Advanced Machine Learning Summer 2019

Part 15 – Latent Variable Models II 06.06.2019

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Course Outline

- Regression Techniques
 - Linear Regression
 - Regularization (Ridge, Lasso)
 - Kernels (Kernel Ridge Regression)
- Deep Reinforcement Learning
- Probabilistic Graphical Models
 - Bayesian Networks
 - Markov Random Fields
 - Inference (exact & approximate)
 - Latent Variable Models
- Deep Generative Models
 - Generative Adversarial Networks
 - Variational Autoencoders

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Topics of This Lecture

- Recap: Mixtures of Gaussians and General EM
 - Mixtures of Gaussians
 - General EM
- Mixtures of Gaussians revisited
 - General EM derivation
- The EM algorithm in general
 - Generalized EM
 - Relation to Variational inference





Recap: Mixtures of Gaussians



Recap: GMMs as Latent Variable Models

- Write GMMs in terms of latent variables \mathbf{z}
 - Marginal distribution of ${\bf x}$

$$p(\mathbf{x}) = \sum_{\mathbf{z}} p(\mathbf{x}, \mathbf{z}) = \sum_{\mathbf{z}} p(\mathbf{z}) p(\mathbf{x} | \mathbf{z}) = \sum_{k=1}^{K} \pi_k \mathcal{N}(\mathbf{x} | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)$$

- Advantage of this formulation
 - We have represented the marginal distribution in terms of latent variables z.
 - Since $p(\mathbf{x}) = \sum_{\mathbf{z}} p(\mathbf{x}, \mathbf{z})$, there is a corresponding latent variable \mathbf{z}_n for each data point \mathbf{x}_n .
 - We are now able to work with the joint distribution $p(\mathbf{x}, \mathbf{z})$ instead of the marginal distribution $p(\mathbf{x})$.
 - \Rightarrow This will lead to significant simplifications...





Recap: Sampling from a Gaussian Mixture

- MoG Sampling
 - We can use ancestral sampling to generate random samples from a Gaussian mixture model.
 - 1. Generate a value $\hat{\mathbf{z}}$ from the marginal distribution $p(\mathbf{z})$.
 - 2. Generate a value $\hat{\mathbf{x}}$ from the conditional distribution $p(\mathbf{x}|\hat{\mathbf{z}})$.



Image source: C.M. Bishop, 2006

Recap: Gaussian Mixtures Revisited

- Applying the latent variable view of EM
 - Goal is to maximize the log-likelihood using the observed data ${\bf X}$

$$\log p(\mathbf{X}|\boldsymbol{\theta}) = \log \left\{ \sum_{\mathbf{Z}} p(\mathbf{X}, \mathbf{Z}|\boldsymbol{\theta}) \right\} \xrightarrow{\pi} \left\{ \sum_{\mathbf{X}_n} p(\mathbf{X}, \mathbf{Z}|\boldsymbol{\theta}) \right\}$$

– Corresponding graphical model:

- Suppose we are additionally given the values of the latent variables ${f Z}$.
- The corresponding graphical model for the complete data now looks like this:
- \Rightarrow Straightforward to marginalize...

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Image source: C.M. Bishop, 2006

Recap: Alternative View of EM

- In practice, however,...
 - We are not given the complete data set { \mathbf{X}, \mathbf{Z} }, but only the incomplete data \mathbf{X} . All we can compute about \mathbf{Z} is the posterior distribution $p(\mathbf{Z}|\mathbf{X}, \boldsymbol{\theta})$.
 - Since we cannot use the complete-data log-likelihood, we consider instead its expected value under the posterior distribution of the latent variables:

$$Q(\theta, \theta^{\text{old}}) = \sum_{\mathbf{Z}} p(\mathbf{Z}|\mathbf{X}, \theta^{\text{old}}) \log p(\mathbf{X}, \mathbf{Z}|\theta)$$

- This corresponds to the E-step of the EM algorithm.
- In the subsequent M-step, we then maximize the expectation to obtain the revised parameter set θ^{new} .

$$oldsymbol{ heta}^{ ext{new}} = rg\max_{oldsymbol{ heta}} ~ \mathcal{Q}(oldsymbol{ heta},oldsymbol{ heta}^{ ext{old}})$$





Recap: General EM Algorithm

- Algorithm
 - 1. Choose an initial setting for the parameters $\theta^{\rm old}$
 - 2. E-step: Evaluate $p(\mathbf{Z}|\mathbf{X}, \boldsymbol{\theta}^{\text{old}})$
 - 3. M-step: Evaluate $\boldsymbol{\theta}^{\mathrm{new}}$ given by

$$oldsymbol{ heta}^{ ext{new}} = rg\max_{oldsymbol{ heta}} \, \mathcal{Q}(oldsymbol{ heta},oldsymbol{ heta}^{ ext{old}})$$

where

$$\mathcal{Q}(\boldsymbol{\theta}, \boldsymbol{\theta}^{\text{old}}) = \sum_{\mathbf{Z}} p(\mathbf{Z} | \mathbf{X}, \boldsymbol{\theta}^{\text{old}}) \log p(\mathbf{X}, \mathbf{Z} | \boldsymbol{\theta})$$

4. While not converged, let $\boldsymbol{\theta}^{\mathrm{old}} \leftarrow \boldsymbol{\theta}^{\mathrm{new}}$ and return to step 2.





Recap: MAP-EM

- Modification for MAP
 - The EM algorithm can be adapted to find MAP solutions for models for which a prior $p(\theta)$ is defined over the parameters.
 - Only changes needed:
 - 2. E-step: Evaluate $p(\mathbf{Z}|\mathbf{X}, \boldsymbol{\theta}^{\text{old}})$
 - 3. M-step: Evaluate $\boldsymbol{\theta}^{\mathrm{new}}$ given by

$$oldsymbol{ heta}^{ ext{new}} = rg\max_{oldsymbol{ heta}} \, \mathcal{Q}(oldsymbol{ heta},oldsymbol{ heta}^{ ext{old}}) + \log p(oldsymbol{ heta})$$

 \Rightarrow Suitable choices for the prior will remove the ML singularities!





Recap: Monte Carlo EM

- EM procedure
 - M-step: Maximize expectation of complete-data log-likelihood

$$Q(\theta, \theta^{\text{old}}) = \int p(\mathbf{Z} | \mathbf{X}, \theta^{\text{old}}) \log p(\mathbf{X}, \mathbf{Z} | \theta) d\mathbf{Z}$$

- For more complex models, we may not be able to compute this analytically anymore...
- Idea

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– Use sampling to approximate this integral by a finite sum over samples $\{\mathbf{Z}^{(l)}\}\$ drawn from the current estimate of the posterior

$$\mathcal{Q}(\boldsymbol{\theta}, \boldsymbol{\theta}^{\text{old}}) \sim \frac{1}{L} \sum_{l=1}^{L} \log p(\mathbf{X}, \mathbf{Z}^{(l)} | \boldsymbol{\theta})$$

- This procedure is called the Monte Carlo EM algorithm.





- Applying the latent variable view of EM
 - Goal is to maximize the log-likelihood using the observed data ${f X}$

$$\log p(\mathbf{X}|\boldsymbol{\theta}) = \log \left\{ \sum_{\mathbf{Z}} p(\mathbf{X}, \mathbf{Z}|\boldsymbol{\theta}) \right\} \xrightarrow{\pi} \left\{ \begin{array}{c} \mathbf{x}_{n} \\ \mathbf{x}_{n} \end{array} \right\}$$

– Corresponding graphical model:

- Suppose we are additionally given the values of the latent variables \mathbf{Z} .
- The corresponding graphical model for the complete data now looks like this:



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- Maximize the likelihood
 - For the complete-data set $\{\mathbf{X},\!\mathbf{Z}\},$ the likelihood has the form

$$p(\mathbf{X}, \mathbf{Z} | \boldsymbol{\mu}, \boldsymbol{\Sigma}, \boldsymbol{\pi}) = \prod_{n=1}^{N} \prod_{k=1}^{K} \pi_k^{z_{nk}} \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)^{z_{nk}}$$

- Taking the logarithm, we obtain

$$\log p(\mathbf{X}, \mathbf{Z} | \boldsymbol{\mu}, \boldsymbol{\Sigma}, \boldsymbol{\pi}) = \sum_{n=1}^{N} \sum_{k=1}^{K} z_{nk} \left\{ \log \pi_k + \log \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k) \right\}$$

- Compared to the incomplete-data case, the order of the sum and logarithm has been interchanged.
- \Rightarrow Much simpler solution to the ML problem.
- Maximization w.r.t. a mean or covariance is exactly as for a single Gaussian, except that it involves only the subset of data points that are "assigned" to that component ($z_{nk} = 1$).





- Maximization w.r.t. mixing coefficients
 - More complex, since the π_k are coupled by the summation constraint

$$\sum_{j=1}^{n} \pi_j = 1$$

- Solve with a Lagrange multiplier

$$\log p(\mathbf{X}, \mathbf{Z} | \boldsymbol{\mu}, \boldsymbol{\Sigma}, \boldsymbol{\pi}) + \lambda \left(\sum_{k=1}^{K} \pi_k - 1 \right)$$

/ TZ

- Solution (after a longer derivation):

$$\pi_k = \frac{1}{N} \sum_{n=1}^N z_{nk}$$

 \Rightarrow The complete-data log-likelihood can be maximized trivially in closed form.

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- In practice, we don't have values for the latent variables
 - Consider the expectation w.r.t. the posterior distribution of the latent variables instead.
 - The posterior distribution takes the form

$$p(\mathbf{Z}|\mathbf{X}, \boldsymbol{\mu}, \boldsymbol{\Sigma}, \boldsymbol{\pi}) \propto \prod_{n=1}^{N} \prod_{k=1}^{K} \left[\pi_k \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k) \right]^{z_{nk}}$$

and factorizes over n, so that the $\{\mathbf{z}_n\}$ are independent under the posterior.

– Expected value of indicator variable z_{nk} under the posterior.

$$\mathbb{E}[z_{nk}] = \frac{\sum_{z_{nk}} z_{nk} \left[\pi_k \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)\right]^{z_{nk}}}{\sum_{z_{nj}} \left[\pi_j \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_j, \boldsymbol{\Sigma}_j)\right]^{z_{nj}}}$$
$$= \frac{\pi_k \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)}{\sum_{j=1}^{K} \pi_j \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_j, \boldsymbol{\Sigma}_j)} = \gamma(z_{nk})$$



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- Continuing the estimation
 - The expected value of the complete-data log-likelihood is therefore

$$\mathbb{E}_{\mathbf{Z}}[\log p(\mathbf{X}, \mathbf{Z} | \boldsymbol{\mu}, \boldsymbol{\Sigma}, \boldsymbol{\pi})] = \sum_{n=1}^{N} \sum_{k=1}^{K} \gamma z_{nk} \{\log \pi_k + \log \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)\}$$

- Putting everything together
 - Start by choosing some initial values for μ^{old} , Σ^{old} , and π^{old} .
 - Use these to evaluate the responsibilities (the E-Step).
 - Keep the responsibilities fixed and maximize the above for μ^{new} , Σ^{new} , and π^{new} (the M-Step).
 - This leads to the familiar closed-form solutions for μ^{new} , Σ^{new} , and π^{new} .
 - ⇒ This is precisely the EM algorithm for Gaussian mixtures as derived before. But we can now also apply it to other distributions.

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The EM Algorithm in General

- General formulation
 - Given a probabilistic model with observed variables X, hidden variables Z and parameters θ .
 - Our goal is to maximize the likelihood given by

$$p(\mathbf{X}|\theta) = \sum_{\mathbf{Z}} p(\mathbf{X}, \mathbf{Z}|\theta)$$

- However, a direct optimization of $p(\mathbf{X}|\mathbf{\theta})$ is often difficult. Optimization of the complete-data log-likelihood $p(\mathbf{X}, \mathbf{Z}|\mathbf{\theta})$ is significantly easier.





The EM Algorithm in General

- Decomposition
 - Introduce a distribution $q(\mathbf{Z})$ over the latent variables. For any choice of $q(\mathbf{Z})$, the following decomposition holds

$$\log p(\mathbf{X}|\theta) = \mathcal{L}(q, \mathbf{\theta}) + KL(q \parallel p)$$

- where

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$$\mathcal{L}(q, \mathbf{\theta}) = \sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{X}, \mathbf{Z} | \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$

$$KL(q \parallel p) = -\sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{Z} \mid \mathbf{X}, \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$

- (Proof on extra slide set)

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Analysis of this Result

- Decomposition
 - For any choice of $q(\mathbf{Z})$, the following decomposition holds

$$\log p(\mathbf{X}|\theta) = \mathcal{L}(q, \theta) + KL(q \parallel p)$$

$$\mathcal{L}(q, \mathbf{\theta}) = \sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{X}, \mathbf{Z} | \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$
$$KL(q \parallel p) = -\sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{Z} | \mathbf{X}, \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$

• Notes (1)

- $-\mathcal{L}(q, \theta)$ is a functional of the distribution $q(\mathbf{Z})$ and a function of the parameters θ .
- A functional is an operator that takes as input a function and outputs again a function.



Analysis of this Result

- Decomposition
 - For any choice of $q(\mathbf{Z})$, the following decomposition holds

$$\log p(\mathbf{X}|\theta) = \mathcal{L}(q, \theta) + KL(q \parallel p)$$

$$\mathcal{L}(q, \mathbf{\theta}) = \sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{X}, \mathbf{Z} | \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$
$$KL(q \parallel p) = -\sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{Z} | \mathbf{X}, \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$

• Notes (2)

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- $KL(q \parallel p)$ is the Kullback-Leibler divergence between the distribution $q(\mathbf{Z})$ and the posterior distribution $p(\mathbf{Z}|\mathbf{X}, \mathbf{\theta})$.
- The KL divergence satisfies $KL(q \parallel p) \ge 0$ with = 0 iff $q(\mathbf{Z}) = p(\mathbf{Z} \mid \mathbf{X}, \mathbf{\theta})$.



Analysis of this Result

- Decomposition
 - For any choice of $q(\mathbf{Z})$, the following decomposition holds

$$\log p(\mathbf{X}|\theta) = \mathcal{L}(q, \theta) + KL(q \parallel p)$$

$$\mathcal{L}(q, \mathbf{\theta}) = \sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{X}, \mathbf{Z} | \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$
$$KL(q \parallel p) = -\sum_{\mathbf{Z}} q(\mathbf{Z}) \log \left\{ \frac{p(\mathbf{Z} | \mathbf{X}, \mathbf{\theta})}{q(\mathbf{Z})} \right\}$$

• Notes (3)

- It therefore follows that $\mathcal{L}(q, \theta) \leq \log p(\mathbf{X}|\theta)$.
- In other words: $\mathcal{L}(q, \theta)$ is a lower bound on $\log p(\mathbf{X}|\theta)$.
- We can now use this result in order to analyze how EM works...





Decomposition

$$\log p(\mathbf{X}|\theta) = \mathcal{L}(q, \boldsymbol{\theta}) + KL(q \parallel p)$$



Interpretation

- $-\mathcal{L}(q, \theta)$ is a lower bound on $\log p(\mathbf{X}|\theta)$.
- The approximation comes from the fact that we use an approximative distribution $q(\mathbf{Z}) = p(\mathbf{Z}|\mathbf{X}, \mathbf{\theta}^{old})$ Instead of the (unknown) real posterior.
- The KL divergence measures the difference between the approximative distribution $q(\mathbf{Z})$ and the real posterior $p(\mathbf{Z}|\mathbf{X}, \mathbf{\theta})$.
- In every EM iteration, we try to make this difference smaller.





Decomposition

$$\log p(\mathbf{X}|\theta) = \mathcal{L}(q, \theta) + KL(q \parallel p)$$



• E-Step

- Suppose the current value of the parameter vector is θ^{old} .
- The E-step maximizes the lower bound $\mathcal{L}(q, \theta)$ w.r.t. $q(\mathbf{Z})$ while holding θ^{old} fixed.
- The solution to this maximization problem of $\log p(\mathbf{X}|\mathbf{\theta}^{old})$ will occur when the KL divergence vanishes, i.e. when $q(\mathbf{Z}) = p(\mathbf{Z}|\mathbf{X}, \mathbf{\theta}^{old})$.
- In this case, the lower bound equals the log-likelihood.





Decomposition

 $\log p(\mathbf{X}|\theta) = \mathcal{L}(q, \mathbf{\theta}) + KL(q \parallel p)$



• M-Step

- In the M-step, the distribution $q(\mathbf{Z})$ is held fixed and the lower bound $\mathcal{L}(q, \mathbf{\theta})$ is maximized w.r.t. $\mathbf{\theta}$ to give some new value $\mathbf{\theta}^{new}$.
- This causes the lower bound \mathcal{L} to increase (unless it is already at maximum), which will cause the log-likelihood to increase.
- Because $q(\mathbf{Z})$ is determined using the old parameter values, it will not equal the posterior distribution $p(\mathbf{Z}|\mathbf{X}, \mathbf{\theta}^{new})$ and there will be a non-zero KL divergence.





 Visualization in the space of parameters



- The EM algorithm alternately
 - Computes a lower bound on the log-likelihood for the current parameters values
 - And then maximizes this bound to obtain the new parameter values.





References and Further Reading

• More information about EM and MoG estimation is available in Chapter 9 of Bishop's book (recommendable to read).

Christopher M. Bishop Pattern Recognition and Machine Learning Springer, 2006



- Additional information
 - A.P. Dempster, N.M. Laird, D.B. Rubin, <u>Maximum-Likelihood from incomplete</u> data via EM algorithm", In J. Royal Statistical Society, Series B. Vol 39, 1977
 - J.A. Bilmes, "<u>A Gentle Tutorial of the EM Algorithm and its Application to</u> <u>Parameter Estimation for Gaussian Mixture and Hidden Markov Models</u>", TR-97-021, ICSI, U.C. Berkeley, CA,USA





