

# Tailoring Differentially Private Bayesian Inference to Distance Between Distributions

Extended Abstract

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## ABSTRACT

Bayesian inference is a statistical method which allows one to derive a *posterior* distribution over a parameter, starting from a *prior* distribution and observed data. Settings where sensitive individual information is involved, different approaches have been taken to make this process differentially private. Dimitrakakis et al. [2], and Wang et al. [6], for instance, proved that, under specific conditions, sampling from the posterior distribution is already differentially private. Other works, e.g., [8], [3], designed differentially private mechanisms that output a representation of the full posterior distribution. Also, accuracy of these mechanisms was measured using metrics on the space of the numeric parameters of the posterior distribution, e.g., the  $\ell_1$ -norm.

In this work we present a new differentially private algorithm to compute the posterior in a Beta-Binomial system and in a Dirichlet-Multinomial system. We apply the exponential mechanism to the smooth sensitivity of a metric on probability distributions; we measured the accuracy of the process by using the same metric. We compared the accuracy of this approach with the ones based on  $\ell_1$ -norm. Experimental results show that when the data size is small or when the parameters of the prior distribution are large, the former outperforms the latter.

## KEYWORDS

Differential privacy, Bayesian inference, Hellinger distance

## 1 AN INFORMAL MOTIVATION

Publishing the posterior distribution inferred from a sensitive dataset can leak information about individuals on the dataset. Differential privacy enables release of an approximate posterior distribution through the introduction of random noise. The amount of the noise that must be introduced depends on the privacy parameter  $\epsilon$  and the sensitivity of the inference to small changes in the data set. Sensitivity can be computed in many different ways based on which metric space is imposed on the output set of the mechanism. In the literature on private Bayesian inference ([7, 8]), it is only measured with respect to the vector of numbers parameterizing the output distribution using e.g. The  $\ell_1$  norm. A more natural approach which we explore here, is to measure sensitivity with respect to a metric on the space of inferred probability distributions. A re-loved question is that of how to measure accuracy. Again, this can be answered in different ways based on the metric imposed

on the output space, and yet again only in few works in literature (e.g. [8]) distances between probability measures have been used for these purposes.

The question that this work aims at answering is whether an approach based on probability metrics can improve on the accuracy of approaches based on metrics over the numeric parameters of the distributions. We will see that in some cases this can happen.

## 2 BAYESIAN INFERENCE BACKGROUND

Given a prior belief  $\Pr(\theta)$  on some parameter  $\theta$ , and an observation  $\mathbf{x}$ , the posterior distribution on  $\theta$  given  $\mathbf{x}$  is computed as:

$$\Pr(\theta|\mathbf{x}) = \frac{\Pr(\mathbf{x}|\theta) \cdot \Pr(\theta)}{\Pr(\mathbf{x})}$$

where the expression  $\Pr(\mathbf{x}|\theta)$  denotes the *likelihood* of observing  $\mathbf{x}$  under a value of  $\theta$ . Since we consider  $\mathbf{x}$  to be fixed, the likelihood is a function of  $\theta$ . For the same reason  $\Pr(\mathbf{x})$  is a constant independent of  $\theta$ . Usually in statistics the prior distribution  $\Pr(\theta)$  is chosen so that it represents the initial belief on  $\theta$ , that is, when no data has been observed. In practice though, prior distributions and likelihood functions are usually chosen so that the posterior belongs to the same *family* of distributions. In this case we say that the prior is conjugate to the likelihood function. Use of a conjugate prior simplifies calculations and allows for inference to be performed in a recursive fashion over the data. In this work we will consider a specific instance of Bayesian inference and one of its generalizations. Specifically, we will consider the situation where  $\theta$  represents the parameter –informally called *bias*– of a Bernoulli distributed random variable, and its immediate generalization where the parameter  $\theta$  represents the vector of parameters of a conjugate distributed random variable. In the former case, the prior distribution over  $\theta \in [0, 1]$  is going to be a beta distribution,  $\text{beta}(\alpha, \beta)$ , with parameters  $\alpha, \beta \in \mathbb{R}^+$ , and with p.d.f:

$$\Pr(\theta) \equiv \frac{\theta^\alpha (1 - \theta)^\beta}{B(\alpha, \beta)}$$

where  $B(\cdot, \cdot)$  is the beta function. The data  $\mathbf{x}$  will be a sequence of  $n \in \mathbb{N}$  binary values, that is  $\mathbf{x} = (x_1, \dots, x_n)$ ,  $x_i \in \{0, 1\}$ , and the likelihood function is:

$$\Pr(\mathbf{x}|\theta) \equiv \theta^{\Delta\alpha} (1 - \theta)^{n - \Delta\alpha}$$

where  $\Delta\alpha = \sum_{i=1}^n x_i$ . From this it can easily be derived that the posterior distribution is:

$$\Pr(\theta|\mathbf{x}) = \text{beta}(\alpha + \Delta\alpha, \beta + n - \Delta\alpha)$$

In the latter case the prior distribution over  $\theta \in [0, 1]^k$  is given by a Dirichlet distribution,  $\text{DL}(\alpha)$ , for  $k \in \mathbb{N}$ , and  $\alpha \in (\mathbb{R}^+)^k$ , with p.d.f:

$$\Pr(\theta) \equiv \frac{1}{B(\alpha)} \cdot \prod_{i=1}^k \theta_i^{\alpha_i-1}$$

where  $B(\cdot)$  is the generalized beta function. The data  $\mathbf{x}$  will be a sequence of  $n \in \mathbb{N}$  values coming from a universe  $\mathcal{X}$ , such that  $|\mathcal{X}| = k$ . The likelihood function will be:

$$\Pr(\mathbf{x}|\theta) \equiv \prod_{a_i \in \mathcal{X}} \theta_i^{\Delta\alpha_i},$$

with  $\Delta\alpha_i = \sum_{j=1}^n [x_j = a_i]$ , where  $[\cdot]$  represents Iverson bracket notation. Denoting by  $\Delta\alpha$  the vector  $(\Delta\alpha_1, \dots, \Delta\alpha_k)$  the posterior distribution over  $\theta$  turns out to be

$$\Pr(\theta|\mathbf{x}) = \text{DL}(\alpha + \Delta\alpha).$$

where  $+$  denotes the component wise sum of vectors of reals.

### 3 THE PROBLEM AND BASELINE APPROACH

We are interested in designing a mechanism for privately releasing the full posterior distributions derived in section 2, as opposed to just sampling from them. It's worth noticing that the posterior distributions are fully characterized by their parameters, and the family (beta, Dirichlet) they belong to. Hence, in case of the Beta-Binomial model we are interested in releasing a private version of the pair of parameters  $(\alpha', \beta') = (\alpha + \Delta\alpha, \beta + n - \Delta\alpha)$ , and in the case of the Dirichlet-Multinomial model we are interested in a private version of  $\alpha' = (\alpha + \Delta\alpha)$ . Zhang et al. [8] and Xiao and Xiong [7] have already attacked this problem by adding independent Laplacian noise to the parameters of the posteriors. That is, in the case of the Beta-Binomial system, the value released would be:  $(\tilde{\alpha}, \tilde{\beta}) = (\alpha + \tilde{\Delta}\alpha, \beta + n - \tilde{\Delta}\alpha)$  where  $\tilde{\Delta}\alpha \sim \text{Lap}(\Delta\alpha, \frac{2}{\epsilon})$ , and where  $\text{Lap}(\mu, \nu)$  denotes a Laplace random variable with mean  $\mu$  and scale  $\nu$ . This mechanism is  $\epsilon$ -differentially private, and the noise is calibrated w.r.t. to a sensitivity of 2 which is derived by using  $\ell_1$  norm over the pair of parameters. Indeed, considering two adjacent<sup>1</sup> data observations  $\mathbf{x}, \mathbf{x}'$ , that, from a unique prior, give rise to two posterior distributions, characterized by the pairs  $(\alpha', \beta')$  and  $(\alpha'', \beta'')$  then  $|\alpha' - \alpha''| + |\beta' - \beta''| \leq 2$ . This argument extends similarly to the Dirichlet-Multinomial system. Also, in previous works, the accuracy of the posterior was measured again with respect to  $\ell_1$  norm. That is, an upper bound was given on

$$\Pr[|\alpha - \tilde{\alpha}| + |\beta - \tilde{\beta}| \geq \gamma]$$

where  $(\alpha, \beta)$ ,  $(\tilde{\alpha}, \tilde{\beta})$  are as defined above. In this work we will use a metric based on a different norm to compute the sensitivity and

<sup>1</sup>Given  $\mathbf{x}, \mathbf{x}'$  we say that  $\mathbf{x}$  and  $\mathbf{x}'$  are adjacent and we write,  $\text{adj}(\mathbf{x}, \mathbf{x}')$ , iff

$$\sum_{i=1}^n [x_i = x'_i] \leq 1.$$

provide guarantees on the accuracy. In particular we will consider a metric over probability measures and not over the parameters that represent them. Specifically, we will use the Hellinger distance  $\mathcal{H}(\cdot, \cdot)$ . Our choice to use Hellinger distance is motivated by two facts, first of all it simplifies calculations in the case of the probabilistic models considered here and second of all it also automatically yields bounds on the total variation distance, which represents also the maximum advantage an unbounded adversary can have in distinguishing two distributions. Given two beta distributions  $\beta_1 = \text{beta}(\alpha_1, \beta_1)$ , and  $\beta_2 = \text{beta}(\alpha_2, \beta_2)$  the following equality holds

$$\mathcal{H}(\beta_1, \beta_2) = \sqrt{1 - \frac{B(\frac{\alpha_1+\alpha_2}{2}, \frac{\beta_1+\beta_2}{2})}{\sqrt{B(\alpha_1, \beta_1)B(\alpha_2, \beta_2)}}}$$

The same change of metric will be applied to the experimental accuracy guarantees.

### 4 OUR APPROACH - EXPONENTIAL MECHANISM WITH SMOOTH SENSITIVITY

Given a prior distribution  $\beta_{\text{prior}} = \text{beta}(\alpha, \beta)$  and a sequence of  $n$  observations  $\mathbf{x} \in \{0, 1\}^n$ , we define the following set:

$$\mathcal{R}_{\text{post}} \equiv \{\text{beta}(\alpha', \beta') \mid \alpha' = \alpha + \Delta\alpha, \beta' = \beta + n - \Delta\alpha\}$$

where  $\Delta\alpha$  is as defined in Section 2. Notice that  $\mathcal{R}_{\text{post}}$  has  $n + 1$  elements, and the Bayesian Inference process will produce an element from  $\mathcal{R}_{\text{post}}$  that we denote by  $\text{BI}(\mathbf{x})$  – we don't explicitly parametrize the result by the prior, which from now on we consider fixed and we denote it by  $\beta_{\text{prior}}$ .

We can now define the mechanism  $\mathcal{M}_{\mathcal{H}}^B$  which, given in input a sequence of observations  $\mathbf{x}$  and parameters  $\epsilon > 0$  and  $\delta > 0$ , produces an element  $r$  in  $\mathcal{R}_{\text{post}}$  with probability:

$$\Pr_{z \sim \mathcal{M}_{\mathcal{H}}^B} [z = r] = \frac{\exp\left(\frac{-\epsilon \cdot \mathcal{H}(\text{BI}(\mathbf{x}), r)}{2 \cdot S(\mathbf{x})}\right)}{\sum_{r \in \mathcal{R}_{\text{post}}} \exp\left(\frac{-\epsilon \cdot \mathcal{H}(\text{BI}(\mathbf{x}), r)}{2 \cdot S(\mathbf{x})}\right)}.$$

This mechanism is based on the basic exponential mechanism [4], with  $\mathcal{R}_{\text{post}}$  as the range and  $\mathcal{H}(\cdot, \cdot)$  as the scoring function. The difference is that in this mechanism we don't calibrate the noise w.r.t. to the global sensitivity of the scoring function but w.r.t. to the smooth sensitivity  $S(\mathbf{x})$  – defined by Nissim et al. [5] – of  $\mathcal{H}(\text{BI}(\mathbf{x}), \cdot)$ . The smooth sensitivity is computed as follows:

$$S(\mathbf{x}) = \max_{\mathbf{x}' \neq \mathbf{x}, \mathbf{x}' \in \{0, 1\}^n} \left\{ \Delta_l \left( \mathcal{H}(\text{BI}(\mathbf{x}'), \cdot) \right) \cdot e^{-\gamma \cdot d(\mathbf{x}, \mathbf{x}')} \right\}, \quad (1)$$

where  $d$  is the Hamming distance between two datasets,  $\gamma = \gamma(\epsilon, \delta)$  is a function of  $\epsilon$  and  $\delta$  to be determined later, and where  $\Delta_l \left( \mathcal{H}(\text{BI}(\mathbf{x}'), \cdot) \right)$  denotes the local sensitivity at  $\text{BI}(\mathbf{x}')$ , or equivalently at  $\mathbf{x}'$ , the scoring function used in our mechanism. That is:

$$\Delta_l \left( \mathcal{H}(\text{BI}(\mathbf{x}'), \cdot) \right) = \max_{\mathbf{x}'' \in \mathcal{X}^n: \text{adj}(\mathbf{x}', \mathbf{x}''), r \in \mathcal{R}_{\text{post}}} |\mathcal{H}(\text{BI}(\mathbf{x}'), r) - \mathcal{H}(\text{BI}(\mathbf{x}''), r)|.$$

This mechanism also extends to the Dirichlet-Multinomial system  $\text{DL}(\alpha)$  by rewriting the Hellinger distance as:

$$\mathcal{H}(\text{DL}(\alpha_1), \text{DL}(\alpha_2)) = \sqrt{1 - \frac{B(\frac{\alpha_1 + \alpha_2}{2})}{\sqrt{B(\alpha_1)B(\alpha_2)}}},$$

and by replacing the  $\mathcal{R}_{\text{post}}$  with set of posterior Dirichlet distributions candidates. Also, the smooth sensitivity  $S(\mathbf{x})$  in (1) will be computed by letting  $\mathbf{x}'$  range over all the elements in  $\mathcal{X}^n$  adjacent to  $\mathbf{x}$ . Notice that  $\mathcal{R}_{\text{post}}$  has  $\binom{n+1}{m-1}$  elements in this case. We will denote by  $\mathcal{M}_{\mathcal{H}}^D$  the mechanism for the Dirichlet-Multinomial system.

The following result guarantees that the mechanisms are indeed differentially private.

**THEOREM 4.1 (PRIVACY).** *For  $\gamma = \ln(1 - \frac{\epsilon}{2 \ln(\frac{\delta}{2(n+1)})})$ , both  $\mathcal{M}_{\mathcal{H}}^B$  and  $\mathcal{M}_{\mathcal{H}}^D$  are  $(\epsilon, \delta)$ -differentially private.*

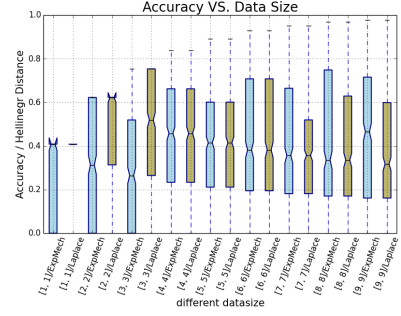
## 5 PRELIMINARY EXPERIMENTAL RESULTS

In this section, we evaluate the accuracy of the mechanisms defined in Section (4) w.r.t. four variables, including data size, dimensions, data variance, prior distribution, and some combinations thereof. Every plot is an average over 1000 runs. In all the experiments we set  $\epsilon = 0.8$ , and  $\delta = 10^{-8}$ .

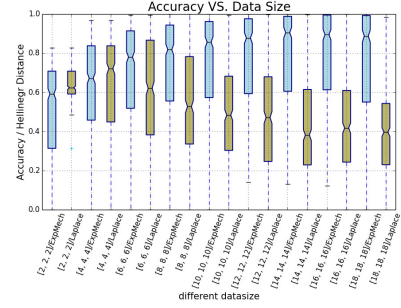
In the following whiskers-plots, the y-axis shows the accuracy (or equivalently, the error) of the mechanisms. Accuracy is measured using Hellinger distance. The x-axis, instead shows one of the previously mentioned independent variables –or a combination thereof. The boxes extend from the lower to the upper quartile values of the data, with a line at the median. A notch on the box around the median is also drawn to give a rough guide to the significance of difference of medians; The whiskers extend from the box to show the range of the data. A blue box in the plots represents our newly designed exponential mechanism behavior, while the yellow box next to it represents the performance of a variation of the basic Laplace mechanism presented in Section (3) with the same settings: that is  $\epsilon, \delta$ , data, prior. The variation considered performs a post processing on the released parameters so that they are consistent. For instance when the sum of the noised parameters is greater than  $n$  we will truncate them so that they sum up to  $n$ .

*Increasing data size with balanced datasets.* In Figure 1 we consider *balanced* datasets of observations. This means that in the Beta-Binomial setting the datasets will be half 1s and half 0s, while in the Dirichlet-Multinomial the data will be equally split in all the  $k = 3$  bins. Figure 1 shows that when the data size increases, the accuracy of  $\mathcal{M}_{\mathcal{H}}^B$  and  $\mathcal{M}_{\mathcal{H}}^D$  decreases. In Figure 1(a), when the data size is smaller than 12,  $\mathcal{M}_{\mathcal{H}}^B$  can outperform the Laplace mechanism but it fails to do so for bigger datasets. The same happens in as in Figure 1(b): the Laplace mechanism starts performing better for datasets bigger than 15.

*Increasing dimensions and data size with balanced dataset.* In Figure 2, in the x-axis are observed data sets of different sizes and different dimensions. The plot shows that increasing the number of dimensions has a similar pejorative effect on  $\mathcal{M}_{\mathcal{H}}^B$ ,  $\mathcal{M}_{\mathcal{H}}^D$ , and the Laplace mechanism. Fixing the number of dimensions and increasing the data size shows that the Laplace mechanism is more



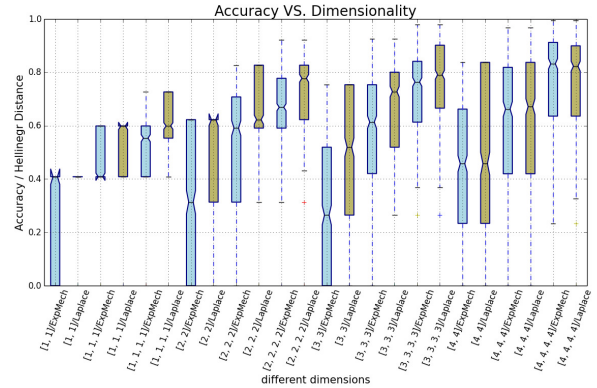
(a) two dimensions with beta(1, 1) prior distribution



(b) three dimensions with DL(1, 1, 1) prior distribution

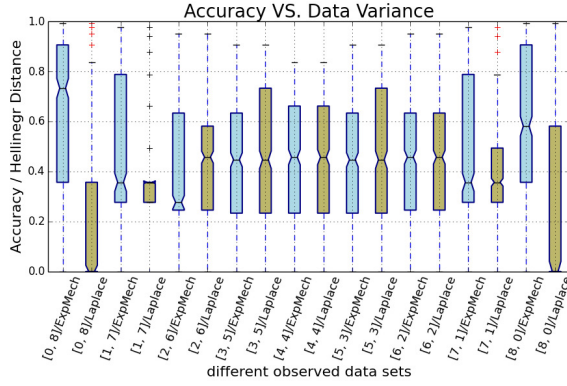
**Figure 1: Balanced datasets**

accurate than both  $\mathcal{M}_{\mathcal{H}}^B$  and  $\mathcal{M}_{\mathcal{H}}^D$ . In other words, dimensions have little influence on whether our mechanisms will beat the Laplace mechanism.



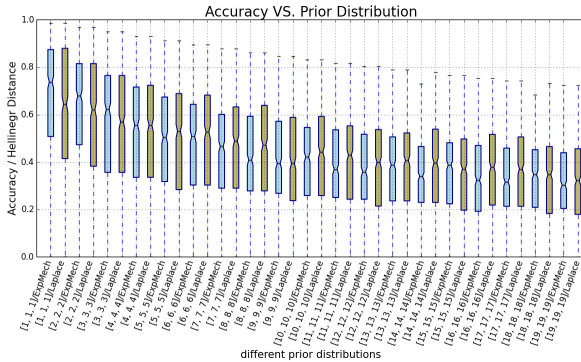
**Figure 2: Increasing dimensions and data size with balanced dataset. Priors have 1s in every dimension.**

*Fixed data size and unbalanced datasets.* In Figure 3, in the x-axis we considered different levels of balance of the datasets. We study this variable only under the two-dimensional beta distribution in order to be more concise. The plot shows that  $\mathcal{M}_{\mathcal{H}}^B$  accuracy is better than the one of Laplace when the dataset is balanced.



**Figure 3: Unbalanced datasets. Prior distributions have 1s in every dimension.**

*Fixed dataset varying balanced priors.* In Figure 4, we fix the dataset to be  $\langle 5, 5, 5 \rangle$ . We also considered balanced priors with increasing values in their dimensions. The plot shows that in the beginning the Laplace mechanism performs better but it is outperformed after a while.



**Figure 4: Observed data set is:  $\langle 5, 5, 5 \rangle$ , varying balanced priors**

## 6 CONCLUSIONS AND FUTURE WORKS

From what we have seen in the previous sections we can obtain some preliminary conclusions. That is, the probability measure approach outperforms the  $\ell_1$ -norm approach in the following cases:

- (1) When the data size is small
- (2) When the observed data is balanced
- (3) When priors parameters increase (eventually).

These results although very motivating, are still not enough for real world applications. Hence, we will continue our work in the following directions:

- (1) For now, we just have a intuitive idea on the accuracy behavior of our mechanisms, and not a precise formula or bound on it. When do our mechanisms perform better than the baseline mechanism and when they don't? How much

influence will elements in Section 5 have on the accuracy? Are there any other important factors we missed? These are all questions w.r.t. the accuracy that we are going to explore next, and in a more principled and formal way.

- (2) Theorem 4.1 provides an upper bound on the privacy loss for  $\mathcal{M}_{\mathcal{H}}^B$  and  $\mathcal{M}_{\mathcal{H}}^D$  but not necessarily a tight one. Indeed, experiments have shown that the actual privacy loss in the experiments can be smaller than  $\epsilon$ . This means that we could improve accuracy, by adding less noise – that is noise proportional to a higher value of  $\epsilon$  – but still achieve  $(\epsilon, \delta)$ -dp.
- (3) The choice of the Hellinger distance might seem quite ad-hoc. Hence, it is worth exploring other distances over distributions. An interesting class of probability metrics is the family of  $f$ -divergences [1].

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