Clustering: K-Means and Mixtures of Gaussians

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Reading: Manning/Raghavan/Schuetze, Chapters 16 (not 16.3) and 17 (http://nlp.stanford.edu/IR-book/)

Outline

- Supervised vs. Unsupervised Learning
- Hierarchical Clustering
 - Hierarchical Agglomerative Clustering (HAC)
- Non-Hierarchical Clustering
 - K-means
 - Mixtures of Gaussians and EM-Algorithm

Non-Hierarchical Clustering

- K-means clustering ("hard")
- Mixtures of Gaussians and training via Expectation maximization Algorithm ("soft")

Clustering Criterion

- Evaluation function that assigns a (usually real-valued) value to a clustering
 - Clustering criterion typically function of
 - within-cluster similarity and
 - between-cluster dissimilarity
- Optimization
 - Find clustering that maximizes the criterion
 - Global optimization (often intractable)
 - Greedy search
 - Approximation algorithms

Centroid-Based Clustering

- Assumes instances are real-valued vectors.
- Clusters represented via centroids (i.e. average of points in a cluster) c:

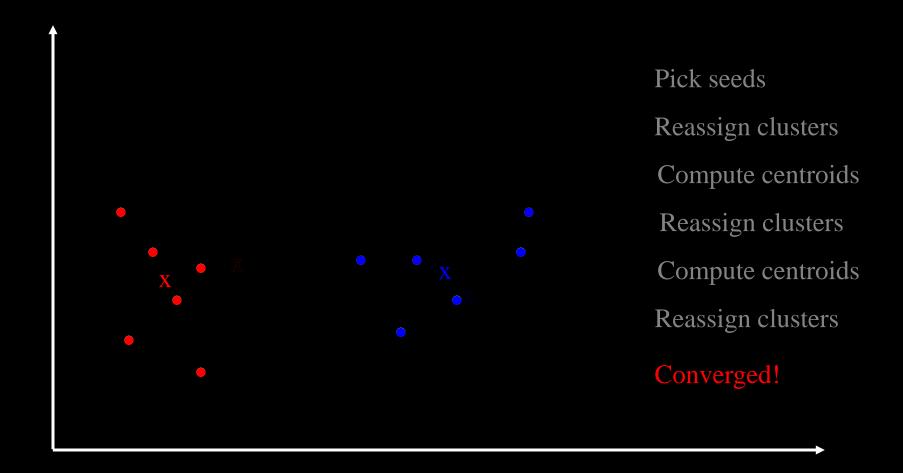
$$\vec{\mu}(c) = \frac{1}{|c|} \sum_{\vec{x} \in c} \vec{x}$$

 Reassignment of instances to clusters is based on distance to the current cluster centroids.

K-Means Algorithm

- Input: k = number of clusters, distance measure d
- Select k random instances $\{s_1, s_2, \dots s_k\}$ as seeds.
- Until clustering converges or other stopping criterion:
 - For each instance x_i :
 - Assign x_i to the cluster c_i such that $d(x_i, s_i)$ is min.
 - For each cluster c_i //update the centroid of each cluster
 - $s_i = \mu(c_i)$

K-means Example (k=2)



Time Complexity

- Assume computing distance between two instances is O(N) where N is the dimensionality of the vectors.
- Reassigning clusters for n points: O(kn) distance computations, or O(knN).
- Computing centroids: Each instance gets added once to some centroid: O(nN).
- Assume these two steps are each done once for i iterations: O(iknN).
- Linear in all relevant factors, assuming a fixed number of iterations, more efficient than HAC.

Buckshot Algorithm

Problem

- Results can vary based on random seed selection, especially for high-dimensional data.
- Some seeds can result in poor convergence rate, or convergence to sub-optimal clusterings.

Idea: Combine HAC and K-means clustering.

- First randomly take a sample of instances of size n^{1/2}
- Run group-average HAC on this sample
- Use the results of HAC as initial seeds for K-means.
- Overall algorithm is efficient and avoids problems of bad seed selection.

Clustering as Prediction

Setup

- Learning Task: P(X)
- I Training Sample: $S=(ec{x}_1$, ... , $ec{x}_n$)
- Hypothesis Space: $H = \{h_1, ..., h_{|H|}\}$ each describes $P(X|h_i)$ where h_i are parameters
- Goal: learn which $P(X|h_i)$ produces the data
- What to predict?
 - Predict where new points are going to fall

Gaussian Mixtures and EM

- Gaussian Mixture Models
 - Assume

$$P(X = \vec{x} | h_i) = \sum_{j=1}^k P(X = \vec{x} | Y = j, h_i) P(Y = j)$$
 where $P(X = \vec{x} | Y = j, h) = N(X = \vec{x} | \vec{\mu}_j, \Sigma_j) = \frac{1}{\sqrt{2\pi\sigma_{ij}^2}} e^{-\frac{1}{2\sigma_{ij}^2}(x - \mu_{ij})^2}$ and $h = (\vec{\mu}_1, ..., \vec{\mu}_k, \Sigma_1, ..., \Sigma_k)$.

- EM Algorithm
 - Assume P(Y) and k known and $\Sigma_i = 1$.
 - REPEAT

•
$$\vec{\mu}_j = \frac{\sum_{i=1}^n P(Y=j|X=\vec{x}_i,\vec{\mu}_1,...,\vec{\mu}_k)\vec{x}_i}{\sum_{i=1}^n P(Y=j|X=\vec{x}_i,\vec{\mu}_1,...,\vec{\mu}_k)}$$

•
$$P(Y = j | X = \vec{x}_i, \vec{\mu}_1, ..., \vec{\mu}_k) = \frac{P(X = \vec{x}_i | Y = j, \vec{\mu}_j) P(Y = j)}{\sum_{l=1}^k P(X = \vec{x}_i | Y = l, \vec{\mu}_l) P(Y = l)} = \frac{e^{-0.5(\vec{x}_i - \vec{\mu}_j)^2} P(Y = j)}{\sum_{l=1}^k e^{-0.5(\vec{x}_i - \vec{\mu}_l)^2} P(Y = l)}$$