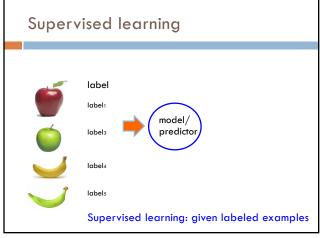
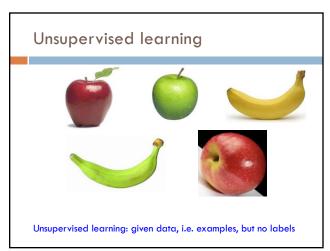


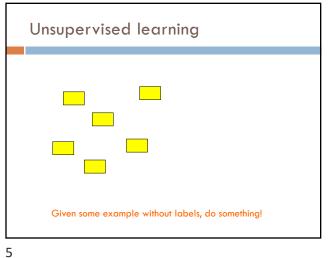
Administrative Final project □ Project proposal feedback soon □ Progress report due next Wednesday

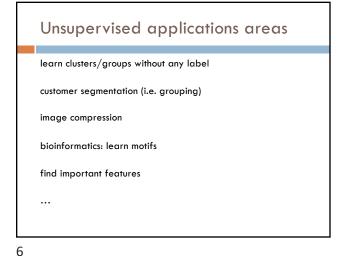
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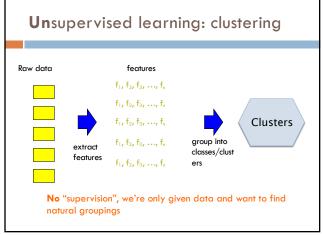




3 4



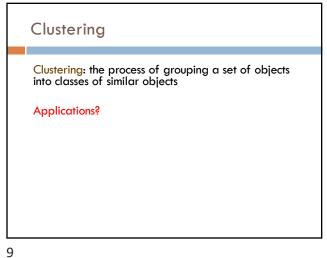




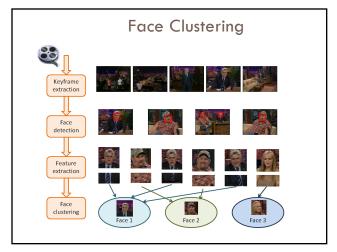
Unsupervised learning: modeling Most frequently, when people think of unsupervised learning they think clustering Another category: learning probabilities/parameters for models without supervision Learn a translation dictionary ■ Learn a grammar for a language Learn the social graph

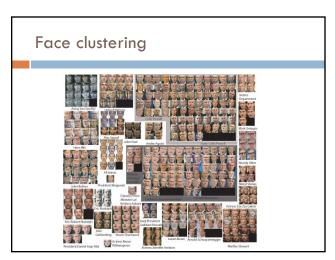
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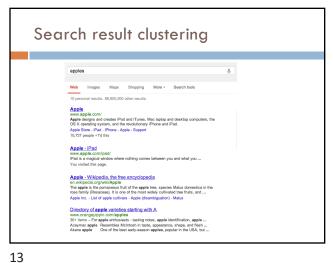
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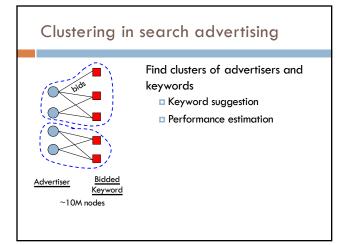


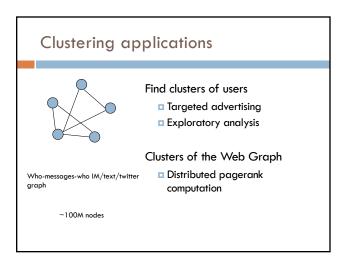


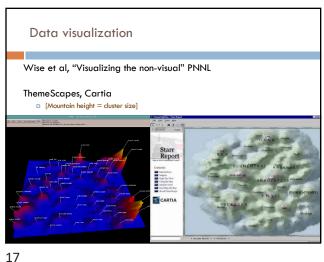


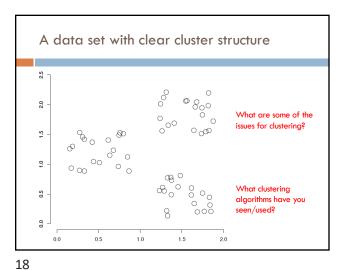








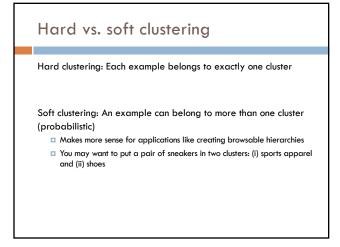


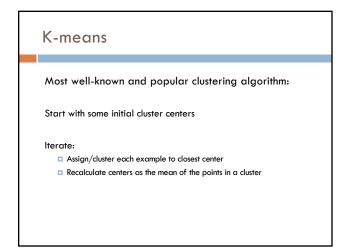


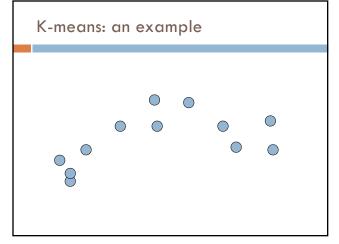
Issues for clustering Representation for clustering ■ How do we represent an example features, etc. □ Similarity/distance between examples Flat clustering or hierarchical Number of clusters □ Fixed a priori □ Data driven?

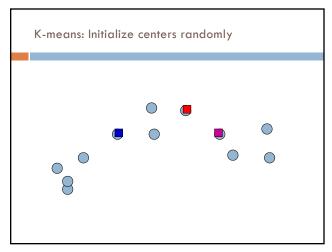
Clustering Algorithms Flat algorithms Usually start with a random (partial) partitioning ■ Refine it iteratively ■ K means clustering ■ Model based clustering ■ Spectral clustering Hierarchical algorithms ■ Bottom-up, agglomerative □ Top-down, divisive

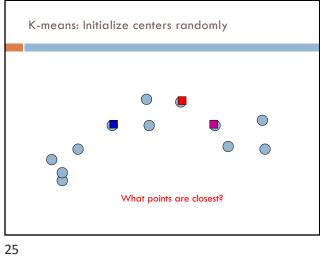
19 20

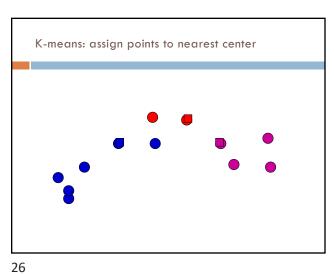


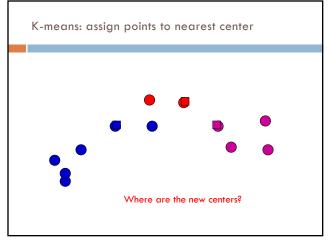


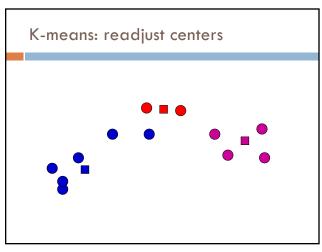


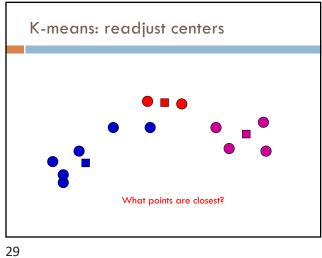


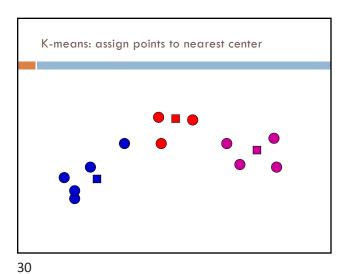


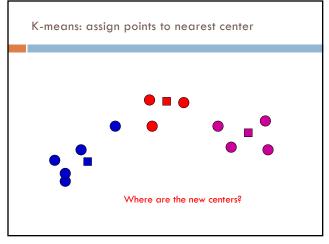


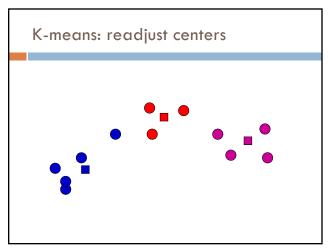


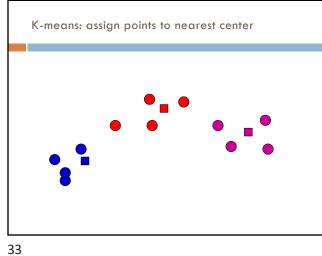


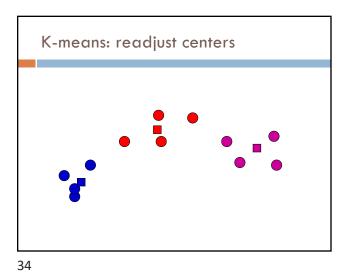


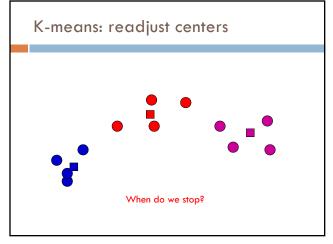


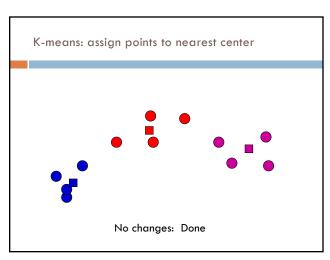


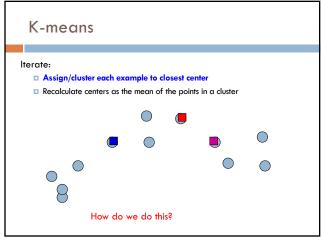












Iterate:

Assign/cluster each example to closest center iterate over each point:

- get distance to each cluster center
- assign to closest center (hard cluster)

Recalculate centers as the mean of the points in a cluster

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Iterate:

• Assign/cluster each example to closest center iterate over each point:

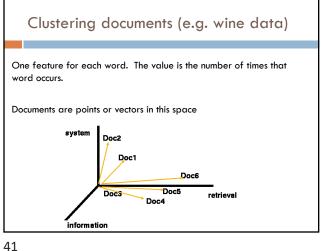
• get distance to each cluster center

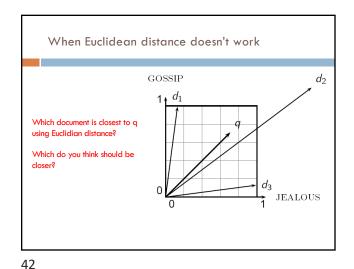
• assign to closest center (hard cluster)

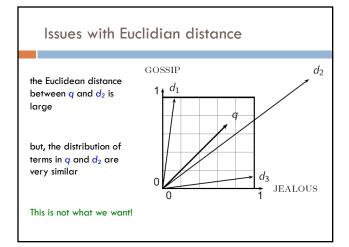
• Recalculate centers as the mean of the points in a cluster

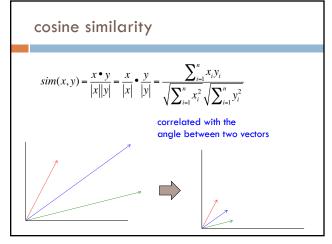
What distance measure should we use?

Distance measures $d(x,y) = \sqrt{\sum_{i=1}^{n} (x_i - y_i)^2}$ good for spatial data









cosine distance

cosine similarity ranges from 0 and 1, with things that are similar 1 and dissimilar θ

cosine distance:

$$d(x, y) = 1 - sim(x, y)$$

- good for text data and many other "real world" data sets
 computationally friendly since we only need to consider
- features that have non-zero values for **both** examples

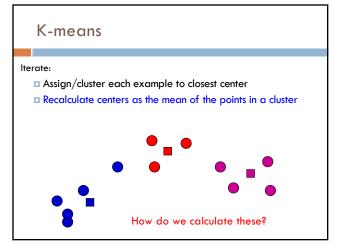
Iterate:

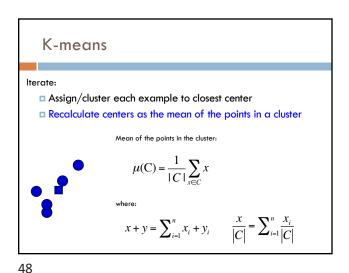
Assign/cluster each example to closest center
Recalculate centers as the mean of the points in a cluster

Where are the cluster centers?

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K-means loss function

K-means tries to minimize what is called the "k-means" loss function:

$$loss = \sum_{i=1}^{n} d(x_i, \mu_k)^2 \text{ where } \mu_k \text{ is cluster center for } x_i$$

the sum of the squared distances from each point to the associated cluster center

Minimizing k-means loss

lterate:

- 1. Assign/cluster each example to closest center
- 2. Recalculate centers as the mean of the points in a cluster

$$loss = \sum_{i=1}^{n} d(x_i, \mu_k)^2 \text{ where } \mu_k \text{ is cluster center for } x_i$$

Does each step of k-means move towards reducing this loss function (or at least not increasing it)?

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Minimizing k-means loss

lterate

- 1. Assign/cluster each example to closest center
- 2. Recalculate centers as the mean of the points in a cluster

$$loss = \sum_{i=1}^{n} d(x_i, \mu_k)^2 \text{ where } \mu_k \text{ is cluster center for } x_i$$

This isn't quite a complete proof/argument, but:

- 1. Any other assignment would end up in a larger loss
- 2. The mean of a set of values minimizes the squared error

Minimizing k-means loss

terate.

- 1. Assign/cluster each example to closest center
- 2. Recalculate centers as the mean of the points in a cluster

$$loss = \sum_{i=1}^{n} d(x_i, \mu_k)^2 \text{ where } \mu_k \text{ is cluster center for } x_i$$

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Minimizing k-means loss Iterate: 1. Assign/cluster each example to closest center 2. Recalculate centers as the mean of the points in a cluster $loss = \sum_{i=1}^{n} d(x_i, \mu_k)^2 \text{ where } \mu_k \text{ is cluster center for } x_i$ NO! It will find a minimum. Unfortunately, the k-means loss function is generally not convex and for most problems has many, many minima

K-means variations/parameters

Start with some initial cluster centers

Iterate:

Assign/cluster each example to closest center
Recalculate centers as the mean of the points in a cluster

What are some other variations/parameters we haven't specified?

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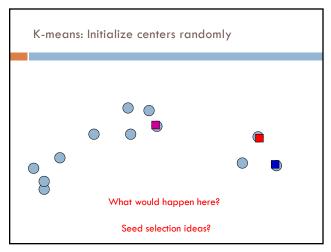
We're only guaranteed to find one of them

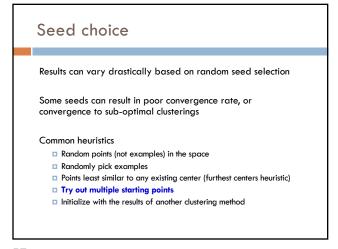
K-means variations/parameters

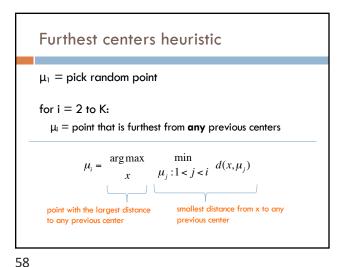
Initial (seed) cluster centers

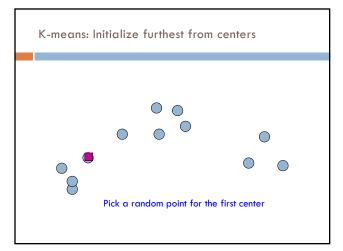
Convergence

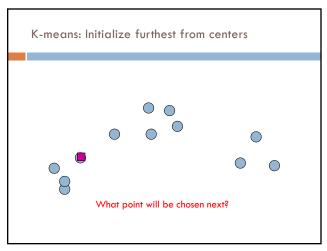
A fixed number of iterations
partitions unchanged
Cluster centers don't change



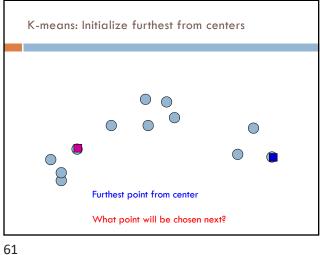


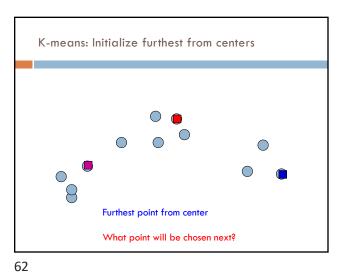


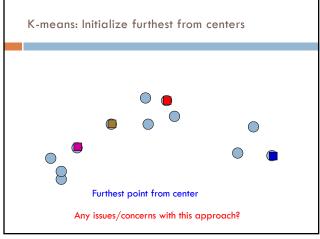


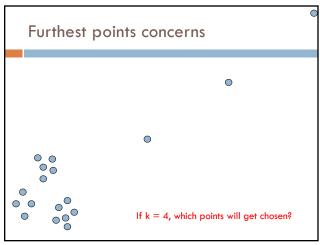


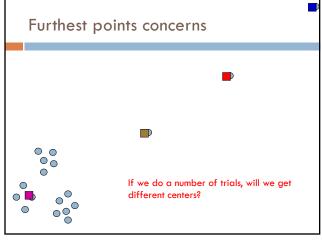
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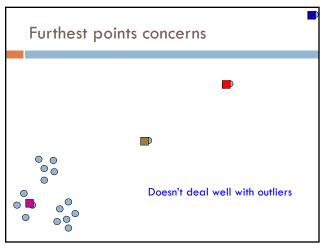






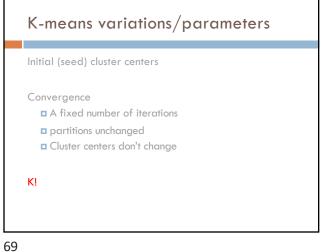


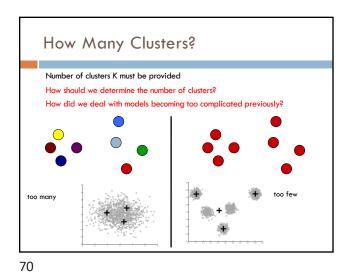


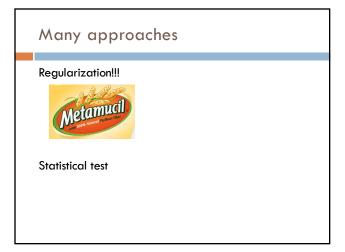


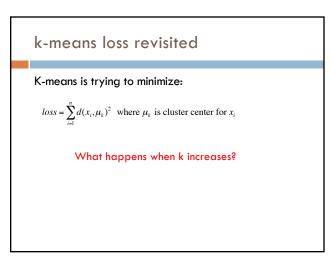
 $\begin{array}{l} \textbf{K-means++} \\ \\ \mu_1 = \text{pick random point} \\ \\ \text{for } k=2 \text{ to } \textbf{K}: \\ \\ \text{for } i=1 \text{ to } \textbf{N}: \\ \\ \\ s_i = \min \ d(x_i, \mu_{1...k-1}) \ / / \text{ smallest distance to any center} \\ \\ \\ \mu_k = \text{randomly pick point proportionate to s} \\ \\ \\ \text{How does this help?} \end{array}$

 $\begin{aligned} & \text{K-means++} \\ & \mu_1 = \text{pick random point} \\ & \text{for } k = 2 \text{ to } \textbf{K}: \\ & \text{for } i = 1 \text{ to } \textbf{N}: \\ & s_i = \min d(x_{i_r} \mu_{1...k-1}) \text{// smallest distance to any center} \\ & \mu_k = \text{randomly pick point } \text{proportionate to s} \\ & \text{- Makes it possible to select other points} \\ & \text{- if } \text{\#points} >> \text{\#outliers, we will pick good points} \\ & \text{- Makes it non-deterministic, which will help with random runs} \\ & \text{- Nice theoretical guarantees!} \end{aligned}$









k-means loss revisited

K-means is trying to minimize:

 $loss = \sum_{i=1}^{n} d(x_i, \mu_k)^2 \text{ where } \mu_k \text{ is cluster center for } x_i$

Loss goes down!

Making the model more complicated allows us more flexibility, but can "overfit" to the data

k-means loss revisited

K-means is trying to minimize:

 $loss_{kmeans} = \sum_{i=1}^{n} d(x_i, \mu_k)^2$ where μ_k is cluster center for x_i



 $loss_{BIC} = loss_{kmeans} + K \log N$ (where N = number of points)

 $loss_{AIC} = loss_{kmeans} + KN$

What effect will this have?
Which will tend to produce smaller k?

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k-means loss revisited

 $loss_{BIC} = loss_{kmeans} + K \log N$ (where N = number of points)

 $loss_{AIC} = loss_{kmeans} + KN$

AIC penalizes increases in K more harshly

Both require a change to the K-means algorithm

Tend to work reasonably well in practice if you don't know ${\sf K}$