### 10707 Deep Learning: Spring 2020

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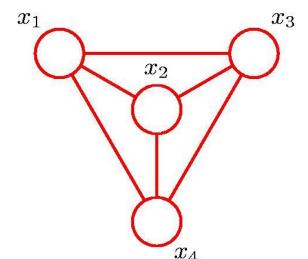
Machine Learning Department

Lecture 11:

Variational methods

### Graphical Models

Recall: graph contains a set of nodes connected by edges.



In a probabilistic graphical model, each node represents a random variable, links represent "probabilistic dependencies" between random variables.

Graph specifies how joint distribution over all random variables decomposes into a **product** of factors, each factor depending on a subset of the variables.

#### Two types of graphical models:

- **Bayesian networks**, also known as Directed Graphical Models (the links have a particular directionality indicated by the arrows)
- **Markov Random Fields**, also known as **Undirected Graphical Models** (the links do not carry arrows and have no directional significance).

## Algorithmic pros/cons of latent-variable models (so far)

#### RBM's

- S Hard to draw samples (In fact, #P-hard provably, even in Ising models)
- Second Easy to sample posterior distribution over latents



#### **Directed models**

S Easy to draw samples



Mard to sample posterior distribution over latents



#### Canonical tasks with graphical models

#### <u>Inference</u>

Given values for the parameters  $\theta$  of the model, *sample/calculate* marginals (e.g. sample  $p_{\theta}(x_1), p_{\theta}(x_4, x_5), p_{\theta}(z|x)$ , etc.)

#### **Learning**

Find values for the parameters  $\theta$  of the model, that give a *high likelihood* for the observed data. (e.g. canonical way is solving maximum likelihood optimization

$$\max_{\theta \in \Theta} \sum_{i=1}^{n} \log p(x_i)$$

Other methods exist, e.g. method of moments (matching moments of model), but less used in deep learning practice.

#### Canonical tasks with graphical models

#### <u>Inference</u>

Inference is hard in undirected fully-observable models (due to partition function); easy in fully-observable Bayesian nets. It's easy for RBM's, hard for latent-variable Bayesian nets (again, implicit normalizing factor is hard.)

#### **Learning**

We will derive "iterative"/"incremental" learning algorithms, using inference algorithms as subroutines.

We will, in particular see how a technique called "variational methods" can be used. (Next time, we see how MCMC methods can be used.)

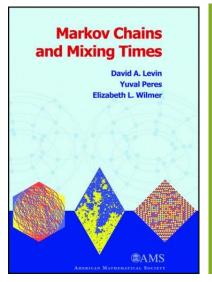
### Inference

### Algorithmic approaches

When faced with a difficult to calculate probabilistic quantity (partition function, difficult posterior), there are two families of approaches:

#### MARKOV CHAIN MONTE CARLO

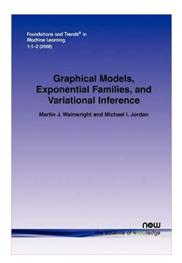
- \*Random walk w/ equilibrium distribution the one we are trying to sample from.
  - ❖ Well studied in TCS.





#### **VARIATIONAL METHODS**

- Based on solving an optimization problem.
- Very popular in practice.
- Comparatively poorly understood



# Part I: Inference in undirected graphical models

Though not always true, calculating marginals is often reducible to calculating partition functions.

**Simple example**: Ising models

$$P_{\theta}(\mathbf{x}) = \frac{1}{\mathcal{Z}(\theta)} \exp\left(\sum_{ij \in E} x_i x_j \theta_{ij} + \sum_{i \in V} x_i \theta_i\right)$$

Partition fn of an appropriately modified Ising model

$$P_{\theta}(x_{k} = 1) = \frac{\sum_{x_{-i} \in \{-1,1\}^{n-1}} \exp(\theta_{k} + \sum_{kj \in E} x_{k} \theta_{j} + \sum_{ij \in E, i \neq k, j \neq k} x_{i} x_{j} \theta_{ij} + \sum_{i \neq k} \theta_{i} x_{i})}{\sum_{x \in \{-1,1\}^{n}} \exp(\sum_{ij} x_{i} x_{j} \theta_{ij} + \sum_{i} \theta_{i} x_{i})}$$

Partition fn of original Ising model

# Part I: Inference in undirected graphical models

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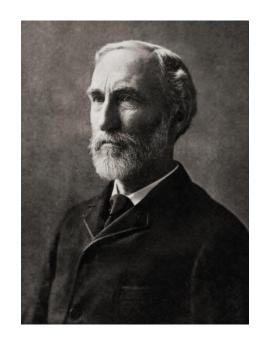
Formal term for when sampling reduces to calculating partition functions: *self-reducible problems*.

How do we calculate/approximate the partition function?

**Gibbs variational principle**: Let  $p(x) = \frac{1}{Z} \exp(E(x))$  be a distribution over a domain  $\mathcal{X}$ . Then, Z is the solution to the following optimization problem:  $\log Z = \max_{p: \text{ distribution over } \mathcal{X}} H(p) + \mathbb{E}_{x \sim p}[E(x)]$ 

Find the distribution that has both high entropy, and high expected energy value

$$H(p) := -\sum_{x \in \mathcal{X}} p(x) \log p(x)$$

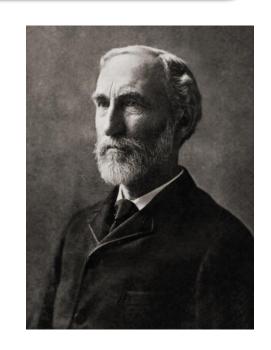


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Proof: 
$$0 \le KL(q \mid\mid p) = \mathbb{E}_q \log q - \mathbb{E}_q \log p$$

$$= -H(q) - \mathbb{E}_{x \sim q}[E(x)] + \log Z$$

$$H(q) + \mathbb{E}_{x \sim q}[E(x)] \le \log Z$$
Hence,  $\max_{q: \text{ distribution over } \mathcal{X}} H(q) + \mathbb{E}_{x \sim q}[E(x)] \le \log Z$ 



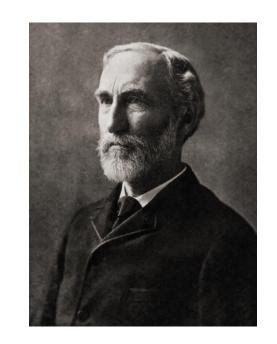
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$$H(q) + \mathbb{E}_{x \sim q}[E(x)] \le \log Z$$
Equality is attained if  $p = q$ :  $KL(q \mid\mid p) = 0$ , so

 $H(q) + \mathbb{E}_{x \sim q}[E(x)] = \log Z$ 



**Gibbs variational principle**: Let  $p(x) = \frac{1}{Z} \exp(E(x))$  be a distribution over a domain  $\mathcal{X}$ . Then, Z is the solution to the following optimization problem:  $\log Z = \max_{q: \text{ distribution over } \mathcal{X}} H(q) + \mathbb{E}_{x \sim q}[E(x)]$ 

Hence, we've reduced calculating partition function to an optimization problem!

But, there is a **serious issue**: how do we solve an optimization over the set of distributions over X?

Even if  $\mathcal{X}$  is a really simple domain, e.g.  $\mathcal{X} = \{\pm 1\}^n$ , the trivial way to solve the problem would involve introducing a variable q(x),  $\forall x \in \{\pm 1\}^n$ : there are  $2^n$  of them.

In fact, you can't be clever – there are results showing this can be #P hard even for Ising models!

**Gibbs variational principle**: Let  $p(x) = \frac{1}{Z} \exp(E(x))$  be a distribution over a domain  $\mathcal{X}$ . Then, Z is the solution to the following optimization problem:  $\log Z = \max_{q: \text{ distribution over } \mathcal{X}} H(q) + \mathbb{E}_{x \sim q}[E(x)]$ 

What can we do to try to approximate this expression?

**Inspiration from physics**: solve a *simpler* optimization problem over a *restricted class* of distributions we can explicitly parametrize.

**Typical example:** mean-field approximation

Consider again  $\mathcal{X} = \{\pm 1\}^n$ . A *product distribution* depends on n parameters only: since  $p(x) = \prod_i p_i(x_i)$ , for each  $i \in [n]$ , we only need to specify  $p_i(x_i) = 1$ .

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Then, we will solve the optimization problem:  $\max_{q=\Pi_i q_i} H(q) + \mathbb{E}_{x\sim q}[E(x)]$ 

It's clear that the number of parameters is small at least.

If we can take gradient wrt variables  $q_i(x_i)$  we can at least do *gradient* descent. Objective in general is *non-convex* though, so technically, this can fail !! (But often works ok.)

**Gibbs variational principle**: Let  $p(x) = \frac{1}{Z} \exp(E(x))$  be a distribution over a domain  $\mathcal{X}$ . Then, Z is the solution to the following optimization problem:  $\log Z = \max_{q: \text{ distribution over } \mathcal{X}} H(q) + \mathbb{E}_{x \sim q}[E(x)]$ 

**Typical example:** mean-field approximation:  $\max_{q=\Pi_i q_i} H(q) + \mathbb{E}_{x\sim q}[E(x)]$ 

Can we take gradients?

(1) Entropy factorizes: 
$$H(q_1q_2) = \sum_{x_1,x_2 \in \{\pm 1\}} q_1(x_1)q_2(x_2) \log(q_1(x_1)q_2(x_2))$$

$$= \sum_{x_1,x_2} q_1(x_1) q_2(x_2) \left(\log q_1(x_1) + \log q_2(x_2)\right) \\ = \sum_{x_2} q_2(x_2) H(q_1) + \sum_{x_1} q_1(x_1) H(q_2)$$

$$= H(q_1) + H(q_2) = q_1(x_1)\log q_1(x_1) + (1 - q_1(x_1))\log(1 - q_1(x_1)) + H(q_2)$$

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**Typical example:** mean-field approximation:  $\max_{q=\Pi_i q_i} H(q) + \mathbb{E}_{x\sim q}[E(x)]$  Can we solve this?

(2)  $\mathbb{E}_{x \sim q}[E(x)]$  is often not a problem:

e.g. for Ising models,  $E(x) = \sum_{ij} J_{ij} x_i x_j$  so  $\mathbb{E}_q[E(x)] = \sum_{ij} J_{ij} \mathbb{E}[x_i] \mathbb{E}[x_j]$  $\mathbb{E}[x_j] = q_i(x_i)$ , so taking a gradient is simple.

**Gibbs variational principle**: Let  $p(x) = \frac{1}{Z} \exp(E(x))$  be a distribution over a domain  $\mathcal{X}$ . Then, Z is the solution to the following optimization problem:  $\log Z = \max_{q: \text{ distribution over } \mathcal{X}} H(q) + \mathbb{E}_{x \sim q}[E(x)]$ 

**Typical example:** mean-field approximation:  $\max_{q=\Pi_i q_i} H(q) + \mathbb{E}_{x\sim q}[E(x)]$  Can we solve this?

(2)  $\mathbb{E}_{x \sim q}[E(x)]$  is often not a problem:

Even if gradients are not explicit, if E(x) is a sum of terms  $\phi_S(x_S)$ , we can estimate the expectations, and estimate the gradient from that. (i.e. use zeroth order optimization method.)

**Gibbs variational principle**: Let  $p(x) = \frac{1}{Z} \exp(E(x))$  be a distribution over a domain  $\mathcal{X}$ . Then, Z is the solution to the following optimization problem:  $\log Z = \max_{q: \text{ distribution over } \mathcal{X}} H(q) + \mathbb{E}_{x \sim q}[E(x)]$ 

**Typical example:** Gaussians with mean/covariance mx as parameters.

(1) Entropy of Gaussian has a closed-form formula:

$$\frac{1}{2}\log((2\pi e)^n\det\Sigma)$$

(2) Expectations wrt to Gaussian can be (often) estimated by drawing samples.

# Part II: Inference in latent-variable Bayesian networks

As we noted last time, the difficulty for calculating posteriors in latent-variable models is of a *similar nature*:

$$p(z|x) = \frac{p(z)p(x|z)}{p(x)}$$
 Integrates over all possible z's, hard normalizing factor

Taking inspiration from previous part, we will approximate posteriors p(z|x) in an analogous way:

# Variational methods for posterior distributions

Let p(z, x) be a joint distribution over latent variables and observables. Then, same as in the *Gibbs principle* calculation:

$$KL(q(z|x)||p(z|x)) = \mathbb{E}_{q(z|x)} \log q(z|x) - \mathbb{E}_{q(z|x)} \log p(z|x)$$
$$= -H(q(z|x)) - \mathbb{E}_{z \sim q}[p(z,x)] + \log p(x)$$

Hence,

$$\operatorname{argmax}_{q(z|x)} \mathit{KL}(q(z|x)) \Big| p(z|x) \Big) = \operatorname{argmax} \{ \mathbb{E}_{q(z|x)} \log q(z|x) - \mathbb{E}_{q(z|x)} \log p(z,x) \}$$

As in the undirected case, if q is a simple distribution (e.g. product distribution, Gaussian), we can optimize this using gradient descent.

## Yet another mean field strategy: coordinate ascent

Consider updating a \*single\* coordinate of the mean-field distribution, that is keep  $q_{-i}(z_i|x)$  fixed, and optimize for  $q_i(z_i|x)$ . Rewriting, we have:

$$= \mathbb{E}_{q(z|x)} \log q(z|x) - \mathbb{E}_{q(z|x)} \log p(z,x)$$

$$= \sum_{i} \mathbb{E}_{q_{i}(z_{i}|x)} \log q_{i}(z_{i}|x) - \mathbb{E}_{q_{i}(z_{i}|x)} \left[ \mathbb{E}_{q_{-i}(Z_{-i}|x)} \log p(z_{i}, z_{-i}, x) \right]$$

$$= \mathbb{E}_{q_i(z_i|x)} \log q_i(z_i|x) - \mathbb{E}_{q_i(z_i|x)} [\log \tilde{p}(z_i,x)] + C$$

Renormalize to make it a distribution

## Yet another mean field strategy: coordinate ascent

Consider updating a \*single\* coordinate of the mean-field distribution, that is keep  $q_{-i}$  ( $z_i|x$ ) fixed, and optimize for  $q_i$  ( $z_i|x$ ). Rewriting, we have:

$$KL(q(z|x)||p(z|x))$$

$$= \mathbb{E}_{q(z|x)} \log q(z|x) - \mathbb{E}_{q(z|x)} \log p(z,x)$$

$$= \sum_{i} \mathbb{E}_{q_{i}(z_{i}|x)} \log q_{i}(z_{i}|x) - \mathbb{E}_{q_{i}(z_{i}|x)} \left[ \mathbb{E}_{q_{-i}(z_{-i}|x)} \log p(z_{i}, z_{-i}, x) \right]$$

$$= \mathbb{E}_{q_{i}(z_{i}|x)} \log q_{i}(z_{i}|x) - \mathbb{E}_{q_{i}(z_{i}|x)} [\log \tilde{p}(z_{i}, x)] + C$$

$$= KL(q_{i}(z_{i}|x)||\tilde{p}(z_{i}, x)) + C$$

$$= \frac{\mathbb{E}_{q_{-i}(z_{-i}|x)} \log p(z_{i}, z_{-i}, x)}{\int_{z_{i}} \mathbb{E}_{q_{-i}(z_{-i}|x)} \log p(z_{i}, z_{-i}, x)}$$

### Yet another mean field strategy: coordinate ascent

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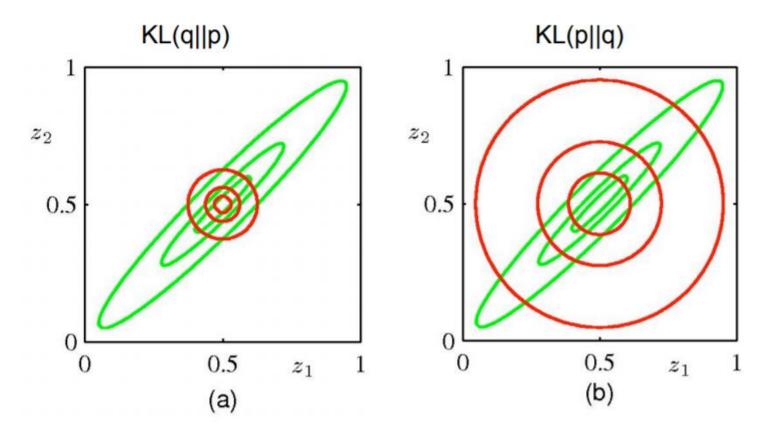
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### What if we changed the order of p, q?



Approximation is too compact.

Approximation is too spread.

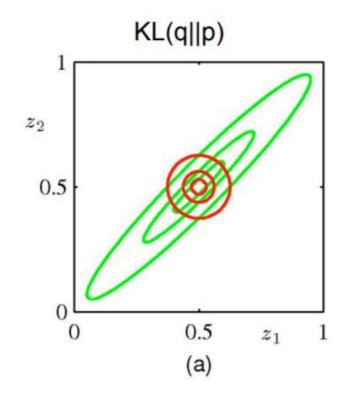
### What if we changed the order of p, q?

$$\mathrm{KL}(q||p) = -\int q(\mathbf{Z}) \ln \frac{p(\mathbf{Z})}{q(\mathbf{Z})} \mathrm{d}\mathbf{Z}.$$

There is a large positive contribution to the KL divergence from regions of Z space in which:

- p(Z) is near zero
- unless q(Z) is also close to zero.

Minimizing KL(q||p) leads to distributions q(Z) that avoid regions in which p(Z) is small.



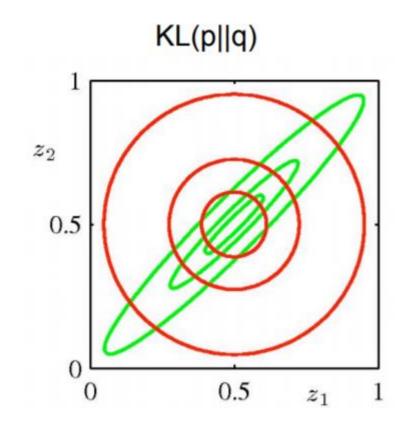
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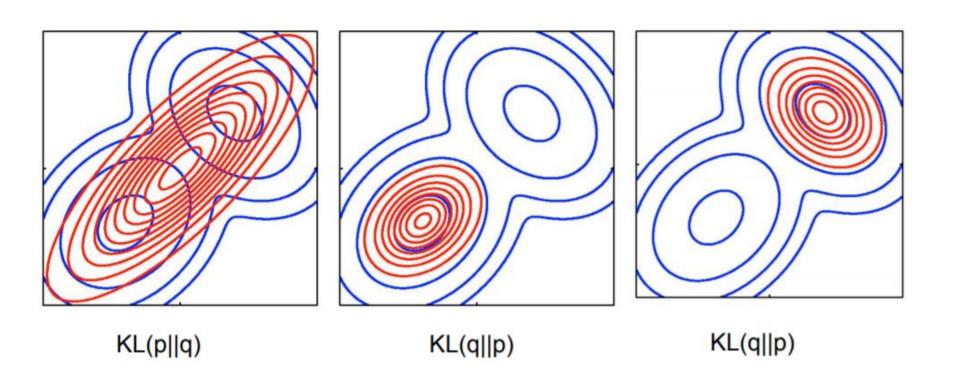
There is a large positive contribution to the KL divergence from regions of Z space in which:

- q(Z) is near zero,
- unless p(Z) is also close to zero.

Minimizing KL(p||q) leads to distributions q(Z) that are nonzero in regions where p(Z) is nonzero.



#### Multimodal distributions



Blue contours show bimodal distribution, red contours single Gaussian distribution that best approximates it.

KL(q||p) will tend to find a single mode, whereas KL(p||q) will average across all of the modes.

### Learning

# Learning latent-variable directed graphical models

How should we try to learn the parameters of a graphical model?

The most obvious strategy: maximum likelihood estimation

Given data  $x_1, x_2, ..., x_n$ , solve the optimization problem

$$\max_{\theta \in \Theta} \sum_{i=1}^{n} \log p(x_i)$$

Latent variables: we will use the variational principle again!

$$\log_{\theta} p(x) = \max_{q(z|x): \text{ distribution over } \mathcal{Z}} H(q(z|x)) + \mathbb{E}_{q(z|x)}[\log p_{\theta}(x,z)]$$

Hence, MLE objective can be written as double maximization:

$$\max_{\theta \in \Theta} \max_{\{q_i(z|x_i)\}} \sum_{i=1}^n H(q_i(z|x_i)) + \mathbb{E}_{q_i(z|x_i)}[\log p_{\theta}(x_i, z)]$$

# Variational methods for posterior distributions

**ELBO (Evidence Lower Bound)**: Let p(z, x) be a joint distribution over latent variables and observables. Then:

$$\log p(x) = \max_{q(z|x): \text{ distribution over } Z} H(q|z) + \mathbb{E}_{q(z|x)}[\log p(x,z)]$$

Write, by Bayes rule,  $p(z|x) = \frac{p(x,z)}{p(x)}$ . Then, the formula above follows by Gibbs variational principle with  $E(x) = \log p(x,z)$ . **Argmax** = p(z|x)!

**Gibbs variational principle**: Let  $p(x) = \frac{1}{Z} \exp(E(x))$  be a distribution over a domain  $\mathcal{X}$ . Then, Z is the solution to the following optimization problem:  $\log Z = \max_{q: \text{ distribution over } \mathcal{X}} H(q) + \mathbb{E}_{x \sim q}[E(x)]$ 

## Expectation-maximization/variational inference

The canonical algorithm for learning a single-layer latent-variable Bayesian network is an iterative algorithm as follows.

Consider the max-likelihood objective, rewritten as in the previous slide:

$$\max_{\theta \in \Theta} \max_{\{q_i(z|x_i)\}} \sum_{i=1}^n H(q_i(z|x_i)) + \mathbb{E}_{q_i(z|x_i)}[\log p_{\theta}(x_i, z)]$$

Algorithm maintains iterates  $\theta^t$ ,  $\{q_i^t(z|x_i)\}$ , and updates them iteratively

#### (1) Expectation (E)-step:

Keep  $\theta^t$  fixed, set  $\{q_i^{t+1}(z|x_i)\}$ , s.t. they maximize the objective above.

#### (2) Maximization (M)-step:

Keep  $\{q_i^t(z|x_i)\}$  fixed, set  $\theta^{t+1}$  s.t. it maximizes the objective above.

Clearly, every step cannot make the objective worse!

Does \*not\* mean it converges to global optimum – could, e.g. get stuck in a local minimum.

### Expectation-maximization/variational inference

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Algorithm maintains iterates  $\theta^t$ ,  $q_i^t(z|x_i)$ , and updates them iteratively

#### (1) Expectation step:

Keep  $\theta^t$  and set  $q_i^{t+1}(z|x_i)$ , s.t. they maximize the objective above.

If the class is infinitely rich, the optimum is  $q_i^{t+1}(z|x_i) = p_{\theta^t}(z|x_i)$ 

This is called **expectation-maximization (EM)**. If class is not infinitely rich, it's called **variational inference**.

#### 1: Mixtures of spherical Gaussians

Consider a mixture of K Gaussians with unknown means  $p = \sum_{i=1}^{K} \frac{1}{K} \mathcal{N}(\mu_i, I_d)$ 

Let's try to calculate the E and M steps.

**E-step**: the optimal  $q_i^{t+1}(z|x_i)$  is  $p_{\theta^t}(z|x_i)$ . Can we calculate this?

By Bayes rule, 
$$p_{\theta^t}(z = k|x_i) \propto p(x_i|z = k) \propto e^{-\left||x_i - \mu_k^t|\right|^2}$$

Writing out the normalizing constant, we have

$$p_{\theta^t}(z = k|x_i) = \frac{e^{-||x_i - \mu_k^t||^2}}{\sum_{k'} e^{-||x_i - \mu_{k'}^t||^2}}$$

"Soft" version of assigning point to nearest cluster

#### 1: Mixtures of spherical Gaussians

Consider a mixture of K Gaussians with unknown means  $p = \sum_{i=1}^{K} \frac{1}{K} \mathcal{N}(\mu_i, I_d)$ 

Let's try to calculate the E and M steps.

**M-step**: given a quess  $q_i^t(z|x_i)$  , we can rewrite the maximization for  $\theta$  as:

$$\max_{\theta \in \Theta} \sum_{i=1}^{n} H(q_i^t(z|x_i)) + \mathbb{E}_{q_i^t(z|x_i)}[\log p_{\theta}(x_i, z)]$$

$$= \mathbb{E}_{q_i^t(z|x_i)}[\log p_{\theta}(x|z)]$$

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**M-step**: given a quess  $q_i^t(z|x_i)$  , we can rewrite the maximization for  $\theta$  as:

$$\max_{\theta} \mathbb{E}_{q_i^t(z|x_i)}[\log p_{\theta}(x|z)] = \max_{\theta} -\sum_{i=1}^n \sum_{k=1}^K q_i^t(z=k|x_i)||x_i - \mu_k||^2$$

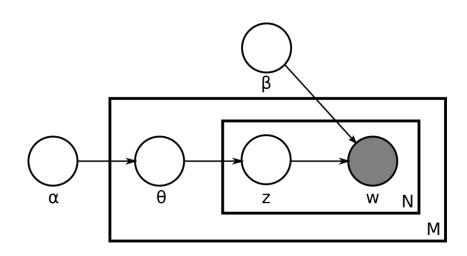
Setting the derivative wrt to  $\mu_k$  to 0, we have:

$$\mu_k^{t+1} = \sum_{i=1}^n \frac{e^{-\left||x_i - \mu_k^t|\right|^2}}{\sum_{k,l} e^{-\left||x_i - \mu_{k,l}^t|\right|^2}} x_i$$

#### 2: Latent Dirichlet Allocation

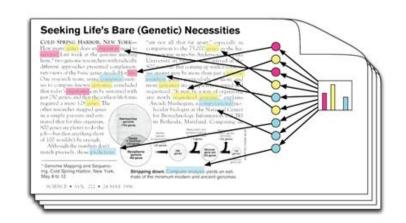
The **parameters** are:  $\{\alpha_i\}_{i=1}^K$  (Dirichlet parameters) and matrix  $\beta \in \mathbb{R}_+^{N \times K}$ , where N is the size of the vocabulary.

The columns of  $\beta$  satisfy  $\sum_{j=1}^{N} \beta_{ij} = 1$  (the distribution of words in a topic i)



#### To produce document:

- First, sample  $\theta \sim \text{Dir}(\cdot | \alpha)$ : this will be the topic proportion vector for the document.
- Each word in the document is generated in order, independently.
- ❖ To generate word i:
  - **Sample topic**  $z_i$  with categorical distribution with parameters  $\theta$
  - Sample word  $w_i$  with categorical distribution with parameters  $\beta_{z_i}$



The E-step cannot be done in closed form:

$$\begin{split} p(\vec{\theta}_{1:D}, z_{1:D,1:N}, \vec{\beta}_{1:K} \mid w_{1:D,1:N}, \alpha, \eta) &= \\ \frac{p(\vec{\theta}_{1:D}, \vec{z}_{1:D}, \vec{\beta}_{1:K} \mid \vec{w}_{1:D}, \alpha, \eta)}{\int_{\vec{\beta}_{1:K}} \int_{\vec{\theta}_{1:D}} \sum_{\vec{z}} p(\vec{\theta}_{1:D}, \vec{z}_{1:D}, \vec{\beta}_{1:K} \mid \vec{w}_{1:D}, \alpha, \eta)} \end{split}$$

(In fact, can be shown to be #P-hard to perform in the worst case.)

The variational family to approximate the posterior is commonly chosen to be a mean-field family:

$$q(\vec{\theta}_{1:D}, z_{1:D,1:N}, \vec{\beta}_{1:K}) = \prod_{k=1}^{K} q(\vec{\beta}_k \mid \vec{\lambda}_k) \prod_{d=1}^{D} \left( q(\vec{\theta}_{dd} \mid \vec{\gamma}_d) \prod_{n=1}^{N} q(z_{d,n} \mid \vec{\phi}_{d,n}) \right)$$

- Probability of topic z given document d:  $q(\theta_d \mid \gamma_d)$  Each document has its own Dirichlet prior  $\gamma_d$
- -Probability of word w given topic z:  $q(\beta_z \mid \lambda_z)$  Each topic has its own Dirichlet prior  $\lambda_z$
- -Probability of topic assignment to word  $w_{d,n}$ :  $q(z_{d,n} \mid \varphi_{d,n})$  Each word position word[d][n] has its own prior  $\varphi_{d,n}$

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#### One iteration of mean field variational inference for LDA

(1) For each topic k and term v:

(8) 
$$\lambda_{k,v}^{(t+1)} = \eta + \sum_{d=1}^{D} \sum_{n=1}^{N} 1(w_{d,n} = v) \phi_{n,k}^{(t)}.$$

- (2) For each document d:
  - (a) Update  $\gamma_d$ :

(9) 
$$\gamma_{d,k}^{(t+1)} = \alpha_k + \sum_{n=1}^N \phi_{d,n,k}^{(t)}.$$

(b) For each word n, update  $\vec{\phi}_{d,n}$ :

(10) 
$$\phi_{d,n,k}^{(t+1)} \propto \exp\left\{\Psi(\gamma_{d,k}^{(t+1)}) + \Psi(\lambda_{k,w_n}^{(t+1)}) - \Psi(\sum_{v=1}^{V} \lambda_{k,v}^{(t+1)})\right\},\,$$

where  $\Psi$  is the digamma function, the first derivative of the  $\log \Gamma$  function.

Parameter updates:

$$\beta_{ij} \propto \sum_{d=1}^{M} \sum_{n=1}^{N_d} \phi_{dni} w_{dn}^j.$$