

# Inference from Privatized Data

## Sampling-based and Variational-based Bayesian Computation

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<https://yifei-xiong.github.io/>

Joint work with Nianqiao Phyllis Ju (Dartmouth College)

# Why inference from privatized data?

- In many modern data-sharing settings, the raw confidential data cannot be released directly.
- Instead, a data curator releases only a *privatized* or *privacy-protected* summary.
- This already appears in settings such as **tech companies**, the **U.S. Census Bureau**, and **social science data releases**.

raw data → summary → add privacy noise → released value

## Central question

How can we perform inference when we only observe a noisy, privatized release rather than the raw data?

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# Differential privacy and the Laplace mechanism

Two datasets  $\mathbf{x}, \mathbf{x}' \in \mathbb{X}^n$  are **neighboring** if they differ in one individual record.

## $\epsilon$ -differential privacy (Dwork, 2006)

A mechanism  $\eta$  is  $\epsilon$ -DP if for any neighboring  $\mathbf{x}, \mathbf{x}'$  and any measurable set  $B$ ,

$$\mathbb{P}(\eta(\mathbf{x}) \in B) \leq e^\epsilon \mathbb{P}(\eta(\mathbf{x}') \in B).$$

For a query  $s : \mathbb{X}^n \rightarrow \mathbb{R}^d$ , define its  $\ell_1$ -sensitivity by

$$\Delta_1(s) := \sup_{d(\mathbf{x}, \mathbf{x}')=1} \|s(\mathbf{x}) - s(\mathbf{x}')\|_1.$$

## Laplace mechanism

Release

$$S_{\text{dp}} = s(\mathbf{X}) + Z, \quad Z_j \stackrel{\text{iid}}{\sim} \text{Laplace} \left( 0, \frac{\Delta_1(s)}{\epsilon} \right).$$

Then  $S_{\text{dp}}$  is  $\epsilon$ -differentially private.

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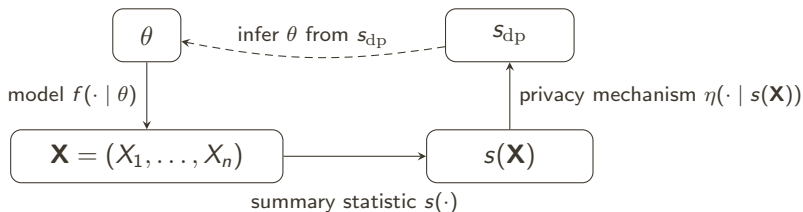
$$S_{\text{dp}} = s(\mathbf{X}) + Z, \quad Z_j \stackrel{\text{iid}}{\sim} \text{Laplace}\left(0, \frac{\Delta_1(s)}{\epsilon}\right).$$

Then  $S_{\text{dp}}$  is  $\epsilon$ -differentially private.

# A generative view of privatized inference

## Notations

- $\theta$ : model parameters
- $\mathbf{X} = (X_1, \dots, X_n)$ : confidential data
- $s(\mathbf{X})$ : summary statistics of  $\mathbf{X}$
- $s_{\text{dp}}$ : privatized query of  $s(\mathbf{X})$ .

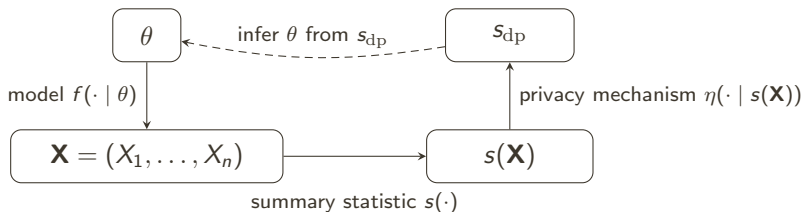


- Posterior:  $\pi(\theta | s_{\text{dp}}) \propto \pi(\theta) f(s_{\text{dp}} | \theta)$
- Likelihood:  $f(s_{\text{dp}} | \theta) = \int_{\mathbf{X}^n} f(\mathbf{x} | \theta) \eta(s_{\text{dp}} | s(\mathbf{x})) d\mathbf{x}$

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# Two computational viewpoints for inference from $s_{\text{dp}}$

Our target is the private posterior  $\pi(\theta | s_{\text{dp}}) \propto \pi(\theta) f(s_{\text{dp}} | \theta)$ .

## Sampling-based methods

Construct a Markov chain whose stationary distribution is

$$\pi(\theta | s_{\text{dp}}) \quad \text{or} \quad \pi(\theta, \mathbf{x} | s_{\text{dp}}),$$

and use posterior samples  $\theta^{(1)}, \dots, \theta^{(M)}$  to approximate expectations under the posterior.

- exact or asymptotically exact
- often computationally expensive

## Variational-based methods

Choose a tractable family

$$q_{\phi}(\theta | s_{\text{dp}})$$

and fit  $\phi$  so that

$$q_{\phi}(\theta | s_{\text{dp}}) \approx \pi(\theta | s_{\text{dp}})$$

as a distribution approximation.

- typically non-exact
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# Sampling-based idea: data augmentation

A natural strategy is to augment the unobserved confidential data

$$\mathbf{X} = (X_1, \dots, X_n)$$

and work with the joint posterior

$$\pi(\theta, \mathbf{x} \mid s_{\text{dp}}) \propto \pi(\theta) f(\mathbf{x} \mid \theta) \eta(s_{\text{dp}} \mid s(\mathbf{x})).$$

If the data model factorizes as

$$f(\mathbf{x} \mid \theta) = \prod_{i=1}^n f(x_i \mid \theta),$$

then

$$\pi(\theta, \mathbf{x} \mid s_{\text{dp}}) \propto \pi(\theta) \left[ \prod_{i=1}^n f(x_i \mid \theta) \right] \eta(s_{\text{dp}} \mid s(\mathbf{x})).$$

## Idea

If we can sample from  $\pi(\theta, \mathbf{x} \mid s_{\text{dp}})$ , then the  $\theta$ -marginal targets the posterior of interest  $\pi(\theta \mid s_{\text{dp}})$ .

# Gibbs update

The augmented posterior suggests a two-block Gibbs strategy.

## Inference step

$$\pi(\theta | \mathbf{x}, s_{\text{dp}}) = \pi(\theta | \mathbf{x}) \propto \pi(\theta) \prod_{i=1}^n f(x_i | \theta).$$

## Imputation step

$$\pi(\mathbf{x} | \theta, s_{\text{dp}}) \propto \left[ \prod_{i=1}^n f(x_i | \theta) \right] \eta(s_{\text{dp}} | s(\mathbf{x})).$$

- The **privacy mechanism** only enters through the imputation step.
- This is the computational bottleneck:  $\mathbf{x}$  is high-dimensional, and the privacy term couples all coordinates through  $s(\mathbf{x})$ .

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# Independent Metropolis-within-Gibbs sampler

- We want to sample from

$$\pi(\mathbf{x} \mid \theta, s_{\text{dp}}) \propto \prod_{i=1}^n f(x_i \mid \theta) \cdot \eta(s_{\text{dp}} \mid s(\mathbf{x})). \quad (1)$$

- Ju et al. (2022) target the DP imputation distribution using a **independent Metropolis-within-Gibbs** (IMwG) sampler.

## Algorithm: IMwG Update

Given current state  $\mathbf{x} = (x_1, \dots, x_n)$ :

- 1 Pick  $I \sim \text{Unif}\{1, \dots, n\}$ .
- 2 Propose  $y \sim q(\cdot)$ , independent of  $x_I$ . (We can choose  $q(\cdot) = f(\cdot \mid \theta)$ )
- 3 Form candidate  $\mathbf{x}' = [\mathbf{x}_{-I}, y]$ .
- 4 Accept with probability

$$\alpha_I^{\text{IMwG}}(y, \mathbf{x}) = 1 \wedge \frac{\pi(\mathbf{x}' \mid \theta, s_{\text{dp}}) q(x_I)}{\pi(\mathbf{x} \mid \theta, s_{\text{dp}}) q(y)}. \quad \left( = 1 \wedge \frac{\eta(s_{\text{dp}} \mid s(\mathbf{x}'))}{\eta(s_{\text{dp}} \mid s(\mathbf{x}))} \geq e^{-\epsilon} \right)$$

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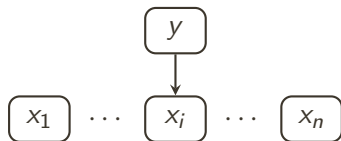
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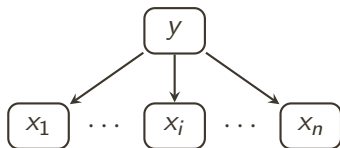
# Idea: reusing proposals

Independent Metropolis (IMwG)



one proposal  $\rightarrow$  one coordinate

SOMA

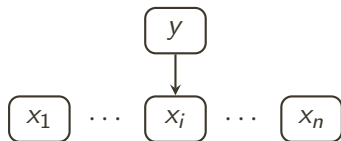


one proposal  $\rightarrow$  many coordinates

- In IMwG, each proposal  $y$  only gets *one chance* to replace a randomly chosen coordinate  $x_i$ .
- Our idea: keep **one** proposal  $y$ , but let it compete across **all** coordinates and move it where it helps the most.
- We let  $y$  “compete” across the slots and update the most promising coordinate. This is SOMA: **Single-Offer-Multiple-Attempts**.

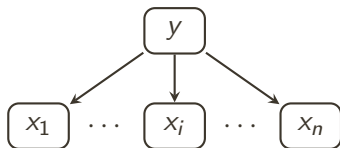
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# SOMA kernel (one iteration)

## One SOMA iteration targeting $\pi(\mathbf{x})$

❶ Propose a single value  $y \sim q(\cdot)$ .

❷ Compute weights

$$w_0(\mathbf{x}) = \frac{\pi(\mathbf{x})}{\prod_{j=1}^n q(x_j)}, \quad w_i(y, \mathbf{x}) = \frac{\pi([\mathbf{x}_{-i}, y])}{q(y) \prod_{j \neq i} q(x_j)},$$

and set  $W(y, \mathbf{x}) = \sum_{i=1}^n w_i(y, \mathbf{x})$ .

❸ Choose a coordinate

$$\mathbb{P}(I = i \mid y, \mathbf{x}) = \frac{w_i(y, \mathbf{x})}{W(y, \mathbf{x})}, \quad i = 1, \dots, n.$$

❹ Accept / reject replacement of  $x_I$  by  $y$  with probability

$$\alpha_I^{\text{SOMA}}(y, \mathbf{x}) = \min \left\{ 1, \frac{W(y, \mathbf{x})}{W(y, \mathbf{x}) + w_0(\mathbf{x}) - w_I(y, \mathbf{x})} \right\}.$$

# Why SOMA is valid and useful

## Reversibility

If the target  $\pi(\mathbf{x})$  is permutation invariant in  $(x_1, \dots, x_n)$ , then SOMA is reversible with respect to  $\pi(\mathbf{x})$ .

For the DP imputation target

$$\pi(\mathbf{x}) \propto \left[ \prod_{i=1}^n f(x_i | \theta) \right] \eta(s_{\text{dp}} | s(\mathbf{x})),$$

this holds whenever  $s(\mathbf{x})$  is permutation invariant.

## Advantage over IMwG

For any state  $\mathbf{x}$ , coordinate  $i$ , and proposal  $y$ ,

$$\alpha_i^{\text{SOMA}}(y, \mathbf{x}) \geq \alpha_i^{\text{IMwG}}(y, \mathbf{x}).$$

Under  $\epsilon$ -DP, the overall acceptance probability satisfied

$$A^{\text{SOMA}}(\mathbf{x}) \geq \frac{n}{n + e^\epsilon - 1}, \quad A^{\text{IMwG}}(\mathbf{x}) \geq e^{-\epsilon}.$$

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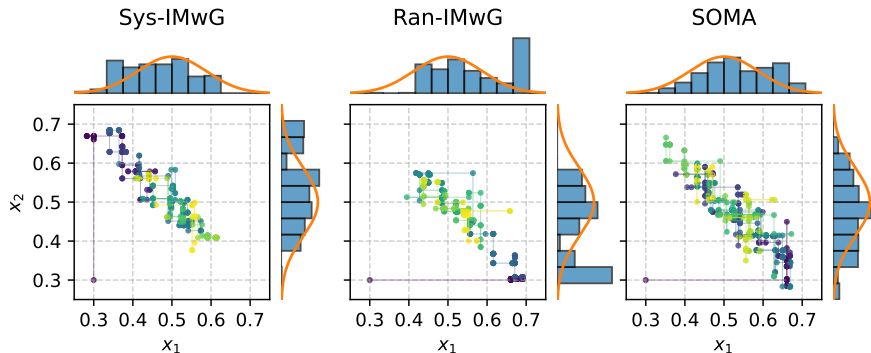
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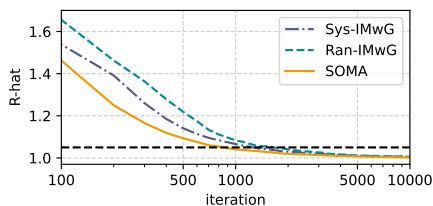
# Synthetic Example: Trace Plots ( $n = 2$ )



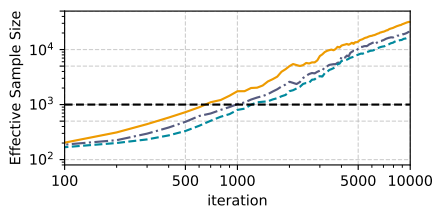
- All chains start from the same initial state  $(0.3, 0.3)$  over 500 iterations.
- SOMA (right) traverses the high-density region of  $\pi$  much more quickly and explores the state space more efficiently.
- Ran-IMwG tends to get stuck in local regions, leading to under-exploration of the posterior support.

# Convergence Diagnostics at $n = 100$

- Privatised Bayesian linear regression (Bernstein and Sheldon, 2019) with  $n = 100$ .
- The state space has dimension  $n \cdot (\rho + 2) + (\rho + 1) = 403$ ; we use standard MCMC diagnostics: Gelman–Rubin  $\hat{R}$  and effective sample size (ESS).



$\hat{R}$  across iterations (dashed line: 1.05)



Bulk ESS across iterations

- $\hat{R}$ : SOMA reaches  $\hat{R} < 1.05$  after about 800 iterations, while both Sys-IMwG and Ran-IMwG need around 1,500 iterations.
- ESS: SOMA yields substantially larger effective sample size than the two IMwG baselines at the same iteration budget.

# Coupling time and Acceptance rate

Table: Coupling time, convergence rate, acceptance rate for  $n = 10$  and  $\epsilon = 3$  and  $\epsilon = 30$ . All values are presented as mean with quartiles.

Method	Coupling Time	Convergence Rate	Acceptance Rate (%)
$n = 10, \epsilon = 3$			
Ran-IMwG	156.04 (46, 216)	0.9937 (0.9933, 0.9941)	94.80 (94.52, 95.11)
Sys-IMwG	<b>99.60 (28, 138)</b>	<b>0.9904 (0.9895, 0.9911)</b>	94.81 (94.55, 95.09)
SOMA	152.98 (42, 211)	0.9937 (0.9932, 0.9942)	<b>99.49 (99.44, 99.55)</b>
$n = 10, \epsilon = 30$			
Ran-IMwG	900.79 (261, 1262)	0.9989 (0.9988, 0.9990)	50.95 (50.64, 51.25)
Sys-IMwG	549.07 (170, 764)	0.9982 (0.9980, 0.9983)	50.94 (50.66, 51.23)
SOMA	<b>120.64 (46, 163)</b>	<b>0.9909 (0.9902, 0.9915)</b>	<b>91.91 (91.80, 92.01)</b>

# Sampling-based approach: takeaway

## What we gained

- Data augmentation turns inference from  $s_{\text{dp}}$  into sampling from

$$\pi(\theta, \mathbf{x} \mid s_{\text{dp}}).$$

- SOMA improves the imputation step by reusing one proposal across coordinates.
- Under DP, acceptance can be controlled explicitly.

## Limitation

Sampling-based methods are exact or asymptotically exact, but they still require iterative MCMC over a high-dimensional latent state.

This motivates a variational approximation to the private posterior.

# Why a variational approximation?

Recall the private posterior

$$\pi(\theta \mid s_{\text{dp}}) \propto \pi(\theta) f(s_{\text{dp}} \mid \theta), \quad f(s_{\text{dp}} \mid \theta) = \int_{\mathbb{X}^n} f(\mathbf{x} \mid \theta) \eta(s_{\text{dp}} \mid s(\mathbf{x})) d\mathbf{x}.$$

## Difficulty

Even if we can simulate from

$$\theta \sim \pi(\theta), \quad \mathbf{X} \sim f(\cdot \mid \theta), \quad s_{\text{dp}} \sim \eta(\cdot \mid s(\mathbf{X})),$$

the private likelihood  $f(s_{\text{dp}} \mid \theta)$  is usually intractable.

- Sampling-based methods target the posterior  $\pi(\theta \mid s_{\text{dp}})$  by MCMC.
- Here we instead learn an approximation  $q_\phi$  from simulated privatized data.

# Two approximation targets: posterior or likelihood

There are two natural objects to approximate.

Privatized posterior estimation  
(PPE)

Approximate

$$\pi(\theta \mid s_{\text{dp}}) \quad \text{by} \quad q_{\phi}(\theta \mid s_{\text{dp}}).$$

Privatized likelihood estimation  
(PLE)

Approximate

$$f(s_{\text{dp}} \mid \theta) \quad \text{by} \quad q_{\phi}(s_{\text{dp}} \mid \theta).$$

- In practice,  $q_{\phi}$  can be any flexible conditional density model, e.g., normalizing flows (Durkan et al., 2019) or diffusion models.
- We do not need its exact architecture here; what matters is the loss function it minimizes.

# PLE: start from a KL objective

We want to approximate the private likelihood

$$f(s_{\text{dp}} | \theta) \quad \text{by} \quad q_{\phi}(s_{\text{dp}} | \theta).$$

A natural population loss is the prior-averaged KL divergence:

$$\ell_{\text{PLE}}(\phi) = \mathbb{E}_{\pi(\theta)} [\text{KL}(f(s_{\text{dp}} | \theta) \| q_{\phi}(s_{\text{dp}} | \theta))].$$

Expanding the KL divergence,

$$\begin{aligned} \ell_{\text{PLE}}(\phi) &= \int \pi(\theta) \int f(s_{\text{dp}} | \theta) \log \frac{f(s_{\text{dp}} | \theta)}{q_{\phi}(s_{\text{dp}} | \theta)} ds_{\text{dp}} d\theta \\ &= \int \pi(\theta) \int f(s_{\text{dp}} | \theta) \log f(s_{\text{dp}} | \theta) ds_{\text{dp}} d\theta \\ &\quad - \int \pi(\theta) \int f(s_{\text{dp}} | \theta) \log q_{\phi}(s_{\text{dp}} | \theta) ds_{\text{dp}} d\theta. \end{aligned}$$

The first term does not depend on  $\phi$ .

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# PLE: from KL to a trainable loss

The first term does not depend on  $\phi$ . Hence

$$\ell_{\text{PLE}}(\phi) \equiv - \int \pi(\theta) \int f(s_{\text{dp}} | \theta) \log q_{\phi}(s_{\text{dp}} | \theta) ds_{\text{dp}} d\theta.$$

Now use

$$f(s_{\text{dp}} | \theta) = \int f(\mathbf{x} | \theta) \eta(s_{\text{dp}} | s(\mathbf{x})) d\mathbf{x}.$$

Substituting this into the loss,

$$\begin{aligned} \ell_{\text{PLE}}(\phi) &\equiv - \int \pi(\theta) \int \left[ \int f(\mathbf{x} | \theta) \eta(s_{\text{dp}} | s(\mathbf{x})) d\mathbf{x} \right] \log q_{\phi}(s_{\text{dp}} | \theta) ds_{\text{dp}} d\theta \\ &= \mathbb{E}_{p(\theta, \mathbf{x})} \left[ - \int \eta(s_{\text{dp}} | s(\mathbf{x})) \log q_{\phi}(s_{\text{dp}} | \theta) ds_{\text{dp}} \right]. \end{aligned}$$

So the trainable loss is

$$\ell_{\text{PLE}}(\phi) = \mathbb{E}_{p(\theta, \mathbf{x})} \left[ - \int \eta(s_{\text{dp}} | s(\mathbf{x})) \log q_{\phi}(s_{\text{dp}} | \theta) ds_{\text{dp}} \right].$$

# PPE: start from a KL objective

We now approximate the private posterior

$$\pi(\theta \mid s_{\text{dp}}) \quad \text{by} \quad q_\phi(\theta \mid s_{\text{dp}}).$$

A natural population loss is the average conditional KL divergence:

$$\ell_{\text{PPE}}(\phi) = \mathbb{E}_{p(s_{\text{dp}})} [\text{KL}(\pi(\theta \mid s_{\text{dp}}) \parallel q_\phi(\theta \mid s_{\text{dp}}))].$$

Expanding the KL divergence,

$$\begin{aligned} \ell_{\text{PPE}}(\phi) &= \int p(s_{\text{dp}}) \int \pi(\theta \mid s_{\text{dp}}) \log \frac{\pi(\theta \mid s_{\text{dp}})}{q_\phi(\theta \mid s_{\text{dp}})} d\theta ds_{\text{dp}} \\ &= \int p(s_{\text{dp}}) \int \pi(\theta \mid s_{\text{dp}}) \log \pi(\theta \mid s_{\text{dp}}) d\theta ds_{\text{dp}} \\ &\quad - \int p(s_{\text{dp}}) \int \pi(\theta \mid s_{\text{dp}}) \log q_\phi(\theta \mid s_{\text{dp}}) d\theta ds_{\text{dp}}. \end{aligned}$$

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$$\ell_{\text{PPE}}(\phi) = - \int p(s_{\text{dp}}) \int \pi(\theta | s_{\text{dp}}) \log q_{\phi}(\theta | s_{\text{dp}}) d\theta ds_{\text{dp}}.$$

Using the joint density

$$p(\theta, s_{\text{dp}}) = p(s_{\text{dp}})\pi(\theta | s_{\text{dp}}),$$

we rewrite this as

$$\ell_{\text{PPE}}(\phi) = - \int p(\theta, s_{\text{dp}}) \log q_{\phi}(\theta | s_{\text{dp}}) d\theta ds_{\text{dp}}.$$

Since

$$p(\theta, s_{\text{dp}}) = \int \pi(\theta) f(\mathbf{x} | \theta) \eta(s_{\text{dp}} | s(\mathbf{x})) d\mathbf{x},$$

we obtain

$$\ell_{\text{PPE}}(\phi) = \mathbb{E}_{p(\theta, \mathbf{x})} \left[ - \int \eta(s_{\text{dp}} | s(\mathbf{x})) \log q_{\phi}(\theta | s_{\text{dp}}) ds_{\text{dp}} \right].$$

So PPE learns the posterior directly, while PLE learns the private likelihood.

# PPE: from KL to a trainable loss

Again, the first term does not depend on  $\phi$ . Hence

$$\ell_{\text{PPE}}(\phi) = - \int p(s_{\text{dp}}) \int \pi(\theta | s_{\text{dp}}) \log q_{\phi}(\theta | s_{\text{dp}}) d\theta ds_{\text{dp}}.$$

Using the joint density

$$p(\theta, s_{\text{dp}}) = p(s_{\text{dp}})\pi(\theta | s_{\text{dp}}),$$

we rewrite this as

$$\ell_{\text{PPE}}(\phi) = - \int p(\theta, s_{\text{dp}}) \log q_{\phi}(\theta | s_{\text{dp}}) d\theta ds_{\text{dp}}.$$

Since

$$p(\theta, s_{\text{dp}}) = \int \pi(\theta) f(\mathbf{x} | \theta) \eta(s_{\text{dp}} | s(\mathbf{x})) d\mathbf{x},$$

we obtain

$$\ell_{\text{PPE}}(\phi) = \mathbb{E}_{p(\theta, \mathbf{x})} \left[ - \int \eta(s_{\text{dp}} | s(\mathbf{x})) \log q_{\phi}(\theta | s_{\text{dp}}) ds_{\text{dp}} \right].$$

So PPE learns the posterior directly, while PLE learns the private likelihood.

# A unified form of the PPE / PLE losses

Both losses have the same nested-integral structure:

$$\ell(\phi) = -\mathbb{E}_{p(\theta, \mathbf{x})} \left[ \int \eta(s_{\text{dp}} | s(\mathbf{x})) g_{\phi}(s_{\text{dp}}, \theta) ds_{\text{dp}} \right].$$

Here

$$g_{\phi}(s_{\text{dp}}, \theta) = \begin{cases} \log q_{\phi}(\theta | s_{\text{dp}}), & \text{for PPE,} \\ \log q_{\phi}(s_{\text{dp}} | \theta), & \text{for PLE.} \end{cases}$$

## Key point

The optimization over  $\phi$  is driven by repeated evaluation of the inner integral

$$l_{\phi}(\theta, \mathbf{x}) = \int \eta(s_{\text{dp}} | s(\mathbf{x})) g_{\phi}(s_{\text{dp}}, \theta) ds_{\text{dp}}.$$

This is exactly where quasi-Monte Carlo can help.

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This is exactly where quasi-Monte Carlo can help.

# Monte Carlo estimator for the inner privacy integral

Suppose the privacy mechanism can be written as

$$s_{\text{dp}} = \tau(u; s(\mathbf{x})), \quad u \sim \text{Unif}([0, 1]^r),$$

where  $r = \dim(\mathbb{S})$  is the dimension of the released private summary.

Then

$$I_\phi(\theta, \mathbf{x}) = \int_{[0,1]^r} g_\phi(\tau(u; s(\mathbf{x})), \theta) \, du.$$

Using i.i.d. uniforms  $u^{(1)}, \dots, u^{(M)}$ , the standard MC estimator is

$$\widehat{I}_{\theta, \mathbf{x}}^{\text{MC}} = \frac{1}{M} \sum_{m=1}^M g_\phi(\tau(u^{(m)}; s(\mathbf{x})), \theta),$$

with root-mean-squared-error (RMSE) typically of order

$$\mathcal{O}(M^{-1/2}).$$

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# RQMC estimator for the inner privacy integral

Now replace the i.i.d. uniforms by randomized low-discrepancy points

$$v^{(1)}, \dots, v^{(M)} \in [0, 1]^r.$$

The RQMC estimator is

$$\widehat{I}_{\theta, \mathbf{x}}^{\text{RQMC}} = \frac{1}{M} \sum_{m=1}^M g_{\phi}(\tau(v^{(m)}; s(\mathbf{x})), \theta).$$

If the integrand

$$\tilde{g}_{\theta, \mathbf{x}}(u) := g_{\phi}(\tau(u; s(\mathbf{x})), \theta)$$

has bounded Hardy–Krause variation, then

$$\mathbb{E} \left[ \left( \widehat{I}_{\theta, \mathbf{x}}^{\text{RQMC}} - I_{\phi}(\theta, \mathbf{x}) \right)^2 \right] = \mathcal{O}(M^{-2+2\delta}),$$

so the RMSE is

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# Why RQMC is especially natural here

- The difficult integral is over the **released private query**  $s_{\text{dp}}$ , not over the full confidential data  $\mathbf{X}$ .
- In many examples,  $\dim(\mathbb{S})$  is small even when  $\dim(\mathbb{X}^n)$  is very large.
- This is the regime where RQMC can substantially reduce variance.

## Computation

For each training pair  $(\theta, \mathbf{x})$ , we approximate

$$l_{\phi}(\theta, \mathbf{x}) = \int \eta(s_{\text{dp}} | s(\mathbf{x})) g_{\phi}(s_{\text{dp}}, \theta) ds_{\text{dp}}$$

by RQMC, and then optimize the empirical loss over  $\phi$ .

# RQMC improves the inner-integral estimation

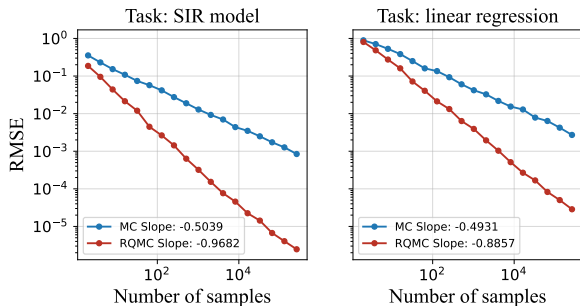


Figure: Rate of convergence.

- MC: RMSE typically scales as  $M^{-1/2}$ .
- RQMC: RMSE can approach  $M^{-1}$ .
- The gain is strongest when the released summary has low dimension.

## Message

In PPE / PLE, the computational bottleneck is the inner privacy integral, so better integration directly improves training.

# Performance comparison

Table: Estimated posterior mean and 95% credible intervals for the privatized Bayesian linear regression example using various methods. Here privacy loss budget is set to  $\epsilon = 10$ .

	$\beta_0$	$\beta_1$	$\beta_2$
Confidential Posterior	-2.15 (-2.68, -1.61)	-2.79 (-3.08, -2.50)	-0.83 (-1.08, -0.58)
Naive posterior	-4.63 (-5.04, -4.22)	-6.23 (-6.57, -5.90)	-5.10 (-5.40, -4.79)
DA-MCMC*	-0.62 (-2.50, 0.99)	-2.72 (-3.74, -0.96)	0.54 (-1.06, 2.46)
SMC-ABC	-0.59 (-2.28, 0.93)	-2.44 (-3.67, -0.38)	0.85 (-0.88, 2.77)
<b>PPE</b>	-0.51 (-2.25, 1.07)	-2.40 (-3.61, -0.30)	0.90 (-0.89, 2.85)
<b>PLE</b>	-0.64 (-2.31, 0.96)	-2.61 (-3.65, -0.98)	0.64 (-0.93, 2.48)
Gibbs-SS*	-0.46 (-2.22, 1.39)	-0.50 (-2.08, 1.28)	0.41 (-1.68, 2.13)
RNPE	-1.40 (-3.59, 0.94)	-2.15 (-3.34, 0.51)	0.29 (-2.11, 2.97)

# Performance comparison

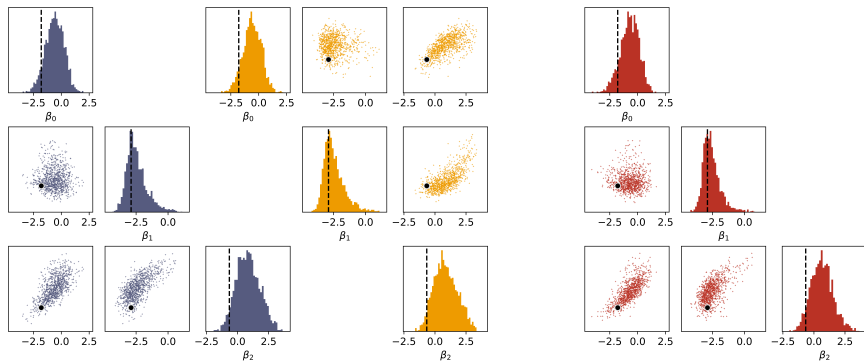


Figure: Posterior comparison on the Bayesian linear regression model. Grey: SMC-ABC; orange: PPE; red: PLE. The vertical lines and black dots indicate true data generating parameters.

# Summary: one problem, two strategies

## Original problem

The target posterior

$$\pi(\theta \mid s_{\text{dp}}) \propto \pi(\theta) f(s_{\text{dp}} \mid \theta)$$

is hard because the private likelihood

$$f(s_{\text{dp}} \mid \theta) = \int_{\mathbb{X}^n} f(\mathbf{x} \mid \theta) \eta(s_{\text{dp}} \mid s(\mathbf{x})) d\mathbf{x}$$

is typically intractable.

## Sampling-based strategy

Augment the latent confidential data and sample from  $\pi(\theta, \mathbf{x} \mid s_{\text{dp}})$  using MCMC; SOMA improves the imputation step.

## Variational-based strategy

Approximate either the posterior (PPE) or the private likelihood (PLE); RQMC improves the numerical evaluation of the resulting loss.

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# Thank you

Thank you!

Questions and comments are welcome.

# Random-Scan and Systematic-Scan IMwG

- To compare SOMA with more familiar samplers, we consider two standard **independent Metropolis-within-Gibbs** schemes targeting the same  $\pi(\mathbf{x})$ .

## Single-coordinate IMwG update at index $i$

- 1 Propose  $y \sim q(\cdot)$ .
- 2 Form  $\mathbf{x}' = [\mathbf{x}_{-i}, y]$ .
- 3 Accept with probability  $\alpha_i^{\text{IMwG}}(y, \mathbf{x}) = 1 \wedge \frac{\pi(\mathbf{x}')q(x_i)}{\pi(\mathbf{x})q(y)}$ .

- **Random-scan IMwG (Ran-scan):**

- at each step, draw  $l \sim \text{Unif}\{1, \dots, n\}$ ;
- perform the IMwG update at coordinate  $l$ .

- **Systematic-scan IMwG (Sys-scan):**

- in each sweep, visit  $i = 1, \dots, n$  in a fixed order;
- at each  $i$  perform the IMwG update above.

- See Qin and Jones (2022) for a detailed comparison of Ran-scan and Sys-scan when  $n = 2$ .

# Theoretical Comparison: SOMA vs IMwG

## Theorem (Acceptance dominance)

For any state  $\mathbf{x}$ , coordinate  $i$  and proposal  $y$ ,

$$\alpha_i^{\text{SOMA}}(y, \mathbf{x}) \geq \alpha_i^{\text{IMwG}}(y, \mathbf{x}). \quad (2)$$

Let  $A^{\text{SOMA}}(\mathbf{x})$  and  $A^{\text{RAN}}(\mathbf{x})$  denote the overall acceptance probabilities from  $\mathbf{x}$  for SOMA and Ran-scan IMwG, respectively. Then

$$A^{\text{SOMA}}(\mathbf{x}) \geq A^{\text{RAN}}(\mathbf{x}) \quad \text{for all } \mathbf{x}. \quad (3)$$

## Corollary (DP-based lower bounds)

If the mechanism is  $\epsilon$ -differentially private, we can show

$$A^{\text{SOMA}}(\mathbf{x}) \geq \frac{n}{n + e^\epsilon - 1}, \quad A^{\text{RAN}}(\mathbf{x}) \geq e^{-\epsilon}. \quad (4)$$

- As  $n \rightarrow \infty$ , SOMA becomes almost rejection-free, while the bound for Ran-scan IMwG stays at  $e^{-\epsilon}$ .

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# Convergence Rate under $\epsilon$ -DP

- We measure convergence by **geometric ergodicity**: for a kernel  $P$  with invariant distribution  $\pi$ ,

$$\|P^t(\cdot | x) - \pi(\cdot)\|_{\text{TV}} \leq C(x) r^t,$$

where  $r \in (0, 1)$  is the **convergence rate** (smaller  $r \Rightarrow$  faster mixing).

## Theorem (Upper bound on convergence rate)

*For the case  $n = 2$ , the convergence rate of the SOMA algorithm is upper bounded by*

$$r_{\text{SOMA}} \leq \frac{e^{2\epsilon} + 3e^\epsilon - 2 + \sqrt{e^{4\epsilon} + 2e^{3\epsilon} + 9e^{2\epsilon} - 8e^\epsilon}}{2(1 + e^\epsilon)^2}. \quad (5)$$

Smaller loss budget  $\epsilon$  (i.e., larger DP noise) leads to a smaller upper bound  $r_{\text{SOMA}}$  and hence faster convergence in this example.

# Computational Cost per iteration

- We compare SOMA with a single IMwG update from the viewpoint of **computational work per proposal**  $y$ .
- Both methods use the same proposal distribution  $q(\cdot)$  and draw one proposal  $y$  per iteration.

## Evaluating the target

- IMwG at index  $i$ : evaluate  $\pi(\mathbf{x})$  and  $\pi([\mathbf{x}_{-i}, y])$  for a *single* coordinate.
  - SOMA: evaluate  $\pi([\mathbf{x}_{-i}, y])$  for *all*  $i = 1, \dots, n$  plus  $\pi(\mathbf{x})$  (to form  $w_0$ ).
- 
- Naively, this is an extra factor of  $n$  target evaluations for SOMA.
  - However these  $n$  weights can be computed in **parallel** (or vectorized). In our DP applications, the wall-clock time per SOMA step is still comparable to a single-try IMwG update.

# SOMA is a good sampler, but not magical

## Drawbacks

- It is still based on an independent Metropolis proposal; performance depends on how well  $q$  matches  $\pi$ .
- Although the time cost may be comparable, the computational cost remains large.

## Good news: when does SOMA shine?

- Target is **permutation-invariant** and naturally written as  $n$  exchangeable coordinates.
- Expensive part of  $\pi(\mathbf{x})$  can be evaluated in **parallel** for all coordinates.
- In this setting, SOMA
  - strictly dominates random-scan IMwG in acceptance probability;
  - can be almost rejection-free when  $n$  is large;
  - empirically mixes faster than both random- and systematic-scan IMwG.

# Normalizing flows

Start from a simple base density

$$z \sim q_0(z), \quad \text{e.g. } q_0 = N(0, I_d),$$

and apply an invertible map

$$x = \tau(z).$$

## Change of variables

If  $\tau$  is invertible and differentiable, then

$$q_\tau(x) = q_0(z) |\det J_\tau(z)|^{-1}, \quad z = \tau^{-1}(x).$$

Hence

$$\log q_\tau(x) = \log q_0(\tau^{-1}(x)) - \log |\det J_\tau(\tau^{-1}(x))|.$$

- A normalizing flow chooses  $\tau$  to be flexible but still easy to invert and differentiate.
- Relatedly, Liu (2024) studies QMC-friendly transport maps inspired by normalizing flows.