

Unsupervised_learning-HW

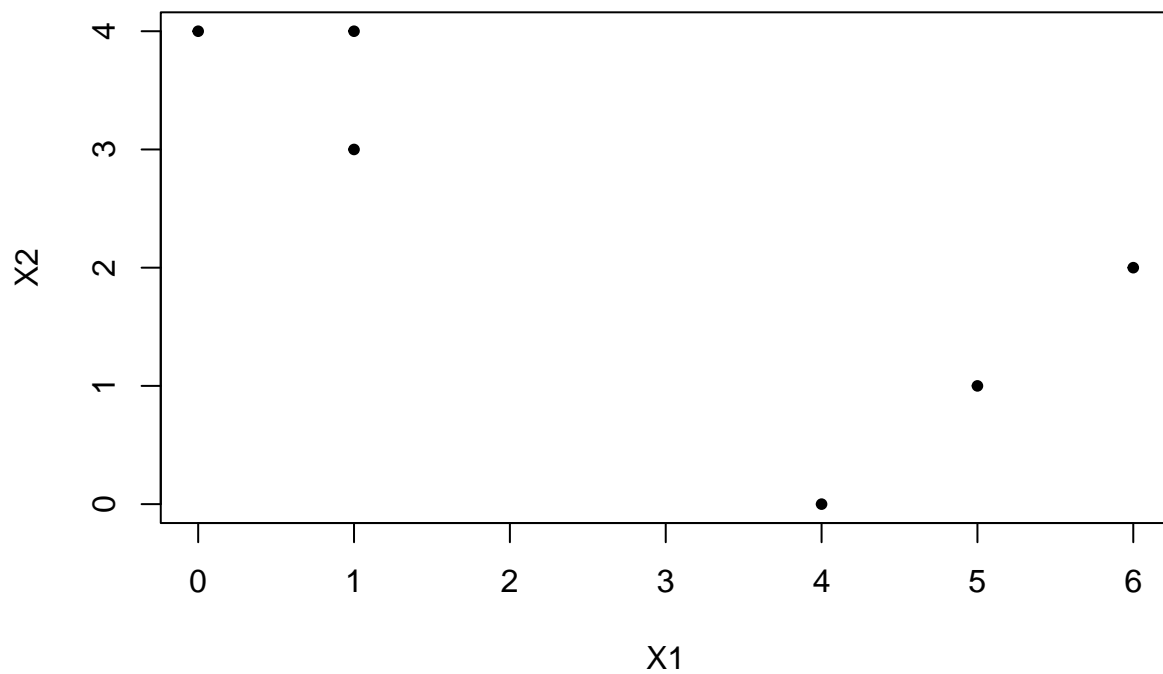
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3/30/2022

12.3

(a)

```
X1 <- c(1, 1, 0, 5, 6, 4)
X2 <- c(4, 3, 4, 1, 2, 0)
dat <- cbind(X1, X2)
plot(dat, pch = 20)
```



(b)

```
set.seed(1108)
label <- sample(c(1, 2), 6, replace = TRUE)
label
```

```
## [1] 1 1 2 2 1 1
```

(c)

```
center1 <- apply(dat[which(label == 1), ], 2, mean)
center2 <- apply(dat[which(label == 2), ], 2, mean)
centroid <- rbind(center1, center2)
```

(d)

```
dist <- Euclidean(vec1 = dat, vec2 = centroid)
label <- ifelse(dist[, 1] > dist[, 2], 1, 2)
```

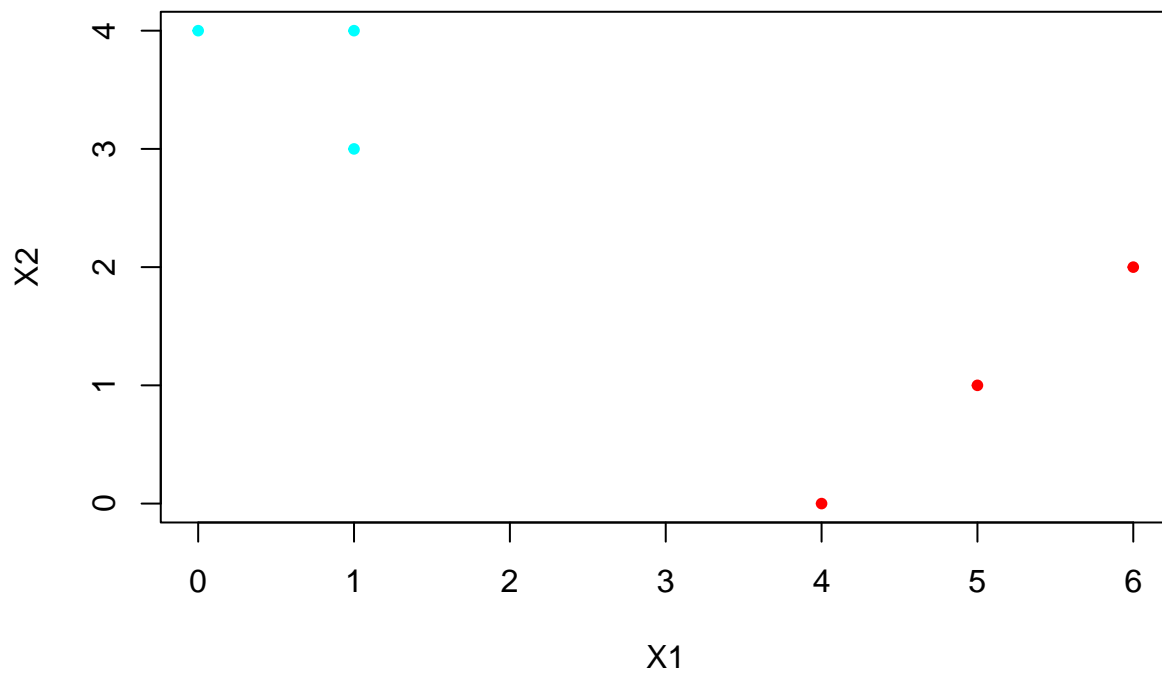
(e)

```
set.seed(1108)
label <- sample(c(1, 2), 6, replace = TRUE)
iter = 10
for (i in 1:iter) {
  cat("iteration", i, ": ", label, "\n")
  centroid <- rbind(apply(dat[which(label == 1), ], 2, mean),
    apply(dat[which(label == 2), ], 2, mean))
  dist <- Euclidean(vec1 = dat, vec2 = centroid)
  label <- ifelse(dist[, 1] > dist[, 2], 2, 1)
}
```

```
## iteration 1 : 1 1 2 2 1 1
## iteration 2 : 2 2 2 1 1 1
## iteration 3 : 2 2 2 1 1 1
## iteration 4 : 2 2 2 1 1 1
## iteration 5 : 2 2 2 1 1 1
## iteration 6 : 2 2 2 1 1 1
## iteration 7 : 2 2 2 1 1 1
## iteration 8 : 2 2 2 1 1 1
## iteration 9 : 2 2 2 1 1 1
## iteration 10 : 2 2 2 1 1 1
```

(f)

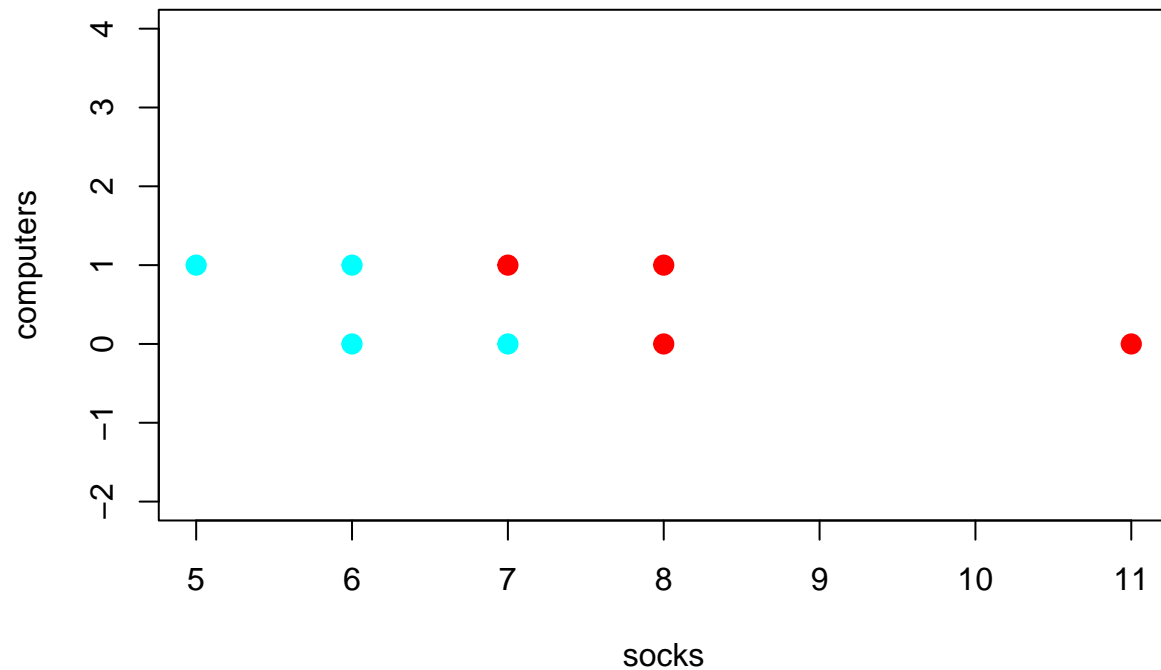
```
plot(dat, pch = 20, col = Cols(label),
      xlab = "X1", ylab = "X2")
```



12.5

(1)

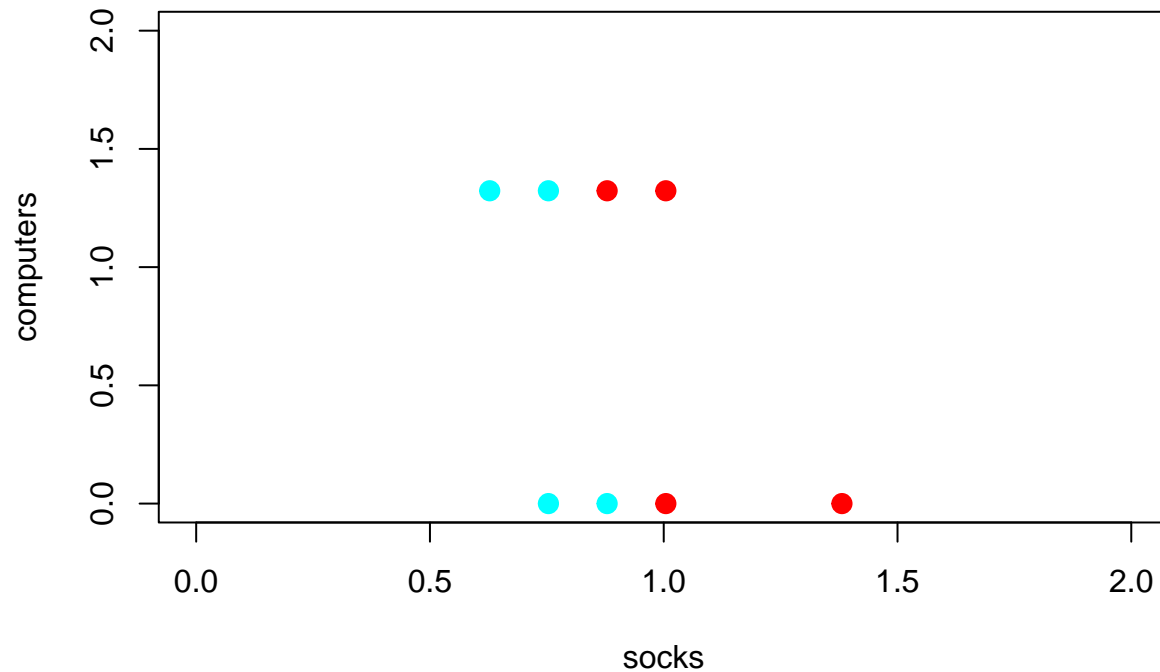
```
socks <- c(8, 11, 7, 6, 5, 6, 7, 8)
computers <- c(0, 0, 0, 0, 1, 1, 1, 1)
dat <- cbind(socks, computers)
labels <- c(1, 1, 2, 2, 2, 2, 1, 1)
plot(dat, col = Cols(labels), pch = 20, cex = 2,
      xlab = "socks", ylab = "computers", xlim = c(5, 11), ylim = c(-2, 4))
```



If take the original data into consideration, the number of socks performs to be a more important factor when clustering the data. I expect to see one cluster with less socks (obs: 3, 4, 5, 6) and another with more socks (obs: 1, 2, 7, 8)

(2)

```
socks      <- scale(c(8, 11, 7, 6, 5, 6, 7, 8), center = FALSE)
computers  <- scale(c(0, 0, 0, 0, 1, 1, 1, 1) , center = FALSE)
dat <- cbind(socks, computers)
labels <- c(1, 1, 2, 2, 2, 2, 1, 1)
plot(dat, col = Cols(labels), pch = 20, cex = 2,
      xlab = "socks", ylab = "computers", xlim = c(0, 2), ylim = c(0, 2))
```



If take the centered data without scale, we prefer to cluster by the number of computers – group 1: obs 1, 2, 3, 4; group 2: obs 5, 6, 7, 8.

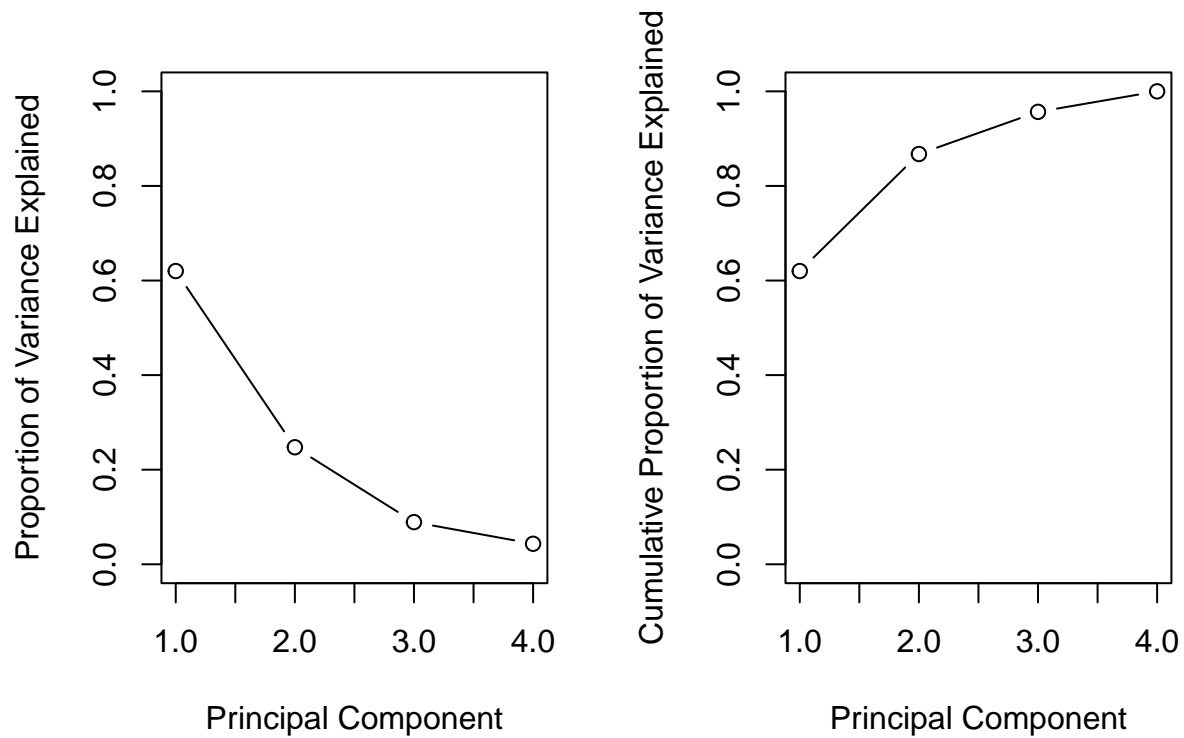
(3)

obviously, if we do the clustering by the number of dollars spent, the factor socks will be ignored and the clustering totally depends on the factor computer.

12.8

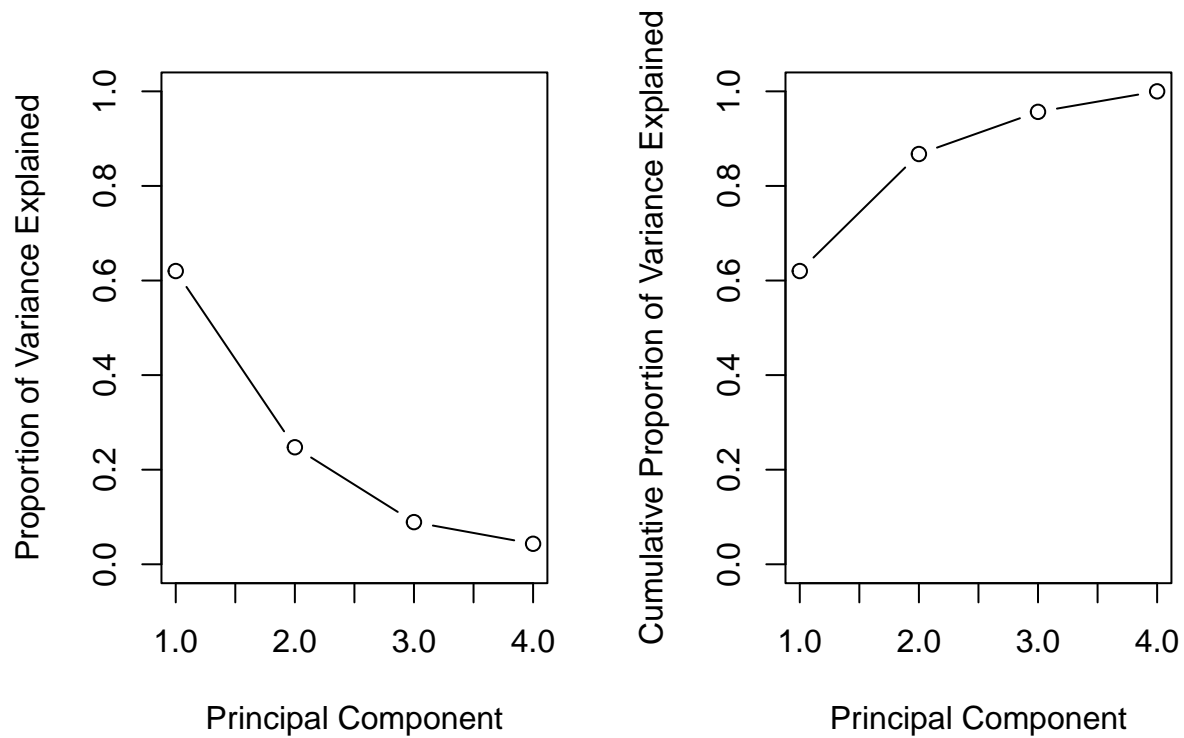
(a)

```
pr.out <- prcomp(USArrests, scale = TRUE)
pr.var <- pr.out$sdev^2
pve <- pr.var / sum(pr.var)
par(mfrow = c(1, 2))
plot(pve, xlab = "Principal Component",
     ylab = "Proportion of Variance Explained", ylim = c(0, 1),
     type = "b")
plot(cumsum(pve), xlab = "Principal Component",
     ylab = "Cumulative Proportion of Variance Explained", ylim = c(0, 1), type = "b")
```



(b)

```
dat <- scale(USArrests)
# apply(dat, 2, mean)
# apply(dat, 2, var)
z_score <- dat %*% pr.out$rotation
pve <- apply(z_score, 2, var)/sum(apply(dat, 2, var)) # eq 12.10
# sum(pve)
par(mfrow = c(1, 2))
plot(pve, xlab = "Principal Component",
     ylab = "Proportion of Variance Explained", ylim = c(0, 1),
     type = "b")
plot(cumsum(pve), xlab = "Principal Component",
     ylab = "Cumulative Proportion of Variance Explained", ylim = c(0, 1), type = "b")
```



12.9

(a)

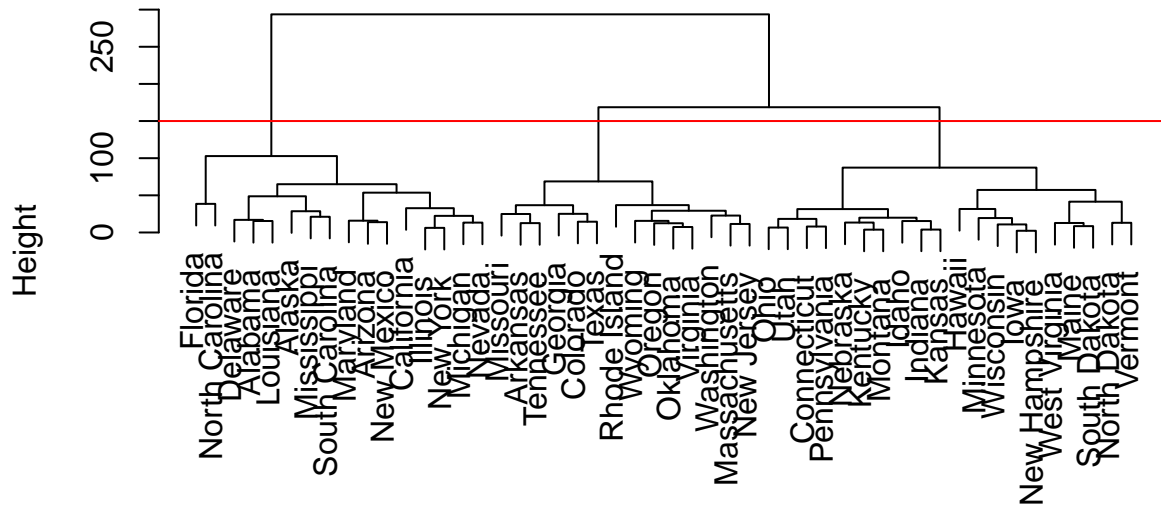
```
hc.complete <- hclust(dist(USArrests), method = "complete") # Euclidean distance, complete linkage
```

(b)

Cut at the height 150, and the clustering results are as follow:

```
hc.cluster <- cutree(hc.complete, 3)
plot(hc.complete)
abline(h = 150, col = "red")
```

Cluster Dendrogram



dist(USArrests)
hclust (*, "complete")

hc.cluster

##	Alabama	Alaska	Arizona	Arkansas	California
##	1	1	1	2	1
##	Colorado	Connecticut	Delaware	Florida	Georgia
##	2	3	1	1	2
##	Hawaii	Idaho	Illinois	Indiana	Iowa
##	3	3	1	3	3
##	Kansas	Kentucky	Louisiana	Maine	Maryland
##	3	3	1	3	1
##	Massachusetts	Michigan	Minnesota	Mississippi	Missouri
##	2	1	3	1	2
##	Montana	Nebraska	Nevada	New Hampshire	New Jersey
##	3	3	1	3	2
##	New Mexico	New York	North Carolina	North Dakota	Ohio
##	1	1	1	3	3
##	Oklahoma	Oregon	Pennsylvania	Rhode Island	South Carolina
##	2	2	3	2	1
##	South Dakota	Tennessee	Texas	Utah	Vermont
##	3	2	2	3	3
##	Virginia	Washington	West Virginia	Wisconsin	Wyoming
##	2	2	3	3	2

(c)


```
hc.complete.scaled <- hclust(dist(dat), method = "complete") # centered and scaled data, Euclidean dist
```

(d)

```
table(cutree(hc.complete, 3), cutree(hc.complete.scaled, 3))
```

```
##
##      1  2  3
##    1  6  9  1
##    2  2  2 10
##    3  0  0 20
```

after scaling the data, complete linkage hierarchical clustering assigns more obs to 3rd cluster. To my consideration, the variables should be scaled beforehand because variables are in different units.

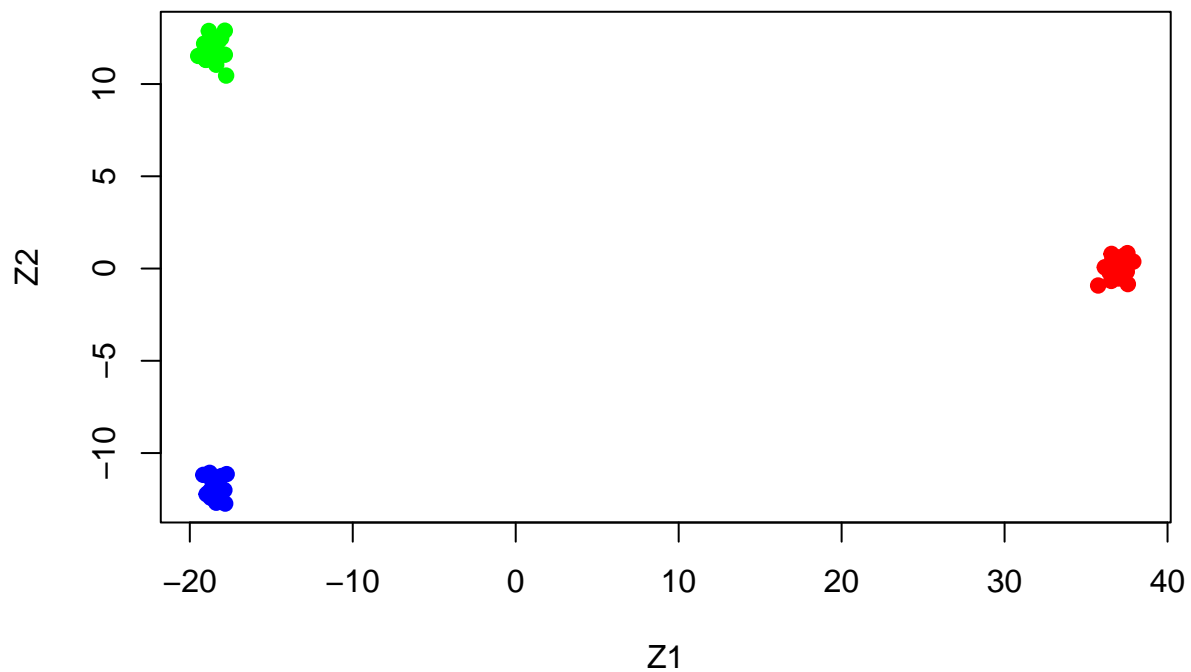
12.10

(a)

```
set.seed(1754)
dat <- matrix(rnorm(20*3*50, sd = .5), ncol = 50)
dat[1:20, 1] <- dat[1:20, 1] + 2
dat[1:20, 2] <- dat[1:20, 2] + 13
dat[21:40, 1] <- dat[21:40, 1] + 22
# dat[21:40, 2] <- dat[21:40, 2] + 33
dat[41:60, 1] <- dat[41:60, 1] + 42
dat[41:60, 2] <- dat[41:60, 2] + 53
labels <- c(rep(3, 20), rep(2, 20), rep(1, 20))
```

(b)

```
pr.out <- prcomp(dat)
plot(pr.out$x[, 1:2], col = Cols(labels),
     xlab = "Z1", ylab = "Z2", pch = 19)
```



(c)

```
set.seed(2)
km.out <- kmeans(dat, 3, nstart = 20)
table(labels, km.out$cluster)
```

```
##
## labels  1  2  3
##      1 20  0  0
##      2  0 20  0
##      3  0  0 20
```

the observations are perfectly clustered

(d)

```
set.seed(2)
km.out <- kmeans(dat, 2, nstart = 20)
table(labels, km.out$cluster)
```

```
##
## labels  1  2
##      1 20  0
##      2  0 20
##      3  0 20
```

(e)

```
set.seed(2)
km.out <- kmeans(dat, 4, nstart = 20)
table(labels, km.out$cluster)
```

```
##
## labels  1  2  3  4
##        1  0  0  0 20
##        2 20  0  0  0
##        3  0 12  8  0
```

(f)

```
set.seed(2)
km.out <- kmeans(pr.out$x[, 1:2], 3, nstart = 20)
table(labels, km.out$cluster)
```

```
##
## labels  1  2  3
##        1 20  0  0
##        2  0 20  0
##        3  0  0 20
```

still perfectly clustered

(g)

```
scales.dat <- scale(dat)
set.seed(2)
km.out <- kmeans(scales.dat, 3, nstart = 20)
table(labels, km.out$cluster)
```

```
##
## labels  1  2  3
##        1  0  6 14
##        2  5 14  1
##        3  8 11  1
```

the result is worse since differences are mitigated after scaling and hard to cluster.