Syllabus



Fri. 21.10.	(1)	0. Introduction
		A. Supervised Learning: Linear Models & Fundamentals
Fri. 27.10.	(2)	A.1 Linear Regression
Fri. 3.11.	(3)	A.2 Linear Classification
Fri. 10.11.	(4)	A.3 Regularization
Fri. 17.11.	(5)	A.4 High-dimensional Data
		B. Supervised Learning: Nonlinear Models
Fri. 24.11.	(6)	B.1 Nearest-Neighbor Models
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Fri. 12.1.	(10)	B.5 A First Look at Bayesian and Markov Networks
		C. Unsupervised Learning
Fri. 19.1.	(11)	C.1 Clustering
Fri. 26.1.	(12)	C.2 Dimensionality Reduction
Fri. 2.2.	(13)	C.3 Frequent Pattern Mining

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Machine Learning

Outline



- 1. k-means & k-medoids
- 2. Gaussian Mixture Models
- 3. Hierarchical Cluster Analysis

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- 1. k-means & k-medoids
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Machine Learning 1. k-means & k-medoids



Partitions

Let X be a set. A set $P \subseteq \mathcal{P}(X)$ of subsets of X is called a partition of X if the subsets

1. are pairwise disjoint:

$$A \cap B = \emptyset$$
, $A, B \in P, A \neq B$

2. cover *X*:

$$\bigcup_{A} A = X, \text{ and }$$

 $A \in P$

3. do not contain the empty set: $\emptyset \notin P$.

Let $X := \{x_1, \dots, x_N\}$ be a finite set. A set $P := \{X_1, \dots, X_K\}$ of subsets $X_k \subseteq X$ is called a **partition of** X if the subsets

1. are pairwise disjoint:

$$X_k \cap X_j = \emptyset, \quad k, j \in \{1, \dots, K\}, k \neq j$$

2. **cover** *X*:

$$\bigcup_{k=1}^K X_k = X, \text{ and }$$

3. do not contain the empty set: $X_k \neq \emptyset$, $k \in \{1, ..., K\}$.

A set X_k is also called a cluster, a partition P a clustering.

The Cluster Analysis Problem (given K)



Given

- ▶ a set \mathcal{X} called **data space**, e.g., $\mathcal{X} := \mathbb{R}^M$,
- ▶ a set $X \subseteq \mathcal{X}$ called data, and
- ► a function

$$D: \bigcup_{X\subseteq \mathcal{X}} \mathsf{Part}(X) o \mathbb{R}_0^+$$

called **distortion measure** where D(P) measures how bad a partition $P \in \text{Part}(X)$ for a data set $X \subseteq \mathcal{X}$ is, and

▶ a number $K \in \mathbb{N}$ of clusters,

find a partition $P = \{X_1, X_2, ... X_K\} \in \text{Part}_{K}(X)$ with K clusters with minimal distortion D(P).

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Machine Learning 1. k-means & k-medoids

k-means: Distortion Sum of Distances to Cluster Centers. Sum of squared distances to cluster centers:

$$D(P) := \sum_{n=1}^{N} \sum_{k=1}^{K} P_{n,k} ||x_n - \mu_k||^2 = \sum_{k=1}^{K} \sum_{\substack{n=1: \ P_{n,k}=1}}^{N} ||x_n - \mu_k||^2$$

with

$$\mu_k := \frac{\sum_{n=1}^{N} P_{n,k} x_n}{\sum_{n=1}^{N} P_{n,k}} = \text{mean } \{x_n \mid P_{n,k} = 1, n \in \{1, \dots, N\}\}$$

Minimizing D over partitions with varying number of clusters leads to singleton clustering with distortion 0; only the cluster analysis problem with given K makes sense.

Minimizing D is not easy as reassigning a point to a different cluster also shifts the cluster centers.

k-means: Minimizing Distances to Cluster Centers



Add cluster centers μ as auxiliary optimization variables:

$$D(P,\mu) := \sum_{n=1}^{N} \sum_{k=1}^{K} P_{n,k} ||x_n - \mu_k||^2$$

Block coordinate descent:

1. fix μ , optimize $P \rightsquigarrow$ reassign data points to clusters:

$$P_{n,k} := \delta(k = \ell_n), \quad \ell_n := \underset{k \in \{1,...,K\}}{\operatorname{arg \, min}} ||x_n - \mu_k||^2$$

2. fix P, optimize $\mu \rightsquigarrow$ recompute cluster centers:

$$\mu_k := \frac{\sum_{n=1}^{N} P_{n,k} x_n}{\sum_{n=1}^{N} P_{n,k}}$$

Iterate until partition is stable.

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Machine Learning 1. k-means & k-medoids

k-means: Initialization



k-means is usually initialized by picking K data points as cluster centers at random:

- 1. pick the first cluster center μ_1 out of the data points at random and then
- 2. sequentially select the data point with the largest sum of distances to already choosen cluster centers as next cluster center

$$\mu_k := x_n, \quad n := \underset{n \in \{1, \dots, N\}}{\operatorname{arg max}} \sum_{\ell=1}^{k-1} ||x_n - \mu_\ell||^2, \quad k = 2, \dots, K$$

Different initializations may lead to different local minima.

- ▶ run k-means with different random initializations and
- ▶ keep only the one with the smallest distortion (random restarts).

k-means Algorithm



```
1: procedure Cluster-kmeans (\mathcal{D} := \{x_1, \dots, x_N\} \subseteq \mathbb{R}^M, K \in \mathbb{N}, \epsilon \in \mathbb{R}^+)
             n_1 \sim \operatorname{unif}(\{1,\ldots,N\}),
 2:
                                                     \mu_1 := x_{n_1}
             for k := 2, ..., K do
 3:
                   n_k := rg \max_{n \in \{1, \dots, N\}} \sum_{j=1}^{k-1} ||x_n - \mu_j||, \quad \mu_k := x_{n_k}
 4:
 5:
             repeat
                  \mu^{\mathrm{old}} := \mu
 6:
                  for n := 1, \ldots, N do
 7:
                         P_n := \operatorname{arg\,min}_{k \in \{1, \dots, K\}} ||x_n - \mu_k||
 8:
                  for k := 1, \ldots, K do
 9:
                         \mu_k := \text{mean } \{x_n \mid P_n = k, n \in \{1, \dots, N\}\}
10:
             until \frac{1}{K}\sum_{k=1}^{K}||\mu_k-\mu_k^{\mathrm{old}}||<\epsilon return P
11:
12:
```

Note: In implementations, the two loops over the data (lines 7 and 10) can be combined in one loop.

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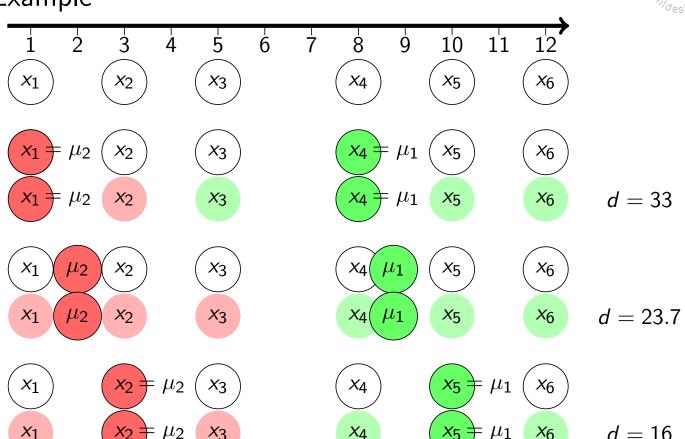




Example

 X_1

 μ_2

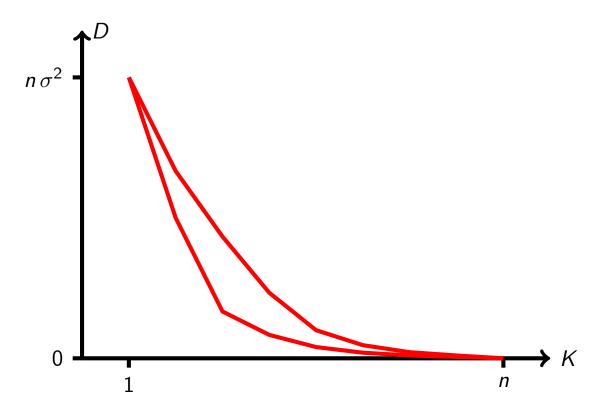


 X_4

d = 16

How Many Clusters K?





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Machine Learning 1. k-means & k-medoids

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k-medoids: k-means for General Distances

One can generalize k-means to general distances d:

$$D(P, \mu) := \sum_{n=1}^{N} \sum_{k=1}^{K} P_{n,k} d(x_n, \mu_k)$$

▶ step 1 assigning data points to clusters remains the same

$$P_{n,k} := \underset{k \in \{1,...,K\}}{\operatorname{arg \, min}} d(x_n, \mu_k)$$

▶ but step 2 finding the best cluster representatives μ_k is not solved by the mean and may be difficult in general.

idea k-medoids: choose cluster representatives out of cluster data points:

$$\mu_k := x_n, \quad n := \underset{n \in \{1, \dots, N\}: P_{n,k} = 1}{\operatorname{arg \, min}} \sum_{\ell=1}^N P_{\ell,k} d(x_\ell, x_n)$$

k-medoids: k-means for General Distances



k-medoids is a "kernel method": it requires no access to the variables, just to the distance measure.

For the Manhattan distance/ L_1 distance, step 2 finding the best cluster representatives μ_k can be solved without restriction to cluster data points:

$$(\mu_k)_m := \mathsf{median}\{(x_n)_m \mid P_{n,k} = 1, n \in \{1, \dots, N\}\}, \quad m = 1, \dots, M$$

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Machine Learning 2. Gaussian Mixture Models



Outline

- 1. k-means & k-medoids
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Soft Partitions: Row Stochastic Matrices



Let $X:=\{x_1,\ldots,x_N\}$ be a finite set. A $N\times K$ matrix

$$P \in [0,1]^{N \times K}$$

is called a **soft partition matrix of** X if it

- 1. is row-stochastic: $\sum_{k=1}^{n} P_{n,k} = 1, \qquad n \in \{1, \dots, N\}$
- 2. does not contain a zero column: $X_{.,k} \neq (0,...,0)^T$, $k \in \{1,...,K\}$

 $P_{n,k}$ is called the

- **▶** membership degree of instance *n* in class *k* or the
- ightharpoonup cluster weight of instance n in cluster k.

 $P_{.,k}$ is called membership vector of class k.

SoftPart(X) denotes the set of all soft partitions of X.

Note: Soft partitions are also called soft clusterings and fuzzy clusterings.

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Machine Learning 2. Gaussian Mixture Models

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The Soft Clustering Problem (with given K)

Given

- ▶ a set \mathcal{X} called **data space**, e.g., $\mathcal{X} := \mathbb{R}^M$,
- ▶ a set $X \subseteq \mathcal{X}$ called **data**, and
- ► a function

$$D: \bigcup_{X\subseteq \mathcal{X}} \mathsf{SoftPart}(X) o \mathbb{R}_0^+$$

called **distortion measure** where D(P) measures how bad a soft partition $P \in \mathsf{SoftPart}(X)$ for a data set $X \subseteq \mathcal{X}$ is, and

▶ a number $K \in \mathbb{N}$ of clusters,

find a soft partition $P \in \text{SoftPart}_K(X) \subseteq [0,1]^{|X| \times K}$ with K clusters with minimal distortion D(P).

Mixture Models



Mixture models assume that there exists an unobserved nominal variable Z with K levels:

$$p(X,Z) = p(Z)p(X \mid Z) = \prod_{k=1}^{K} (\pi_k p(X \mid Z = k)^{\delta(Z=k)})$$

The complete data loglikelihood of the completed data (X, Z) then is

$$\ell(\Theta; X, Z) := \sum_{n=1}^{N} \sum_{k=1}^{K} \delta(Z_n = k) (\ln \pi_k + \ln p(X = x_n \mid Z = k; \theta_k))$$
with $\Theta := (\pi_1, \dots, \pi_K, \theta_1, \dots, \theta_K)$

 ℓ cannot be computed because z_n 's are unobserved.

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Mixture Models: Expected Loglikelihood



Given an estimate $\Theta^{(t-1)}$ of the parameters, mixtures aim to optimize the **expected complete data loglikelihood**:

$$Q(\Theta; \Theta^{(t-1)}) := \mathbb{E}[\ell(\Theta; X, Z) \mid \Theta^{(t-1)}]$$

$$= \sum_{n=1}^{N} \sum_{k=1}^{K} \mathbb{E}[\delta(Z_n = k) \mid x_n, \Theta^{(t-1)}] (\ln \pi_k + \ln p(X = x_n \mid Z = k; \theta_k))$$

which is relaxed to

$$Q(\Theta, r; \Theta^{(t-1)}) = \sum_{n=1}^{N} \sum_{k=1}^{K} r_{n,k} (\ln \pi_k + \ln p(X = x_n \mid Z = k; \theta_k)) + (r_{n,k} - \mathbb{E}[\delta(Z_n = k) \mid x_n, \Theta^{(t-1)}])^2$$

Mixture Models: Expected Loglikelihood



Block coordinate descent (EM algorithm): alternate until convergence

1. expectation step:

$$r_{n,k}^{(t-1)} := \mathbb{E}[\delta(Z_n = k) \mid x_n, \Theta^{(t-1)}] = p(Z = k \mid X = x_n; \Theta^{(t-1)})$$

$$= \frac{p(X = x_n \mid Z = k; \Theta^{(t-1)}) p(Z = k; \Theta^{(t-1)})}{\sum_{k'=1}^{K} p(X = x_n \mid Z = k'; \Theta^{(t-1)}) p(Z = k'; \Theta^{(t-1)})}$$

$$= \frac{p(X = x_n \mid Z = k; \theta_k^{(t-1)}) \pi_k^{(t-1)}}{\sum_{k'=1}^{K} p(X = x_n \mid Z = k'; \theta_k^{(t-1)}) \pi_k^{(t-1)}}$$

$$(0)$$

2. maximization step:

$$\begin{split} \Theta^{(t)} &:= \argmax_{\Theta} Q(\Theta, r^{(t-1)}; \Theta^{(t-1)}) \\ &= \argmax_{\pi_1, \dots, \pi_K, \theta_1, \dots, \theta_K} \sum_{n=1}^N \sum_{k=1}^K r_{n,k} (\ln \pi_k + \ln p(X = x_n \mid Z = k; \theta_k)) \end{split}$$

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Machine Learning 2. Gaussian Mixture Models

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Mixture Models: Expected Loglikelihood

2. maximization step:

$$\Theta^{(t)} = \underset{\pi_1, ..., \pi_K, \theta_1, ..., \theta_K}{\arg \max} \sum_{n=1}^{N} \sum_{k=1}^{K} r_{n,k} (\ln \pi_k + \ln p(X = x_n \mid Z = k; \theta_k))$$

$$\rightsquigarrow \quad \pi_k^{(t)} = \frac{\sum_{n=1}^N r_{n,k}}{N} \tag{1}$$

$$\sum_{n=1}^{N} \frac{r_{n,k}}{p(X=x_n \mid Z=k; \theta_k)} \frac{\partial p(X=x_n \mid Z=k; \theta_k)}{\partial \theta_k} = 0, \quad \forall k$$
 (*)

(*) needs to be solved for specific cluster specific distributions p(X|Z).

Gaussian Mixtures



Gaussian mixtures:

▶ use Gaussians for p(X|Z):

$$p(X = x \mid Z = k) = \frac{1}{\sqrt{(2\pi)^{M}|\Sigma_{k}|}} e^{-\frac{1}{2}(x-\mu_{k})^{T}\Sigma_{k}^{-1}(x-\mu_{k})}, \quad \theta_{k} := (\mu_{k}, \Sigma_{k})$$

$$\Rightarrow \quad \mu_{k}^{(t)} = \frac{\sum_{n=1}^{N} r_{n,k}^{(t-1)} x_{n}}{\sum_{n=1}^{N} r_{n,k}^{(t-1)}}$$

$$\Sigma_{k}^{(t)} = \frac{\sum_{n=1}^{N} r_{n,k}^{(t-1)} (x_{n} - \mu_{k}^{(t)})^{T} (x_{n} - \mu_{k}^{(t)})}{\sum_{n=1}^{N} r_{n,k}^{(t-1)}}$$

$$= \frac{\sum_{n=1}^{N} r_{n,k}^{(t-1)} x_{n}^{T} x_{n} - \mu_{k}^{(t)} T \mu_{k}^{(t)}}{\sum_{n=1}^{N} r_{n,k}^{(t-1)} x_{n}^{T} x_{n} - \mu_{k}^{(t)} T \mu_{k}^{(t)}}$$

$$(3)$$

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Machine Learning 2. Gaussian Mixture Models



Gaussian Mixtures: EM Algorithm, Summary

1. expectation step: $\forall n, k$

$$\tilde{r}_{n,k}^{(t-1)} = \pi_k^{(t-1)} \frac{1}{\sqrt{(2\pi)^M |\Sigma_k^{(t-1)}|}} e^{-\frac{1}{2}(x_n - \mu_k^{(t-1)})^T \Sigma_k^{(t-1) - 1} (x_n - \mu_k^{(t-1)})} \quad (0a)$$

$$r_{n,k}^{(t-1)} = \frac{\tilde{r}_{n,k}^{(t-1)}}{\sum_{k'=1}^{K} \tilde{r}_{n,k'}^{(t-1)}}$$
(0b)

2. maximization step: $\forall k$

$$\pi_k^{(t)} = \frac{\sum_{n=1}^{N} r_{n,k}^{(t-1)}}{N} \tag{1}$$

$$\mu_k^{(t)} = \frac{\sum_{n=1}^{N} r_{n,k}^{(t-1)} x_n}{\sum_{n=1}^{N} r_{n,k}^{(t-1)}}$$
(2)

$$\Sigma_{k}^{(t)} = \frac{\sum_{n=1}^{N} r_{n,k}^{(t-1)} x_{n}^{T} x_{n} - \mu_{k}^{(t)} T \mu_{k}^{(t)}}{\sum_{n=1}^{N} r_{n,k}^{(t-1)}}$$
(3)

Gaussian Mixtures for Soft Clustering



▶ The **responsibilities** $r \in [0,1]^{N \times K}$ are a soft partition.

$$P := r$$

▶ The negative expected loglikelihood can be used as cluster distortion:

$$D(P) := -\max_{\Theta} Q(\Theta, r)$$

 \blacktriangleright To optimize D, we simply can run EM.

For hard clustering:

assign points to the cluster with highest responsibility (hard EM):

$$r_{n,k}^{(t-1)} = \delta(k = \underset{k'=1,...,K}{\operatorname{arg max}} \, \tilde{r}_{n,k'}^{(t-1)})$$
 (0b')

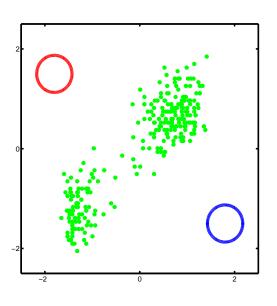
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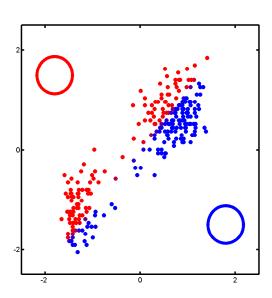
Machine Learning 2. Gaussian Mixture Models

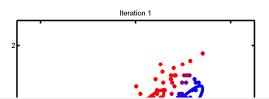
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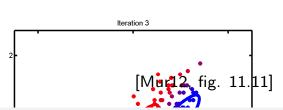


Gaussian Mixtures for Soft Clustering / Example









Model-based Cluster Analysis



Different parametrizations of the covariance matrices Σ_k restrict possible cluster shapes:

- full Σ: all sorts of ellipsoid clusters.
- ► diagonal Σ: ellipsoid clusters with axis-parallel axes
- unit Σ: spherical clusters.

One also distinguishes

- ▶ cluster-specific Σ_k : each cluster can have its own shape.
- ▶ shared $\Sigma_k = \Sigma$: all clusters have the same shape.

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Machine Learning 2. Gaussian Mixture Models

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k-means: Hard EM with spherical clusters

1. expectation step: $\forall n, k$

$$\begin{split} \tilde{r}_{n,k}^{(t-1)} &= \frac{1}{\sqrt{(2\pi)^{M}|\Sigma_{k}^{(t-1)}|}} e^{-\frac{1}{2}(x_{n} - \mu_{k}^{(t-1)})^{T} \Sigma_{k}^{(t-1) - 1}(x_{n} - \mu_{k}^{(t-1)})}} \quad \text{(0a)} \\ &= \frac{1}{\sqrt{(2\pi)^{M}}} e^{-\frac{1}{2}(x_{n} - \mu_{k}^{(t-1)})^{T}(x_{n} - \mu_{k}^{(t-1)})} \\ r_{n,k}^{(t-1)} &= \delta(k = \underset{k'=1,\dots,K}{\text{arg max }} \tilde{r}_{n,k'}^{(t-1)}) \\ \arg\max_{k'=1,\dots,K} \tilde{r}_{n,k'}^{(t-1)} &= \underset{k'=1,\dots,K}{\text{arg max }} \frac{1}{\sqrt{(2\pi)^{M}}} e^{-\frac{1}{2}(x_{n} - \mu_{k}^{(t-1)})^{T}(x_{n} - \mu_{k}^{(t-1)})} \\ &= \underset{k'=1,\dots,K}{\text{arg min }} ||x_{n} - \mu_{k}^{(t-1)}||^{2} \\ &= \underset{k'=1,\dots,K}{\text{arg min }} ||x_{n} - \mu_{k}^{(t-1)}||^{2} \end{split}$$

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Machine Learning 3. Hierarchical Cluster Analysis



Hierarchies

Let X be a set.

A tree (H, E), $E \subseteq H \times H$ edges pointing towards root

- lacktriangle with leaf nodes h corresponding bijectively to elements $x_h \in X$
- ▶ plus a surjective map L : $H \rightarrow \{0, \ldots, d\}, d \in \mathbb{N}$ with
 - ► L(root) = 0 and
 - ▶ L(h) = d for all leaves $h \in H$ and
 - ▶ $L(h) \le L(g)$ for all $(g, h) \in E$

called level map

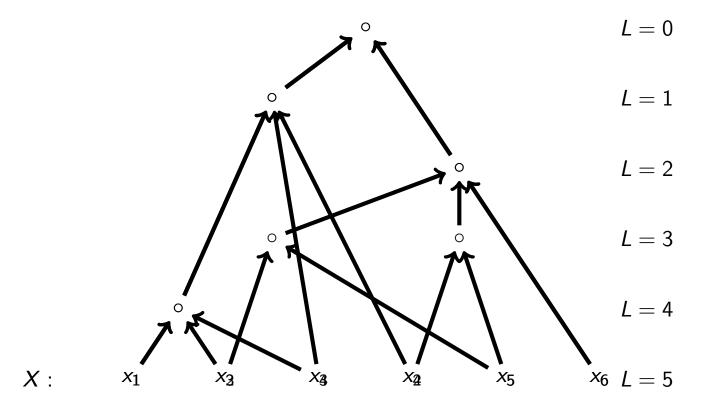
is called an **hierarchy over** X.

d is called the **depth** of the hierarchy.

Hier(X) denotes the set of all hierarchies over X.

Hierarchies / Example





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Machine Learning 3. Hierarchical Cluster Analysis

Ners/E



Hierarchies: Nodes Correspond to Subsets

Let (H, E) be such an hierarchy:

- ▶ nodes of an hierarchy correspond to subsets of X:
 - ▶ leaf nodes *h* correspond to a singleton subset by definition.

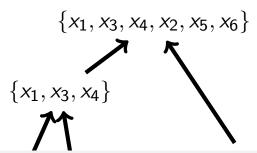
$$subset(h) := \{x_h\}, x_h \in X \text{ corresponding to leaf } h$$

▶ interior nodes *h* correspond to the union of the subsets of their children:

$$subset(h) := \bigcup_{\substack{g \in H \\ (g,h) \in E}} subset(g)$$

▶ thus the root node *h* corresponds to the full set *X*:

$$subset(h) = X$$



Hierarchies: Levels Correspond to Partitions

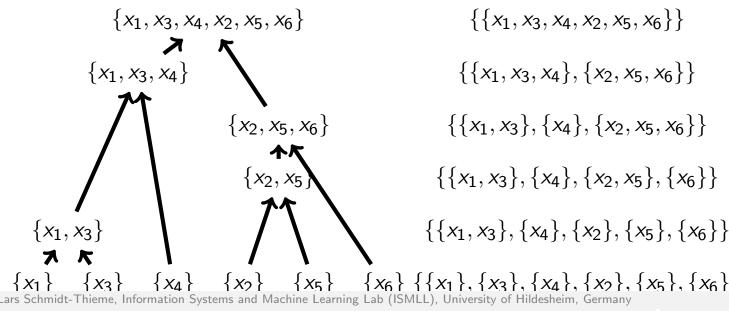


Let (H, E) be such an hierarchy:

▶ levels $\ell \in \{0, ..., d\}$ correspond to partitions

$$P_{\ell}(H, L) := \{ h \in H \mid L(h) \ge \ell, \not\exists g \in H : L(g) \ge \ell,$$

 $\mathsf{subset}(h) \subsetneq \mathsf{subset}(g) \}$



Machine Learning 3. Hierarchical Cluster Analysis

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The Hierarchical Cluster Analysis Problem



Given

- ▶ a set \mathcal{X} called **data space**, e.g., $\mathcal{X} := \mathbb{R}^M$,
- ▶ a set $X \subseteq \mathcal{X}$ called **data** and
- ► a function

$$D: \bigcup_{X\subseteq \mathcal{X}} \operatorname{Hier}(X) o \mathbb{R}_0^+$$

called **distortion measure** where D(P) measures how bad a hierarchy $H \in \text{Hier}(X)$ for a data set $X \subseteq \mathcal{X}$ is,

find a hierarchy $H \in Hier(X)$ with minimal distortion D(H).

Distortions for Hierarchies



Examples for distortions for hierarchies:

$$D(H) := \sum_{K=1}^{N} \tilde{D}(P_K(H))$$

where

- $ightharpoonup P_K(H)$ denotes the partition at level K-1 (with K classes) and
- ullet $ilde{D}$ denotes a distortion for partitions.

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Machine Learning 3. Hierarchical Cluster Analysis

Agglomerative and Divisive Hierarchical Clustering

Hierarchies are usually learned by greedy search level by level:

- agglomerative clustering:
 - 1. start with the singleton partition P_N :

$$P_N := \{X_k \mid k = 1, \dots, N\}, \quad X_k := \{x_k\}, \quad k = 1, \dots, N$$

2. in each step $K = N, \ldots, 2$ build P_{K-1} by joining the two clusters $k, \ell \in \{1, \dots, K\}$ that lead to the minimal distortion

$$D(\{X_1,\ldots,\widehat{X_k},\ldots,\widehat{X_\ell},\ldots,X_K,X_k\cup X_\ell))$$

- divisive clustering:
 - 1. start with the all partition P_1 :

$$P_1 := \{X\}$$

2. in each step K = 1, N-1 build P_{K+1} by splitting one cluster X_k in two clusters X_k', X_ℓ' that lead to the minimal distortion

$$D(\{X_1,\ldots,\widehat{X_k},\ldots,X_K,X_k',X_\ell'),\quad X_k=X_k'\cup X_\ell'$$
 Note: $\widehat{X_k}$ denotes that the class X_k is omitted from the partition.

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Class-wise Defined Partition Distortions



If the partition distortion can be written as a sum of distortions of its classes,

$$D(\{X_1,\ldots,X_K\})=\sum_{k=1}^K \tilde{D}(X_k)$$

then the optimal pair does only depend on X_k, X_ℓ :

$$D(\{X_1,\ldots,\widehat{X_k},\ldots,\widehat{X_\ell},\ldots,X_K,X_k\cup X_\ell)=\tilde{D}(X_k\cup X_\ell)-(\tilde{D}(X_k)+\tilde{D}(X_\ell))$$

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Machine Learning 3. Hierarchical Cluster Analysis

Ners/F.

Closest Cluster Pair Partition Distortions



For a cluster distance

$$ilde{d}: \mathcal{P}(X) imes \mathcal{P}(X) o \mathbb{R}_0^+$$
 with $ilde{d}(A \cup B, C) \geq \min \{ ilde{d}(A, C), ilde{d}(B, C)\}, \quad A, B, C \subseteq X$

a partition can be judged by the closest cluster pair it contains:

$$D(\{X_1,\ldots,X_K\}) = \min_{k,\ell=1,K\atop k\neq\ell} \tilde{d}(X_k,X_\ell)$$

Such a distortion has to be maximized.

To increase it, the closest cluster pair has to be joined.

Single Link Clustering



$$d_{\mathsf{sl}}(A,B) := \min_{x \in A, y \in B} d(x,y), \quad A,B \subseteq X$$

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Complete Link Clustering



$$d_{\mathsf{cl}}(A,B) := \max_{x \in A, y \in B} d(x,y), \quad A,B \subseteq X$$

Average Link Clustering



$$d_{\mathsf{al}}(A,B) := \frac{1}{|A||B|} \sum_{x \in A, y \in B} d(x,y), \quad A,B \subseteq X$$

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Recursion Formulas for Cluster Distances

$$\begin{split} d_{\text{sl}}(X_{i} \cup X_{j}, X_{k}) &:= \min_{x \in X_{i} \cup X_{j}, y \in X_{k}} d(x, y) \\ &= \min \{ \min_{x \in X_{i}, y \in X_{k}} d(x, y), \min_{x \in X_{j}, y \in X_{k}} d(x, y) \} \\ &= \min \{ d_{\text{sl}}(X_{i}, X_{k}), d_{\text{sl}}(X_{j}, X_{k}) \} \\ d_{\text{cl}}(X_{i} \cup X_{j}, X_{k}) &:= \max_{x \in X_{i} \cup X_{j}, y \in X_{k}} d(x, y) \\ &= \max \{ \max_{x \in X_{i}, y \in X_{k}} d(x, y), \max_{x \in X_{j}, y \in X_{k}} d(x, y) \} \\ &= \max \{ d_{\text{cl}}(X_{i}, X_{k}), d_{\text{cl}}(X_{j}, X_{k}) \} \\ d_{\text{al}}(X_{i} \cup X_{j}, X_{k}) &:= \frac{1}{|X_{i} \cup X_{j}|} \sum_{x \in X_{i} \cup X_{j}, y \in X_{k}} d(x, y) \\ &= \frac{|X_{i}|}{|X_{i} \cup X_{j}|} \frac{1}{|X_{i}||X_{k}|} \sum_{x \in X_{i} \cup X_{j}, y \in X_{k}} d(x, y) \end{split}$$

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Conclusion (1/2)



- ► Cluster analysis aims at **detecting latent groups** in data, without labeled examples (↔ **record linkage**).
- ► Latent groups can be described in three different granularities:
 - ▶ partitions segment data into K subsets (hard clustering).
 - ► hierarchies structure data into an hierarchy, in a sequence of consistent partitions (hierarchical clustering).
 - ► soft clusterings / row-stochastic matrices build overlapping groups to which data points can belong with some membership degree (soft clustering).
- ▶ k-means finds a K-partition by finding K cluster centers with smallest Euclidean distance to all their cluster points.
- ▶ k-medoids generalizes k-means to general distances; it finds a K-partition by selecting K data points as cluster representatives with smallest distance to all their cluster points.

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Conclusion (2/2)

- hierarchical single link, complete link and average link methods
 - ▶ find a hierarchy by greedy search over consistent partitions,
 - starting from the singleton parition (agglomerative)
 - being efficient due to recursion formulas,
 - ► requiring only a distance matrix.
- ▶ Gaussian Mixture Models find soft clusterings by modeling data by a class-specific multivariate Gaussian distribution $p(X \mid Z)$ and estimating expected class memberships (expected likelihood).
- ► The Expectation Maximiation Algorithm (EM) can be used to learn Gaussian Mixture Models via block coordinate descent.
- ▶ k-means is a special case of a Gaussian Mixture Model
 - ▶ with hard/binary cluster memberships (hard EM) and
 - spherical cluster shapes.

Readings



- ▶ k-means:
 - ► [HTFF05], ch. 14.3.6, 13.2.3, 8.5 [Bis06], ch. 9.1, [Mur12], ch. 11.4.2
- ► hierarchical cluster analysis:
 - ► [HTFF05], ch. 14.3.12, [Mur12], ch. 25.5. [PTVF07], ch. 16.4.
- ► Gaussian mixtures:
 - ► [HTFF05], ch. 14.3.7, [Bis06], ch. 9.2, [Mur12], ch. 11.2.3, [PTVF07], ch. 16.1.

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