisson et al. 2007, Toni et al. 2008, Beaumont et al. 2009, Del Moral et al. 2011, Drovandi et al. 2011, ... Choose a sequence of tolerances $\epsilon_{1}>\epsilon_{2}>\ldots>\epsilon_{T}$ and let $\pi_{t}$ be the ABC approximation when using tolerance $\epsilon_{t}$.
We aim to sample $N$ particles successively from

$$
\pi_{1}(\theta), \ldots, \pi_{T}(\theta)=\text { target }
$$



At each stage $t$, we aim to construct a weighted sample of particles that approximates $\pi_{t}(\theta, x)$.


Picture from Toni and Stumpf 2010

## Model selection

W. 2007, Grelaud et al. 2009

Often we want to compare models $\rightarrow$ Bayes factors

$$
B_{12}=\frac{\pi\left(D \mid M_{1}\right)}{\pi\left(D \mid M_{2}\right)}
$$

where $\pi\left(D \mid M_{i}\right)=\int \mathbb{I}_{\rho(D, X) \leq \epsilon} \pi\left(x \mid \theta, M_{i}\right) \pi(\theta) \mathrm{d} x \mathrm{~d} \theta$.

## Model selection

W. 2007, Grelaud et al. 2009

Often we want to compare models $\rightarrow$ Bayes factors

$$
B_{12}=\frac{\pi\left(D \mid M_{1}\right)}{\pi\left(D \mid M_{2}\right)}
$$

where $\pi\left(D \mid M_{i}\right)=\int \mathbb{I}_{\rho(D, X) \leq \epsilon} \pi\left(x \mid \theta, M_{i}\right) \pi(\theta) \mathrm{d} x \mathrm{~d} \theta$.
For rejection $A B C$

$$
\pi(D \mid M) \approx \frac{1}{N} \sum \mathbb{I}_{\rho\left(D, X_{i}\right) \leq \epsilon}
$$

where $X_{i} \sim M\left(\theta_{i}\right)$ with $\theta_{i} \sim \pi(\theta)$.

## Summary Statistics

## References:

- Blum, Nunes, Prangle and Sisson 2012
- Joyce and Marjoram 2008
- Nunes and Balding 2010
- Fearnhead and Prangle 2012
- Robert et al. 2011


## Choosing summary statistics

Blum, Nunes, Prangle, Fearnhead 2012

If $S(D)=s_{o b s}$ is sufficient for $\theta$, i.e., $s_{o b s}$ contains all the information contained in $D$ about $\theta$

$$
\pi\left(\theta \mid s_{o b s}\right)=\pi(\theta \mid D)
$$

then using summaries has no detrimental effect

## Choosing summary statistics

Blum, Nunes, Prangle, Fearnhead 2012

If $S(D)=s_{\text {obs }}$ is sufficient for $\theta$, i.e., $s_{\text {obs }}$ contains all the information contained in $D$ about $\theta$

$$
\pi\left(\theta \mid s_{o b s}\right)=\pi(\theta \mid D)
$$

then using summaries has no detrimental effect
However, low-dimensional sufficient statistics are rarely available. How do we choose good low dimensional summaries?

## Error trade-off

Fearnhead and Prangle 2012
The error in the $A B C$ approximation can be broken into two parts
(1) Choice of summary:

$$
\pi(\theta \mid D) \stackrel{?}{\approx} \pi\left(\theta \mid s_{o b s}\right)
$$

## Error trade-off

Fearnhead and Prangle 2012
The error in the $A B C$ approximation can be broken into two parts
(1) Choice of summary:

$$
\pi(\theta \mid D) \stackrel{?}{\approx} \pi\left(\theta \mid s_{o b s}\right)
$$

(2) Use of ABC acceptance kernel:

$$
\pi\left(\theta \mid s_{o b s}\right) \stackrel{?}{\approx} \pi_{A B C}\left(\theta \mid s_{o b s}\right)
$$

## Error trade-off

## Fearnhead and Prangle 2012

The error in the $A B C$ approximation can be broken into two parts
(1) Choice of summary:

$$
\pi(\theta \mid D) \stackrel{?}{\approx} \pi\left(\theta \mid s_{o b s}\right)
$$

(2) Use of ABC acceptance kernel:

$$
\pi\left(\theta \mid s_{o b s}\right) \stackrel{?}{\approx} \pi_{A B C}\left(\theta \mid s_{o b s}\right)
$$

The first approximation allows the matching between $S(D)$ and $S(X)$ to be done in a lower dimension. There is a trade-off

- $\operatorname{dim}(S)$ small: $\pi\left(\theta \mid s_{o b s}\right) \approx \pi_{A B C}\left(\theta \mid s_{o b s}\right)$, but $\pi\left(\theta \mid s_{o b s}\right) \not \approx \pi(\theta \mid D)$
- $\operatorname{dim}(S)$ large: $\pi\left(\theta \mid s_{o b s}\right) \approx \pi(\theta \mid D)$ but $\pi\left(\theta \mid s_{o b s}\right) \not \approx \pi_{A B C}\left(\theta \mid s_{o b s}\right)$ as curse of dimensionality forces us to use larger $\epsilon$


## Error trade-off

## Fearnhead and Prangle 2012

The error in the $A B C$ approximation can be broken into two parts
(1) Choice of summary:

$$
\pi(\theta \mid D) \stackrel{?}{\approx} \pi\left(\theta \mid s_{o b s}\right)
$$

(2) Use of ABC acceptance kernel:

$$
\pi\left(\theta \mid s_{o b s}\right) \stackrel{?}{\approx} \pi_{A B C}\left(\theta \mid s_{o b s}\right)
$$

The first approximation allows the matching between $S(D)$ and $S(X)$ to be done in a lower dimension. There is a trade-off

- $\operatorname{dim}(S)$ small: $\pi\left(\theta \mid s_{o b s}\right) \approx \pi_{A B C}\left(\theta \mid s_{o b s}\right)$, but $\pi\left(\theta \mid s_{o b s}\right) \not \approx \pi(\theta \mid D)$
- $\operatorname{dim}(S)$ large: $\pi\left(\theta \mid s_{o b s}\right) \approx \pi(\theta \mid D)$ but $\pi\left(\theta \mid s_{o b s}\right) \not \approx \pi_{A B C}\left(\theta \mid s_{o b s}\right)$ as curse of dimensionality forces us to use larger $\epsilon$
Optimal (in some sense) to choose $\operatorname{dim}(s)=\operatorname{dim}(\theta)$


## Machine learning invasion

ML algorithms are good at classification, often better than humans.
ABC can be done via classification, albeit at the cost of abandoning the Bayesian interpretation.

## Machine learning invasion

ML algorithms are good at classification, often better than humans.
ABC can be done via classification, albeit at the cost of abandoning the Bayesian interpretation.
E.g. 1)Use random forests, (C)NNs etc to generate a summary
(1) Train a ML model, $m(X)$, to predict $\theta$ from $D$ using a large number of simulator runs $\left\{\theta_{i}, X_{i}\right\}$
(2) ABC then simulates $\theta$ from the prior and $X$ from the simulator, and accepts $\theta$ if $m(X) \approx m\left(D_{o b s}\right)$

## Machine learning invasion

ML algorithms are good at classification, often better than humans.
ABC can be done via classification, albeit at the cost of abandoning the Bayesian interpretation.
E.g. 1)Use random forests, (C)NNs etc to generate a summary
(1) Train a ML model, $m(X)$, to predict $\theta$ from $D$ using a large number of simulator runs $\left\{\theta_{i}, X_{i}\right\}$
(2) ABC then simulates $\theta$ from the prior and $X$ from the simulator, and accepts $\theta$ if $m(X) \approx m\left(D_{o b s}\right)$
E.g. 2) Generative Adversarial Networks (GANs) play a game between a generator and a discriminative classifier. The classifier tries to distinguish between data and simulation, and the generator tries to trick the classifier.

## Machine learning invasion

ML algorithms are good at classification, often better than humans.
ABC can be done via classification, albeit at the cost of abandoning the Bayesian interpretation.
E.g. 1)Use random forests, (C)NNs etc to generate a summary
(1) Train a ML model, $m(X)$, to predict $\theta$ from $D$ using a large number of simulator runs $\left\{\theta_{i}, X_{i}\right\}$
(2) ABC then simulates $\theta$ from the prior and $X$ from the simulator, and accepts $\theta$ if $m(X) \approx m\left(D_{o b s}\right)$
E.g. 2) Generative Adversarial Networks (GANs) play a game between a generator and a discriminative classifier. The classifier tries to distinguish between data and simulation, and the generator tries to trick the classifier.
E.g. 3) Park et al. 2016, . ., suggested using the kernel mean embedding of the distribution (MMD) in an RKHS - inference is then simply projection in the RKHS.

## Machine learning invasion

ML algorithms are good at classification, often better than humans.
ABC can be done via classification, albeit at the cost of abandoning the Bayesian interpretation.
E.g. 1)Use random forests, (C)NNs etc to generate a summary
(1) Train a ML model, $m(X)$, to predict $\theta$ from $D$ using a large number of simulator runs $\left\{\theta_{i}, X_{i}\right\}$
(2) ABC then simulates $\theta$ from the prior and $X$ from the simulator, and accepts $\theta$ if $m(X) \approx m\left(D_{o b s}\right)$
E.g. 2) Generative Adversarial Networks (GANs) play a game between a generator and a discriminative classifier. The classifier tries to distinguish between data and simulation, and the generator tries to trick the classifier.
E.g. 3) Park et al. 2016, ..., suggested using the kernel mean embedding of the distribution (MMD) in an RKHS - inference is then simply projection in the RKHS.

All work well in simulation studies where the model is well specified and there is a true $\theta \ldots$

- Warning: beware of all automated summary selection approaches if misspecified


## Inference for misspecified models



## Inference under discrepancy

How should we do inference if the model is imperfect?
Data generating process

$$
y \sim G
$$

Model (complex simulator, finite dimensional parameter)

$$
\mathcal{F}=\left\{F_{\theta}: \theta \in \Theta\right\}
$$



If $G=F_{\theta_{0}} \in \mathcal{F}$ then we know what to $\mathrm{do}^{2}$.
How should we proceed if

$$
G \notin \mathcal{F}
$$

- $G$

[^1]
## Maximum likelihood

Maximum likelihood estimator

$$
\hat{\theta}_{n}=\arg \max _{\theta} I(y \mid \theta)
$$

If $G=F_{\theta_{0}} \in \mathcal{F}$, then (under some conditions)

$$
\begin{aligned}
\hat{\theta}_{n} & \rightarrow \theta_{0} \text { almost surely as } n \rightarrow \infty \\
\sqrt{n}\left(\hat{\theta}_{n}-\theta_{0}\right) & \stackrel{d}{\Rightarrow} N\left(0, \mathcal{I}^{-1}\left(\theta_{0}\right)\right)
\end{aligned}
$$

Asymptotic consistency, efficiency, normality.

## Maximum likelihood

Maximum likelihood estimator

$$
\hat{\theta}_{n}=\arg \max _{\theta} I(y \mid \theta)
$$

If $G=F_{\theta_{0}} \in \mathcal{F}$, then (under some conditions)

$$
\begin{aligned}
\hat{\theta}_{n} & \rightarrow \theta_{0} \text { almost surely as } n \rightarrow \infty \\
\sqrt{n}\left(\hat{\theta}_{n}-\theta_{0}\right) & \stackrel{d}{\Rightarrow} N\left(0, \mathcal{I}^{-1}\left(\theta_{0}\right)\right)
\end{aligned}
$$

Asymptotic consistency, efficiency, normality. If $G \notin \mathcal{F}$

$$
\begin{aligned}
\hat{\theta}_{n} \rightarrow \theta^{*} & =\arg \min _{\theta} D_{K L}\left(G, F_{\theta}\right) \text { almost surely } \\
& =\arg \min _{\theta} \int \log \frac{d G}{d F_{\theta}} d G \\
\sqrt{n}\left(\hat{\theta}_{n}-\theta_{0}\right) & \stackrel{d}{\Rightarrow} N\left(0, V^{-1}\right)
\end{aligned}
$$



## Bayes

Bayesian posterior

$$
\pi(\theta \mid y) \propto \pi(y \mid \theta) \pi(\theta)
$$

If $G=F_{\theta_{0}} \in \mathcal{F}$

$$
\pi(\theta \mid y) \stackrel{d}{\Rightarrow} N\left(\theta_{0}, n^{-1} \mathcal{I}^{-1}\left(\theta_{0}\right)\right) \text { as } \quad n \rightarrow \infty
$$

Bernstein-von Mises theorem: we forget the prior, and get asymptotic concentration and normality (under some conditions).

## Bayes

Bayesian posterior

$$
\pi(\theta \mid y) \propto \pi(y \mid \theta) \pi(\theta)
$$

If $G=F_{\theta_{0}} \in \mathcal{F}$

$$
\pi(\theta \mid y) \stackrel{d}{\Rightarrow} N\left(\theta_{0}, n^{-1} \mathcal{I}^{-1}\left(\theta_{0}\right)\right) \text { as } \quad n \rightarrow \infty
$$

Bernstein-von Mises theorem: we forget the prior, and get asymptotic concentration and normality (under some conditions).

If $G \notin \mathcal{F}$, we still get asymptotic concentration (and possibly normality) but to $\theta^{*}$ (the pseudo-true value).
"there is no obvious meaning for Bayesian analysis in this case"

## Bayes

Bayesian posterior

$$
\pi(\theta \mid y) \propto \pi(y \mid \theta) \pi(\theta)
$$

If $G=F_{\theta_{0}} \in \mathcal{F}$

$$
\pi(\theta \mid y) \stackrel{d}{\Rightarrow} N\left(\theta_{0}, n^{-1} \mathcal{I}^{-1}\left(\theta_{0}\right)\right) \text { as } \quad n \rightarrow \infty
$$

Bernstein-von Mises theorem: we forget the prior, and get asymptotic concentration and normality (under some conditions).

If $G \notin \mathcal{F}$, we still get asymptotic concentration (and possibly normality) but to $\theta^{*}$ (the pseudo-true value).
"there is no obvious meaning for Bayesian analysis in this case"
Often with non-parametric models (eg GPs), we don't even get this convergence to the pseudo-true value due to lack of identifiability.

## An appealing idea: model the discrepancy

## Kennedy an O'Hagan 2001

Can we expand the class of models by adding a Gaussian process (GP) to our simulator?

If $f_{\theta}(x)$ is our simulator, $y$ the observation, then perhaps we can correct $f$ by modelling

$$
y=f_{\theta^{*}}(x)+\delta(x) \quad \text { where } \quad \delta \sim G P
$$



## An appealing idea：model the discrepancy

## Kennedy an O＇Hagan 2001

Can we expand the class of models by adding a Gaussian process（GP）to our simulator？

If $f_{\theta}(x)$ is our simulator，$y$ the observation，then perhaps we can correct $f$ by modelling

$$
y=f_{\theta^{*}}(x)+\delta(x) \quad \text { where } \quad \delta \sim G P
$$

This greatly expands $\mathcal{F}$ into a non－parametric world．


## An appealing, but flawed, idea

Kennedy and O'Hagan 2001, Brynjarsdottir and O'Hagan 2014
Simulator
Reality

$$
f_{\theta}(x)=\theta x
$$

$$
g(x)=\frac{\theta x}{1+\frac{x}{a}} \quad \theta=0.65, a=20
$$

Solid=model with true theta, dashed=truth


## An appealing, but flawed, idea

Bolting on a GP can correct your predictions, but won't necessarily fix your inference:

- No discrepancy:

$$
\begin{gathered}
y=f_{\theta}(x)+N\left(0, \sigma^{2}\right) \\
\theta \sim N(0,100), \sigma^{2} \sim \Gamma^{-1}(0.001,0.001)
\end{gathered}
$$

- GP discrepancy:

$$
\begin{aligned}
y=f_{\theta}(x) & +\delta(x)+N\left(0, \sigma^{2}\right) \\
\delta(\cdot) & \sim G P(\cdot, \cdot) \text { with objective priors }
\end{aligned}
$$




## Dangers of non-parametric model extensions

There are (at least) two problems with this approach:

- We may still find $G \notin \mathcal{F}$
- Identifiability


## Dangers of non-parametric model extensions

There are (at least) two problems with this approach:

- We may still find $G \notin \mathcal{F}$
- Identifiability
- A GP is an incredibly complex infinite dimensional model, which is not necessarily identified even asymptotically. The posterior can concentrate not on a point, but on some sub manifold of parameter space, and the projection of the prior on this space continues to impact the posterior even as more and more data are collected.
ie We never forget the prior, but the prior is too complex to understand


## Dangers of non-parametric model extensions

There are (at least) two problems with this approach:

- We may still find $G \notin \mathcal{F}$
- Identifiability
- A GP is an incredibly complex infinite dimensional model, which is not necessarily identified even asymptotically. The posterior can concentrate not on a point, but on some sub manifold of parameter space, and the projection of the prior on this space continues to impact the posterior even as more and more data are collected.
ie We never forget the prior, but the prior is too complex to understand
- Brynjarsdottir and O'Hagan 2014 try to model their way out of trouble with prior information - which is great if you have it.
- We can also have problems finding the true optima for the hyperparameters, even in 1d problems:


－We can also have problems finding the true optima for the hyperparameters，even in 1d problems：


－Wong et al 2017 impose identifiability（for $\delta$ and $\theta$ ）by giving up and identifying

$$
\theta^{*}=\arg \min _{\theta} \int\left(\zeta(x)-f_{\theta}(x)\right)^{2} d \pi(x)
$$

J．R．Statist．Soc．B（2017）
79，Part 2，pp．635－648

## A frequentist approach to computer model calibration

## History matching

ABC was proposed as a method of last resort, but there is evidence it works particularly well for mis-specified models.

## History matching

ABC was proposed as a method of last resort, but there is evidence it works particularly well for mis-specified models.
History matching was designed for inference in mis-specified models. It seeks to find a NROY set

$$
\mathcal{P}_{\theta}=\left\{\theta: S_{H M}\left(\hat{F}_{\theta}, y\right) \leq 3\right\}
$$

where

$$
S_{H M}\left(F_{\theta}, y\right)=\frac{\left|\mathbb{E}_{F_{\theta}}(Y)-y\right|}{\sqrt{\operatorname{Var}_{F_{\theta}}(Y)}}
$$

## History matching

ABC was proposed as a method of last resort, but there is evidence it works particularly well for mis-specified models.
History matching was designed for inference in mis-specified models. It seeks to find a NROY set

$$
\mathcal{P}_{\theta}=\left\{\theta: S_{H M}\left(\hat{F}_{\theta}, y\right) \leq 3\right\}
$$

where

$$
S_{H M}\left(F_{\theta}, y\right)=\frac{\left|\mathbb{E}_{F_{\theta}}(Y)-y\right|}{\sqrt{\operatorname{Var}_{F_{\theta}}(Y)}}
$$

ABC approximates the posterior as

$$
\pi_{\epsilon}(\theta) \propto \pi(\theta) \mathbb{E}\left(\mathbb{I}_{S\left(\hat{F}_{\theta}, y\right) \leq \epsilon}\right)
$$

for some choice of $S$ (typically $S\left(\hat{F}_{\theta}, y\right)=\rho\left(\eta(y), \eta\left(y^{\prime}\right)\right)$ where $y^{\prime} \sim F_{\theta}$ ) and $\epsilon$.

## History matching

ABC was proposed as a method of last resort, but there is evidence it works particularly well for mis-specified models.
History matching was designed for inference in mis-specified models. It seeks to find a NROY set

$$
\mathcal{P}_{\theta}=\left\{\theta: S_{H M}\left(\hat{F}_{\theta}, y\right) \leq 3\right\}
$$

where

$$
S_{H M}\left(F_{\theta}, y\right)=\frac{\left|\mathbb{E}_{F_{\theta}}(Y)-y\right|}{\sqrt{\operatorname{Var}_{F_{\theta}}(Y)}}
$$

ABC approximates the posterior as

$$
\pi_{\epsilon}(\theta) \propto \pi(\theta) \mathbb{E}\left(\mathbb{I}_{S\left(\hat{F}_{\theta}, y\right) \leq \epsilon}\right)
$$

for some choice of $S$ (typically $S\left(\hat{F}_{\theta}, y\right)=\rho\left(\eta(y), \eta\left(y^{\prime}\right)\right)$ where $y^{\prime} \sim F_{\theta}$ ) and $\epsilon$.
They have thresholding of a score in common and are algorithmically comparable (thresholding).

## History matching and $A B C$

These methods (anecdotally) seem to work better in mis-specified situations.
Why?

## History matching and $A B C$

These methods (anecdotally) seem to work better in mis-specified situations.
Why?
They differ from likelihood based approaches in that

- They only use some aspect of the simulator output
- Typically we hand pick which simulator outputs to compare, and weight them on a case by case basis.
- Potentially use generalised scores/loss-functions
- The thresholding type nature potentially makes them somewhat conservative
- Bayes/Max-likelihood estimates usually concentrate asymptotically. If $G \notin \mathcal{F}$ can we hope to learn precisely about $\theta$ ?
- We should use methods that limit the amount of learning that is possible about $\theta$.


## History matching and $A B C$

These methods (anecdotally) seem to work better in mis-specified situations.
Why?
They differ from likelihood based approaches in that

- They only use some aspect of the simulator output
- Typically we hand pick which simulator outputs to compare, and weight them on a case by case basis.
- Potentially use generalised scores/loss-functions
- The thresholding type nature potentially makes them somewhat conservative
- Bayes/Max-likelihood estimates usually concentrate asymptotically. If $G \notin \mathcal{F}$ can we hope to learn precisely about $\theta$ ?
- We should use methods that limit the amount of learning that is possible about $\theta$.


## Conclusions

ABC allows inference in models for which it would otherwise be impossible.

- not a silver bullet - if likelihood methods possible, use them instead

Algorithms and post-hoc regression can greatly improve computational efficiency, but computation is still usually the limiting factor.

- Challenge is to develop more efficient methods to allow inference in more expensive models.

Scoring of parameter values needs careful thought

- Likelihood isn't always fit for purpose.


## Conclusions

ABC allows inference in models for which it would otherwise be impossible.

- not a silver bullet - if likelihood methods possible, use them instead

Algorithms and post-hoc regression can greatly improve computational efficiency, but computation is still usually the limiting factor.

- Challenge is to develop more efficient methods to allow inference in more expensive models.

Scoring of parameter values needs careful thought

- Likelihood isn't always fit for purpose.

Thank you for listening!


[^0]:    ${ }^{1}$ Unless your model is wrong...

[^1]:    ${ }^{2}$ Even if we can't agree about it!

