STAT406- Methods of Statistical Learning Lecture 19

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Unsupervised learning

- Unsupervised ≠ Supervised
- High-"density" regions (w/o model)
- Agglomerative / hierarchical methods
- High-"density" regions (with a model)EM-algorithm
- Dimension reduction (PCA, MDS, etc.)

Clustering - Problem

Data: p features / variables per "unit"

$$\mathbf{X} = \begin{pmatrix} X_1 \\ X_2 \\ \vdots \\ X_p \end{pmatrix}$$

• X_1, \ldots, X_n

- Goal: find regions where X_i's are "clustered"
- Goal: find regions where P(X) is relatively high
- These regions are sometimes modeled

- Lower dimensional subspaces (linear manifolds)
 - **Principal Components**
- Convex regions with high P(X)
 K-means / K-medoids Hierarchical methods

- Intrinsically different from classification
- There is no clear performance measure to evaluate "success"
- Hence the name: "unsupervised learning"

Example 1 – 9 Breweries - 26 attributes

```
> a
   [,1] [,2] [,3] [,4] [,5] [,6] [,7] [,8] [,9] [...]
V1 3.51 3.41 3.20 2.73 2.35 3.03 2.21 3.91 3.07 [...]
V2 4.43 4.05 3.66 5.25 3.88 4.23 3.27 2.71 4.08 [...]
V3 4.76 3.42 4.22 2.44 4.18 2.47 3.67 4.59 4.74 [...]
V4 3.68 3.78 3.07 2.75 2.78 3.12 2.49 3.91 3.34 [...]
V5 4.77 1.04 3.86 5.28 3.86 4.24 3.40 4.23 4.23 [...]
[\ldots]
   [,15] [,16] [,17] [,18] [,19] [,20] [,21] [,22] [...]
V1 3.07 3.45 2.53 3.12 2.93 2.24 2.41 3.32 [...]
V2 3.82 4.29 4.71 3.58 3.27
                                3.11 3.14 3.74 [...]
V3 4.17 4.44 4.53 4.10 4.13
                                4.12 3.43 4.32 [...]
V4 3.21 3.74 2.83 3.14 2.80 2.39 2.40 3.32 [...]
V5 3.94 4.47 4.83 3.82 3.46
                                 3.39 3.22 4.01 [...]
[\ldots]
```

$$\mathbf{X}_1, \ \mathbf{X}_2, \ \dots \ \mathbf{X}_9 \in \mathbb{R}^{26}$$

Do they appear grouped / clustered?

UN Votes

- From http://hdl.handle.net/1902.1/12379
- UN, founded 1946, 193 members
- "important" votes (U.S. State Department)
- Votes: Yes (1), Abstain (2), No (3), Absent (8), Not a Member (9)
- 368 important votes, 77 countries voted > 95% of these

UN Votes

- Do voting patterns reflect political alignments?
- Do countries vote along known political blocks?
- Data: X_i: votes for country i

$$X_i \in \mathbb{R}^{368}$$
, $i = 1, ..., 77$ (countries)

What groups are there?

Cancer example

- From [HTF09], details in script
- Gene expression for 64 samples
- There are 6830 genes
- $X_1, X_2, \ldots, X_{64} \in \mathbb{R}^{6830}$
- We know tissue type for ea. sample
- Really: "feature selection"

- Look for convex sets of relative high density
- The number of sets K is specified a priori (but we'll come back to this)
- Since "high density" is related to "closeness"

$$\min \sum_{r=1}^{\mathbf{K}} \sum_{i,j \in \mathcal{C}_r} d^2 \left(\mathbf{X}_i, \mathbf{X}_j \right)$$

minimize over all partitions C_1, \ldots, C_K

Note that

$$\begin{split} &\sum_{i=1}^{n} \sum_{j=1}^{n} d^{2} \left(\mathbf{X}_{i}, \mathbf{X}_{j} \right) = \sum_{r=1}^{K} \sum_{i \in \mathcal{C}_{r}} \sum_{j=1}^{n} d^{2} \left(\mathbf{X}_{i}, \mathbf{X}_{j} \right) \\ &= \sum_{r=1}^{K} \sum_{i \in \mathcal{C}_{r}} \left[\sum_{j \in \mathcal{C}_{r}} d^{2} \left(\mathbf{X}_{i}, \mathbf{X}_{j} \right) + \sum_{j \notin \mathcal{C}_{r}} d^{2} \left(\mathbf{X}_{i}, \mathbf{X}_{j} \right) \right] \\ &\sum_{r=1}^{K} \sum_{i,j \in \mathcal{C}_{r}} d^{2} \left(\mathbf{X}_{i}, \mathbf{X}_{j} \right) + \sum_{r=1}^{K} \sum_{i \in \mathcal{C}_{r}} \sum_{j \notin \mathcal{C}_{r}} d^{2} \left(\mathbf{X}_{i}, \mathbf{X}_{j} \right) \\ &T = W + B \end{split}$$

When
$$d^{2}(\mathbf{X}_{i}, \mathbf{X}_{j}) = \|\mathbf{X}_{i} - \mathbf{X}_{j}\|^{2}$$

$$W = \sum_{r=1}^{K} \sum_{i,j \in \mathcal{C}_{r}} \|\mathbf{X}_{i} - \mathbf{X}_{j}\|^{2} = \sum_{r=1}^{K} \sum_{i \in \mathcal{C}_{r}} \|\mathbf{X}_{i} - \bar{\mathbf{X}}_{r}\|^{2}$$

• Given C_1, C_2, \ldots, C_K , assign X_i to the cluster C_i with closest mean

$$\mathbf{X}_{i} \leftarrow \arg\min_{1 \leq i \leq K} \left\| \mathbf{X}_{i} - \bar{\mathbf{X}}_{j} \right\|^{2}$$

Note that

$$\bar{\mathbf{X}}_r = \hat{\mu}_r = \arg\min_{\mu} \sum_{i \in \mathcal{C}_r} \|\mathbf{X}_i - \mu\|^2$$

• Given $\hat{\mu}_1, \ldots, \hat{\mu}_K$

$$\min_{\mathcal{C}_1, \dots, \mathcal{C}_k} \sum_{r=1}^k \sum_{i \in \mathcal{C}_r} \|\mathbf{x}_i - \hat{\boldsymbol{\mu}}_r\|^2$$

is attained with

$$\mathbf{X}_i \leftarrow \arg\min_{1 \leq i \leq K} \left\| \mathbf{X}_i - \hat{\boldsymbol{\mu}}_j \right\|^2$$

• And, given C_1, \ldots, C_K

$$\min_{\hat{\boldsymbol{\mu}}_1, \dots, \hat{\boldsymbol{\mu}}_K} \sum_{r=1}^k \sum_{i \in \mathcal{C}_r} \|\mathbf{x}_i - \hat{\boldsymbol{\mu}}_r\|^2$$

is attained with

$$\hat{\boldsymbol{\mu}}_r \leftarrow \bar{\mathbf{X}}_r = \frac{1}{n_r} \sum_{i \in \mathcal{C}_r} \mathbf{x}_i$$

This suggests a simple iterative (and greedy) algorithm.

Remarks

- Algorithm is greedy
- Answer depends on the initial configuration
- It needs to be started from many initial configurations

Cancer data

```
> set.seed(31)
> nci.km <- kmeans(nci, centers=8)</pre>
> table(nci.km$cluster)
 1 2 3 4 5 6 7 8
8 6 6 14 3 8 4 15
> set.seed(311)
> nci.km <- kmeans(nci, centers=8)</pre>
> table(nci.km$cluster)
 1 2 3 4 5 6 7 8
4 12 6 9 4 8 19 2
```

Need **more** starting points...

```
> set.seed(31)
> nci.km <- kmeans(nci, centers=8, iter.max = 5000,
 nstart=1000)
> table(nci.km$cluster)
 1 2 3 4 5 6 7 8
 3 8 5 14 6 15 9 4
> set.seed(311)
> nci.km <- kmeans(nci, centers=8, iter.max = 5000,</pre>
 nstart=1000)
> table(nci.km$cluster)
 4 5 8 9 14 3 15 6
```

These clusters are the same

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UN Votes example

Not all countries voted each time

```
> dim(X)
[1] 368 77
> sum(complete.cases(X))
[1] 145
```

Only use resolutions with full votes

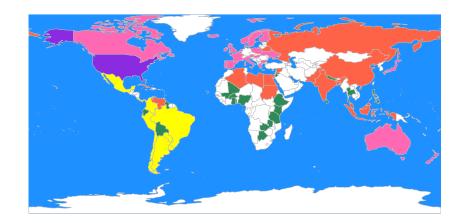
```
X2 <- X[complete.cases(X),]</pre>
```

• Use kmeans in R

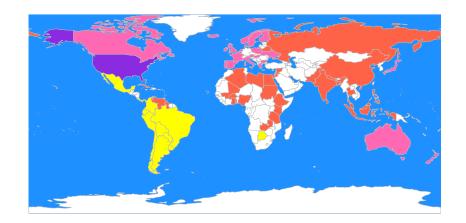
UN Votes example

```
> set.seed(123)
> b <- kmeans(t(X2), centers=5,
                   iter.max=20, nstart=1)
> table(b$cluster)
1 2 3 4 5
18 2 7 19 31
> b <- kmeans(t(X2), centers=5,
                   iter.max=20, nstart=1)
> table(b$cluster)
1 2 3 4 5
27 12 13 7 18
```

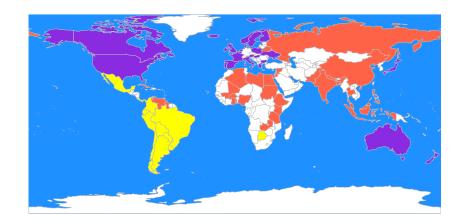
UN Votes example - K=5



UN Votes example - K=4



UN Votes example - K=3



K-means++

- A cleverly chosen set of initial centres
- K-means++
 - Pick a centre c₁ at random (from data)
 - Then for j in 2:k
 - Compute weights

$$w_i = \min\left(d^2(\boldsymbol{x}_i,\boldsymbol{c}_1),\dots,d^2(\boldsymbol{x}_i,\boldsymbol{c}_{j-1})\right)\,,$$

- ▶ Pick next centre \mathbf{c}_j from data with prob $\propto d_i$
- Implemented in flexclust::kcca

Choosing *K*

For each cluster C_r , let

$$W\left(\mathcal{C}_{r}\right) = \sum_{i,i \in \mathcal{C}_{r}} d^{2}\left(\mathbf{X}_{i},\mathbf{X}_{j}\right) \qquad r = 1,\ldots,\mathbf{K}$$

and

$$W_{\mathbf{K}} = \sum_{i=1}^{\mathbf{K}} W(\mathcal{C}_{r})$$

Choosing *K*

- Note that selecting K to minimize W_K does not generally work
- W_K typically decreases with K
- A simple example follows

Selecting the number *K* of clusters

For each cluster C_r , let

$$W\left(\mathcal{C}_{r}\right) = \sum_{i,i \in \mathcal{C}_{r}} d\left(\mathbf{X}_{i}, \mathbf{X}_{j}\right) \qquad r = 1, \dots, \mathbf{K}$$

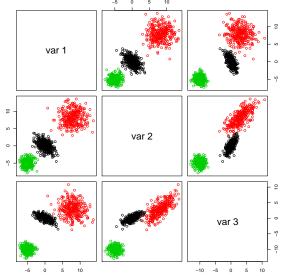
and

$$W_{\mathbf{K}} = \sum_{i=1}^{\mathbf{K}} W(\mathcal{C}_r)$$

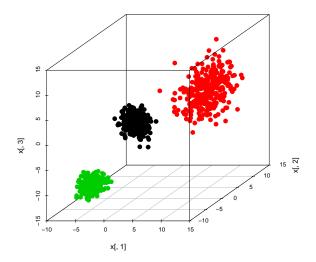
Selecting the number *K* of clusters

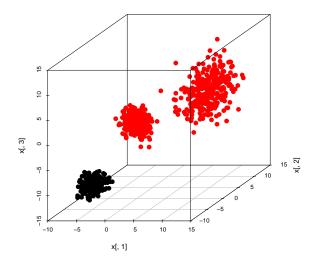
- Note that selecting K to minimize W_K does not generally work
- W_K typically decreases with K
- A simple example follows

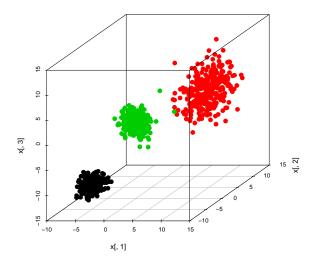
Pairs plot - Easy case

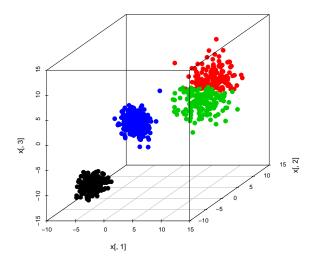


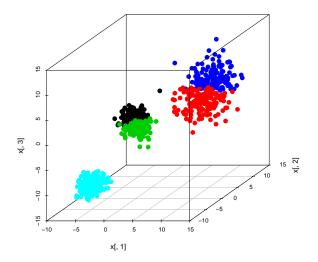
Pairs plot - Easy case

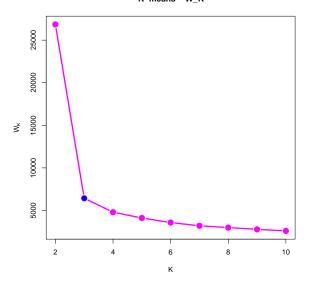




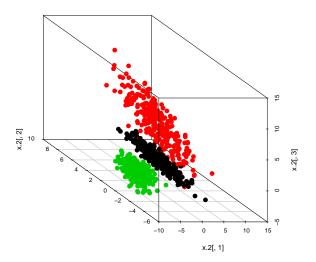


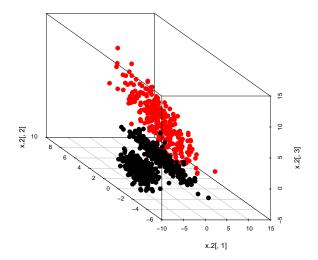


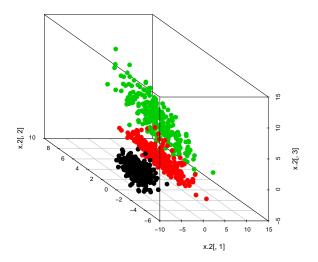


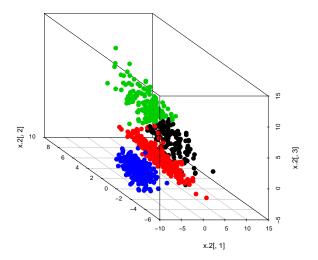


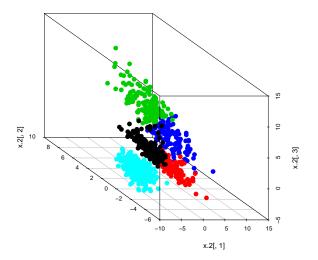
Pairs plot - K-means

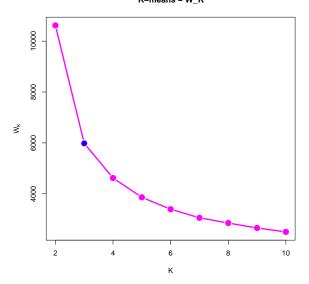












GAP Statistic

GAP Statistic (Tibshirani, Walther and Hastie, 2001)

Consider

$$G(\mathbf{K}) = E[\log(W_{\mathbf{K}})] - \log(W_{\mathbf{K}})$$

where $E[\log(W_{\mathbf{K}})]$ is the expected value under a certain reference distribution

Clest algorithm

Clest algorithm

Idea - select the value of **K** that produces classes that are best predicted by your favourite classification method.

Dudoit, Fridlyand, 2002, A prediction-based resapmling method for estimating the number of clusters in a dataset, Genome Biology **3(7)**: research0036.1 - 0036.21

Other approaches to select **K**

Dudoit, Fridlyand, 2003, Bagging to improve the accuracy of a clustering procedure, Bioinformatics, **19**, 1090-1099

Note that in K-means

- We used $d^2(\mathbf{X}_i, \mathbf{X}_j) = \|\mathbf{X}_i \mathbf{X}_j\|^2$
- The cluster "centers" may not be actual observations
- Need to manipulate the "features" (X_i)
- Can we use different distance measures?
- Can we work with the dissimilarities only?

A slightly different algorithm is

• Given C_1 , C_2 , ..., C_K , for each cluster C_r find

$$\mathbf{j}_{r}^{*} = \arg\min_{i \in \mathcal{C}_{r}} \sum_{j \in \mathcal{C}_{r}} d\left(\mathbf{X}_{i}, \mathbf{X}_{j}\right)$$

and let $m_r = \mathbf{X}_{j_r^*}$

• Given $m_1, m_2, ..., m_K$, assign X_i to the cluster C_i with closest centre:

$$\mathbf{X}_{i} \leftarrow \arg\min_{1 \leq i \leq K} d\left(\mathbf{X}_{i}, \mathbf{m}_{j}\right)$$

- 1. Find K initial cluster centres
- 2. Given centres m_{ℓ} , assign points to the cluster C_i with closest centre:

$$\mathbf{X}_i \leftarrow \arg\min_{1 \leq i \leq K} d\left(\mathbf{X}_i, \mathbf{m}_j\right)$$

- 3. Explore all possible swaps between centres and non-centres.
- 4. If there's improvement, go to step 2

Note that now

- We can use any distance robustness?
- The cluster representatives / prototypes are actual observations
- We do not need the observations, only the dissimilarities

Beers - 9 beers with 26 attributes

```
> a <- read.table('breweries.dat', header=FALSE)</pre>
> a < - t(a)
> a.dis <- dist(a, method='manhattan')</pre>
>
> brew.pam <- pam(a.dis, k=3)
>
> brew.pam
Medoids:
     ID
[1,] "7" "V7"
[2,] "2" "V2"
[3,] "6" "V6"
Clustering vector:
V1 V2 V3 V4 V5 V6 V7 V8 V9
 1 2 3 1 2 3 1 3 2
```

Silhouette plot

• For each unit $\mathbf{X}_i \in \mathcal{C}_\ell$

$$a_i = \frac{1}{n_\ell} \sum_{\mathbf{X}_i \in \mathcal{C}_\ell} d\left(\mathbf{X}_i, \mathbf{X}_i\right)$$

Dissimilarity with other clusters

$$d(i, C_r) = \frac{1}{n_r} \sum_{\mathbf{X}_i \in C_r} d(\mathbf{X}_i, \mathbf{X}_j)$$

Silhouette plot

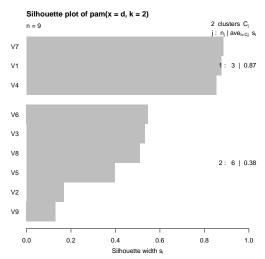
Then, dissimilarity to closest cluster

$$b_i = \min_{r \neq \ell} d(i, C_r)$$

Silhouette

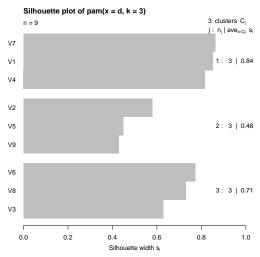
$$s_i = (b_i - a_i)/\max(a_i, b_i)$$

Breweries - K=2



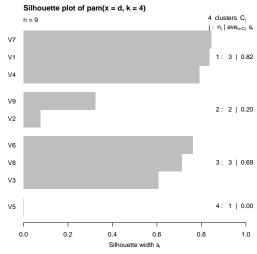
Average silhouette width: 0.54

Breweries - K=3



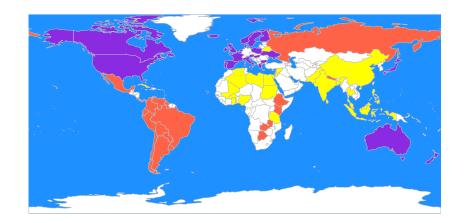
Average silhouette width: 0.68

Breweries - K=4



Average silhouette width: 0.55

UN Votes PAM - K=3



UN Votes Kmeans - K=3

