

Probabilistic & Unsupervised Learning

The EM algorithm

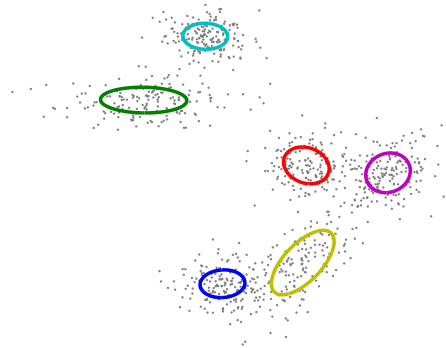
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Term 1, Autumn 2013

Example: mixture of Gaussians



Data: $\mathcal{X} = \{\mathbf{x}_1 \dots \mathbf{x}_N\}$

Latent process:
 $s_i \stackrel{\text{iid}}{\sim} \text{Disc}[\boldsymbol{\pi}]$

Component distributions:
 $\mathbf{x}_i | (s_i = m) \sim \mathcal{P}_m[\theta_m] = \mathcal{N}(\boldsymbol{\mu}_m, \boldsymbol{\Sigma}_m)$

Marginal distribution:
$$P(\mathbf{x}_i) = \sum_{m=1}^k \pi_m \mathcal{P}_m(\mathbf{x}; \theta_m)$$

Log-likelihood:

$$\ell(\{\boldsymbol{\mu}_m\}, \{\boldsymbol{\Sigma}_m\}, \boldsymbol{\pi}) = \sum_{i=1}^n \log \sum_{m=1}^k \frac{\pi_m}{\sqrt{|2\pi\boldsymbol{\Sigma}_m|}} e^{-\frac{1}{2}(\mathbf{x}_i - \boldsymbol{\mu}_m)^\top \boldsymbol{\Sigma}_m^{-1}(\mathbf{x}_i - \boldsymbol{\mu}_m)}$$

Log-likelihoods

▶ Exponential family models: $p(\mathbf{x}|\boldsymbol{\theta}) = f(\mathbf{x})e^{\boldsymbol{\theta}^\top \mathbf{T}(\mathbf{x})} / Z(\boldsymbol{\theta})$

$$\ell(\boldsymbol{\theta}) = \boldsymbol{\theta}^\top \sum_n \mathbf{T}(\mathbf{x}_n) - N \log Z(\boldsymbol{\theta}) \quad (+ \text{constants})$$

- ▶ Concave function.
- ▶ Maximum may be closed-form.
- ▶ If not, numerical optimisation is still generally straightforward.

▶ Latent variable models: $p(\mathbf{x}|\boldsymbol{\theta}_x, \boldsymbol{\theta}_y) = \int d\mathbf{y} f_x(\mathbf{x}) \frac{e^{\boldsymbol{\phi}(\boldsymbol{\theta}_x, \mathbf{y})^\top \mathbf{T}_x(\mathbf{x})}}{Z_x(\boldsymbol{\phi}(\boldsymbol{\theta}_x, \mathbf{y}))} f_y(\mathbf{y}) \frac{e^{\boldsymbol{\theta}_y^\top \mathbf{T}_y(\mathbf{y})}}{Z_y(\boldsymbol{\theta}_y)}$

$$\ell(\boldsymbol{\theta}_x, \boldsymbol{\theta}_y) = \sum_n \log \int d\mathbf{y} f_x(\mathbf{x}) \frac{e^{\boldsymbol{\phi}(\boldsymbol{\theta}_x, \mathbf{y})^\top \mathbf{T}_x(\mathbf{x})}}{Z_x(\boldsymbol{\phi}(\boldsymbol{\theta}_x, \mathbf{y}))} f_y(\mathbf{y}) \frac{e^{\boldsymbol{\theta}_y^\top \mathbf{T}_y(\mathbf{y})}}{Z_y(\boldsymbol{\theta}_y)}$$

- ▶ Usually no closed form optimum.
- ▶ Often multiple local maxima.
- ▶ Direct numerical optimisation may be possible but infrequently easy.

The joint-data likelihood

▶ For many models, maximisation might be straightforward if \mathbf{y} were not latent, and we could just maximise the joint-data likelihood:

$$\ell(\boldsymbol{\theta}_x, \boldsymbol{\theta}_y) = \sum_n \boldsymbol{\phi}(\boldsymbol{\theta}_x, \mathbf{y}_n)^\top \mathbf{T}_x(\mathbf{x}_n) + \boldsymbol{\theta}_y^\top \sum_n \mathbf{T}_y(\mathbf{y}_n) - \sum_n \log Z_x(\boldsymbol{\phi}(\boldsymbol{\theta}_x, \mathbf{y}_n)) - N \log Z_y(\boldsymbol{\theta}_y)$$

- ▶ Conversely, if we knew $\boldsymbol{\theta}$, we could compute (the posterior over) the values of \mathbf{y} .
- ▶ **Idea:** update $\boldsymbol{\theta}$ and (the distribution on) \mathbf{y} in alternation, converging to a self-consistent answer.
- ▶ Will this yield the right answer?
- ▶ Typically, it will (as we shall see). This is the **Expectation Maximisation (EM)** algorithm.

The Expectation Maximisation (EM) algorithm

The EM algorithm (Dempster, Laird & Rubin, 1977; but significant earlier precedents) finds a (local) maximum of a latent variable model likelihood. It starts from arbitrary values of the parameters, and iterates two steps:

E step: Fill in values of latent variables according to posterior given data.

M step: Maximise likelihood as if latent variables were not hidden.

- ▶ Useful in models where learning would be easy if hidden variables were, in fact, observed (e.g. MoGs).
- ▶ Decomposes difficult problems into series of tractable steps.
- ▶ No learning rate.
- ▶ Framework lends itself to principled approximations.
- ▶ How does it work?

The lower bound for EM – “free energy”

Observed data $\mathcal{X} = \{\mathbf{x}_i\}$; Latent variables $\mathcal{Y} = \{\mathbf{y}_i\}$; Parameters $\theta = \{\theta_x, \theta_y\}$.

Log-likelihood:

$$\ell(\theta) = \log P(\mathcal{X}|\theta) = \log \int d\mathcal{Y} P(\mathcal{Y}, \mathcal{X}|\theta)$$

By Jensen, any distribution, $q(\mathcal{Y})$, over the latent variables generates a lower bound:

$$\ell(\theta) = \log \int d\mathcal{Y} q(\mathcal{Y}) \frac{P(\mathcal{Y}, \mathcal{X}|\theta)}{q(\mathcal{Y})} \geq \int d\mathcal{Y} q(\mathcal{Y}) \log \frac{P(\mathcal{Y}, \mathcal{X}|\theta)}{q(\mathcal{Y})} \stackrel{\text{def}}{=} \mathcal{F}(q, \theta).$$

Now,

$$\begin{aligned} \int d\mathcal{Y} q(\mathcal{Y}) \log \frac{P(\mathcal{Y}, \mathcal{X}|\theta)}{q(\mathcal{Y})} &= \int d\mathcal{Y} q(\mathcal{Y}) \log P(\mathcal{Y}, \mathcal{X}|\theta) - \int d\mathcal{Y} q(\mathcal{Y}) \log q(\mathcal{Y}) \\ &= \int d\mathcal{Y} q(\mathcal{Y}) \log P(\mathcal{Y}, \mathcal{X}|\theta) + \mathbf{H}[q], \end{aligned}$$

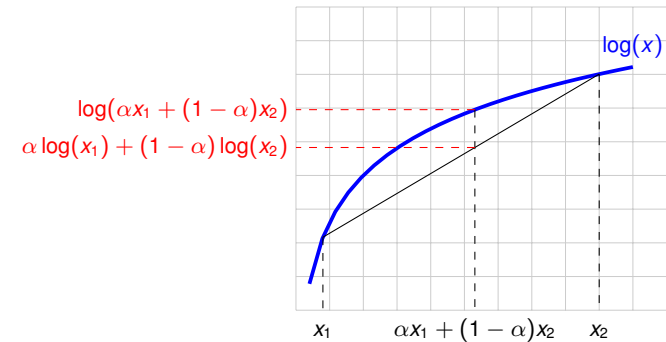
where $\mathbf{H}[q]$ is the entropy of $q(\mathcal{Y})$.

So:

$$\mathcal{F}(q, \theta) = \langle \log P(\mathcal{Y}, \mathcal{X}|\theta) \rangle_{q(\mathcal{Y})} + \mathbf{H}[q]$$

Jensen's inequality

One view: EM iteratively refines a lower bound on the log-likelihood.



In general:

For $\alpha_i \geq 0$, $\sum \alpha_i = 1$ (and $\{x_i > 0\}$):

For probability measure α and concave f

$$\log \left(\sum_i \alpha_i x_i \right) \geq \sum_i \alpha_i \log(x_i)$$

$$f(E_\alpha [x]) \geq E_\alpha [f(x)]$$

Equality (if and only if $f(x)$ is almost surely constant or linear on (convex) support of α).

The E and M steps of EM

The lower bound on the log likelihood is given by:

$$\mathcal{F}(q, \theta) = \langle \log P(\mathcal{Y}, \mathcal{X}|\theta) \rangle_{q(\mathcal{Y})} + \mathbf{H}[q],$$

EM alternates between:

- ▶ **E step:** optimize $\mathcal{F}(q, \theta)$ wrt distribution over hidden variables holding parameters fixed:

$$q^{(k)}(\mathcal{Y}) := \operatorname{argmax}_{q(\mathcal{Y})} \mathcal{F}(q(\mathcal{Y}), \theta^{(k-1)}).$$

- ▶ **M step:** maximize $\mathcal{F}(q, \theta)$ wrt parameters holding hidden distribution fixed:

$$\theta^{(k)} := \operatorname{argmax}_{\theta} \mathcal{F}(q^{(k)}(\mathcal{Y}), \theta) = \operatorname{argmax}_{\theta} \langle \log P(\mathcal{Y}, \mathcal{X}|\theta) \rangle_{q^{(k)}(\mathcal{Y})}$$

The second equality comes from the fact $\mathbf{H}[q^{(k)}(\mathcal{Y})]$ does not depend directly on θ .

The E Step

The free energy can be re-written

$$\begin{aligned}\mathcal{F}(q, \theta) &= \int q(\mathcal{Y}) \log \frac{P(\mathcal{Y}, \mathcal{X}|\theta)}{q(\mathcal{Y})} d\mathcal{Y} \\ &= \int q(\mathcal{Y}) \log \frac{P(\mathcal{Y}|\mathcal{X}, \theta)P(\mathcal{X}|\theta)}{q(\mathcal{Y})} d\mathcal{Y} \\ &= \int q(\mathcal{Y}) \log P(\mathcal{X}|\theta) d\mathcal{Y} + \int q(\mathcal{Y}) \log \frac{P(\mathcal{Y}|\mathcal{X}, \theta)}{q(\mathcal{Y})} d\mathcal{Y} \\ &= \ell(\theta) - \mathbf{KL}[q(\mathcal{Y})||P(\mathcal{Y}|\mathcal{X}, \theta)]\end{aligned}$$

The second term is the Kullback-Leibler divergence.

This means that, for fixed θ , \mathcal{F} is bounded above by ℓ , and achieves that bound when $\mathbf{KL}[q(\mathcal{Y})||P(\mathcal{Y}|\mathcal{X}, \theta)] = 0$.

But $\mathbf{KL}[q||p]$ is zero if and only if $q = p$ (see appendix.)

So, the E step sets

$$q^{(k)}(\mathcal{Y}) = P(\mathcal{Y}|\mathcal{X}, \theta^{(k-1)})$$

and, after an E step, the free energy equals the likelihood.

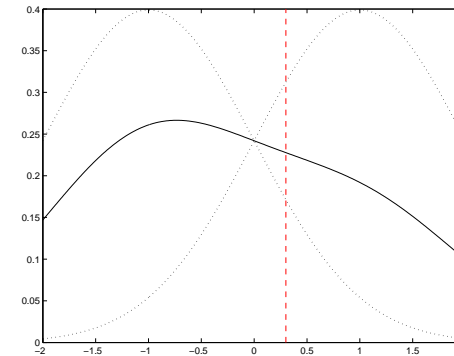
Coordinate Ascent in \mathcal{F} (Demo)

To visualise, we consider a one parameter / one latent mixture:

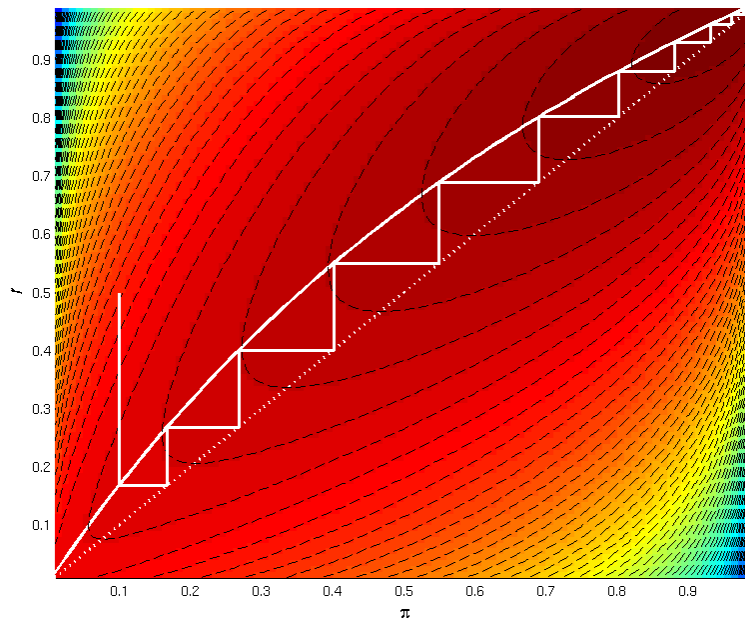
$$\begin{aligned}s &\sim \text{Bernoulli}[\pi] \\ x|s=0 &\sim \mathcal{N}[-1, 1] \quad x|s=1 \sim \mathcal{N}[1, 1].\end{aligned}$$

Single data point $x_1 = .3$.

$q(s)$ is a distribution on a single binary latent, and so is represented by $r_1 \in [0, 1]$.



Coordinate Ascent in \mathcal{F} (Demo)



EM Never Decreases the Likelihood

The E and M steps together never decrease the log likelihood:

$$\ell(\theta^{(k-1)}) \stackrel{\text{E step}}{=} \mathcal{F}(q^{(k)}, \theta^{(k-1)}) \leq \mathcal{F}(q^{(k)}, \theta^{(k)}) \stackrel{\text{Jensen}}{\leq} \ell(\theta^{(k)}),$$

- ▶ The E step brings the free energy to the likelihood.
- ▶ The M-step maximises the free energy wrt θ .
- ▶ $\mathcal{F} \leq \ell$ by Jensen – or, equivalently, from the non-negativity of KL

If the M-step is executed so that $\theta^{(k)} \neq \theta^{(k-1)}$ iff \mathcal{F} increases, then the overall EM iteration will step to a new value of θ iff the likelihood increases.

Can also show that fixed points of EM (generally) correspond to maxima of the likelihood (see appendices).

Partial M steps and Partial E steps

Partial M steps: The proof holds even if we just *increase* \mathcal{F} wrt θ rather than maximize. (Dempster, Laird and Rubin (1977) call this the generalized EM, or GEM, algorithm).

In fact, immediately after an E step

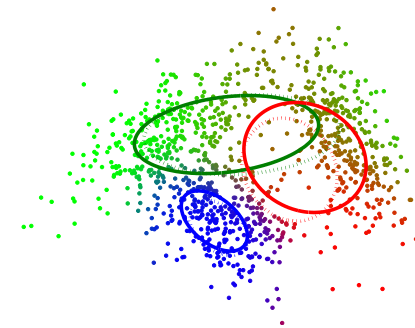
$$\left. \frac{\partial}{\partial \theta} \right|_{\theta^{(k-1)}} \langle \log P(\mathcal{X}, \mathcal{Y} | \theta) \rangle_{q^{(k)}(\mathcal{Y}) [= P(\mathcal{Y} | \mathcal{X}, \theta^{(k-1)})]} = \left. \frac{\partial}{\partial \theta} \right|_{\theta^{(k-1)}} \log P(\mathcal{X} | \theta)$$

So E-step (inference) can be used to construct other gradient-based optimisation schemes (e.g. "Expectation Conjugate Gradient", Salakhutdinov et al. *ICML* 2003).

Partial E steps: We can also just *increase* \mathcal{F} wrt to some of the q_s .

For example, sparse or online versions of the EM algorithm would compute the posterior for a subset of the data points or as the data arrives, respectively. One might also update the posterior over a subset of the hidden variables, while holding others fixed...

EM for MoGs



► Evaluate responsibilities

$$r_{im} = \frac{P_m(\mathbf{x})\pi_m}{\sum_{m'} P_{m'}(\mathbf{x})\pi_{m'}}$$

► Update parameters

$$\begin{aligned} \mu_m &\leftarrow \frac{\sum_i r_{im} \mathbf{x}_i}{\sum_i r_{im}} \\ \Sigma_m &\leftarrow \frac{\sum_i r_{im} (\mathbf{x}_i - \mu_m)(\mathbf{x}_i - \mu_m)^T}{\sum_i r_{im}} \\ \pi_m &\leftarrow \frac{\sum_i r_{im}}{N} \end{aligned}$$

The Gaussian mixture model (E-step)

In a univariate Gaussian mixture model, the density of a data point x is:

$$p(x|\theta) = \sum_{m=1}^k p(s=m|\theta)p(x|s=m, \theta) \propto \sum_{m=1}^k \frac{\pi_m}{\sigma_m} \exp\left\{-\frac{1}{2\sigma_m^2}(x - \mu_m)^2\right\},$$

where θ is the collection of parameters: means μ_m , variances σ_m^2 and mixing proportions $\pi_m = p(s=m|\theta)$.

The hidden variable s_i indicates which component generated observation x_i .

The E-step computes the posterior for s_i given the current parameters:

$$\begin{aligned} q(s_i) &= p(s_i|x_i, \theta) \propto p(x_i|s_i, \theta)p(s_i|\theta) \\ r_{im} &\stackrel{\text{def}}{=} q(s_i=m) \propto \frac{\pi_m}{\sigma_m} \exp\left\{-\frac{1}{2\sigma_m^2}(x_i - \mu_m)^2\right\} \quad (\text{responsibilities}) \quad \leftarrow \langle \delta_{s_i=m} \rangle_q \end{aligned}$$

with the normalization such that $\sum_m r_{im} = 1$.

The Gaussian mixture model (M-step)

In the M-step we optimize the sum (since s is discrete):

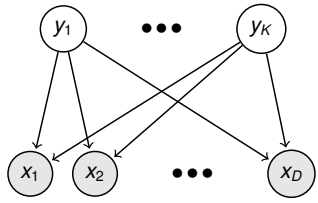
$$\begin{aligned} E &= \langle \log p(x, s|\theta) \rangle_{q(s)} = \sum q(s) \log[p(s|\theta) p(x|s, \theta)] \\ &= \sum_{i,m} r_{im} \left[\log \pi_m - \log \sigma_m - \frac{1}{2\sigma_m^2}(x_i - \mu_m)^2 \right]. \end{aligned}$$

Optimum is found by setting the partial derivatives of E to zero:

$$\begin{aligned} \frac{\partial}{\partial \mu_m} E &= \sum_i r_{im} \frac{(x_i - \mu_m)}{2\sigma_m^2} = 0 \Rightarrow \mu_m = \frac{\sum_i r_{im} x_i}{\sum_i r_{im}}, \\ \frac{\partial}{\partial \sigma_m} E &= \sum_i r_{im} \left[-\frac{1}{\sigma_m} + \frac{(x_i - \mu_m)^2}{\sigma_m^3} \right] = 0 \Rightarrow \sigma_m^2 = \frac{\sum_i r_{im} (x_i - \mu_m)^2}{\sum_i r_{im}}, \\ \frac{\partial}{\partial \pi_m} E &= \sum_i r_{im} \frac{1}{\pi_m}, \quad \frac{\partial E}{\partial \pi_m} + \lambda = 0 \Rightarrow \pi_m = \frac{1}{n} \sum_i r_{im}, \end{aligned}$$

where λ is a Lagrange multiplier ensuring that the mixing proportions sum to unity.

EM for Factor Analysis



The model for \mathbf{x} :

$$p(\mathbf{x}|\theta) = \int p(\mathbf{y}|\theta)p(\mathbf{x}|\mathbf{y}, \theta)d\mathbf{y} = \mathcal{N}(0, \Lambda\Lambda^T + \Psi)$$

Model parameters: $\theta = \{\Lambda, \Psi\}$.

E step: For each data point \mathbf{x}_n , compute the posterior distribution of hidden factors given the observed data: $q_n(\mathbf{y}_n) = p(\mathbf{y}_n|\mathbf{x}_n, \theta_t)$.

M step: Find the θ_{t+1} that maximises $\mathcal{F}(q, \theta)$:

$$\begin{aligned} \mathcal{F}(q, \theta) &= \sum_n \int q_n(\mathbf{y}_n) [\log p(\mathbf{y}_n|\theta) + \log p(\mathbf{x}_n|\mathbf{y}_n, \theta) - \log q_n(\mathbf{y}_n)] d\mathbf{y}_n \\ &= \sum_n \int q_n(\mathbf{y}_n) [\log p(\mathbf{y}_n|\theta) + \log p(\mathbf{x}_n|\mathbf{y}_n, \theta)] d\mathbf{y}_n + c. \end{aligned}$$

The M step for Factor Analysis

M step: Find θ_{t+1} by maximising $\mathcal{F} = \sum_n (\log p(\mathbf{y}_n|\theta) + \log p(\mathbf{x}_n|\mathbf{y}_n, \theta))_{q_n(\mathbf{y}_n)} + c$

$$\begin{aligned} &\log p(\mathbf{y}_n|\theta) + \log p(\mathbf{x}_n|\mathbf{y}_n, \theta) \\ &= c - \frac{1}{2}\mathbf{y}_n^T\mathbf{y}_n - \frac{1}{2}\log|\Psi| - \frac{1}{2}(\mathbf{x}_n - \Lambda\mathbf{y}_n)^T\Psi^{-1}(\mathbf{x}_n - \Lambda\mathbf{y}_n) \\ &= c' - \frac{1}{2}\log|\Psi| - \frac{1}{2}[\mathbf{x}_n^T\Psi^{-1}\mathbf{x}_n - 2\mathbf{x}_n^T\Psi^{-1}\Lambda\mathbf{y}_n + \mathbf{y}_n^T\Lambda^T\Psi^{-1}\Lambda\mathbf{y}_n] \\ &= c' - \frac{1}{2}\log|\Psi| - \frac{1}{2}[\mathbf{x}_n^T\Psi^{-1}\mathbf{x}_n - 2\mathbf{x}_n^T\Psi^{-1}\Lambda\mathbf{y}_n + \text{Tr}[\Lambda^T\Psi^{-1}\Lambda\mathbf{y}_n\mathbf{y}_n^T]] \end{aligned}$$

Taking expectations wrt $q_n(\mathbf{y}_n)$:

$$= c' - \frac{1}{2}\log|\Psi| - \frac{1}{2}[\mathbf{x}_n^T\Psi^{-1}\mathbf{x}_n - 2\mathbf{x}_n^T\Psi^{-1}\Lambda\boldsymbol{\mu}_n + \text{Tr}[\Lambda^T\Psi^{-1}\Lambda(\boldsymbol{\mu}_n\boldsymbol{\mu}_n^T + \Sigma)]]$$

Note that we don't need to know everything about $q(\mathbf{y}_n)$, just the moments $\langle \mathbf{y}_n \rangle$ and $\langle \mathbf{y}_n\mathbf{y}_n^T \rangle$. These are the **expected sufficient statistics**.

The E step for Factor Analysis

E step: For each data point \mathbf{x}_n , compute the posterior distribution of hidden factors given the observed data: $q_n(\mathbf{y}_n) = p(\mathbf{y}_n|\mathbf{x}_n, \theta) = p(\mathbf{y}_n, \mathbf{x}_n|\theta)/p(\mathbf{x}_n|\theta)$

Tactic: write $p(\mathbf{y}_n, \mathbf{x}_n|\theta)$, consider \mathbf{x}_n to be fixed. What is this as a function of \mathbf{y}_n ?

$$\begin{aligned} p(\mathbf{y}_n, \mathbf{x}_n) &= p(\mathbf{y}_n)p(\mathbf{x}_n|\mathbf{y}_n) \\ &= (2\pi)^{-\frac{K}{2}} \exp\{-\frac{1}{2}\mathbf{y}_n^T\mathbf{y}_n\} |2\pi\Psi|^{-\frac{1}{2}} \exp\{-\frac{1}{2}(\mathbf{x}_n - \Lambda\mathbf{y}_n)^T\Psi^{-1}(\mathbf{x}_n - \Lambda\mathbf{y}_n)\} \\ &= c \times \exp\{-\frac{1}{2}[\mathbf{y}_n^T\mathbf{y}_n + (\mathbf{x}_n - \Lambda\mathbf{y}_n)^T\Psi^{-1}(\mathbf{x}_n - \Lambda\mathbf{y}_n)]\} \\ &= c' \times \exp\{-\frac{1}{2}[\mathbf{y}_n^T(I + \Lambda^T\Psi^{-1}\Lambda)\mathbf{y}_n - 2\mathbf{y}_n^T\Lambda^T\Psi^{-1}\mathbf{x}_n]\} \\ &= c'' \times \exp\{-\frac{1}{2}[\mathbf{y}_n^T\Sigma^{-1}\mathbf{y}_n - 2\mathbf{y}_n^T\Sigma^{-1}\boldsymbol{\mu}_n + \boldsymbol{\mu}_n^T\Sigma^{-1}\boldsymbol{\mu}_n]\} \end{aligned}$$

So $\Sigma = (I + \Lambda^T\Psi^{-1}\Lambda)^{-1} = I - \beta\Lambda$ and $\boldsymbol{\mu}_n = \Sigma\Lambda^T\Psi^{-1}\mathbf{x}_n = \beta\mathbf{x}_n$. Where $\beta = \Sigma\Lambda^T\Psi^{-1}$. Note that $\boldsymbol{\mu}_n$ is a linear function of \mathbf{x}_n and Σ does not depend on \mathbf{x}_n .

The M step for Factor Analysis (cont.)

$$\mathcal{F} = c' - \frac{N}{2}\log|\Psi| - \frac{1}{2}\sum_n [\mathbf{x}_n^T\Psi^{-1}\mathbf{x}_n - 2\mathbf{x}_n^T\Psi^{-1}\Lambda\boldsymbol{\mu}_n + \text{Tr}[\Lambda^T\Psi^{-1}\Lambda(\boldsymbol{\mu}_n\boldsymbol{\mu}_n^T + \Sigma)]]$$

Taking derivatives wrt Λ and Ψ^{-1} , using $\frac{\partial \text{Tr}[AB]}{\partial B} = A^T$ and $\frac{\partial \log|A|}{\partial A} = A^{-T}$:

$$\frac{\partial \mathcal{F}}{\partial \Lambda} = \Psi^{-1}\sum_n \mathbf{x}_n\boldsymbol{\mu}_n^T - \Psi^{-1}\Lambda\left(N\Sigma + \sum_n \boldsymbol{\mu}_n\boldsymbol{\mu}_n^T\right) = 0$$

$$\Rightarrow \hat{\Lambda} = \left(\sum_n \mathbf{x}_n\boldsymbol{\mu}_n^T\right) \left(N\Sigma + \sum_n \boldsymbol{\mu}_n\boldsymbol{\mu}_n^T\right)^{-1}$$

$$\frac{\partial \mathcal{F}}{\partial \Psi^{-1}} = \frac{N}{2}\Psi - \frac{1}{2}\sum_n [\mathbf{x}_n\mathbf{x}_n^T - \Lambda\boldsymbol{\mu}_n\mathbf{x}_n^T - \mathbf{x}_n\boldsymbol{\mu}_n^T\Lambda^T + \Lambda(\boldsymbol{\mu}_n\boldsymbol{\mu}_n^T + \Sigma)\Lambda^T]$$

$$\Rightarrow \hat{\Psi} = \frac{1}{N}\sum_n [\mathbf{x}_n\mathbf{x}_n^T - \Lambda\boldsymbol{\mu}_n\mathbf{x}_n^T - \mathbf{x}_n\boldsymbol{\mu}_n^T\Lambda^T + \Lambda(\boldsymbol{\mu}_n\boldsymbol{\mu}_n^T + \Sigma)\Lambda^T]$$

$$\hat{\Psi} = \Lambda\Sigma\Lambda^T + \frac{1}{N}\sum_n (\mathbf{x}_n - \Lambda\boldsymbol{\mu}_n)(\mathbf{x}_n - \Lambda\boldsymbol{\mu}_n)^T \quad (\text{squared residuals})$$

Note: we should actually only take derivatives w.r.t. Ψ_{dd} since Ψ is diagonal. As $\Sigma \rightarrow 0$ these become the equations for ML linear regression

Mixtures of Factor Analysers

Simultaneous clustering and dimensionality reduction.

$$p(\mathbf{x}|\theta) = \sum_k \pi_k \mathcal{N}(\boldsymbol{\mu}_k, \Lambda_k \Lambda_k^\top + \Psi)$$

where π_k is the mixing proportion for FA k , $\boldsymbol{\mu}_k$ is its centre, Λ_k is its “factor loading matrix”, and Ψ is a common sensor noise model. $\theta = \{\{\pi_k, \boldsymbol{\mu}_k, \Lambda_k\}_{k=1\dots K}, \Psi\}$

We can think of this model as having *two* sets of hidden latent variables:

- ▶ A discrete indicator variable $s_n \in \{1, \dots, K\}$
- ▶ For each factor analyzer, a continous factor vector $\mathbf{y}_{n,k} \in \mathcal{R}^{D_k}$

$$p(\mathbf{x}|\theta) = \sum_{s_n=1}^K p(s_n|\theta) \int p(\mathbf{y}|s_n, \theta) p(\mathbf{x}_n|\mathbf{y}, s_n, \theta) d\mathbf{y}$$

As before, an EM algorithm can be derived for this model:

E step: We need moments of $p(\mathbf{y}_n, s_n|\mathbf{x}_n, \theta)$, specifically: $\langle \delta_{s_n=m} \rangle$, $\langle \delta_{s_n=m} \mathbf{y}_n \rangle$ and $\langle \delta_{s_n=m} \mathbf{y}_n \mathbf{y}_n^\top \rangle$.

M step: Similar to M-step for FA with responsibility-weighted moments.

See <http://www.learning.eng.cam.ac.uk/zoubin/papers/tr-96-1.pdf>

References

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EM for exponential families

EM is often applied to models whose **joint** over $\mathbf{z} = (\mathbf{y}, \mathbf{x})$ has exponential-family form:

$$p(\mathbf{z}|\theta) = f(\mathbf{z}) \exp\{\theta^\top \mathbf{T}(\mathbf{z})\} / Z(\theta)$$

(with $Z(\theta) = \int f(\mathbf{z}) \exp\{\theta^\top \mathbf{T}(\mathbf{z})\} d\mathbf{z}$) but whose marginal $p(\mathbf{x}) \notin \text{ExpFam}$.

The free energy dependence on θ is given by:

$$\begin{aligned} \mathcal{F}(q, \theta) &= \int q(\mathbf{y}) \log p(\mathbf{y}, \mathbf{x}|\theta) d\mathbf{y} - \mathbf{H}[q] \\ &= \int q(\mathbf{y}) [\theta^\top \mathbf{T}(\mathbf{z}) - \log Z(\theta)] d\mathbf{y} + \text{const wrt } \theta \\ &= \theta^\top \langle \mathbf{T}(\mathbf{z}) \rangle_{q(\mathbf{y})} - \log Z(\theta) + \text{const wrt } \theta \end{aligned}$$

So, in the **E step** all we need to compute are the **expected sufficient statistics** under q .

We also have:

$$\begin{aligned} \frac{\partial}{\partial \theta} \log Z(\theta) &= \frac{1}{Z(\theta)} \frac{\partial}{\partial \theta} Z(\theta) = \frac{1}{Z(\theta)} \frac{\partial}{\partial \theta} \int f(\mathbf{z}) \exp\{\theta^\top \mathbf{T}(\mathbf{z})\} \\ &= \int \frac{1}{Z(\theta)} f(\mathbf{z}) \exp\{\theta^\top \mathbf{T}(\mathbf{z})\} \cdot \mathbf{T}(\mathbf{z}) = \langle \mathbf{T}(\mathbf{z}) \rangle_{\theta} \end{aligned}$$

Thus, the **M step** solves: $\frac{\partial \mathcal{F}}{\partial \theta} = \langle \mathbf{T}(\mathbf{z}) \rangle_{q(\mathbf{y})} - \langle \mathbf{T}(\mathbf{z}) \rangle_{\theta} = 0$

Proof of the Matrix Inversion Lemma

$$(A + XBX^\top)^{-1} = A^{-1} - A^{-1}X(B^{-1} + X^\top A^{-1}X)^{-1}X^\top A^{-1}$$

Need to prove:

$$(A^{-1} - A^{-1}X(B^{-1} + X^\top A^{-1}X)^{-1}X^\top A^{-1})(A + XBX^\top) = I$$

Expand:

$$I + A^{-1}XBX^\top - A^{-1}X(B^{-1} + X^\top A^{-1}X)^{-1}X^\top - A^{-1}X(B^{-1} + X^\top A^{-1}X)^{-1}X^\top A^{-1}XBX^\top$$

Regroup:

$$\begin{aligned} &= I + A^{-1}X \left(BX^\top - (B^{-1} + X^\top A^{-1}X)^{-1}X^\top - (B^{-1} + X^\top A^{-1}X)^{-1}X^\top A^{-1}XBX^\top \right) \\ &= I + A^{-1}X \left(BX^\top - (B^{-1} + X^\top A^{-1}X)^{-1}B^{-1}BX^\top - (B^{-1} + X^\top A^{-1}X)^{-1}X^\top A^{-1}XBX^\top \right) \\ &= I + A^{-1}X \left(BX^\top - (B^{-1} + X^\top A^{-1}X)^{-1}(B^{-1} + X^\top A^{-1}X)BX^\top \right) \\ &= I + A^{-1}X(BX^\top - BX^\top) = I \end{aligned}$$

KL[q(x)||p(x)] ≥ 0, with equality iff ∀x : p(x) = q(x)

First consider discrete distributions; the Kullback-Liebler divergence is:

$$\text{KL}[q||p] = \sum_i q_i \log \frac{q_i}{p_i}.$$

To minimize wrt distribution q we need a **Lagrange multiplier** to enforce normalisation:

$$E \stackrel{\text{def}}{=} \text{KL}[q||p] + \lambda(1 - \sum_i q_i) = \sum_i q_i \log \frac{q_i}{p_i} + \lambda(1 - \sum_i q_i)$$

Find conditions for stationarity

$$\left. \begin{aligned} \frac{\partial E}{\partial q_i} &= \log q_i - \log p_i + 1 - \lambda = 0 \Rightarrow q_i = p_i \exp(\lambda - 1) \\ \frac{\partial E}{\partial \lambda} &= 1 - \sum_i q_i = 0 \Rightarrow \sum_i q_i = 1 \end{aligned} \right\} \Rightarrow q_i = p_i.$$

Check sign of curvature (Hessian):

$$\frac{\partial^2 E}{\partial q_i \partial q_i} = \frac{1}{q_i} > 0, \quad \frac{\partial^2 E}{\partial q_i \partial q_j} = 0,$$

so unique stationary point $q_i = p_i$ is indeed a minimum. Easily verified that at that minimum, $\text{KL}[q||p] = \text{KL}[p||p] = 0$.

A similar proof holds for continuous densities, using functional derivatives.

Maxima in \mathcal{F} correspond to maxima in ℓ

Let θ^* now be the parameter value at a local maximum of \mathcal{F} (and thus at a fixed point)

Differentiating the previous expression wrt θ again we find

$$\frac{d^2}{d\theta^2} \ell(\theta) = \frac{d^2}{d\theta^2} \langle \log P(\mathcal{Y}, \mathcal{X}|\theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} - \frac{d^2}{d\theta^2} \langle \log P(\mathcal{Y}|\mathcal{X}, \theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)}$$

The first term on the right is negative (a maximum) and the second term is positive (a minimum). Thus the curvature of the likelihood is negative and

θ^* is a maximum of ℓ .

[... as long as the derivatives exist. They sometimes don't (zero-noise ICA)].

Fixed Points of EM are Stationary Points in ℓ

Let a fixed point of EM occur with parameter θ^* . Then:

$$\left. \frac{\partial}{\partial \theta} \langle \log P(\mathcal{Y}, \mathcal{X} | \theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} \right|_{\theta^*} = 0$$

$$\begin{aligned} \text{Now, } \ell(\theta) &= \log P(\mathcal{X}|\theta) = \langle \log P(\mathcal{X}|\theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} \\ &= \left\langle \log \frac{P(\mathcal{Y}, \mathcal{X}|\theta)}{P(\mathcal{Y}|\mathcal{X}, \theta)} \right\rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} \\ &= \langle \log P(\mathcal{Y}, \mathcal{X}|\theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} - \langle \log P(\mathcal{Y}|\mathcal{X}, \theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} \end{aligned}$$

$$\text{so, } \frac{d}{d\theta} \ell(\theta) = \frac{d}{d\theta} \langle \log P(\mathcal{Y}, \mathcal{X}|\theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} - \frac{d}{d\theta} \langle \log P(\mathcal{Y}|\mathcal{X}, \theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)}$$

The second term is 0 at θ^* if the derivative exists (minimum of $\text{KL}[\cdot||\cdot]$), and thus:

$$\left. \frac{d}{d\theta} \ell(\theta) \right|_{\theta^*} = \left. \frac{d}{d\theta} \langle \log P(\mathcal{Y}, \mathcal{X}|\theta) \rangle_{P(\mathcal{Y}|\mathcal{X}, \theta^*)} \right|_{\theta^*} = 0$$

So, EM converges to a stationary point of $\ell(\theta)$.