



POLITECNICO
MILANO 1863

Coupled Markov chains with applications to Approximate Bayesian Computation for model based clustering

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11 november 2021

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- find a way to speed up computation time by parallelizing a Monte Carlo Markov chain method;
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⇒ Unbiased Markov chain Monte Carlo methods with couplings;
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⇒ Approximate Bayesian Computation



Unbiased Markov chain Monte Carlo methods with couplings

The road to parallelization: coupling of Markov chains

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What is the coupling of Markov chains? The coupling of two probability distributions μ and ν refers to the construction of a bivariate probability distribution whose marginals are the original distributions μ and ν . Markov chain coupling allow to **reduce the convergence time**.

Rhee–Glynn estimator I

The goal is to estimate

$$\mathbb{E}_{\pi}[h(X)] = \int h(x)\pi(\mathrm{d}x).$$

The estimator we are going to construct is based on a coupled pair of Markov chains, $(X_t)_{t \geq 0}$ and $(Y_t)_{t \geq 0}$, which marginally start from π_0 and evolve accordingly to P .

We consider some assumptions:

1 as $t \rightarrow \infty$,

$$\mathbb{E}[h(X_t)] \rightarrow \mathbb{E}_\pi[h(X)];$$

and there exists $\eta > 0$ and $D < \infty$ such that $\mathbb{E}[|h(X_t)|^{2+\eta}] \leq D$ for all $t \geq 0$;

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- ② the chains are such that the meeting time

$$\tau = \inf\{t \geq 1 : X_t = Y_{t-1}\}$$

satisfies $\mathbb{P}(\tau > t) \leq C\delta^t$ for all $t \geq 0$, for some constants $C < \infty$ and $\delta \in (0, 1)$;

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- ③ the chains stay together after meeting:

$$X_t = Y_{t-1} \text{ for all } t \geq \tau.$$

Rhee–Glynn estimator II

Thanks to the previous assumptions we can prove that:

$$\mathbb{E}_\pi[h(X)] = \mathbb{E}[h(X_k) + \sum_{t=k+1}^{\tau-1} \{h(X_t) - h(Y_{t-1})\}];$$

and we define the Rhee–Glynn estimator as:

$$H_k(X, Y) = h(X_k) + \sum_{t=k+1}^{\tau-1} \{h(X_t) - h(Y_{t-1})\}$$

which is **unbiased** by construction.

Time-averaged estimator

In such form, Rhee-Glynn is not computationally feasible, the time-averaged estimator keeps the assumptions allowing the computation:

$$H_{k:m} = \frac{1}{m-k+1} \sum_{l=k}^m h(X_l) + \sum_{l=k+1}^{\tau-1} \min(1, \frac{l-k}{m-k+1}) \{h(X_l) - h(Y_{l-1})\}$$

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- $MCMC_{k:m}$ is the standard MCMC average;
- $BC_{k:m}$ is the bias correction;
- $k-1$ number of burn-in iterations;
- m is a fixed integer of number of maximum iterations;
- τ is a random variable representing the meeting time.

Application to Metropolis-Hasting algorithm I

The MH **algorithm** adapted with couplings:

- ① draw X_0 and Y_0 from an initial distribution π_0 and draw $X_1 \sim P(X_0, \cdot)$;
- ② set $t = 1$: while $t < \max\{m, \tau\}$ and:
 - a draw $(X_{t+1}, Y_t) \sim \bar{P}\{(X_t, Y_{t-1}), \cdot\}$;
 - b set $t \leftarrow t + 1$;
- ③ compute

$$H_{k:m}(X, Y)$$

with the time-averaged estimator.

Application to Metropolis-Hasting algorithm II

The following is the **algorithm** to calculate the coupled kernel

$\bar{P}\{(X_t, Y_{t-1}), \cdot\}$ via MH:

- 1 sample $(X^*, Y^*) | (X_t, Y_{t-1})$ from a maximal coupling of $q(X_t, \cdot)$ and $q(Y_{t-1}, \cdot)$;
- 2 sample $U \sim \mathcal{U}([0, 1])$;

3 if

$$U \leq \min \left\{ 1, \frac{\pi(X^*)q(X^*, X_t)}{\pi(X_t)q(X_t, X^*)} \right\}$$

then $X_{t+1} = X^*$; otherwise $X_t = X_{t-1}$;

4 if

$$U \leq \min \left\{ 1, \frac{\pi(Y^*)q(Y^*, Y_t)}{\pi(Y_t)q(Y_t, Y^*)} \right\}$$

then $Y_{t+1} = Y^*$; otherwise $Y_t = Y_{t-1}$.



Approximate Bayesian Computation

To solve this issue we can use methods based on the **approximation of the likelihood function**, called *Likelihood-free methods*. Here the **algorithm**: *Inputs*:

- a target posterior density $\pi(\theta|y_{obs}) \propto p(y_{obs}|\theta)\pi(\theta)$, consisting of a prior distribution $\pi(\theta)$ and a procedure of generating data under the model $p(y_{obs}|\theta)$;
- a proposal density $g(\theta)$, with $g(\theta) > 0$ if $\pi(\theta|y_{obs}) > 0$;
- an integer $N > 0$.

Sampling for $i = 1, \dots, N$:

- ① generate $\theta^{(i)} \sim g(\theta)$ from sampling density g ;
- ② generate $y \sim p(y|\theta^{(i)})$ from the likelihood;
- ③ if $y = y_{obs}$, then accept $\theta^{(i)}$ with probability $\frac{\pi(\theta^{(i)})}{Kg(\theta^{(i)})}$, where
 $K \geq \max_{\theta} \frac{\pi(\theta)}{g(\theta)}$; else go to 1.

Output:

- a set of parameter vectors $\theta^{(1)}, \dots, \theta^{(N)}$ which are samples from $\pi(\theta|y_{obs})$.

Is this an efficient method for complex analysis?

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3 If $\|y - y_{obs}\| \leq h$, then accept $\theta^{(i)}$ with probability $\frac{\pi(\theta^{(i)})}{Kg(\theta^{(i)})}$, where

$$K \geq \max_{\theta} \frac{\pi(\theta)}{g(\theta)}.$$

Else go to 1.

Approximate Bayesian Computation

We focused on a particular case of the Likelihood-free methods: the *Approximate Bayesian Computation (ABC)*.

The aim: find a practical way of performing Bayesian analysis, while keeping the approximation and the computation to a minimum.

The likelihood-free rejection algorithm is sampling from the joint distribution

$$\propto \mathbb{I}(\|y - y_{obs}\| \leq h) p(y|\theta) \pi(\theta)$$

\implies replace the indicator function with a standard smoothing kernel function $K_h(u)$, with $u = \|y - y_{obs}\|$:

$$K_h(u) = \frac{1}{h} K\left(\frac{u}{h}\right)$$

Hence:

$$\pi_{ABC}(\theta, y|y_{obs}) \propto K_h(u) p(y|\theta) \pi(\theta)$$

Is this feasible in practice?

In practice: difficult to have $y \approx y_{obs}$ from $p(y|\theta)$, unless y_{obs} very low dimensional or $p(y|\theta)$ factorises into low-dimensional components.

Thus we should use a large h , obtaining a poor posterior approximation!

\implies use summary statistics $s = S(y)$

$$\pi_{ABC}(\theta|s_{obs})$$

Critical decision: choice of summary statistics

Dimension of summary statistics:

- large enough to contain as much as information about observed data as possible
- low enough to avoid curse of dimensionality of matching s and s_{obs}

⇒ choose sufficient statistics, such that:

$$\pi(\theta | s_{obs}) \equiv \pi(\theta | y_{obs})$$

Distance measure: substantial impact on ABC algorithm efficiency

$$\| \mathbf{s} - \mathbf{s}_{obs} \| = (\mathbf{s} - \mathbf{s}_{obs})^\top \Sigma^{-1} (\mathbf{s} - \mathbf{s}_{obs})$$

- $\Sigma = \text{identity matrix} \rightarrow \text{Euclidean distance}$
- $\Sigma = \text{diagonal matrix of non-zero weights} \rightarrow \text{Weighted Euclidean distance}$
- $\Sigma = \text{full covariance matrix of } \mathbf{s} \rightarrow \text{Mahalanobis distance}$

The following is an ABC **algorithm**: *Inputs*:

- a target posterior density $\pi(\theta|y_{obs}) \propto p(y_{obs}|\theta)\pi(\theta)$, consisting of a prior distribution $\pi(\theta)$ and a procedure of generating data under the model $p(y_{obs}|\theta)$;
- a proposal density $g(\theta)$, with $g(\theta) > 0$ if $\pi(\theta|y_{obs}) > 0$;
- an integer $N > 0$;
- a kernel function $K_h(u)$ and a scale parameter $h > 0$;
- a low dimensional vector of summary statistics $s = S(y)$.

ABC rejection sampling Algorithm

Sampling for $i = 1, \dots, N$:

- ① generate $\theta^{(i)} \sim g(\theta)$ from sampling density g ;
- ② generate $y \sim p(y|\theta^{(i)})$ from the likelihood;
- ③ compute summary statistic $s = S(y)$;
- ④ accept $\theta^{(i)}$ with probability $\frac{K_h(\|s - s_{obs}\|)\pi(\theta^{(i)})}{Kg(\theta^{(i)})}$, where
 $K \geq K_h(0) \max_{\theta} \frac{\pi(\theta)}{g(\theta)}$; else go to 1.

Output:

- a set of parameter vectors $\theta^{(1)}, \dots, \theta^{(N)} \sim \pi_{ABC}(\theta|S_{obs})$.

Possibly, can add a stopping rule.



Conclusions

Our focus till now was to understand the fundamental concepts and collect the missing information.

The next step will be a **simple and separate implementation** of both solution to be tested on simulated data.

Further steps will consider the **integration** of both solution into a single implementation and the testing on more complex data.

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