The EM Algorithm

Further Statistical Methods, Lecture 6 Hilary Term 2004

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The EM algorithm

The EM algorithm is an alternative to Newton-Raphson or the method of scoring for computing MLE in cases where the complications in calculating the MLE are due to *incomplete observation* and data are *MAR*, missing at random, with *separate parameters* for observation and the missing data mechanism, so the missing data mechanism can be ignored.

Data (X, Y) are the *complete data* whereas only incomplete data Y = y are observed. (Rubin uses $Y = Y_{obs}$ and $X = Y_{mis}$).

The complete data log-likelihood is:

 $l(\theta) = \log L(\theta; x, y) = \log f(x, y; \theta).$

The marginal log-likelihood or incomplete data log-likelihood is based on y alone and is equal to

$$l_y(\theta) = \log L(\theta; y) = \log f(y; \theta).$$

We wish to maximize l_y in θ but l_y is typically quite unpleasant:

$$l_y(\theta) = \log \int f(x, y; \theta) \, dx.$$

The EM algorithm is a method of maximizing the latter iteratively and alternates between two steps, one known as the *E-step* and one as the *M-step*, to be detailed below.

We let θ^* be and arbitrary but fixed value, typically the value of θ at the current iteration.

The E-step calculates the expected complete data log-likelihood ratio $q(\theta | \theta^*)$:

$$q(\theta \mid \theta^*) = \mathbf{E}_{\theta^*} \left[\log \frac{f(X, y; \theta)}{f(X, y; \theta^*)} \mid Y = y \right]$$
$$= \int \log \frac{f(x, y; \theta)}{f(x, y; \theta^*)} f(x \mid y; \theta^*) \, dx.$$

The M-step maximizes $q(\theta\,|\,\theta^*)$ in θ for for fixed $\theta^*,$ i.e. calculates

$$\theta^{**} = \arg \max_{\theta} q(\theta \mid \theta^*).$$

We will show that after an E-step and subsequent M-step, the likelihood function has never decreased.

Kullback-Leibler divergence

The *KL divergence* between f and g is $KL(f:g) = \int f(x) \log \frac{f(x)}{a(x)} dx.$

Also known as *relative entropy* of g with respect to f. Since $-\log x$ is a convex function, Jensen's inequality gives $KL(f:g) \ge 0$ and KL(f:g) = 0 if and only if f = g, since

$$KL(f:g) = \int f(x) \log \frac{f(x)}{g(x)} dx \ge -\log \int f(x) \frac{g(x)}{f(x)} dx = 0,$$

so KL divergence defines an (asymmetric) distance measure between probability distributions.

Expected and marginal log-likelihood

Since $f(x \mid y; \overline{\theta}) = f\{(x, y); \theta\} / f(y; \theta)$ we have $q(\theta \mid \theta^*) = \int \log \frac{f(y; \theta) f(x \mid y; \theta)}{f(y; \theta^*) f(x \mid y; \theta^*)} f(x \mid y; \theta^*) dx$ $= \log f(y; \theta) - \log f(y; \theta^*)$

$$+ \int \log \frac{f(x \mid y; \theta)}{f(x \mid y; \theta^*)} f(x \mid y; \theta^*) dx$$
$$= l_y(\theta) - l_y(\theta^*) - KL(f_{\theta^*}^y : f_{\theta}^y).$$

Since the KL-divergence is minimized for $\theta = \theta^*$, differentiation of the above expression yields

$$\frac{\partial}{\partial \theta} q(\theta \mid \theta^*) \bigg|_{\theta = \theta^*} = \left. \frac{\partial}{\partial \theta} l_y(\theta) \right|_{\theta = \theta^*}$$

Let now $\theta_0 = \theta^*$ and define the iteration

$$\theta_{n+1} = \arg \max_{\theta} q(\theta \mid \theta_n).$$

Then

$$l_y(\theta_{n+1}) = l_y(\theta_n) + q(\theta_{n+1} | \theta_n) + KL(f_{\theta_{n+1}}^y : f_{\theta_n}^y)$$

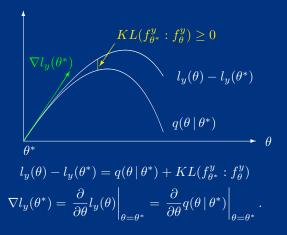
$$\geq l_y(\theta_n) + 0 + 0.$$

So the log-likelihood never decreases after a combined E-step and M-step.

It follows that any limit point must be a saddle point or a local maximum of the likelihood function.

The picture on the next overhead should show it all.

Expected and complete data likelihood



Mixtures

Consider a sample $Y = (Y_1, \ldots, Y_n)$ from individual densities

$$f(y;\alpha,\mu) = \{\alpha\phi(y-\mu) + (1-\alpha)\phi(y)\}$$

where ϕ is the normal density

$$\phi(y) = \frac{1}{\sqrt{2\pi}} e^{-y^2/2}$$

and α and μ are both unknown, $0 < \alpha < 1$.

This corresponds to a fraction α of the observations being contaminated, or originating from a different population.

Incomplete observation

The likelihood function becomes

$$L_y(\alpha,\mu) = \prod_i \left\{ \alpha \phi(y_i - \mu) + (1 - \alpha)\phi(y_i) \right\}$$

is quite unpleasant, although both Newton–Raphson and the method of scoring can be used.

But suppose we knew which observations came from which population?

In other words, let $X = (X_1, \ldots, X_n)$ be i.i.d. with $P(X_i = 1) = \alpha$ and suppose that the conditional distribution of Y_i given $X_i = 1$ was $\mathcal{N}(\mu, 1)$ whereas given $X_i = 0$ it was $\mathcal{N}(0, 1)$, i.e. that X_i was indicating whether Y_i was contaminated or not.

Then the marginal distribution of Y is precisely the mixture distribution and the 'complete data likelihood' is

$$L_{x,y}(\alpha,\mu) = \prod_{i} \alpha^{x_{i}} \phi(y_{i}-\mu)^{x_{i}} (1-\alpha)^{1-x_{i}} \phi(y_{i})^{1-x_{i}}$$

$$\propto \alpha^{\sum x_{i}} (1-\alpha)^{n-\sum x_{i}} \prod_{i} \phi(y_{i}-\mu)^{x_{i}}$$

so taking logarithms we get (ignoring a constant) that

$$l_{x,y}(\alpha,\mu) = \sum_{i} x_i \log \alpha + \left(n - \sum_{i} x_i\right) \log(1-\alpha) \\ - \sum_{i} x_i (y_i - \mu)^2 / 2.$$

If we did not know how to maximize this explicitly,

differentiation easily leads to:

$$\hat{\alpha} = \sum x_i/n, \quad \hat{\mu} = \sum x_i y_i / \sum x_i.$$

Thus, when complete data are available the frequency of contaminated observations is estimated by the observed frequency and the mean μ of these is estimated by the average among the contaminated observations.

E-step and M-step

By taking expectations, we get the E-step as

$$q(\alpha, \mu \,|\, \alpha^*, \mu^*) = \mathbf{E}_{\alpha^*, \mu^*} \{ l_{X, y}(\alpha, \mu) \,|\, Y = y \} \\ = \sum_{i} x_i^* \log \alpha + \left(n - \sum_{i} x_i^* \right) \log(1 - \alpha) \\ - \sum_{i} x_i^* (y_i - \mu)^2 / 2$$

where

$$x_i^* = \mathbf{E}_{\alpha^*,\mu^*}(X_i \mid Y_i = y_i) = P_{\alpha^*,\mu^*}(X_i = 1 \mid Y_i = y_i).$$

Since this has the same form as the complete data likelihood, just with x_i^* replacing x_i , the M-step simply

becomes

$$\alpha^{**} = \sum x_i^*/n, \quad \mu^{**} = \sum x_i^* y_i / \sum x_i^*,$$

i.e. here the mean of the contaminated observations is estimated by a weighted average of all the observations, the weight being proportional to the probability that this observation is contaminated. In effect, x_i^* act as *imputed values* of x_i .

The imputed values x_i^* needed in the E-step are calculated as follows:

$$\begin{aligned} x_i^* &= \mathbf{E}(X_i \mid Y_i = y_i) = P(X_i = 1 \mid Y_i = y_i) \\ &= \frac{\alpha^* \phi(y_i - \mu^*)}{\alpha^* \phi(y_i - \mu^*) + (1 - \alpha^*) \phi(y_i)}. \end{aligned}$$