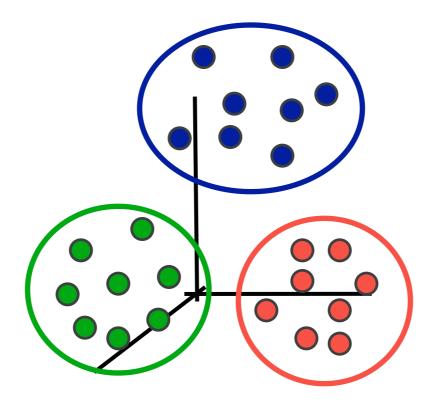
CS 565: Data mining

- Clustering: David Arthur, Sergei Vassilvitskii. *k-means* ++: The Advantages of Careful Seeding. In SODA 2007
- Thanks A. Gionis and S. Vassilvitskii for the slides

What is clustering?

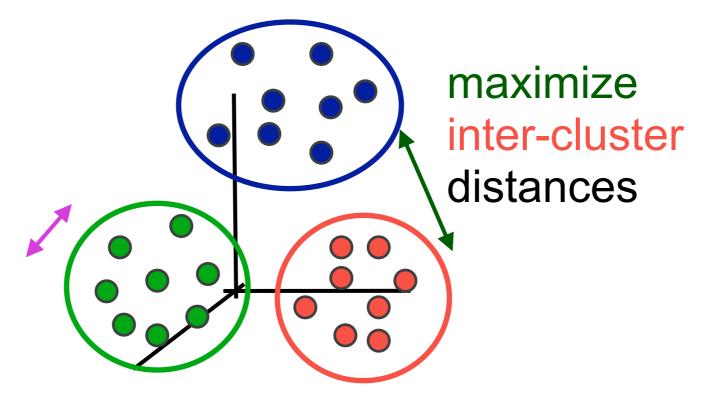
 a grouping of data objects such that the objects within a group are similar (or near) to one another and dissimilar (or far) from the objects in other groups



How to capture this objective?

a grouping of data objects such that the objects within a group are similar (or near) to one another and dissimilar (or far) from the objects in other groups

minimize intra-cluster distances



The clustering problem

- Given a collection of data objects
- Find a grouping so that
 - similar objects are in the same cluster
 - dissimilar objects are in different clusters
- Why we care ?
- stand-alone tool to gain insight into the data
 - visualization
- preprocessing step for other algorithms
 - indexing or compression often relies on clustering

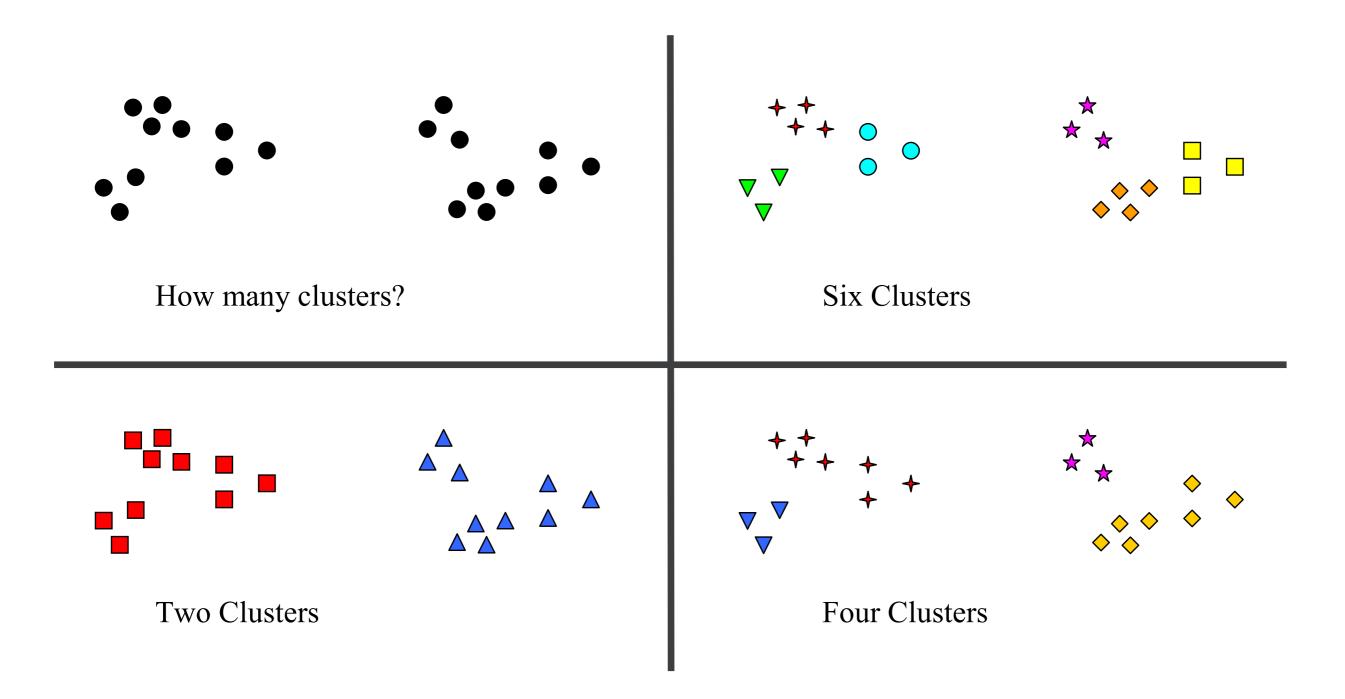
Applications of clustering

- image processing
 - cluster images based on their visual content
- web mining
 - cluster groups of users based on their access patterns on webpages
 - cluster webpages based on their content
- bioinformatics
 - cluster similar proteins together (similarity wrt chemical structure and/or functionality etc)
- many more...

The clustering problem

- Given a collection of data objects
- Find a grouping so that
 - similar objects are in the same cluster
 - dissimilar objects are in different clusters
- Basic questions:
 - what does similar mean?
 - what is a good partition of the objects?
 i.e., how is the quality of a solution measured?
 - how to find a good partition?

Notion of a cluster can be ambiguous



Types of clusterings

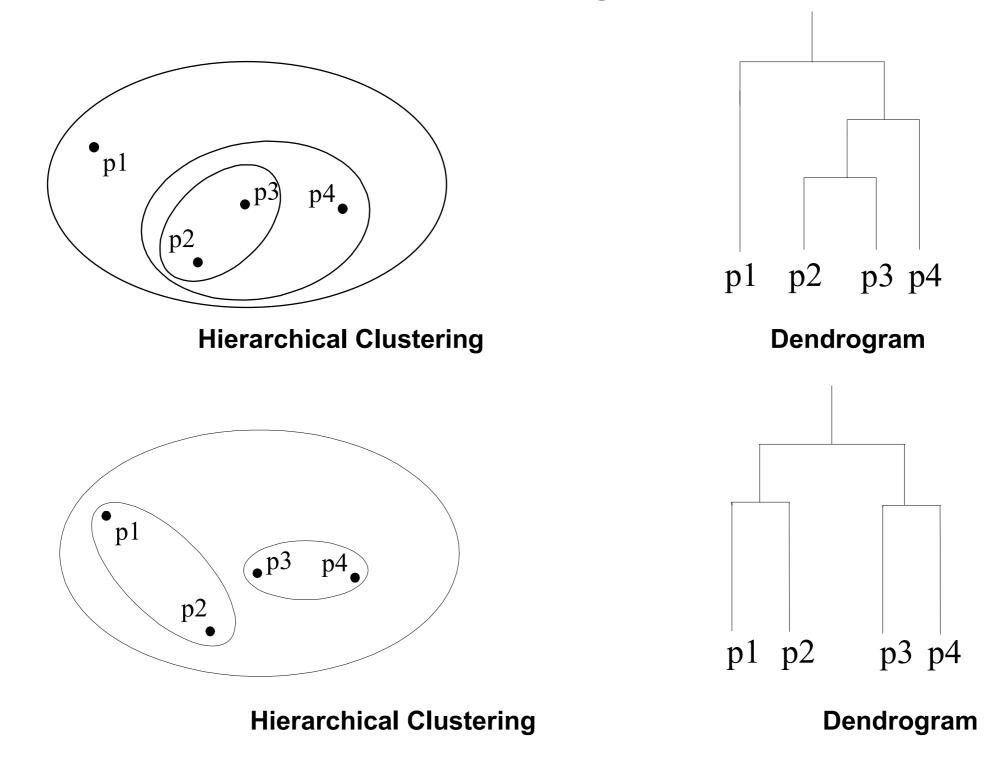
Partitional

each object belongs in exactly one cluster

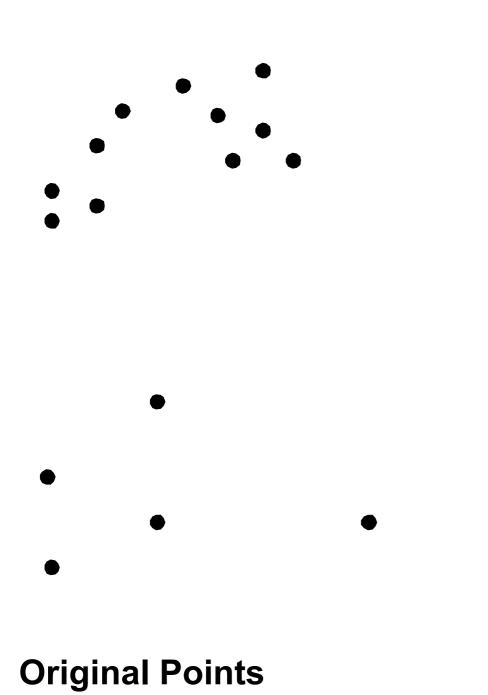
Hierarchical

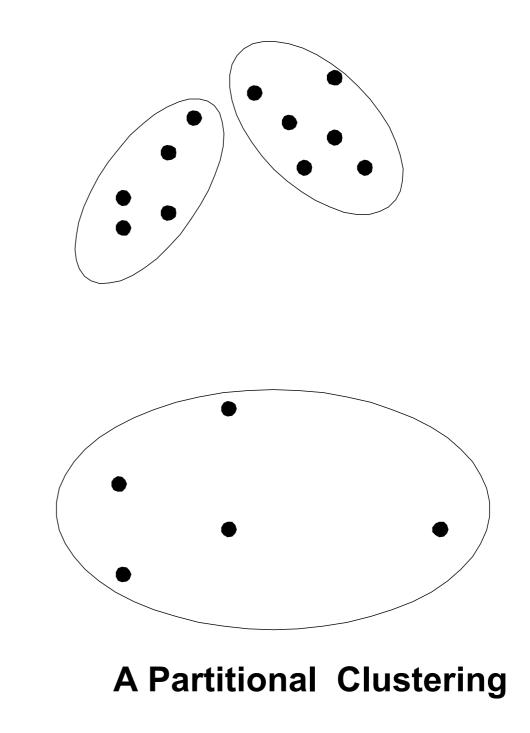
• a set of nested clusters organized in a tree

Hierarchical clustering



Partitional clustering





Partitional algorithms

partition the n objects into k clusters

- each object belongs to exactly one cluster
- the number of clusters k is given in advance

The k-means problem

- consider set X={x₁,...,x_n} of n points in R^d
- assume that the number k is given
- problem:
 - find k points c₁,...,c_k (named centers or means) so that the cost

$$\sum_{i=1}^{n} \min_{j} \left\{ L_2^2(x_i, c_j) \right\} = \sum_{i=1}^{n} \min_{j} ||x_i - c_j||_2^2$$

is minimized

The k-means problem

- consider set $X = \{x_1, ..., x_n\}$ of n points in \mathbb{R}^d
- assume that the number k is given
- problem:
 - find k points c₁,...,c_k (named centers or means)
 - and partition X into {X₁,...,X_k} by assigning each point x_i in X to its nearest cluster center,
 - so that the cost

$$\sum_{i=1}^{n} \min_{j} ||x_i - c_j||_2^2 = \sum_{j=1}^{k} \sum_{x \in X_j} ||x - c_j||_2^2$$

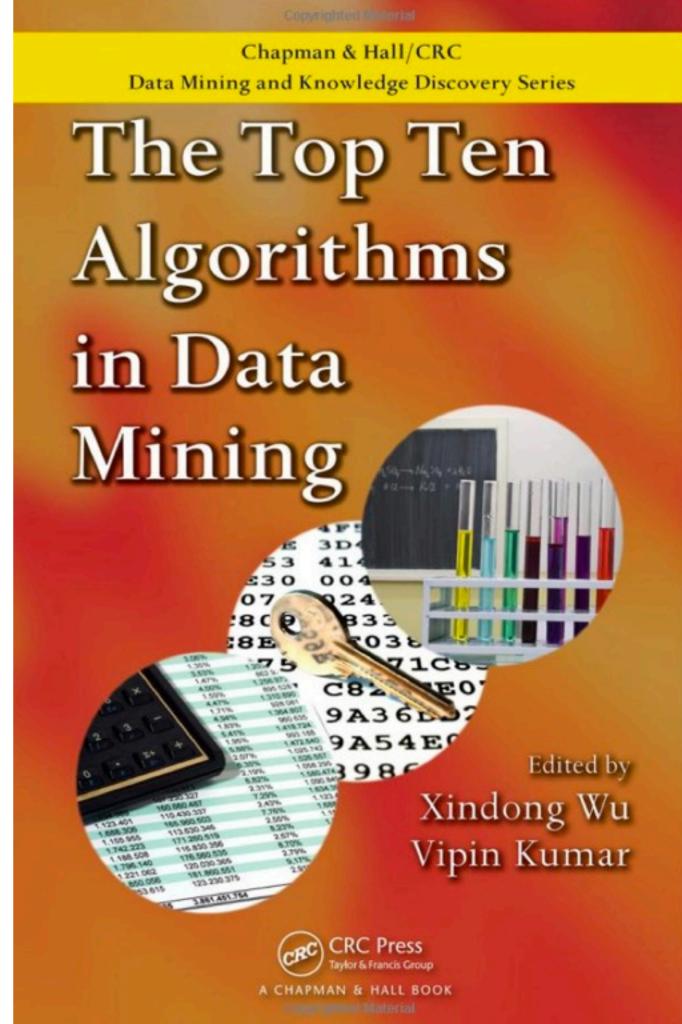
is minimized

The k-means problem

- k=1 and k=n are easy special cases (why?)
- an NP-hard problem if the dimension of the data is at least 2 (d≥2)
 - for d≥2, finding the optimal solution in polynomial time is infeasible
- for d=1 the problem is solvable in polynomial time
- in practice, a simple iterative algorithm works quite well

The k-means algorithm

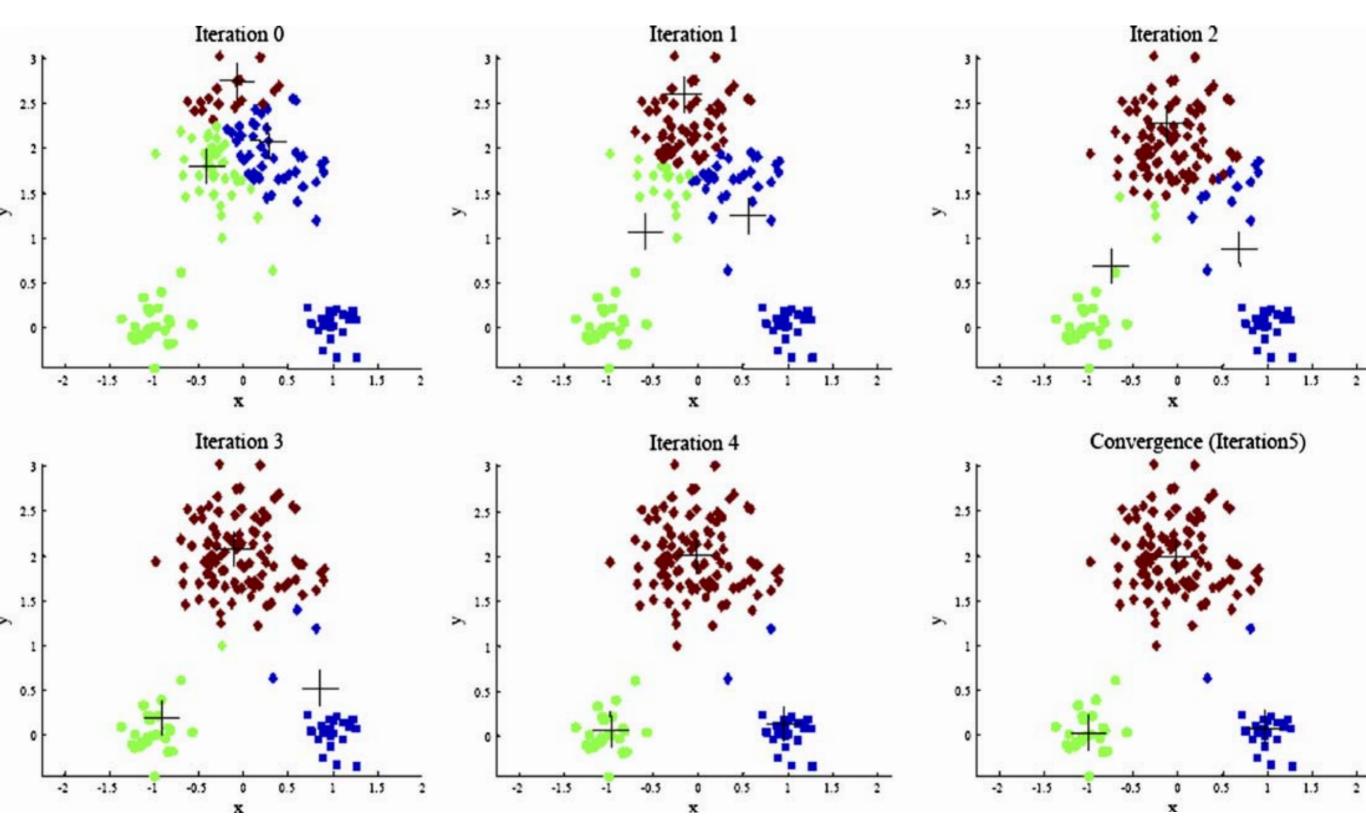
- voted among the top-10 algorithms in data mining
- one way of solving the kmeans problem



The k-means algorithm

- 1.randomly (or with another method) pick k cluster centers {c₁,...,c_k}
- 2.for each j, set the cluster X_j to be the set of points in X that are the closest to center c_j
- 3.for each j let c_j be the center of cluster X_j (mean of the vectors in X_j)
- 4.repeat (go to step 2) until convergence

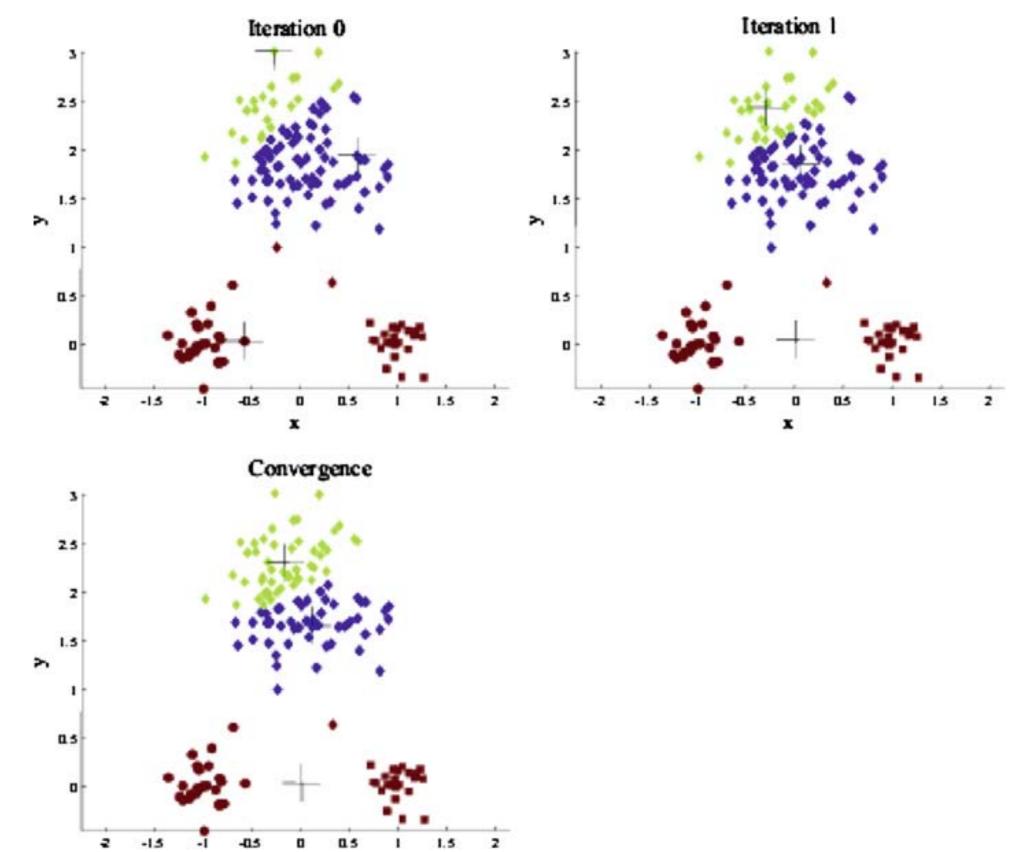
Sample execution



Properties of the k-means algorithm

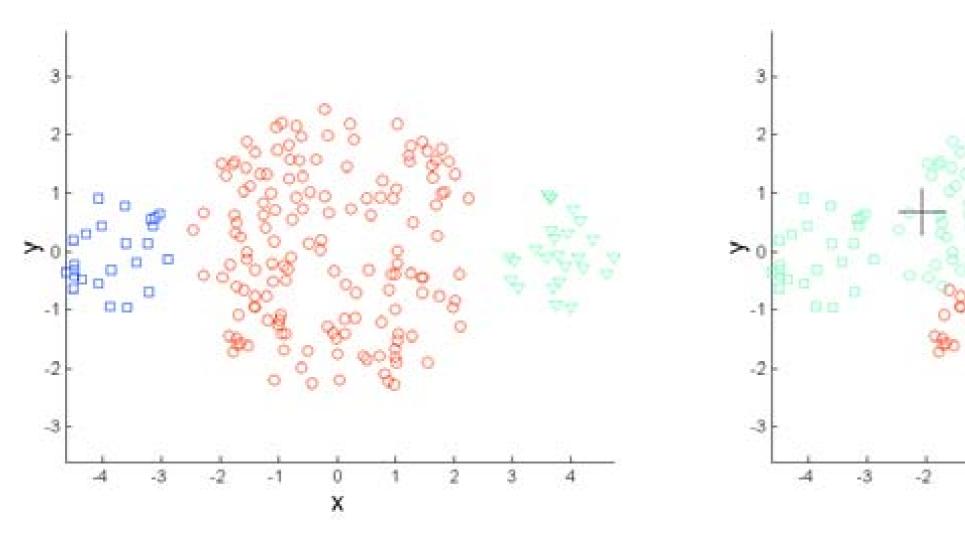
- finds a local optimum
- often converges quickly but not always
- the choice of initial points can have large influence in the result

Effects of bad initialization



x

Limitations of k-means: different sizes



Original Points

K-means (3 Clusters)

0

Х

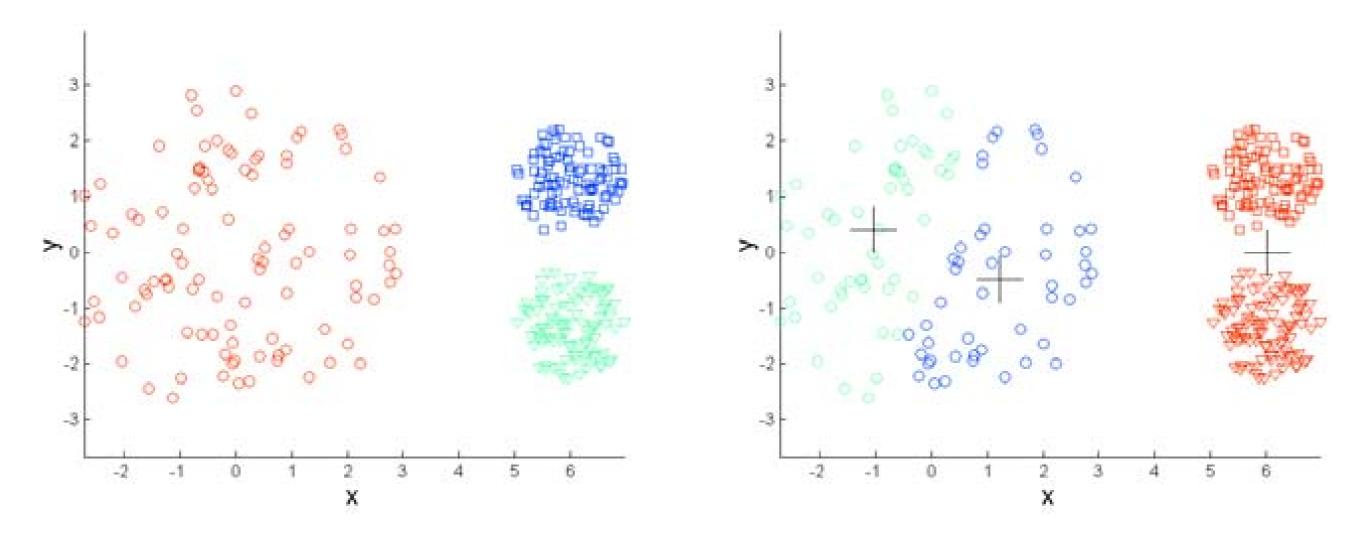
2

3

4

-1

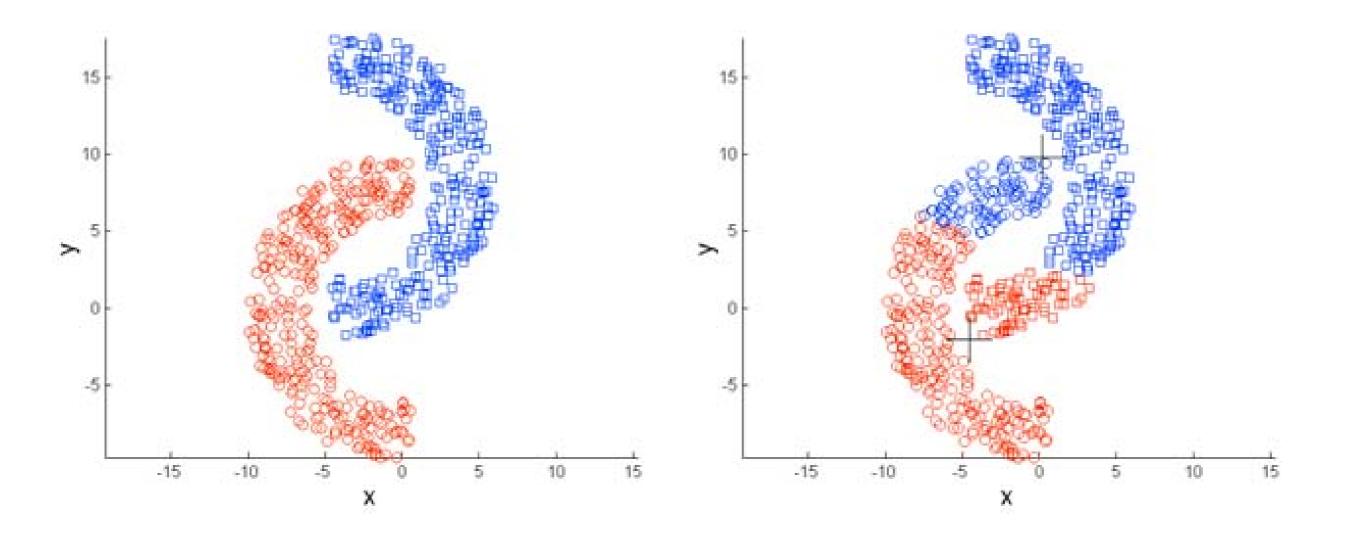
Limitations of k-means: different density



Original Points

K-means (3 Clusters)

Limitations of k-means: non-spherical shapes



Original Points

K-means (2 Clusters)

Discussion on the k-means algorithm

- finds a local optimum
- often converges quickly

but not always

- the choice of initial points can have large influence in the result
- tends to find spherical clusters
- outliers can cause a problem
- different densities may cause a problem

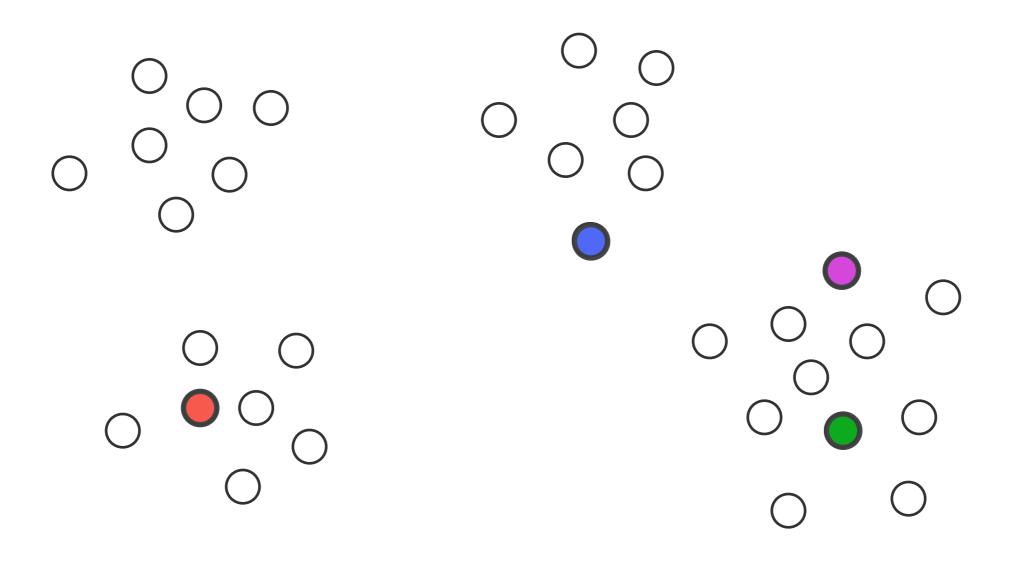
Initialization

- random initialization
- random, but repeat many times and take the best solution
 - helps, but solution can still be bad
- pick points that are distant to each other
 - k-means++
 - provable guarantees

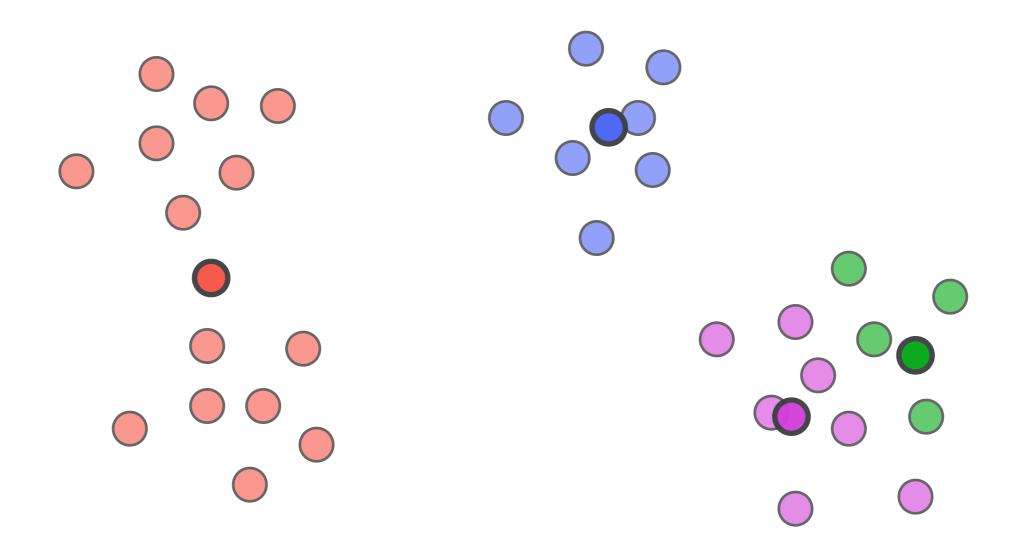
k-means++

David Arthur and Sergei Vassilvitskii k-means++: The advantages of careful seeding SODA 2007

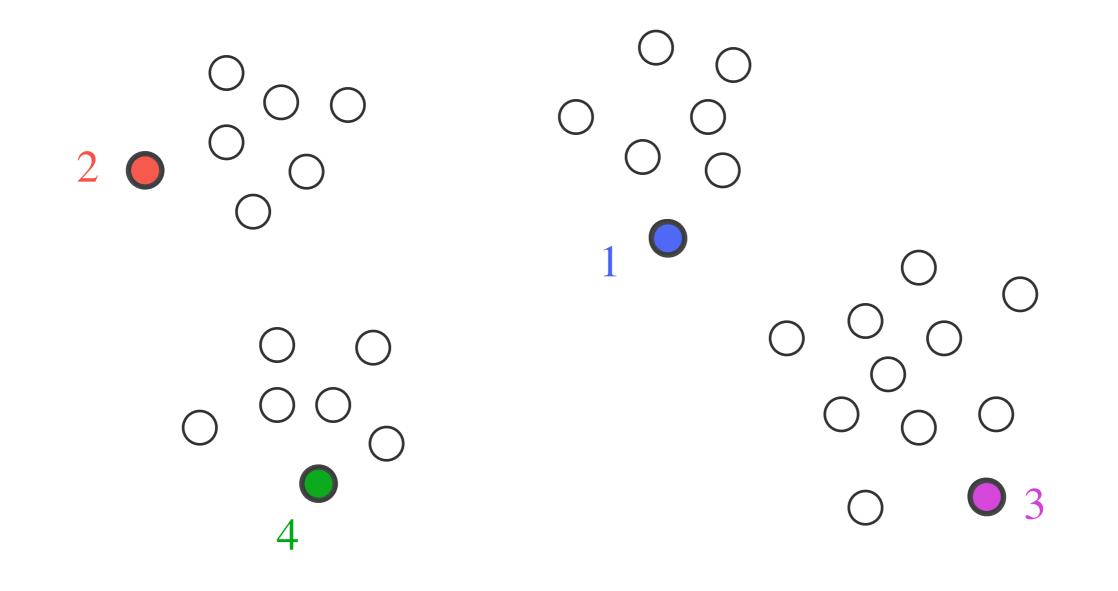
k-means algorithm: random initialization



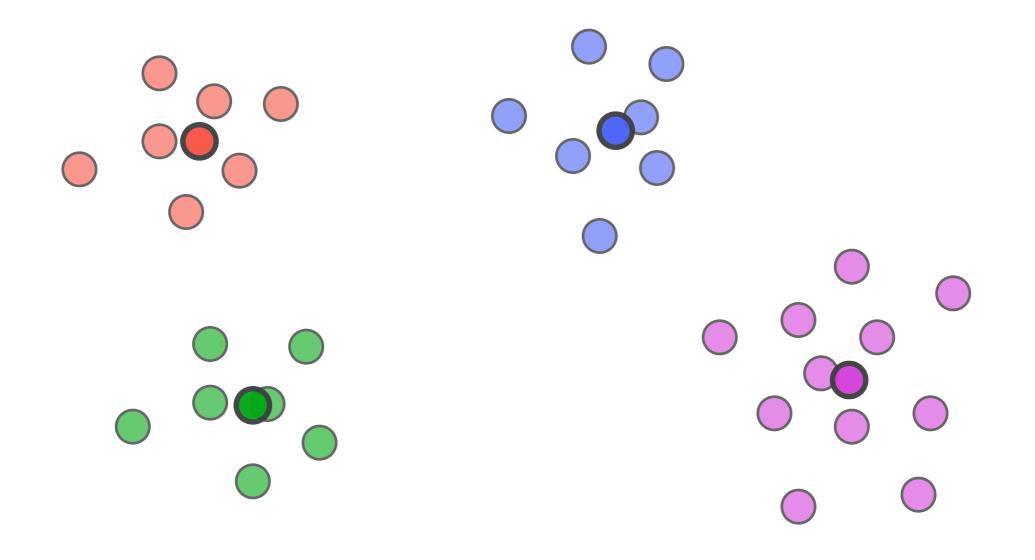
k-means algorithm: random initialization



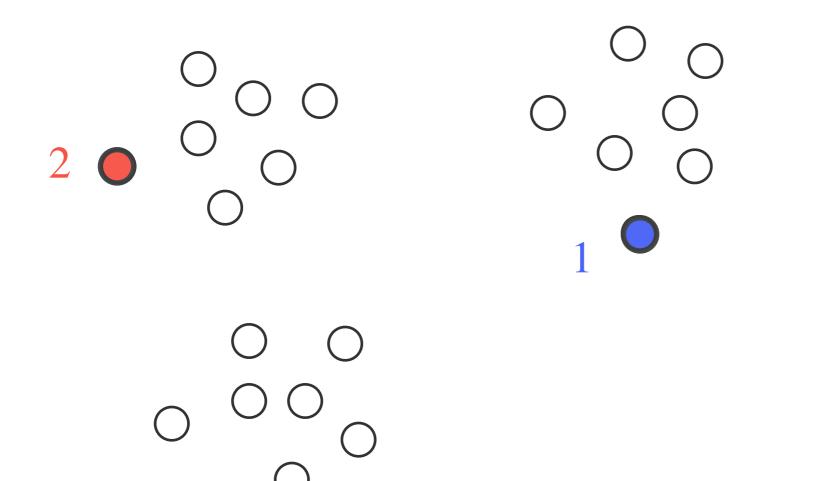
k-means algorithm: initialization with further-first traversal



k-means algorithm: initialization with further-first traversal

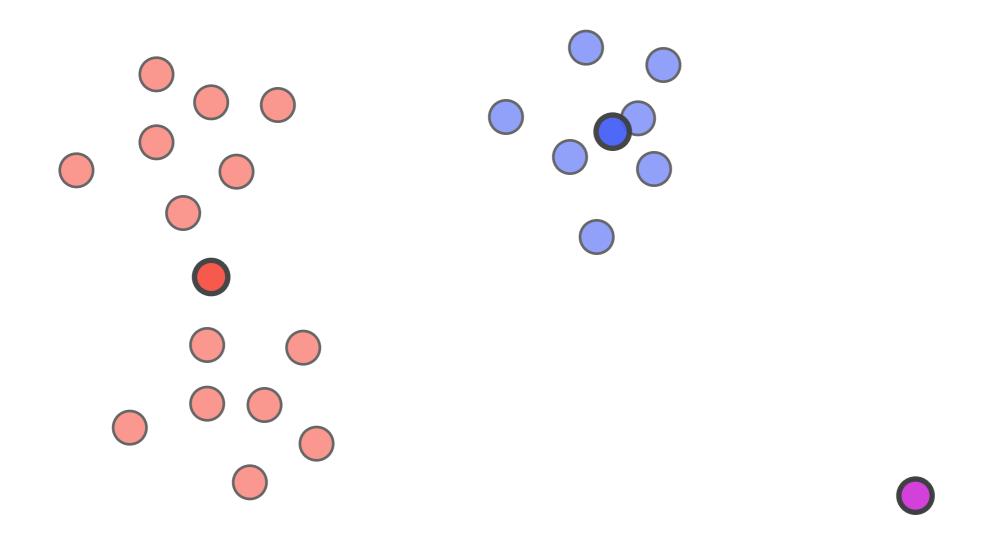


but... sensitive to outliers

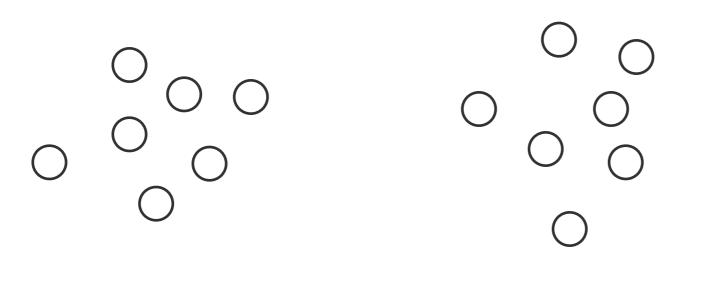




but... sensitive to outliers



Here random may work well



С

k-means++ algorithm

- interpolate between the two methods
- let D(x) be the distance between x and the nearest center selected so far
- choose next center with probability proportional to

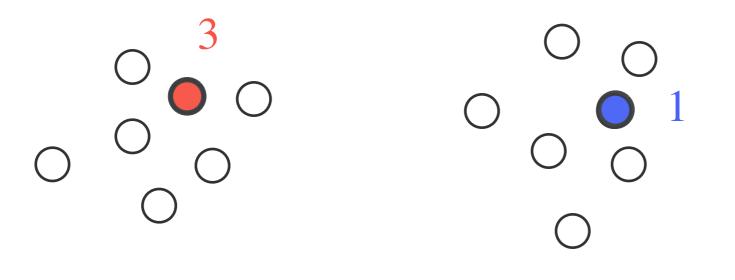
 $(D(x))^{a} = D^{a}(x)$

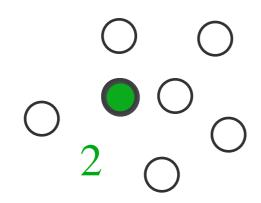
- + a = 0 random initialization
- * $a = \infty$ furthest-first traversal
- + a = 2 k-means++

k-means++ algorithm

- initialization phase:
 - choose the first center uniformly at random
 - choose next center with probability proportional to $D^2(x)$
- iteration phase:
 - iterate as in the k-means algorithm until convergence

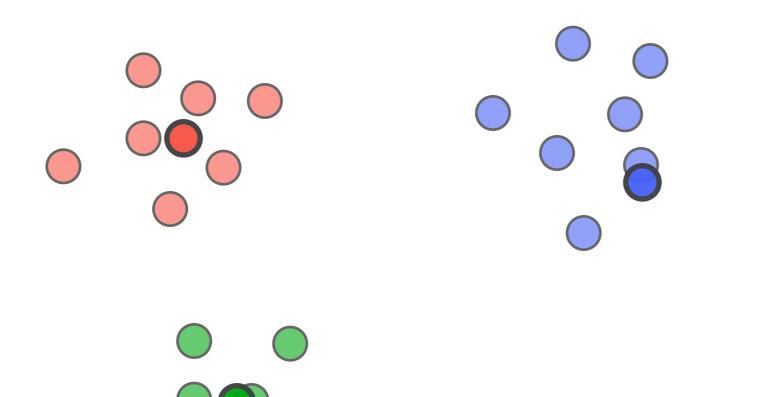
k-means++ initialization





С

k-means++ result





k-means++ provable guarantee

Theorem:

k-means++ is O(logk) approximate in expectation

k-means++ provable guarantee

- approximation guarantee comes just from the first iteration (initialization)
- subsequent iterations can only improve cost

k-means++ analysis

- consider optimal clustering C^{*}
- assume that k-means++ selects a center from a new optimal cluster

• then

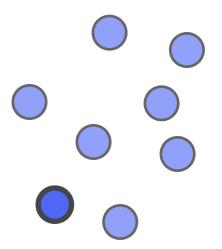
- k-means++ is 8-approximate in expectation
- intuition: if no points from a cluster are picked, then it probably does not contribute much to the overall error
- an inductive proof shows that the algorithm is O(logk) approximate

k-means++ proof : first cluster

- fix an optimal clustering C^{*}
- first center is selected uniformly at random
- bound the total error of the points in the optimal cluster of the first center

k-means++ proof : first cluster

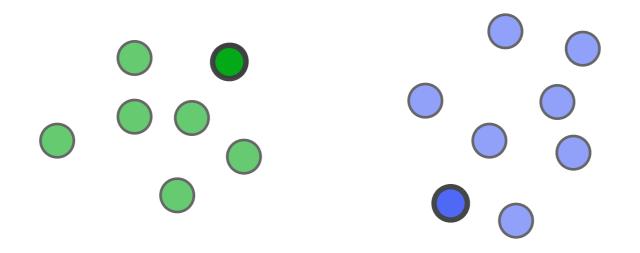
- let A be the first cluster
- each point a₀ ∈ A is equally likely to be selected as center



expected error:

$$E[\phi(A)] = \sum_{a_0 \in A} \frac{1}{|A|} \sum_{a \in A} ||a - a_0||^2$$
$$= 2 \sum_{a \in A} ||a - \bar{A}||^2 = 2\phi^*(A)$$

k-means++ proof : other clusters

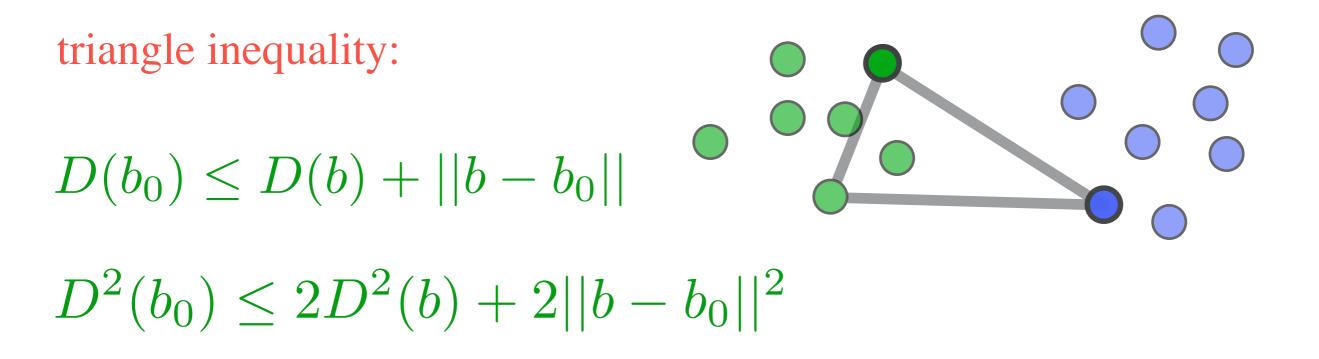


- suppose next center is selected from a new cluster in the optimal clustering C^{*}
- bound the total error of that cluster

k-means++ proof : other clusters

let B be the second cluster and b₀ the center selected

$$E[\phi(B)] = \sum_{b_0 \in B} \frac{D^2(b_0)}{\sum_{b \in B} D^2(b)} \sum_{b \in B} \min\{D(b), ||b - b_0||^2\}$$



k-means++ proof : other clusters

$$D^{2}(b_{0}) \leq 2D^{2}(b) + 2||b - b_{0}||^{2}$$

average over all points b in B

 \blacklozenge

$$D^{2}(b_{0}) \leq \frac{2}{|B|} \sum_{b \in B} D^{2}(b) + \frac{2}{|B|} \sum_{b \in B} ||b - b_{0}||^{2}$$
recall

$$E[\phi(B)] = \sum_{b_0 \in B} \frac{D^2(b_0)}{\sum_{b \in B} D^2(b)} \sum_{b \in B} \min\{D(b), ||b - b_0||^2\}$$
$$\leq 4 \sum_{b \in B} \frac{1}{|B|} \sum_{b_0 \in B} ||b - b_0||^2 = 4 \sum_{b \in B} 2||b - \bar{B}||^2 = 8\phi^*(B)$$

k-means++ analysis

- if that k-means++ selects a center from a new optimal cluster
- then
 - k-means++ is 8-approximate in expectation
- an inductive proof shows that the algorithm is O(logk) approximate

Lesson learned

no reason to use k-means and not k-means++

• k-means++ :

- easy to implement
- provable guarantee
- works well in practice

The k-median problem

- consider set $X = \{x_1, \dots, x_n\}$ of n points in \mathbb{R}^d
- assume that the number k is given
- problem:
 - find k points c₁,...,c_k (named medians)
 - and partition X into {X₁,...,X_k} by assigning each point x_i in X to its nearest cluster median,
 - so that the cost

$$\sum_{i=1}^{n} \min_{j} ||x_{i} - c_{j}||_{2} = \sum_{j=1}^{k} \sum_{x \in X_{j}} ||x - c_{j}||_{2}$$
 is minimized

the k-medoids algorithm

or PAM (partitioning around medoids)

- 1.randomly (or with another method) choose k medoids {c₁,...,c_k} from the original dataset X
- 2.assign the remaining n-k points in X to their closest medoid c_j
- 3.for each cluster, replace each medoid by a point in the cluster that improves the cost
- 4.repeat (go to step 2) until convergence

Discussion on the k-medoids algorithm

- very similar to the k-means algorithm
- same advantages and disadvantages
- how about efficiency?

The Local-kMedian algorithm

- Pick a random set of k cluster centers $S = \{c_1, \ldots, c_k\}$
- \mathbb{P}_S : partition induced by assigning each data point to its closest point in S
- Repeat
 - Find S' ``**similar**" to S
 - If $kMedian-Cost(\mathbb{P}_{S'}) < kMedian-Cost(\mathbb{P}_S)$ then S = S'
- Until convergence
- Similar S' is find via swaps or p-swaps.

The Local-kMedian algorithm

• Proposition: If \mathbb{P}_S is the partition output by the LocalkMedian with single swaps and \mathbb{P}_{S^*} is the optimal partition for the k-Median problem, then

kMedian-Cost(\mathbb{P}_S) $\leq 5 \times \text{kMedian-Cost}(\mathbb{P}_{S^*})$

The Local-kMedian algorithm

• Proposition: If \mathbb{P}_S is the partition output by the LocalkMedian with p-swaps and \mathbb{P}_{S^*} is the optimal partition for the k-Median problem, then

kMedian-Cost(\mathbb{P}_S) $\leq (3 + 2/p) \times kMedian-Cost(\mathbb{P}_{S^*})$

The k-center problem

- consider set $X = \{x_1, ..., x_n\}$ of n points in \mathbb{R}^d
- assume that the number k is given
- problem:
 - find k points c₁,...,c_k (named centers)
 - and partition X into {X₁,...,X_k} by assigning each point x_i in X to its nearest cluster center,
 - so that the cost

is minimized
$$\max_{i=1}^{n} \min_{j=1}^{k} ||x_i - c_j||_2$$

Properties of the k-center problem

- NP-hard for dimension d≥2
- for d=1 the problem is solvable in polynomial time (how?)
- a simple combinatorial algorithm works well

The k-center problem

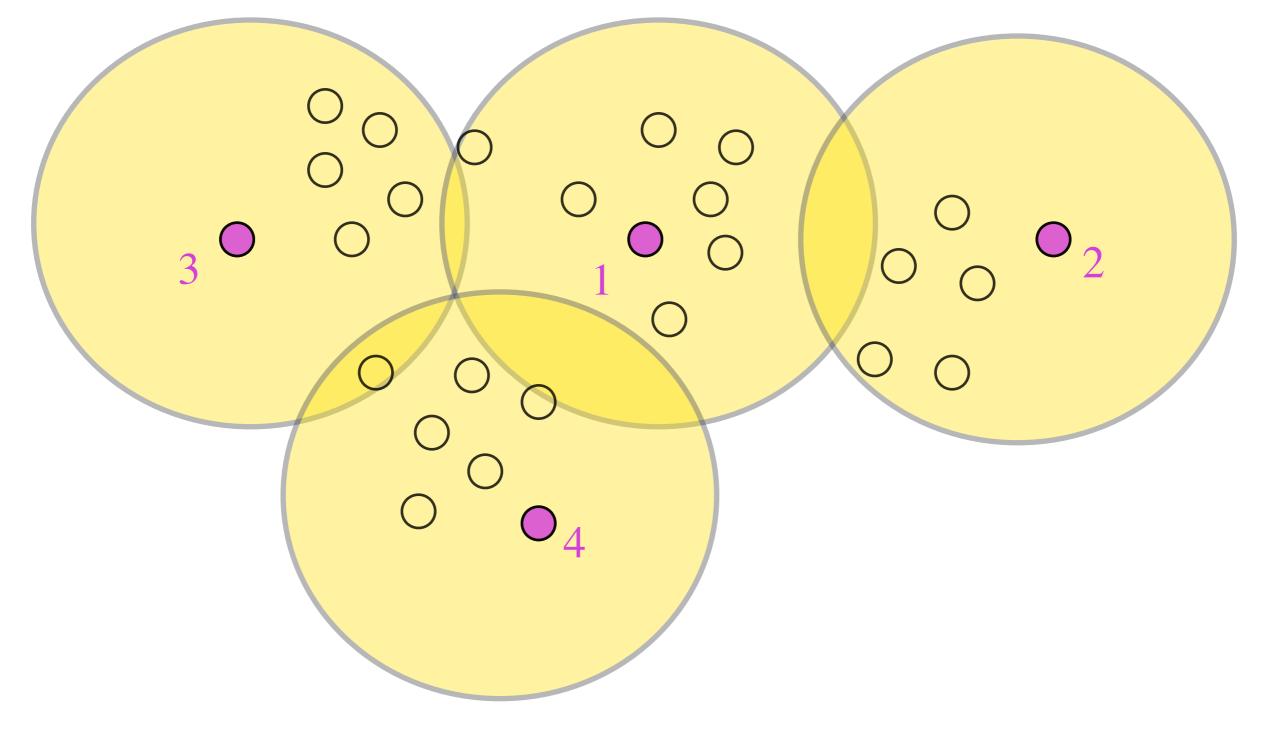
- consider set $X = \{x_1, ..., x_n\}$ of n points in \mathbb{R}^d
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 - find k points c₁,...,c_k (named centers)
 - and partition X into {X₁,...,X_k} by assigning each point x_i in X to its nearest cluster center,
 - so that the cost

is minimized
$$\max_{i=1}^{n} \min_{j=1}^{k} ||x_i - c_j||_2$$

Furthest-first traversal algorithm

- pick any data point and label it 1
- for i=2,...,k
 - find the unlabeled point that is furthest from {1,2,...,i-1}
 - // use d(x,S) = min y∈S d(x,y)
 - label that point i
- assign the remaining unlabeled data points to the closest labeled data point

Furthest-first traversal algorithm: example



Furthest-first traversal algorithm

 furthest-first traversal algorithm gives a factor 2 approximation

Furthest-first traversal algorithm

- pick any data point and label it 1
- for i=2,...,k
 - find the unlabeled point that is furthest from {1,2,...,i-1}
 - // use d(x,S) = min y∈S d(x,y)
 - label that point i
 - $p(i) = \operatorname{argmin}_{j \le i} d(i,j)$
 - R_i = d(i,p(i))
- assign the remaining unlabeled data points to the closest labeled data point

Analysis

• Claim 1: $R_1 \ge R_2 \ge ... \ge R_k$

• proof:

•
$$R_j = d(j,p(j))$$

= $d(j,\{1,2,...,j-1\})$
 $\leq d(j,\{1,2,...,i-1\}) // j > i$
 $\leq d(i,\{1,2,...,i-1\}) = R_i$

Analysis

- Claim 2:
 - let C be the clustering produced by the FFT algorithm
 - let R(C) be the cost of that clustering
 - then R(C) = R_{k+1}
- proof:
 - for any i>k we have :

 $d(i,\!\{1,\!2,\!\ldots,\!k\}) \leq d(k\!+\!1,\!\{1,\!2,\!\ldots,\!k\}) = R_{k\!+\!1}$

Analysis

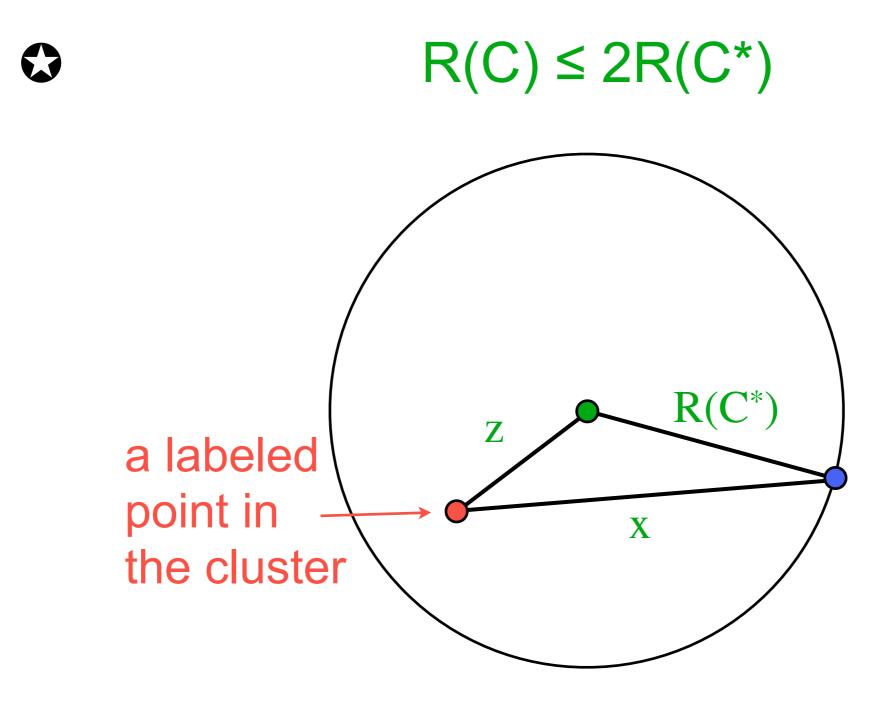
Theorem

- let C be the clustering produced by the FFT algorithm
- let C* be the optimal clustering
- then $R(C) \leq 2R(C^*)$
- proof:
 - let C_{1}^{*} ,..., C_{k}^{*} be the clusters of the optimal k-clustering
 - if these clusters contain points {1,...,k} then

 $R(C) \leq 2R(C^*)$

- otherwise suppose that one of these clusters contains two or more of the points in {1,...,k}
- these points are at distance at least R_k from each other
- this (optimal) cluster must have radius

 $\frac{1}{2} R_k \ge \frac{1}{2} R_{k+1} = \frac{1}{2} R(C)$



 $\mathsf{R}(\mathsf{C}) \le \mathsf{x} \le \mathsf{z} + \mathsf{R}(\mathsf{C}^*) \le 2\mathsf{R}(\mathsf{C}^*)$