# Class 7: Machine Learning 1

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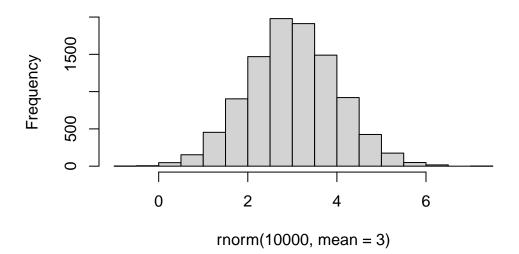
Today we will explore unsupervised machine learning methods starting with clustering and dimensionality reduction.

## Clustering

To start let's make up some data to cluster where we know what the answer should be. The <code>rnorm()</code> function will help us here.

```
hist(rnorm(10000, mean=3))
```

# Histogram of rnorm(10000, mean = 3)



Return 30 numbers centered on -3

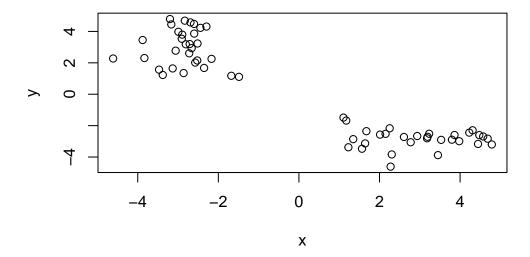
```
[1,] -2.991297
                 3.978344
 [2,] -3.834980
                 2.304359
 [3,] -2.691131
                 4.573042
 [4,] -2.603694
                 4.476778
 [5,] -2.598901
                 3.861289
 [6,] -3.470747
                 1.565748
 [7,] -3.166711
                4.445206
 [8,] -2.806797
                 3.179574
 [9,] -2.168453
                 2.252648
[10,] -3.877895
                 3.450803
[11,] -2.661681
                 2.935304
[12,] -2.573306
                 2.014116
[13,] -2.906548
                 3.530135
[14,] -2.861164 1.347440
```

- [15,] -2.710405 3.193785
- [16,] -2.353080 1.674420
- [17,] -2.895673 3.797106
- [18,] -3.058580 2.773551
- [19,] -2.515713 3.231854
- [20,] -2.831384 4.684583
- [21,] -2.523303 2.150658
- [22,] -1.487225 1.106359
- [23,] -3.132889 1.642689
- [24,] -2.722205 2.606078
- [25,] -2.444967 4.227515
- [26,] -3.378592 1.227771
- [27,] -4.610479 2.275412
- [28,] -1.677803 1.174826
- [29,] -3.200248 4.787185
- [30,] -2.295523 4.310386
- [31,] 4.310386 -2.295523
- [32,] 4.787185 -3.200248
- [33,] 1.174826 -1.677803
- [34,] 2.275412 -4.610479
- [35,] 1.227771 -3.378592
- [36,] 4.227515 -2.444967
- [37,] 2.606078 -2.722205
- [38,] 1.642689 -3.132889
- [39,] 1.106359 -1.487225
- [40,] 2.150658 -2.523303
- [41,] 4.684583 -2.831384
- [42,]3.231854 -2.515713
- [43,]2.773551 -3.058580
- [44,]3.797106 -2.895673
- [45,]1.674420 -2.353080
- [46,] 3.193785 -2.710405
- [47,]1.347440 -2.861164
- [48,] 3.530135 -2.906548
- [49,]2.014116 -2.573306
- [50,] 2.935304 -2.661681
- [51,] 3.450803 -3.877895
- [52,] 2.252648 -2.168453
- [53,] 3.179574 -2.806797
- [54,] 4.445206 -3.166711
- [55,] 1.565748 -3.470747
- [56,] 3.861289 -2.598901
- [57,] 4.476778 -2.603694

```
[58,] 4.573042 -2.691131
[59,] 2.304359 -3.834980
[60,] 3.978344 -2.991297
```

Make a plot of x

### plot(x)



#### K-means

The main function in "base" R for K-means clustering is called kmeans():

```
km <- kmeans(x, centers=2)
km</pre>
```

K-means clustering with 2 clusters of sizes 30, 30

Cluster means:

Clustering vector:

Within cluster sum of squares by cluster:

[1] 51.89665 51.89665

(between\_SS / total\_SS = 90.7 %)

Available components:

[1] "cluster" "centers" "totss" "withinss" "tot.withinss"

[6] "betweenss" "size" "iter" "ifault"

The kmeans() function return a "list" with 9 components. You can see the named components of any list with the attributes() function.

#### attributes(km)

#### \$names

[1] "cluster" "centers" "totss" "withinss" "tot.withinss"

[6] "betweenss" "size" "iter" "ifault"

#### \$class

[1] "kmeans"

Q. How many points in each cluster?

#### km\$size

[1] 30 30

Q. Cluster assignment/memebership vector?

#### km\$cluster

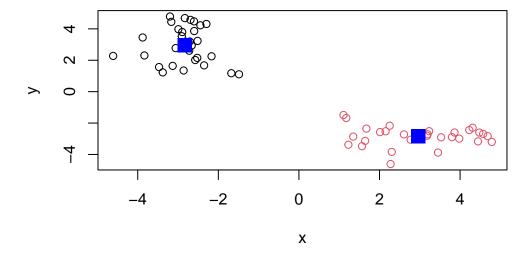
Q. Cluster centers?

#### km\$centers

```
x y
1 -2.835046 2.959299
2 2.959299 -2.835046
```

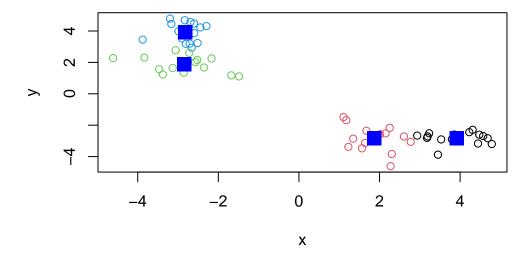
Q. Make a plot of our kmeans() results showing cluster assignment using different colors for each cluster/group of points and cluster centers in blue.

```
plot(x, col=km$cluster)
points(km$centers, col="blue", pch=15, cex=2)
```



Q. Run kmeans() again on x and this cluster into 4 groups/clusters and plot the same result figure as above.

```
km4 <- kmeans(x, centers=4)
plot(x, col=km4$cluster)
points(km4$centers, col="blue", pch=15, cex=2)</pre>
```



**key-point**: K-means clustering is super popular but can be miss-used. One big limitation is that it can impose a clustering pattern on your data even if clear natural grouping don't exist - i.e. it does what you tell it to do in terms of centers.

#### **Hierarchical Clustering**

The main function in "base" R for hierarchical clustering is called hclust().

You can't just pass our dataset as is into hclust(). You must give "distancee matrix" as input. We can get this from the dist() function in R.

```
d <- dist(x)
hc <- hclust(d)
hc</pre>
```

# Call: hclust(d = d)

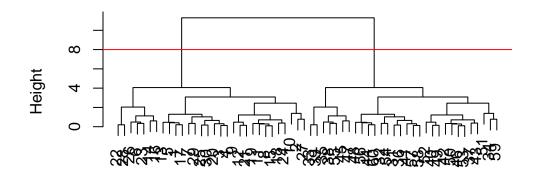
Cluster method : complete
Distance : euclidean

Number of objects: 60

The results of hclust() don't have a useful print() method but do have a special plot() method.

```
plot(hc)
abline(h=8, col="red")
```

# **Cluster Dendrogram**

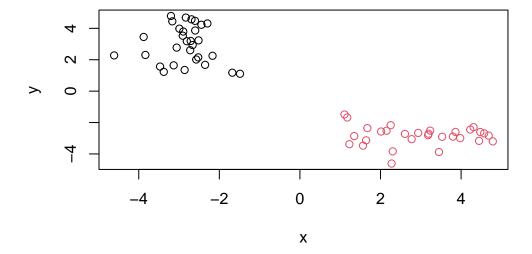


d hclust (\*, "complete")

To get our main cluster assignment (membership vector), we need to "cut" the tree at the big goal posts...

```
grps <- cutree(hc, h=8)
grps</pre>
```

table(grps)



Hierarchical Clustering is distinct in that the dendrogram (tree figure) can reveal the potential grouping in your data (unlike K-means).

## **Principal Component Analysis (PCA)**

PCA is a common and highly useful dimensionality reduction technique used in many fields - particularly bioinformatics.

Here we will analyze some data from the UK on food consumption.

### Data import

```
url <- "https://tinyurl.com/UK-foods"
x <- read.csv(url)
head(x)</pre>
```

```
X England Wales Scotland N.Ireland
1
          Cheese
                     105
                           103
                                    103
                                               66
2 Carcass_meat
                     245
                           227
                                    242
                                              267
3
     Other_meat
                     685
                           803
                                    750
                                              586
4
            Fish
                           160
                                    122
                                               93
                     147
5 Fats_and_oils
                     193
                           235
                                    184
                                              209
          Sugars
                     156
                           175
                                    147
                                              139
```

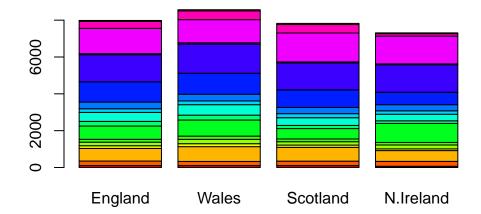
```
rownames(x) <- x[,1]
x <- x[,-1]
head(x)
```

	England	Wales	${\tt Scotland}$	N.Ireland
Cheese	105	103	103	66
Carcass_meat	245	227	242	267
Other_meat	685	803	750	586
Fish	147	160	122	93
Fats_and_oils	193	235	184	209
Sugars	156	175	147	139

```
x <- read.csv (url, row.names=1)
head(x)</pre>
```

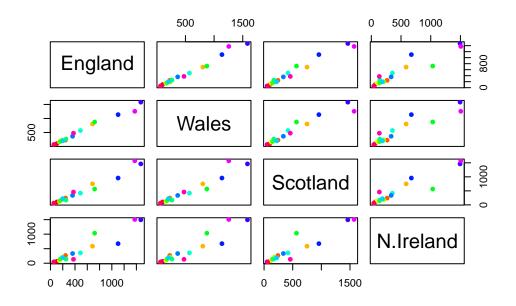
	England	Wales	${\tt Scotland}$	N.Ireland
Cheese	105	103	103	66
Carcass_meat	245	227	242	267
Other_meat	685	803	750	586
Fish	147	160	122	93
Fats_and_oils	193	235	184	209
Sugars	156	175	147	139

```
barplot(as.matrix(x), beside=F, col=rainbow(nrow(x)))
```



One conventional plot that can be useful is called a "paris" plot.

## pairs(x, col=rainbow(nrow(x)), pch=16)



#### PCA to the rescue

The main function in base R for PCA is called prcomp().

```
pca <- prcomp(t(x))
summary(pca)</pre>
```

Importance of components:

```
PC1 PC2 PC3 PC4
Standard deviation 324.1502 212.7478 73.87622 3.176e-14
Proportion of Variance 0.6744 0.2905 0.03503 0.000e+00
Cumulative Proportion 0.6744 0.9650 1.00000 1.000e+00
```

The prcomp() function returns a list object of our result with five attributes/components.

```
attributes(pca)
```

```
$names
```

```
[1] "sdev" "rotation" "center" "scale" "x"
```

#### \$class

[1] "prcomp"

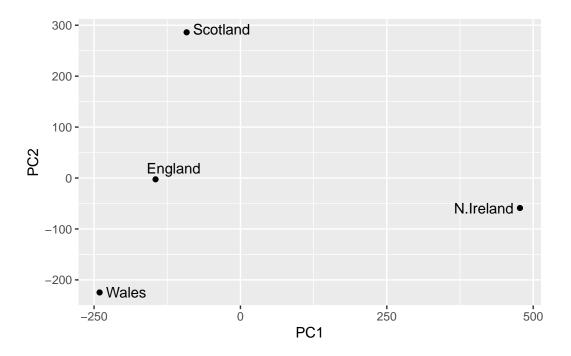
The two main "results" in here are pca\$x and pca\$rotation. The first of these (pca\$x) contains the scores of the data on the new PC axis - we use these to make our "PCA plot".

```
pca$x
```

```
PC1 PC2 PC3 PC4
England -144.99315 -2.532999 105.768945 -4.894696e-14
Wales -240.52915 -224.646925 -56.475555 5.700024e-13
Scotland -91.86934 286.081786 -44.415495 -7.460785e-13
N.Ireland 477.39164 -58.901862 -4.877895 2.321303e-13
```

```
library(ggplot2)
library(ggrepel)

# Make a plot of pca$x with PC1 vs. PC2
ggplot(pca$x) +
   aes (PC1, PC2, label=rownames(pca$x)) +
   geom_point() +
   geom_text_repel()
```

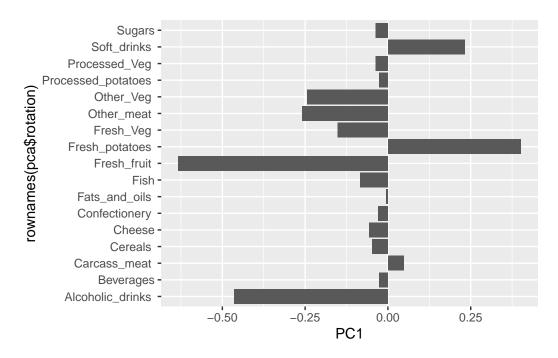


Notes: Using the ggplot2 packages, this results in a figure describing lists of vectors and the matrix between four different countries, to mainly represent the variation: how similar and dissimilar they are from each other. For example, Scotland, England, and Wales are on same area (the negative PC1), while N.Ireland are plotted in the other area (the positive PC1)

So first we will need to take whatever it is we want to plot and convert it to a data.frame with the as.data.frame() function. Then to make our plotting life easier we will also add the food labels as a column (called "Food") to this data frame with the rownames\_to\_column() function from the tibble package (you might need to install this):

The second major result is contained in the pca\$rotation object or component. Let's plot this to see what PCA is picking up...

```
ggplot(pca$rotation) +
aes(PC1, rownames(pca$rotation)) +
geom_col()
```



Notes: based on the graph above, this plot specifically describes the proportion of each component of food in bar representation, based on PC1 value. Therefore, for example, fresh fruit and alcoholic drinks, being represented significantly high toward the negative PC1, is likely being consumed in three countries, Scotland, England, and Wales. On the other hand, soft drinks and fresh potatos, being represented significantly high toward the positive PC1, is likely consumed in N.Ireland.