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Online but Accurate Inference for Latent Variable Models with Local Gibbs Sampling

Christophe Dupuy
Technicolor
INRIA
Departement d’Informatique de l’ENS, Ecole normale supérieure, CNRS, PSL Research University
Paris, France.

Francis Bach
INRIA
Departement d’Informatique de l’ENS, Ecole normale supérieure, CNRS, PSL Research University
Paris, France.

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Abstract

We study parameter inference in large-scale latent variable models. We first propose a unified treatment of online inference for latent variable models from a non-canonical exponential family, and draw explicit links between several previously proposed frequentist or Bayesian methods. We then propose a novel inference method for the frequentist estimation of parameters, that adapts MCMC methods to online inference of latent variable models with the proper use of local Gibbs sampling. Then, for latent Dirichlet allocation, we provide an extensive set of experiments and comparisons with existing work, where our new approach outperforms all previously proposed methods. In particular, using Gibbs sampling for latent variable inference is superior to variational inference in terms of test log-likelihoods. Moreover, Bayesian inference through variational methods perform poorly, sometimes leading to worse fits with latent variables of higher dimensionality.

Keywords: Latent Variables Models, Online Learning, Gibbs Sampling, Topic Modelling, Latent Dirichlet Allocation

1. Introduction

Probabilistic graphical models provide general modelling tools for complex data, where it is natural to include assumptions on the data generating process by adding latent variables in the model. Such latent variable models are adapted to a wide variety of unsupervised learning tasks (Koller and Friedman, 2009; Murphy, 2012). In this paper, we focus on parameter inference in such latent variable models where the main operation needed for the standard expectation-maximization (EM) algorithm is intractable, namely dealing with conditional distributions over latent variables given the observed variables; latent Dirichlet allocation (LDA) (Blei et al., 2003) is our motivating example, but many hierarchical models exhibit this behavior, e.g., ICA with heavy-tailed priors. For such models, there exist two main classes of methods to deal efficiently with intractable exact inference in large-scale situations: sampling methods or variational methods.

Sampling methods can handle arbitrary distributions and lead to simple inference algorithms while converging to exact inference. However it may be slow to converge and non scalable to
big datasets in practice. In particular, although efficient implementations have been developed, for example for LDA (Zhao et al., 2014; Yan et al., 2009), MCMC methods may not deal efficiently yet with continuous streams of data for our general class of models.

On the other hand, **variational inference** builds an approximate model for the posterior distribution over latent variables—called variational—and infer parameters of the true model through this approximation. The fitting of this variational distribution is formulated as an optimization problem where efficient (deterministic) iterative techniques such as gradient or coordinate ascent methods apply. This approach leads to scalable inference schemes (Hoffman et al., 2013), but due to approximations, there always remains a gap between the variational posterior and the true posterior distribution, inherent to algorithm design, and that will not vanish when the number of samples and the number of iterations increase.

Beyond the choice of approximate inference techniques for latent variables, parameter inference may be treated either from the frequentist point of view, e.g., using maximum likelihood inference, or a Bayesian point of view, where the posterior distribution of the parameter given the observed data is approximated. With massive numbers of observations, this posterior distribution is typically peaked around the maximum likelihood estimate, and the two inference frameworks should not differ much (Van der Vaart, 2000).

In this paper, we focus on methods that make a single pass over the data to estimate parameters. We make the following contributions:

1. We review and compare existing methods for online inference for latent variable models from a non-canonical exponential family in Section 2, and draw explicit links between several previously proposed frequentist or Bayesian methods. Given the large number of existing methods, our unifying framework allows to understand differences and similarities between all of them.

2. We propose in Section 3 a novel inference method for the frequentist estimation of parameters, that adapts MCMC methods to online inference of latent variable models with the proper use of “local” Gibbs sampling. In our online scheme, we apply Gibbs sampling to the current observation, which is “local”, as opposed to “global” batch schemes where Gibbs sampling is applied to the entire dataset.

3. After formulating LDA as a non-canonical exponential family in Section 4, we provide an extensive set of experiments in Section 6, where our new approach outperforms all previously proposed methods. In particular, using Gibbs sampling for latent variable inference is superior to variational inference in terms of test log-likelihoods. Moreover, Bayesian inference through variational methods perform poorly, sometimes leading to worse fits with latent variables of higher dimensionality.

## 2. Online EM

We consider an exponential family model on random variables \((X, h)\) with parameter \(\eta \in \mathcal{E} \subseteq \mathbb{R}^d\) and with density (Lehmann and Casella, 1998):

\[
p(X, h|\eta) = a(X, h) \exp \left[ \phi(\eta), S(X, h) - \psi(\eta) \right].
\]

(1)

We assume that \(h\) is hidden and \(X\) is observed. The vector \(\phi(\eta) \in \mathbb{R}^d\) represents the natural parameter, \(S(X, h) \in \mathbb{R}^d\) is the vector of sufficient statistics, \(\psi(\eta)\) is the log-normalizer, and \(a(X, h)\)
is the underlying base measure. We consider a non-canonical family as in many models (such as LDA), the natural parameter \( \phi(\eta) \) does not coincide with the model parameter \( \eta \), that is, \( \phi(\eta) \neq \eta \); we however assume that \( \phi \) is injective.

We consider \( N \) i.i.d. observations \( (X_i)_{i=1,...,N} \) from a distribution \( t(X) \), which may be of the form \( P(X|\eta^*) = \int_{\phi} p(X, h|\eta^*) dh \) for our model above and a certain \( \eta^* \in \mathcal{E} \) (well-specified model) or not (misspecified model). Our goal is to obtain a predictive density \( r(X) \) built from the data and using the model defined in (1), with the maximal expected log-likelihood \( \mathbb{E}_{t(X)} \log r(X) \).

### 2.1 Maximum Likelihood Estimation

In the frequentist perspective, the predictive distribution \( r(X) \) is of the form \( p(X|\hat{\eta}) \), for a well-defined estimator \( \hat{\eta} \in \mathcal{E} \). The most common method is the EM algorithm (Dempster et al., 1977), which is an algorithm that aims at maximizing the likelihood of the observed data, that is,

\[
\max_{\eta \in \mathcal{E}} \sum_{i=1}^{N} \log p(X_i|\eta).
\]

More precisely, the EM algorithm is an iterative process to find the maximum likelihood (ML) estimate given observations \( (X_i)_{i=1,...,N} \) associated to hidden variables \( (h_i)_{i=1,...,N} \). It may be seen as the iterative construction of lower bounds of the log-likelihood function (Bishop, 2006). In the exponential family setting (1), we have, by Jensen’s inequality, given the model defined by \( \eta' \in \mathcal{E} \) from the previous iteration, and for any parameter \( \eta \in \mathcal{E} \):

\[
\log p(X_i|\eta) = \log \int p(X_i, h_i|\eta) dh_i \\
\geq \int p(h_i|X_i, \eta') \log \frac{p(X_i, h_i|\eta)}{p(h_i|X_i, \eta')} dh_i \\
= \int p(h_i|X_i, \eta') \left( \langle \phi(\eta), S(X_i, h_i) \rangle - \psi(\eta) \right) dh_i - C_i(\eta') \\
= \langle \phi(\eta), \mathbb{E}_{p(h_i|X_i, \eta')} [S(X_i, h_i)] \rangle - \psi(\eta) - C_i(\eta'),
\]

for a certain constant \( C_i(\eta') \), with equality if \( \eta' = \eta \). Thus, EM-type algorithms build locally tight lower bounds of the log-likelihood in (2), which are equal to

\[
\langle \phi(\eta), \sum_{i=1}^{N} s_i \rangle - N\psi(\eta) + \text{cst},
\]

for appropriate values of \( s_i \in \mathbb{R}^d \) obtained by computing conditional expectations with the distribution of \( h_i \) given \( X_i \) for the current model defined by \( \eta' \) (E-step), i.e., \( s_i = \mathbb{E}_{p(h_i|X_i, \eta')} [S(X_i, h_i)] \). Then this function of \( \eta \) is maximized to obtain the next iterate (M-step). In standard EM applications, these two steps are assumed tractable. In Section 3, we will only assume that the M-step is tractable while the E-step is intractable.

Standard EM will consider \( s_i = \mathbb{E}_{p(h_i|X_i, \eta')} [S(X_i, h)] \) for the previous value of the parameter \( \eta \) for all \( i \), and hence, at every iteration, all observations \( X_i, i = 1, \ldots, N \) are considered for latent variable inference, leading to a slow “batch” algorithm for large \( N \).

Incremental EM (Neal and Hinton, 1998) will only update a single element \( s_i \) coming from a single observation \( X_i \) and update the corresponding part of the sum \( \sum_{j=1}^{N} s_j \) without changing
other elements. In the extreme case where a single pass over the data is made, then the M-step at iteration \( i \) maximizes

\[
\langle \phi(\eta), \sum_{j=1}^{i} \mathbb{E}_{p(h_j|X_j, \eta_{j-1})} [S(X_j, h_j)] \rangle - i\psi(\eta),
\]

with respect to \( \eta \). In the next section, we provide a (known) other interpretation of this algorithm.

### 2.2 Stochastic Approximation

Given our frequentist objective \( \mathbb{E}_{t(X)} \log p(X|\eta) \) to maximize defined as an expectation, we may consider two forms of stochastic approximation (Kushner and Yin, 2003), where observations \( X_t \) sampled from \( t(X) \) are processed only once. The first one is stochastic gradient ascent, of the form

\[
\eta_i = \eta_{i-1} + \rho_i \frac{\partial \log p(X_i|\eta)}{\partial \eta},
\]

or appropriately renormalized version thereof, i.e., \( \eta_i = \eta_{i-1} + \rho_i H^{-1} \frac{\partial \log p(X_i|\eta)}{\partial \eta} \), with several possibilities for the \( d \times d \) matrix \( H \), such as the negative Hessian of the partial or the full log-likelihood, or the negative covariance matrix of gradients, which can be seen as versions of natural gradient—see Titterington (1984); Delyon et al. (1999); Cappé and Moulines (2009). This either leads to slow convergence (without \( H \)) or expensive iterations (with \( H \)), with the added difficulty of choosing a proper scale and decay for the step-size \( \rho_i \).

A key insight of Delyon et al. (1999); Cappé and Moulines (2009) is to use a different formulation of stochastic approximation, not explicitly based on stochastic gradient ascent. Indeed, they consider the stationary equation \( \mathbb{E}_{t(X)} \left[ \frac{\partial \log p(X|\eta)}{\partial \eta} \right] = 0 \) and expand it using the exponential family model (1) as follows:

\[
\frac{\partial \log p(X|\eta)}{\partial \eta} = \frac{\partial \log \int p(X, h|\eta) dh}{\partial \eta} = \phi'(\eta) \mathbb{E}_{p(h|X, \eta)} [S(X, h)] - \psi'(\eta).
\]

Given standard properties of the exponential family, namely

\[
\psi'(\eta) = \phi'(\eta) \mathbb{E}_{p(h, X|\eta)} [S(X, h)]^{-1},
\]

and assuming invertibility of \( \phi'(\eta) \), this leads to the following stationary equation:

\[
\mathbb{E}_{t(X)} \left[ \mathbb{E}_{p(h|X, \eta)} [S(X, h)] \right] = \mathbb{E}_{p(h, X|\eta)} [S(X, h)].
\]

---

**Proof:** Given (1), \( \int_{X,h} p(X, h|\eta) d(X, h) = 1 \) \( \Rightarrow \psi(\eta) = \log \left[ \int_{X,h} a(X, h)e^{\phi(\eta)S(X, h)} d(X, h) \right] \). We then derive this identity with respect to \( \eta \), which gives:

\[
\psi'(\eta) = \frac{\int_{X,h} \phi'(\eta) S(X, h) a(X, h) e^{\phi(\eta)S(X, h)} d(X, h)}{\int_{X,h} a(X, h) e^{\phi(\eta)S(X, h)} d(X, h)} = \phi'(\eta) \int_{X,h} S(X, h) a(X, h) e^{\phi(\eta)S(X, h)} d(X, h) = \phi'(\eta) \int_{X} S(X, h) p(X, h|\eta) d(X, h) = \phi'(\eta) \mathbb{E}_{p(h, X|\eta)} [S(X, h)].
\]
This stationary equation states that at optimality the sufficient statistics have the same expectation for the full model $p(h, X|\eta)$ and the joint “model/data” distribution $t(X)p(h|X, \eta)$.

Another important insight of Delyon et al. (1999); Cappe and Moulines (2009) is to consider the change of variable $s(\eta) = \mathbb{E}_{p(h, X|\eta)}[S(X, h)]$ on sufficient statistics, which is equivalent to

$$\eta = \eta^*(s) \in \text{arg max } \langle \phi(\eta), s \rangle - \psi(\eta),$$

(which is the usual M-step update). See Cappe and Moulines (2009) for detailed assumptions allowing this inversion. We may then rewrite the equation above as

$$\mathbb{E}_{t(X)} (\mathbb{E}_{p(h|X, \eta^*(s))}[S(X, h)]) = s.$$

This is a non-linear equation in $s \in \mathbb{R}^d$, with an expectation with respect to $t(X)$ which is only accessed through i.i.d. samples $X_i$, and thus a good candidate for the Robbins-Monro algorithm to solve stationary equations (and not to minimize functions) (Kushner and Yin, 2003), which takes the simple form:

$$s_i = s_{i-1} - \rho_i (s_{i-1} - \mathbb{E}_{p(h_i|X_i, \eta_{i-1})}[S(X_i, h_i)])$$

with a step-size $\rho_i$. It may be rewritten as

$$\begin{cases} s_i = (1 - \rho_i) s_{i-1} + \rho_i \mathbb{E}_{p(h_i|X_i, \eta_i-1)}[S(X_i, h_i)] \\ \eta_i = \eta^*(s_i), \end{cases} \tag{3}$$

which has a particularly simple interpretation: instead of computing the expectation for all observations as in full EM, this stochastic version keeps tracks of old sufficient statistics through the variable $s_{i-1}$ which is updated towards the current value $\mathbb{E}_{p(h_i|X_i, \eta_{i-1})}[S(X_i, h_i)]$. The parameter $\eta$ is then updated to the value $\eta^*(s_i)$. Cappe and Moulines (2009) show that this update is asymptotically equivalent to the natural gradient update with three main improvements: (a) no matrix inversion is needed, (b) the algorithm may be accelerated through Polyak-Ruppert averaging (Polyak and Juditsky, 1992), i.e., using the average $\bar{\eta}_N$ of all $\eta_i$ instead of the last iterate $\eta_N$, and (c) the step-size is particularly simple to set, as we are taking convex combinations of sufficient statistics, and hence only the decay rate of $\rho_i$ has to be chosen, i.e., of the form $\rho_i = i^{-\kappa}$, for $\kappa \in (0, 1]$, without any multiplicative constant.

2.2.1 Incremental View

For the specific stepsize $\rho_i = 1/i$, the online EM algorithm (3) corresponds exactly to the incremental EM presented above (Neal and Hinton, 1998), as then

$$s_i = \frac{1}{i} \sum_{j=1}^{i} \mathbb{E}_{p(h_j|X_j, \eta_{j-1})}[S(X_j, h_j)].$$

See Mairal (2014) for a detailed convergence analysis of incremental algorithms, in particular showing that step-sizes larger than $1/i$ are preferable (we observe this in practice in Section 6).
2.2.2 Monte Carlo Methods

There exist alternative methods to the EM algorithm based on Monte Carlo sampling to compute the maximum likelihood. For instance, the Monte Carlo EM method (MCEM) (Wei and Tanner, 1990) is a general Bayesian approach (i.e., \( \eta \) is a random variable) to approximate the maximizer of the posterior distribution \( p(\eta|X, h) \) with Monte Carlo sampling. More precisely, in the MCEM method, similarly to EM, a surrogate function of the log-likelihood is used, given by:

\[
Q(\eta, \eta_t) = \int_h \log[p(\eta|X, h)]p(h|X, \eta_t)dh.
\]

The function \( Q \) is approximated by sampling the latent variables \( h \) from the current conditional \( p(h|X, \eta_t) \):

\[
\hat{Q}(\eta, \eta_t) = \sum_{i=1}^{P} \log p(\eta|X, h^i),
\]

where \((h^i)_{i=1,...,P}\) are the samples drawn from the conditional \( p(h|X, \eta_t) \). The approximation \( \hat{Q} \) is then maximize with respect to \( \eta \). Note that this method is a batch method, namely, samples are drawn over all the dataset.

Other sequential Monte Carlo methods (SMC) use importance sampling to estimate the conditional distributions \( p(h|X) \) using an auxiliary density function. This auxiliary distribution is then reweighted at each iteration. In our case where the conditional distribution \( p(h|X) \) is intractable to compute, these methods fail to accurately approximate this distribution (Cappé et al., 2005; Kantas et al., 2015) and our Gibbs sampling scheme is more relevant. The SMC methods are also more adapted to models with dependency between observations.

The two Monte Carlo methods mentioned above also consist in sufficient statistics updates for the class of models considered here.

3. Online EM with Intractable Models

The online EM updates in (3) lead to a scalable algorithm for optimization when the local E-step is tractable. However, in many latent variable models—e.g., LDA, hierarchical Dirichlet processes (Teh et al., 2006), or ICA (Hyvärinen et al., 2004)—it is intractable to compute the conditional expectation \( \mathbb{E}_{p(h|X, \eta)}[S(X, h)] \).

Following Rohde and Cappé (2011), we propose to leverage the scalability of online EM updates (3) and locally approximate the conditional distribution \( p(h|X, \eta) \) in the case this distribution is intractable to compute. We will however consider different approximate methods, namely Gibbs sampling or variational inference. Our method is thus restricted to models where the hidden variable \( h \) may naturally be splitted in two or more groups of simple random variables. Our algorithm is described in Algorithm 1 and may be instantiated with two approximate inference schemes which we now describe.
Algorithm 1 Gibbs / Variational online EM

Input: $\eta_0$, $s_0$, $\kappa \in (0, 1]$.  
\textbf{for} $i = 1, \ldots, N$ \textbf{do}  
\begin{itemize} 
\item Collect observation $X_i$,  
\item Estimate $p(h_i | X_i, \eta_{i-1})$ with sampling ($G$--OEM) or variational inference ($V$--OEM),  
\item Apply (3) to sufficient statistics $s_i$ and parameter $\eta_i$ with $\rho_i = 1/i^\kappa$, 
\end{itemize}  
\textbf{end for}  
Output: $\bar{\eta}_N = \frac{1}{N} \sum_{i=1}^{N} \eta_i$ or $\eta_N$.

3.1 Variational Inference: $V$--OEM

While variational inference had been considered before for online estimation of latent variable models, in particular for LDA for incremental EM (Sato et al., 2010), using it for online EM (which is empirically faster) had not been proposed and allows to use bigger step-sizes (e.g., $\kappa = 1/2$). These methods are based on maximizing the negative variational “free-energy”

$$ \mathbb{E}_{q(h|\eta)} \left[ \log \frac{p(X, h|\eta)}{q(h|\eta)} \right], $$

with respect to $q(h|\eta)$ having a certain factorized form adapted to the model at hand, so that efficient coordinate ascent may be used. See, e.g., Hoffman et al. (2013). We now denote online EM with variational approximation of the conditional distribution $p(h|X, \eta)$ as $V$--OEM.

3.2 Sampling Methods: $G$--OEM

MCMC methods to approximate the conditional distribution of latent variables with online EM have been considered by Rohde and Cappé (2011), who apply locally the Metropolis-Hasting (M-H) algorithm (Metropolis et al., 1953; Hastings, 1970), and show results on simple synthetic datasets. While Gibbs sampling is widely used for many models such as LDA due to its simplicity and lack of external parameters, M-H requires a proper proposal distribution with frequent acceptance and fast mixing, which may be hard to find in high dimensions. We provide a different simpler local scheme based on Gibbs sampling (thus adapted to a wide variety of models), and propose a thorough favorable comparison on synthetic and real datasets with existing methods.

The Gibbs sampler is used to estimate posterior distributions by alternatively sampling parts of the variables given the other ones (see Casella and George 1992, for details), and is standard and easy to use in many common latent variable models. In the following, the online EM method with Gibbs estimation of the conditional distribution $p(h|X, \eta)$ is denoted $G$--OEM.

As mentioned above, the online EM updates correspond to a stochastic approximation algorithm and thus are robust to random noise in the local E-step. As a result, our sampling method is particularly adapted as it is a random estimate of the E-step—see a theoretical analysis by Rohde and Cappé (2011), and thus we only need to compute a few Gibbs samples for the estimation of $p(h_i | X_i, \eta_{i-1})$. A key contribution of our paper is to reuse sampling techniques that have proved competitive in the batch set-up and to compare them to existing variational approaches.
3.3 “Boosted” Inference

As the variational and MCMC estimations of \( p(h|X_i, \eta_{i-1}) \) are done with iterative methods, we can boost the inference of Algorithm 1 by applying the update in the parameter \( \eta \) in (3) after each iteration of the estimation of \( p(h|X_i, \eta_{i-1}) \). In the context of LDA, this was proposed by Sato et al. (2010) for incremental EM and we extend it to all versions of online EM. With this boost, we expect that the global parameters \( \eta \) converge faster, as they are updated more often. In the following, we denote by G-OEM++ (resp. V-OEM++) the method G-OEM (resp. V-OEM) augmented with this boost.

3.4 Variational Bayesian Estimation

In the Bayesian perspective where \( \eta \) is seen as a random variable, we either consider a distribution based on model averaging, e.g., \( r(X) = \int p(X|\eta)q(\eta)d\eta \) where \( q(\eta) \propto \prod_{i=1}^N p(X_i|\eta)p(\eta) \) is the posterior distribution, or

\[
q(\eta) \propto \prod_{i=1}^N p(X_i|\eta),
\]

where \( \bar{\eta} \) is the summary (e.g., the mean) of the posterior distribution \( q(\eta) \), or of an approximation, which is usually done in practice (see, e.g., Hoffman and Blei 2015) and is asymptotically equivalent when \( N \) tends to infinity.

The main problem is that, even when the conditional distribution of latent variables is tractable, it is intractable to manipulate the joint posterior distribution over the latent variables \( h_1, \ldots, h_N \), and the parameter \( \eta \). Variational inference techniques consider an approximation where hidden variables are independent of the parameter \( \eta \), i.e., such that

\[
p(\eta, h_1, \ldots, h_N|X_1, \ldots, X_N) \approx q(\eta)\prod_{i=1}^N q(h_i),
\]

which corresponds to the maximization of the following lower bound—called Evidence Lower BOund (ELBO)—on the log-likelihood \( \log p(X_1, \ldots, X_n) \) (Bishop 2006):

\[
\int q(\eta)\prod_{i=1}^N q(h_i) \log \frac{p(\eta)\prod_{i=1}^n p(X_i, h_i|\eta)}{q(\eta)\prod_{i=1}^N q(h_i)} d\eta dh_1 \cdots dh_N.
\]

The key insight from Hoffman et al. (2010); Broderick et al. (2013) is to consider the variational distribution \( q(\eta) \) as the global parameter, and the cost function above as a sum of local functions that depend on the data \( X_i \) and the variational distribution \( q(h_i) \). Once the local variational distribution \( q(h_i) \) is maximized out, the sum structure may be leveraged in similar ways than for frequentist estimation, either by direct (natural) stochastic gradient (Hoffman et al., 2010) or incremental techniques that accumulate sufficient statistics (Broderick et al. 2013). A nice feature of these techniques is that they extend directly to models with intractable latent variable inference, by making additional assumptions on \( q(h_i) \) (see for example the LDA situation in Section 4).

In terms of actual updates, they are similar to online EM in Section 3.1 with a few changes, but which turn out to lead to significant differences in practice. The similarity comes from the expansion of the ELBO as

\[
\mathbb{E}_{q(\eta)} \left[ \sum_{i=1}^N \mathbb{E}_{q(h_i)} \log \frac{p(X_i, h_i|\eta)}{q(h_i)} \right] + \mathbb{E}_{q(\eta)} \left[ \log \frac{p(\eta)}{q(\eta)} \right].
\]
The left hand side has the same structure than the variational EM update in (4), thus leading to similar updates, while the right hand side corresponds to the “Bayesian layer”, and the maximization with respect to $q(\eta)$ is similar to the M-step of EM (where $\eta$ is seen as a parameter).

Like online EM techniques presented in Section 3, approximate inference for latent variable is used, but, when using Bayesian stochastic variational inference techniques, there are two additional sources of inefficiencies: (a) extra assumptions regarding the independence of $\eta$ and $h_1, \ldots, h_N$, and (b) the lack of explicit formulation as the minimization of an expectation, which prevents the simple use of the most efficient stochastic approximation techniques (together with their guarantees). While (b) can simply slow down the algorithm, (a) may lead to results which are far away from exact inference, even for large numbers of samples (see examples in Section 6).

Beyond variational inference, Gibbs sampling has been recently considered by Gao et al. (2016): their method consists in sampling hidden variables for the current document given current parameters, but (a) only some of the new parameters are updated by incrementally aggregating the samples of the current document with current parameters, and (b) the method is slower than G-OEM (see Section 6).

4. Application to LDA

LDA (Blei et al., 2003) is a probabilistic model that infers hidden topics given a text corpus where each document of the corpus can be represented as topic probabilities. In particular, the assumption behind LDA is that each document is generated from a mixture of topics and the model infers the hidden topics and the topic proportions of each document. In practice, inference is done using Bayesian variational EM (Blei et al. 2003), Gibbs sampling (Griffiths and Steyvers, 2004; Wallach, 2006) or stochastic variational inference (Hoffman et al., 2010; Broderick et al., 2013; Sato et al., 2010).

4.0.1 Hierarchical Probabilistic Model.

Let $\mathcal{C} = \{X_1, \ldots, X_D\}$ be a corpus of $D$ documents, $V$ the number of words in our vocabulary and $K$ the number of latent topics in the corpus. Each topic $\beta^k$ corresponds to a discrete distribution on the $V$ words (that is an element of the simplex in $V$ dimensions). A hidden discrete distribution $\theta_i$ over the $K$ topics (that is an element of the simplex in $K$ dimensions) is attached to each document $X_i$. LDA is a generative model applied to a corpus of text documents which assumes that each word of the $i^{th}$ document $X_i$ is generated as follows:

- Choose $\theta_i \sim \text{Dirichlet}(\alpha)$,
- For each word $x_n \in X_i = (x_1, \ldots, x_{N_{X_i}})$:
  - Choose a topic $z_n \sim \text{Multinomial}(\theta_i)$,
  - Choose a word $x_n \sim \text{Multinomial}(\beta^{z_n})$.

In our settings, an observation is a document $X_i = (x_1, \ldots, x_{N_{X_i}})$ where for all $1 \leq n \leq N_{X_i}$, $x_n \in \{0, 1\}^V$ and $\sum_{v=1}^V x_{nv} = 1$. Each observation $X_i$ is associated with the hidden variables $h_i$, with $h_i \equiv (Z_i = (z_1, \ldots, z_{N_{X_i}}), \theta_i)$. The vector $\theta_i$ represents the topic proportions of document $X_i$ and $Z_i$ is the vector of topic assignments of each word of $X_i$. The variable $h_i$ is local, i.e., attached.
to one observation \( X_i \). The parameters of the model are global, represented by \( \eta \equiv (\beta, \alpha) \), where \( \beta \) represents the topic matrix and \( \alpha \) represents the Dirichlet prior on topic proportions.

We derive the LDA model in Section 4.1 to find \( \phi, S, \psi \) and \( a \) such that the joint probability \( p(Z, \theta|X, \alpha, \beta) \) is in a non-canonical exponential family (1).

We may then readily apply all algorithms from Section 3 by estimating the conditional expectation \( E_{Z,\theta|X,\alpha,\beta}[S(X, Z, \theta)] \) with either variational inference (V-OEM) or Gibbs sampling (G-OEM). See Sections 4.2 and 4.3 for online EM derivations. Note that the key difficulty of LDA is the presence of two interacting hidden variables \( Z \) and \( \theta \).

### 4.1 LDA and Exponential Families

An observation \( X \) is a document of length \( N_X \), where \( X = (x_1, \ldots, x_{N_X}) \), each word is represented by \( x_n \in \{0, 1\}^V \) with \( \sum_{v=1}^V x_{nv} = 1 \). Our corpus \( \mathcal{C} \) is a set of \( D \) observations \( \mathcal{C} = (X^1, \ldots, X^D) \).

For each document \( X^i \) a hidden variable \( \theta^i \) is associated, corresponding to the topic distribution of document \( X^i \). For each word \( x_n \) of document \( X^i \) a hidden variable \( z_n \in \{0, 1\}^K \) is attached, corresponding to the topic assignment of word \( x_n \). We want to find \( \phi, S, \psi \) and \( a \) such that, the joint probability is in the exponential family (1):

\[
p(X, Z, \theta|\beta, \alpha) = a(X, Z, \theta) \exp [\langle \phi(\beta, \alpha), S(X, Z, \theta) \rangle - \psi(\beta, \alpha)],
\]

given an observation \( X \) and hidden variables \( Z \) and \( \theta \).

For the LDA model, we have:

\[
p(X, Z, \theta|\beta, \alpha) = \prod_{n=1}^{N_X} p(x_n|z_n, \beta)p(z_n|\theta)p(\theta|\alpha) = \prod_{n=1}^{N_X} \prod_{k=1}^{K} \prod_{v=1}^{V} (\beta^k_{v})^{x_{nv}z_{nk}} \theta^k \]

which we can expand as:

\[
p(X, Z, \theta|\beta, \alpha) = \exp \left[ \sum_{n=1}^{N_X} \sum_{k=1}^{K} z_{nk} \log \theta^k \right] \times \exp \left[ \sum_{n=1}^{N_X} \sum_{k=1}^{K} \sum_{v=1}^{V} x_{nv}z_{nk} \log \beta^k_{v} \right] \times \exp \left[ \sum_{k=1}^{K} (\alpha_k - 1) \log \theta^k + B(\alpha) \right],
\]

with \( B(\alpha) = \log \left[ \Gamma \left( \sum_{i=1}^{K} \alpha_i \right) \right] - \sum_{i=1}^{K} \log[\Gamma(\alpha_i)] \), where \( \Gamma \) is the gamma function. We deduce the non-canonical exponential family setting \( \phi, S, \psi \) and \( a \):

\[
S(X, Z, \theta) = \begin{pmatrix}
S_{1k}^{\ell} & \equiv & \left[ \sum_{n=1}^{N_X} z_{nk}x_{nv} \right]_{kv} \\
S_{2k} & \equiv & [\log \theta^k]_{k}
\end{pmatrix}, \quad \phi(\beta, \alpha) = \begin{pmatrix}
\phi_{1k} & \equiv & [\log \beta^k_{v}]_{k} \\
\phi_{2k} & \equiv & [\alpha_k]_{k}
\end{pmatrix},
\]
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with \( S^1, \phi^1 \in \mathbb{R}^{K \times V} \) and \( S^2, \phi^2 \in \mathbb{R}^K \),

\[
\psi(\beta, \alpha) = \sum_{i=1}^{K} \log[\Gamma(\alpha_i)] - \log \left[ \Gamma \left( \sum_{i=1}^{K} \alpha_i \right) \right],
\]

(7)

and

\[
\alpha(X, Z, \theta) = \exp \left[ \sum_{k=1}^{K} \left( \sum_{n=1}^{N_X} z_{nk} - 1 \right) \log \theta_k \right].
\]

The one-to-one mapping between the sufficient statistics \( s = (s^1, s^2) \) and \((\beta, \alpha)\) is defined by:

\[
(\beta, \alpha)^*[s] = \left\{ \begin{array}{ll}
\arg \max_{\beta \geq 0, \alpha \geq 0} & \langle \phi(\beta, \alpha), s \rangle - \psi(\beta, \alpha) \\
\text{s.t.} & \beta^\top 1 = 1,
\end{array} \right.
\]

where \( 1 \) denotes the vector whose all entries equal 1. The objective \( \langle \phi(\beta, \alpha), s \rangle - \psi(\beta, \alpha) \) is concave in \( \beta \) from the concavity of \( \log \) and concave in any \( \alpha_k \) for \( \alpha \geq 0 \) as the function \( B(\alpha) \) is concave as the negative log-partition of the Dirichlet distribution. We use the Lagrangian method for \( \beta \):

\[
L(\beta, \lambda) = \sum_{k=1}^{K} \sum_{v=1}^{V} s^1_{kv} \log \beta^k_v + \lambda^\top (\beta^\top 1 - 1),
\]

with \( \lambda \in \mathbb{R}^K \). The derivative of \( L \) is set to zero when:

\[
\forall (k, v), s^1_{kv} \beta^k_v + \lambda_k = 0 \Rightarrow \lambda_k = -\sum_{v=1}^{V} s^1_{kv},
\]

as \( \sum_{v=1}^{V} \beta^k_v = 1 \). We then have \( (\beta^*(s))_{kv} = \frac{s^1_{kv}}{\sum_{j} s^1_{kj}} \). This mapping satisfies the constraint \( \beta \geq 0 \) because for any observation \( X \) and hidden variable \( Z \), we have \( S^1(X, Z)_{kv} \geq 0 \). This comes from (5) and the fact that \( \forall (n, k, v), (x_{nv}, z_{nk}) \in \{0, 1\}^2 \). We find the condition on \( \alpha \) by setting the derivatives to 0, which gives \( \forall k \in [1, K] \):

\[
s^2_k - \Psi([\alpha^*(s)]_k) + \Psi \left( \sum_{i=1}^{K} [\alpha^*(s)]_i \right) = 0,
\]

where \( \Psi : x \mapsto \frac{\partial}{\partial x} [\log \Gamma](x) \) is the digamma function. Finally, \( (\alpha^*(s), \beta^*(s)) \) satisfies \( \forall (k, v) \):

\[
\begin{cases}
(\beta^*(s))_{kv} = \left[ \frac{s^1_{kv}}{\sum_{j} s^1_{kj}} \right]_{kv} \\
\Psi([\alpha^*(s)]_k) - \Psi \left( \sum_{i=1}^{K} [\alpha^*(s)]_i \right) = s^2_k,
\end{cases}
\]

(8)

The parameter \( \alpha^* \) is usually estimated with gradient ascent (Blei et al., 2003; Hoffman et al., 2010). We can also estimate \( \alpha \) with the fixed point iteration (Minka, 2000) which consists in repeating the following update until convergence:

\[
\alpha_k^{\text{new}} = \Psi^{-1} \left( \Psi \left( \sum_{i=1}^{K} \alpha_i^{\text{old}} \right) + s^2_k \right).
\]

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We use the fixed point iteration to estimate $\alpha^*$ as it is more stable in practice. We study different updates for $\alpha$ in Appendix C.

We can now apply Algorithm 1 to LDA. The only missing step is the estimation of the conditional expectation $E_{Z,\theta|X,\alpha_t,\beta_t}[S(X, Z, \theta)]$, with $X = (x_1, \ldots, x_{N_X})$ and $Z = (z_1, \ldots, z_{N_Z})$. We explain how to approximate this expectation with variational inference and Gibbs sampling.

### 4.2 Variational Online EM Applied to LDA (V-OEM)

In this section we explain how to approximate $E_{Z,\theta|X,\alpha_t,\beta_t}[S(X, Z, \theta)]$ with variational inference, in the frequentist setting. See Hoffman et al. (2013) for detailed derivations of variational inference for LDA in the Bayesian setting (from which the updates in the frequentist setting may be easily obtained). The idea behind variational inference is to maximize the Evidence Lower BOund (ELBO), a lower bound on the probability of the observations:

$$p(X) \geq \text{ELBO}(X, p, q),$$

where $q$ represents the variational model. In the case of LDA, the variational model is often set with a Dirichlet($\gamma$) prior on $\theta$ and a multinomial prior on $Z$ (Hoffman et al., 2013):

$$q(Z, \theta) = q(\theta|\gamma) \prod_{n=1}^{N_X} q(z_n|\zeta_n).$$

(9)

We then maximize the ELBO with respect to $\gamma$ and $\zeta$, which is equivalent to minimizing the Kullback-Leibler (KL) divergence between the variational posterior and the true posterior:

$$\max_{\gamma,\zeta} \text{ELBO}(X, p, q) \Leftrightarrow \min_{\gamma,\zeta} \text{KL}[p(Z, \theta|X)||q(\theta, Z)].$$

(10)

We solve this problem with block coordinate descent, which leads to iteratively updating $\gamma$ and $\zeta$ as follows:

$$\zeta_{nk} \propto \prod_{v=1}^{V} \left( \beta_{kv} \right)^{x_{nv}} \exp \left[ \Psi(\gamma_k) \right],$$

(11)

$$\gamma_k = \alpha_k + \sum_{n=1}^{N_X} \zeta_{nk}. \quad (12)$$

We then approximate $E_{Z,\theta|X,\alpha_t,\beta_t}[S(X, Z, \theta)]$ with the variational posterior. Given (5) and (9), we have:

$$E_{p(Z,\theta|X)}[S(X, Z, \theta)] \approx E_{q(Z,\theta)}[S(X, Z, \theta)] = \left[ \begin{array}{c} \left( \sum_{n=1}^{N_X} \zeta_{nk} x_{nv} \right)_{kv} \\ \Psi(\gamma_k) - \Psi \left( \sum_{j=1}^{K} \gamma_j \right) \end{array} \right].$$

(13)

The variational approximation of $E_{p(Z,\theta|X)}[S(X, Z, \theta)]$ is then done in two steps:

1. Iteratively update $\zeta$ with (11) and $\gamma$ with (12),
2. $E_{p(Z,\theta|X)}[S(X, Z, \theta)] \leftarrow E_{q(Z,\theta|\gamma,\zeta)}[S(X, Z, \theta)]$ with equation (13).

As $\gamma$ and $\zeta$ are set to minimize the distance between the variational posterior and the true posterior (10) we expect that this approximation is close to the true expectation. However, as the variational model is a simplified version of the true model, there always remains a gap between the true posterior and the variational posterior.
4.3 Gibbs Online EM Applied to LDA (G-OEM)

In this section we explain how to approximate $E_{Z,\theta|X,\alpha,\beta}(S(X, Z))$ with Gibbs sampling.

4.3.1 Expectation of $S^1$.

Given (5), we have $\forall k \in [1, K], \forall v \in [1, V]$:

$$E_{Z,\theta|X,\alpha,\beta} \left( (S^1(X, Z))_{kv} \right) = E_{Z,\theta|X,\alpha,\beta} \left( \sum_{n=1}^{N_X} z_{nk} x_{nv} \right) = \sum_{n=1}^{N_X} \int_{Z,\theta} z_{nk} x_{nv} p(z_n, \theta|X, \beta, \alpha) d\theta dz = \sum_{n=1}^{N_X} x_{nv} p(z_{nk} = 1|X, \beta, \alpha).$$

We see that we only need the probability of $z$, and can thus use collapsed Gibbs sampling (Griffiths and Steyvers, 2004). We have, following Bayes rule:

$$p(z_{nk} = 1|z_{-n}, X, \beta, \alpha) \propto p(x_n|z_{nk} = 1, \beta) p(z_{nk} = 1|z_{-n}, \alpha),$$

where $z_{-n}$ is the topic assignments except index $n$. In the LDA model, each word $x_n$ is drawn from a multinomial with parameter $\beta^{z_n}$, which gives:

$$p(x_n|z_{nk} = 1, \beta) = \sum_{v=1}^{V} x_{nv} \beta_{kv}.$$

In the following, we use the notation $\beta_{x_n}^k \equiv \sum_{v=1}^{V} x_{nv} \beta_{kv}$ for the sake of simplicity. We then use the fact that the topic proportions $\theta$ has a Dirichlet($\alpha$) prior, which implies that $Z|\alpha$ follows a Dirichlet-multinomial distribution (or multivariate Pólya distribution). As a result, the conditional distribution is:

$$p(z_{nk} = 1|z_{-n}, \alpha) = \frac{N_{-n,k} + \alpha_k}{(N_X - 1) + \sum_j \alpha_j},$$

with $N_{-n,k}$ the number of words assigned to topic $k$ in the current document, except index $n$. Finally, we have the following relation (Griffiths and Steyvers, 2004):

$$p(z_{nk} = 1|z_{-n}, X, \beta, \alpha) \propto \beta_{x_n}^k \times \frac{N_{-n,k} + \alpha_k}{(N_X - 1) + \sum_j \alpha_j}.$$ (14)

We estimate $p(z_{nk} = 1|X, \beta, \alpha)$ with Gibbs sampling by iteratively sampling topic assignments $z_n$ for each word, as detailed in Algorithm 2. We average over the last quarter of samples to reduce noise in the final output. We then incorporate the output in Algorithm 1.

4.3.2 Expectation of $S^2$.

Given (5), we also have $\forall k \in [1, K], \forall v \in [1, V]$:

$$E_{Z,\theta|X,\alpha,\beta} \left( (S^2(X, Z))_k \right) = E_{Z,\theta|X,\alpha,\beta}[\log \theta_k].$$
On the one hand, we have:

\[
p(Z, \theta | X, \beta, \alpha) = p(Z | \theta, X, \beta, \alpha) p(\theta | X, \beta, \alpha) = C(\alpha) \prod_{k=1}^{K} \theta_k^{\left(\sum_{n=1}^{N_X} z_{nk}\right) + \alpha_k - 1},
\]

with \(C(\alpha) = \frac{\Gamma\left(\sum_{i=1}^{K} \alpha_i\right)}{\prod_{i=1}^{K} \Gamma(\alpha_i)}\). On the other hand:

\[
p(Z, \theta | X, \beta, \alpha) \propto p(\theta | Z, \alpha) p(Z | X, \beta).
\]

We deduce from the two identities:

\[
p(\theta | Z, \alpha) \propto K \prod_{k=1}^{K} \theta_k^{\left(\sum_{n=1}^{N_X} z_{nk}\right) + \alpha_k - 1} \Rightarrow \theta | Z, \alpha \sim \text{Dirichlet} \left(\alpha + \sum_{n=1}^{N_X} z_n\right).
\]

Finally, the expectation is:

\[
\mathbb{E}_{Z, \theta | X, \beta, \alpha} \left( [S^2(X, Z)]_k \right) = \mathbb{E}_{Z, \theta | X, \beta, \alpha} \left( \log \theta_k \right)
\]

\[
= \mathbb{E}_{Z | X, \beta, \alpha} \left[ \mathbb{E}_{\theta | Z, \alpha} \left[ \log \theta_k \right] \right]
\]

\[
= \mathbb{E}_{Z | X, \beta, \alpha} \left[ \Psi \left( [\alpha(s)]_k + \sum_{n=1}^{N_X} z_{nk} \right) - \Psi \left( \sum_{i=1}^{K} [\alpha(s)]_i + N_X \right) \right],
\]

as the distribution of \(\theta | Z, \alpha\) is \(\text{Dirichlet} \left(\alpha + \sum_{n=1}^{N_X} z_n\right)\). We use the values of \(z\) sampled with Algorithm 2 to estimate this expectation. More precisely, keeping notations of Algorithm 2:

\[
\mathbb{E}_{Z | X, \beta, \alpha} \left[ \Psi \left( [\alpha(s)]_k + \sum_{n=1}^{N_X} z_{nk} \right) \right] \approx \frac{1}{P} \sum_{t=1}^{P} \Psi \left( [\alpha(s)]_k + \sum_{n=1}^{N_X} z_{nk}^{(t)} \right).
\]

### 4.4 Bayesian Approach

In a Bayesian setting, we consider \(\beta\) as a random variable, with \(\beta \sim \text{Dirichlet}(b 1)\), with \(b \in \mathbb{R}\) and \(1 \in \mathbb{R}^V\) denotes the vector whose all entries equal 1. The variational distribution of the global parameter \(\beta\) is then set to \(q(\beta^k | \lambda^k) = \text{Dirichlet}(\lambda^k)\), with \(\lambda^k \in \mathbb{R}^V \forall k = 1, \ldots, K\).

The main difference with the frequentist methods above (G-OEM and V-OEM) is to optimize the ELBO with respect to the variational parameters \((\lambda^k)_k\). In practice, it is equivalent to replace \(\beta^k_v\) by \(\exp \left[ \mathbb{E}_q[\log \beta^k_v] \right]\) in all the updates above (i.e., in Equation (11) for V-OEM and in Equation (14) for G-OEM). The variational parmater \(\lambda^k \in \mathbb{R}^V\) is updated with stochastic gradient on the ELBO, which gives, at iteration \(t\):

\[
\lambda^k(t + 1) = \rho_t \lambda^k(t) + (1 - \rho_t) \hat{\lambda}^k,
\]

with \(\hat{\lambda}^k \in \mathbb{R}^V\), \(\lambda^k_0 = b + D \mathbb{E}_q \left[ S^1_{kv} \right]\), where \(D\) is the total number of documents in the dataset and \(b\) is the prior on \(\beta^k\) (Hoffman et al., 2013).
Algorithm 2: Gibbs sampling scheme to approximate $p(z_{nk} = 1|X, \beta, \alpha)$

**Input:** $\beta$, $\alpha$, $X$.

**Initialization:** $Z_0^n \sim \text{Mult}([\beta^k_{xn}]_{k=1,\ldots,K})$, with $\beta^k_{xn} = \frac{\beta^k_n}{\sum_j \beta^k_j} \forall n \in [1, N_X]$.

**for** $t = 1, 2, \ldots, P$ **do**

Compute random permutation $\sigma^t$ on $[1, N_X]$,

**for** $n = 0, 1, \ldots, N_X$ **do**

- Set $Z^t_{-n} = \{(z^t_{\sigma^t(i)})_{1 \leq i < n}, (z^t_{\sigma^t(i)})_{n \leq i \leq N_X}\}$,
- Compute $\forall k, p(z_{\sigma^t(n)k} = 1|Z^t_{-n}, X, \beta, \alpha)$ with Equation (14),
- Sample $z^t_{\sigma^t(n)} \sim \text{Mult} \big[p(z^t_{\sigma^t(n)}|Z^t_{-n}, X, \beta, \alpha)\big]$,

**end for**

**end for**

**for** $n = 0, 1, \ldots, N_X$ **do**

Set $Z^t_{-n} = \{(z^t_i)_{1 \leq i < n}, (z^t_i)_{n \leq i \leq N_X}\}$ for $t \geq \frac{3}{4}P$,

$p(z_{nk} = 1|X, \beta, \alpha) \leftarrow \frac{4}{P} \sum_{t = \frac{3}{4}P}^P p(z_{nk} = 1|Z^t_{-n}, X, \beta, \alpha)$

**end for**

**Output:** $\forall k, \forall n: (z^t_{nk})_{t=1,\ldots,P}, p(z_{nk} = 1|X, \beta, \alpha)$.

5. Application to Hierarchical Dirichlet Process (HDP) (Teh et al., 2006)

The HDP model is a generative process to model documents from an infinite set of topics $\beta_k$, $k = 1, 2, 3, \ldots$. Each topic is a discrete distribution of size $V$, the size of the vocabulary. Each topic is associated to a weight $\pi_k \in [0, 1]$, representing the importance of the topic in the corpus. For each document $d$, the (infinite) topic proportions $\nu_d$ are drawn from $\nu_d \sim \text{Dirichlet}(b\pi)$. We then generate words with a similar scheme to LDA scheme. More formally a corpus is generated as follows:

1. Draw an infinite number of topics $\beta_k \sim \text{Dirichlet}(\eta)$, for $k \in \{1, 2, 3, \ldots\}$;

2. Draw corpus breaking proportions $\bar{\pi}_k \sim \text{Beta}(1, \alpha)$, for $k = 1, 2, 3, \ldots$; with $\pi_k = \sigma_k(\bar{\pi})$;

3. For each document $d$:

   (a) Draw document-level topic proportions: $\nu_d \sim \text{Dirichlet}(b\pi)$;

   (b) For each word $n$ in $d$:
   i. Draw topic assignment $z_{dn} \sim \text{Multinomial}(\nu_d)$;
   ii. Draw word $w_n \sim \text{Multinomial}(\beta_{z_{dn}})$.

In practice, we set the initial number of topics to $T = 2$. We then increase the number of topics used in the corpus using Gibbs sampling and $p(z_{dn} > T|X, \eta) \propto b(1 - \sum_{i=1}^T \pi_i)$. See Section 5.2 for details.
5.1 HDP and Exponential Families

We consider an exponential family model on random variables \((X, h)\) with parameter \(\eta \in \mathcal{E} \subseteq \mathbb{R}^d\) and with density:

\[
p(X, h|\eta) = a(X, h) \exp \left[ \langle \phi(\eta), S(X, h) \rangle - \Psi(\eta) \right].
\]

In the case of HDP, an observation \(X\) is a document of length \(N_X\), where \(X = (x_1, \ldots, x_{N_X})\), \(x_n \in \{0, 1\}^V\) and \(\sum_{v=1}^{V} x_{nv} = 1\). In the frequentist approach, the parameters of the model are global, represented by \(\eta \equiv (\beta, \pi)\), where \(\beta\) represents the corpus topics, \(\pi\) represents the corpus breaking proportions. Our corpus \(\mathcal{C}\) is a set of \(D\) observations \(\mathcal{C} = (X_1, \ldots, X_D)\). For each document \(X^d\), the associated hidden variables are \(\nu_d \in [0, 1]^K\) corresponding to document-level topic proportions. For each word \(x_n\) of document \(X^d\), a hidden variable \(z_n \in \{0, 1\}^T\) is attached, corresponding to the topic assignment of word \(x_n\).

We want to find \(\phi, S, \psi\) and \(a\) such that, the joint probability is in the exponential family:

\[
p(X, Z, \nu|\beta, \pi) = a(X, Z, \nu) \exp \left[ \langle \phi(\beta, \pi), S(X, Z, \nu) \rangle - \psi(\beta, \pi) \right],
\]

given an observation \(X\) and hidden variables \(Z\) and \(\nu\). For the HDP model, we have:

\[
p(X, Z, \nu|\beta, \pi) = p(\nu|\pi) \prod_{n=1}^{N_X} p(x_n|z_n, \beta)p(z_n|\nu)
\]

\[
= W(\pi) \prod_{k \in \mathbb{N}^*} \prod_{n=1}^{N_X} (\nu_k)^{b \pi_k - 1} \prod_{k} \prod_{v} (\nu_k)^{z_{nk}} \prod_{v} (\beta_{k,v})^{x_{nv}z_{nk}}
\]

\[
= \exp \left[ -\psi(\pi) \right] \exp \left[ \sum_k \log \nu_k \left( \sum_{n=1}^{N_X} z_{nk} - 1 \right) \right]
\]

\[
\times \exp \left[ \sum_k (b \pi_k) \log \nu_k \right]
\]

\[
\times \exp \left[ \sum_{k,v} \log \beta_{k,v} \sum_{n=1}^{N_X} x_{nv}z_{nk} \right],
\]

with \(\psi(\pi) = \sum_k \log \Gamma(b \pi_k) - \log \Gamma(b)\). We deduce the exponential family setting \(\phi, S, a\):

\[
S(X, Z, \nu) = \begin{pmatrix}
S^1_k \equiv [\log \nu_k]_k \\
S^2_{kv} \equiv \left[ \sum_{n=1}^{N_X} z_{nk}x_{nv} \right]_{kv}
\end{pmatrix}
\]

\[
\phi(\beta, \pi) = \begin{pmatrix}
\phi^1_k \equiv [b \pi_k]_k \\
\phi^2_{kv} \equiv [\log \beta_{k,v}]_{kv}
\end{pmatrix}
\]

with

\[
a(X, Z, \nu) = \exp \left[ \sum_k \log \nu_k \left( \sum_{n=1}^{N_X} z_{nk} - 1 \right) \right].
\]
The one-to-one mapping between the sufficient statistics \( s = (s^1, s^2)\) and \((\beta, \pi)\) is defined by:

\[
(\beta, \pi)^*[s] = \begin{cases} 
\arg \max_{\beta \geq 0, \pi \geq 0} \phi(\beta, \pi), s - \psi(\beta, \pi) \\
\text{s.t. } \beta^\top \mathbf{1} = 1,
\end{cases}
\]

where \( \mathbf{1} \) denotes the vector whose all entries equal 1.

With the same computation than LDA, \( \beta^*(s)_{kv} \equiv \left[ s^2_{kv} \sum_{j} s^2_{kj} \right] \). We find \( \pi^*(s) \) by solving:

\[
\pi^*(s) = \arg \max_{1 \geq \pi \geq 0} \sum_{k=1}^{K} (b\pi_k s^1_k - \log \Gamma(b\pi_k)) + \log \Gamma(b \sum_k \pi_k),
\]

which gives:

\[
\Psi(b\pi^*(s)_k) - \Psi(b \sum_i \pi^*(s)_i) = s^1_k.
\]

where \( \Psi : x \mapsto \frac{\partial}{\partial x} \log \Gamma(x) \) is the digamma function. We estimate \((b\pi)^*\) with the fixed point iteration which consists in repeating the following update until convergence:

\[
(b\pi)^{\text{new}}_k = \Psi^{-1} \left( \Psi \left( \sum_i (b\pi_i)^{\text{old}} \right) + s^1_k \right).
\]

Finally, \((\beta, \pi)^*[s]\) satisfies \forall (k, v):

\[
\begin{aligned}
\beta^*(s)_{kv} &= \left[ s^2_{kv} \sum_{j} s^2_{kj} \right] \\
\Psi(b\pi^*(s)_k) - \Psi(b \sum_i \pi^*(s)_i) &= s^1_k.
\end{aligned}
\]

5.2 Inference with Online EM

In this section, we explain how to approximate \( \mathbb{E}_{Z, \nu | X, \eta} [S(X, Z, \nu)] \) with Gibbs sampling from a frequentist and a Bayesian perspective. In particular, as the total number of topics is infinite, we need to keep track of the previously used topics and iteratively extend the number of topics considered.

5.2.1 Gibbs Online EM (G–OEM)

In our frequentist G–OEM approach, \( \eta \) is a parameter. The Gibbs sampling scheme to approximate \( \mathbb{E}_{Z, \nu | X, \eta} [S(X, Z, \nu)] \) is different from LDA and a probability of adding a new topic to the current list is computed at each iteration, as explained below.

5.2.2 Expectation of \( S^1 \).

We have:

\[
\mathbb{E}_{Z, \nu | X, \eta} [S^1(X, Z, \nu)]_k = \mathbb{E}_{Z, \nu | X, \eta} [\log \nu_k] \\
= \mathbb{E}_{Z | X, \eta} \left[ \psi \left( b\pi_k + \sum_{n=1}^{N_X} z_{nk} \right) \right] - \psi \left( b \sum \pi_i + N_X \right),
\]
and we use the values of $z$ sampled with Gibbs sampling to compute:

$$
\mathbb{E}_{Z|X,\eta} \left[ \psi \left( b\pi_k + \sum_{n=1}^{N_X} z_{nk} \right) \right] \approx \frac{1}{P} \sum_{t=1}^{P} \mathbb{E}_{Z|X,\eta} \left[ \sum_{n=1}^{N_X} z_{nk} \right]
$$

5.2.3 **EXPECTATION OF $S^2$.**

We have:

$$
E_{Z,\nu|X,\eta}[S^2(X,\nu)]_{kv} = E_{Z,\nu|X,\eta} \left[ \sum_{n=1}^{N_X} z_{nk}x_{nv} \right] = \sum_{n=1}^{N_X} x_{nv}p(z_{nk} = 1|X,\eta)
$$

5.2.4 **SAMPLING $z|X,\eta$.**

If $T$ is the current number of topics, we have:

$$
p(z_{nk} = 1|z^{-n},X,\eta) \propto \left( N_{k}^{-n} + b\pi_k \right) \times \left( N_{k}^{-n} + b\pi_k \right) + \beta_{k,x_n},
$$

and the probability of sampling a new topic is given by:

$$
p(z_n > T|z^{-n},X,\eta) \propto b \left( 1 - \sum_{t=1}^{T} \pi_k \right) / V.
$$

When a new topic is generated, we initialize the probability $\pi_{T+1}$ with $\bar{\pi}_{T+1} \sim \text{Beta}(1,\alpha)$ and $\pi_{T+1} = \bar{\pi}_{T+1} \prod_{t=1}^{T} (1 - \bar{\pi}_t)$.

5.2.5 **BAYESIAN APPROACH: VARGIBBS** (WANG AND BLEI, 2012)

In a Bayesian settings where $\beta^k \sim \text{Dirichlet}(\eta)$; $q(\beta^k|\lambda) = \text{Dirichlet}(\lambda^k)$ and $\pi_k \sim \text{Beta}(1,a)$; $q(\pi_k|a_k,b_k) = \text{Beta}(a_k,b_k)$, the sampling scheme is different as we also sample $\pi$ and an auxiliary variable $s_{dk}$ corresponding to the number of “tables” serving “dish” $k$ in “restaurant” $d$ (in the formulation of HDP as a Chinese restaurant process; see Wang and Blei (2012) for details).

**Sampling $z$:**

$$
p(z_{nk} = 1|z^{-n},\lambda,\pi) \propto \left( N_{dk}^{-n} + b\pi_k \right) \frac{N_{k}^{-n} + \lambda_{kx_n}}{N_{k}^{-n} + \sum_v \lambda_{kv}}.
$$

**Sampling $s$:**

$$
p(s_{dk}|N_{dk},b\pi_k) = \frac{\Gamma(b\pi_k)}{\Gamma(b\pi_k + N_{dk})} S(N_{dk},s_{dk})(b\pi_k)^{s_{dk}},
$$

with $S(n,m)$ are unsigned Stirling number of the first kind.

**Sampling $\pi$:**

$$
p(\bar{\pi}_k) \propto \bar{\pi}_k^{\alpha_k - 1 + \sum_{d \in S} s_{dk}} (1 - \bar{\pi}_k)^{b_k - \alpha + \sum_{d \in S} \sum_{j=k+1}^{\infty} s_{dj}}
$$
We then set:

\[
\begin{align*}
\hat{\lambda}_{kv} &= \eta + D \sum_{n=1}^{N} z_{nk} x_{nv} \\
\hat{a}_k &= 1 + D s_{dk} \\
\hat{b}_k &= \alpha + D \sum_{j=k+1}^{\infty} s_{dj}
\end{align*}
\]

(18)

and \( \lambda^{t+1} = (1 - \rho_t) \lambda^t + \rho_t \hat{\lambda}; \ a^{t+1} = (1 - \rho_t) a^t + \rho_t \hat{a}; \ b^{t+1} = (1 - \rho_t) b^t + \rho_t \hat{b} \).

In practice, for each document we sample the hidden variables \( z \) for each word and compute the topic counts \( N_{dk} \) for topic \( k \) in document \( d \), then we sample the variable \( s \). Finally, we perform the online EM algorithm by making the approximation \( \mathbb{E}_{p(h|X,\eta)}[S(X, h)] \approx \mathbb{E}_{q(h)}[S(X, h)] \), which corresponds to equation (18). Note that in this Bayesian approach, the parameters \( (\lambda, a, b) \) represent the distribution parameters of the random variables \( \beta \) and \( \pi \).

### 6. Evaluation

We evaluate our method by computing the likelihood on held-out documents, that is \( p(X|\beta, \alpha) \) for any test document \( X \). For LDA, the likelihood is intractable to compute. We approximate \( p(X|\beta, \alpha) \) with the “left-to-right” evaluation algorithm (Wallach et al., 2009) applied to each test document. This algorithm is a mix of particle filtering and Gibbs sampling. On any experiments, this leads essentially to the same log-likelihood than Gibbs sampling with sufficiently enough samples—e.g., 200. In the following, we present results in terms of log-perplexity, defined as the opposite of the log-likelihood \(-\log p(X|\eta)\). The lower the log-perplexity, the better the corresponding model. In our experiments, we compute the average test log-perplexity on \( N_t \) documents. We compare eight different methods:

- **G-OEM** (our main algorithm): Gibbs online EM. Online EM algorithm with Gibbs estimation of the conditional distribution \( p(h|X, \eta) \) (Algorithm 2). Frequentist approach and step-size \( \rho_i = 1/\sqrt{i} \);
- **V-OEM++**: variational online EM (also a new algorithm). Online EM algorithm with variational estimation of the conditional distribution \( p(h|X, \eta) \), augmented with inference boosting from Section 3.3. Frequentist approach and step-size \( \rho_i = 1/\sqrt{i} \);
- **OLDA**: online LDA (Hoffman et al., 2010). Bayesian approach which maximizes the ELBO from Section 3.4, with natural stochastic gradient ascent and a step-size \( \rho_i = 1/\sqrt{i} \);
- **VarGibbs**: Sparse stochastic inference for LDA (Mimno et al., 2012). This method also maximizes the ELBO but estimates the variational expectations \( q(Z, \theta) \) with Gibbs sampling instead of iterative maximization of variational parameters—see Section 4.2;
- **SVB**: streaming variational Bayes (Broderick et al., 2013). A variational Bayesian equivalent of V-OEM with step-size \( \rho_i = 1/i \);
• **SPLDA**: single pass LDA (Sato et al., 2010). The difference with V-OEM++ is that \( \rho_i = 1/i \) and the updates in \( \alpha \) done with a Gamma prior (see Appendix C);

• **SGS**: streaming Gibbs sampling (Gao et al. 2016). This method is related to G-OEM with \( \rho_i = 1/i \). In this method, \( \alpha \) is not optimized and set to a constant \( C \alpha \). For comparison purposes, for each dataset, we set \( C \alpha \) to be the averaged final parameter \( \hat{\alpha} \) obtained with G-OEM on the same dataset: \( C \alpha = \frac{1}{K} \sum_k \hat{\alpha}_k \). For each observation, only the last Gibbs sample is considered, leading to extra noise in the output;

• **LDS**: Stochastic gradient Riemannian Langevin dynamics sampler (Patterson and Teh. 2013). The authors use the Langevin Monte Carlo methods on probability simplex and apply their online algorithm to LDA. For this method and only this method, we set to \( P = 200 \) the number of internal updates.

For existing variational methods—OLDA, SVB, SPLDA—\( \beta \) is a random variable with prior \( q(\beta) \). We estimate the likelihood \( p(X|\hat{\beta}, \alpha) \) with the “left-to-right” algorithm by setting \( \hat{\beta} = E_q[\beta] \) for Bayesian methods. For simplicity, we only present our results obtained with G-OEM and V-OEM++. Indeed, the inference boost presented in Section 4 is only beneficial for V-OEM. A detailed analysis is presented in Appendix A.1.

### 6.1 Explicit Links for LDA

In this section, we propose to make the links between the methods listed above explicit, using the framework described in Section 4 for the particular LDA model. We present in Table 1 a summary of the compared method.

#### 6.1.1 Category

In the frequentist approach, \( \beta \) is a parameter and is updated with Equation (8), as the “M-step” in online EM.

In a Bayesian setting, \( \beta \) is a random variable with prior \( \beta \sim \text{Dirichlet}(b1) \), with \( b \in \mathbb{R} \) and \( 1 \in \mathbb{R}^V \) denotes the vector whose all entries equal 1. The variational distribution of the global parameter \( \beta \) is then set to \( q(\beta|\lambda^k) = \text{Dirichlet}(\lambda^k) \), with \( \lambda^k \in \mathbb{R}^V \forall k = 1, \ldots, K \). The variational parameter \( \lambda^k \) is updated by maximizing the ELBO with stochastic gradient ascent (Equation (15)).

<table>
<thead>
<tr>
<th>Method</th>
<th>Category</th>
<th>Gibbs Sampling</th>
<th>Step-size</th>
<th>Update for ( \alpha )</th>
</tr>
</thead>
<tbody>
<tr>
<td>G-OEM</td>
<td>Frequentist</td>
<td>Freq</td>
<td>Fixed</td>
<td>Fixed point</td>
</tr>
<tr>
<td>V-OEM</td>
<td>Frequentist</td>
<td>Varyational</td>
<td>Free</td>
<td>Fixed point</td>
</tr>
<tr>
<td>OLDA</td>
<td>Bayesian</td>
<td>Varyational</td>
<td>Free</td>
<td>Gradient ascent</td>
</tr>
<tr>
<td>VARGIBBS</td>
<td>Bayesian</td>
<td>Gibbs Sampling</td>
<td>Free</td>
<td>( \alpha ) fixed</td>
</tr>
<tr>
<td>SVB</td>
<td>Bayesian</td>
<td>Varyational</td>
<td>Fixed: 1/t</td>
<td>Gradient ascent</td>
</tr>
<tr>
<td>SPLDA</td>
<td>Frequentist</td>
<td>Varyational</td>
<td>Fixed: 1/t</td>
<td>Gamma prior</td>
</tr>
<tr>
<td>SGS</td>
<td>Frequentist</td>
<td>Gibbs Sampling</td>
<td>Fixed: 1/t</td>
<td>( \alpha ) fixed</td>
</tr>
</tbody>
</table>

Table 1: Comparison of existing methods for LDA.
6.1.2 Estimation of $\mathbb{E}_{Z|X,\eta}[S(X, Z)]$

For LDA, the expectation $\mathbb{E}_{Z|X,\eta}[S(X, Z)]$ can either be estimated with Gibbs sampling—using Equation (14)—or with variational approximation—using Equation (10).

6.1.3 Step-size

Some of the methods listed above (SVB, SPLDA and SGS) are incremental, which means the sufficient statistics are incrementally aggregated $s_t = s_{t-1} + \mathbb{E}_{Z|X_t,\eta}[S(X_t, Z_t)]$. For LDA, it exactly corresponds to a step-size $\rho_t = 1/t$ in the online EM setting, even though the link is not explicit in the corresponding papers.

For the other listed methods, the step-size exponent $\kappa$ is chosen arbitrarily in $[0.5, 1)$, with $\rho_t = 1/t^\kappa$. However, results are mostly presented with $\kappa = 1/2$ and $\rho_t = 1/\sqrt{t}$.

6.2 General Settings

6.2.1 Initialization

We initialize randomly $\eta \equiv (\beta, \alpha)$. For a given experiment, we initialize all the methods with the same values of $(\beta, \alpha)$ for fair comparison, except SPLDA that has its own initialization scheme—see Sato et al. (2010) for more details.

6.2.2 Minibatch

We consider minibatches of size 100 documents for each update in order to reduce noise (Liang and Klein 2009). In the case of online EM in Equation (3), we estimate an expectation for each observation of the minibatch. We update the new sufficient statistics $s$ towards the average of the expectations over the minibatch. We do the same averaging for all the presented methods.

6.2.3 Number of Local Updates

For all the presented methods, we set the number of passes through each minibatch to $P = 20$. For $\text{G-OEM}$, this means that we perform 20 Gibbs sampling for each word of the minibatch. All other methods access each document 20 times (e.g., 20 iterations of variational inference on each document). For $\text{G-OEM}$, inference with larger values for $P$ (e.g., $P = 50$ or $P = 100$) leads to very similar results.

6.2.4 Datasets

We apply the methods on six different datasets, summarized in Table 2 ($\overline{N_X}$ is the average length of documents). Following Blei et al. (2003), the synthetic dataset has been generated from 10 topics and the length of each document drawn from a Poisson(60). The 10 topics are inferred with online LDA (Hoffman et al. 2010) from 50,000 reviews of the IMDB dataset with a vocabulary size of 10,000. We only consider the entries of the 1,000 most frequent words of this dataset that we normalize to satisfy the constraint $\sum_v \beta_{kv} = 1$.

The words in the datasets IMDB, Wikipedia, New York Times, Pubmed and Amazon movies are filtered by removing the stop-words and we select the most frequent words of the datasets. For the synthetic dataset, IMDB, Pubmed and Amazon movies, the size of the test sets is $N_t = 5,000$ documents. For Wikipedia and New York Times, the test sets contain $N_t = 2,000$ documents.
We run the methods on 11 different train/test splits of each dataset. For all the presented results, we plot the median from the 11 experiments as a line—solid or dashed. For the sake of readability, we only present the same plots with error bars between the third and the seventh decile in Appendix D and Appendix E.

6.2.5 Computation Time

For each presented method and dataset, the computational time is reported in Table 3. Although all methods have the same running-time complexities, coded in Python, sampling methods (G-OEM, VarGibbs and SGS) need an actual loop over all documents while variational methods (OLDA, SVB, SPLDA and V-OEM++) may use vector operations, and may thus be up to twice faster. This could be mitigated by using efficient implementations of Gibbs sampling on minibatches (Yan et al., 2009; Zhao et al., 2014; Gao et al., 2016). Note also that to attain a given log-likelihood, our method G-OEM is significantly faster and often attains log-likelihoods not attainable by other methods (e.g., for the dataset New York Times).

6.2.6 Step-size

In the following, we compare the results of our methods G-OEM and V-OEM++ with \( \kappa = 1/2 \), i.e., the step-size \( \rho_t = 1/\sqrt{t} \), without averaging. Detailed analysis of different settings of our method can be found in Appendix A. In particular, we compare different step-sizes and the effect of averaging over all iterates. We also compare the performance of OLDA with different step-sizes in Appendix A.2 and observe that results are very similar for all the step-sizes that we try. Note that for incremental methods (SVB, SPLDA, SGS), the step-size is fixed to \( \rho_t = 1/t \). For LDS, we run the method with parameters as close as possible to our method for fair comparison.

6.3 Results on LDA

Results obtained with the presented methods applied to LDA on different datasets for different values of the number \( K \) of topics are presented in Figure 1. Performance through iterations (i.e., as the number of documents increases) is presented in Figure 4. We first observe that for all experiments, our new method G-OEM performs better—often significantly—than all existing methods. In particular, it is highly robust to diversity of datasets.

---

2. Dataset described in Diao et al. (2014)
3. Data from Leskovec and Krevl (2014)
4. UCI dataset (Lichman, 2013)
6.3.1 Influence of the number of topics $K$

As shown in Figure 1, for synthetic data in plot (a), although the true number of topics is $K^* = 10$, SPLDA, OLDA, VarGibbs and SGS perform slightly better with $K = 20$, while G-OEM has the better fit for the correct value of $K$; moreover, SVB has very similar performances for any value of $K$, which highlights the fact that this method does not capture more information with a higher value of $K$. LDS performs very poorly on this dataset—for any value of $K$ the log-perplexity is around 400—and is not displayed in Figure 1 (a) for clarity.

On non-synthetic datasets in plots (b)-(f), while the log-perplexity of frequentist methods—G-OEM, V-OEM++ and SPLDA—decreases with $K$, the log-perplexity of variational Bayesian methods—OLDA and SVB—does not decrease significantly with $K$. As explained below, our interpretation is that the actual maximization of the ELBO does not lead to an improvement in log-likelihood. The hybrid Bayesian method VarGibbs—which uses Gibbs sampling for local updates ($\theta, z$) and variational updates for global parameters ($\beta, \alpha$)—performs much better than the variational Bayesian methods. Our interpretation is that the objective function maximized with VarGibbs is a much better approximation of the log-likelihood than the ELBO.

In terms of robustness, G-OEM and LDS are the only methods that do not display overfitting on any dataset. However, LDS is only competitive for the highest values of $K$—$K \geq 500$.

6.3.2 Performance Through Iterations

As shown in Figure 4 for synthetic data in plot (a), after only few dozens of iterations—few thousands of documents seen—G-OEM, V-OEM++ and VarGibbs outperform the other presented methods. Variational Bayesian methods again do converge but to a worse parameter value. On real datasets in plots (b)-(f), G-OEM and VarGibbs are significantly faster; we can indeed still observe that after around 100 iterations—10,000 documents seen—G-OEM and VarGibbs perform better than other methods on all the datasets except Pubmed, where the performances of G-OEM, V-OEM++, VarGibbs and SPLDA are similar. Note that

6.3.3 Variational vs. Sampling

Our method G-OEM directly optimizes the likelihood with a consistent approximation, and performs better than its variational counterparts SPLDA and V-OEM++ in all experiments. The hybrid method VarGibbs is less robust than G-OEM as it performs either similarly to G-OEM—for the datasets

<table>
<thead>
<tr>
<th>Method</th>
<th>IMDB</th>
<th>WIKIPEDIA</th>
<th>NYT</th>
<th>PUBMED</th>
</tr>
</thead>
<tbody>
<tr>
<td>G-OEM</td>
<td>13H</td>
<td>55H</td>
<td>30H</td>
<td>58H</td>
</tr>
<tr>
<td>V-OEM++</td>
<td>9H</td>
<td>37H</td>
<td>20H</td>
<td>54H</td>
</tr>
<tr>
<td>OLDA</td>
<td>7H</td>
<td>33H</td>
<td>8H</td>
<td>30H</td>
</tr>
<tr>
<td>VarGibbs</td>
<td>12H</td>
<td>50H</td>
<td>28H</td>
<td>54H</td>
</tr>
<tr>
<td>SVB</td>
<td>7H</td>
<td>34H</td>
<td>9H</td>
<td>30H</td>
</tr>
<tr>
<td>SPLDA</td>
<td>9H</td>
<td>37H</td>
<td>20H</td>
<td>54H</td>
</tr>
<tr>
<td>SGS</td>
<td>11H</td>
<td>48H</td>
<td>27H</td>
<td>50H</td>
</tr>
<tr>
<td>LDS</td>
<td>7H</td>
<td>17H</td>
<td>12H</td>
<td>40H</td>
</tr>
</tbody>
</table>

Table 3: Average computational time (in hours) for each method—$K = 128$. 

Online Inference for Latent Variable Models
Figure 1: Perplexity on different test sets as a function of $K$, the number of topics inferred. Best seen in color.
Figure 2: OLDA. Perplexity on different test sets as a function of $K$ for OLDA with $P = 10$ (red) and $P = 100$ (black) internal updates.


6.3.4 Frequentist vs. Bayesian

In all our experiments we observe that frequentist methods—G-OEM, V-OEM++ and SPLDA—outperform variational Bayesian methods—OLDA and SVB. As described in Section 3.4, variational Bayesian methods maximize the ELBO, which makes additional strong independence assumptions and here leads to poor results. For example, as the number $K$ of topics increases, the log-likelihood goes down for some datasets. In order to investigate if this is an issue of slow convergence, we show on Figure 2 (dotted black line) that running $P = 100$ internal updates in OLDA to get a finer estimate of the ELBO for each document may deteriorate the performance. Moreover, Figure 3 presents the evolution of the ELBO, which does always increase when $K$ increases, showing that the online methods do optimize correctly the ELBO (while not improving the true log-likelihood). See Appendix B for additional results on the convergence of the ELBO. The results are mitigated for the hybrid Bayesian method VarGibbs. The performance of this method is either similar to G-OEM and V-OEM++ or significantly worse than both G-OEM and V-OEM++.

6.3.5 Small Step-sizes vs. Large Step-sizes

SPLDA is also a variational method which is equivalent to V-OEM++, but with a step-size $1/t$, which is often slower than bigger step-sizes (Mairal 2014), which we do observe—see Appendix A.2 for a further analysis on the effect of the choice of step-sizes as $1/i^\kappa$ on G-OEM. Note that we run all the methods on a fixed (finite) number of observations. If we were to extend to infinite datasets, the difference between the step-sizes should be the speed of convergence. However, even if the number of observations is large, the gap between the step-sizes is still significant to justify the use of $1/\sqrt{t}$ for the step-size. Indeed, when considering large datasets, the contribution of each
iteration at the end of the pass over the data is squeezed by the step-size in $1/t$. When the number of observations is large enough to prevent the use of batch algorithms but still insufficient for an online algorithm to converge in one pass, a possible solution could be to consider constant step-sizes in order to converge even faster to a local maxima. As proposed, we do not have any guarantee for our methods to converge with constant step-sizes, but previous works have shown the benefits of using constant step-sizes under certain assumptions (e.g., Bach and Moulines (2013)).

### 6.4 Empirical Analysis

In this section we provide a qualitative empirical analysis on the topics extracted with the different methods. We note this is clearly a subjective analysis but it stresses the benefits of a “better” inference mechanism in terms of log-likelihood (Chang et al., 2009). Examples of eight topics extracted with G-OEM and OLDA on the IMDB dataset of movie reviews are presented in Table 4 page 30.

We first compute the KL divergence between the $K = 128$ topics extracted with G-OEM and the $K = 128$ topics extracted with OLDA. We run the Hungarian algorithm on the resulting distance matrix to assign each topic extracted with G-OEM to a single topic of OLDA. We choose manually eight topics extracted with G-OEM that are representative of the usual behavior, and display the eight corresponding topics of OLDA assigned with the above method.

We observe that the topics extracted with G-OEM are more consistent than topics extracted with OLDA: topics of G-OEM precisely describe only one aspect of the reviews while the topics of OLDA tend to mix several aspects in each topic. For instance, the words of topic 1 extracted with G-OEM are related to horror movies. The words of the corresponding topic extracted with OLDA mix horror movies—e.g., horror, scary—and ghost movies—e.g., ghost, haunt. In this OLDA topic 1, we can also observe less relevant words, like effective, mysterious, which are not directly linked with horror and ghost vocabularies. We can make the same remarks with topic 2 and topic 3, respectively related to comedy movies and romantic comedy movies. In topic 2 extracted with G-OEM, the least related

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**Figure 3:** Dataset: IMDB. Evidence Lower BOund (ELBO) computed on test sets (20 internal iterations and 1 pass over the dataset). *Left:* ELBO through iterations with $K = 128$. *Right:* ELBO as a function of the number of topics $K$. 

---

[26]
(a) Synthetic, $K = 10$

(b) IMDB, $K = 128$

(c) Wikipedia, $K = 128$

(d) New York Times, $K = 128$

(e) Pubmed, $K = 128$

(f) Amazon movies, $K = 128$

Figure 4: Perplexity through iterations on different test sets with the presented methods. Best seen in color.
words to *comedy* are names of characters/actors—i.e., *steve* and *seth*—while the words not related to *comedy* in topic 25 of OLDA are more general, belonging to a different lexical field—e.g., *sport, site, progress, brave, definition*. In topic 3 of G-OEM, all the presented words are related to *romantic comedy* while in topic 3 of OLDA, the words *old, hard* and *review* are not related to this genre.

We also observe that G-OEM extracts strongly “qualitative” topics—topic 4 and topic 5—which is not done with OLDA. Indeed, it is difficult to group the top words of topic 4 or topic 5 of OLDA in the same lexical field. Except *dialogue* and *suppose*, all the top words of topic 4 of G-OEM are negative words. These two words may appear in a lot of negative sentences, leading to a high weight in this topic. In topic 5 of G-OEM, the words *absolutely* and *visual* are non strictly positive words while the thirteen other words in this topic convey a positive opinion. The word *absolutely* is an adverb much more employed in positive sentences than negative or neutral sentences, which can explain its high weight in topic 5.

The topic 6 of both G-OEM and OLDA can be considered as a “junk” topic, as for both method, most of its top words are contractions of modal verbs or frequent words—e.g., *didn’t, isn’t, wait, bad*. The contractions are not filtered when removing the stop words as they are not included in the list of words removed\(^1\).

For both G-OEM and OLDA, the top words of topic 7 are general words about movies. These words are usually employed to describe a movie as a whole—e.g., *narrative, filmmaker*.

Finally, the top words of topic 8 of G-OEM are related to the situation of the scenes. We could not find such topic in the other presented methods and we can see that the top words of topic 8 of OLDA—supposedly close to topic 8 of G-OEM—are related to *family* movies. Each word of topic 8 of G-OEM—except *group* and *beautiful*—are related to a *spatial location*, and may help answer the question “*where does the scene take place?*”.

### 6.5 Results on HDP

For the HDP model, we compare our G-OEM method to the Bayesian VarGibbs (Wang and Blei, 2012) method. We set the initial number of topics to \(T = 2\). We present in Figure 5 results obtained with G-OEM and VarGibbs applied to both LDA and HDP. Results with error bars are presented in Appendix F. For both LDA and HDP, G-OEM outperforms the Bayesian method VarGibbs.

### 7. Conclusion

We have developed an online inference scheme to handle intractable conditional distributions of latent variables, with a proper use of local Gibbs sampling within online EM, that leads to significant improvements over variational methods and Bayesian estimation procedures. Note that all methods for the same problem are similar (in fact a few characters away from each other); ours is based on a proper stochastic approximation maximum likelihood framework and is empirically the most robust. It would be interesting to explore distributed large-scale settings (Broderick et al., 2013; Yan et al. 2009; Gao et al. 2016) and potentially larger (e.g., constant) step-sizes that have proved efficient in supervised learning (Bach and Moulines, 2013).

---

1. See NLTK toolbox (Bird et al., 2009) for the exhaustive list of stop words.
Figure 5: Perplexity through iterations on different test sets with G-OEM and VarGibbs applied to both LDA and HDP. Best seen in color.

Acknowledgments

We would like to thank David Blei, Olivier Cappé, Nicolas Flammarion and John Canny for helpful discussions related to this work.
Table 4: Comparison of topics extracted on IMDB dataset, $K = 128$—15 top words of eight topics extracted with G-OEM and OLDA.

<table>
<thead>
<tr>
<th>#</th>
<th>Topic 1</th>
<th>Topic 2</th>
<th>Topic 3</th>
<th>Topic 4</th>
<th>Topic 5</th>
<th>Topic 6</th>
<th>Topic 7</th>
<th>Topic 8</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>VIOLENCE</td>
<td>COMEDY</td>
<td>ROMANTIC</td>
<td>BAD</td>
<td>BRILLIANT</td>
<td>DIDN’T</td>
<td>NARRATIVE</td>
<td>TOWN</td>
</tr>
<tr>
<td>2</td>
<td>VIOLENT</td>
<td>FUNNY</td>
<td>COMEDY</td>
<td>WORST</td>
<td>PERFECT</td>
<td>I’VE</td>
<td>ULTIMATE</td>
<td>LOCAL</td>
</tr>
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Appendix A. Gibbs/Variational Online EM Analysis

In this section we evaluate the proposed methods G-OEM and V-OEM with different settings in terms of step-sizes, averaging outputs and boosting internal updates.

A.1 Effect of Inference Boosting on G-OEM and V-OEM

The effect of the inference boost as described in Section 3.3 on G-OEM and V-OEM with synthetic and IMDB datasets is presented in Figure 6 and in Figure 7. It leads to a minor improvement for G-OEM++ and a significant one for V-OEM++.

A.2 Step-sizes and Averaging

We apply G-OEM with different stepsizes $\rho_i = \frac{1}{i^\kappa}$. Note that because we average sufficient statistics, there is no needed proportionality constants. We first compare the performance of the last iterate $\eta_N$ (without averaging) and the average of the iterates $\bar{\eta}_N = \frac{1}{N} \sum_{i=0}^{N} \eta_i$ (with averaging) for different values of $\kappa$.

Results are presented in Figure 8 on the synthetic data and in Figure 9 on the IMDB dataset. For $\kappa \in [0, \frac{1}{2}]$, averaging improves the performance while for $\kappa \in [\frac{1}{2}, 1]$, averaging deteriorates the performance. For $\kappa = \frac{1}{2}$, averaging is only slightly beneficial on IMDB dataset. For constant stepsizes $\kappa = 0$ the averaging improves significantly the performance, as the iterates do not converge and tend to oscillate around a local optimum (Bach and Moulines, 2013). We can expect the same effect for $\kappa \in [0, \frac{1}{2}]$ as the function $n \mapsto \frac{1}{n^\kappa}$ deacreases slowly for such values of $\kappa$. For $\kappa \in [\frac{1}{2}, 1]$, the stochastic gradient ascent scheme is guaranteed to converge to a local optimum (Bottou, 1998). The averaging then deteriorates the performance as it incorporates the first iterates, which gets the last iterate away from local optimum. However, the stepsize $1/i$ ($\kappa = 1$) is not competitive. The performance with $\kappa = 0.75$ is only slightly better on IMDB dataset. The setting $\kappa = \frac{1}{2}$ represents a good balance between first and last iterates. For this step-size, performances with or without averaging are similar but results without averaging seem to be more stable, hence our choice for all our other simulations.

We also apply OLDA with different step-sizes $\rho_t = \tau/t^\kappa$ for different values of $\tau, \kappa$. Results are presented in Figure 10 without error bars and in Figure 11 with error bars. For OLDA, results are very similar for any step-size.
Figure 6: G-OEM. Perplexity on different test sets as a function of the number of topics $K$ for regular EM and boosted EM ($++$). We observe that for almost all datasets, there is no significant improvement when boosting the inference. Our interpretation is that each Gibbs sample is noisy and does not provide a stable boost. Best seen in color.
Figure 7: V-OEM. Perplexity on different test sets as a function of the number of topics $K$ for regular EM and boosted EM (++)

We observe that boosting inference improves significantly the results on all the datasets excepted on Wikipedia where V-OEM and V-OEM++ have similar performances. The variational estimation of the posterior is finer and finer through iterations. When updating the parameters at each iteration of the posterior estimation, the inference is indeed boosted. Best seen in color.
Figure 8: Dataset: synthetic. Perplexity on different test sets as a function of the exponent $\kappa$—the corresponding stepsize is $\rho_i = \frac{1}{i^{\kappa}}$—for G-OEM with averaging (left) and without averaging (right). The number of topics inferred $K$ goes from 5 (the lightest) to 20 (the darkest). Best seen in color.

Figure 9: Dataset: IMDB. Perplexity on different test sets as a function of the exponent $\kappa$—the corresponding stepsize is $\rho_i = \frac{1}{i^{\kappa}}$—for G-OEM with averaging (left) and without averaging (right). The number of topics inferred $K$ goes from 8 (the lightest) to 128 (the darkest). Best seen in color.
Figure 10: Evolution of perplexity on different test sets as a function of the number of documents analyzed. For OLDA, we compare the performance with different step-sizes \( \rho_t = \tau / t^\kappa \) for different values of \( \tau, \kappa \). Solid line: \( \kappa = 1/2 \); Dashed line: \( \kappa = 1 \).

Figure 11: Evolution of perplexity on different test sets as a function of the number of documents analyzed, with error bars. For OLDA, we compare the performance with different step-sizes \( \rho_t = \tau / t^\kappa \) for different values of \( \tau, \kappa \). Solid line: \( \kappa = 1/2 \); Dashed line: \( \kappa = 1 \).
Appendix B. Evolution of the ELBO

Figure 12 presents the evolution of the ELBO for online LDA (OLDA) and SVB on different test sets. We compute the ELBO on test documents as described by Hoffman et al. (2010). This plot helps us to observe that even if the ELBO reaches a local maximum (i.e., it stabilizes), the quality of the model in terms of perplexity is not controllable. We can also see in Figure 12 that the ELBO is much better optimized with $K = 128$ than with other values of $K$ for both SVB and OLDA, that is, as expected, latent variables of higher dimensionality lead to better fits for the cost function which is optimized. However, for several datasets the performance in terms of perplexity is better with low values of $K$ ($K = 8$ or $K = 16$) than with high dimensional variables ($K = 64$ or $K = 128$).

Figure 12: Evidence Lower BOund (ELBO) computed on different test sets. Top: ELBO through iterations, with 4 passes over each dataset and 200 internal iterations. Bottom: ELBO as a function of the number of topics $K$, with 20 internal iterations and 1 pass over each dataset. Best seen in color.
In order to check if more internal iterations could help variational Bayesian methods, we present in Table 5 the values of perplexity reached by OLDA when running 4 passes over each dataset with \( P = 200 \) internal iterations 1 pass over each dataset with \( P = 20 \) internal iterations. We observe that the ELBO converges quickly to a local optimum and doing ten times more internal iterations does not change significantly the final performance.

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<th>( P = 200, 4 ) passes</th>
<th>( P = 20, 1 ) pass</th>
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<td>SVB</td>
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Table 5: Comparison of log-perplexity levels reached with OLDA and SVB on IMDB dataset.

### Appendix C. Updates in \( \alpha \)

In this section we compare the different types of updates for \( \alpha \). Figure 13 presents results obtained on synthetic dataset for fixed point iteration algorithm (Minka, 2000) and by putting a gamma prior on \( \alpha \) (Sato et al., 2010). We observe that the fixed point method leads to better performance for \( G\text{-OEM} \) and \( G\text{-OEM++} \). For \( V\text{-OEM} \), the gamma updates better perform for \( \kappa = \frac{1}{2} \). The performances of the gamma updates and the fixed point method are very similar for \( V\text{-OEM++} \). Note that the algorithm \( V\text{-OEM++} \) with \( \kappa = 1 \) and gamma updates on \( \alpha \) is exactly equivalent to SPLDA (Sato et al., 2010). The performance of this method can be improved by setting \( \kappa = \frac{1}{2} \) with any update on \( \alpha \).

We also observe that fixing \( \alpha \) to \( \alpha_{\text{true}} \) that generated the data does not necessarily lead to better performance.
Figure 13: Dataset: Synthetic, $K = 10$. Perplexity on different test sets for different types of updates for $\alpha$; for boosted methods, we use the same inference for $\alpha$ for local and global updates. NO: $\alpha$ is fixed and set to $\alpha_{true}$ that generated the data; FP: fixed point iteration; Gam: gamma prior on $\alpha$ (Sato et al., 2010). Best seen in color.
Appendix D. Performance with Different $K$, with Error Bars

The performance of the presented methods for different values of $K$ on the different datasets is presented in Figure 14. We plot the median from the 11 experiments as a line—solid or dashed—and a shaded region between the third and the seventh decile.

Appendix E. Performance Through Iterations, with Error Bars

The performance through iterations of the presented methods on the different datasets is presented in Figure 15. We plot the median from the 11 experiments as a line—solid or dashed—and a shaded region between the third and the seventh decile.

Appendix F. Results on HDP, with Error Bars

The performance through iterations of the G-OEM and VarGibss applied to both LDA and HDP is presented in Figure 15. We plot the median from the 11 experiments as a line—solid or dashed—and a shaded region between the third and the seventh decile.
Figure 14: Perplexity on different test sets as a function of $K$, the number of topics inferred. Same as Figure 1, but with error bars. Best seen in colors.
Figure 15: Perplexity through iterations on different test sets with the presented methods. Same as Figure 4, but with error bars. Best seen in colors.
Figure 16: Perplexity through iterations on different test sets with G-OEM and VarGibbs applied to both LDA and HDP. Best seen in color.
References


