# Class 7: Machine Learning 1

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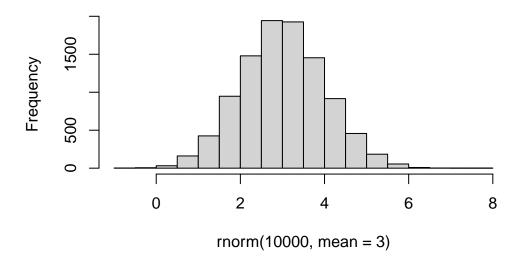
Today we will explore unsupervised machine learning methods starting with clustering and dimensionality reduction.

### 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.

```
tmp <- c(rnorm(30, mean=-3), rnorm(30, mean=3))
x <- cbind(x=tmp, y=rev(tmp))
x</pre>
```

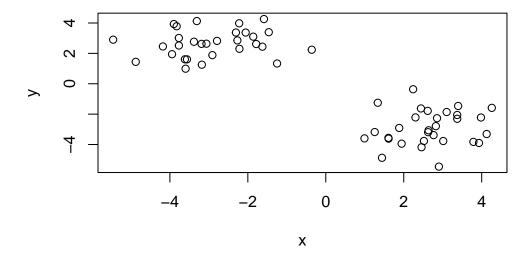
```
Х
 [1,] -4.1792452
                  2.4568767
 [2,] -1.8611433
                  3.1012253
 [3,] -3.6145729
                  1.6109763
 [4,] -3.7684518
                  2.5178595
 [5,] -1.4628548
                  3.3955168
 [6,] -3.1796362
                  1.2557467
 [7,] -1.2496618 1.3337774
 [8,] -1.6263705
                  2.4410630
 [9,] -3.5617735
                  1.6070464
[10,] -3.8267480
                  3.7865929
[11,] -2.0555639
                  3.3714719
[12,] -0.3600657
                  2.2407225
[13,] -1.5863887
                  4.2595145
[14,] -3.8977488
                  3.9308435
```

- [15,] -3.1875920 2.6254494
- [16,] -2.2159012 2.3007914
- [17,] -3.7698507 3.0134074
- [18,] -2.2682917 2.8545262
- [19,] -1.7851870 2.6147537
- [20,] -4.8754301 1.4418009
- [21,] -3.3824479 2.7656617
- [22,] -3.9445149 1.9460432
- [23,] -2.9097697 1.8836764
- [24,] -2.7905928 2.8234348
- [25,] -2.3058212 3.3724875
- [26,] -3.0626125 2.6377440
- [27,] -5.4557729 2.9019691
- [27,] 5.4557729 2.9019091
- [28,] -3.3088316 4.1273693
- [29,] -2.2218544 3.9811237
- [30,] -3.5969278 0.9921391
- [31,] 0.9921391 -3.5969278
- [32,] 3.9811237 -2.2218544
- [33,] 4.1273693 -3.3088316
- [34,] 2.9019691 -5.4557729
- [35,] 2.6377440 -3.0626125
- [36,] 3.3724875 -2.3058212
- [37,] 2.8234348 -2.7905928
- [38,] 1.8836764 -2.9097697
- [39,] 1.9460432 -3.9445149
- [40,] 2.7656617 -3.3824479
- [41,] 1.4418009 -4.8754301
- [42,] 2.6147537 -1.7851870
- [43,] 2.8545262 -2.2682917
- [44,] 3.0134074 -3.7698507
- [45,] 2.3007914 -2.2159012
- [46,] 2.6254494 -3.1875920
- [47,] 3.9308435 -3.8977488
- [48,] 4.2595145 -1.5863887
- [49,] 2.2407225 -0.3600657
- [50,] 3.3714719 -2.0555639
- [51,] 3.7865929 -3.8267480
- ,-[FO ] 4 COZO4C4 O FC4770F
- [52,] 1.6070464 -3.5617735
- [53,] 2.4410630 -1.6263705 [54,] 1.3337774 -1.2496618
- [[[]]]
- [55,] 1.2557467 -3.1796362
- [56,] 3.3955168 -1.4628548
- [57,] 2.5178595 -3.7684518

```
[58,] 1.6109763 -3.6145729
[59,] 3.1012253 -1.8611433
[60,] 2.4568767 -4.1792452
```

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] 60.94494 60.94494

(between\_SS / total\_SS = 88.4 %)

Available components:

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

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

The kmeans() function returns 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 are in each cluster?

#### km\$size

[1] 30 30

Cluster assignment/membership vector?

#### km\$cluster

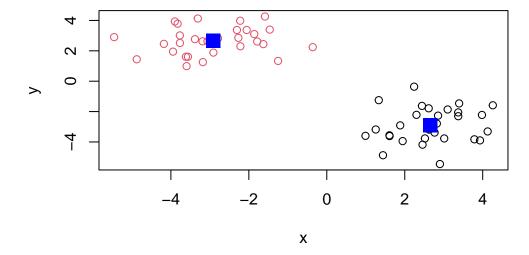
Q. Cluster centers?

#### km\$centers

```
x y
1 2.653054 -2.910387
2 -2.910387 2.653054
```

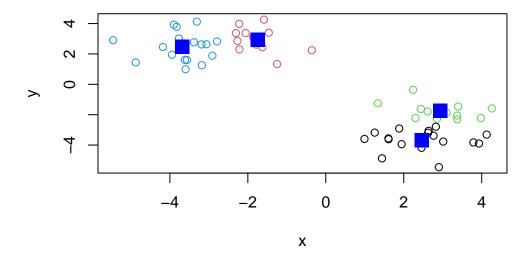
Q. Make a plot of our kmeans() results showing cluster assignment using different colors for each cluster/group of points and cluster centers in blue.

```
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.

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



**key-point**: K-means clustering is super popular but can be misused. 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 "distance 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)

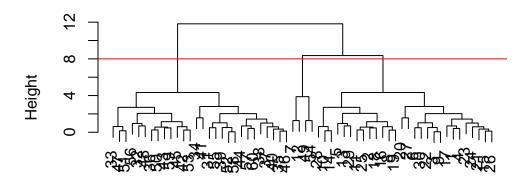
Cluster method : complete
Distance : euclidean

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**



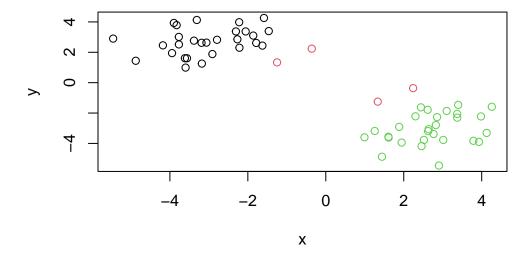
d hclust (\*, "complete")

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)
```

grps 1 2 3 28 4 28



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

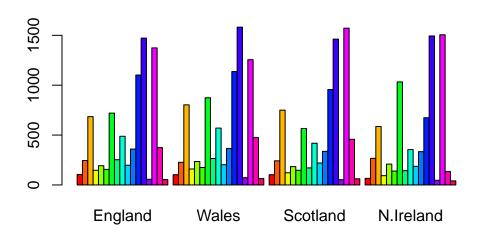
```
url <- "https://tinyurl.com/UK-foods"
x <- read.csv(url)
head(x)</pre>
```

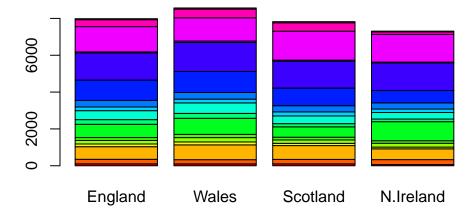
X England Wales Scotland N.Ireland Cheese Carcass\_meat Other\_meat Fish 5 Fats\_and\_oils Sugars 

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

	England	Wales	${\tt Scotland}$	${\tt N.Ireland}$
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

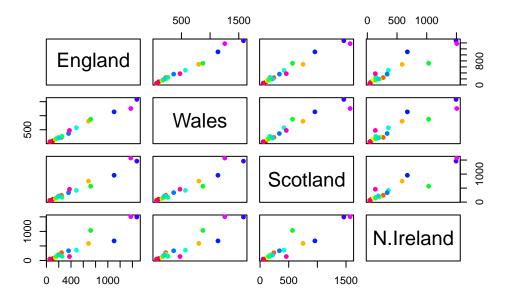
barplot(as.matrix(x), beside=T, col=rainbow(nrow(x)))





One conventional plot that can be useful is called a "pairs" 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 results with five attributes/components.

```
attributes(pca)
```

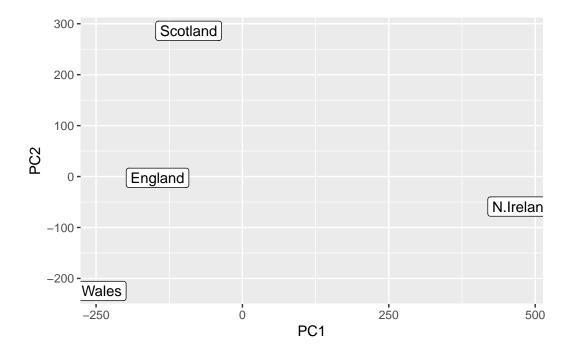
```
$names
[1] "sdev"          "rotation" "center"          "scale"          "x"
$class
[1] "prcomp"
```

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
```

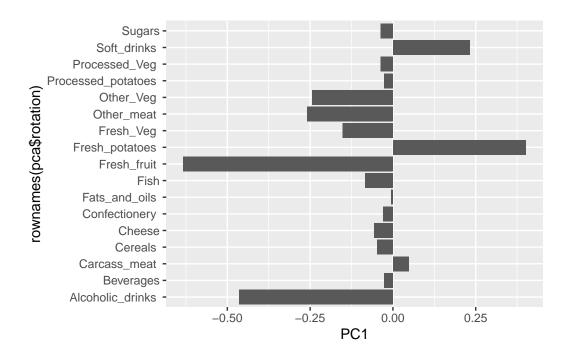
```
# Make a plot of pca$x with PC1 vs PC2
# ggrepel , geom_text_repel
ggplot(pca$x) +
   aes(PC1, PC2, label=rownames(pca$x)) +
   geom_point() +
   geom_label()
```



The graph above shows that the results from N. Ireland are most different compared to the other three countries.

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()
```



The graph above shows that, based on the previous graph, soft drinks and fresh potatoes are more likely to be associated with N. Ireland and every other food are more likely to be associated with the other three countries.