

Mixture Models & EM

Rishabh Iyer
University of Texas at Dallas

based on the slides of Nick Rouzzi and Vibhav Gogate

Previously...



- We looked at k-means and hierarchical clustering as mechanisms for unsupervised learning
 - k-means was simply a block coordinate descent scheme for a specific objective function
- Today: how to learn probabilistic models for unsupervised learning problems

EM: Soft Clustering



- Clustering (e.g., k-means) typically assumes that each instance is given a "hard" assignment to exactly one cluster
- Does not allow uncertainty in class membership or for an instance to belong to more than one cluster
 - Problematic because data points that lie roughly midway between cluster centers are assigned to one cluster
- Soft clustering gives probabilities that an instance belongs to each of a set of clusters

Probabilistic Clustering

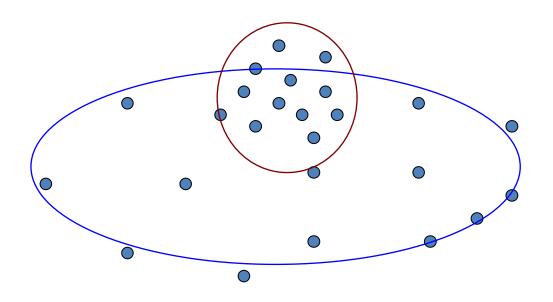


- Try a probabilistic model!
 - Allows overlaps, clusters of different size, etc.
- Can tell a generative story for data
 - p(x|y) p(y)
- Challenge: we need to estimate model parameters without labeled y's (i.e., in the unsupervised setting)

Z	X ₁	X ₂
??	0.1	2.1
??	0.5	-1.1
??	0.0	3.0
??	-0.1	-2.0
??	0.2	1.5
•••	•••	•••

Probabilistic Clustering





- Clusters of different shapes and sizes
- Clusters can overlap! (k-means doesn't allow this)

Finite Mixture Models



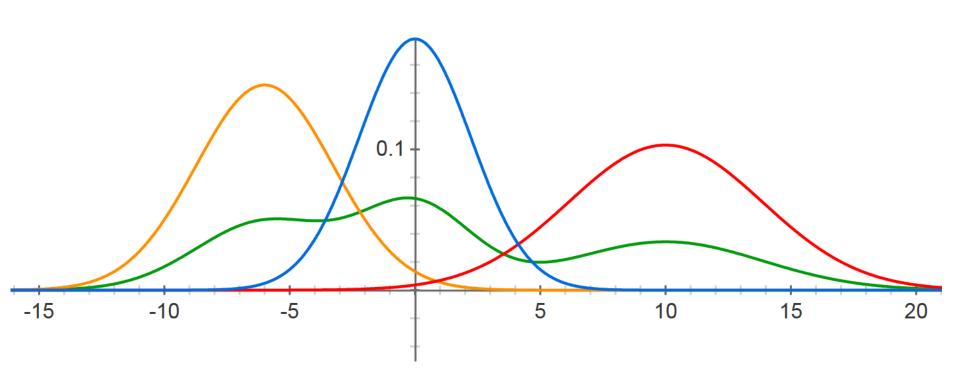
- Given a dataset: $x^{(1)}, \dots, x^{(N)}$
- Mixture model: $\Theta = \{\lambda_1, ..., \lambda_k, \theta_1, ..., \theta_k\}$

$$p(x|\Theta) = \sum_{y=1}^{k} \lambda_y p_y(x|\theta_y)$$

- $p_y(x|\theta_y)$ is a mixture component from some family of probability distributions parameterized by θ_y and $\lambda \geq 0$ such that $\sum_{\nu} \lambda_{\nu} = 1$ are the mixture weights
- We can think of $\lambda_y = p(Y = y | \Theta)$ for some random variable Y that takes values in $\{1, \dots, k\}$

Finite Mixture Models





Uniform mixture of 3 Gaussians

Multivariate Gaussian



• A d-dimensional multivariate Gaussian distribution is defined by a $d \times d$ covariance matrix Σ and a mean vector μ

$$p(x|\mu,\Sigma) = \frac{1}{\sqrt{(2\pi)^{\mathrm{d}}\mathrm{det}(\Sigma)}} \exp\left(-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)\right)$$

- The covariance matrix describes the degree to which pairs of variables vary together
 - The diagonal elements correspond to variances of the individual variables

Multivariate Gaussian



• A d-dimensional multivariate Gaussian distribution is defined by a $d \times d$ covariance matrix Σ and a mean vector μ

$$p(x|\mu,\Sigma) = \frac{1}{\sqrt{(2\pi)^{\mathrm{d}} \mathrm{det}(\Sigma)}} \exp\left(-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)\right)$$

- The covariance matrix must be a symmetric positive definite matrix in order for the above to make sense
 - Positive definite: all eigenvalues are positive & matrix is invertible
 - Ensures that the quadratic form is concave

Gaussian Mixture Models (GMMs)



• We can define a GMM by choosing the k^{th} component of the mixture to be a Gaussian density with parameters

$$\theta_k = \{\mu_k, \Sigma_k\}$$

$$p(x|\mu_k, \Sigma_k) = \frac{1}{\sqrt{(2\pi)^d \det(\Sigma_k)}} \exp\left(-\frac{1}{2}(x - \mu_k)^T \Sigma_k^{-1}(x - \mu_k)\right)$$

We could cluster by fitting a mixture of k Gaussians to our data

How do we learn these kinds of models?

Learning Gaussian Parameters



MLE for supervised univariate Gaussian

$$\mu_{MLE} = \frac{1}{N} \sum_{i=1}^{N} \chi^{(i)}$$

$$\sigma_{MLE}^{2} = \frac{1}{N} \sum_{i=1}^{N} (x^{(i)} - \mu_{MLE})^{2}$$

MLE for supervised multivariate Gaussian

$$\mu_{MLE} = \frac{1}{N} \sum_{i=1}^{N} x^{(i)}$$

$$\Sigma_{\text{MLE}} = \frac{1}{N} \sum_{i=1}^{N} (x^{(i)} - \mu_{MLE}) (x^{(i)} - \mu_{MLE})^{T}$$

Learning Gaussian Parameters



• MLE for supervised multivariate mixture of k Gaussian distributions

$$\mu_{MLE}^k = \frac{1}{|M_k|} \sum_{i \in M_k} x^{(i)}$$

$$\Sigma_{MLE}^{k} = \frac{1}{|M_k|} \sum_{i \in M_k} (x^{(i)} - \mu_{MLE}^k) (x^{(i)} - \mu_{MLE}^k)^T$$

Sums are over the observations that were generated by the k^{th} mixture component (this requires that we know which points were generated by which distribution!)

The Unsupervised Case



- What if our observations do not include information about which of the k mixture components generated them?
- Consider a joint probability distribution over data points, $x^{(i)}$, and mixture assignments, $y \in \{1, ..., k\}$

$$\arg \max_{\Theta} \prod_{i=1}^{N} p(x^{(i)}|\Theta) = \arg \max_{\Theta} \prod_{i=1}^{N} \sum_{y=1}^{k} p(x^{(i)}, Y = y|\Theta)$$

$$= \arg \max_{\Theta} \prod_{i=1}^{N} \sum_{y=1}^{k} p(x^{(i)}|Y = y, \Theta) p(Y = y|\Theta)$$

The Unsupervised Case



- What if our observations do not include information about which of the k mixture components generated them?
- Consider a joint probability distribution over data points, $x^{(i)}$, and mixture assignments, $y \in \{1, ..., k\}$

$$\arg \max_{\Theta} \prod_{i=1}^{N} p(x^{(i)}|\Theta) = \arg \max_{\Theta} \prod_{i=1}^{N} \sum_{y=1}^{k} p(x^{(i)}, Y = y|\Theta)$$

$$= \arg \max_{\Theta} \prod_{i=1}^{N} \sum_{y=1}^{k} p(x^{(i)}|Y = y, \Theta) p(Y = y|\Theta)$$

We only know how to compute the probabilities for each mixture component

The Unsupervised Case



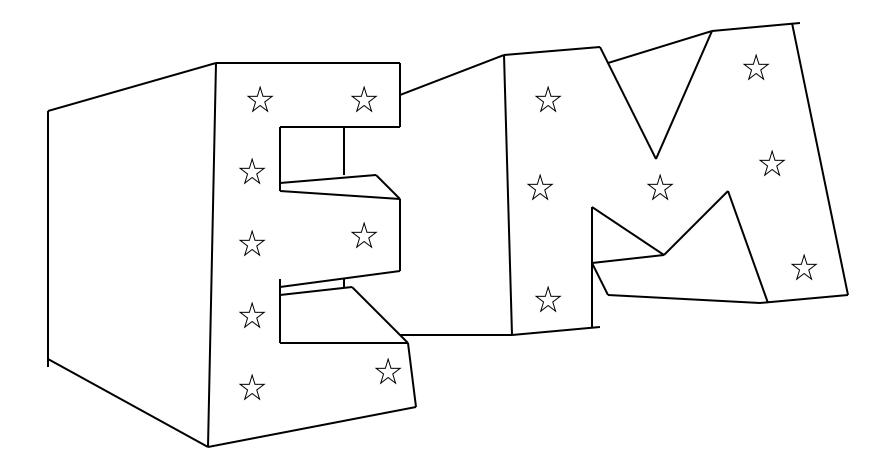
In the case of a Gaussian mixture model

$$p(x^{(i)}|Y = y, \Theta) = N(x^{(i)}|\mu_y, \Sigma_y)$$
$$p(Y = y|\Theta) = \lambda_y$$

The solution: modify the objective to make the optimization easier

Expectation Maximization





Jensen's Inequality



For a convex function $f: \mathbb{R}^n \to \mathbb{R}$, any $a_1, \dots, a_k \in [0,1]$ such that $a_1 + \dots + a_k = 1$, and any $x^{(1)}, \dots, x^{(k)} \in \mathbb{R}^n$,

$$a_1 f(x^{(1)}) + \dots + a_k f(x^{(k)}) \ge f(a_1 x^{(1)} + \dots + a_k x^{(k)})$$

Inequality is reversed for concave functions



$$\log \ell(\Theta) = \sum_{i=1}^{N} \log \sum_{y=1}^{k} p(x^{(i)}, Y = y | \Theta)$$

$$= \sum_{i=1}^{N} \log \sum_{y=1}^{k} \frac{q_i(y)}{q_i(y)} p(x^{(i)}, Y = y | \Theta)$$

$$\geq \sum_{i=1}^{N} \sum_{y=1}^{k} q_i(y) \log \frac{p(x^{(i)}, Y = y | \Theta)}{q_i(y)}$$

$$\equiv F(\Theta, q)$$



$$\log \ell(\Theta) = \sum_{i=1}^{N} \log \sum_{y=1}^{k} p(x^{(i)}, Y = y | \Theta)$$

$$= \sum_{i=1}^{N} \log \sum_{y=1}^{k} \frac{q_i(y)}{q_i(y)} p(x^{(i)}, Y = y | \Theta)$$

$$\geq \sum_{i=1}^{N} \sum_{y=1}^{k} q_i(y) \log \frac{p(x^{(i)}, Y = y | \Theta)}{q_i(y)} \text{ arbitrary positive probability distribution}$$

$$\equiv F(\Theta, q)$$



$$\log \ell(\Theta) = \sum_{i=1}^{N} \log \sum_{y=1}^{k} p(x^{(i)}, Y = y | \Theta)$$

$$= \sum_{i=1}^{N} \log \sum_{y=1}^{k} \frac{q_i(y)}{q_i(y)} p(x^{(i)}, Y = y | \Theta)$$

$$\geq \sum_{i=1}^{N} \sum_{y=1}^{k} q_i(y) \log \frac{p(x^{(i)}, Y = y | \Theta)}{q_i(y)}$$

$$= F(\Theta, q)$$
Jensen's ineq.



$$\arg\max_{\Theta,q_1,\dots,q_N} \sum_{i=1}^N \sum_{v=1}^k q_i(y) \log \frac{p(x^{(i)}, Y = y | \Theta)}{q_i(y)}$$

- This objective is not jointly concave in Θ and $q_1, ..., q_N$
 - Best we can hope for is a local maxima (and there could be A LOT of them)
- The EM algorithm is a block coordinate ascent scheme that finds a local optimum of this objective
 - Start from an initialization Θ^0 and q_1^0, \dots, q_N^0



• E step: with the θ 's fixed, maximize the objective over q

$$q^{t+1} \in \arg\max_{q_1,...,q_N} \sum_{i=1}^{N} \sum_{y=1}^{k} q_i(y) \log \frac{p(x^{(i)}, Y = y | \Theta^t)}{q_i(y)}$$

• Using the method of Lagrange multipliers for the constraint that $\sum_{y} q_i(y) = 1$ gives

$$q_i^{t+1}(y) = p(Y = y | X = x^{(i)}, \Theta^t)$$



• M step: with the q's fixed, maximize the objective over Θ

$$\theta^{t+1} \in \arg\max_{\Theta} \sum_{i=1}^{N} \sum_{y=1}^{k} q_i^{t+1}(y) \log \frac{p(x^{(i)}, Y = y | \theta)}{q_i^{t+1}(y)}$$

- For the case of GMM, we can compute this update in closed form
 - This is not necessarily the case for every model
 - May require gradient ascent



- Start with random parameters
- E-step maximizes a lower bound on the log-sum for fixed parameters
- M-step solves a MLE estimation problem for fixed probabilities
- Iterate between the E-step and M-step until convergence

EM for Gaussian Mixtures



• E-step:

$$q_i^t(y) = \frac{\lambda_y^t \cdot p(x^{(i)} \middle| \mu_y^t, \Sigma_y^t)}{\sum_{y'} \lambda_{y'}^t \cdot p(x^{(i)} \middle| \mu_{y'}^t, \Sigma_{y'}^t)}$$
Probability of $x^{(i)}$ under the appropriate

M-step:

$$\mu_y^{t+1} = \frac{\sum_{i=1}^{N} q_i^t(y) x^{(i)}}{\sum_{i=1}^{N} q_i^t(y)}$$

Probability of $x^{(i)}$ under the appropriate multivariate normal distribution

$$\Sigma_{y}^{t+1} = \frac{\sum_{i=1}^{N} q_{i}^{t}(y) (x^{(i)} - \mu_{y}^{t+1}) (x^{(i)} - \mu_{y}^{t+1})^{T}}{\sum_{i=1}^{N} q_{i}^{t}(y)}$$

$$\lambda_{y}^{t+1} = \frac{1}{N} \sum_{i=1}^{N} q_{i}^{t}(y)$$

EM for Gaussian Mixtures



• E-step:

$$q_i^t(y) = \frac{\lambda_y^t \cdot p\left(x^{(i)} \middle| \mu_y^t, \Sigma_y^t\right)}{\sum_{y'} \lambda_{y'}^t \cdot p\left(x^{(i)} \middle| \mu_{y'}^t, \Sigma_{y'}^t\right)}$$
Probability of x(i) under the mixture model

M-step:

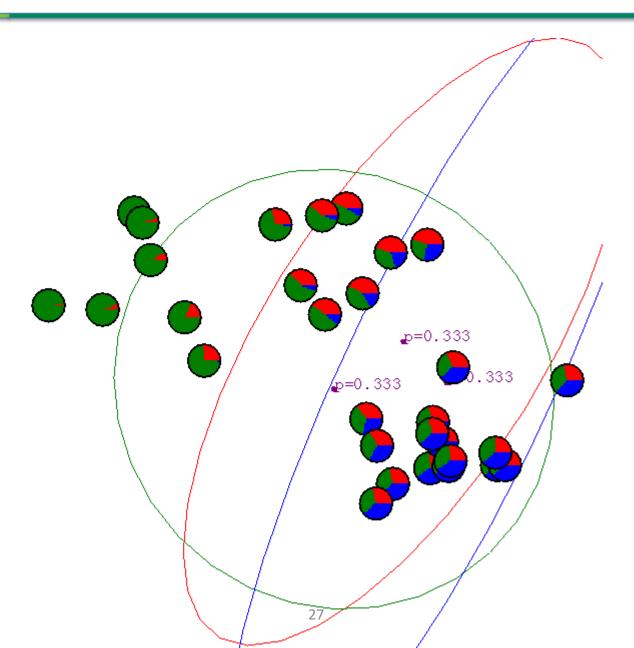
$$\mu_y^{t+1} = \frac{\sum_{i=1}^{N} q_i^t(y) x^{(i)}}{\sum_{i=1}^{N} q_i^t(y)}$$

$$\Sigma_{y}^{t+1} = \frac{\sum_{i=1}^{N} q_{i}^{t}(y) (x^{(i)} - \mu_{y}^{t+1}) (x^{(i)} - \mu_{y}^{t+1})^{T}}{\sum_{i=1}^{N} q_{i}^{t}(y)}$$

$$\lambda_{y}^{t+1} = \frac{1}{N} \sum_{i=1}^{N} q_{i}^{t}(y)$$

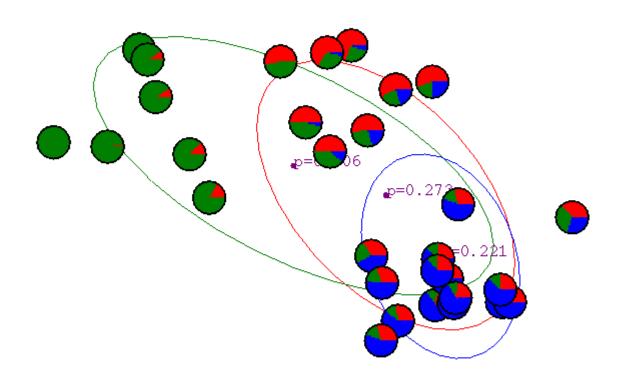
Gaussian Mixture Example: Start





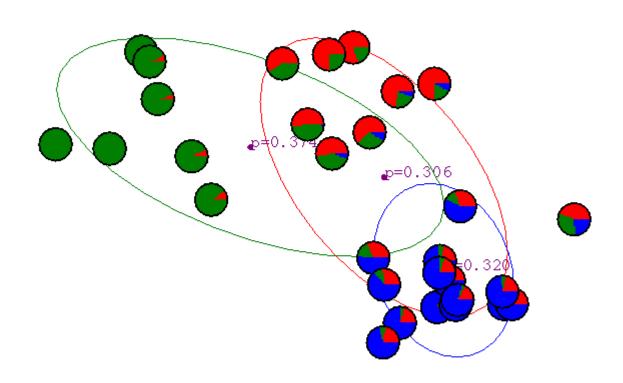
After first iteration





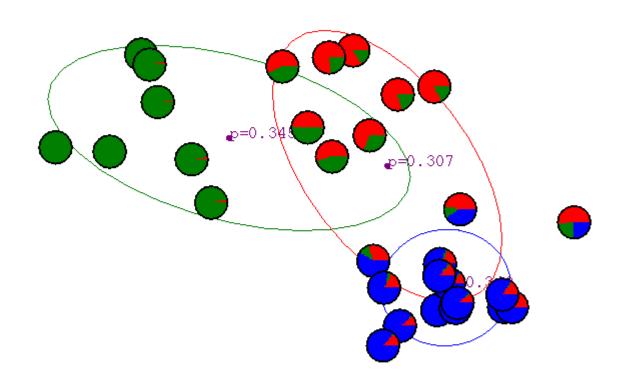
After 2nd iteration





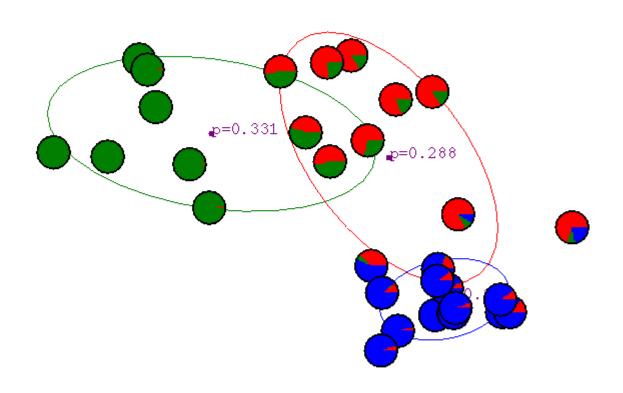
After 3rd iteration





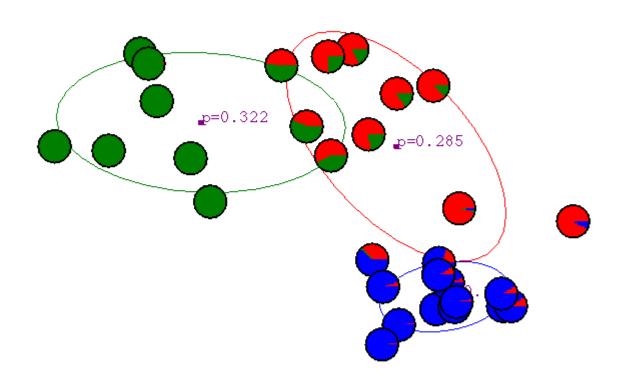
After 4th iteration





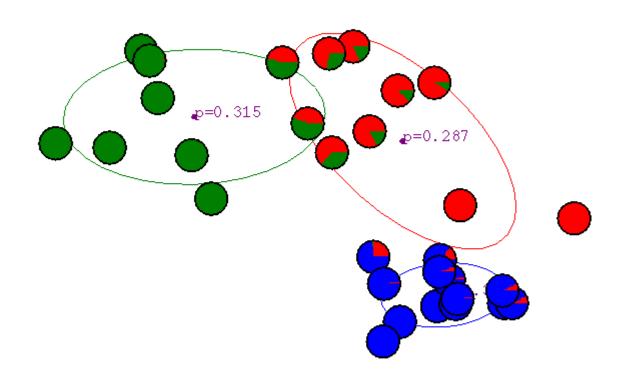
After 5th iteration





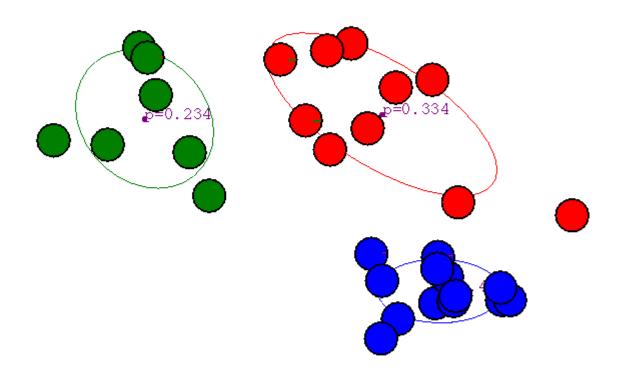
After 6th iteration





After 20th iteration





Properties of EM



- EM converges to a local optimum
 - This is because each iteration improves the log-likelihood
 - Proof same as k-means (just block coordinate ascent)
 - E-step can never decrease likelihood
 - M-step can never decrease likelihood
- If we make hard assignments instead of soft ones, algorithm is equivalent to k-means!