Untitled

Problem 7.1

Function helping to make right choice about number of K

```
wssplot <- function(data, nc=15, seed=1234){
            wss <- (nrow(data)-1)*sum(apply(data,2,var))
            for (i in 2:nc){
                set.seed(seed)
                wss[i] <- sum(kmeans(data, centers=i)$withinss)}
            plot(1:nc, wss, type="b", xlab="Number of Clusters",
                ylab="Within groups sum of squares")}
```
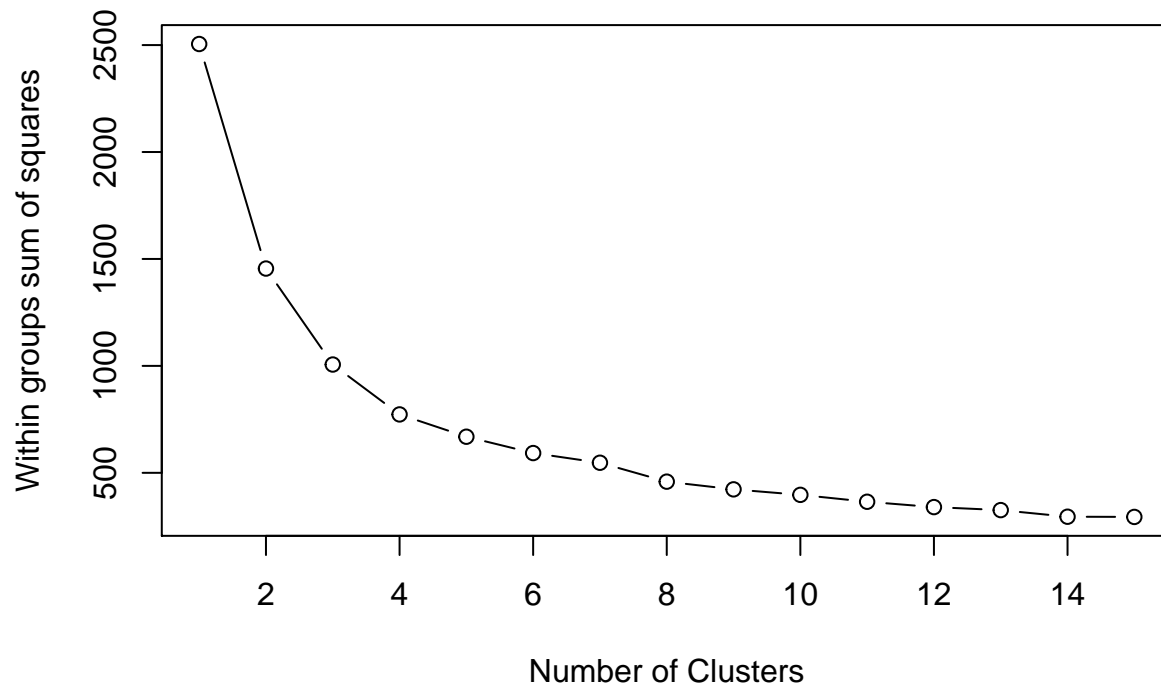
Data that we want to use for clusterization

```
kmeans.data <- data[4:6]
head(kmeans.data)
```

```
##     Gm Age Salary
## 1   74  28   9.50
## 2   77  29   9.25
## 3   83  29   9.00
## 4   93  29   8.75
## 5   81  29   8.75
## 6  108  24   8.00
```

Scaling the data

```
kmeans.data.scaled <- scale(kmeans.data)
wssplot(kmeans.data.scaled)
```

Let's do some clusterization

3 clusters

```r
# K-Means Cluster Analysis
fit3 <- kmeans(kmeans.data.scaled, 3)
# get cluster means
aggregate(kmeans.data.scaled,by=list(fit3$cluster),FUN=mean)
```

```
##   Group.1         Gm        Age     Salary
## 1       1 -1.2750839 -0.4807131 -0.6641629
## 2       2  0.6240421  0.9326274  1.1868775
## 3       3  0.5830986 -0.3564442 -0.4084625
```
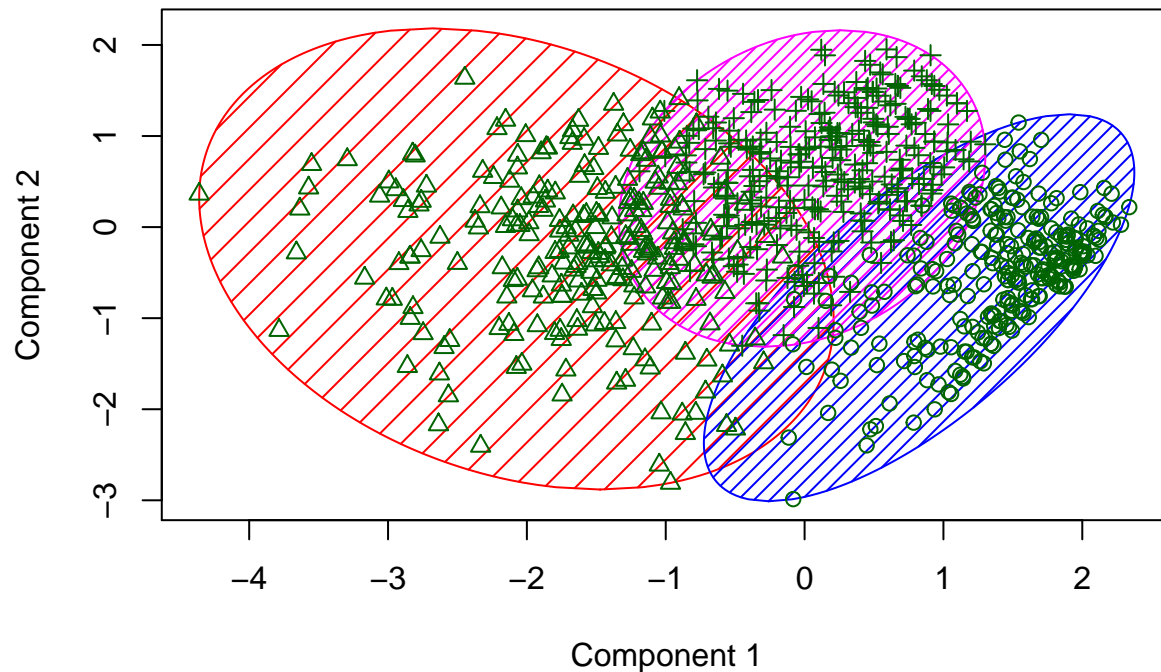
```r
# append cluster assignment
fit3$tot.withinss #total within-cluster sum of squares
```

```
## [1] 1006.333
```

Cluster Plot against 1st 2 principal components

```r
# vary parameters for most readable graph
library(cluster)
clusplot(kmeans.data.scaled, fit3$cluster, color=TRUE, shade=TRUE)
```

CLUSPLOT(kmeans.data.scaled)

Component 1

These two components explain 84.33 % of the point variability.

4 clusters

```
# K-Means Cluster Analysis
fit4 <- kmeans(kmeans.data.scaled, 4)
# get cluster means
aggregate(kmeans.data.scaled,by=list(fit4$cluster),FUN=mean)
```

```
##   Group.1        Gm        Age      Salary
## 1       1 -1.2910727 -0.5847154 -0.67308221
## 2       2  0.2674915  1.3058219  0.08614001
## 3       3  0.6108817 -0.5813090 -0.38003721
## 4       4  0.7990348  0.5224068  1.73816663
```
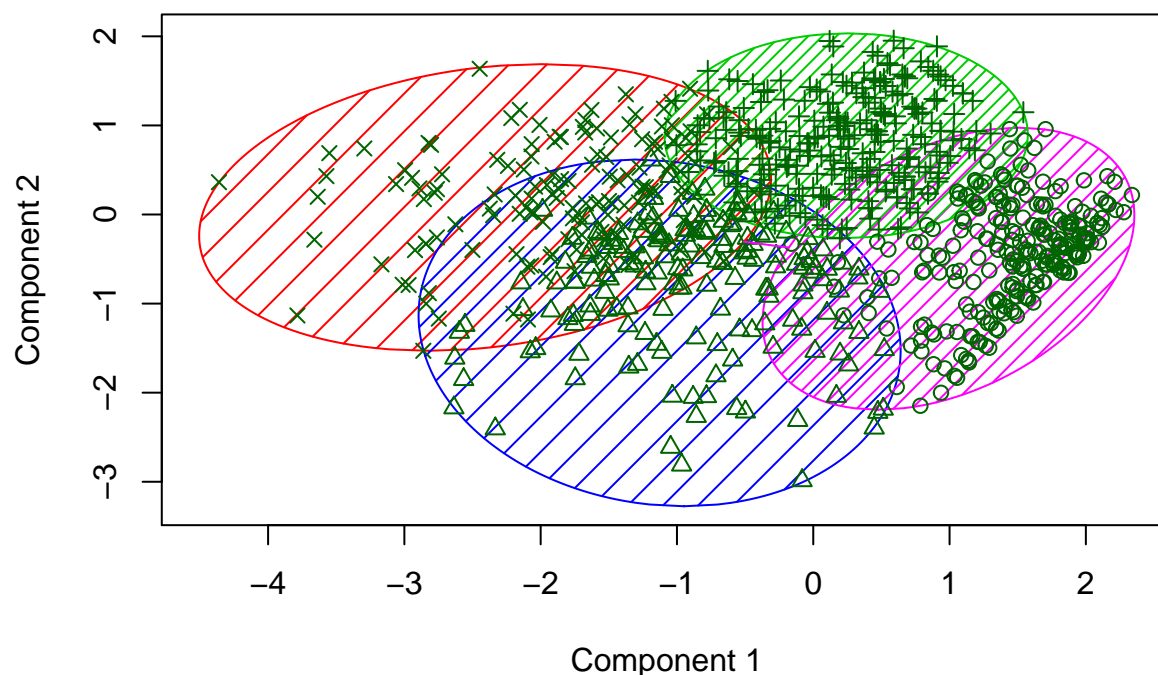
```
# append cluster assignment
fit4$tot.withinss #total within-cluster sum of squares
```

```
## [1] 772.9491
```

Cluster Plot against 1st 2 principal components

```
# vary parameters for most readable graph
clusplot(kmeans.data.scaled, fit4$cluster, color=TRUE, shade=TRUE)
```

CLUSPLOT(kmeans.data.scaled)

Component 1
These two components explain 84.33 % of the point variability.

7 clusters

```r
# K-Means Cluster Analysis
fit7 <- kmeans(kmeans.data.scaled, 7)
# get cluster means
aggregate(kmeans.data.scaled,by=list(fit7$cluster),FUN=mean)
```

```
##   Group.1         Gm          Age     Salary
## 1       1  0.7198899  0.08273893  1.0259907
## 2       2  0.9529056  0.75135856  2.7388054
## 3       3  0.5314117  1.54295287  0.7870524
## 4       4 -1.5100957 -0.63245174 -0.6874792
## 5       5  0.8137360 -0.63694294 -0.4563627
## 6       6 -0.3322552 -0.77335084 -0.6681983
## 7       7 -0.1372308  0.99114365 -0.4660662
```
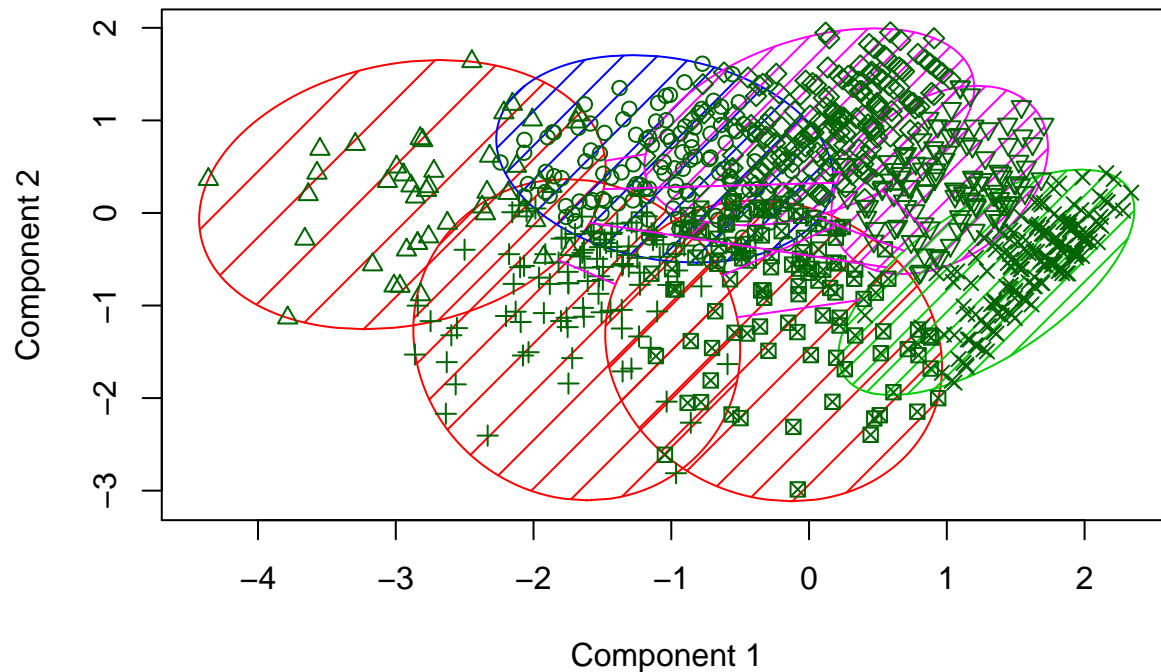
```r
# append cluster assignment
fit7$tot.withinss #total within-cluster sum of squares
```

```
## [1] 545.4409
```

Cluster Plot against 1st 2 principal components

```r
# vary parameters for most readable graph
clusplot(kmeans.data.scaled, fit7$cluster, color=TRUE, shade=TRUE)
```

CLUSPLOT(kmeans.data.scaled)

Component 1
These two components explain 84.33 % of the point variability.

comparing 2 cluster solutions 3-means and 4-means

```
library(fpc)
d <- dist(kmeans.data.scaled)
cluster.stats(d, fit3$cluster, fit4$cluster, compareonly = TRUE)
```

```
## $corrected.rand
## [1] 0.7083329
##
## $vi
## [1] 0.6987167
```

comparing 2 cluster solutions 4-means and 7-means

```
library(fpc)
d <- dist(kmeans.data.scaled)
cluster.stats(d, fit4$cluster, fit7$cluster, compareonly = TRUE)
```

```
## $corrected.rand
## [1] 0.5015822
##
## $vi
## [1] 1.306123
```

The correct choice of k is often ambiguous, with interpretations depending on the shape and scale of the distribution of points in a data set and the desired clustering resolution of the user. In addition, increasing k without penalty will always reduce the amount of error in the

resulting clustering, to the extreme case of zero error if each data point is considered its own cluster (i.e., when k equals the number of data points, n). Intuitively then, the optimal choice of k will strike a balance between maximum compression of the data using a single cluster, and maximum accuracy by assigning each data point to its own cluster.

So, in our case we see that the total within-cluster sum of squares is redusing as k is getting bigger.

Problem 7.2

Build a minimum spanning tree connecting all points and delete 6 maximum edges step by step.

```
stdData <- kmeans.data.scaled
dis <- dist(stdData) #dissimilarity data
tr <- spantree(dis, toolong = 0) #build an MST
plot(tr, cmdscale(dis)) #plot an MST
```
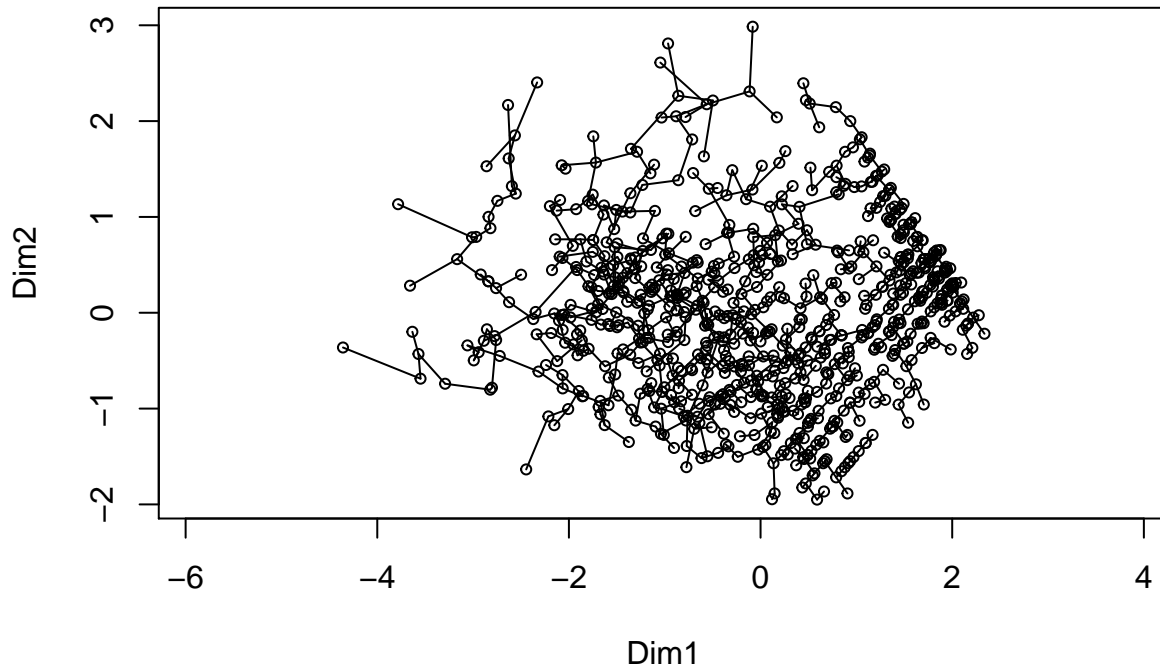


```
for(i in 1:6) #loop for finding 6 maximum edges in MST
{
  m <- which.max(tr$dist)
  tr$dist[m] <- NA
}
m <- which.max(tr$dist)
tres <- tr$dist[m] + 0.00001 #set a treshold
j <- distconnected(dis, toolong = tres) #find connected components in MST after deleting 6 edges
```

```
## Connectivity of distance matrix with threshold dissimilarity 0.758652
## Data are disconnected: 7 groups
## Groups sizes
##    1   2   3   4   5   6   7
## 828   1   3   1   1   1   1
```

```
clusplot(stdData, j, color=TRUE, shade=TRUE) #plot clusters
```
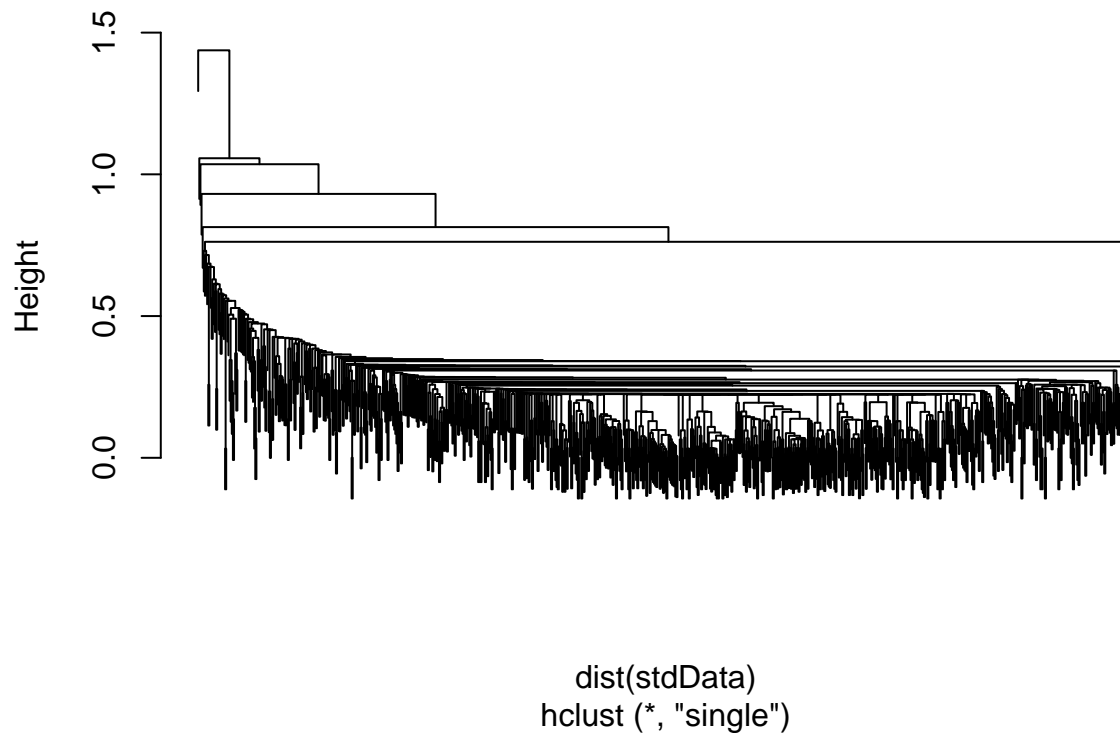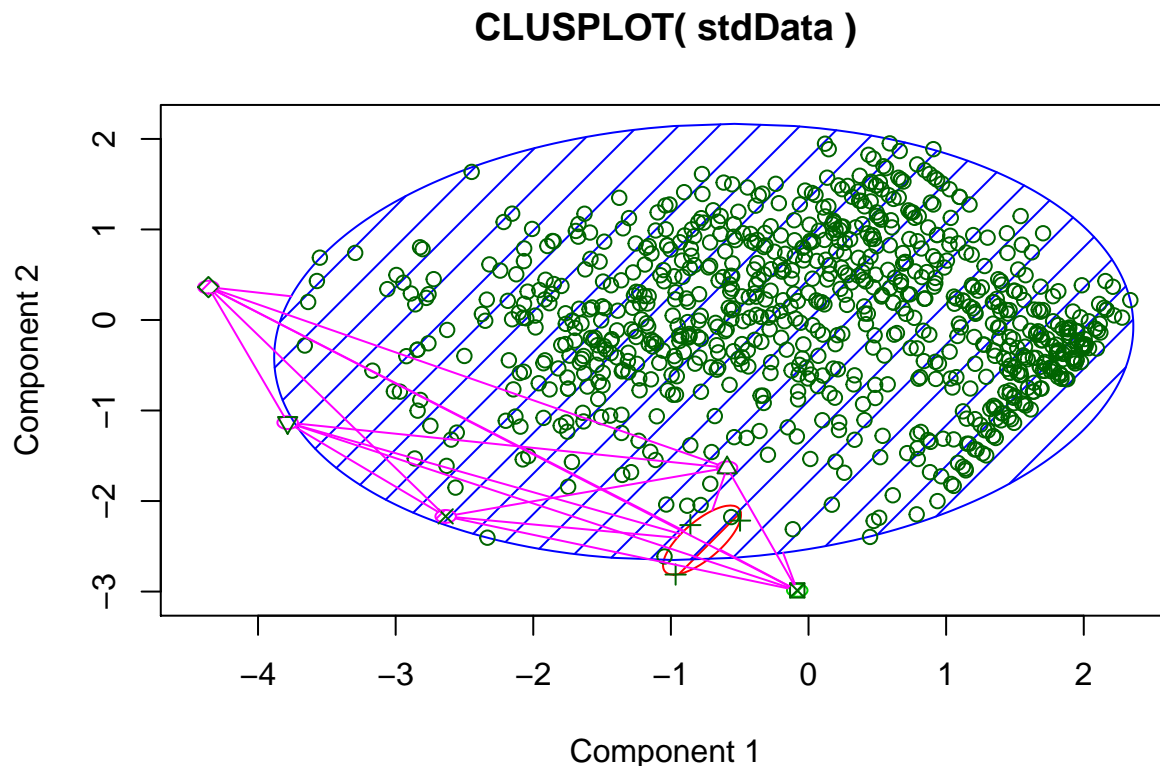
7

CLUSPLOT(stdData)
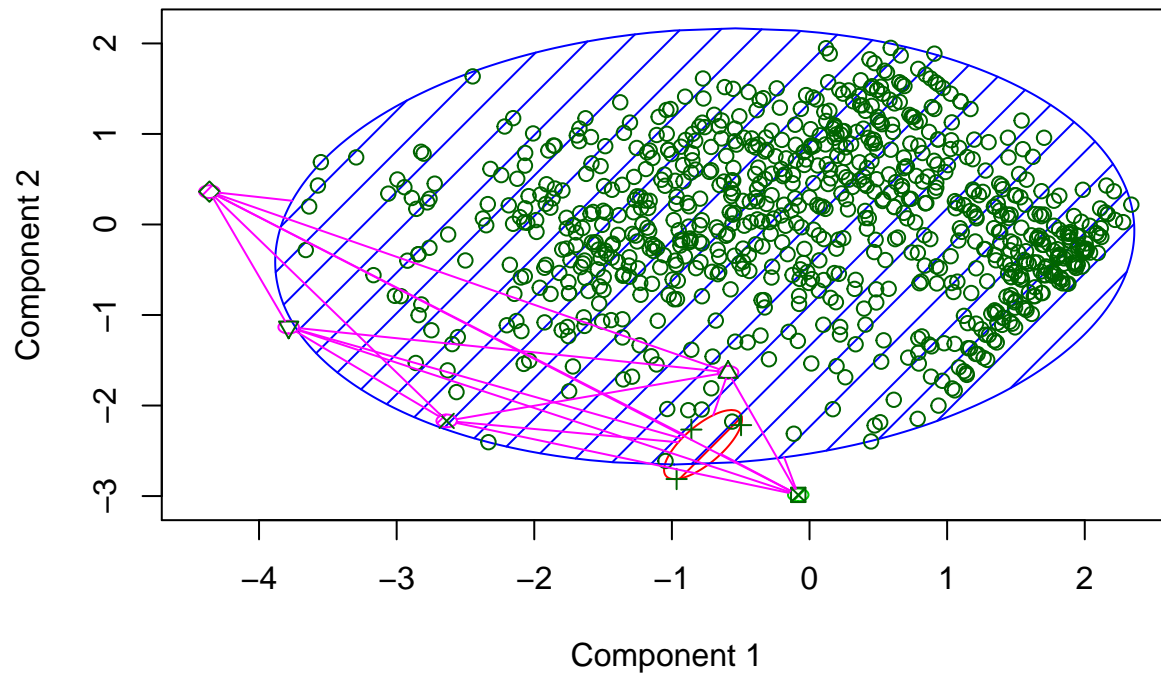
Component 1
These two components explain 84.33 % of the point variability.

```r
hc <- hclust(dist(stdData), method = "single")
plot(hc, labels = FALSE)
```

Cluster Dendrogram

dist(stdData)
hclust (*, "single")

```
memb <- cutree(hc, k = 7)
clusplot(stdData, memb, color=TRUE, shade=TRUE)
```

CLUSPLOT(stdData)

Component 1
These two components explain 84.33 % of the point variability.

Consider a set of 2D points presented. Those in problem 2 clustered by using the single link approach, whereas those in problem 1, by using the square error criterion of K-Means.

Overall, this example demonstrates the major difference between conventional clustering and single link clustering: the latter finds elongated structures whereas the former cuts out convex parts. Sometimes, especially in the analysis of results of physical processes or experiments over real-world particles, the elongated structures do capture the essence of the data and are of great interest. In other cases, especially when entities/features have no intuitive geometric meaning - think of hockey players, for example, convex clusters make much more sense as groupings around their centroids.

By using MST (with single linkage) we get 1 big clusters and 6 small cluster consisting of 1/3 player/s. This is probably can be players that differs very much from other players and doesn't get to any big cluster. However, K-means placed these player to some more general cluster, because in this case clusters are more convex. The shape of such clusters lets the objects (that are standing far away from others in terms of features' distances) to get into them.