

Assignment_4

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Due 3/20/2022

Following is the link to my GitHub account:

https://github.com/Kgardner22/64060_-kgardner

IMPORT AND PREPARE DATA:

Import the Pharmaceuticals.csv file

```
Pharmaceuticals <- read.table('C:/R/MyData/Pharmaceuticals.csv', header = T,
sep = ',')

summary(Pharmaceuticals)
```

##	Symbol	Name	Market_Cap	Beta	
##	Length:21	Length:21	Min. : 0.41	Min. :0.1800	
##	Class :character	Class :character	1st Qu.: 6.30	1st Qu.:0.3500	
##	Mode :character	Mode :character	Median : 48.19	Median :0.4600	
##			Mean : 57.65	Mean :0.5257	
##			3rd Qu.: 73.84	3rd Qu.:0.6500	
##			Max. :199.47	Max. :1.1100	
##	PE_Ratio	ROE	ROA	Asset_Turnover	Leverage
##	Min. : 3.60	Min. : 3.9	Min. : 1.40	Min. :0.3	Min. :0.0000
##	1st Qu.:18.90	1st Qu.:14.9	1st Qu.: 5.70	1st Qu.:0.6	1st Qu.:0.1600
##	Median :21.50	Median :22.6	Median :11.20	Median :0.6	Median :0.3400
##	Mean :25.46	Mean :25.8	Mean :10.51	Mean :0.7	Mean :0.5857
##	3rd Qu.:27.90	3rd Qu.:31.0	3rd Qu.:15.00	3rd Qu.:0.9	3rd Qu.:0.6000
##	Max. :82.50	Max. :62.9	Max. :20.30	Max. :1.1	Max. :3.5100

```
##      Rev_Growth      Net_Profit_Margin      Median_Recommendation      Location
##      Min.      : -3.17      Min.      : 2.6      Length:21      Length:21
##      1st Qu.: 6.38      1st Qu.:11.2      Class :character      Class :character
##      Median : 9.37      Median :16.1      Mode  :character      Mode  :character
##      Mean   :13.37      Mean   :15.7
##      3rd Qu.:21.87      3rd Qu.:21.1
##      Max.   :34.21      Max.   :25.5
##      Exchange
##      Length:21
##      Class :character
##      Mode  :character
##
##
##
```

Load required libraries

```
library(tidyverse) #for data manipulation
library(factoextra) #for clustering and visualization
library(flexclust)
```

Use cluster analysis to explore and analyze the given dataset as follows:

REQUIREMENTS A and B:

Use only the numerical variables (1 to 9) to cluster the 21 firms. Justify the various choices made in conducting the cluster analysis, such as weights for different variables, the specific clustering algorithm(s) used, the number of clusters formed, and so on.

First, we create the data.frame with only the numerical variables 1 to 9

```
set.seed(64060)

df <- Pharmaceuticals[,c(3:11)]

summary(df)

##      Market_Cap      Beta      PE_Ratio      ROE
##      Min.      : 0.41      Min.      :0.1800      Min.      : 3.60      Min.      : 3.9
##      1st Qu.: 6.30      1st Qu.:0.3500      1st Qu.:18.90      1st Qu.:14.9
##      Median : 48.19      Median :0.4600      Median :21.50      Median :22.6
##      Mean   : 57.65      Mean   :0.5257      Mean   :25.46      Mean   :25.8
##      3rd Qu.: 73.84      3rd Qu.:0.6500      3rd Qu.:27.90      3rd Qu.:31.0
##      Max.   :199.47      Max.   :1.1100      Max.   :82.50      Max.   :62.9
##      ROA      Asset_Turnover      Leverage      Rev_Growth
##      Min.      : 1.40      Min.      :0.3      Min.      :0.0000      Min.      : -3.17
##      1st Qu.: 5.70      1st Qu.:0.6      1st Qu.:0.1600      1st Qu.: 6.38
##      Median :11.20      Median :0.6      Median :0.3400      Median : 9.37
##      Mean   :10.51      Mean   :0.7      Mean   :0.5857      Mean   :13.37
##      3rd Qu.:15.00      3rd Qu.:0.9      3rd Qu.:0.6000      3rd Qu.:21.87
##      Max.   :20.30      Max.   :1.1      Max.   :3.5100      Max.   :34.21
```

```
## Net_Profit_Margin
## Min. : 2.6
## 1st Qu.:11.2
## Median :16.1
## Mean :15.7
## 3rd Qu.:21.1
## Max. :25.5
```

Before we can begin cluster analysis, we must first scale the data.

```
# Scaling the data frame (z-score)
```

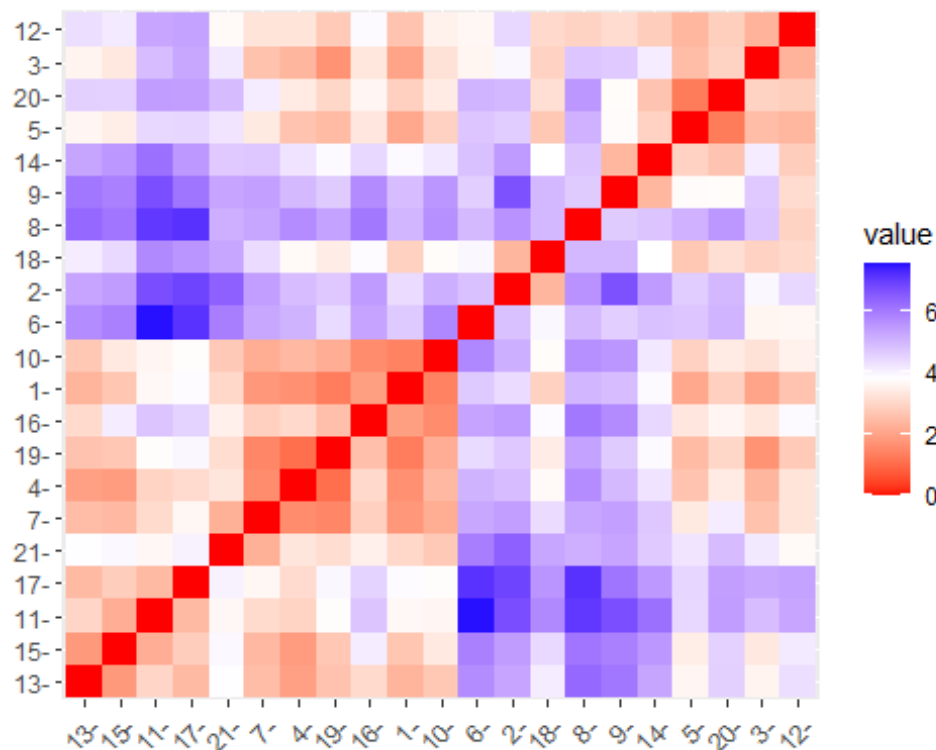
```
df <- scale(df)
```

```
summary(df)
```

```
##      Market_Cap      Beta      PE_Ratio      ROE
## Min.   :-0.9768  Min.   :-1.3466  Min.   :-1.3404  Min.   :-1.4515
## 1st Qu.:-0.8763  1st Qu.:-0.6844  1st Qu.:-0.4023  1st Qu.:-0.7223
## Median :-0.1614  Median :-0.2560  Median :-0.2429  Median :-0.2118
## Mean   : 0.0000  Mean   : 0.0000  Mean   : 0.0000  Mean   : 0.0000
## 3rd Qu.: 0.2762  3rd Qu.: 0.4841  3rd Qu.: 0.1495  3rd Qu.: 0.3450
## Max.    : 2.4200  Max.    : 2.2758  Max.    : 3.4971  Max.    : 2.4597
##      ROA      Asset_Turnover      Leverage      Rev_Growth
## Min.   :-1.7128  Min.   :-1.8451  Min.   :-0.74966  Min.   :-1.4971
## 1st Qu.:-0.9047  1st Qu.:-0.4613  1st Qu.:-0.54487  1st Qu.:-0.6328
## Median : 0.1289  Median :-0.4613  Median :-0.31449  Median :-0.3621
## Mean   : 0.0000  Mean   : 0.0000  Mean   : 0.00000  Mean   : 0.0000
## 3rd Qu.: 0.8430  3rd Qu.: 0.9225  3rd Qu.: 0.01828  3rd Qu.: 0.7693
## Max.    : 1.8389  Max.    : 1.8451  Max.    : 3.74280  Max.    : 1.8862
## Net_Profit_Margin
## Min.   :-1.99560
## 1st Qu.:-0.68504
## Median : 0.06168
## Mean   : 0.00000
## 3rd Qu.: 0.82364
## Max.    : 1.49416
```

We'll compute and visualize the distance matrix between rows using `get_dist()` and `fviz_dist()`

```
distance <- get_dist(df)
fviz_dist(distance)
```



The above graph

shows the distance between firms.

I'll run the k-means algorithm to cluster the firms, choosing an initial random value of $k = 4$.

```
set.seed(64060)
k4 <- kmeans(df, centers = 4, nstart = 25) # k = 4, number of restarts = 25

# the following will help us Visualize the output

k4$centers # output the centers

##      Market_Cap      Beta  PE_Ratio      ROE      ROA Asset_Turnover
## 1 -0.03142211 -0.4360989 -0.3172485  0.1950459  0.4083915  1.729746e-01
## 2  1.69558112 -0.1780563 -0.1984582  1.2349879  1.3503431  1.153164e+00
## 3 -0.82617719  0.4775991 -0.3696184 -0.5631589 -0.8514589 -9.994088e-01
## 4 -0.52462814  0.4451409  1.8498439 -1.0404550 -1.1865838  1.480297e-16
##      Leverage Rev_Growth Net_Profit_Margin
## 1 -0.2744931 -0.7041516      0.5569544
## 2 -0.4680782  0.4671788      0.5912425
## 3  0.8502201  0.9158889     -0.3319956
## 4 -0.3443544 -0.5769454     -1.6095439

k4$size # Number of firms in each cluster

## [1] 8 4 6 3

fviz_cluster(k4, data = df) # Visualize the output
```



This produces

similar size clusters (8,4,6,3)

Other Distances

I'll rerun the example using other distance measures to compare the results

```
set.seed(64060)

k4M = kcca(df, k=4, kccaFamily("kmedians")) #kmedians uses Manhattan
distance
k4M

## kcca object of family 'kmedians'
##
## call:
## kcca(x = df, k = 4, family = kccaFamily("kmedians"))
##
## cluster sizes:
##
##  1  2  3  4
##  3 12  5  1

k4E = kcca(df, k=4, kccaFamily("kmeans")) #kmeans uses Euclidean distance
k4E

## kcca object of family 'kmeans'
##
```

```
## call:
## kcca(x = df, k = 4, family = kccaFamily("kmeans"))
##
## cluster sizes:
##
##  1  2  3  4
## 12  4  2  3

k4A = kcca(df, k=4, kccaFamily("angle")) #angle uses angle between
observation and centroid
k4A

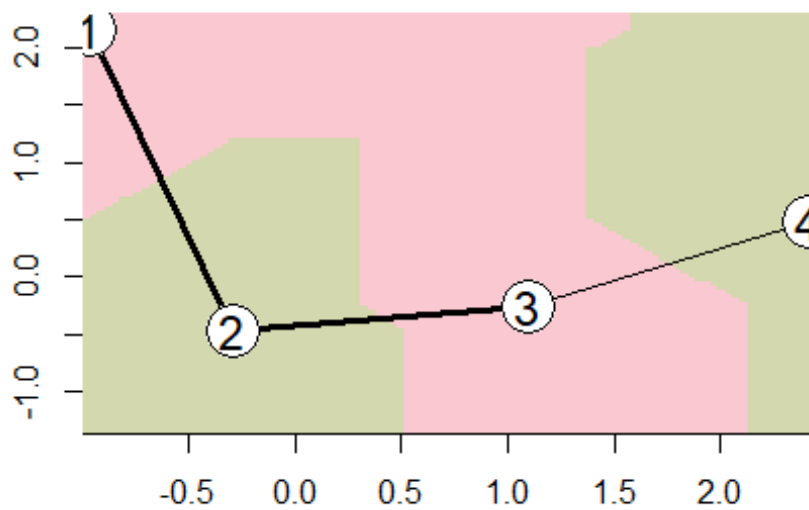
## kcca object of family 'angle'
##
## call:
## kcca(x = df, k = 4, family = kccaFamily("angle"))
##
## cluster sizes:
##
##  1  2  3  4
##  4  4 11  2

# We won't use Jaccard distance as this is primarily used for categorical
data which is not applicable.
```

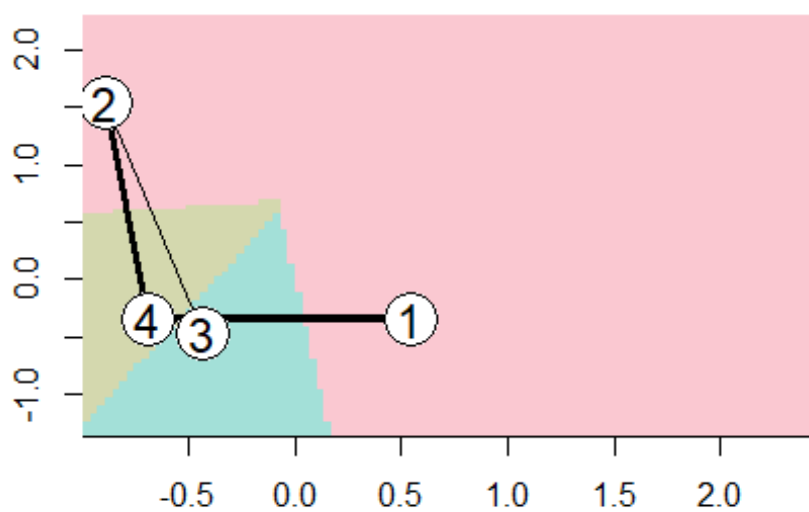
In reviewing these results, we see the cluster sizes using kmedians (Manhattan), kmeans (Euclidean) and angle produce similar results.

Let's take a look at the images of each of these results:

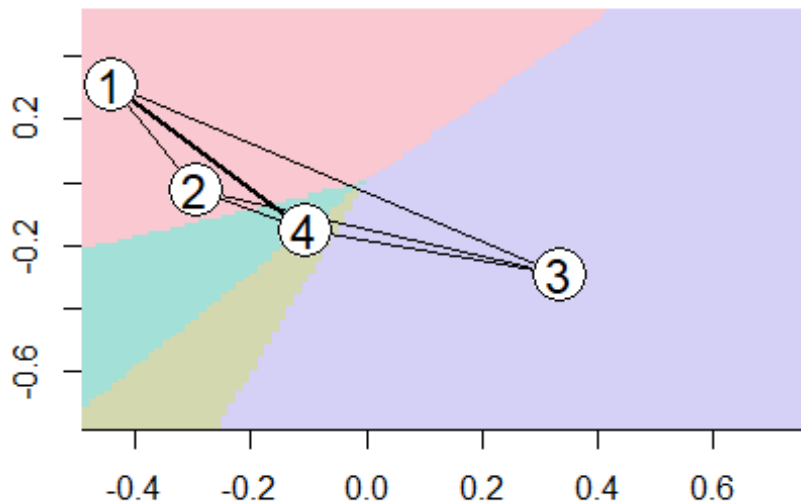
```
image(k4M) # Manhattan distance
```



```
image(k4E) # Euclidean distance
```



```
image(k4A) # angle
```



These images clearly show the various distance measures make a huge difference in the clustering results. It's clear that k4M, which is using Manhattan distance (kmedians), produces better results since the distance between the centroids is maximized in comparison to the other results.

This is also confirmed by looking at the centers:

```
dist(k4M@centers)

##           1           2           3
## 2 3.838654
## 3 5.233264 2.754097
## 4 6.025978 4.677079 2.397111

dist(k4E@centers)

##           1           2           3
## 2 4.169374
## 3 4.306065 4.093207
## 4 3.319190 3.018190 3.847236

dist(k4A@centers)

##           1           2           3
## 2 1.253136
## 3 1.872258 1.823885
## 4 1.158538 1.373808 1.743565
```

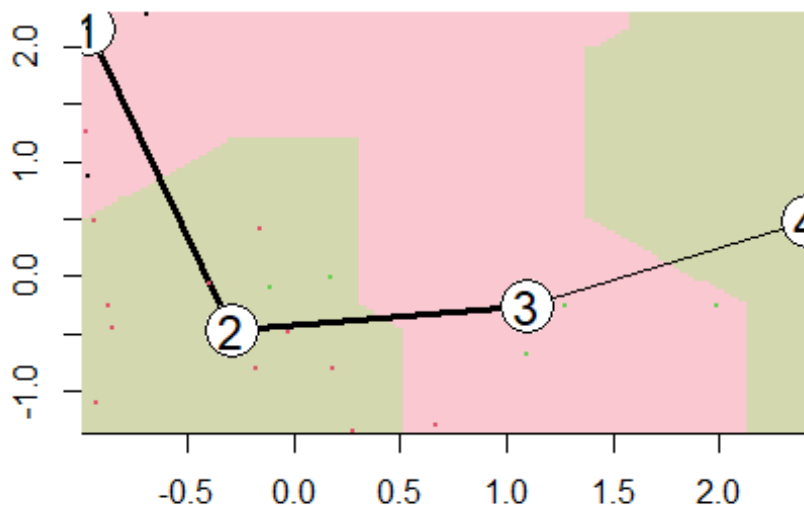

As shown in this data, the distance measure producing the maximum distance between centroids is Manhattan distance (k4M)

I'll now apply the predict function to k4M which uses Manhattan Distance

```
set.seed(64060)
clusters_index4 <- predict(k4M)
dist(k4M@centers)

##           1           2           3
## 2 3.838654
## 3 5.233264 2.754097
## 4 6.025978 4.677079 2.397111

image(k4M)
points(df, col=clusters_index4, pch=19, cex=0.3)
```



But is a K of 4 really the best choice? After all, this was just a random choice. Let's use a K of 3 and examine the results.

```
set.seed(64060)
k3 <- kmeans(df, centers = 3, nstart = 25) # k = 3, number of restarts = 25

# the following will help us Visualize the output

k3$centers # output the centers
```

```
##      Market_Cap      Beta    PE_Ratio      ROE      ROA Asset_Turnover
## 1  0.6733825 -0.3586419 -0.2763512  0.6565978  0.8344159    0.4612656
## 2 -0.8261772  0.4775991 -0.3696184 -0.5631589 -0.8514589   -0.9994088
## 3 -0.6125361  0.2698666  1.3143935 -0.9609057 -1.0174553    0.2306328
##      Leverage Rev_Growth Net_Profit_Margin
## 1 -0.3331068 -0.2902163      0.6823310
## 2  0.8502201  0.9158889     -0.3319956
## 3 -0.3592866 -0.5757385     -1.3784169
```

```
k3$size # Number of firms in each cluster
```

```
## [1] 11  6  4
```

```
fviz_cluster(k3, data = df) # Visualize the output
```



Other Distances

I'll rerun the example using other distance measures to compare the results

```
set.seed(64060)
```

```
k3M = kcca(df, k=3, kccaFamily("kmedians")) #kmedians uses Manhattan
distance
```

```
k3M
```

```
## kcca object of family 'kmedians'
```

```
##
```

```
## call:
```

```
## kcca(x = df, k = 3, family = kccaFamily("kmedians"))
```

```
##
```

```

## cluster sizes:
##
##  1  2  3
##  3 12  6

k3E = kcca(df, k=3, kccaFamily("kmeans")) #kmeans uses Euclidean distance
k3E

## kcca object of family 'kmeans'
##
## call:
## kcca(x = df, k = 3, family = kccaFamily("kmeans"))
##
## cluster sizes:
##
##  1  2  3
## 10  7  4

k3A = kcca(df, k=3, kccaFamily("angle")) #angle uses angle between
observation and centroid
k3A

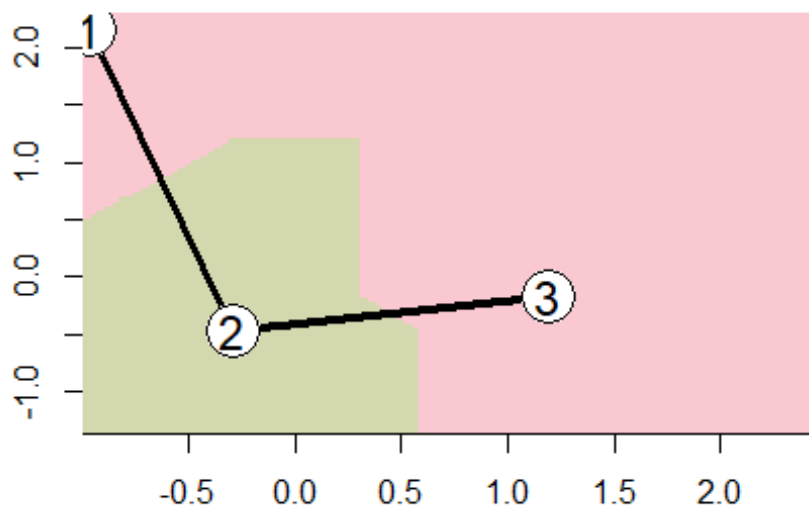
## kcca object of family 'angle'
##
## call:
## kcca(x = df, k = 3, family = kccaFamily("angle"))
##
## cluster sizes:
##
##  1  2  3
##  6  4 11

# We won't use Jaccard distance as this is primarily used for categorical
data which is not applicable.

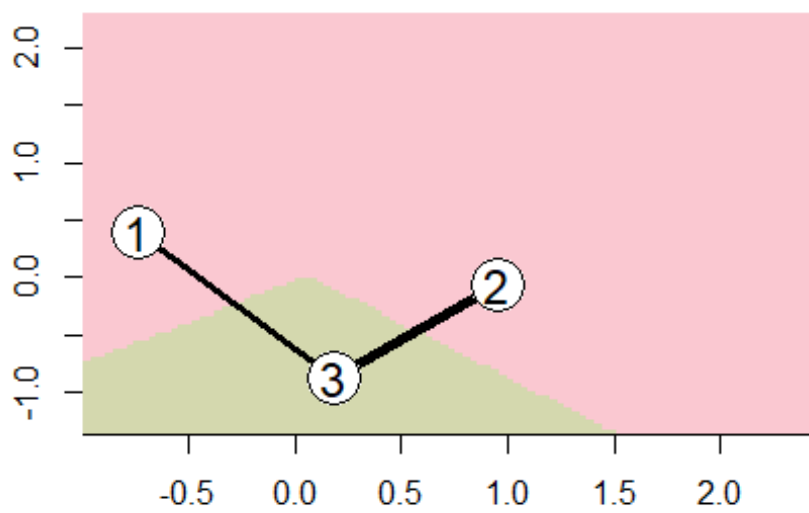
```

Let's take a look at the images of each of these results:

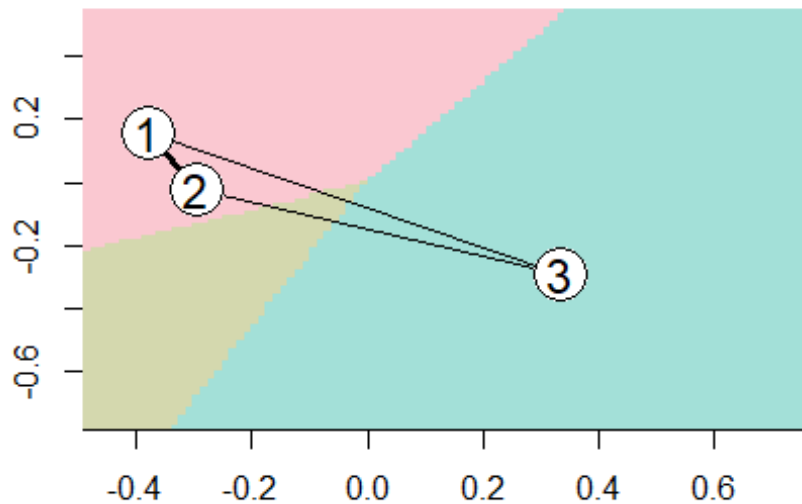
```
image(k3M) # Manhattan distance
```



```
image(k3E) # Euclidean distance
```



```
image(k3A) # angle
```



Once again, using Manhattan distance produces better results with the clustering. There is a greater distance between centroids of the clusters than using Euclidean or Angle.

This is also confirmed by looking at the centers:

```
dist(k3M@centers)
##           1           2
## 2 3.838654
## 3 5.334439 2.990655

dist(k3E@centers)
##           1           2
## 2 3.921756
## 3 2.918545 2.236172

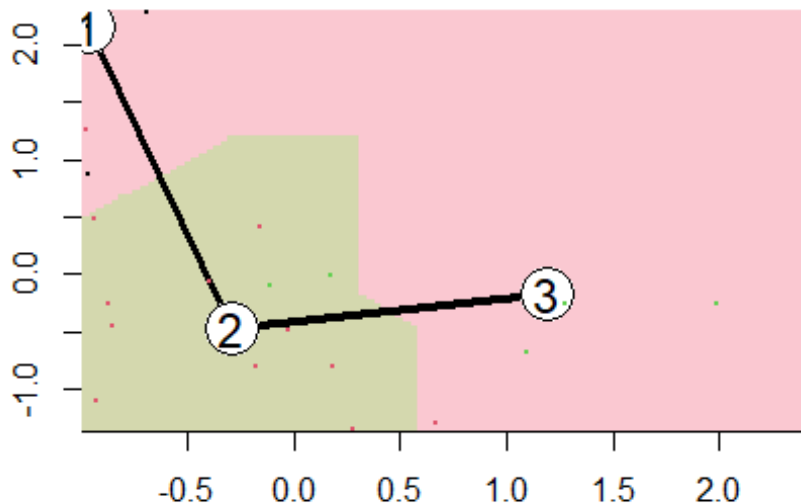
dist(k3A@centers)
##           1           2
## 2 1.276399
## 3 1.898562 1.823885
```

Let's apply the predict function to k3M which uses Manhattan Distance

```
set.seed(64060)
clusters_index3 <- predict(k3M)
dist(k3M@centers)
```

```
##          1          2
## 2 3.838654
## 3 5.334439 2.990655

image(k3M)
points(df, col=clusters_index3, pch=19, cex=0.3)
```



Before we make any conclusions with these results, let's try a K of 5 and analyze the results:

```
set.seed(64060)
k5 <- kmeans(df, centers = 5, nstart = 25) # k = 5, number of restarts = 25
```

the following will help us Visualize the output

```
k5$centers # output the centers
```

##	Market_Cap	Beta	PE_Ratio	ROE	ROA	Asset_Turnover
## 1	-0.87051511	1.3409869	-0.05284434	-0.6184015	-1.1928478	-0.4612656
## 2	-0.43925134	-0.4701800	2.70002464	-0.8349525	-0.9234951	0.2306328
## 3	-0.76022489	0.2796041	-0.47742380	-0.7438022	-0.8107428	-1.2684804
## 4	-0.03142211	-0.4360989	-0.31724852	0.1950459	0.4083915	0.1729746
## 5	1.69558112	-0.1780563	-0.19845823	1.2349879	1.3503431	1.1531640

##	Leverage	Rev_Growth	Net_Profit_Margin
## 1	1.36644699	-0.6912914	-1.320000179
## 2	-0.14170336	-0.1168459	-1.416514761
## 3	0.06308085	1.5180158	-0.006893899

```
## 4 -0.27449312 -0.7041516      0.556954446
## 5 -0.46807818  0.4671788      0.591242521

k5$size # Number of firms in each cluster

## [1] 3 2 4 8 4

fviz_cluster(k5, data = df) # Visualize the output
```



Other Distances

I'll rerun the example using other distance measures to compare the results

```
set.seed(64060)

k5M = kcca(df, k=5, kccaFamily("kmedians")) #kmedians uses Manhattan
distance
k5M

## kcca object of family 'kmedians'
##
## call:
## kcca(x = df, k = 5, family = kccaFamily("kmedians"))
##
## cluster sizes:
##
```

```

## 1 2 3 4 5
## 3 6 5 1 6

k5E = kcca(df, k=5, kccaFamily("kmeans")) #kmeans uses Euclidean distance
k5E

## kcca object of family 'kmeans'
##
## call:
## kcca(x = df, k = 5, family = kccaFamily("kmeans"))
##
## cluster sizes:
##
## 1 2 3 4 5
## 5 1 9 1 5

k5A = kcca(df, k=5, kccaFamily("angle")) #angle uses angle between
observation and centroid
k5A

## kcca object of family 'angle'
##
## call:
## kcca(x = df, k = 5, family = kccaFamily("angle"))
##
## cluster sizes:
##
## 1 2 3 4 5
## 8 4 4 3 2

# We won't use Jaccard distance as this is primarily used for categorical
data which is not applicable.

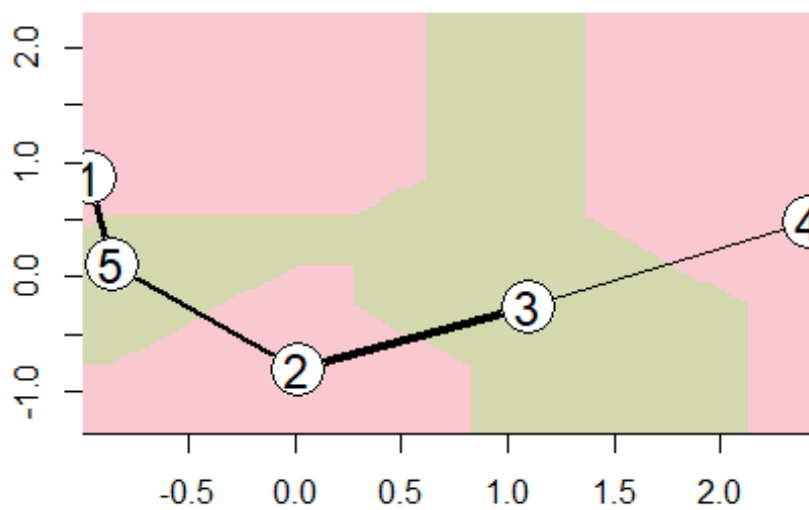
```

Let's take a look at the images of each of these results:

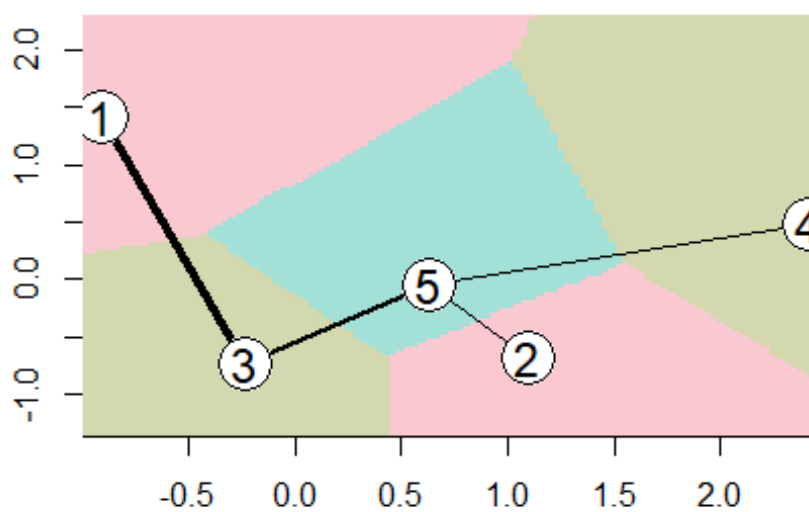
```

image(k5M) # Manhattan distance

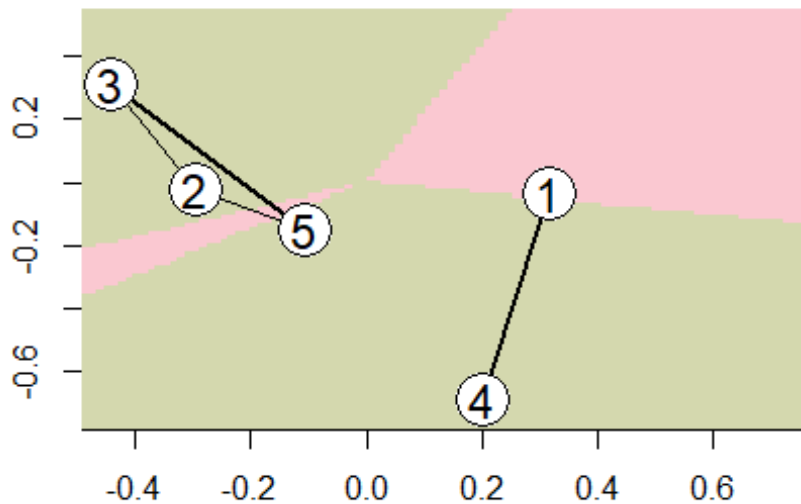
```

```
image(k5E) # Euclidean distance
```



```
image(k5A) # angle
```



This is a little more interesting. It seems with the K of 5, the Euclidean distance may be producing a better result

Let's take a closer look at the centers:

```
dist(k5M@centers)

##          1          2          3          4
## 2 3.721628
## 3 4.689876 2.194249
## 4 5.698767 3.925905 2.397111
## 5 2.931609 2.762659 3.804627 5.718298

dist(k5E@centers)

##          1          2          3          4
## 2 6.108448
## 3 3.091631 4.400111
## 4 5.789244 2.447177 4.465604
## 5 4.177287 2.502227 2.448221 2.791316

dist(k5A@centers)

##          1          2          3          4
## 2 1.822650
## 3 1.840575 1.253136
## 4 1.163136 1.635111 1.732785
## 5 1.776303 1.373808 1.158538 1.508880
```

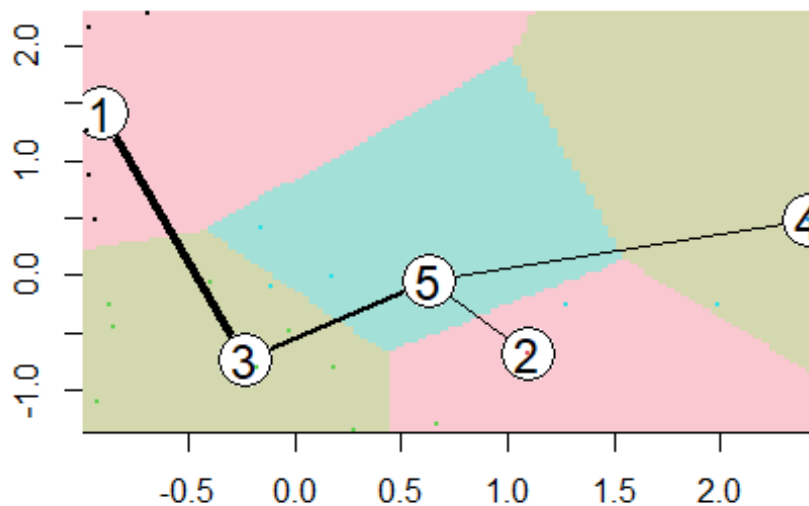
It appears that the Euclidean distance (k5E) is producing better results as can be seen in both the image and the data for the centers. This seems to maximize the distance between the cluster centroids.

Let's apply the predict function to k5E which uses Euclidean Distance

```
set.seed(64060)
clusters_index5 <- predict(k5E)
dist(k5E@centers)

##           1           2           3           4
## 2 6.108448
## 3 3.091631 4.400111
## 4 5.789244 2.447177 4.465604
## 5 4.177287 2.502227 2.448221 2.791316

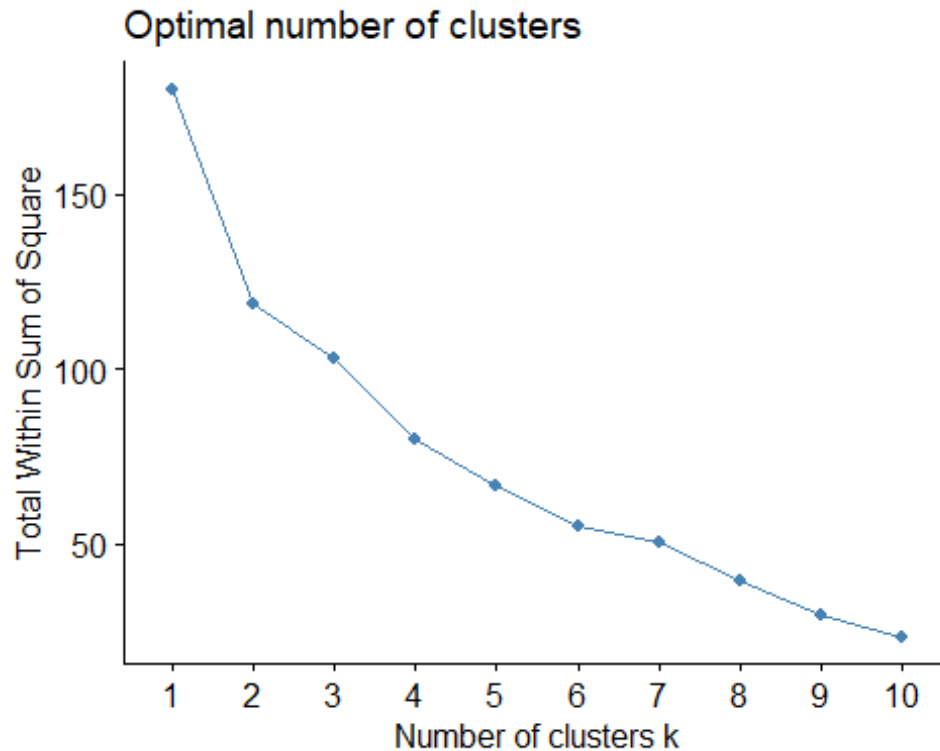
image(k5E)
points(df, col=clusters_index5, pch=19, cex=0.3)
```



CHOOSING THE BEST K Let's use some tools to help us determine the best K value

We'll first review an "elbow chart" to determine k

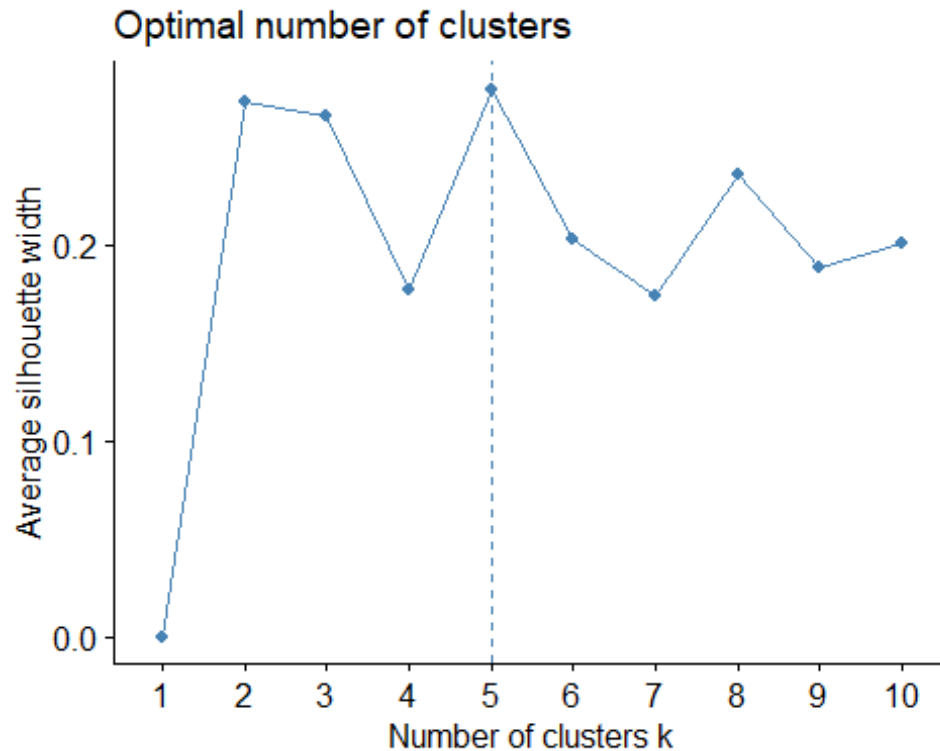
```
fviz_nbclust(df, kmeans, method = "wss")
```



In reviewing the “elbow chart” (WSS), it is a little unclear as to what the optimal number of clusters (k) should be since there is not a clearly visible “knee point” in the results plot. The total WSS has a substantial drop from 1 to 2, less of a drop from 2 to 3, and then another larger drop from 3 to 4. From 4 to 5 and 5 to 6, the decrease is similar and then from 6 to 7, there is little drop. My first thought is the knee point is either at 4 or 6. Honestly, it is difficult to make a reliable determination of the optimal number of clusters (k) using this method (WSS) for this particular set of data. Therefore, I need to confirm this using a different method (Silhouette Method).

Next, we’ll use the Silhouette Method to determine the number of clusters (k)

```
fviz_nbclust(df, kmeans, method = "silhouette")
```



In using the Silhouette Method, it is clear that the optimal number of clusters (k) is 5.

SUMMARY OF REQUIREMENT A:

When clustering, our objective is to minimize the similarity within the cluster and maximize the dissimilarity between the clusters. Meaning, we want the clusters to be as tight as possible with the distance between the clusters to be as great as possible. Also, it is preferable to have the size of the clusters similar.

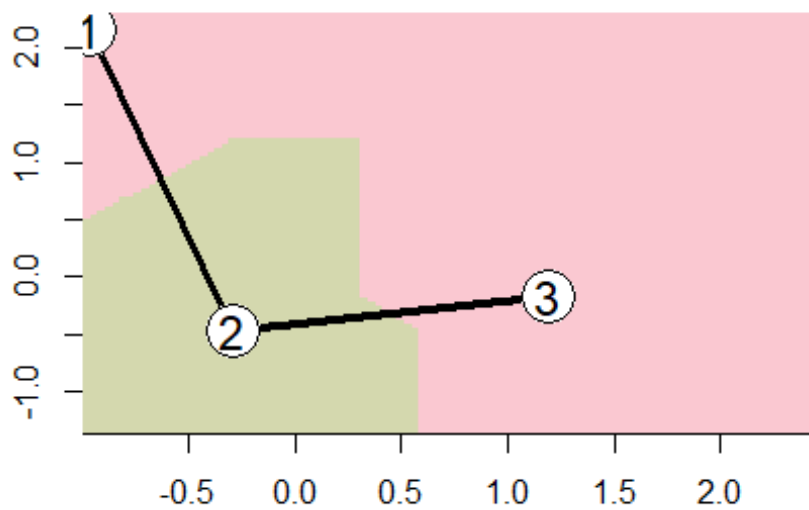
```
k3$size
## [1] 11 6 4

k4$size
## [1] 8 4 6 3

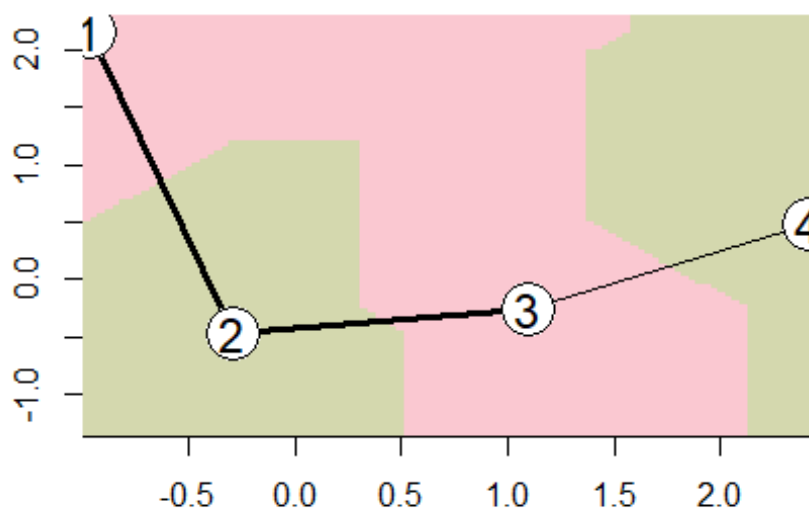
k5$size
## [1] 3 2 4 8 4
```

We have 21 observations. For 3 clusters, the average size would be 7. For 4 clusters, the average size would be 5.25. For 5 clusters, the average size would be 4.2. In reviewing the results of our cluster sizes, the results of our 5 clusters seem to remain closer to the average.

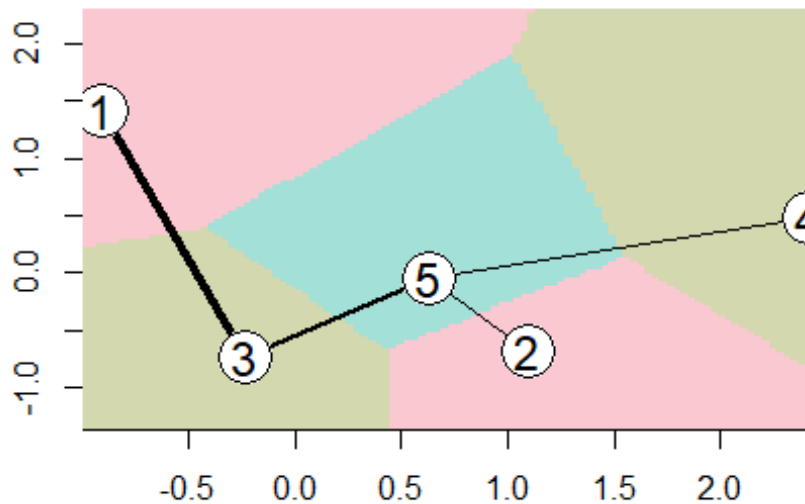
```
image(k3M)
```



image(k4M)



image(k5E)



In reviewing these images, the image with the k of 5 seems to be producing more desirable results. The clusters are “tighter” (more tightly grouped) with the distance between clusters maximized.

```
dist(k3M@centers)
##          1          2
## 2 3.838654
## 3 5.334439 2.990655

dist(k4M@centers)
##          1          2          3
## 2 3.838654
## 3 5.233264 2.754097
## 4 6.025978 4.677079 2.397111

dist(k5E@centers)
##          1          2          3          4
## 2 6.108448
## 3 3.091631 4.400111
## 4 5.789244 2.447177 4.465604
## 5 4.177287 2.502227 2.448221 2.791316
```

The above statement is also confirmed by reviewing the centers

Using the silhouette method in determining k, it confirmed the optimal k value is 5.

WEIGHTING THE VARIABLES In reviewing the data, it's logical to think Market_Cap and PE_Ratio would be more substantial in differentiating the various firms. Going with this assumption, let's place more weight on these variables than on the others.

Let's create the data frame we'll use for the weighted results

```
set.seed(64060)

df_weighted <- Pharmaceuticals[,c(3:11)]

summary(df_weighted)
```

##	Market_Cap	Beta	PE_Ratio	ROE
##	Min. : 0.41	Min. :0.1800	Min. : 3.60	Min. : 3.9
##	1st Qu.: 6.30	1st Qu.:0.3500	1st Qu.:18.90	1st Qu.:14.9
##	Median : 48.19	Median :0.4600	Median :21.50	Median :22.6
##	Mean : 57.65	Mean :0.5257	Mean :25.46	Mean :25.8
##	3rd Qu.: 73.84	3rd Qu.:0.6500	3rd Qu.:27.90	3rd Qu.:31.0
##	Max. :199.47	Max. :1.1100	Max. :82.50	Max. :62.9

##	ROA	Asset_Turnover	Leverage	Rev_Growth
##	Min. : 1.40	Min. :0.3	Min. :0.0000	Min. : -3.17
##	1st Qu.: 5.70	1st Qu.:0.6	1st Qu.:0.1600	1st Qu.: 6.38
##	Median :11.20	Median :0.6	Median :0.3400	Median : 9.37
##	Mean :10.51	Mean :0.7	Mean :0.5857	Mean :13.37
##	3rd Qu.:15.00	3rd Qu.:0.9	3rd Qu.:0.6000	3rd Qu.:21.87
##	Max. :20.30	Max. :1.1	Max. :3.5100	Max. :34.21

##	Net_Profit_Margin
##	Min. : 2.6
##	1st Qu.:11.2
##	Median :16.1
##	Mean :15.7
##	3rd Qu.:21.1
##	Max. :25.5

We need to scale this data frame before we can proceed.

```
# Scaling the data frame (z-score)
df_weighted <- scale(df_weighted)

summary(df_weighted)
```

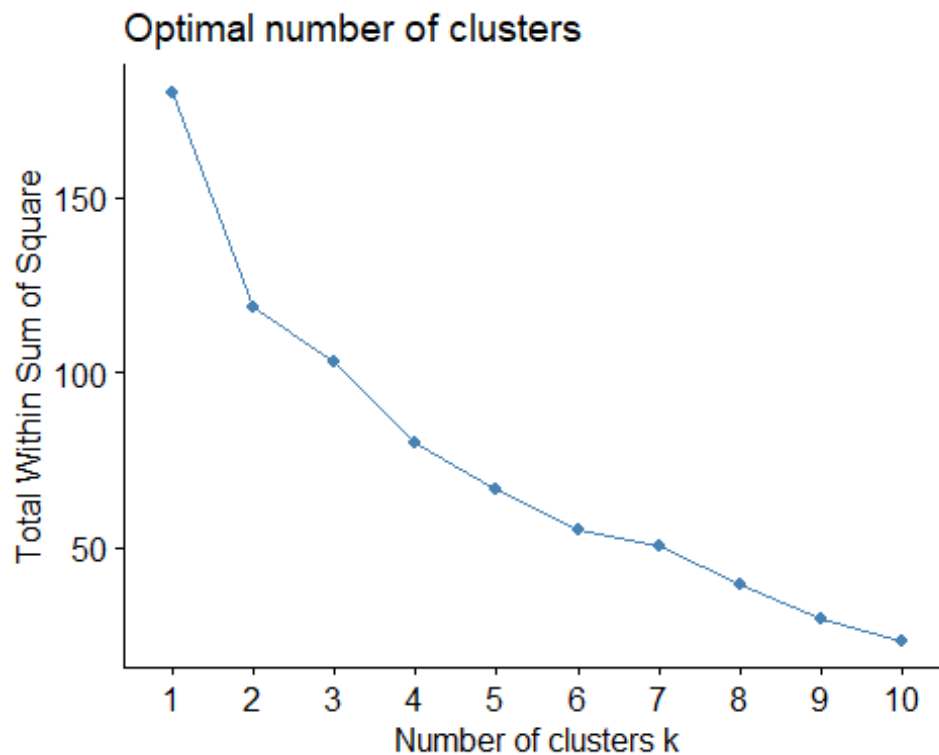
##	Market_Cap	Beta	PE_Ratio	ROE
##	Min. : -0.9768	Min. : -1.3466	Min. : -1.3404	Min. : -1.4515
##	1st Qu.: -0.8763	1st Qu.: -0.6844	1st Qu.: -0.4023	1st Qu.: -0.7223
##	Median : -0.1614	Median : -0.2560	Median : -0.2429	Median : -0.2118
##	Mean : 0.0000	Mean : 0.0000	Mean : 0.0000	Mean : 0.0000
##	3rd Qu.: 0.2762	3rd Qu.: 0.4841	3rd Qu.: 0.1495	3rd Qu.: 0.3450
##	Max. : 2.4200	Max. : 2.2758	Max. : 3.4971	Max. : 2.4597

##	ROA	Asset_Turnover	Leverage	Rev_Growth
##	Min. : -1.7128	Min. : -1.8451	Min. : -0.74966	Min. : -1.4971
##	1st Qu.: -0.9047	1st Qu.: -0.4613	1st Qu.: -0.54487	1st Qu.: -0.6328


```
## Median : 0.1289    Median :-0.4613    Median :-0.31449    Median :-0.3621
## Mean   : 0.0000    Mean   : 0.0000    Mean   : 0.00000    Mean   : 0.0000
## 3rd Qu.: 0.8430    3rd Qu.: 0.9225    3rd Qu.: 0.01828    3rd Qu.: 0.7693
## Max.   : 1.8389    Max.   : 1.8451    Max.   : 3.74280    Max.   : 1.8862
## Net_Profit_Margin
## Min.    :-1.99560
## 1st Qu.:-0.68504
## Median  : 0.06168
## Mean    : 0.00000
## 3rd Qu.: 0.82364
## Max.    : 1.49416
```

CHOOSING THE BEST K Let's review an "elbow chart" to determine k

```
fviz_nbclust(df_weighted, kmeans, method = "wss")
```

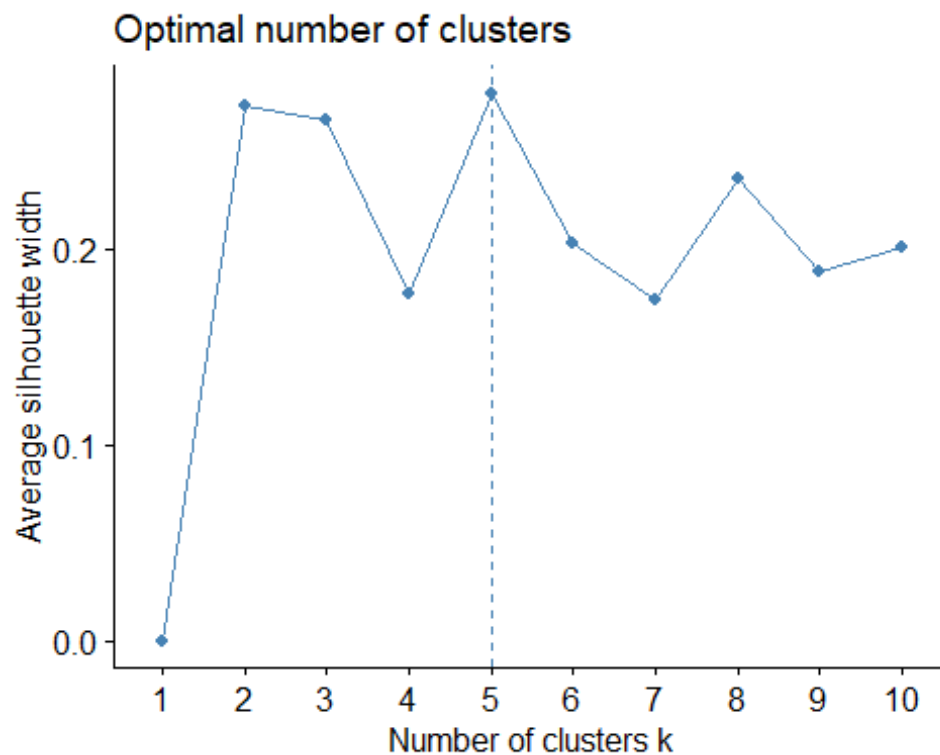


Again, these results

are a little unclear, so let's defer to the Silhouette method.

Now, let's use the Silhouette Method to determine the number of clusters (k)

```
fviz_nbclust(df_weighted, kmeans, method = "silhouette")
```



This shows the optimal number of clusters is 5

Now, we'll place more weight on Market_Cap (1st variable) and PE_Ratio (3rd variable)

```
set.seed(64060)

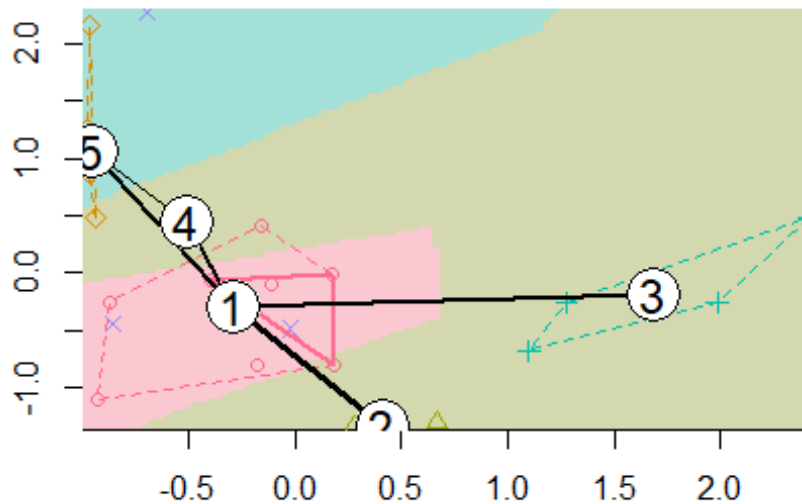
k5_weighted <- cclust(df_weighted, k=5, save.data=TRUE, weights =
c(1,0.5,1,0.5,0.5,0.5,0.5,0.5,0.5), method = "hardcl")

k5_weighted

## kcca object of family 'kmeans'
##
## call:
## cclust(x = df_weighted, k = 5, method = "hardcl", weights = c(1,
## 0.5, 1, 0.5, 0.5, 0.5, 0.5, 0.5, 0.5), save.data = TRUE)
##
## cluster sizes:
##
## 1 2 3 4 5
## 8 2 4 3 4
```

Let's now visualize our results:

```
image(k5_weighted)
```



```
dist(k5_weighted@centers)
##           1           2           3           4
## 2  1.903477
## 3  2.887223  3.278427
## 4  3.498740  4.355677  5.326741
## 5  3.045587  4.188404  5.101772  3.548789
```

CONCLUSIONS: My hypothesis regarding the weight placed on each variable may be incorrect. When reviewing the images of the original k5E and the weighted k5_weighted, k5E produces better cluster results which doesn't seem entirely logical. To get better results with the weighted variables, we would need better estimations of the actual weighted importance of each variable.

REQUIREMENT C:

Is there a pattern in the clusters with respect to the non-numerical variables (10 to 12)?

First, we'll add a column to Pharmaceuticals called "Cluster_No" and set the values equal to the cluster which each observation has been assigned to.

```
Pharmaceuticals$Cluster_No = k5E@cluster
```

We'll first compare the Cluster_No to the Median_Recommendation (variable 10)

```
table(Cluster=Pharmaceuticals$Cluster_No,
Median_Recommendation=Pharmaceuticals$Median_Recommendation)
```

```
##           Median_Recommendation
## Cluster Hold Moderate Buy Moderate Sell Strong Buy
##      1    2           2           1           0
##      2    1           0           0           0
##      3    4           3           1           1
##      4    0           1           0           0
##      5    2           1           2           0
```

In reviewing the distribution of observations of the Median_Recommendation compared to each Cluster, we'll disregard clusters 2 and 4 since these have only 1 observation.

```
Cluster 1, 40% are Hold, 40% are Moderate Buy, and 20% are Moderate Sell.
Cluster 3, 44% are Hold, 33% are Moderate Buy, and 11% are Moderate Sell.
Cluster 5, 40% are Hold, 20% are Moderate Buy, and 40% are Moderate Sell.
```

Therefore, there does NOT seem to be any strong correlation between this variable (variable 10) and the clusters to which they were assigned.

Next, we'll compare the Cluster_No to the Location (variable 11)

```
table(Cluster=Pharmaceuticals$Cluster_No, Location=Pharmaceuticals$Location)
```

```
##           Location
## Cluster CANADA FRANCE GERMANY IRELAND SWITZERLAND UK US
##      1      0      0      1      1           0  0  3
##      2      0      0      0      0           0  1  0
##      3      1      1      0      0           1  1  5
##      4      0      0      0      0           0  0  1
##      5      0      0      0      0           0  1  4
```

In reviewing the distribution of observations of the Location compared to each Cluster, again, we'll disregard Clusters 2 and 4 since these have only 1 observation.

```
Cluster 1, 20% are Germany, 20% are Ireland, and 60% are US.
Cluster 3, 11% are Canada, 11% are France, 11% are Switzerland, 11% are UK,
and 55% are US
Cluster 5, 20% are UK and 80% are US
```

From the raw data, we know that 62% of the firms in the observations are from the US.

While a larger percentage of firms in each Cluster are from the US, the percentage of US firms in each cluster is not much different than the overall average number of firms in the US. Therefore, there does not seem to be a strong correlation between Cluster and Location.

Now, we'll compare the Cluster_No to the Exchange (variable 12)

```
table(Cluster=Pharmaceuticals$Cluster_No, Exchange=Pharmaceuticals$Exchange)
```

```
##           Exchange
## Cluster AMEX NASDAQ NYSE
##      1    1      1      3
```

```
##      2      0      0      1
##      3      0      0      9
##      4      0      0      1
##      5      0      0      5
```

In reviewing the distribution of observations of the Exchange compared to each Cluster, again, we'll disregard Clusters 2 and 4 since these have only 1 observation.

```
Cluster 1, 20% are AMEX, 20% are NASDAQ, and 60% are NYSE
Cluster 3, 100% are NYSE
Cluster 5, 100% are NYSE
```

However, we know from the raw data that of the 21 firms listed, only 1 are on the AMEX and only 1 are on the NASDAQ. All remaining 19 firms are on the NYSE.

Since the one firm on the AMEX and the one firm on the NASDAQ are both listed in the same cluster (Cluster 1) together with 3 firms on the NYSE, it seems there is no correlation between Cluster and Exchange.

REQUIREMENT D:

Provide an appropriate name for each cluster using any or all of the variables in the dataset.

For this, I'll use 2 key variables for simplification: Market_Cap and PE_Ratio

```
set.seed(64060)

d_df <- Pharmaceuticals[,c(3,5)]

summary(d_df)

##      Market_Cap      PE_Ratio
## Min.   : 0.41   Min.   : 3.60
## 1st Qu.: 6.30   1st Qu.:18.90
## Median :48.19   Median :21.50
## Mean   :57.65   Mean    :25.46
## 3rd Qu.:73.84   3rd Qu.:27.90
## Max.   :199.47   Max.    :82.50
```

We'll scale the data frame

```
# Scaling the data frame (z-score)
d_df <- scale(d_df)

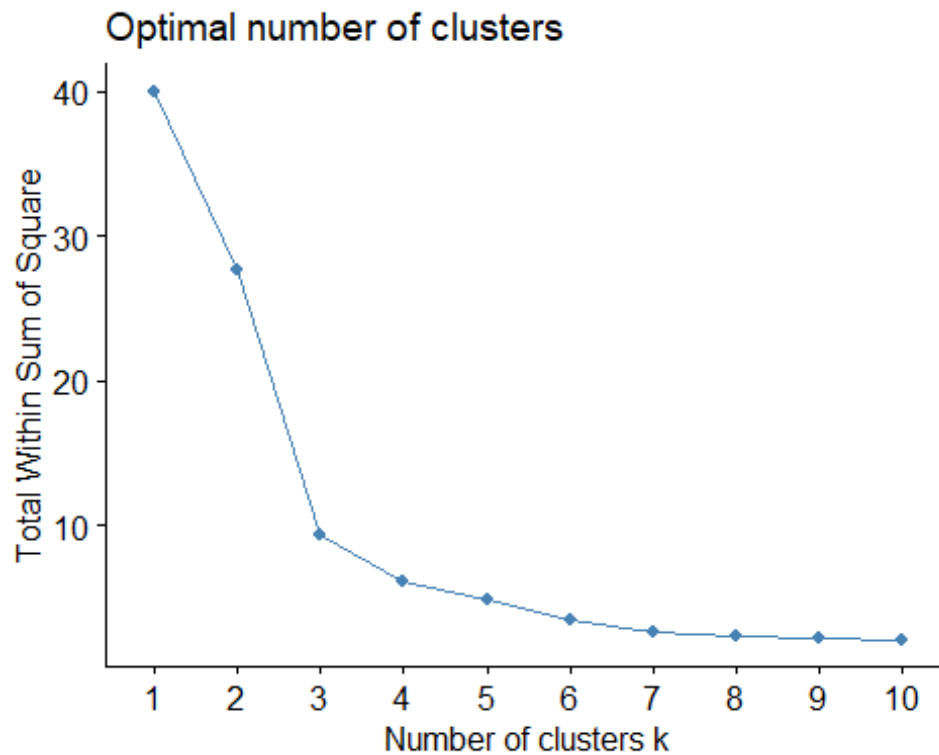
summary(d_df)

##      Market_Cap      PE_Ratio
## Min.   :-0.9768   Min.    :-1.3404
## 1st Qu.: -0.8763   1st Qu.: -0.4023
## Median : -0.1614   Median : -0.2429
## Mean    : 0.0000   Mean     : 0.0000
```

```
## 3rd Qu.: 0.2762    3rd Qu.: 0.1495
## Max.    : 2.4200    Max.    : 3.4971
```

CHOOSING THE BEST K Let's review an "elbow chart" to determine k

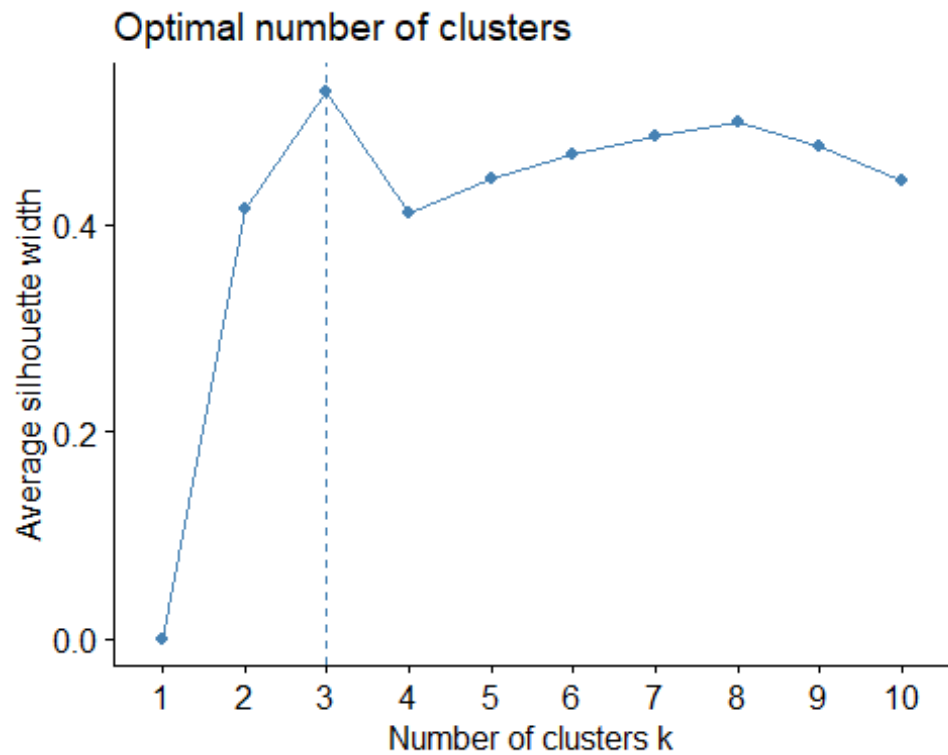
```
fviz_nbclust(d_df, kmeans, method = "wss")
```



This chart clearly shows the knee point at 3, indicating the optimal number of clusters is 3

Now, let's use the Silhouette Method to determine the number of clusters (k)

```
fviz_nbclust(d_df, kmeans, method = "silhouette")
```



This is also

showing the optimal number of clusters is 3.

I'll run the k-means algorithm to cluster the firms and then visualize the output

```
set.seed(64060)

d_k3 <- kmeans(d_df, centers = 3, nstart = 25) # k = 3, number of restarts = 25

fviz_cluster(d_k3, data = d_df) # Visualize the output
```

