

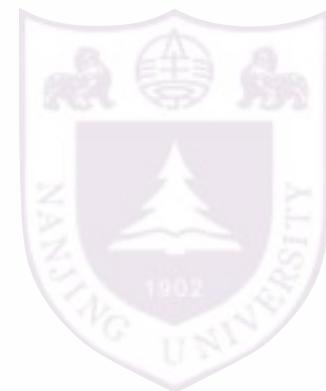
# Lecture 7: Unsupervised Learning

## density estimation and clustering

[http://cs.nju.edu.cn/yuy/course\\_dm13ms.ashx](http://cs.nju.edu.cn/yuy/course_dm13ms.ashx)

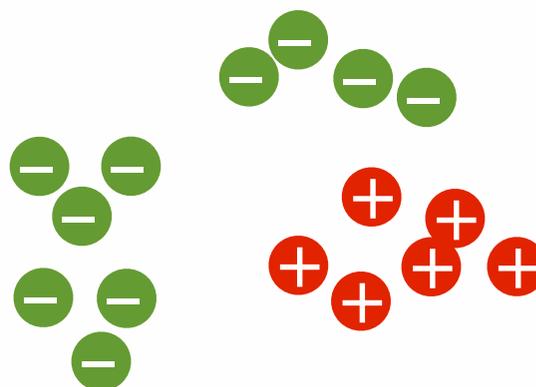


# Unsupervised learning



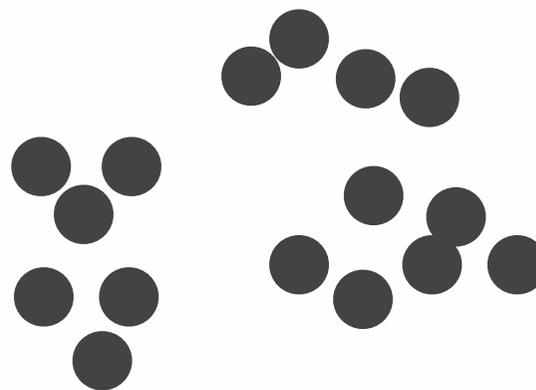
data for supervised learning

target: find a mapping  $h : \mathcal{X} \rightarrow \mathcal{Y}$



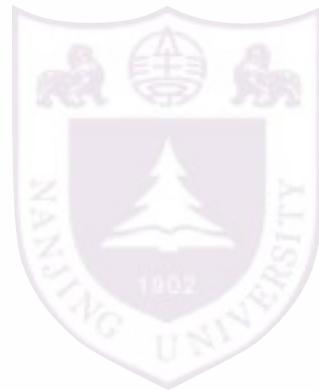
data for unsupervised learning:

target: find structures of the data



*what structures ?*

# Unsupervised learning

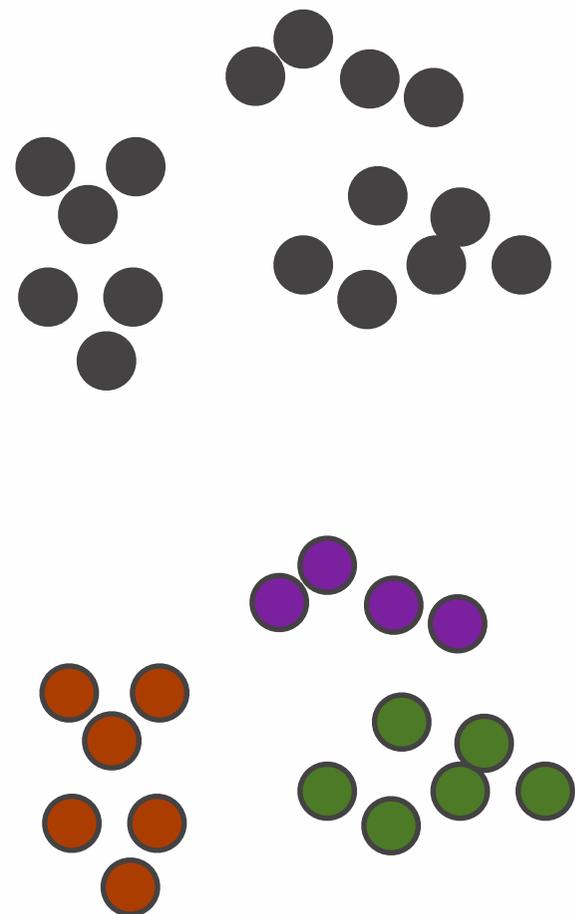


why unsupervised learning?

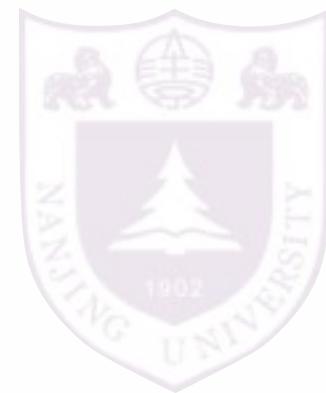
natural need of discovery  
of structures in data

act as a preprocessing step  
to help supervised learning

...



# Density estimation

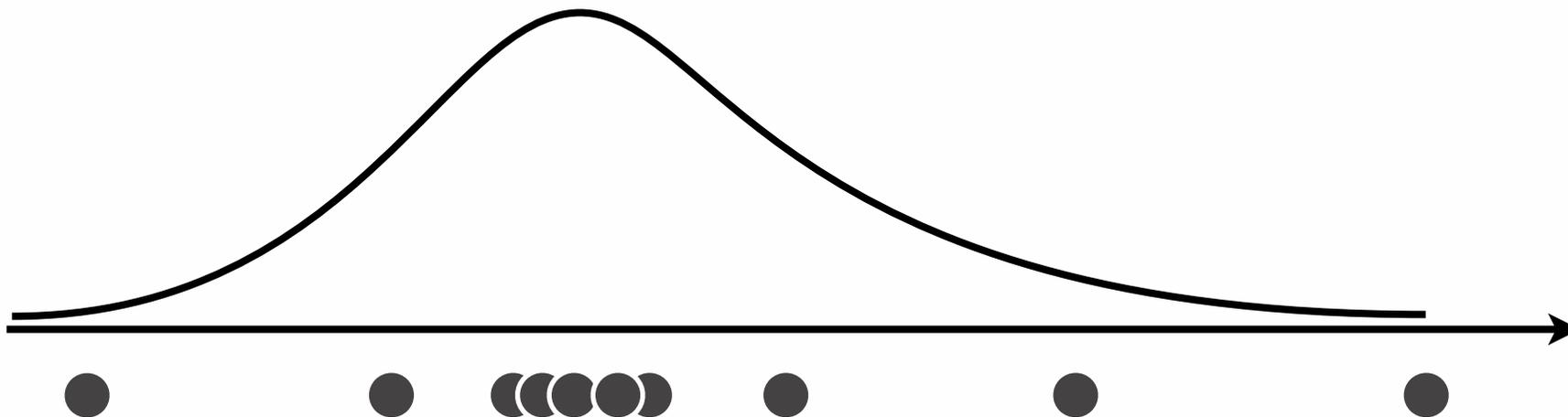


There exists a probability *density* function  $p$   
a data set  $D$  sampled i.i.d. from  $p$

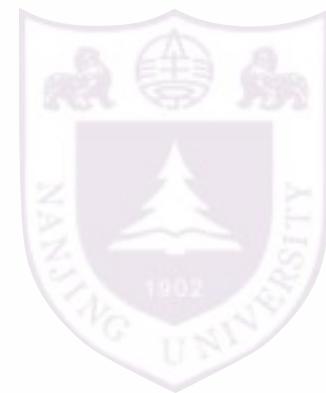
*how large is the density at  $x$ , i.e.,  $p(x)$ ?*

reconstruct  $p$  from  $D$

estimate the density of any instance



# Parametric methods



Assume the family of the density function,  
estimate the parameters

Normal distribution/Gaussian model:

$$p(\mathbf{x}) = \frac{1}{(2\pi)^{k/2} |\Sigma|^{1/2}} e^{-\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu})^\top \Sigma^{-1} (\mathbf{x} - \boldsymbol{\mu})}$$

Estimation:

$\boldsymbol{\mu}$  is data mean

$\Sigma$  is data covariance matrix

$$\Sigma = \frac{1}{n-1} \sum_{i=1}^n (\mathbf{x}_i - \bar{\mathbf{x}})(\mathbf{x}_i - \bar{\mathbf{x}})^\top$$

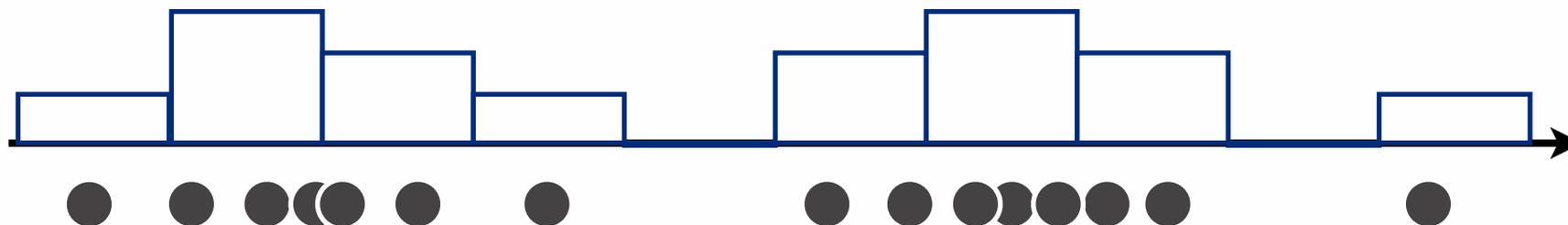
# Nonparametric methods



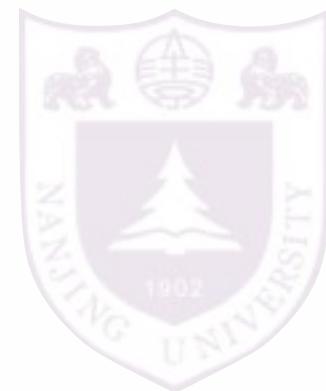
## Histogram estimator

divide the input space into bins  
count the frequency of instances in each bin

$$p(x) = \frac{\# \text{ instances in bin}(x)}{m \times \text{bin-width}}$$



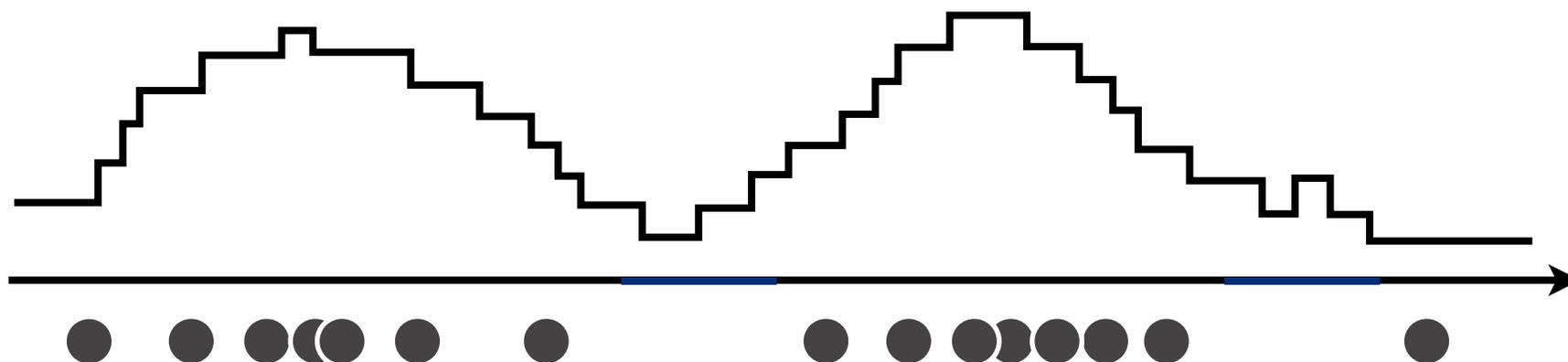
# Nonparametric methods



## Naive estimator

for each position, count instances in the neighbor range

$$p(x) = \frac{\# \text{ instances in } [x - h, x + h]}{m \times 2h}$$



# Nonparametric methods

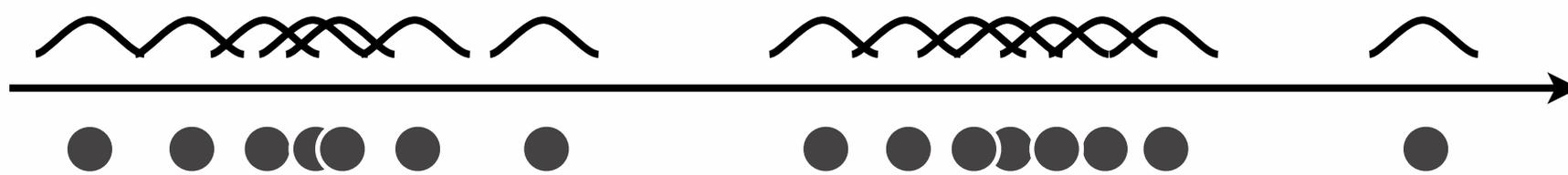


Kernel estimator/Parzen window

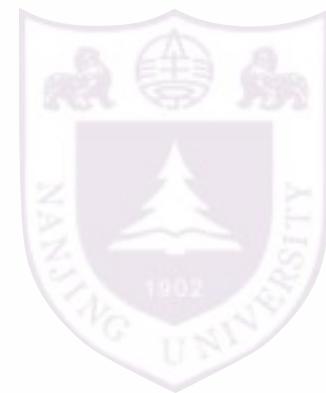
for each position, the influence of an instance decreases with the distance

$$p(x) = \frac{1}{mh} \sum_{i=1}^m K\left(\frac{x - x_i}{h}\right)$$

Gaussian kernel:  $K(\Delta) = \frac{1}{\sqrt{2\pi}} e^{-\Delta^2/2}$



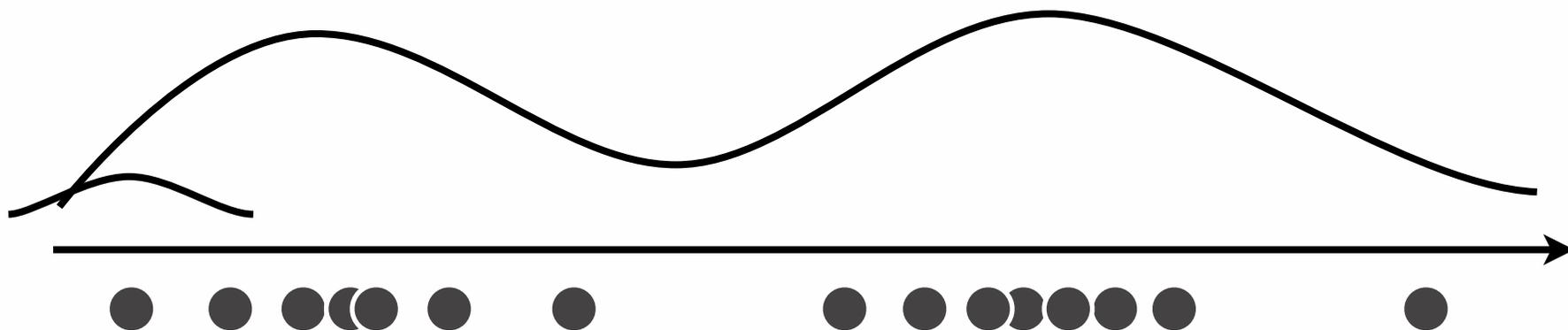
# Nonparametric methods



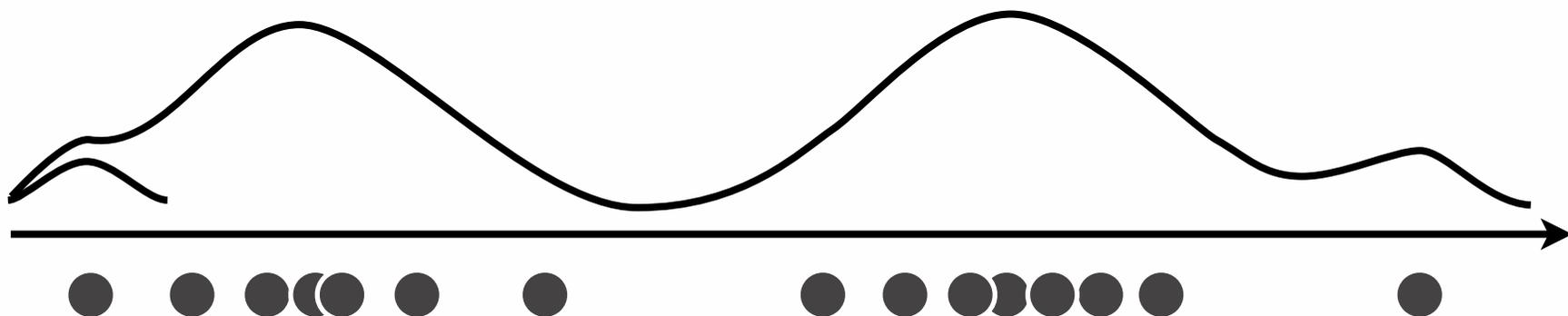
$$p(x) = \frac{1}{mh} \sum_{i=1}^m K\left(\frac{x - x_i}{h}\right)$$

$$\text{Gaussian kernel: } K(\Delta) = \frac{1}{\sqrt{2\pi}} e^{-\Delta^2/2}$$

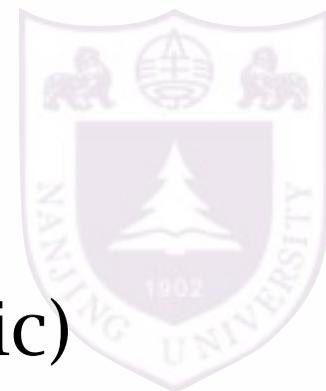
larger  $h$



smaller  $h$



# Nonparametric methods



random partition based method (non-metric)

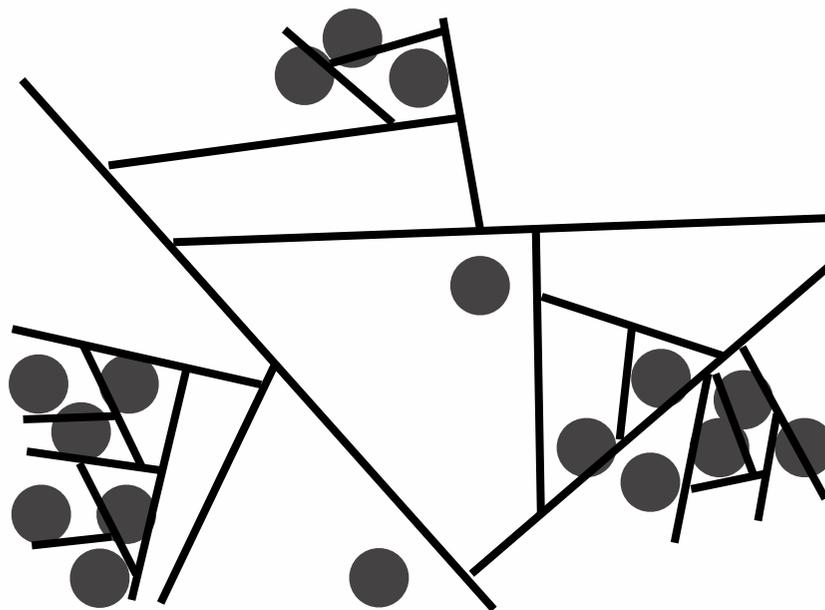
instance in low density region is easily separated

1. grow a full complete random oblique decision tree

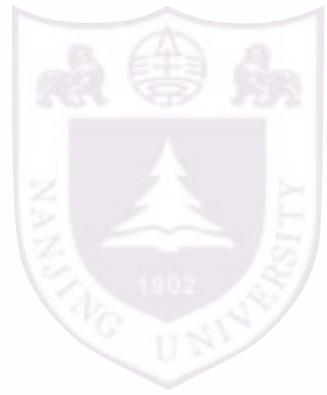
2. the leave depth implies the density

3. build and average many trees to smooth

(normalization is needed)



# Clustering



Clustering is to find clusters in the data

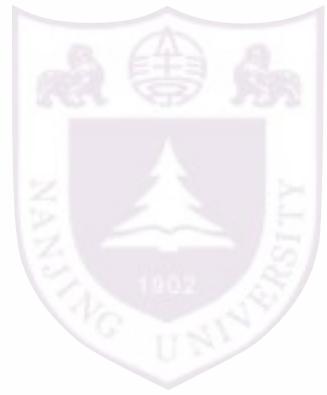


Unfortunately, there is no clear definition of what should be in a cluster



*the subjectivity of clustering*

# Clustering



hierarchical methods

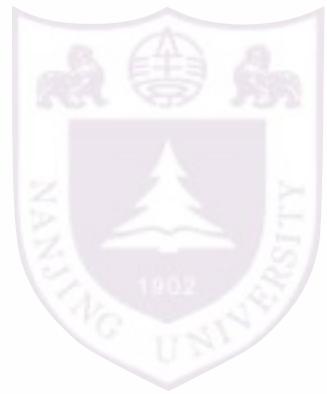
density-based methods

centroid-based methods

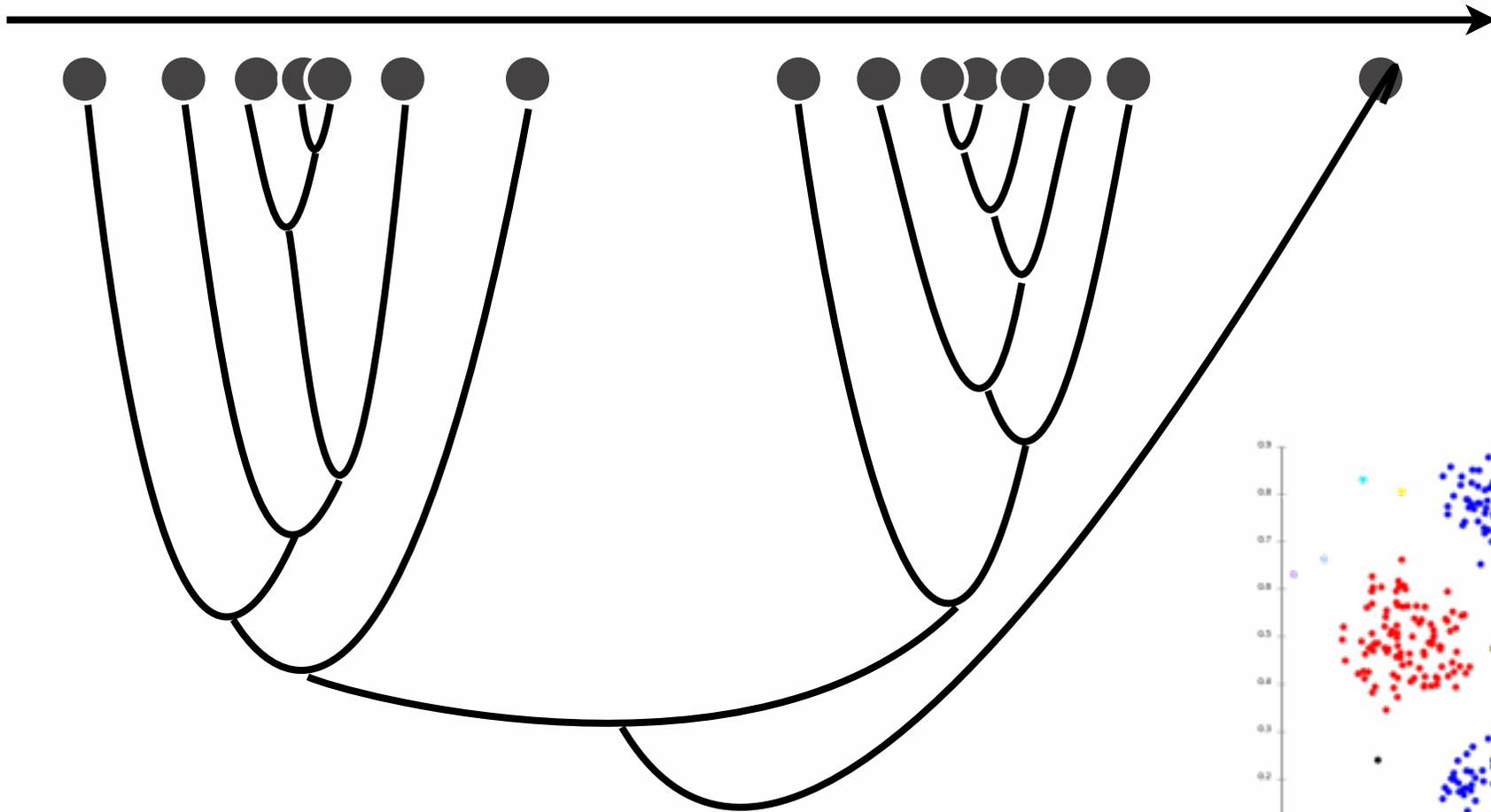
model-based methods

...

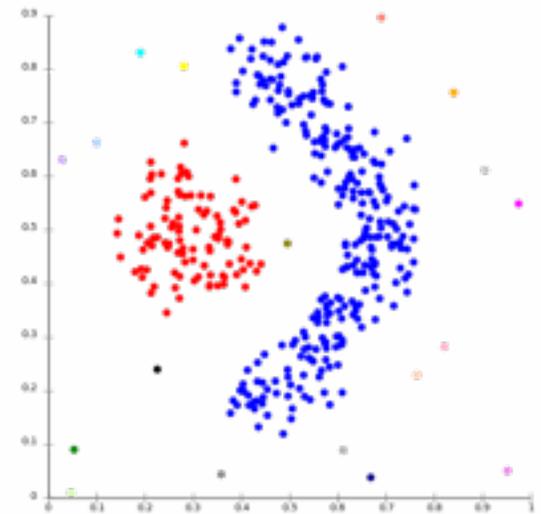
# Hierarchical methods



bottom-up: single-link clustering

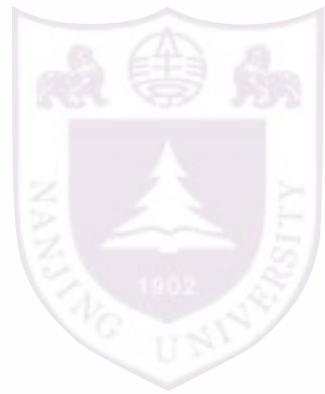


stop at a minimum distance threshold  
e.g. average distance

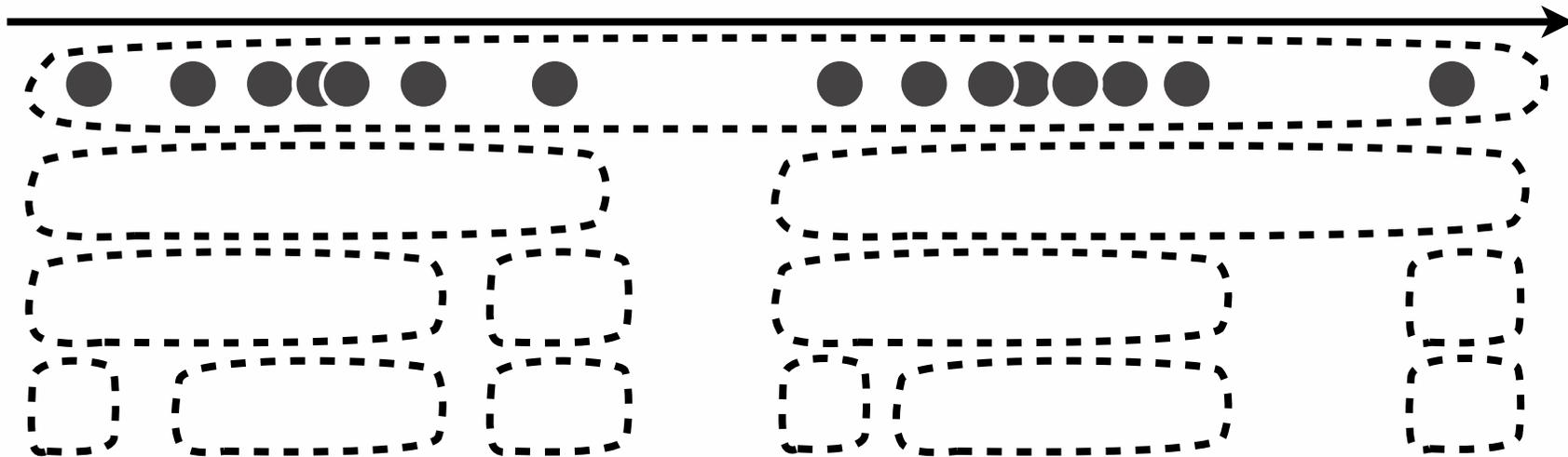


[from wikipedia]

# Hierarchical methods



top-down: divisive clustering



separate data into two groups by maximizing the inter-group distance

expensive in each level



# Density-based methods

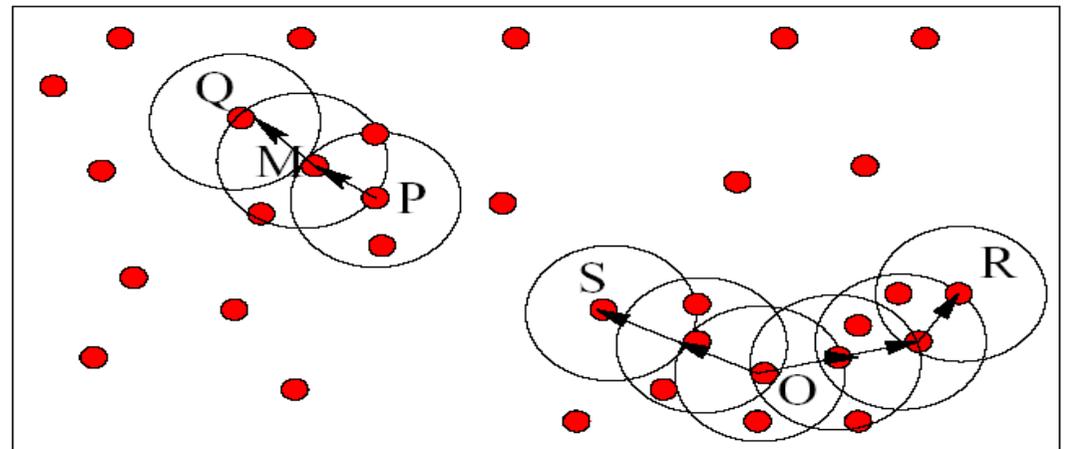
## DBSCAN

focus on dense instances, clustering by connectivity

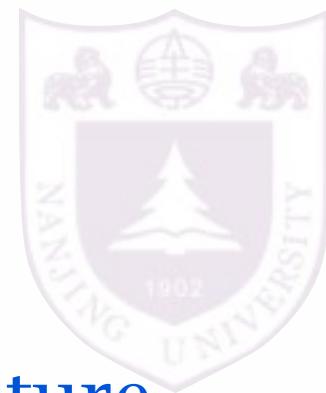
key concepts:

- an object  $P$  whose  $\varepsilon$ -neighborhood containing no less than  $MinPts$  number of objects is a **core object** with respect to  $\varepsilon$  and  $MinPts$
- an object  $M$  is **directly density-reachable** from object  $P$  with respect to  $\varepsilon$  and  $MinPts$  if  $M$  is within the  $\varepsilon$ -neighborhood of  $P$  which contains at least a minimum number of points,  $MinPts$
- an object  $Q$  is **density-reachable** from object  $P$  with respect to  $\varepsilon$  and  $MinPts$  if there is a chain of objects  $p_1, \dots, p_n, p_1 = P$  and  $p_n = Q, p_{i+1}$  is directly density-reachable from  $p_i$  with respect to  $\varepsilon$  and  $MinPts$
- an object  $S$  is **density-connected** to object  $R$  with respect to  $\varepsilon$  and  $MinPts$  if there is an object  $O$  such that both  $S$  and  $R$  are density-reachable from  $O$  with respect to  $\varepsilon$  and  $MinPts$

strictly not a clustering algorithm, leaving instances unclustered



# Density-based methods



## OPTICS

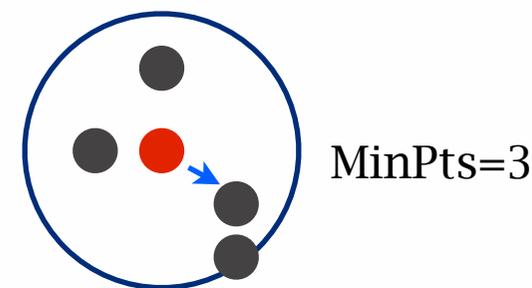
order instances to identify the cluster structure

for each core object, calculate **core-distance** to be the distance to the *MinPts*-th nearest instance

for instance  $p$ , calculate **reachability-distance** to a core object to be

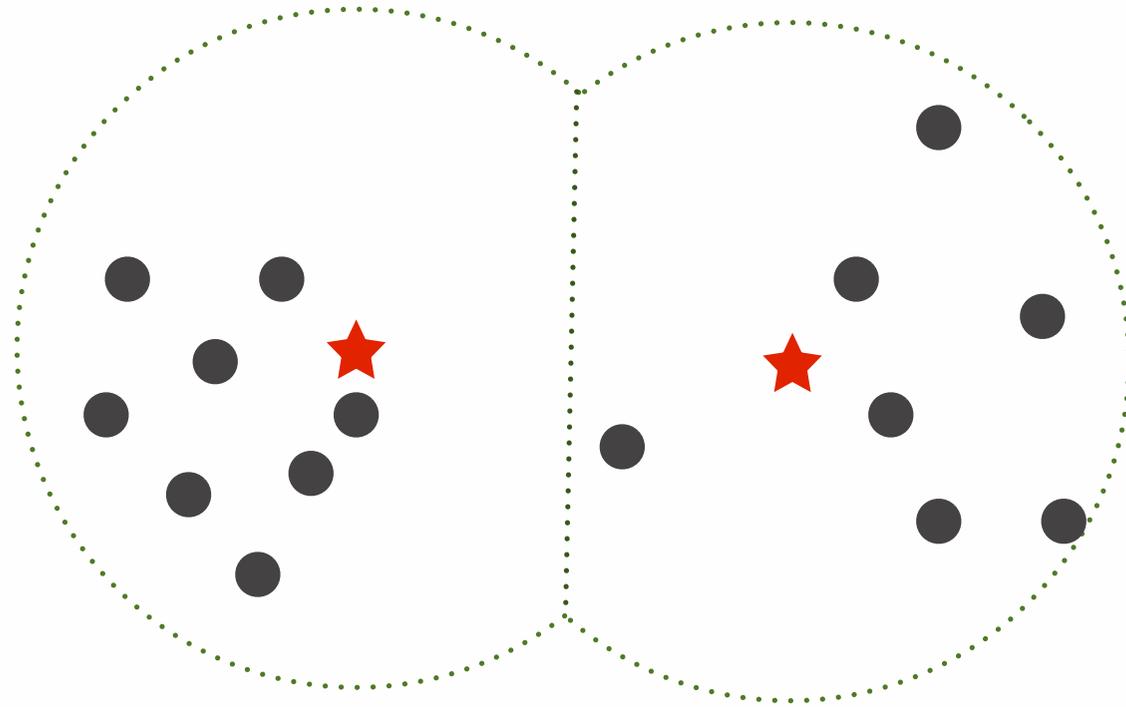
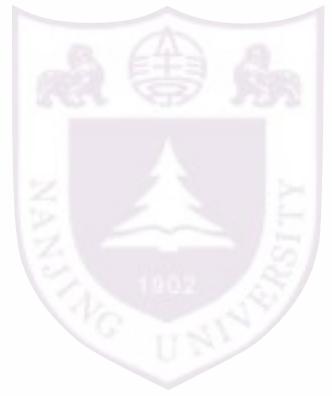
$$\max\{\text{core-distance}(o), \text{distance}(o,p)\}$$

similar to DBSCAN, but adjust the scanning order so that closer instances are ordered closer



[Ankerst et al., SIGMOD99]

# Centroid-based methods



# Centroid-based methods



## $k$ -means

Step1: randomly generate  $k$  centers

Step2: for each instance, assign it to the cluster whose center is the nearest to the instance

Step3: compute the means of the cluster and regard them as the centers

Step4: if there is no change, exit. otherwise go to Step2

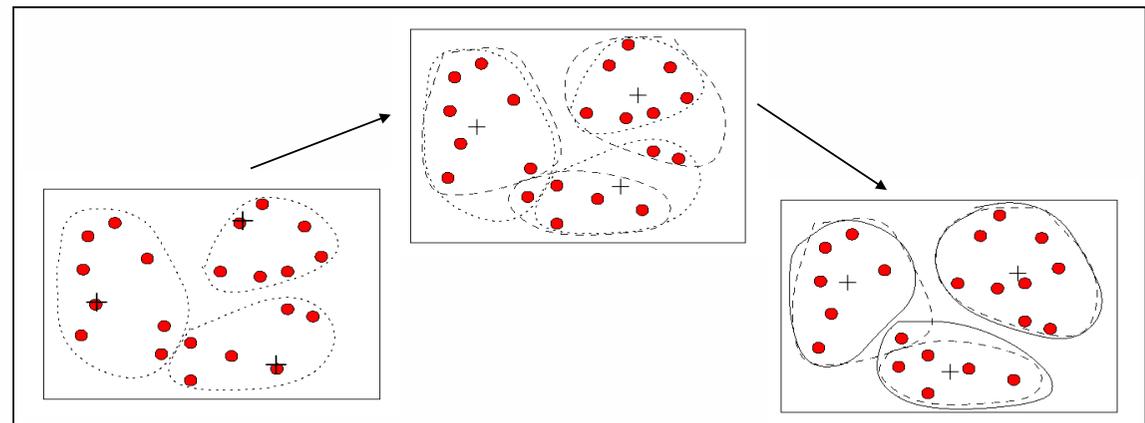
fix centers, update clusters

fix clusters, update centers

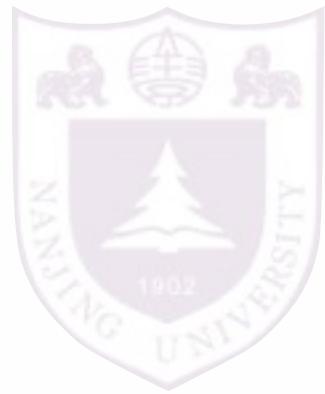
objective:

$$\arg \min_{\mathcal{S}} \sum_{i=1}^k \sum_{\mathbf{x} \in S_i} \|\mathbf{x} - \mu_i\|^2$$

converge to local optimal



# Centroid-based methods



## $k$ -medoids

Step1: randomly select  $k$  objects as the centers of the clusters

Step2: for each remaining object, assign it to the cluster whose center is the nearest to the object

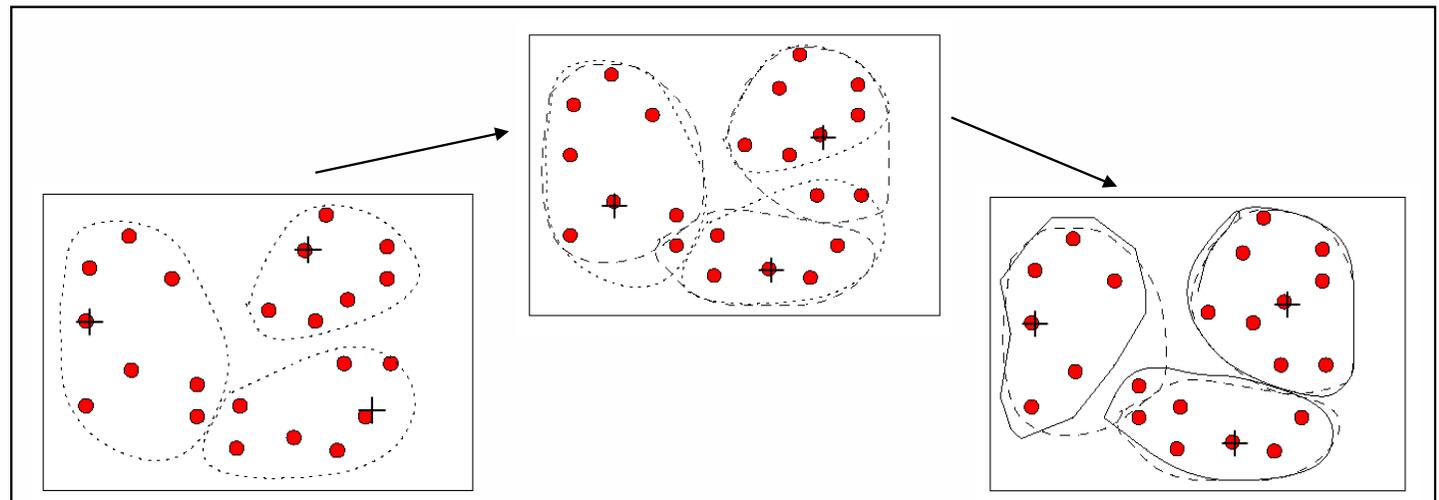
fix centers, update clusters

Step3: compute the means of the cluster, and assign the instance nearest to the mean as the centers

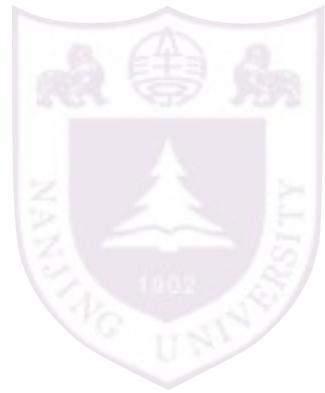
fix clusters, update centers

Step4: if there is no improvement, exit. otherwise go to Step2

$$\arg \min_{\mathcal{S}} \sum_{i=1}^k \sum_{\mathbf{x} \in S_i} \|\mathbf{x} - \mu_i\|^2$$



# Centroid-based methods

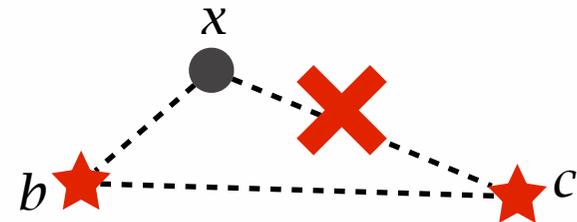


accelerate *k*-means [Elkan, ICML03]

in the original *k*-means algorithm the later iterations do not utilize earlier information

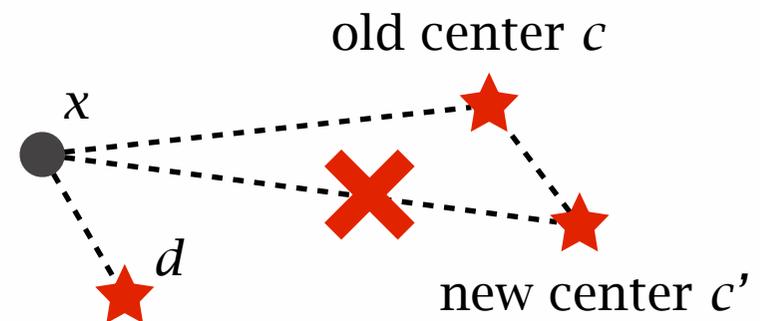
**Lemma 1:** Let  $x$  be a point and let  $b$  and  $c$  be centers. If  $d(b, c) \geq 2d(x, b)$  then  $d(x, c) \geq d(x, b)$ .

when  $d(x, b)$  is calculated, we don't need to calculate  $d(x, c)$  in order to know  $x$  is closer to  $b$  than  $c$ .



**Lemma 2:** Let  $x$  be a point and let  $b$  and  $c$  be centers. Then  $d(x, c) \geq \max\{0, d(x, b) - d(b, c)\}$ .

when we know  $d(x, c)$  and that the new center moves a distance  $\Delta$ , we know  $d(x, c')$  is at least  $d(x, c) - \Delta$  (or 0) without calculate the exact distance.



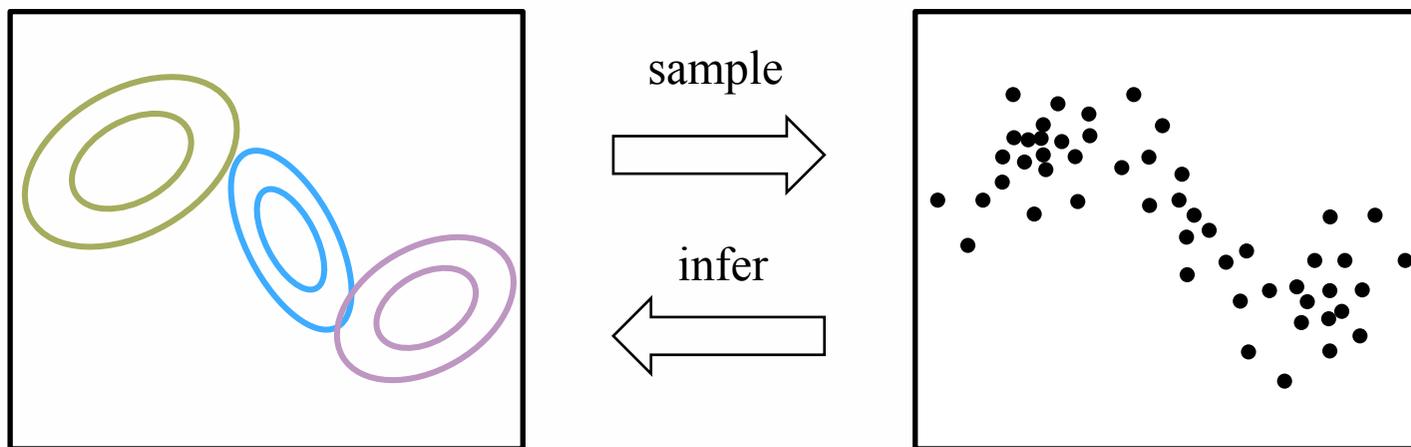
# Model-based methods



## Gaussian-mixture model

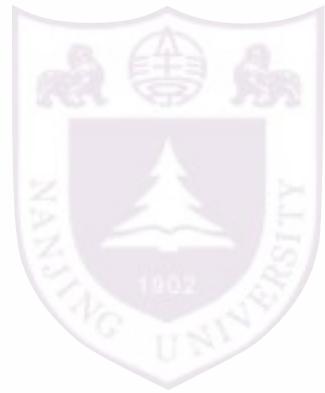
A perspective of dealing with unlabeled data is to imagine how the data is *generated*

assume that the data were generated from multiple Gaussian components



Clustering: To infer the Gaussian components from data

# Model-based methods



Gaussian models:

**Gaussian model has two parameters:**  $\mathcal{N}(\boldsymbol{\mu}, \boldsymbol{\Sigma})$

**Density function:**

$$p(\boldsymbol{x}) = \frac{1}{(2\pi)^{k/2} |\boldsymbol{\Sigma}|^{1/2}} e^{-\frac{1}{2}(\boldsymbol{x} - \boldsymbol{\mu})^\top \boldsymbol{\Sigma}^{-1}(\boldsymbol{x} - \boldsymbol{\mu})}$$

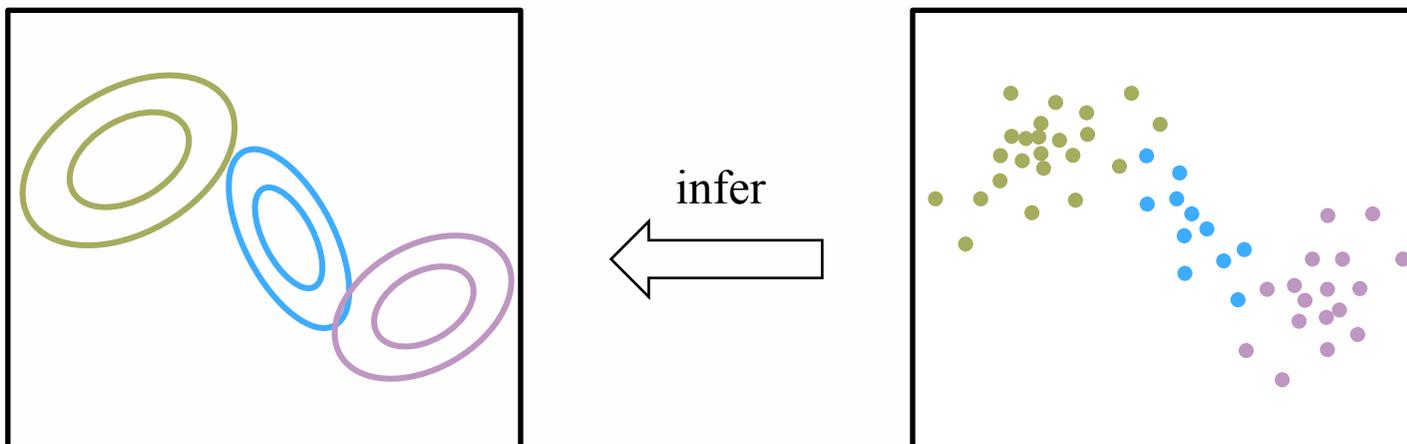
**Log-likelihood function:**

$$\ln p(\boldsymbol{x} | \boldsymbol{\mu}, \boldsymbol{\Sigma}) = -\frac{1}{2} \left( k \ln(2\pi) + \ln |\boldsymbol{\Sigma}| + (\boldsymbol{x} - \boldsymbol{\mu})^\top \boldsymbol{\Sigma}^{-1}(\boldsymbol{x} - \boldsymbol{\mu}) \right)$$

# Model-based methods



When data clusters are known:



**We know that there are three Gaussians models**

**for each model, calculate its parameters by**

**maximizing the log-likelihood function:  $\sum_x \ln p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma})$**

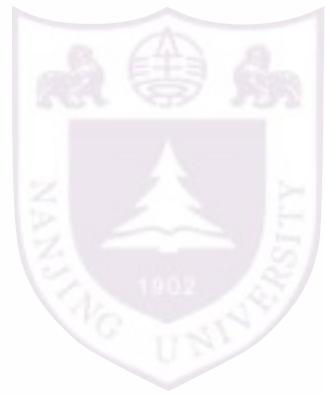
$$\begin{cases} \partial \sum_x \ln p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) / \partial \boldsymbol{\mu} = 0 \\ \partial \sum_x \ln p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) / \partial \boldsymbol{\Sigma} = 0 \end{cases}$$

$$N = \sum_x p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma})$$

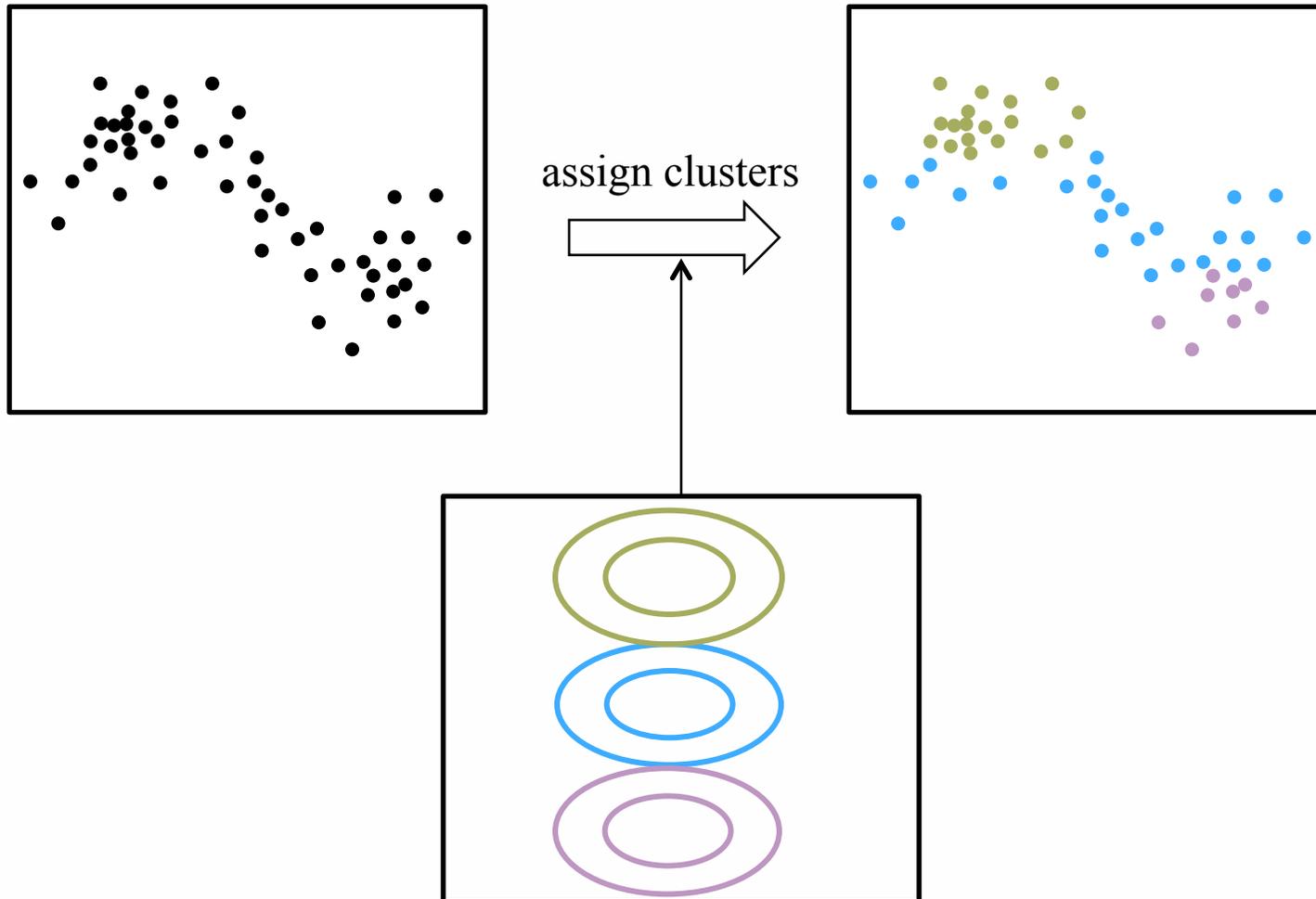
$$\boldsymbol{\mu} = \frac{1}{N} \sum_x p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) \cdot \mathbf{x} \quad \text{(data mean)}$$

$$\boldsymbol{\Sigma} = \frac{1}{N} \sum_x p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) (\mathbf{x} - \boldsymbol{\mu})(\mathbf{x} - \boldsymbol{\mu})^\top \quad \text{(data covariance)}$$

# Model-based methods



When data clusters are unknown:



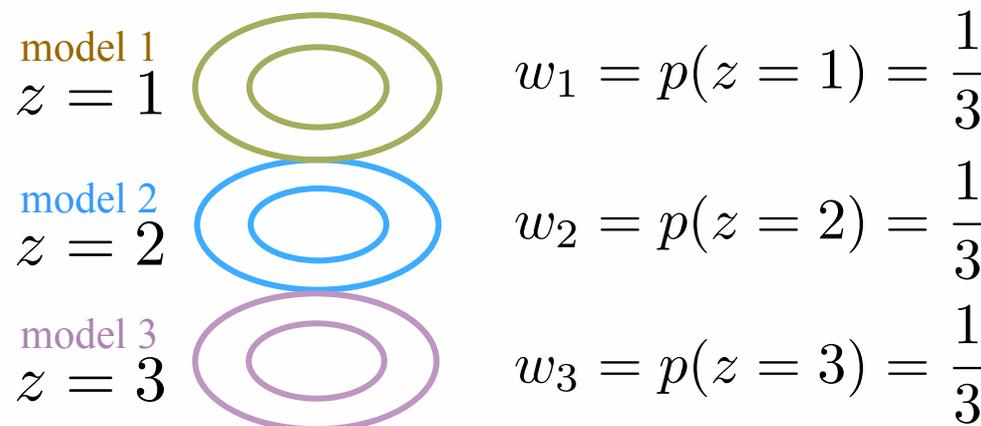
**Guess the model at first!**

# Model-based methods



How to assign clusters to data:

**Assume the models and their prior probabilities**



**Bayes rule:**

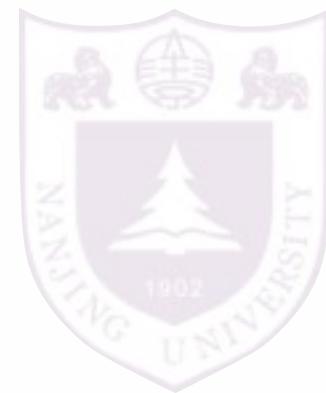
$$p(z | \mathbf{x}) = \frac{p(\mathbf{x} | z)p(z)}{p(\mathbf{x})}$$

density function  $\rightarrow$   $p(\mathbf{x} | z)$       prior probability  $\rightarrow$   $p(z)$

**Assign the cluster of the largest posterior probability**

$$c(\mathbf{x}) = \arg \max_{i=1,2,3} p(\mathbf{x} | \boldsymbol{\mu}_i, \boldsymbol{\Sigma}_i) \cdot w_i$$

# Model-based methods



## EM algorithm:

The original EM approach [Dempster et al, J Royal Statistical Society'77]

1. Initial guess of models (with equal prior probabilities)

$$(\boldsymbol{\mu}_1, \boldsymbol{\Sigma}_1, w_1 = \frac{1}{k}), \dots, (\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k, w_k = \frac{1}{k})$$

2. Assign clusters to data

$$c(\mathbf{x}) = \arg \max_{i=1, \dots, k} p(\mathbf{x} | \boldsymbol{\mu}_i, \boldsymbol{\Sigma}_i) \cdot w_i$$

**Expectation**

complete the data

3. Re-estimate model parameters from data

$$\boldsymbol{\mu}_i = \frac{1}{N_i} \sum_{\mathbf{x}} p(\mathbf{x} | \boldsymbol{\mu}_i, \boldsymbol{\Sigma}_i) \mathbf{x}$$

$$\boldsymbol{\Sigma}_i = \frac{1}{N_i} \sum_{\mathbf{x}} p(\mathbf{x} | \boldsymbol{\mu}_i, \boldsymbol{\Sigma}_i) (\mathbf{x} - \boldsymbol{\mu}_i)(\mathbf{x} - \boldsymbol{\mu}_i)^\top$$

$$w_i = N_i / N$$

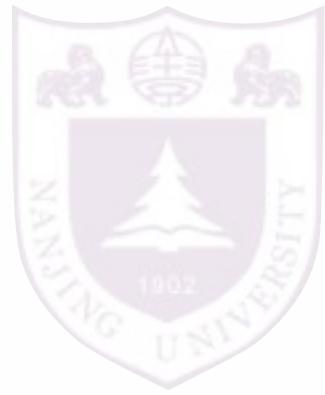
**Maximization**

complete the model

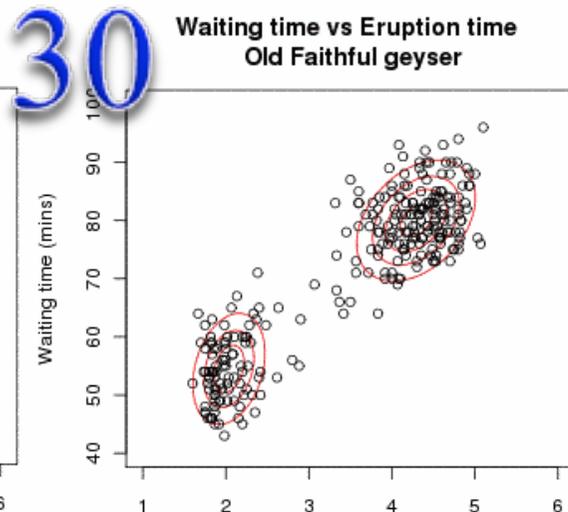
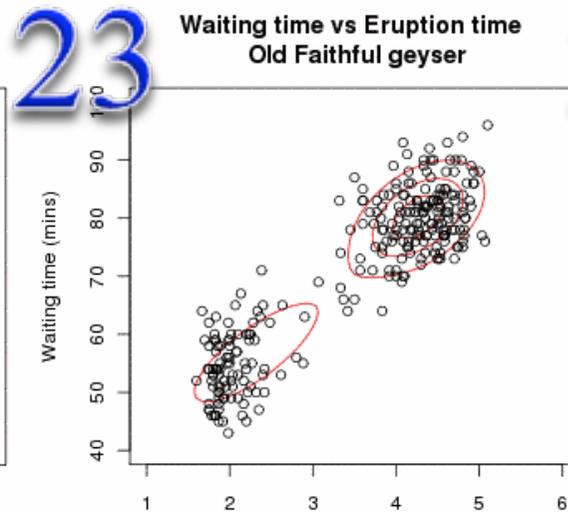
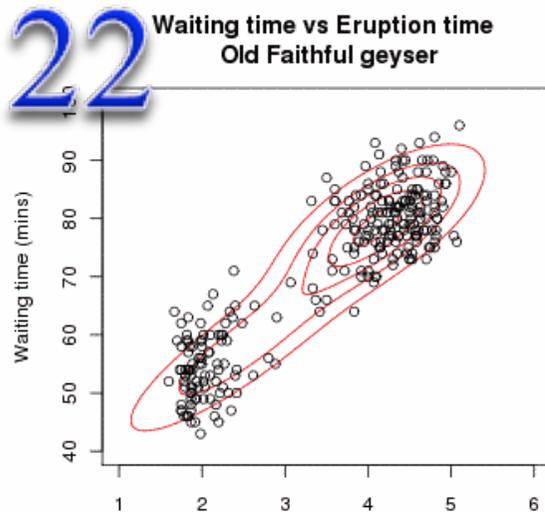
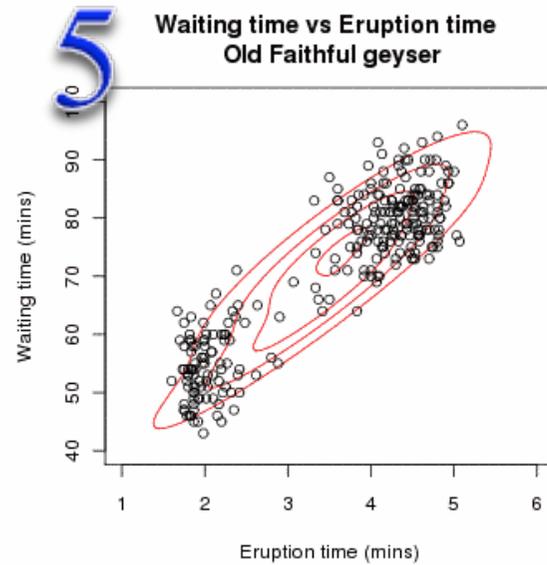
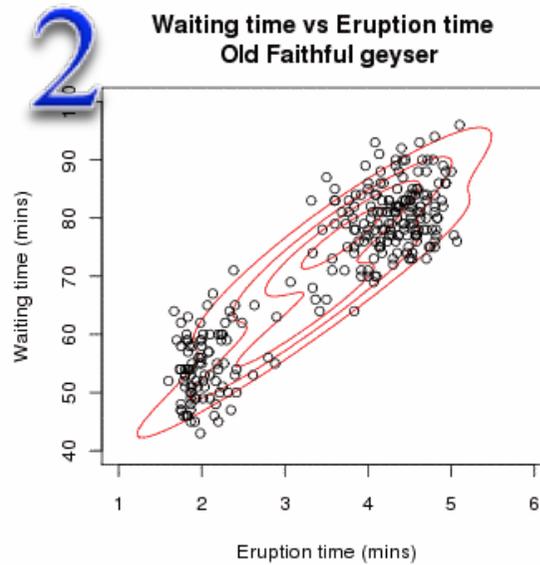
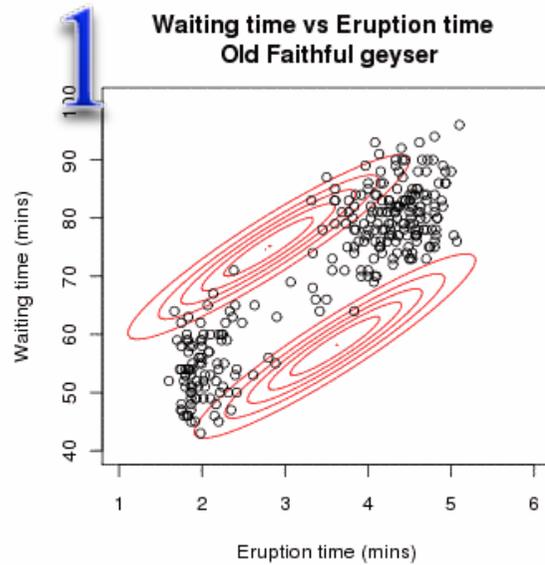
$$N_i = \sum_{\mathbf{x}} p(\mathbf{x} | \boldsymbol{\mu}_i, \boldsymbol{\Sigma}_i)$$

4. Go to 2 if not *converged*

# Model-based methods

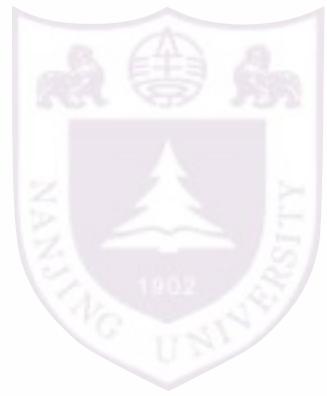


## GMM example:

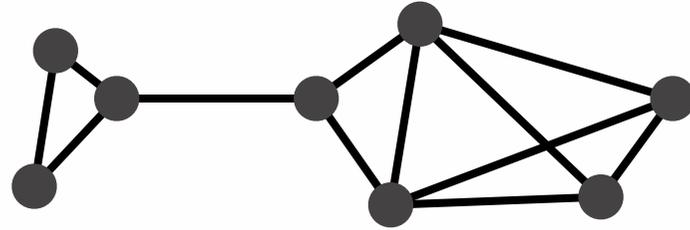


(from  
wikipedia)

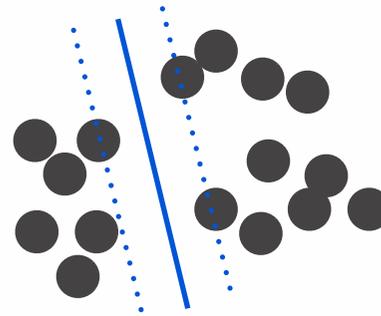
# Some other methods



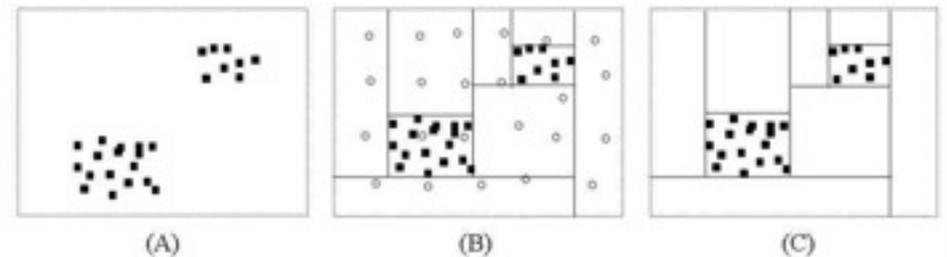
spectral clustering [Shi and Malik, PAMI00]:



maximum margin clustering [Xu *et al.*, NIPS05]:



decision tree-based clustering [Liu *et al.*, FADM05]:



# Determine the number of clusters



Rule of thumb

$$k = \sqrt{n/2}$$

Cross-validation

leave a subset of data as *test data*

try different number of clusters to maximize the performance on the test data

Using density based method

Use OPTICS to find the number of clusters, then run *k*-means

...

# 习题



使用核密度估计(kernel estimator)方法是否会受到距离函数的影响?

k-means 聚类算法的停止条件是什么?

k-means 聚类算法的优化目标是什么?

阐述k-means聚类算法的执行过程和关键步骤。