## Data Clustering: K-means and Hierarchical Clustering

Piyush Rai

#### CS5350/6350: Machine Learning

October 4, 2011

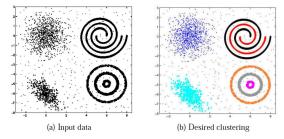
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• Data Clustering is an unsupervised learning problem

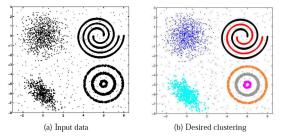
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- Data Clustering is an unsupervised learning problem
- Given: N unlabeled examples  $\{x_1, \ldots, x_N\}$ ; the number of partitions K
- Goal: Group the examples into K partitions



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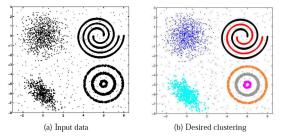
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- Clustering groups examples based of their mutual similarities

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- The only information clustering uses is the similarity between examples
- Clustering groups examples based of their mutual similarities
- A good clustering is one that achieves:
  - High within-cluster similarity
  - Low inter-cluster similarity

Picture courtesy: "Data Clustering: 50 Years Beyond K-Means", A.K. Jain (2008)

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#### Notions of Similarity

- Choice of the similarity measure is very important for clustering
- Similarity is inversely related to distance

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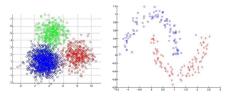
#### Notions of Similarity

- Choice of the similarity measure is very important for clustering
- Similarity is inversely related to distance
- Different ways exist to measure distances. Some examples:
  - Euclidean distance:  $d(\mathbf{x}, \mathbf{z}) = ||\mathbf{x} \mathbf{z}|| = \sqrt{\sum_{d=1}^{D} (x_d z_d)^2}$
  - Manhattan distance:  $d(\mathbf{x}, \mathbf{z}) = \sum_{d=1}^{D} |x_d z_d|$
  - Kernelized (non-linear) distance:  $d(\mathbf{x}, \mathbf{z}) = ||\phi(\mathbf{x}) \phi(\mathbf{z})||$

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- For the left figure above, Euclidean distance may be reasonable
- For the right figure above, kernelized distance seems more reasonable

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## Similarity is Subjective

• Similarity is often hard to define



• Different similarity criteria can lead to different clusterings

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## Data Clustering: Some Real-World Examples

- Clustering images based on their perceptual similarities
- Image segmentation (clustering pixels)



- Clustering webpages based on their content
- Clustering web-search results
- Clustering people in social networks based on user properties/preferences
- ... and many more..

Picture courtesy: http://people.cs.uchicago.edu/~pff/segment/

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### Types of Clustering

- Solutional clustering (K-means, Gaussian mixture models, etc.)
  - Partitions are independent of each other



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## Types of Clustering

- Flat or Partitional clustering (K-means, Gaussian mixture models, etc.)
  - Partitions are independent of each other



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- Partitions can be visualized using a tree structure (a dendrogram)
- Does not need the number of clusters as input
- Possible to view partitions at different levels of granularities (i.e., can refine/coarsen clusters) using different K



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• Input: N examples  $\{\mathbf{x}_1, \ldots, \mathbf{x}_N\}$   $(\mathbf{x}_n \in \mathbb{R}^D)$ ; the number of partitions K

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- Initialize: K cluster centers  $\mu_1, \ldots, \mu_K$ . Several initialization options:
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- Iterate:
  - Assign each of example x<sub>n</sub> to its closest cluster center

$$C_k = \{n: \quad k = \arg\min_k ||\mathbf{x}_n - \mu_k||^2\}$$

 $(\mathcal{C}_k \text{ is the set of examples closest to } \mu_k)$ 

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• Recompute the new cluster centers  $\mu_k$  (mean/centroid of the set  $C_k$ )

$$\mu_k = \frac{1}{|\mathcal{C}_k|} \sum_{n \in \mathcal{C}_k} \mathsf{x}_n$$

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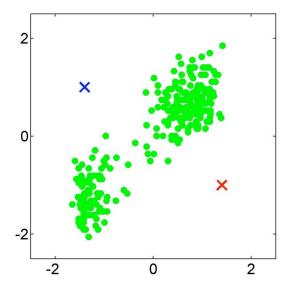
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- Repeat while not converged
- A possible convergence criteria: cluster centers do not change anymore

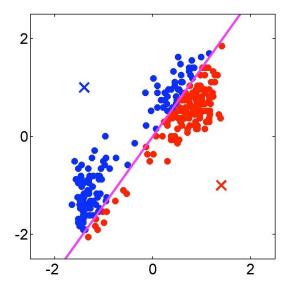
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#### K-means: Initialization (assume K = 2)



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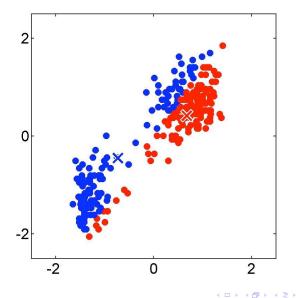
#### K-means iteration 1: Assigning points



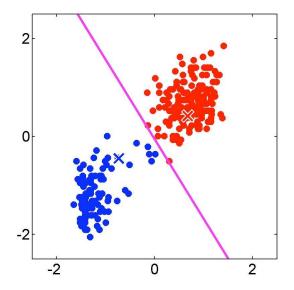
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#### K-means iteration 1: Recomputing the cluster centers

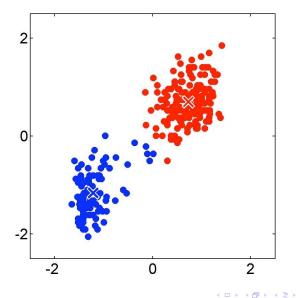


#### K-means iteration 2: Assigning points

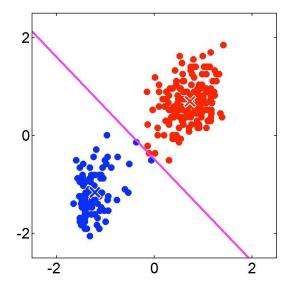


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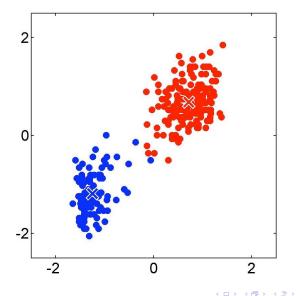


#### K-means iteration 3: Assigning points

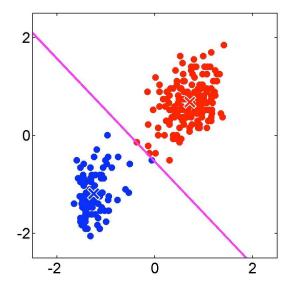


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#### K-means iteration 3: Recomputing the cluster centers

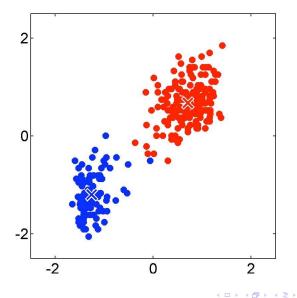


#### K-means iteration 4: Assigning points



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#### K-means iteration 4: Recomputing the cluster centers



### K-means: The Objective Function

The K-means objective function

- Let  $\mu_1, \ldots, \mu_K$  be the K cluster centroids (means)
- Let  $r_{nk} \in \{0,1\}$  be indicator denoting whether point  $\mathbf{x}_n$  belongs to cluster k
- *K*-means objective minimizes the total distortion (sum of distances of points from their cluster centers)

$$J(\mu, r) = \sum_{n=1}^{N} \sum_{k=1}^{K} r_{nk} ||\mathbf{x}_{n} - \mu_{k}||^{2}$$

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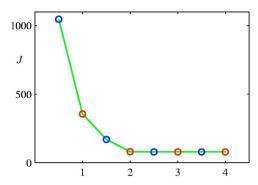
$$J(\mu, \mathbf{r}) = \sum_{n=1}^{N} \sum_{k=1}^{K} r_{nk} ||\mathbf{x}_n - \mu_k||^2$$

- Note: Exact optimization of the K-means objective is NP-hard
- The K-means algorithm is a heuristic that converges to a local optimum

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#### K-means: Choosing the number of clusters K

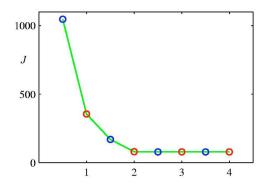
• One way to select K for the K-means algorithm is to try different values of K, plot the K-means objective versus K, and look at the "elbow-point" in the plot



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• For the above plot, K = 2 is the elbow point



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  - Try multiple initializations and choose the best result
  - Other smarter initialization schemes (e.g., look at the *K*-means++ algorithm by Arthur and Vassilvitskii)

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#### K-means: Limitations

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  - K-medians algorithm is a more robust alternative for data with outliers
  - Reason: Median is more robust than mean in presence of outliers

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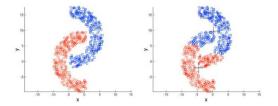
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  - K-medians algorithm is a more robust alternative for data with outliers
  - Reason: Median is more robust than mean in presence of outliers
- Works well only for round shaped, and of roughtly equal sizes/density clusters
- Does badly if the clusters have non-convex shapes
  - Spectral clustering or kernelized K-means can be an alternative

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# K-means Limitations Illustrated

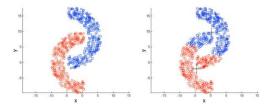
Non-convex/non-round-shaped clusters: Standard K-means fails!



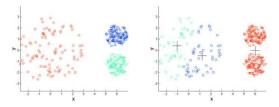
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# K-means Limitations Illustrated

Non-convex/non-round-shaped clusters: Standard K-means fails!



Clusters with different densities

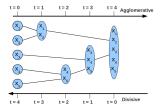


Picture courtesy: Christof Monz (Queen Mary, Univ. of London)

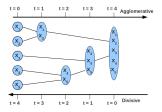
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Data Clustering

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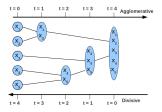


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#### Agglomerative (bottom-up) Clustering

- Start with each example in its own singleton cluster
- At each time-step, greedily merge 2 most similar clusters
- Stop when there is a single cluster of all examples, else go to 2

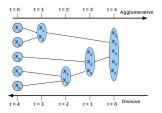


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#### Divisive (top-down) Clustering

- Start with all examples in the same cluster
- At each time-step, remove the "outsiders" from the least cohesive cluster
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#### • Divisive (top-down) Clustering

- Start with all examples in the same cluster
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- Stop when each example is in its own singleton cluster, else go to 2
- Agglomerative is more popular and simpler than divisive (but less accurarate)

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- How to compute the dissimilarity between two clusters R and S?

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- Min-link or single-link: results in chaining (clusters can get very large)

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• Max-link or complete-link: results in small, round shaped clusters

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$$d(R,S) = \max_{\mathbf{x}_R \in R, \mathbf{x}_S \in S} d(\mathbf{x}_R, \mathbf{x}_S)$$

• Average-link: compromise between single and complexte linkage

$$d(R,S) = \frac{1}{|R||S|} \sum_{\mathbf{x}_R \in R, \mathbf{x}_S \in S} d(\mathbf{x}_R, \mathbf{x}_S)$$

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- We know how to compute the dissimilarity  $d(\mathbf{x}_i, \mathbf{x}_j)$  between two examples
- How to compute the dissimilarity between two clusters R and S?
- Min-link or single-link: results in chaining (clusters can get very large)

$$d(R,S) = \min_{\mathbf{x}_R \in R, \mathbf{x}_S \in S} d(\mathbf{x}_R, \mathbf{x}_S)$$

• Max-link or complete-link: results in small, round shaped clusters

$$d(R,S) = \max_{\mathbf{x}_R \in R, \mathbf{x}_S \in S} d(\mathbf{x}_R, \mathbf{x}_S)$$

• Average-link: compromise between single and complexte linkage

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$$d(R, S) = \frac{1}{|R||S|} \sum_{\mathbf{x}_R \in R, \mathbf{x}_S \in S} d(\mathbf{x}_R, \mathbf{x}_S)$$
(a) MIN (single link.) (b) MAX (complete link.) (c) Group average.

# Flat vs Hierarchical Clustering

- Flat clustering produces a single partitioning
- Hierarchical Clustering can give different partitionings depending on the level-of-resolution we are looking at
- Flat clustering needs the number of clusters to be specified
- Hierarchical clustering doesn't need the number of clusters to be specified
- Flat clustering is usually more efficient run-time wise
- Hierarchical clustering can be slow (has to make several merge/split decisions)
- No clear consensus on which of the two produces better clustering

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