

# Advanced Machine Learning

## Expectation Maximization

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Lecture 5, August 11, 2009

Reading:



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## Clustering



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# Unobserved Variables



- A variable can be unobserved (latent) because:
  - it is an imaginary quantity meant to provide some simplified and abstractive view of the data generation process
    - e.g., speech recognition models, mixture models ...
  - it is a real-world object and/or phenomena, but difficult or impossible to measure
    - e.g., the temperature of a star, causes of a disease, evolutionary ancestors ...
  - it is a real-world object and/or phenomena, but sometimes wasn't measured, because of faulty sensors; or was measure with a noisy channel, etc.
    - e.g., traffic radio, aircraft signal on a radar screen,
- Discrete latent variables can be used to partition/cluster data into sub-groups (mixture models, forthcoming).
- Continuous latent variables (factors) can be used for dimensionality reduction (factor analysis, etc., later lectures).

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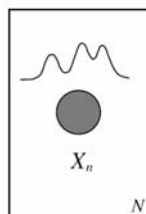
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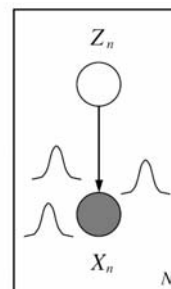
# Mixture Models



- A density model  $p(x)$  may be multi-modal.
- We may be able to model it as a mixture of uni-modal distributions (e.g., Gaussians).
- Each mode may correspond to a different sub-population (e.g., male and female).



(a)



(b)

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# Gaussian Mixture Models (GMMs)

- Consider a mixture of  $K$  Gaussian components:

- $Z$  is a latent class indicator vector:

$$p(z_n) = \text{multi}(z_n : \pi) = \prod_k (\pi_k)^{z_n^k}$$

- $X$  is a conditional Gaussian variable with a class-specific mean/covariance

$$p(x_n | z_n^k = 1, \mu, \Sigma) = \frac{1}{(2\pi)^{m/2} |\Sigma_k|^{1/2}} \exp\left\{-\frac{1}{2} (x_n - \mu_k)^T \Sigma_k^{-1} (x_n - \mu_k)\right\}$$

- The likelihood of a sample:

$$\begin{aligned} p(x_n | \mu, \Sigma) &= \sum_k p(z^k = 1 | \pi) p(x, | z^k = 1, \mu, \Sigma) \\ &= \sum_k \prod_k (\pi_k)^{z_n^k} \mathcal{N}(x_n : \mu_k, \Sigma_k) = \sum_k \pi_k \mathcal{N}(x, | \mu_k, \Sigma_k) \end{aligned}$$

mixture proportion      mixture component

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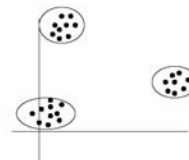
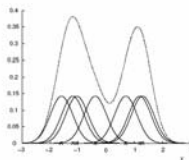
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# Gaussian Mixture Models (GMMs)

- Consider a mixture of  $K$  Gaussian components:

$$p(x_n | \mu, \Sigma) = \sum_k \pi_k \mathcal{N}(x, | \mu_k, \Sigma_k)$$



- This model can be used for unsupervised clustering.

- This model (fit by AutoClass) has been used to discover new kinds of stars in astronomical data, etc.

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## Learning mixture models

- Given data

$$\mathcal{D} = \{\mathbf{x}_n\}_{n=1}^N$$

- Likelihood:

$$L(\pi, \mu, \Sigma; D) = \prod_n p(x_n | \pi, \mu, \Sigma) = \prod_n \left( \sum_k \pi_k N(x_n | \mu_k, \Sigma_k) \right)$$

$$\{\pi^*, \mu^*, \Sigma^*\} = \arg \max L(\pi, \mu, \Sigma, D)$$

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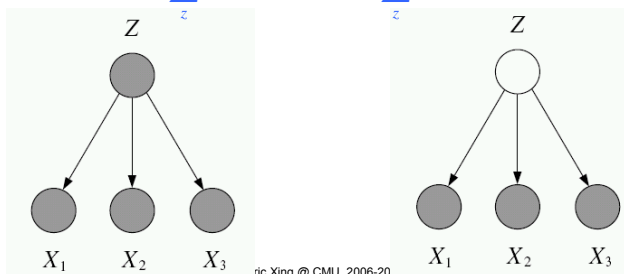
## Why is Learning Harder?

- In fully observed iid settings, the log likelihood decomposes into a sum of local terms.

$$\ell_c(\theta; D) = \log p(x, z | \theta) = \log p(z | \theta_z) + \log p(x | z, \theta_x)$$

- With latent variables, all the parameters become coupled together via *marginalization*

$$\ell_c(\theta; D) = \log \sum_z p(x, z | \theta) = \log \sum_z p(z | \theta_z) p(x | z, \theta_x)$$



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# Gradient Learning for mixture models



- We can learn mixture densities using gradient descent on the log likelihood. The gradients are quite interesting:

$$\begin{aligned}\mathcal{L}(\theta) &= \log p(\mathbf{x} | \theta) = \log \sum_k \pi_k p_k(\mathbf{x} | \theta_k) \\ \frac{\partial \mathcal{L}}{\partial \theta} &= \frac{1}{p(\mathbf{x} | \theta)} \sum_k \pi_k \frac{\partial p_k(\mathbf{x} | \theta_k)}{\partial \theta} \\ &= \sum_k \frac{\pi_k}{p(\mathbf{x} | \theta)} p_k(\mathbf{x} | \theta_k) \frac{\partial \log p_k(\mathbf{x} | \theta_k)}{\partial \theta} \\ &= \sum_k \pi_k \frac{p_k(\mathbf{x} | \theta_k)}{p(\mathbf{x} | \theta)} \frac{\partial \log p_k(\mathbf{x} | \theta_k)}{\partial \theta_k} = \sum_k r_k \frac{\partial \mathcal{L}_k}{\partial \theta_k}\end{aligned}$$

- In other words, the gradient is the responsibility weighted sum of the individual log likelihood gradients.
- Can pass this to a conjugate gradient routine.

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# Parameter Constraints



- Often we have constraints on the parameters, e.g.  $\sum_k \pi_k = 1$ ,  $\Sigma$  being symmetric positive definite (hence  $\Sigma_{ii} > 0$ ).
- We can use constrained optimization, or we can reparameterize in terms of unconstrained values.
  - For normalized weights, use the softmax transform:
  - For covariance matrices, use the Cholesky decomposition:

$$\Sigma^{-1} = \mathbf{A}^T \mathbf{A}$$

where  $\mathbf{A}$  is upper diagonal with positive diagonal:

$$\mathbf{A}_{ii} = \exp(\lambda_i) > 0 \quad \mathbf{A}_{ij} = \eta_{ij} \quad (j > i) \quad \mathbf{A}_{ij} = 0 \quad (j < i)$$

the parameters  $\gamma, \lambda, \eta_{ij} \in \mathbb{R}$  are unconstrained.

- Use chain rule to compute  $\frac{\partial \mathcal{L}}{\partial \pi}, \frac{\partial \mathcal{L}}{\partial \mathbf{A}}$ .

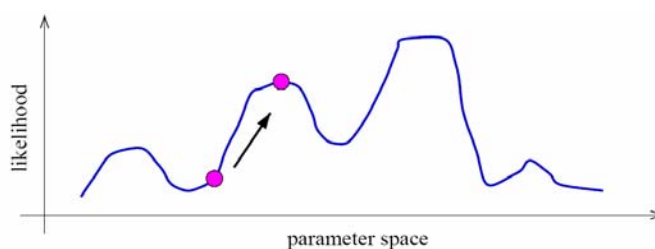
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# Identifiability

- A mixture model induces a multi-modal likelihood.
- Hence gradient ascent can only find a local maximum.
- Mixture models are unidentifiable, since we can always switch the hidden labels without affecting the likelihood.
- Hence we should be careful in trying to interpret the “meaning” of latent variables.



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# Toward the EM algorithm

- E.g., A mixture of K Gaussians:

- $Z$  is a latent class indicator vector

$$p(z_n) = \text{multi}(z_n : \pi) = \prod_k (\pi_k)^{z_n^k}$$

- $X$  is a conditional Gaussian variable with a class-specific mean/covariance

$$p(x_n | z_n^k = 1, \mu, \Sigma) = \frac{1}{(2\pi)^{m/2} |\Sigma_k|^{1/2}} \exp\left\{-\frac{1}{2}(x_n - \mu_k)^T \Sigma_k^{-1} (x_n - \mu_k)\right\}$$

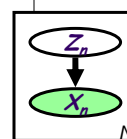
- The likelihood of a sample:

$$\begin{aligned} p(x_n | \mu, \Sigma) &= \sum_k p(z_n^k = 1 | \pi) p(x_n | z_n^k = 1, \mu, \Sigma) \\ &= \sum_{z_n} \prod_k \left( (\pi_k)^{z_n^k} N(x_n : \mu_k, \Sigma_k)^{z_n^k} \right) = \sum_k \pi_k N(x_n | \mu_k, \Sigma_k) \end{aligned}$$

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## Toward the EM algorithm



- Recall MLE for **completely observed data**

- Data log-likelihood

$$\begin{aligned}\mathcal{L}(\theta; D) &= \log \prod_n p(z_n, x_n) = \log \prod_n p(z_n | \pi) p(x_n | z_n, \mu, \sigma) \\ &= \sum_n \log \prod_k \pi_k^{z_n^k} + \sum_n \log \prod_k N(x_n; \mu_k, \sigma)^{z_n^k} \\ &= \sum_n \sum_k z_n^k \log \pi_k - \sum_n \sum_k \frac{1}{2\sigma^2} (x_n - \mu_k)^2 + C\end{aligned}$$

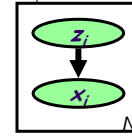
- MLE  $\hat{\pi}_{k,MLE} = \arg \max_{\pi} \mathcal{L}(\theta; D),$

$$\hat{\mu}_{k,MLE} = \arg \max_{\mu} \mathcal{L}(\theta; D)$$

$$\hat{\sigma}_{k,MLE} = \arg \max_{\sigma} \mathcal{L}(\theta; D)$$

$$\Rightarrow \hat{\mu}_{k,MLE} = \frac{\sum_n z_n^k x_n}{\sum_n z_n^k}$$

- What if we do not know  $z_n$ ?



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## Expectation-Maximization (EM) Algorithm



- EM is an optimization strategy for objective functions that can be interpreted as likelihoods in the presence of missing data.
- It is much simpler than gradient methods:
  - No need to choose step size.
  - Enforces constraints automatically.
  - Calls inference and fully observed learning as subroutines.
- EM is an Iterative algorithm with two linked steps:
  - E-step: fill-in hidden values using inference,  $p(z|x, \theta)$ .
  - M-step: update parameters  $t+1$  using standard MLE/MAP method applied to completed data
- We will prove that this procedure monotonically improves (or leaves it unchanged). Thus it always converges to a local optimum of the likelihood.

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# K-means

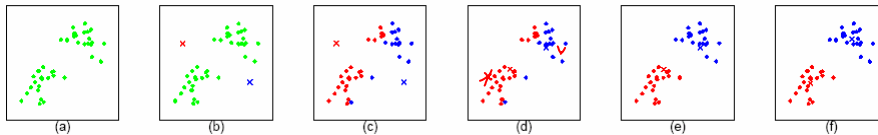
- Start:
  - "Guess" the centroid  $\mu_k$  and covariance  $\Sigma_k$  of each of the K clusters
- Loop
  - For each point  $n=1$  to  $N$ ,  
compute its cluster label:

$$z_n^{(t)} = \arg \max_k (x_n - \mu_k^{(t)})^T \Sigma_k^{-1(t)} (x_n - \mu_k^{(t)})$$

- For each cluster  $k=1:K$

$$\mu_k^{(t+1)} = \frac{\sum_n \delta(z_n^{(t)}, k) x_n}{\sum_n \delta(z_n^{(t)}, k)}$$

$$\Sigma_k^{(t+1)} = \dots$$



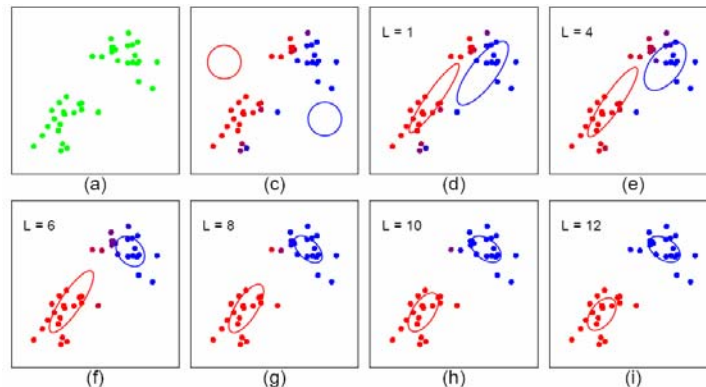
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# Expectation-Maximization

- Start:
  - "Guess" the centroid  $\mu_k$  and covariance  $\Sigma_k$  of each of the K clusters
- Loop



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## How is EM derived?

- A mixture of K Gaussians:

- $Z$  is a latent class indicator vector

$$p(\mathbf{z}_n) = \text{multi}(\mathbf{z}_n : \pi) = \prod_k (\pi_k)^{z_n^k}$$

- $X$  is a conditional Gaussian variable with a class-specific mean/covariance

$$p(x_n | z_n^k = 1, \mu, \Sigma) = \frac{1}{(2\pi)^{m/2} |\Sigma_k|^{1/2}} \exp\left\{-\frac{1}{2} (x_n - \mu_k)^T \Sigma_k^{-1} (x_n - \mu_k)\right\}$$

- The likelihood of a sample:

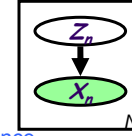
$$\begin{aligned} p(x_n | \mu, \Sigma) &= \sum_k p(z_n^k = 1 | \pi) p(x_n | z_n^k = 1, \mu, \Sigma) \\ &= \sum_{z_n} \prod_k (\pi_k)^{z_n^k} N(x_n; \mu_k, \Sigma_k)^{z_n^k} = \sum_k \pi_k N(x_n | \mu_k, \Sigma_k) \end{aligned}$$

- The “complete” likelihood

$$p(x_n, z_n^k = 1 | \mu, \Sigma) = p(z_n^k = 1 | \pi) p(x_n | z_n^k = 1, \mu, \Sigma) = \pi_k N(x_n | \mu_k, \Sigma_k)$$

$$p(x_n, z_n | \mu, \Sigma) = \prod_k [\pi_k N(x_n | \mu_k, \Sigma_k)]^{z_n^k}$$

But this is itself a random variable! Not good as objective function



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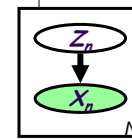
## How is EM derived?

- The complete log likelihood:

$$\begin{aligned} \mathcal{L}(\theta; D) &= \log \prod_n p(z_n, x_n) = \log \prod_n p(z_n | \pi) p(x_n | z_n, \mu, \sigma) \\ &= \sum_n \log \prod_k \pi_k^{z_n^k} + \sum_n \log \prod_k N(x_n; \mu_k, \sigma)^{z_n^k} \\ &= \sum_n \sum_k z_n^k \log \pi_k - \sum_n \sum_k z_n^k \frac{1}{2\sigma^2} (x_n - \mu_k)^2 + C \end{aligned}$$

- The expected complete log likelihood

$$\begin{aligned} \langle \mathcal{L}_c(\theta; \mathbf{x}, \mathbf{z}) \rangle &= \sum_n \langle \log p(z_n | \pi) \rangle_{p(\mathbf{z}|\mathbf{x})} + \sum_n \langle \log p(x_n | z_n, \mu, \Sigma) \rangle_{p(\mathbf{z}|\mathbf{x})} \\ &= \sum_n \sum_k \langle z_n^k \rangle \log \pi_k - \frac{1}{2} \sum_n \sum_k \langle z_n^k \rangle (x_n - \mu_k)^T \Sigma_k^{-1} (x_n - \mu_k) + \log |\Sigma_k| + C \end{aligned}$$



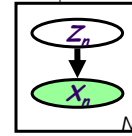
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## E-step

- We maximize  $\langle \ell_c(\theta) \rangle$  iteratively using the following iterative procedure:



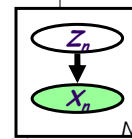
- Expectation step:** computing the expected value of the sufficient statistics of the hidden variables (i.e.,  $z$ ) given current est. of the parameters (i.e.,  $\pi$  and  $\mu$ ).

$$\tau_n^{k(t)} = \langle z_n^k \rangle_{q^{(t)}} = p(z_n^k = 1 | x_n, \mu^{(t)}, \Sigma^{(t)}) = \frac{\pi_k^{(t)} \mathcal{N}(x_n | \mu_k^{(t)}, \Sigma_k^{(t)})}{\sum_i \pi_i^{(t)} \mathcal{N}(x_n | \mu_i^{(t)}, \Sigma_i^{(t)})}$$

- Here we are essentially doing **inference**

## M-step

- We maximize  $\langle \ell_c(\theta) \rangle$  iteratively using the following iterative procedure:



- Maximization step:** compute the parameters under current results of the expected value of the hidden variables

$$\begin{aligned} \pi_k^* &= \arg \max \langle \ell_c(\theta) \rangle, \quad \Rightarrow \frac{\partial}{\partial \pi_k} \langle \ell_c(\theta) \rangle = 0, \forall k, \quad \text{s.t. } \sum_k \pi_k = 1 \\ &\Rightarrow \pi_k^* = \frac{\sum_n \tau_n^{k(t)} \langle z_n^k \rangle_{q^{(t)}}}{\sum_n \tau_n^{k(t)}} = \frac{\sum_n \tau_n^{k(t)}}{\sum_n 1} = \frac{\langle n_k \rangle}{N} \end{aligned}$$

$$\mu_k^* = \arg \max \langle \ell_c(\theta) \rangle, \quad \Rightarrow \mu_k^{(t+1)} = \frac{\sum_n \tau_n^{k(t)} x_n}{\sum_n \tau_n^{k(t)}}$$

$$\Sigma_k^* = \arg \max \langle \ell_c(\theta) \rangle, \quad \Rightarrow \Sigma_k^{(t+1)} = \frac{\sum_n \tau_n^{k(t)} (x_n - \mu_k^{(t+1)})(x_n - \mu_k^{(t+1)})^T}{\sum_n \tau_n^{k(t)}}$$

Fact:

$$\frac{\partial \log |A^{-1}|}{\partial A^{-1}} = A^T$$

$$\frac{\partial x^T A x}{\partial A} = x x^T$$

- This is isomorphic to **MLE** except that the variables that are hidden are replaced by their expectations (in general they will be replaced by their corresponding "**sufficient statistics**").

## Compare: K-means

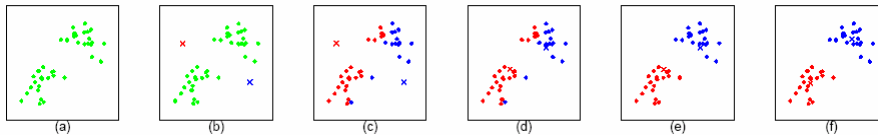


- The EM algorithm for mixtures of Gaussians is like a "soft version" of the K-means algorithm.
- In the K-means "E-step" we do hard assignment:

$$z_n^{(t)} = \arg \max_k (x_n - \mu_k^{(t)})^T \Sigma_k^{-1(t)} (x_n - \mu_k^{(t)})$$

- In the K-means "M-step" we update the means as the weighted sum of the data, but now the weights are 0 or 1:

$$\mu_k^{(t+1)} = \frac{\sum_n \delta(z_n^{(t)}, k) x_n}{\sum_n \delta(z_n^{(t)}, k)}$$



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## Theory underlying EM



- What are we doing?
- Recall that according to MLE, we intend to learn the model parameter that would have maximize the likelihood of the data.
- But we do not observe  $z$ , so computing

$$\ell_c(\theta; D) = \log \sum_z p(x, z | \theta) = \log \sum_z p(z | \theta_z) p(x | z, \theta_x)$$

is difficult!

- What shall we do?

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# Complete & Incomplete Log Likelihoods



- Complete log likelihood

Let  $X$  denote the observable variable(s), and  $Z$  denote the latent variable(s).

If  $Z$  could be observed, then

$$\ell_c(\theta; x, z) \stackrel{\text{def}}{=} \log p(x, z | \theta)$$

- Usually, optimizing  $\ell_c()$  given both  $z$  and  $x$  is straightforward (c.f. MLE for fully observed models).
- Recalled that in this case the objective for, e.g., MLE, decomposes into a sum of factors, the parameter for each factor can be estimated separately.
- **But given that  $Z$  is not observed,  $\ell_c()$  is a random quantity, cannot be maximized directly.**

- Incomplete log likelihood

With  $z$  unobserved, our objective becomes the log of a marginal probability:

$$\ell_c(\theta; x) = \log p(x | \theta) = \log \sum_z p(x, z | \theta)$$

- **This objective won't decouple**

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# Expected Complete Log Likelihood



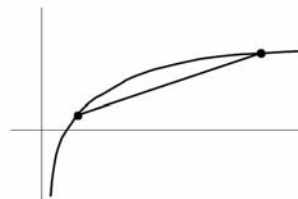
- For **any** distribution  $q(z)$ , define **expected complete log likelihood**:

$$\langle \ell_c(\theta; x, z) \rangle_q \stackrel{\text{def}}{=} \sum_z q(z | x, \theta) \log p(x, z | \theta)$$

- A deterministic function of  $\theta$
- Linear in  $\ell_c()$  --- inherit its factorizability
- Does maximizing this surrogate yield a maximizer of the likelihood?

- Jensen's inequality

$$\begin{aligned} \ell(\theta; x) &= \log p(x | \theta) \\ &= \log \sum_z p(x, z | \theta) \\ &= \log \sum_z q(z | x) \frac{p(x, z | \theta)}{q(z | x)} \\ &\geq \sum_z q(z | x) \log \frac{p(x, z | \theta)}{q(z | x)} \end{aligned} \Rightarrow \ell(\theta; x) \geq \langle \ell_c(\theta; x, z) \rangle_q + H_q$$



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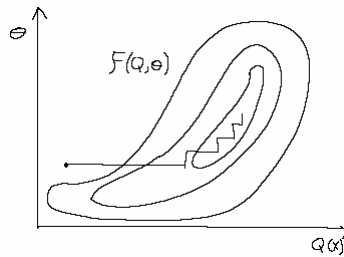
## Lower Bounds and Free Energy

- For fixed data  $x$ , define a functional called the free energy:

$$F(q, \theta) \stackrel{\text{def}}{=} \sum_z q(z | x) \log \frac{p(x, z | \theta)}{q(z | x)} \leq \ell(\theta; x)$$

- The EM algorithm is coordinate-ascent on  $F$ :

- E-step:**  $q^{t+1} = \arg \max_q F(q, \theta^t)$
- M-step:**  $\theta^{t+1} = \arg \max_{\theta} F(q^{t+1}, \theta)$



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## E-step: maximization of expected $\ell_c$ w.r.t. $q$

- Claim:  $q^{t+1} = \arg \max_q F(q, \theta^t) = p(z | x, \theta^t)$ 
  - This is the posterior distribution over the latent variables given the data and the parameters. Often we need this at test time anyway (e.g. to perform classification).

- Proof (easy): this setting attains the bound  $\ell(\theta; x) \geq F(q, \theta)$

$$\begin{aligned} F(p(z|x, \theta^t), \theta^t) &= \sum_z p(z|x, \theta^t) \log \frac{p(x, z | \theta^t)}{p(z|x, \theta^t)} \\ &= \sum_z p(z|x, \theta^t) \log p(x | \theta^t) \\ &= \log p(x | \theta^t) = \ell(\theta^t; x) \end{aligned}$$

- Can also show this result using variational calculus or the fact that  $\ell(\theta; x) - F(q, \theta) = \text{KL}(q \| p(z | x, \theta))$

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## E-step $\equiv$ plug in posterior expectation of latent variables



- Without loss of generality: assume that  $p(\mathbf{x}, \mathbf{z} | \theta)$  is a generalized exponential family distribution:

$$p(\mathbf{x}, \mathbf{z} | \theta) = \frac{1}{Z(\theta)} h(\mathbf{x}, \mathbf{z}) \exp \left\{ \sum_i \theta_i f_i(\mathbf{x}, \mathbf{z}) \right\}$$

- Special cases: if  $p(\mathbf{x} | \mathbf{z})$  are GLIMs, then  $f_i(\mathbf{x}, \mathbf{z}) = \eta_i^T(\mathbf{z}) \xi_i(\mathbf{x})$

- The expected complete log likelihood under  $q^{t+1} = p(\mathbf{z} | \mathbf{x}, \theta^t)$  is

$$\begin{aligned} \langle \ell_c(\theta^t; \mathbf{x}, \mathbf{z}) \rangle_{q^{t+1}} &= \sum_{\mathbf{z}} q(\mathbf{z} | \mathbf{x}, \theta^t) \log p(\mathbf{x}, \mathbf{z} | \theta^t) - A(\theta) \\ &= \sum_i \theta_i^t \langle f_i(\mathbf{x}, \mathbf{z}) \rangle_{q(\mathbf{z} | \mathbf{x}, \theta^t)} - A(\theta) \\ &\stackrel{p\text{-GLIM}}{=} \sum_i \theta_i^t \langle \eta_i(\mathbf{z}) \rangle_{q(\mathbf{z} | \mathbf{x}, \theta^t)} \xi_i(\mathbf{x}) - A(\theta) \end{aligned}$$

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## M-step: maximization of expected $\ell_c$ w.r.t. $\theta$



- Note that the free energy breaks into two terms:

$$\begin{aligned} F(q, \theta) &= \sum_{\mathbf{z}} q(\mathbf{z} | \mathbf{x}) \log \frac{p(\mathbf{x}, \mathbf{z} | \theta)}{q(\mathbf{z} | \mathbf{x})} \\ &= \sum_{\mathbf{z}} q(\mathbf{z} | \mathbf{x}) \log p(\mathbf{x}, \mathbf{z} | \theta) - \sum_{\mathbf{z}} q(\mathbf{z} | \mathbf{x}) \log q(\mathbf{z} | \mathbf{x}) \\ &= \langle \ell_c(\theta; \mathbf{x}, \mathbf{z}) \rangle_q + H_q \end{aligned}$$

- The first term is the expected complete log likelihood (energy) and the second term, which does not depend on  $\theta$ , is the entropy.
- Thus, in the M-step, maximizing with respect to  $\theta$  for fixed  $q$  we only need to consider the first term:

$$\theta^{t+1} = \arg \max_{\theta} \langle \ell_c(\theta; \mathbf{x}, \mathbf{z}) \rangle_{q^{t+1}} = \arg \max_{\theta} \sum_{\mathbf{z}} q(\mathbf{z} | \mathbf{x}) \log p(\mathbf{x}, \mathbf{z} | \theta)$$

- Under optimal  $q^{t+1}$ , this is equivalent to solving a standard MLE of fully observed model  $p(\mathbf{x}, \mathbf{z} | \theta)$ , with the **sufficient statistics** involving  $\mathbf{z}$  replaced by their expectations w.r.t.  $p(\mathbf{z} | \mathbf{x}, \theta)$ .

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## Summary: EM Algorithm



- A way of maximizing likelihood function for latent variable models. Finds MLE of parameters when the original (hard) problem can be broken up into two (easy) pieces:
  1. Estimate some “missing” or “unobserved” data from observed data and current parameters.
  2. Using this “complete” data, find the maximum likelihood parameter estimates.
- Alternate between filling in the latent variables using the best guess (posterior) and updating the parameters based on this guess:
  - E-step:  $q^{t+1} = \arg \max_q F(q, \theta^t)$
  - M-step:  $\theta^{t+1} = \arg \max_{\theta} F(q^{t+1}, \theta)$
- In the M-step we optimize a lower bound on the likelihood. In the E-step we close the gap, making bound=likelihood.

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## A Report Card for EM



- Some good things about EM:
  - no learning rate (step-size) parameter
  - automatically enforces parameter constraints
  - very fast for low dimensions
  - each iteration guaranteed to improve likelihood
- Some bad things about EM:
  - can get stuck in local minima
  - can be slower than conjugate gradient (especially near convergence)
  - requires expensive inference step
  - is a maximum likelihood/MAP method

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