# Assignment 5

#### **MUNERAH**

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```
## Load the file
library(readx1)
Cereals <- read.csv("C:/Users/mnooo/Desktop/Datasets/Cereals.csv")</pre>
# Loading the libraries
library(tidyverse)
## -- Attaching packages ----- tidyverse 1.3.1 --
## v ggplot2 3.3.5
                    v purrr
                              0.3.4
## v tibble 3.1.6 v dplyr 1.0.7
## v tidyr 1.1.4 v stringr 1.4.0
## v readr 2.0.1
                     v forcats 0.5.1
## Warning: package 'tibble' was built under R version 4.1.2
## Warning: package 'tidyr' was built under R version 4.1.2
## -- Conflicts ------ tidyverse conflicts() --
## x dplyr::filter() masks stats::filter()
## x dplyr::lag() masks stats::lag()
library(corrplot)
## corrplot 0.90 loaded
library(tidymodels)
## Warning: package 'tidymodels' was built under R version 4.1.2
## Registered S3 method overwritten by 'tune':
##
    method
                            from
    required_pkgs.model_spec parsnip
##
## -- Attaching packages ------ tidymodels 0.1.4 --
```

```
## v broom
                 0.7.9
                            v rsample
                                           0.1.1
## v dials
                 0.0.10
                                           0.1.6
                            v tune
## v infer
                 1.0.0
                            v workflows
                                           0.2.4
## v modeldata
                            v workflowsets 0.1.0
                 0.1.1
## v parsnip
                 0.1.7
                            v yardstick
                                          0.0.9
## v recipes
                 0.1.17
## Warning: package 'dials' was built under R version 4.1.2
## Warning: package 'infer' was built under R version 4.1.2
## Warning: package 'modeldata' was built under R version 4.1.2
## Warning: package 'parsnip' was built under R version 4.1.2
## Warning: package 'rsample' was built under R version 4.1.2
## Warning: package 'tune' was built under R version 4.1.2
## Warning: package 'workflows' was built under R version 4.1.2
## Warning: package 'workflowsets' was built under R version 4.1.2
## Warning: package 'yardstick' was built under R version 4.1.2
## -- Conflicts ----- tidymodels_conflicts() --
## x scales::discard() masks purrr::discard()
## x dplyr::filter() masks stats::filter()
## x recipes::fixed() masks stringr::fixed()
## x dplyr::lag()
                  masks stats::lag()
## x yardstick::spec() masks readr::spec()
## x recipes::step() masks stats::step()
## * Use suppressPackageStartupMessages() to eliminate package startup messages
library(cluster)
library(factoextra)
## Welcome! Want to learn more? See two factoextra-related books at https://goo.gl/ve3WBa
library(knitr)
```

```
## Data pre-processing
# Transfer the cereal name into a row name
Cereals <- data.frame(Cereals,row.names = 'name')

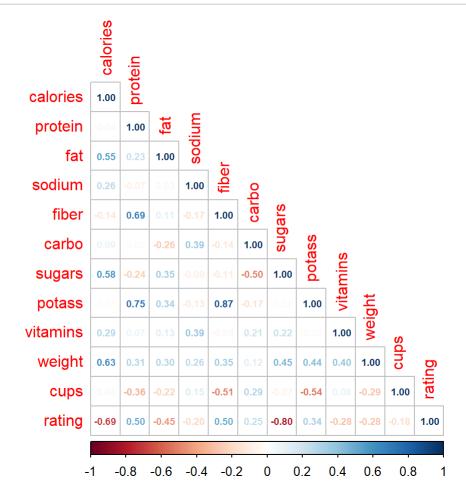
# Remove missing values
Cereals <- Cereals %>%
        filter(complete.cases(.)) %>%
        select(-c(mfr,type,shelf))

#Normalization
scaled <-data.frame(scale(Cereals))</pre>
```

### Exploratory analysis

In this part, we explore the datasets using visualization. We use the correlation plot as well as scatter plots to visualize relationships.

Below is a correlation matrix plot showing the relationship between pairs of variables.



We notice that rating is highly affected negatively by sugars and calories while protein and fiber has positive significant effect on rating.

Apply hierarchical clustering to the data using Euclidean distance to the normalised mesurements.

```
#Hierarchical Clustering
metd<-c('single','complete','average','ward')
names(metd)<- c('single','complete','average','ward')

# we write a function which computes agglomerative clustering
AGNES<- function(x){
        agnes(scaled, method = x)$ac
}</pre>
```

#calculate agglomerative coefficient for each clustering linkage method sapply(metd, AGNES)

```
## single complete average ward
## 0.6072384 0.8469328 0.7881955 0.9087265
```

We notice that Ward's minimum variance method provides the highest agglomerative coefficient. We will use it as our final model for hierarchical clustering.

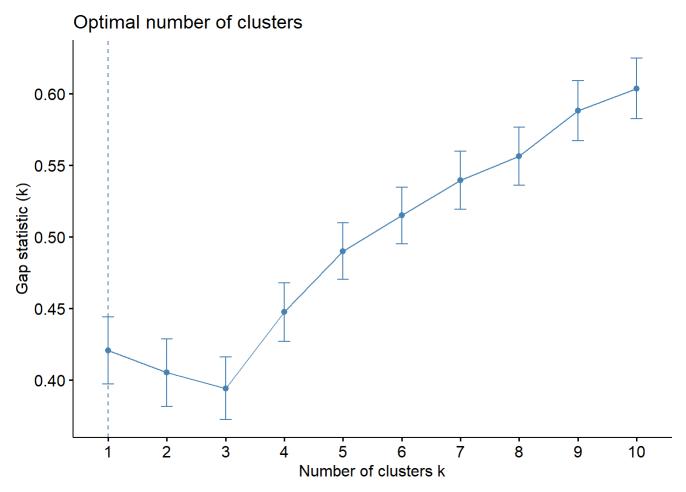
```
set.seed(100)
#Compute th euclidean distance matrix
dm<- dist(scaled, method = "euclidean")

#We perform hierarchical clustering using ward's method
wClust<- hclust(dm, method = "ward.D2")
wClust</pre>
```

```
##
## Call:
## hclust(d = dm, method = "ward.D2")
##
## Cluster method : ward.D2
## Distance : euclidean
## Number of objects: 74
```

# How many clusters will you choose?

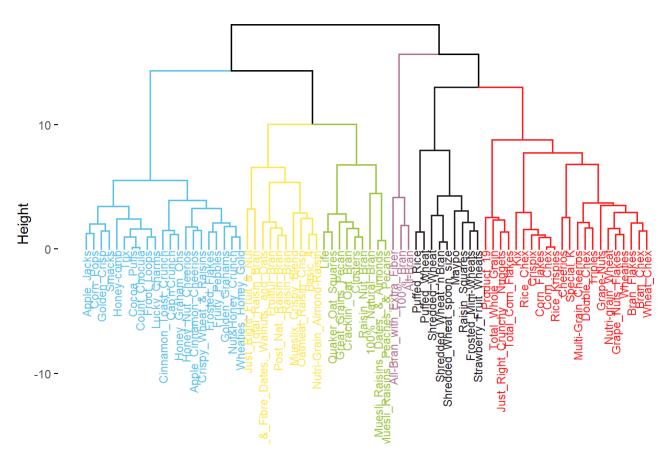
```
#Selecting number of clusters K using gap-stat
fviz_nbclust(scaled, hcut, method = "gap_stat") # K = 7
```



The gap-statistic does not suggest an ideal number of cluster as it indicates its optimal number to be 1. We will not use this but we will rather use 6 clusters.

```
## Warning: `guides(<scale> = FALSE)` is deprecated. Please use `guides(<scale> =
## "none")` instead.
```

### Dendogram for AGNES [ward method]



We notice how these clusters are being partitioned with different colours from the dendrogram above.

# Comment on the structure of the clusters and their stability.

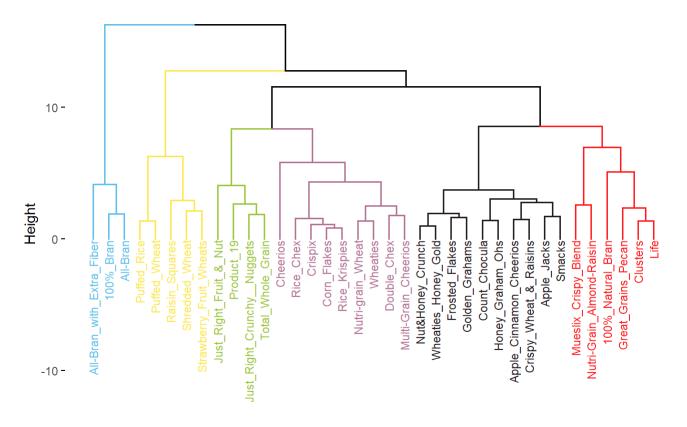
In order to understand the structure of the cluster, we split the data into two. We perform Hierarchical Clustering on both partitions using AGNES Ward method and plotted the dendrograms.

```
#For partA
set.seed(100)
dmA<- dist(partA, method = 'euclidean')
aClust<- hclust(dmA, method = 'ward.D2')</pre>
```

We then plot a dendrogram for the first partition.

```
## Warning: `guides(<scale> = FALSE)` is deprecated. Please use `guides(<scale> =
## "none")` instead.
```

## Dendogram for partition A



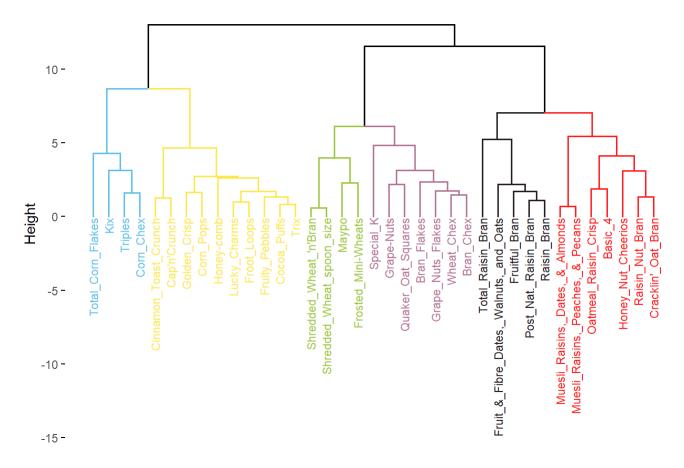
We do a similar clustering on the second partition.

```
#For partB
set.seed(100)
dmB<- dist(partB, method = 'euclidean')
bClust<- hclust(dmB, method = 'ward.D2')</pre>
```

We then plot a dendrogram for the first partition.

```
## Warning: `guides(<scale> = FALSE)` is deprecated. Please use `guides(<scale> =
## "none")` instead.
```

#### Dendogram for partition B



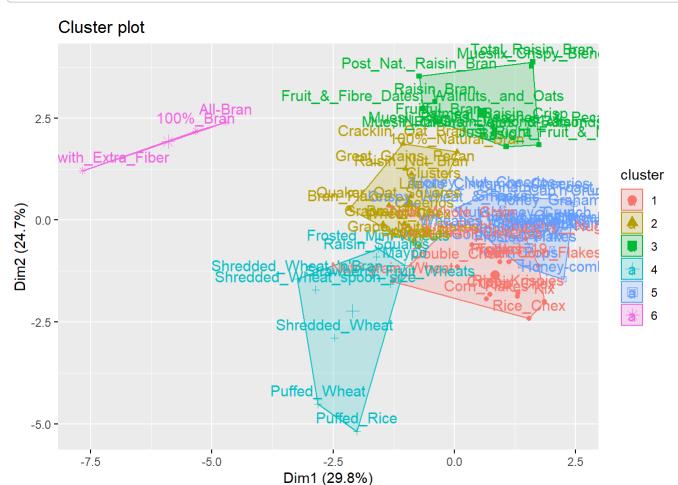
We discovered that the clusters maintain similar structure for more than 90% even when the data was randomly split. We may observe this by looking into the dendrogram for partition A and Partition B.

# Stability of the cluster

A cluster is said to be stable if the partitioning remains similar when the dataset are randomised or the clustering process changes.

We use K means clustering to assess stability of the clusters.

```
set.seed(100)
KM<-kmeans(scaled,6, nstart = 25)
#Plot the clusters
fviz_cluster(KM, data = scaled)</pre>
```



We can see an exact clustering pattern if we examine both the dendrogram and the cluster plot for the K means algorithm.

# Finding a cluster of "healthy cereals."

To identify the cluster with healthy cereals, we may be interested in using rating but we have to identify the features which influences rating through our correlation plot.

From our correlation plot, sugar and calories are having significant negative effect on ratings and protein and fiber has positive effect on rating. We may use these variables or rating to select the cluster with the best nutritional content.

```
# assign cluster to data
cuts<-cutree(wClust, k = 6)
clusterData<- cbind(Cereals, cluster = cuts)

centroids<- clusterData %>%
    group_by(cluster) %>%
    summarise_all(funs(mean))
```

```
## Warning: `funs()` was deprecated in dplyr 0.8.0.
## Please use a list of either functions or lambdas:
##
     # Simple named list:
##
##
     list(mean = mean, median = median)
##
##
     # Auto named with `tibble::lst()`:
     tibble::lst(mean, median)
##
##
##
     # Using lambdas
     list(~ mean(., trim = .2), ~ median(., na.rm = TRUE))
##
## This warning is displayed once every 8 hours.
## Call `lifecycle::last_lifecycle_warnings()` to see where this warning was generated.
```

```
kable(round(centroids,2))
```

cluster	calories	protein	fat	sodium	fiber	carbo	sugars	potass	vitamins	weight	cups	rating
1	63.33	4.00	0.67	176.67	11.00	6.67	3.67	310.00	25.00	1.00	0.39	73.84
2	117.78	3.44	2.67	115.56	2.61	12.50	7.56	131.67	22.22	1.00	0.67	40.72
3	110.95	1.52	1.00	168.57	0.57	12.43	11.48	47.38	25.00	1.00	0.88	28.66
4	132.00	3.00	1.40	192.00	3.65	15.55	10.90	172.50	40.00	1.34	0.71	36.26
5	103.64	2.68	0.55	225.91	1.68	18.32	3.45	74.55	38.64	1.00	0.93	46.18
6	82.22	2.44	0.11	1.67	2.11	15.33	2.33	90.56	11.11	0.87	0.85	63.02

Cluster 1 has the highest rating with an average of 73.8. This cluster has cereals with the highest Protein and fibre but with the least calories. It has almost close to the least sugar on average too. Cluster 1 has 100% Bran, All Bran and All Bran with extra fibre as the cereals in it.

#### Should the data be normalised

For this goal, we use the main dataset which is not normalised so as to be able to properly interpret the result but we use normalised data to obtain clusters since each variable has different measurement scale.

# How should they be used in the cluster analysis?

In order to use this in the cluster analysis, we assigned the cluster values to each cereal in the original dataset and computed the mean for each features based on the clusters assigned. This makes it easy to analyse and advise on which cluster has the healthy cereals.