# Global consensus Monte Carlo

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#### THEORY

## Introduction

For problems involving large data sets, it may be convenient or necessary to partition the data across multiple cores or machines.

Consider a target density given by

$$\pi(z) \propto \mu(z) \prod_{j=1}^b f_j(z)$$

where  $f_j$  is computable on the jth machine, and depends on  $y_j$ , the jth subset of the data. We wish to generate samples distributed according to  $\pi$ .

Existing approaches include:

- Scott et al. (2016), who propose running one MCMC chain on each computing node, with target densities proportional to  $\mu(z)^{1/b}f_j(z)$ . The samples are combined in a way that implicitly assumes approximate Gaussianity.
- Xu et al. (2014), who approximate each  $f_j$  by a density from an exponential family.

We propose a novel procedure, motivated by the global variable consensus optimisation algorithm of Boyd et al. (2011), itself based upon ideas of Bertsekas and Tsitsiklis (1989).

#### The instrumental model

We introduce an instrumental hierarchical model by associating a local 'proxy' variable  $x_j$  with each subset of the data, and introducing a top-level parameter  $\lambda$  (see DAG, above right).

Specifically, on an extended state space we define

$$\pi_{\lambda}(z,x_{1:b}) \propto \mu(z) \prod_{j=1}^{b} K_{\lambda}(z,x_{j}) f_{j}(x_{j}),$$

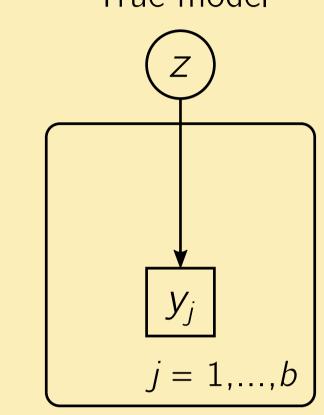
where  $\{K_{\lambda}: \lambda > 0\}$  is a family of Markov transition densities.

Assume that  $f_j$  is bounded, and that  $\int K_{\lambda}(z,x)f_j(x)dx \to f_j(z)$  pointwise as  $\lambda \to 0$ . Then the z-marginal of  $\pi_{\lambda}$  converges in total variation to  $\pi$ , so that for bounded functions  $\varphi$ ,

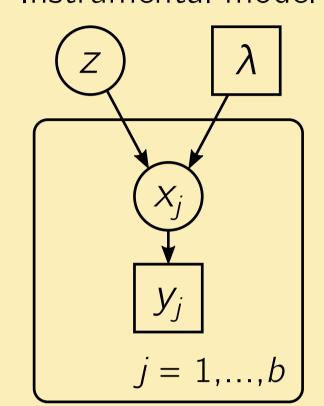
$$\int arphi(z)\pi_\lambda(z) \mathrm{d}z o \int arphi(z)\pi(z) \mathrm{d}z.$$

We look to generate samples distributed according to the artificial joint distribution  $\pi_{\lambda}$ ; for  $\lambda$  sufficiently close to 0, the z-marginal forms an approximation to the true target distribution.

#### True model



#### Instrumental model



## Gibbs sampler

For given  $\lambda$ , a  $\pi_{\lambda}$ -reversible Markov chain is obtained by considering the full conditional densities:

$$egin{align} \pi_{\lambda}(z\mid x_{1:b}) &\propto \mu(z) \prod_{j=1}^b \mathsf{K}_{\lambda}(z,x_j), \ \pi_{\lambda}(x_i\mid z) &\propto \mathsf{K}_{\lambda}(z,x_i) f_i(x_i). \end{aligned}$$

A two-variable Gibbs sampler may be constructed, where the two variables are z and  $x_{1:b}$ . In distributed settings, the drawing of a new  $x'_j \sim \pi_{\lambda}(\cdot \mid z)$  may occur on the jth computing node; the new values  $x'_{1:b}$  may then be sent to a central node, in order to

## **SMC** sampler

draw a new  $z' \sim \pi_{\lambda}(\cdot \mid x'_{1:h})$ .

The parameter  $\lambda$  may be chosen to balance computational tractability with fidelity to the true model, in a form of bias—variance tradeoff. When  $\lambda$  is close to 0, MCMC chains targeting  $\pi_{\lambda}(z, x_{1:b})$  may mix poorly.

To this end, one may approximate a sequence of distributions  $\pi_{\lambda_0}, \pi_{\lambda_1}, \ldots$  using an SMC sampler. By choosing a decreasing sequence of values  $\lambda_p$ , such an approach is observed to produce low-variance estimators for small values of  $\lambda$ .

Work is ongoing on a procedure to specify  $\lambda$  in an automated manner, by adaptively specifying this sequence, and using SMC variance estimators to assess the performance of each  $\lambda$  value.

#### **► EXAMPLES**

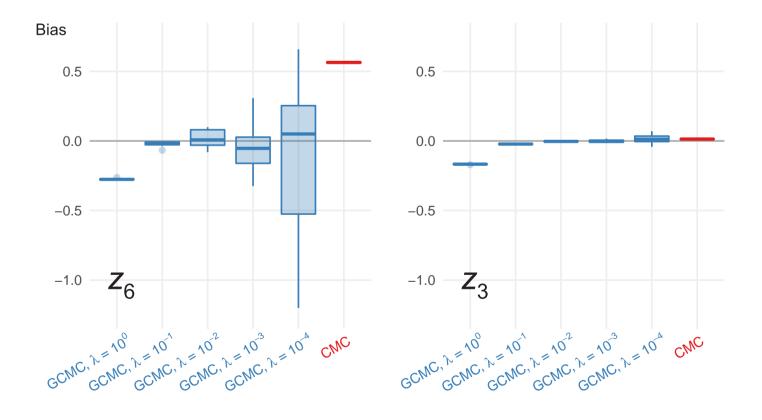
# Binary regression

Binary logistic regression models are commonly used with A/B testing data — in web design for example, to determine which content choices lead to maximised user interaction (such as the user clicking on a link to a product for sale).

- Data set formed of responses  $\eta_i \in \{-1,1\}$  and vectors  $\xi_i$  of binary covariates (originally belonging to  $\{0,1\}^d$ , but centred in a pre-processing step). The data are split into b subsets;  $f_j(z) = \prod_i \sigma(\eta_i z^T \xi_i)$ , where the product is taken over those indices i included in the jth data subset, and  $\sigma$  is the logistic function.
- We use a zero-mean normal prior  $\mu$ .
- For our algorithm, we use normal transition kernels:  $K_{\lambda}(z,x) = \mathcal{N}(x;z,\lambda)$ .

We compare our algorithm (GCMC) with the consensus Monte Carlo algorithm (CMC) proposed by Scott et al. (2016). We aim to estimate the posterior mean  $\int z\pi(z)dz$ .

We demonstrate on a simple data set with d=6 covariates, split into b=8 subsets, each comprising 512 data. We use various choices of  $\lambda$  for GCMC.



These boxplots show the biases of estimates of the posterior mean of two components, as obtained from 10 runs of each setup.

When estimating some parameters, CMC can introduce a far larger bias than GCMC, as the latter is more robust to deviations from Gaussianity, though must  $\lambda$  be chosen appropriately to control the variance. In the case above left, the corresponding covariate is rarely observed in some of the data subsets, so the  $f_j$  are skewed in this dimension, and are poorly approximated by Gaussians.

For parameters corresponding to more frequently-observed covariates, GCMC performs comparably to CMC if  $\lambda$  is chosen suitably (above right).

# Log-normal toy model

Let  $\mathcal{LN}(x; \mu, \sigma^2)$  denote the density at x of a log-normal distribution with parameters  $(\mu, \sigma^2)$ .

- Let  $\mu(z) = \mathcal{LN}(z; \mu_0, \sigma_0^2)$ .
- Let  $f_j(z) = \mathcal{LN}(y_j; \log(z), \sigma_j^2)$ .
- For our algorithm, we use log-normal transition kernels:  $K_{\lambda}(z,x) = \mathcal{LN}(x;\log(z),\lambda)$ .

Although the CMC algorithm of Scott et al. is exact for Gaussian models, this does not apply to reparametrisations, as is the case here. In GCMC, for small enough  $\lambda$  we would still expect the z-marginal of  $\pi_{\lambda}$  to be a good approximation of the true target.

We demonstrate on an example with b=32; in each case,  $10^6$  samples were used, following burn-in. Estimating the first moment  $\int z\pi(z)\mathrm{d}z\approx 1.131$ :

	Bias squared	Variance
GCMC, $\lambda = 1$	$1.32 \times 10^{-3}$	$4.66 \times 10^{-5}$
GCMC, $\lambda=10^{-1}$	$5.45 \times 10^{-5}$	$2.32 \times 10^{-4}$
GCMC, $\lambda = 10^{-2}$	$8.32 \times 10^{-5}$	$6.03 \times 10^{-4}$
GCMC, $\lambda=10^{-3}$	$1.52 \times 10^{-3}$	$1.26 \times 10^{-2}$
CMC	$2.37 \times 10^{-2}$	$1.41 \times 10^{-3}$

When estimating higher moments, GCMC results in far smaller biases than CMC, if a small enough  $\lambda$  is used. For example, estimating the fifth moment  $\int z^5 \pi(z) dz \approx 3.445$ :

	Bias squared	Variance
GCMC, $\lambda = 1$	$1.53 \times 10^{+1}$	$6.57 \times 10^{-2}$
GCMC, $\lambda=10^{-1}$	$1.15 \times 10^{-1}$	$9.84 \times 10^{-2}$
GCMC, $\lambda=10^{-2}$	$7.34 \times 10^{-2}$	$2.07 \times 10^{-1}$
GCMC, $\lambda=10^{-3}$	$1.49 \times 10^{-1}$	$9.45 \times 10^{0}$
CMC	$3.49 \times 10^{+3}$	$3.49 \times 10^{+3}$

#### **▼** REFERENCES

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