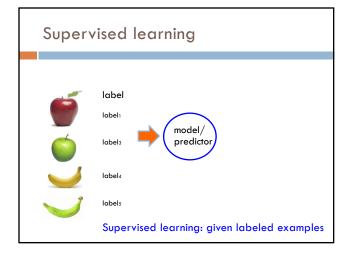
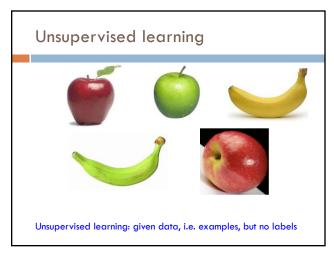
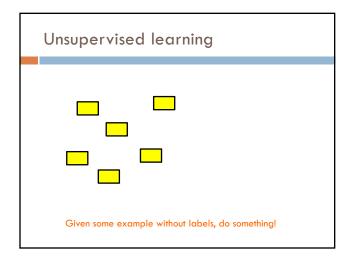
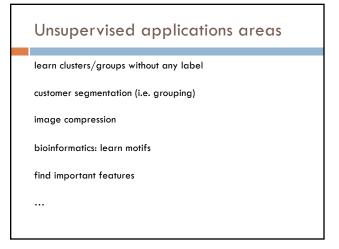


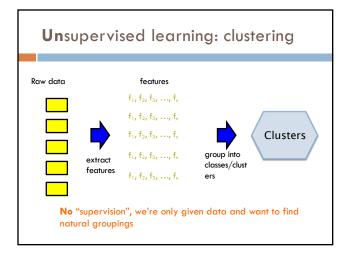
# Final project Project proposal due tomorrow (Wednesday) Progress report due next Wednesday Grading Assignment 8 back soon Assignment 9 next week

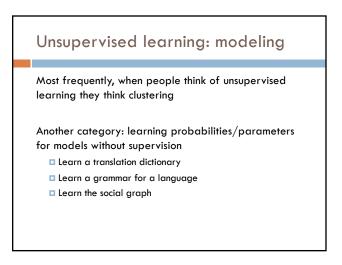


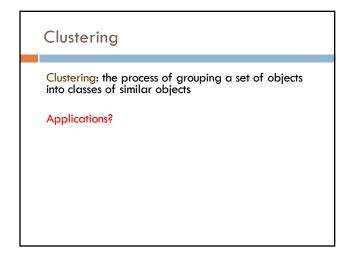




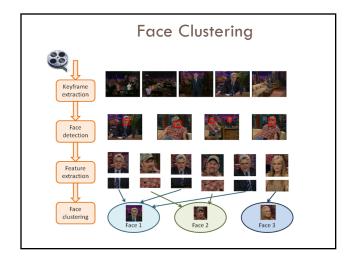


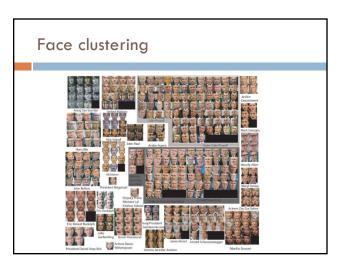


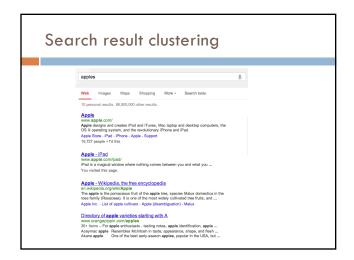




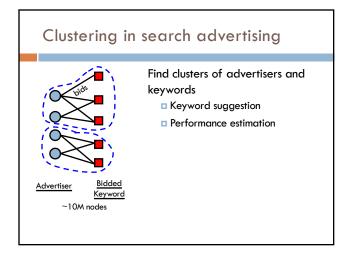


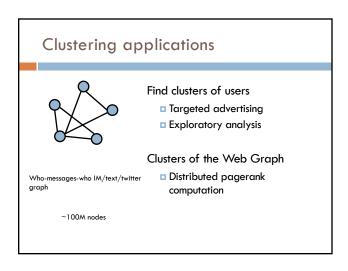


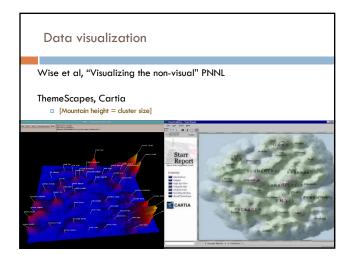


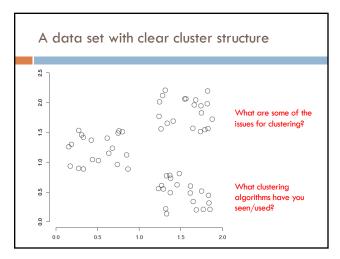


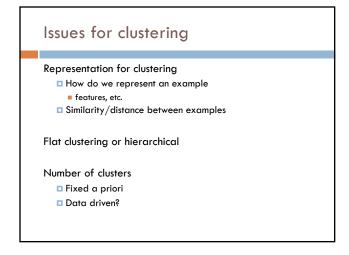


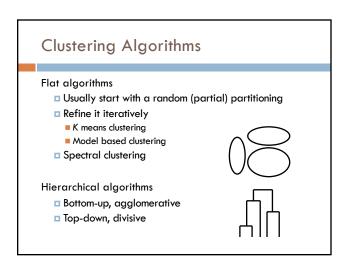




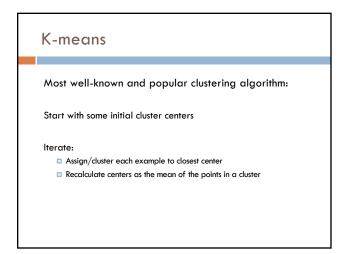


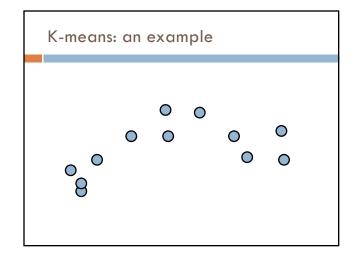


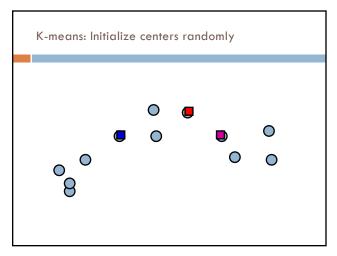


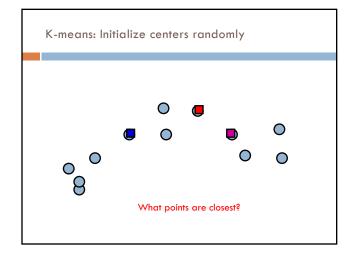


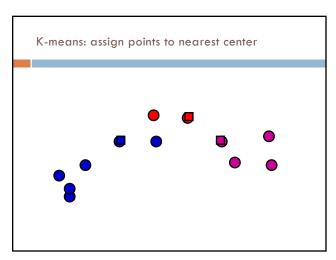
# Hard vs. soft clustering Hard clustering: Each example belongs to exactly one cluster Soft clustering: An example can belong to more than one cluster (probabilistic) Makes more sense for applications like creating browsable hierarchies You may want to put a pair of sneakers in two clusters: (i) sports apparel and (ii) shoes

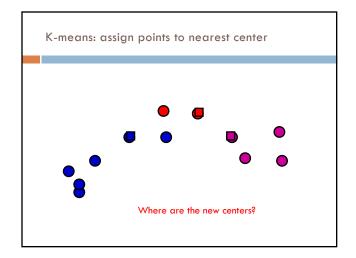


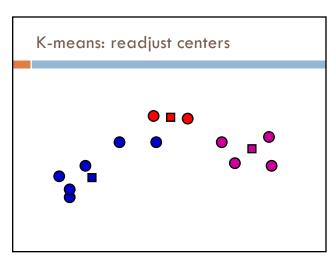


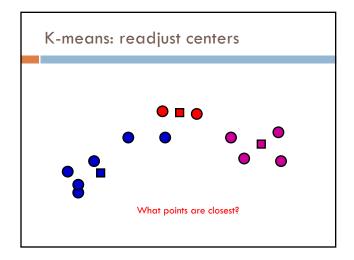


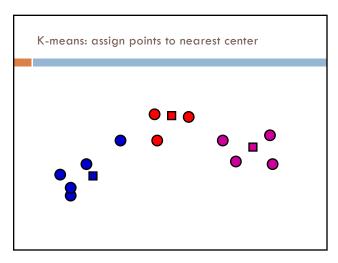


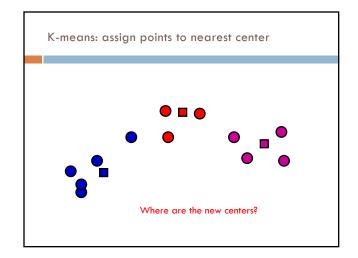


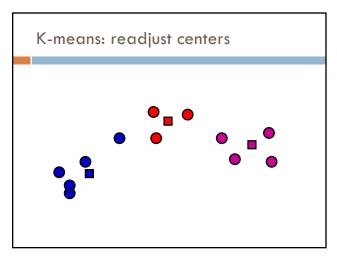


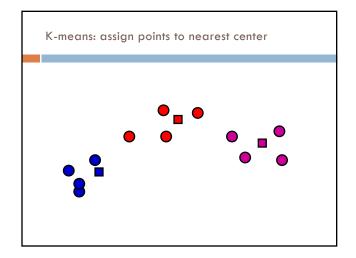


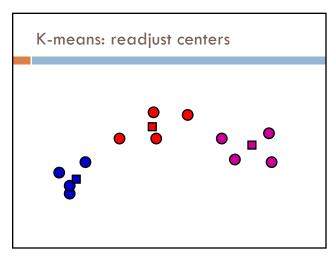


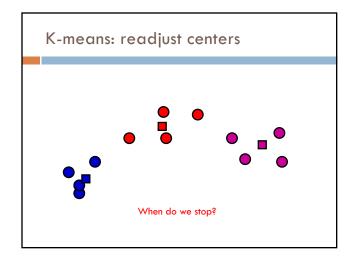


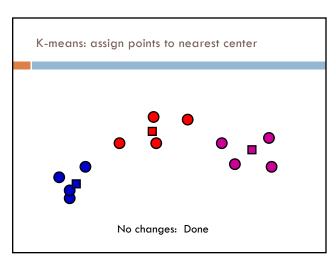


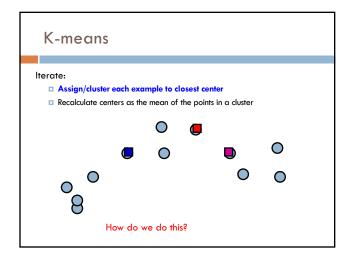


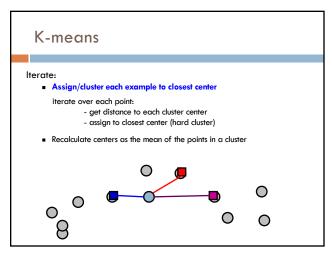


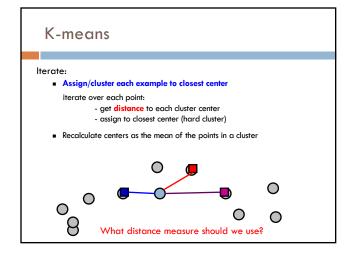


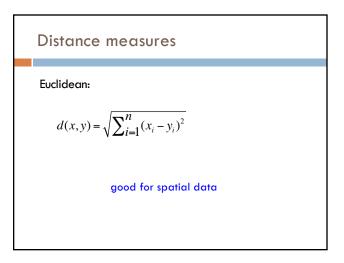


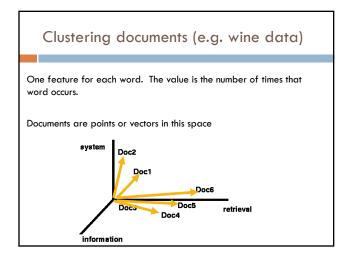


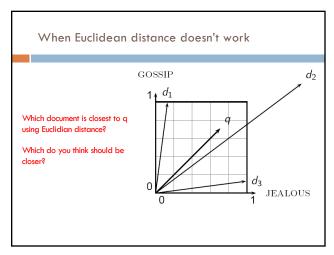


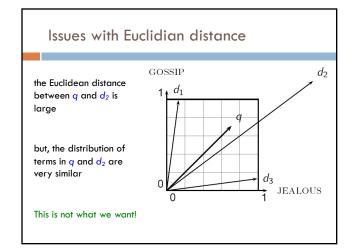


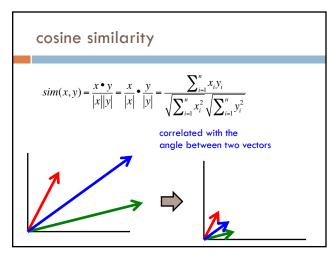












## cosine distance

cosine similarity ranges from 0 and 1, with things that are similar 1 and dissimilar  $\mathbf{0}$ 

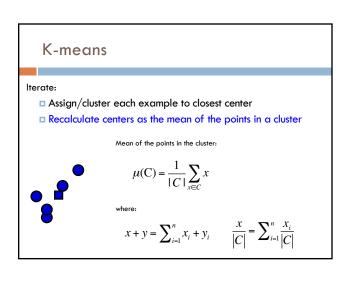
cosine distance:

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

- 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?

# Iterate: Assign/cluster each example to closest center Recalculate centers as the mean of the points in a cluster How do we calculate these?



## 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

### Iterate

- 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)?

# Minimizing k-means loss

### lterate

- 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

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

Does this mean that k-means will always find the minimum loss/clustering?

## 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

We're only guaranteed to find one of them

# 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?

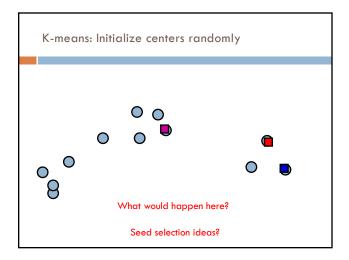
# K-means variations/parameters

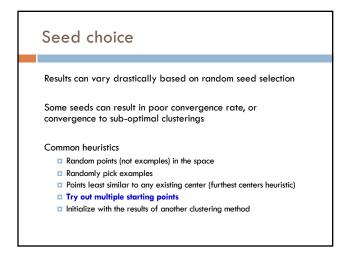
Initial (seed) cluster centers

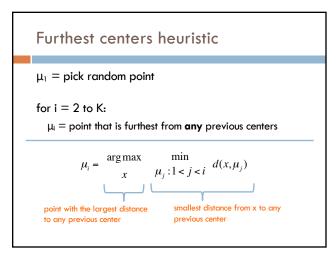
### Convergence

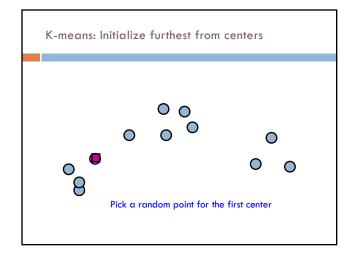
- A fixed number of iterations
- partitions unchanged
- Cluster centers don't change

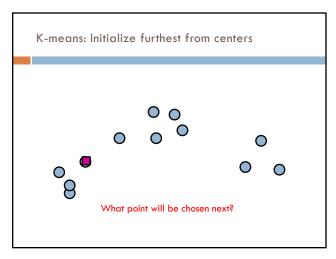
K!

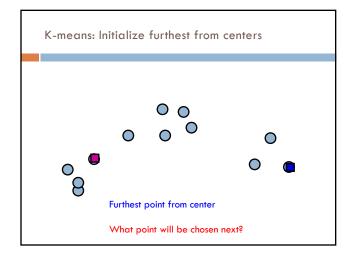


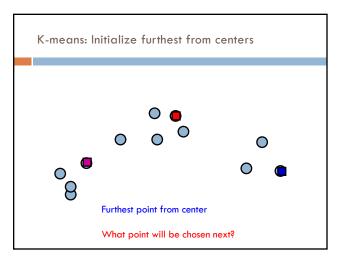


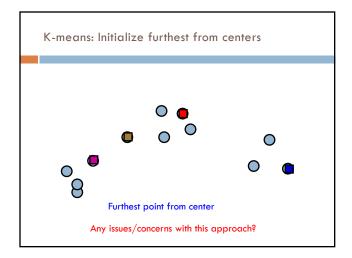


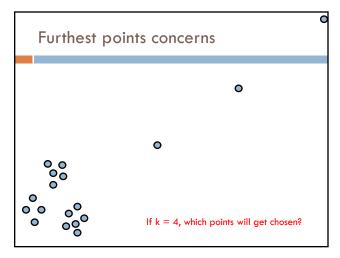


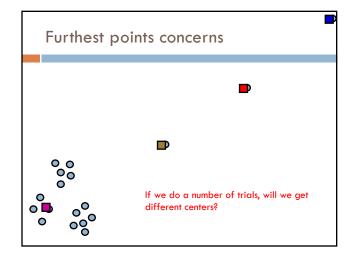


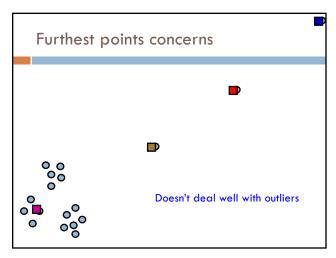




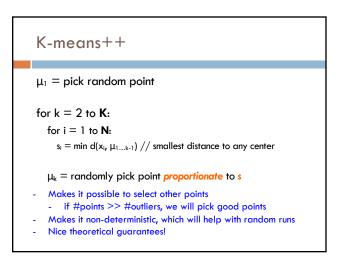


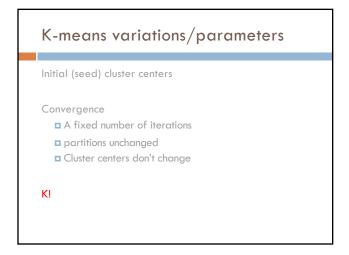


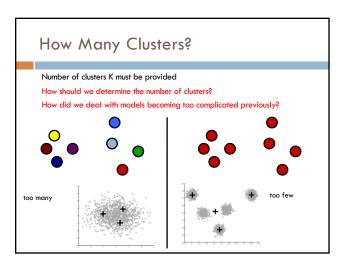


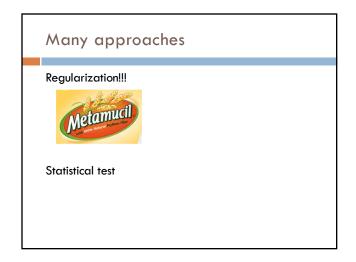


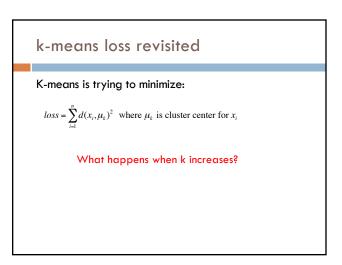
# $\begin{aligned} & \text{K-means++} \\ & \mu_1 = \text{pick random point} \\ & \text{for } k = 2 \text{ to } \textbf{K}: \\ & \text{for } i = 1 \text{ to } \textbf{N}: \\ & s_i = \min d(x_{i_r} \mu_{1...k\cdot 1}) \text{// smallest distance to any center} \\ & \mu_k = \text{randomly pick point } \textit{proportionate}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











## 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  $\mu_k$  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?

### 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  ${\sf K}$ 

### Midterm

Mean: 84%

Q1:79%

Median: 85%

Q3: 90%