Statistics 452: Statistical Learning and Prediction Chapter 10, part 2: Cluster Analysis

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Clustering

- Clustering is a set of techniques for finding subgroups, or clusters, in a data set.
- ▶ A good set of clusters is a partition of the data such that observations are *similar* within clusters and *dissimilar* between clusters.
- Example: Suppose we have n observations of p features for cancer patients.
 - We can cluster to look for cancer sub-types.
 - However, these clusters will be highly dependent on how we measure similarity/dissimilarity.
 - Plus, we don't know that clusters correspond to cancer sub-types.

Clustering as Dimension Reduction

- ► We can view clustering as an exploratory method to understand possible structure in our data.
- ▶ Similar in spirit to PCA, but a different mechanism
 - ► PCA finds a low-dimensional representation of observations explaining a good proportion of the variance
 - Clustering looks to find homogeneous sub-groups.

Clustering Methods

- ▶ There are many.
- ▶ We focus on *K*-means/medoids and hierarchical clustering, and assume all features are quantitative.
- ▶ In K-means we partition into a pre-specified number of clusters.
- In hierarchical clustering we successively group observations.
 - Represent the nested partitions as a tree-like structure called a dendrogram.

K-means Clustering

- Choose the desired number of clusters, K. The K-means clustering algorithm will assign each observation to exactly one of the K clusters.
- Let C_k be the set of indices for observations assigned to cluster k, k = 1, ..., K.
- K-means clustering chooses clusters to solve

$$\min_{C_1,\ldots,C_K}\sum_{k=1}^K W(C_k),$$

where $W(C_k)$ is a measure of the amount by which observations within cluster C_k differ from one another.

K-means Clustering: Choice of $W(C_k)$

▶ For quantitative variables, the standard choice for $W(C_k)$ is the squared Euclidean distance,

$$W(C_k) = \frac{1}{|C_k|} \sum_{i,i' \in C_k} \sum_{j=1}^{p} (x_{ij} - x_{i'j})^2$$

where $|C_k|$ denotes the number of observations in cluster k.

▶ It turns out that

$$\frac{1}{|C_k|} \sum_{i,i' \in C_k} \sum_{j=1}^p (x_{ij} - x_{i'j})^2 = 2 \sum_{i \in C_k} \sum_{j=1}^p (x_{ij} - \bar{x}_{kj})^2$$

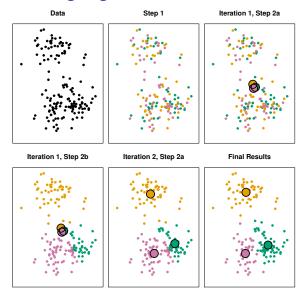
where the cluster centroid $\bar{x}_{kj} = \frac{1}{|C_k|} \sum_{i \in C_k} x_{ij}$.

▶ We can see where the name "K-means" comes from.

K-means Clustering Algorithm

- 1. Randomly assign a number, from $1, \ldots, K$ to each of the observations. They serve as the initial cluster assignments for the observations.
- Iterate the following steps until the cluster assignments stop changing:
 - (a) For each of the *K* clusters, compute the cluster centroid. The *k*th cluster centroid is the vector of the *p* feature means for the observations in the *k*th cluster for continuous variables.
 - (b) Assign each observation to the cluster whose centroid is closest in terms of Euclidean distance.

K-means Clustering Algorithm, Simulated Data Example



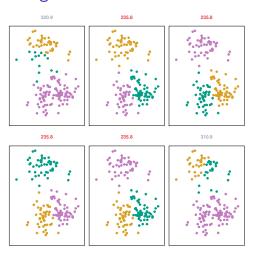
Text, Figure 10.6. Progress of the K-means algorithm with K=3.

See the video demo at http://www.youtube.com/watch?v=zaKjh2N8jN4.

K-means Clustering and Local Minima

- ▶ The algorithm is guaranteed to decrease the value of $\sum_k W(C_k)$ until no further improvement is possible, resulting in a *local* minimum.
- ► However, the local min depends on the random initialization. Hence, in practice we re-run the algorithm several times and keep the solution that achieves the lowest overall criterion value.

K-means Clustering with Different Initialization



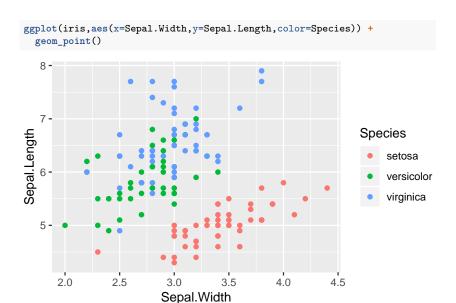
Text, Fig 10.7. K-means clustering performed six times on the same data with K=3 different random initializations. Above each plot is the value of the objective function. Three different local optimums were found. The overall minimum, found four times, is 235.8. Note: Cluster labels different in 3 of 4.

K Means Clustering of the Iris Data

► We know there are three species of iris. Ignore the species labels and do the clustering.

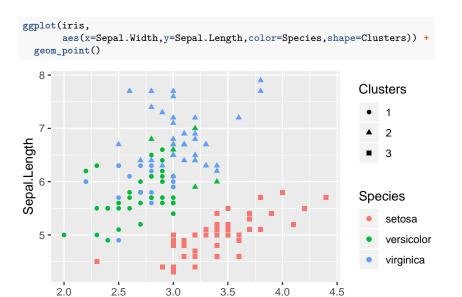
```
library(ggplot2)
data(iris) # help(iris)
head(iris) # plot sepal length vs width with labels
```

```
Sepal.Length Sepal.Width Petal.Length Petal.Width Species
##
## 1
              5.1
                          3.5
                                       1.4
                                                  0.2
                                                        setosa
              4.9
                                       1.4
                                                  0.2 setosa
## 2
                         3.0
## 3
             4.7
                         3.2
                                       1.3
                                                  0.2 setosa
             4.6
## 4
                         3.1
                                      1.5
                                                  0.2 setosa
             5.0
                         3.6
                                      1.4
                                                  0.2 setosa
## 5
## 6
             5.4
                         3.9
                                      1.7
                                                  0.4
                                                        setosa
```



```
library(dplyr)
set.seed(1)
irisX <- iris %>% select(-Species) %>% scale()
kout3 <- kmeans(irisX,centers=3,nstart=10)
iris <- data.frame(iris,Clusters=factor(kout3$cluster))
with(iris,table(Species,Clusters))</pre>
```

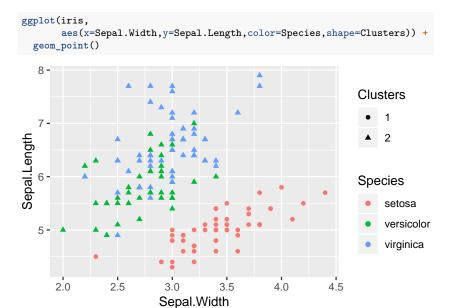
```
## Clusters
## Species 1 2 3
## setosa 0 0 50
## versicolor 39 11 0
## virginica 14 36 0
```



Sepal.Width

```
kout2 <- kmeans(irisX,centers=2,nstart=10)
iris$Clusters <- factor(kout2$cluster)
with(iris,table(Species,Clusters))</pre>
### Clusters
```

Species 1 2 ## setosa 50 0 ## versicolor 0 50 ## virginica 0 50



Scaling Variables

► Whether or not to scale variables depends on the context, but usually we will.

Choosing K

- ▶ Like PCA, there is no "best" method for choosing *K*.
- Cross-validation is not an option because there is no outcome.
- ► For small *p*, can visualize the clusters, but this becomes difficult as *p* grows.
- The silhouette plot is a graphical approach.
 - Discussed after the PAM algorithm below.

Sensitivity to Outliers

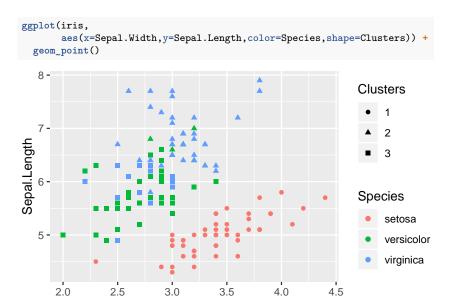
- Means and Euclidean distances are sensitive to outliers and K-means depends on both
 - Means are cluster centroids.
 - Criterion to minimize is a sum of squared Euclidean distances.

K-Medoids Clustering

- ▶ An alternative to *K*-means is *K*-medoids clustering.
- ► Cluster centres are medoids, which are observations chosen to represent each cluster.
- ► There is flexibility in choosing the dissimilarity measure
 - Can choose a more robust measure, such as ℓ_1 (so-called Manhattan) distance.
- ▶ An implementation of *K*-medoids is the PAM (partitioning around medoids) algorithm.

PAM on the Iris Data

```
library(cluster)
pout3 <- pam(irisX,k=3)
iris$Clusters <- factor(pout3$cluster)
with(iris,table(Species,Clusters))</pre>
```



Sepal.Width

```
pout2 <- pam(irisX,k=2)
iris$Clusters <- factor(pout2$cluster)
with(iris,table(Species,Clusters))

## Clusters
## Species 1 2</pre>
```

##

##

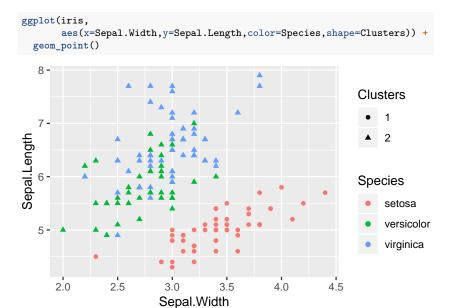
setosa

versicolor 0 50

virginica

50 0

0 50



The Silhouette Plot

- For each observation we compute a measure of cluster certainty, called the silhouette width, that takes values in [-1, 1].
 - ► Values near 1 indicate certainty that the observation is in the right cluster and
 - ▶ Values near −1 indicate that the observation is in the wrong cluster.
- Formula for silhouette width is given in on the next slide.
- ▶ If the silhouette values in a cluster are, say, below average, this suggests the cluster can be merged with another.

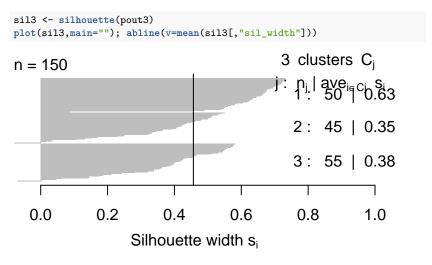
Silhouette Widths

- For observation i in cluster k, the silhouette width s(i) is defined as follows.
- Let a(i) be the average dissimilarity between i and all other observations in **cluster** k.
- For any other cluster C (i not in C), let d(i, C) be the average dissimilarity between i and all observations in cluster C.
- Let $b(i) = \min_C d(i, C)$ (minimum over all C such that i not in C).
- ► Then

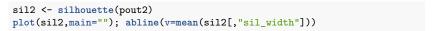
$$s(i) = \frac{b(i) - a(i)}{\max(a(i), b(i))}$$

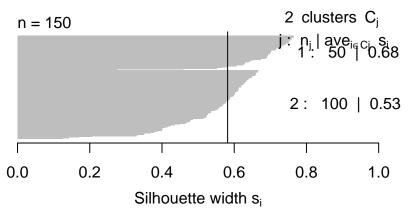
Note: a(i) is a dissimilarity measure, and should be $\ll b(i)$ if i confidently in cluster k.

Silhouette Plots for the Iris Data

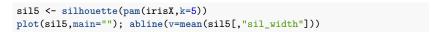


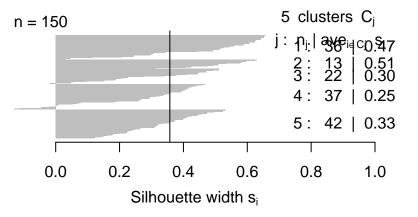
Average silhouette width: 0.46





Average silhouette width: 0.58





Average silhouette width: 0.36

ightharpoonup Had to go to about K=15 before silhouettes suggested merging clusters (not shown).

Clustering with categorical data.

Use the HUI data

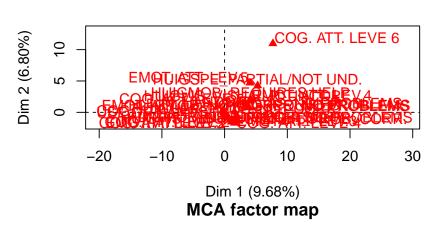
[1] "HUIDCOG" "HUIGDEX" "HUIDEMO" "HUIGHER" "HUIGMOB" "HUIGSPE" "HUIGVIS"

Extract PCs from MCA

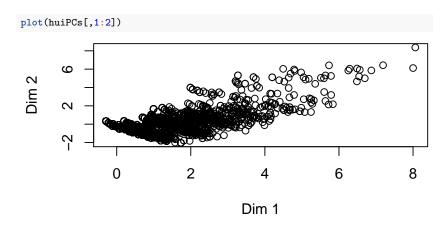
library(FactoMineR)

res.mca <- MCA(hsub,row.w = hui\$WTS_M)#huiPCs <- res.mca\$ind\$coord

MCA factor map







Cluster on PCs

▶ Will work with a small subset (first 1000 people) to keep computation and overplotting down.

```
n <- 1000
hk <- kmeans(huiPCs,centers=4,nstart=10)
# plot for a small sample of the dataset
huiPCs <- data.frame(huiPCs[1:n,])
huiPCs$Clusters <- factor(hk$cluster[1:n])
huiPCs$age <- hui$DHHGAGE[1:n]
huiPCs$sex <- hui$DHH_SEX[1:n]</pre>
```

```
library(ggplot2)
ggplot(huiPCs,
        aes(x=Dim.1,y=Dim.2,shape=Clusters)) +
  geom_point(alpha=.4)
    4 -
                                                                    Clusters
Dim.2
                                             ш
                                                                         2
                                                                          3
                                                 ш
   -2 -
                                                             5
```

Dim.1

