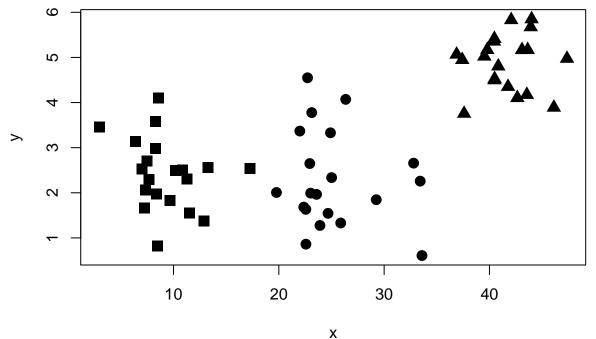
# Exploring the Number of Clusters

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#### Create synthetic data

First we create some synthetic data using the rnorm() function. We create 3 distributions with centers (10, 3), (27, 2) and (41, 5). These are the "true" clusters but the regions overlap a little. We plot the unclustered data with different shapes for each distribution.

```
set.seed(1234)
x <- rep(0, 60)
y <- rep(0, 60)
x[1:20] <- rnorm(20, mean=10, sd=3)
y[1:20] <- rnorm(20, mean=3, sd=1)
x[21:40] <- rnorm(20, mean=27, sd=4)
y[21:40] <- rnorm(20, mean=2, sd=1)
x[41:60] <- rnorm(20, mean=41, sd=3)
y[41:60] <- rnorm(20, mean=5, sd=1)
# uncomment the next two lines to see what happens
# with a more uniform distribution
#x <- rnorm(60, mean=30, sd=10)
#y <- rnorm(60, mean=3, sd=2)
true <- c(rep(1,20), rep(2,20), rep(3,20))
plot(x, y, cex=1.5, pch=c(15, 16, 17)[true])</pre>
```



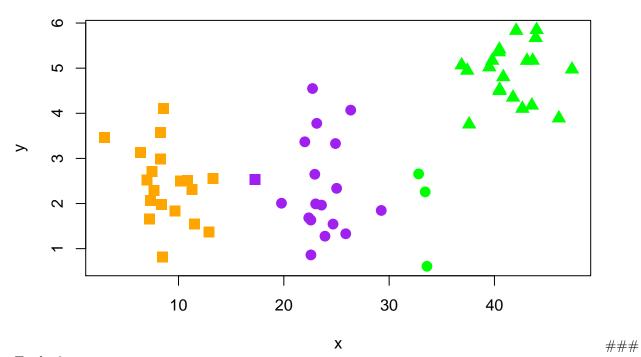
#### k-means: one iteration

Apply the k-means algoirthm with only one iteration and one start.

k-means: unlimited iterations

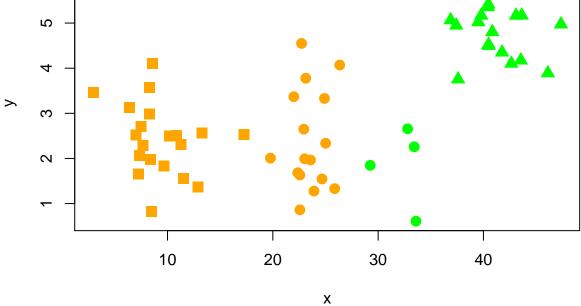
Although when we ran one iteration we got a warning message that it did not converge, we se no change when we let it run as many iterations as needed. Typing res2\$iter at the console shows that it only ran 2 iterations.

```
set.seed(1234)
res3 <- kmeans(df, 3, nstart=1 )
plot(x, y, col=c("orange", "green", "purple")[res3$cluster], cex=1.5, pch=c(15, 16, 17)[true])</pre>
```



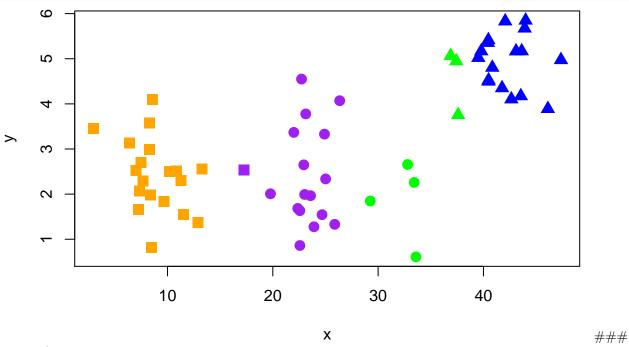
Try k=2

```
set.seed(1234)
res2 <- kmeans(df, 2, nstart= 5)
plot(x, y, col=c("orange", "green", "purple", "blue")[res2$cluster], cex=1.5, pch=c(15, 16, 17)[true])</pre>
```



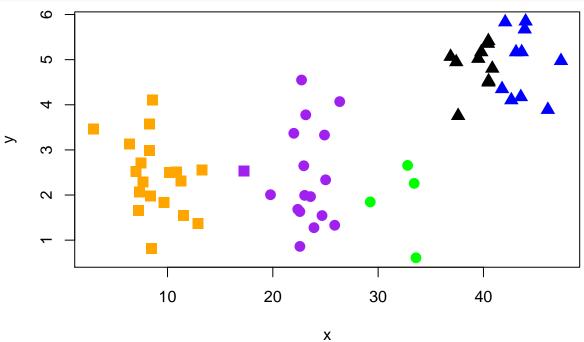
Try k=4

```
set.seed(1234)
res4 <- kmeans(df, 4, nstart= 5)
plot(x, y, col=c("orange", "green", "purple", "blue")[res4$cluster], cex=1.5, pch=c(15, 16, 17)[true])</pre>
```



 ${\rm Try}\ 5\ {\rm clusters}$ 

```
set.seed(1234)
res5 <- kmeans(df, 5, nstart= 5)
plot(x, y, col=c("orange", "green", "purple", "blue", "black")[res5$cluster], cex=1.5, pch=c(15, 16, 17)</pre>
```



withinss

Our goal is to reduce within sum of squares, this means our clusters are more homogenous. Let's compare the withinss for k=2, k=4 and k=5.

###

It seems there is a dramatic drop from k=2 to k=3 then it gradually decreases. It makes sense that the larger

the number of clusters, the smaller the withinss. After all, if k=n then withinss would be 0.

```
print(paste("k=2: ", sum(res2$withinss)))

## [1] "k=2: 2530.74905628694"

print(paste("k=3: ", sum(res3$withinss)))

## [1] "k=3: 611.695448018061"

print(paste("k=4: ", sum(res4$withinss)))

## [1] "k=4: 373.0421592289"

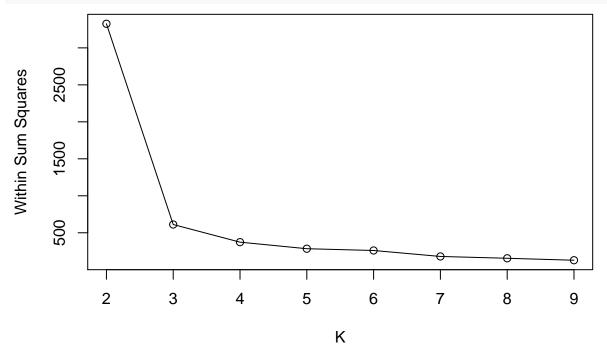
print(paste("k=5: ", sum(res5$withinss)))

## [1] "k=5: 284.426492959912"
```

### Finding k with a function

We can write a function to randomly try different k values and plot the within sum of squares.

```
plot_withinss <- function(df, max_clusters){
  withinss <- rep(0, max_clusters-1)
  for (i in 2:max_clusters){
    set.seed(1234)
    withinss[i] <- sum(kmeans(df, i)$withinss)
  }
  plot(2:max_clusters, withinss[2:max_clusters], type="o", xlab="K", ylab="Within Sum Squares")
}
plot_withinss(df, 9)</pre>
```



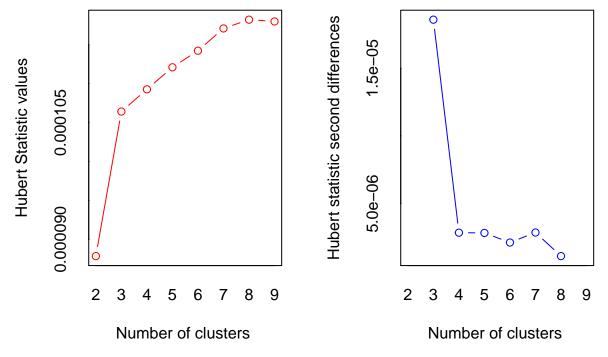
# NbClust()

##

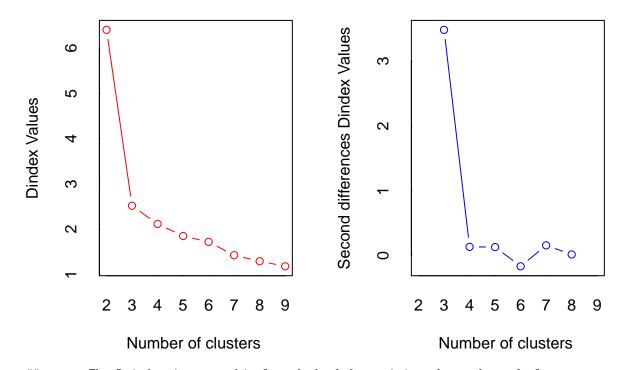
Next we try the NbClust() function to find the best number of clusters.

```
library(NbClust)
set.seed(1234)
nc <- NbClust(df, min.nc=2, max.nc=9, method="kmeans")</pre>
```

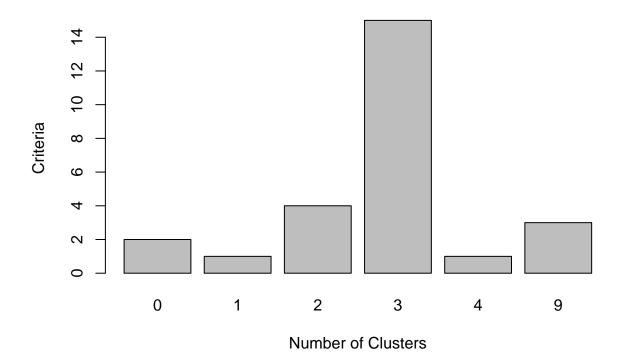
## Warning in pf(beale, pp, df2): NaNs produced



## \*\*\* : The Hubert index is a graphical method of determining the number of clusters.
## In the plot of Hubert index, we seek a significant knee that corresponds to a
## significant increase of the value of the measure i.e the significant peak in Hubert
## index second differences plot.



```
*** : The D index is a graphical method of determining the number of clusters.
                  In the plot of D index, we seek a significant knee (the significant peak in Dindex
##
##
                  second differences plot) that corresponds to a significant increase of the value of
##
                  the measure.
##
  ************************
##
## * Among all indices:
## * 4 proposed 2 as the best number of clusters
## * 15 proposed 3 as the best number of clusters
## * 1 proposed 4 as the best number of clusters
## * 3 proposed 9 as the best number of clusters
##
##
                     **** Conclusion ****
##
## * According to the majority rule, the best number of clusters is 3
##
##
t <- table(nc$Best.n[1,])
t
##
##
         2
           3
                  9
      1
               4
      1
         4 15
              1
barplot(t, xlab="Number of Clusters", ylab = "Criteria")
```



# Plot for the book

