

Tipologia y ciclo de vida de los datos - PRA2

Autor: Eduardo Diaz Villanueva

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Contents

Descripción del dataset.	2
Integración y selección de los datos de interés	2
Limpieza de los datos	3
Elementos vacíos	3
Valores extremos	4
Análisis	16
Selección grupo de datos	27
Comprobación de la normalidad y homogeneidad de la varianza.	27
Aplicación de pruebas estadísticas para comparar los grupos de datos	27
Representación de los resultados	27
Resolución del problema	27

Lectura de datos

```
library(stringr)
library(dplyr)
```

```
##
## Attaching package: 'dplyr'
##
## The following objects are masked from 'package:stats':
##
##   filter, lag
##
## The following objects are masked from 'package:base':
##
##   intersect, setdiff, setequal, union
```

```
library(ggplot2)

# Limpiamos la aplicación de datos anteriores y cargamos el fichero.
rm(list = ls())
datos <- read.csv("winequality-red.csv", sep=",")
datos_originales <- datos
head(datos,5)
```

```
##   fixed.acidity volatile.acidity citric.acid residual.sugar chlorides
## 1          7.4           0.70         0.00           1.9       0.076
## 2          7.8           0.88         0.00           2.6       0.098
```

```
## 3      7.8      0.76      0.04      2.3      0.092
## 4     11.2      0.28      0.56      1.9      0.075
## 5      7.4      0.70      0.00      1.9      0.076
##  free.sulfur.dioxide total.sulfur.dioxide density  pH sulphates alcohol
## 1             11             34 0.9978 3.51      0.56      9.4
## 2             25             67 0.9968 3.20      0.68      9.8
## 3             15             54 0.9970 3.26      0.65      9.8
## 4             17             60 0.9980 3.16      0.58      9.8
## 5             11             34 0.9978 3.51      0.56      9.4
##  quality
## 1      5
## 2      5
## 3      5
## 4      6
## 5      5
```

Descripción del dataset.

El dataset seleccionado contiene 11 variables que describen las propiedades químicas de un vino, como puede ser la acidez, pH nivel de azúcar, etc... estas variables tendrán influencia en la calidad final del vino.

Con este ejercicio queremos estudiar que variables son más representativas y encontrar modelos que puedan predecir la calidad del vino.

Si pensamos por ejemplo en una industria, podríamos reducir el tiempo y coste reduciendo el número de pruebas de calidad a las variables más significativas. Incluso mejorar la calidad del producto final, focalizando esfuerzos y recursos a reducir la variabilidad de las variables que más contribuyan a la calidad final.

Integración y selección de los datos de interés

Realizaremos un primer análisis estadístico para familiarizarnos con las variables.

```
summary(datos)
```

```
## fixed.acidity  volatile.acidity  citric.acid  residual.sugar
## Min.   : 4.60  Min.   :0.1200  Min.   :0.000  Min.   : 0.900
## 1st Qu.: 7.10  1st Qu.:0.3900  1st Qu.:0.090  1st Qu.: 1.900
## Median : 7.90  Median :0.5200  Median :0.260  Median : 2.200
## Mean   : 8.32  Mean   :0.5278  Mean   :0.271  Mean   : 2.539
## 3rd Qu.: 9.20  3rd Qu.:0.6400  3rd Qu.:0.420  3rd Qu.: 2.600
## Max.   :15.90  Max.   :1.5800  Max.   :1.000  Max.   :15.500
## chlorides      free.sulfur.dioxide total.sulfur.dioxide density
## Min.   :0.01200  Min.   : 1.00      Min.   : 6.00      Min.   :0.9901
## 1st Qu.:0.07000  1st Qu.: 7.00      1st Qu.: 22.00     1st Qu.:0.9956
## Median :0.07900  Median :14.00     Median : 38.00     Median :0.9968
## Mean   :0.08747  Mean   :15.87     Mean   : 46.47     Mean   :0.9967
## 3rd Qu.:0.09000  3rd Qu.:21.00     3rd Qu.: 62.00     3rd Qu.:0.9978
## Max.   :0.61100  Max.   :72.00     Max.   :289.00     Max.   :1.0037
## pH             sulphates          alcohol          quality
## Min.   :2.740  Min.   :0.3300  Min.   : 8.40      Min.   :3.000
## 1st Qu.:3.210  1st Qu.:0.5500  1st Qu.: 9.50      1st Qu.:5.000
## Median :3.310  Median :0.6200  Median :10.20     Median :6.000
## Mean   :3.311  Mean   :0.6581  Mean   :10.42     Mean   :5.636
## 3rd Qu.:3.400  3rd Qu.:0.7300  3rd Qu.:11.10     3rd Qu.:6.000
## Max.   :4.010  Max.   :2.0000  Max.   :14.90     Max.   :8.000
```

```
str(datos)
```

```
## 'data.frame': 1599 obs. of 12 variables:
## $ fixed.acidity : num 7.4 7.8 7.8 11.2 7.4 7.4 7.9 7.3 7.8 7.5 ...
## $ volatile.acidity : num 0.7 0.88 0.76 0.28 0.7 0.66 0.6 0.65 0.58 0.5 ...
## $ citric.acid : num 0 0 0.04 0.56 0 0 0.06 0 0.02 0.36 ...
## $ residual.sugar : num 1.9 2.6 2.3 1.9 1.9 1.8 1.6 1.2 2 6.1 ...
## $ chlorides : num 0.076 0.098 0.092 0.075 0.076 0.075 0.069 0.065 0.073 0.071 ...
## $ free.sulfur.dioxide : num 11 25 15 17 11 13 15 15 9 17 ...
## $ total.sulfur.dioxide: num 34 67 54 60 34 40 59 21 18 102 ...
## $ density : num 0.998 0.997 0.997 0.998 0.998 ...
## $ pH : num 3.51 3.2 3.26 3.16 3.51 3.51 3.3 3.39 3.36 3.35 ...
## $ sulphates : num 0.56 0.68 0.65 0.58 0.56 0.56 0.46 0.47 0.57 0.8 ...
## $ alcohol : num 9.4 9.8 9.8 9.8 9.4 9.4 9.4 10 9.5 10.5 ...
## $ quality : int 5 5 5 6 5 5 5 7 7 5 ...
```

```
#Tipo de dato asignado a cada campo
sapply(datos, function(x) class(x))
```

```
##      fixed.acidity    volatile.acidity    citric.acid
##      "numeric"        "numeric"          "numeric"
##      residual.sugar    chlorides    free.sulfur.dioxide
##      "numeric"        "numeric"        "numeric"
## total.sulfur.dioxide    density          pH
##      "numeric"        "numeric"        "numeric"
##      sulphates        alcohol          quality
##      "numeric"        "numeric"        "integer"
```

Limpieza de los datos

Elementos vacios

Analizamos los valores para detectar falta o ausencia de daos

```
# Analizamos la existencia de datos NA
colSums(is.na(datos))
```

```
##      fixed.acidity    volatile.acidity    citric.acid
##      0              0              0
##      residual.sugar    chlorides    free.sulfur.dioxide
##      0              0              0
## total.sulfur.dioxide    density          pH
##      0              0              0
##      sulphates        alcohol          quality
##      0              0              0
```

```
# Analizamos la existencia de datos vacios
colSums(datos=="")
```

```
##      fixed.acidity    volatile.acidity    citric.acid
##      0              0              0
##      residual.sugar    chlorides    free.sulfur.dioxide
##      0              0              0
## total.sulfur.dioxide    density          pH
##      0              0              0
##      sulphates        alcohol          quality
```

```
##          0          0          0
```

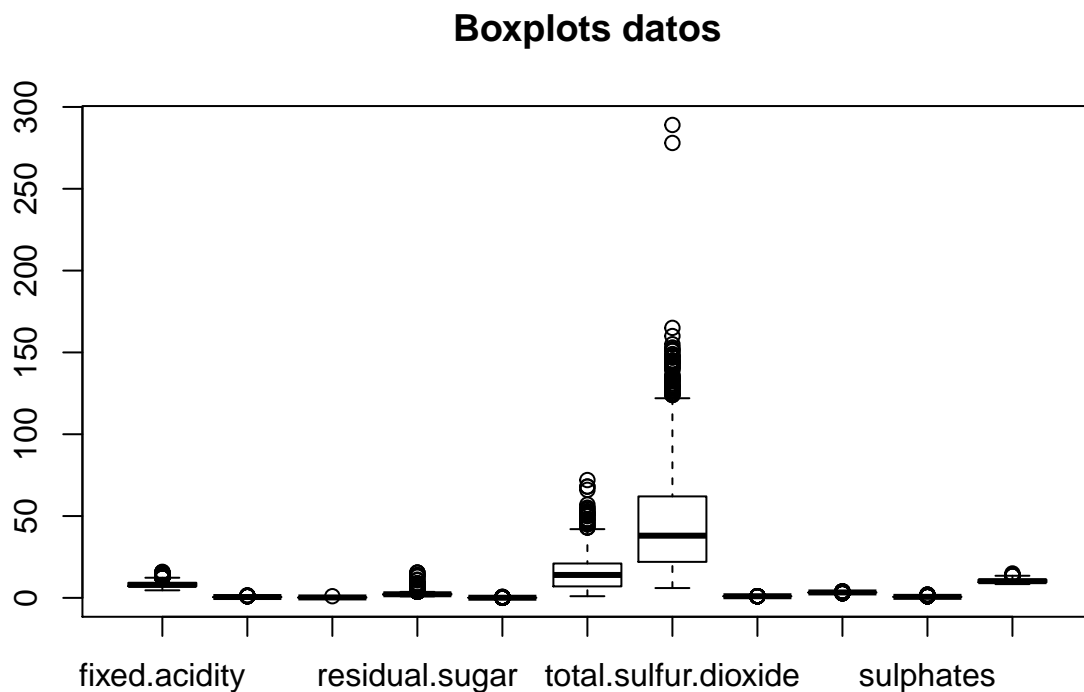
```
# Analizamos la existencia de datos con valor 0
colSums(datos==0)
```

```
##      fixed.acidity    volatile.acidity      citric.acid
##           0           0           132
##      residual.sugar      chlorides  free.sulfur.dioxide
##           0           0           0
## total.sulfur.dioxide      density           pH
##           0           0           0
##      sulphates      alcohol      quality
##           0           0           0
```

Observamos la variable Citric.acid con una gran cantidad de valores 0.

Valores extremos

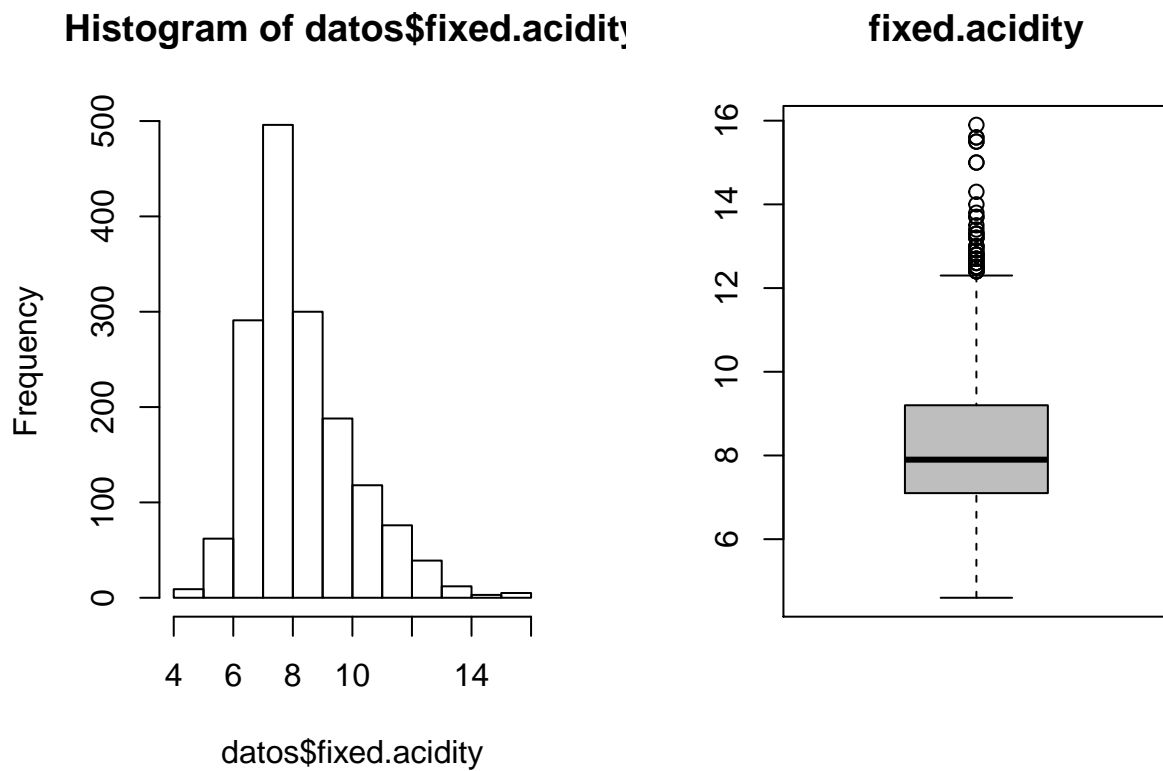
```
boxplot(x=datos[,1:11],main="Boxplots datos")
```



```
#boxplot(x=datos[,1:5],main="Boxplots datos")
#boxplot(x=datos[,6:7],main="Boxplots datos")
#boxplot(x=datos[,8:12],main="Boxplots datos")
```

```
par(mfrow=c(1,2))
hist(datos$fixed.acidity)
```

```
boxplot(datos$fixed.acidity,main="fixed.acidity", col="gray")
```

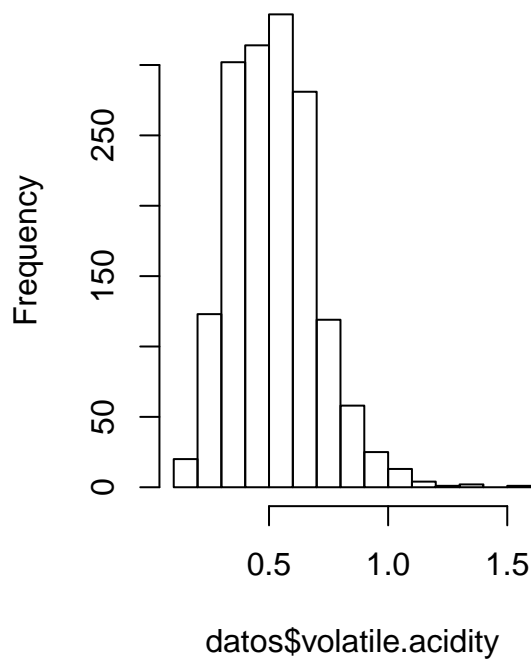


```
boxplot.stats(datos$fixed.acidity)$out
```

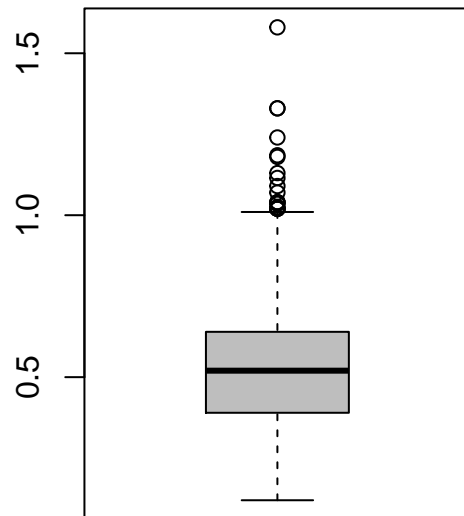
```
## [1] 12.8 12.8 15.0 15.0 12.5 13.3 13.4 12.4 12.5 13.8 13.5 12.6 12.5 12.8 12.8
## [16] 14.0 13.7 13.7 12.7 12.5 12.8 12.6 15.6 12.5 13.0 12.5 13.3 12.4 12.5 12.9
## [31] 14.3 12.4 15.5 15.5 15.6 13.0 12.7 13.0 12.7 12.4 12.7 13.2 13.2 13.2 15.9
## [46] 13.3 12.9 12.6 12.6
```

```
par(mfrow=c(1,2))
hist(datos$volatile.acidity)
boxplot(datos$volatile.acidity,main="volatile.acidity", col="gray")
```

Histogram of datos\$volatile.acidi



volatile.acidity

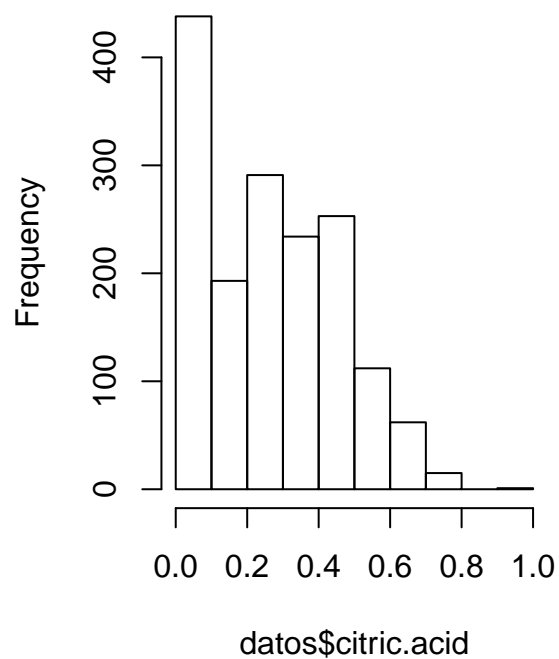


```
boxplot.stats(datos$volatile.acidity)$out
```

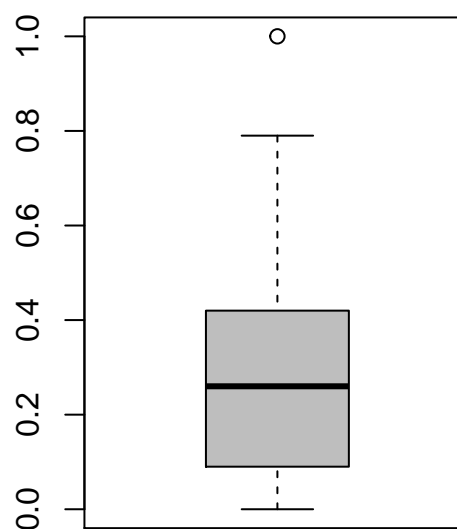
```
## [1] 1.130 1.020 1.070 1.330 1.330 1.040 1.090 1.040 1.240 1.185 1.020 1.035
## [13] 1.025 1.115 1.020 1.020 1.580 1.180 1.040
```

```
par(mfrow=c(1,2))
hist(datos$citric.acid )
boxplot(datos$citric.acid ,main="citric.acid ", col="gray")
```

Histogram of datos\$citric.acid



citric.acid

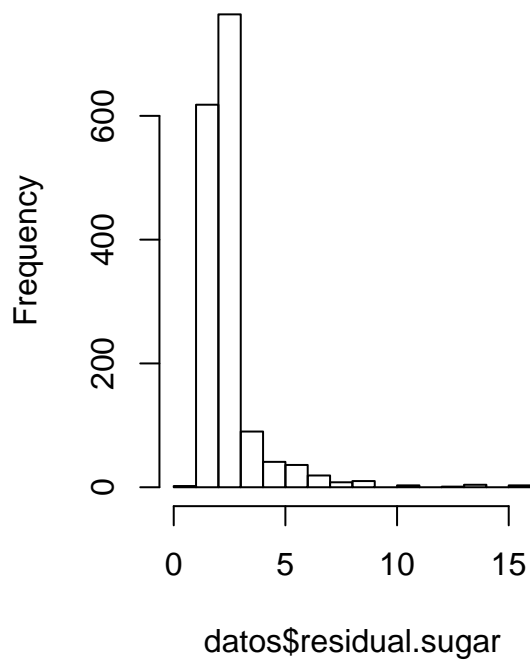


```
boxplot.stats(datos$citric.acid )$out
```

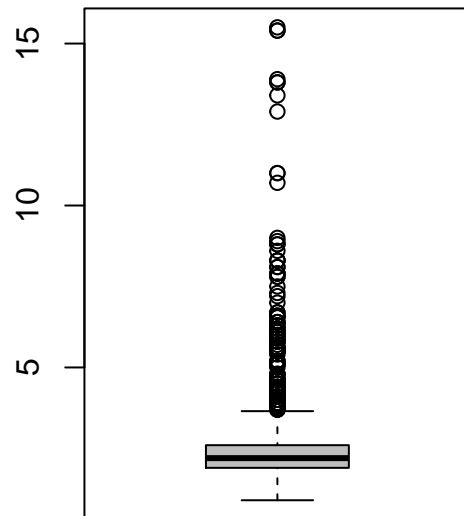
```
## [1] 1
```

```
par(mfrow=c(1,2))  
hist(datos$residual.sugar)  
boxplot(datos$residual.sugar,main="residual.sugar", col="gray")
```

Histogram of datos\$residual.sug



residual.sugar

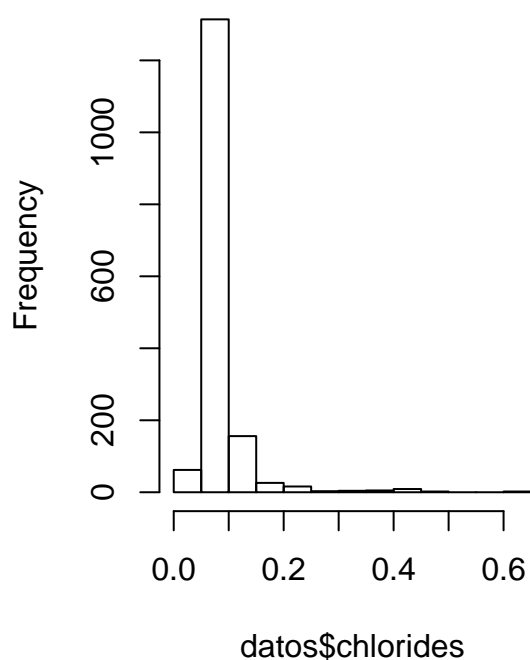


```
boxplot.stats(datos$residual.sugar)$out
```

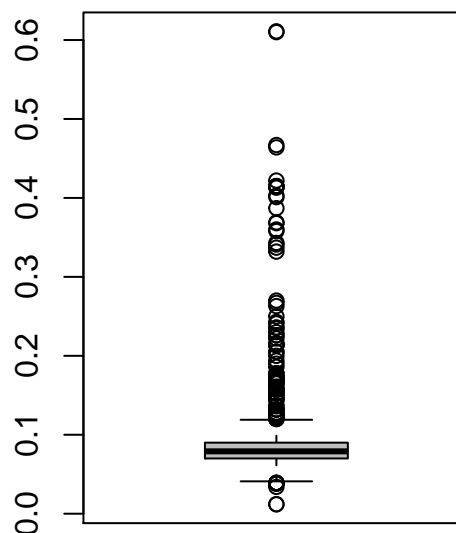
```
## [1] 6.10 6.10 3.80 3.90 4.40 10.70 5.50 5.90 5.90 3.80 5.10 4.65
## [13] 4.65 5.50 5.50 5.50 5.50 7.30 7.20 3.80 5.60 4.00 4.00 4.00
## [25] 4.00 7.00 4.00 4.00 6.40 5.60 5.60 11.00 11.00 4.50 4.80 5.80
## [37] 5.80 3.80 4.40 6.20 4.20 7.90 7.90 3.70 4.50 6.70 6.60 3.70
## [49] 5.20 15.50 4.10 8.30 6.55 6.55 4.60 6.10 4.30 5.80 5.15 6.30
## [61] 4.20 4.20 4.60 4.20 4.60 4.30 4.30 7.90 4.60 5.10 5.60 5.60
## [73] 6.00 8.60 7.50 4.40 4.25 6.00 3.90 4.20 4.00 4.00 4.00 6.60
## [85] 6.00 6.00 3.80 9.00 4.60 8.80 8.80 5.00 3.80 4.10 5.90 4.10
## [97] 6.20 8.90 4.00 3.90 4.00 8.10 8.10 6.40 6.40 8.30 8.30 4.70
## [109] 5.50 5.50 4.30 5.50 3.70 6.20 5.60 7.80 4.60 5.80 4.10 12.90
## [121] 4.30 13.40 4.80 6.30 4.50 4.50 4.30 4.30 3.90 3.80 5.40 3.80
## [133] 6.10 3.90 5.10 5.10 3.90 15.40 15.40 4.80 5.20 5.20 3.75 13.80
## [145] 13.80 5.70 4.30 4.10 4.10 4.40 3.70 6.70 13.90 5.10 7.80
```

```
par(mfrow=c(1,2))
hist(datos$chlorides)
boxplot(datos$chlorides,main="chlorides", col="gray")
```


Histogram of datos\$chlorides



chlorides

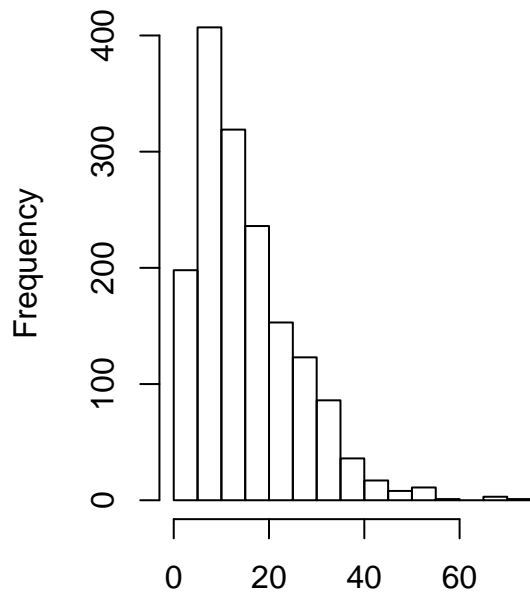


```
boxplot.stats(datos$chlorides)$out
```

```
## [1] 0.176 0.170 0.368 0.341 0.172 0.332 0.464 0.401 0.467 0.122 0.178 0.146
## [13] 0.236 0.610 0.360 0.270 0.039 0.337 0.263 0.611 0.358 0.343 0.186 0.213
## [25] 0.214 0.121 0.122 0.122 0.128 0.120 0.159 0.124 0.122 0.122 0.174 0.121
## [37] 0.127 0.413 0.152 0.152 0.125 0.122 0.200 0.171 0.226 0.226 0.250 0.148
## [49] 0.122 0.124 0.124 0.143 0.222 0.039 0.157 0.422 0.034 0.387 0.415 0.157
## [61] 0.157 0.243 0.241 0.190 0.132 0.126 0.038 0.165 0.145 0.147 0.012 0.012
## [73] 0.039 0.194 0.132 0.161 0.120 0.120 0.123 0.123 0.414 0.216 0.171 0.178
## [85] 0.369 0.166 0.166 0.136 0.132 0.132 0.123 0.123 0.123 0.403 0.137 0.414
## [97] 0.166 0.168 0.415 0.153 0.415 0.267 0.123 0.214 0.214 0.169 0.205 0.205
## [109] 0.039 0.235 0.230 0.038
```

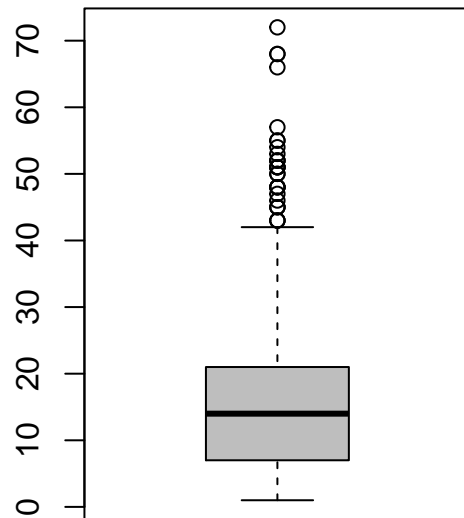
```
par(mfrow=c(1,2))
hist(datos$free.sulfur.dioxide)
boxplot(datos$free.sulfur.dioxide,main="free.sulfur.dioxide", col="gray")
```

Histogram of datos\$free.sulfur.dioxide



datos\$free.sulfur.dioxide

free.sulfur.dioxide

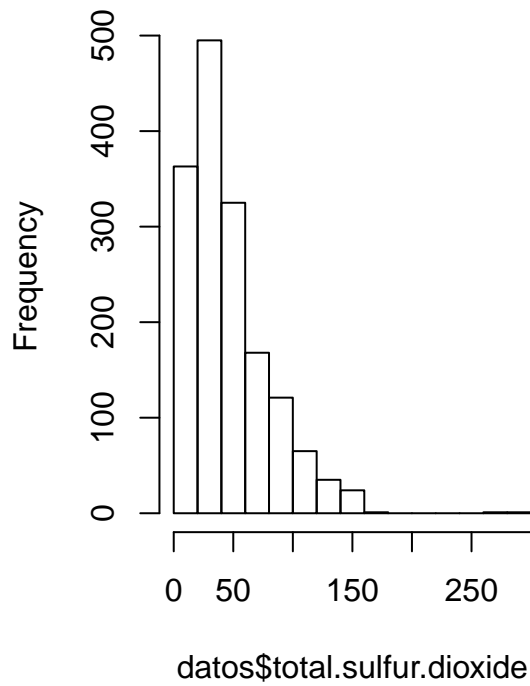


```
boxplot.stats(datos$free.sulfur.dioxide)$out
```

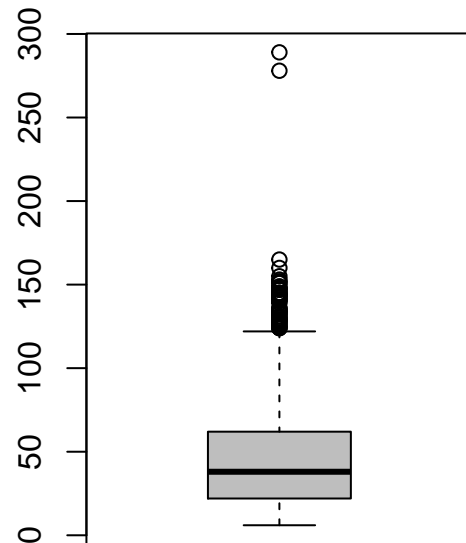
```
## [1] 52 51 50 68 68 43 47 54 46 45 53 52 51 45 57 50 45 48 43 48 72 43 51 51 52
## [26] 55 55 48 48 66
```

```
par(mfrow=c(1,2))
hist(datos$total.sulfur.dioxide)
boxplot(datos$total.sulfur.dioxide,main="total.sulfur.dioxide", col="gray")
```

Histogram of datos\$total.sulfur.dio



total.sulfur.dioxide

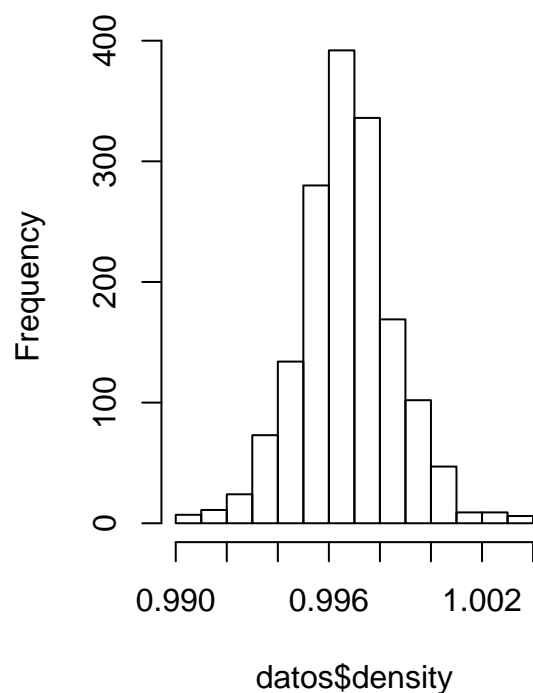


```
boxplot.stats(datos$total.sulfur.dioxide)$out
```

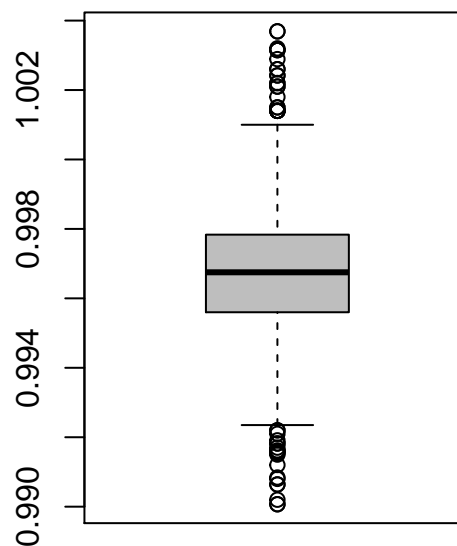
```
## [1] 145 148 136 125 140 136 133 153 134 141 129 128 129 128 143 144 127 126 145
## [20] 144 135 165 124 124 134 124 129 151 133 142 149 147 145 148 155 151 152 125
## [39] 127 139 143 144 130 278 289 135 160 141 141 133 147 147 131 131 131
```

```
par(mfrow=c(1,2))
hist(datos$density)
boxplot(datos$density,main="density", col="gray")
```

Histogram of datos\$density



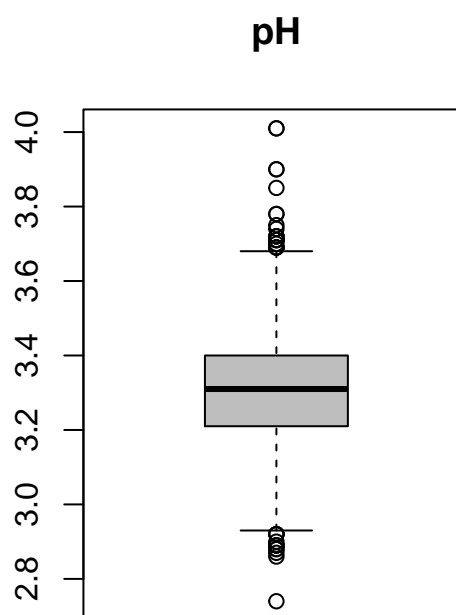
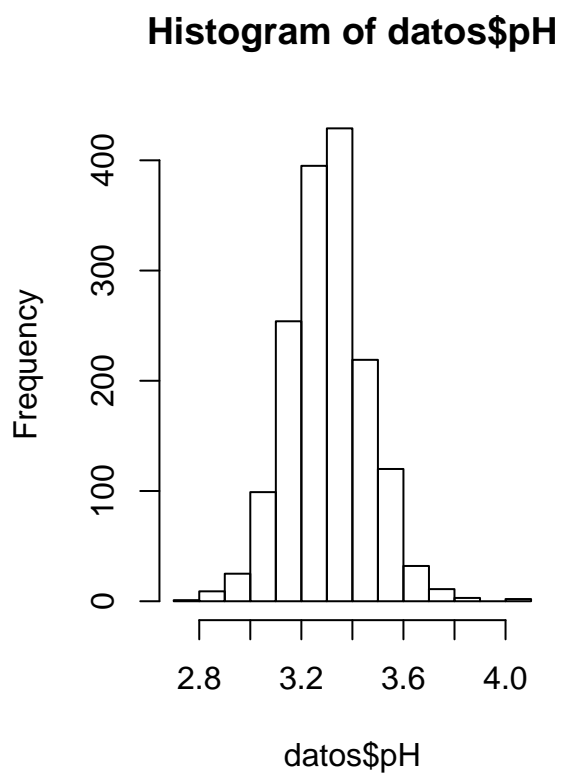
density



```
boxplot.stats(datos$density)$out
```

```
## [1] 0.99160 0.99160 1.00140 1.00150 1.00150 1.00180 0.99120 1.00220 1.00220
## [10] 1.00140 1.00140 1.00140 1.00140 1.00320 1.00260 1.00140 1.00315 1.00315
## [19] 1.00315 1.00210 1.00210 0.99170 0.99220 1.00260 0.99210 0.99154 0.99064
## [28] 0.99064 1.00289 0.99162 0.99007 0.99007 0.99020 0.99220 0.99150 0.99157
## [37] 0.99080 0.99084 0.99191 1.00369 1.00369 1.00242 0.99182 1.00242 0.99182
```

```
par(mfrow=c(1,2))
hist(datos$pH)
boxplot(datos$pH,main="pH", col="gray")
```

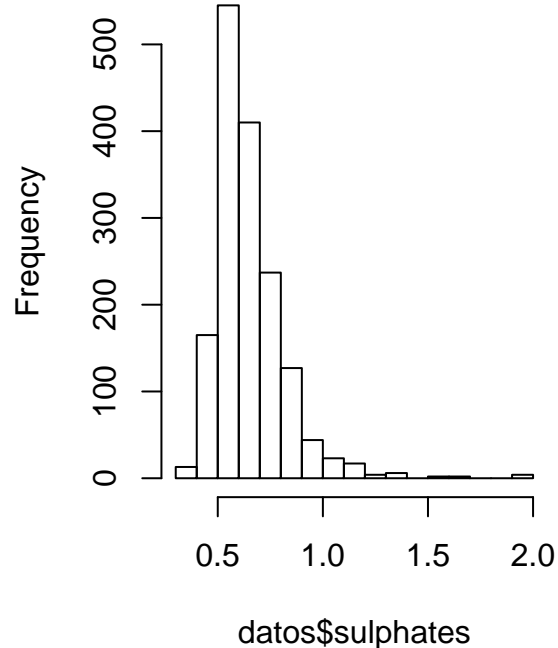


```
boxplot.stats(datos$pH)$out
```

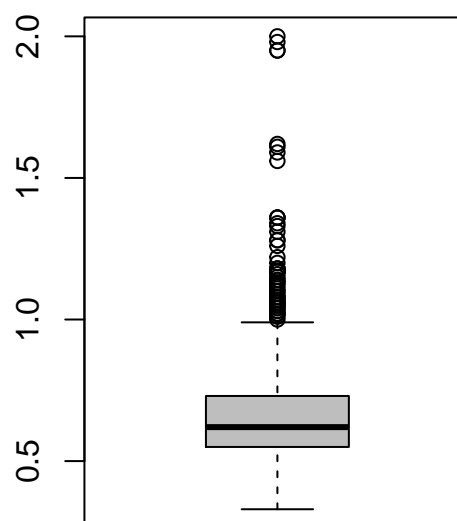
```
## [1] 3.90 3.75 3.85 2.74 3.69 3.69 2.88 2.86 3.74 2.92 2.92 2.92 3.72 2.87 2.89
## [16] 2.89 2.92 3.90 3.71 3.69 3.69 3.71 3.71 2.89 2.89 3.78 3.70 3.78 4.01 2.90
## [31] 4.01 3.71 2.88 3.72 3.72
```

```
par(mfrow=c(1,2))
hist(datos$sulphates)
boxplot(datos$sulphates,main="sulphates", col="gray")
```

Histogram of datos\$sulphates



sulphates

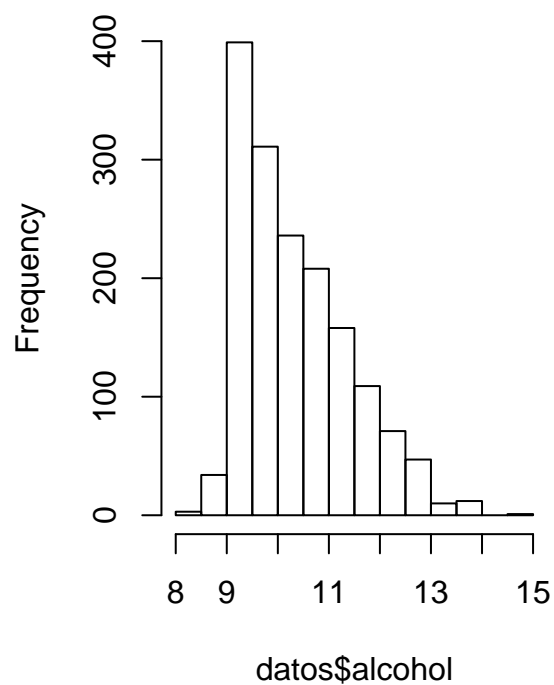


```
boxplot.stats(datos$sulphates)$out
```

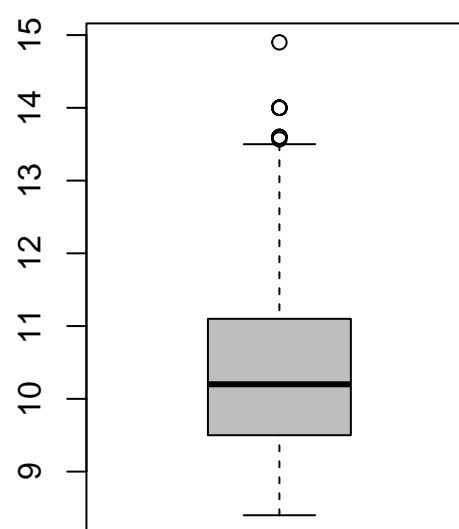
```
## [1] 1.56 1.28 1.08 1.20 1.12 1.28 1.14 1.95 1.22 1.95 1.98 1.31 2.00 1.08 1.59
## [16] 1.02 1.03 1.61 1.09 1.26 1.08 1.00 1.36 1.18 1.13 1.04 1.11 1.13 1.07 1.06
## [31] 1.06 1.05 1.06 1.04 1.05 1.02 1.14 1.02 1.36 1.36 1.05 1.17 1.62 1.06 1.18
## [46] 1.07 1.34 1.16 1.10 1.15 1.17 1.17 1.33 1.18 1.17 1.03 1.17 1.10 1.01
```

```
par(mfrow=c(1,2))
hist(datos$alcohol)
boxplot(datos$alcohol,main="alcohol", col="gray")
```

Histogram of datos\$alcohol



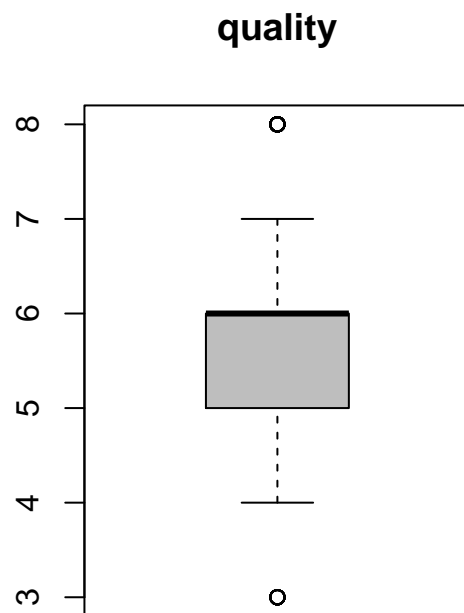
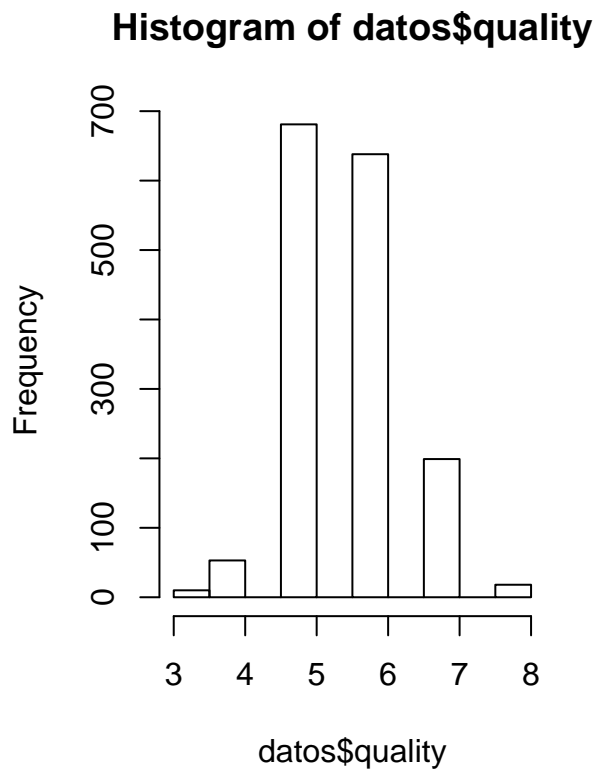
alcohol



```
boxplot.stats(datos$alcohol)$out
```

```
## [1] 14.00000 14.00000 14.00000 14.00000 14.90000 14.00000 13.60000 13.60000
## [9] 13.60000 14.00000 14.00000 13.56667 13.60000
```

```
par(mfrow=c(1,2))
hist(datos$quality)
boxplot(datos$quality,main="quality", col="gray")
```



```
boxplot.stats(datos$quality)$out
```

```
## [1] 8 8 8 8 8 3 8 8 8 3 8 3 8 3 3 8 8 8 8 8 3 3 8 8 3 3 3 8
```

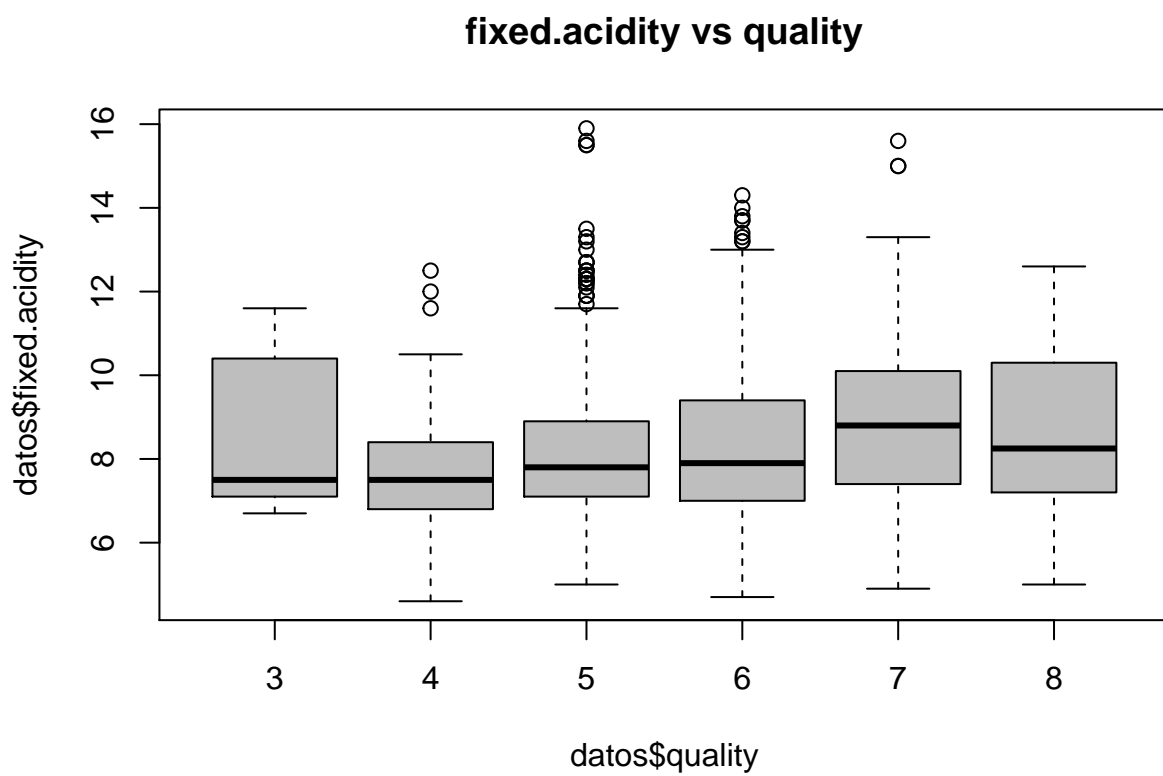
Analisis

Antes de comenzar con el analisis guardaremos una copia de los datos despues del proceso de limpieza

```
# Exportación de los datos limpios en .csv
write.csv(datos, "datos_data_clean.csv")
```

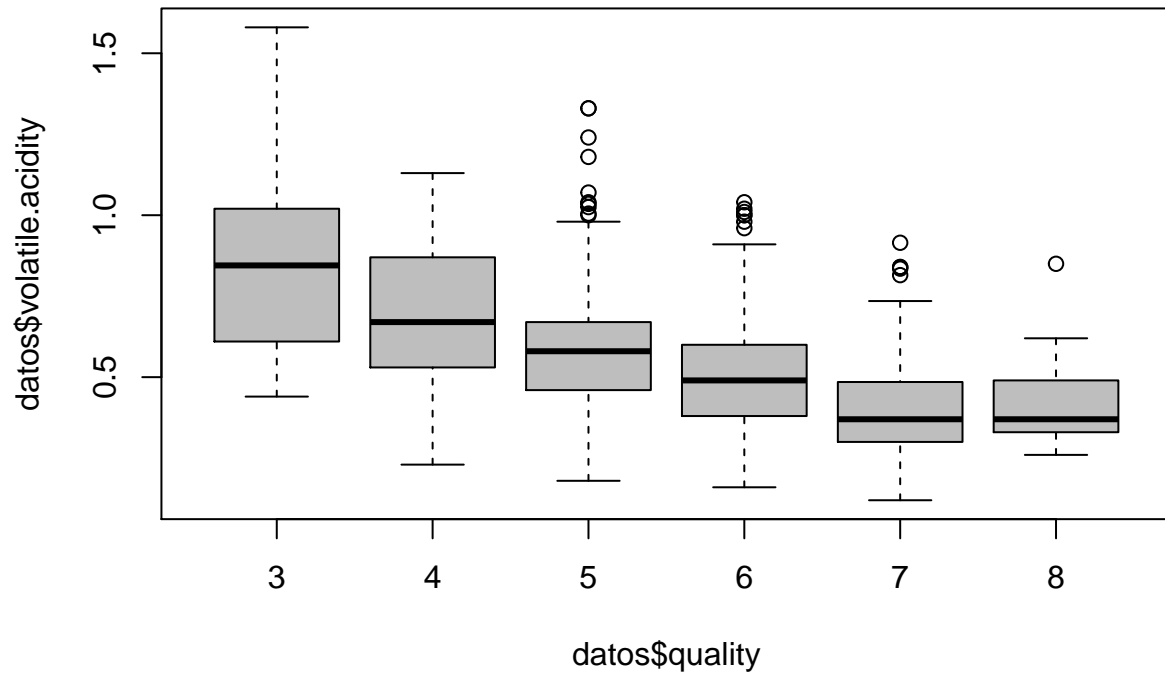
Analizaremos las variables frente a la calidad para decidir cuaes utilizar en el resto del analisis

```
#boxplot(datos$pH,main="quality", col="gray")
boxplot(formula = datos$fixed.acidity ~ datos$quality, main="fixed.acidity vs quality", col="gray")
```

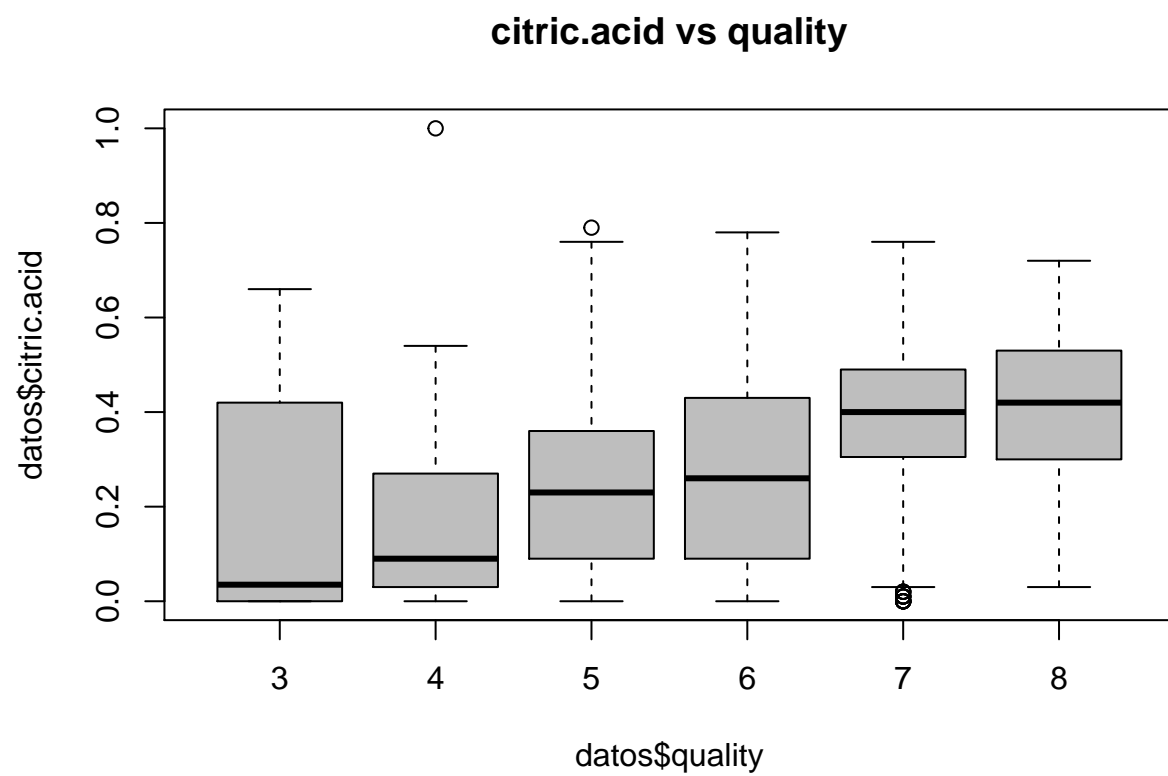



```
boxplot(formula = datos$volatile.acidity ~ datos$quality, main="volatile.acidity vs quality", col="gray"
```

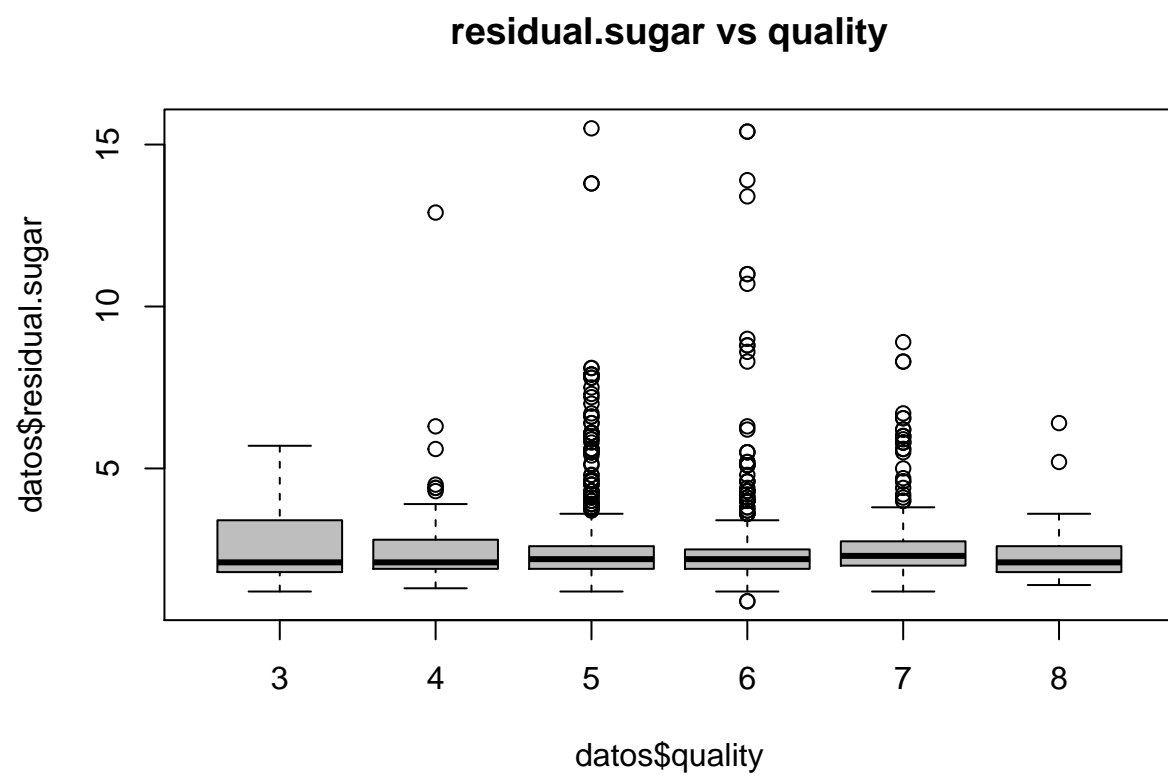
volatile.acidity vs quality



```
boxplot(formula = datos$volatile.acidity ~ datos$quality, main="volatile.acidity vs quality", col="gray")
```

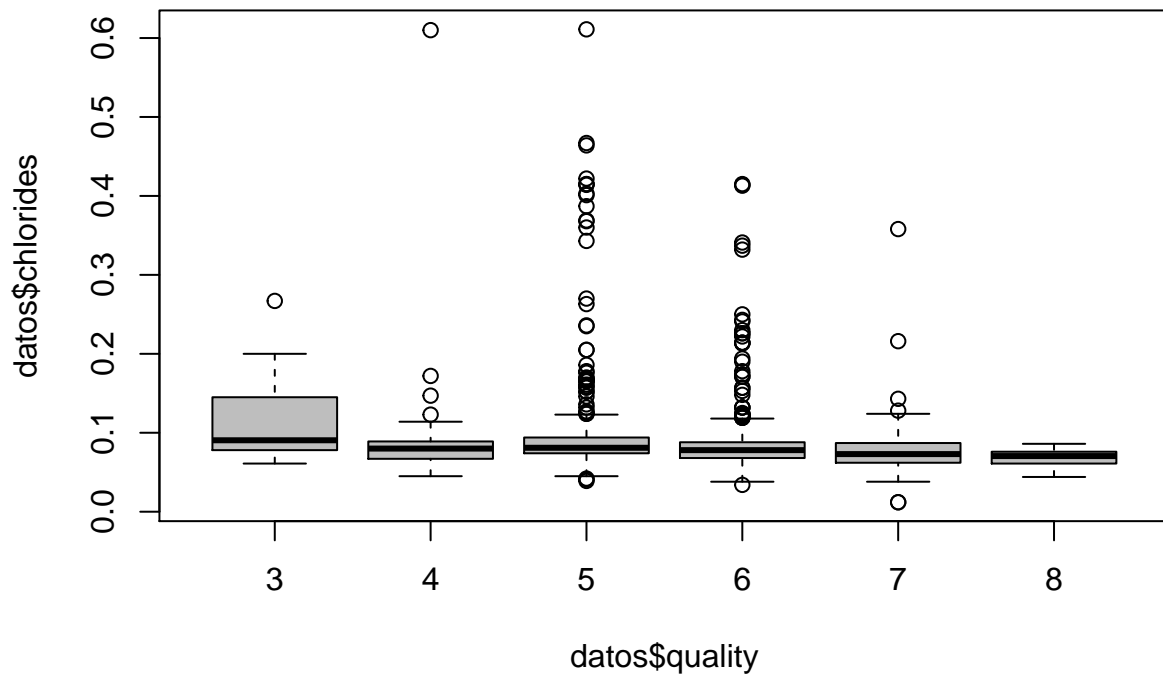


```
boxplot(formula = datos$residual.sugar ~ datos$quality, main="residual.sugar vs quality", col="gray")
```



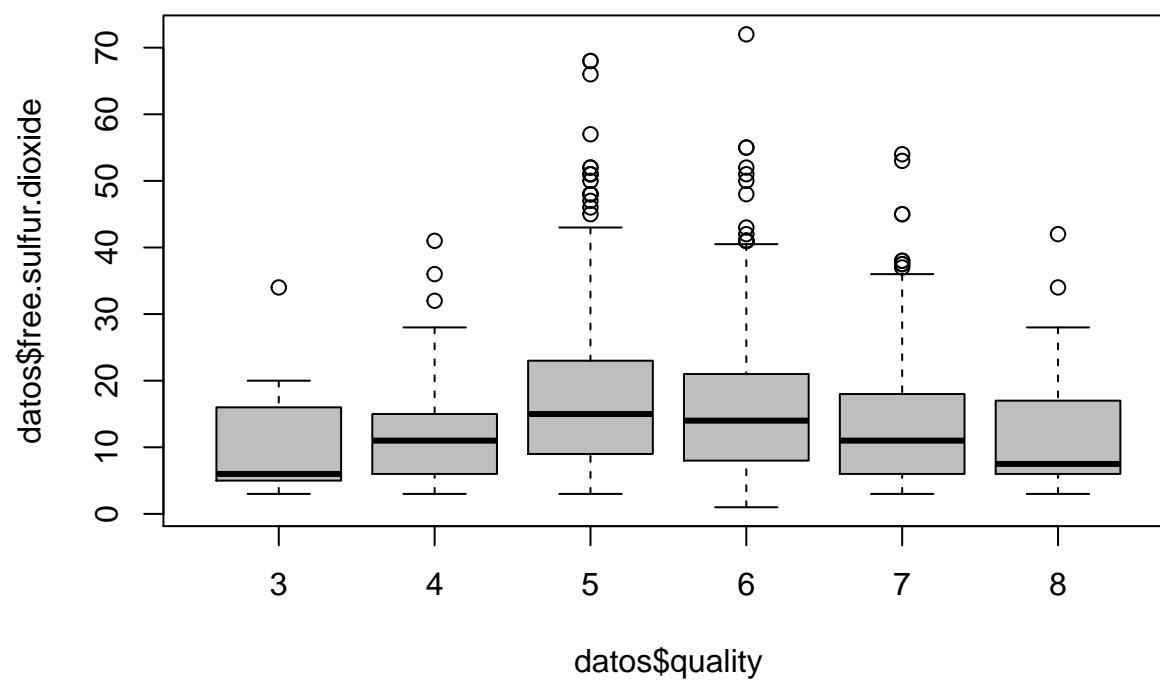
```
boxplot(formula = datos$chlorides ~ datos$quality, main="chlorides vs quality", col="gray")
```

chlorides vs quality

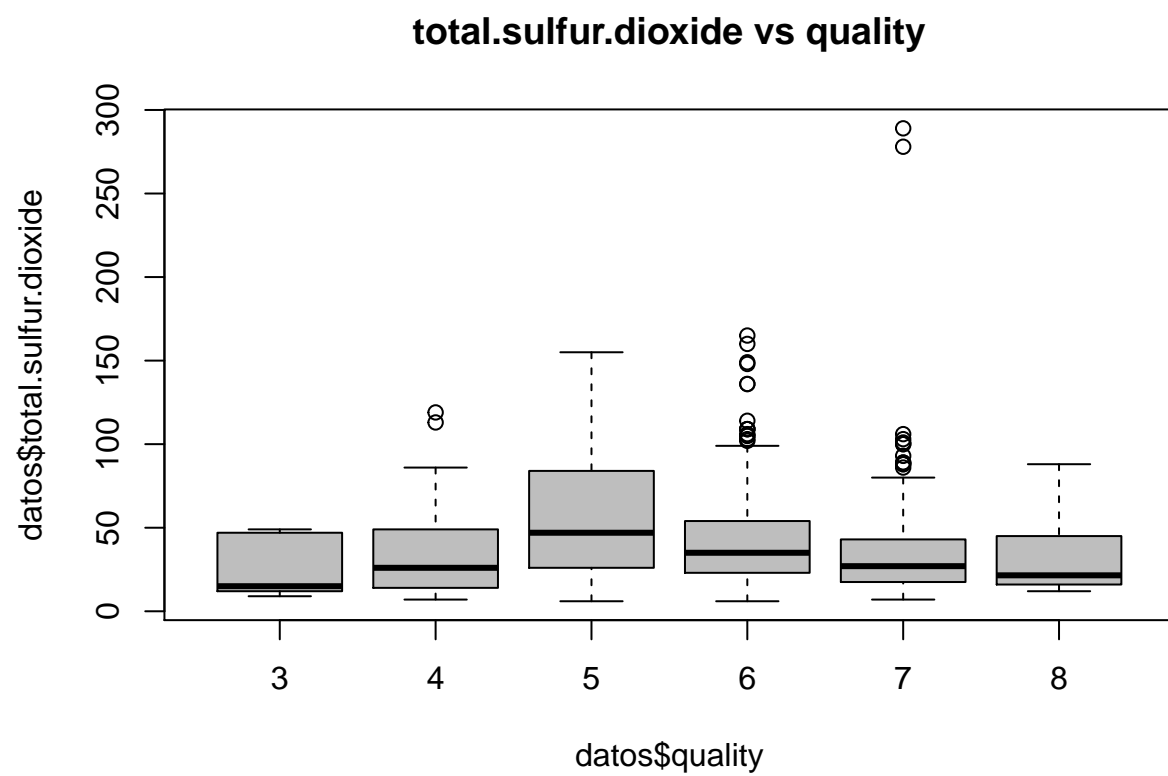


```
boxplot(formula = datos$free.sulfur.dioxide ~ datos$quality, main="free.sulfur.dioxide vs quality", col="gray")
```

free.sulfur.dioxide vs quality

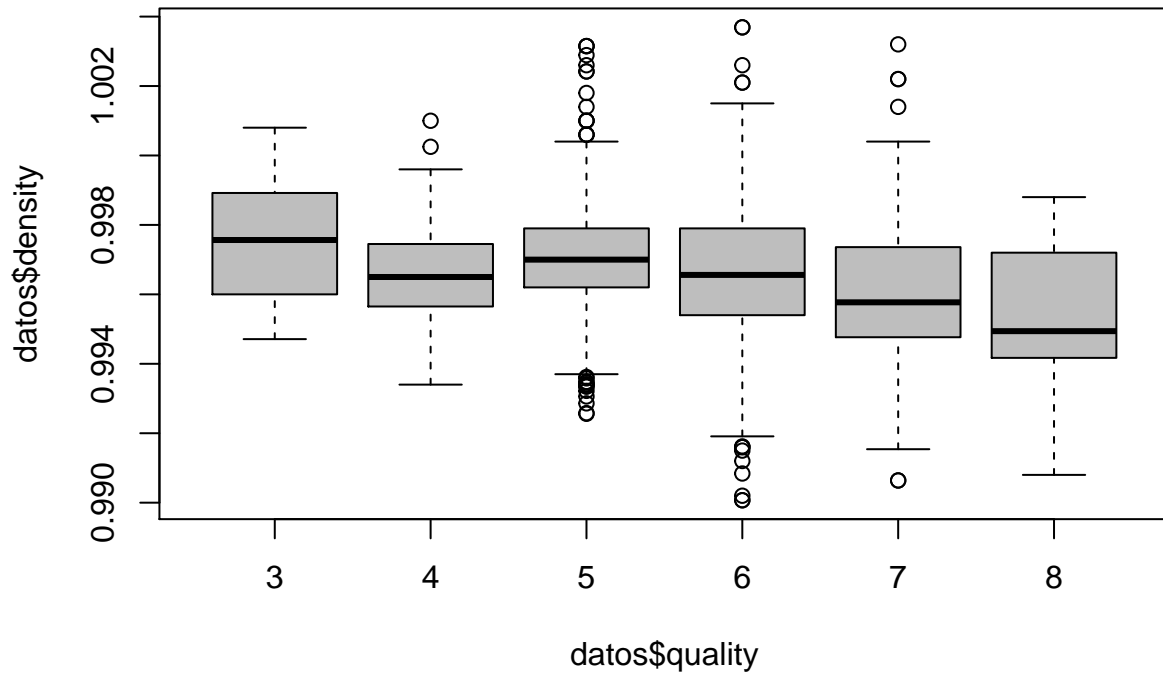


```
boxplot(formula = datos$total.sulfur.dioxide ~ datos$quality, main="total.sulfur.dioxide vs quality", col = "gray")
```

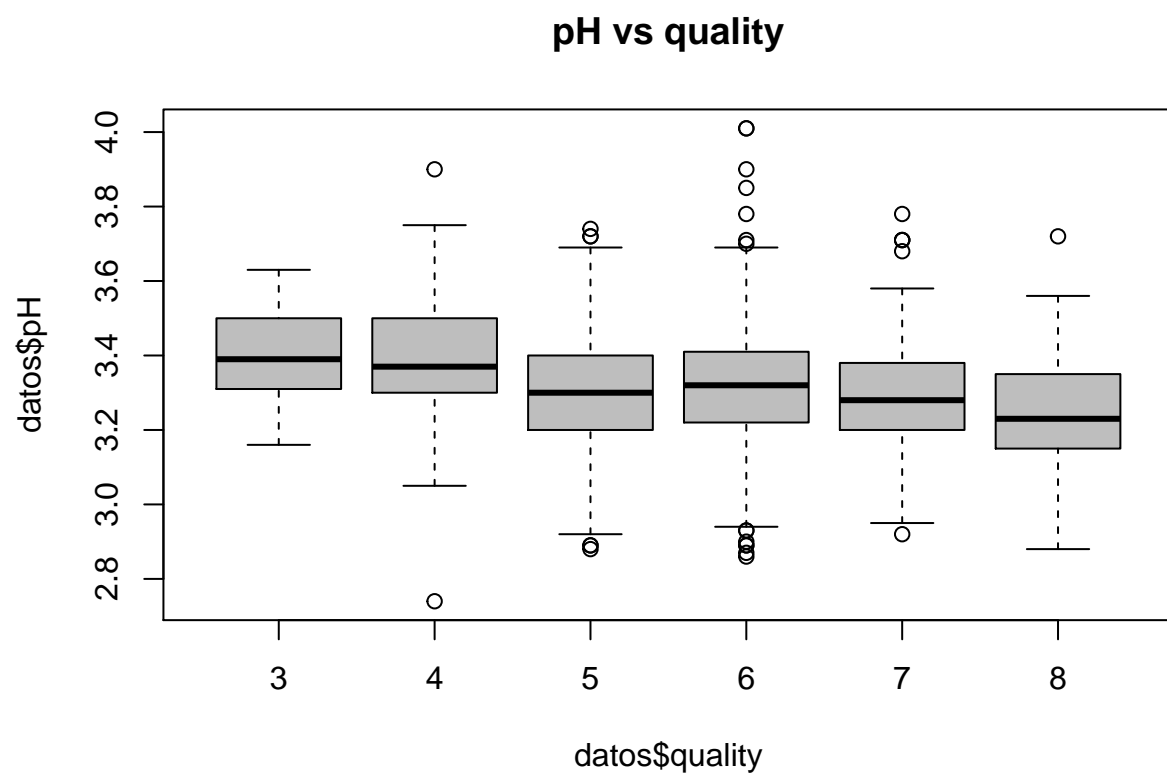


```
boxplot(formula = datos$density ~ datos$quality, main="density vs quality", col="gray")
```

density vs quality

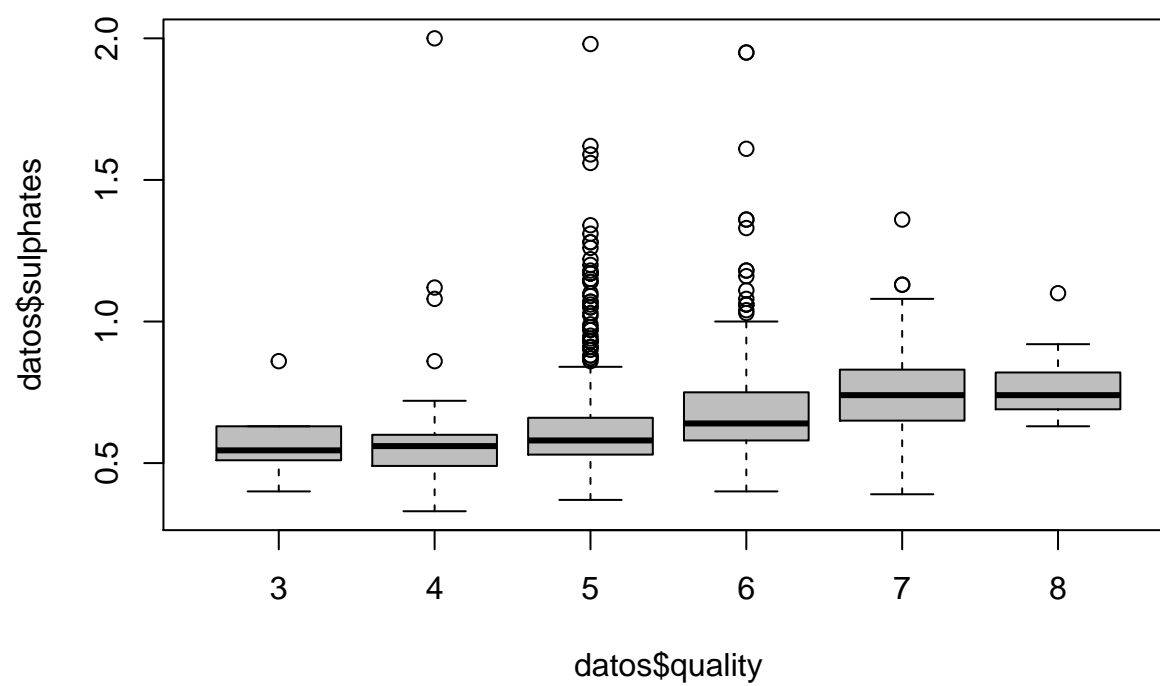


```
boxplot(formula = datos$pH ~ datos$quality, main="pH vs quality", col="gray")
```

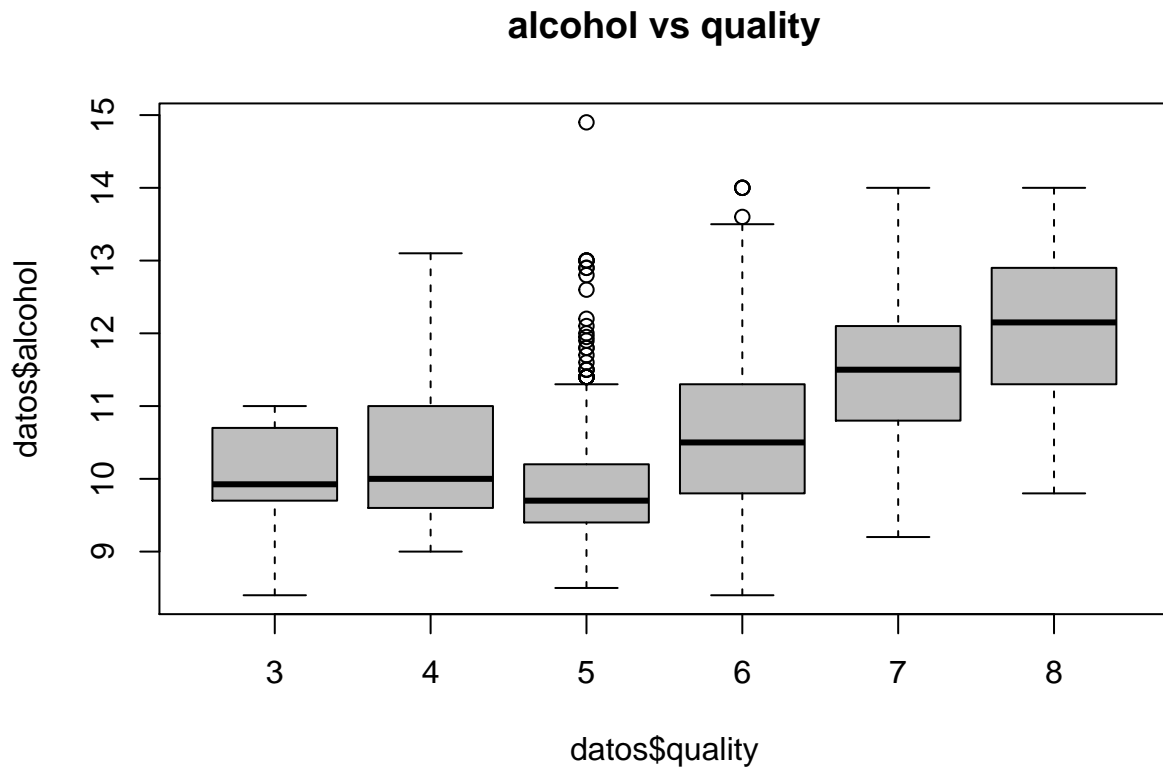



```
boxplot(formula = datos$sulphates ~ datos$quality, main="sulphates vs quality", col="gray")
```

sulphates vs quality



```
boxplot(formula = datos$alcohol ~ datos$quality, main="alcohol vs quality", col="gray")
```



Selección grupo de datos

Comprobación de la normalidad y homogeneidad de la varianza.

```
shapiro.test(datos$fixed.acidity)
```

```
##
##  Shapiro-Wilk normality test
##
## data:  datos$fixed.acidity
## W = 0.94203, p-value < 2.2e-16
```

El test nos indica que ninguna variable está normalizada, ya que el p-valor es inferior al coeficiente 0.05, por lo que se puede rechazar la hipótesis nula y entender que no es normal.

Que no sea normal no quiere decir que no pueda ser normalizable, ya que según el teorema del límite central al tener más de 30 elementos en las observaciones podemos aproximarla como una distribución normal de media 0 y desviación estándar 1.

Aplicación de pruebas estadísticas para comparar los grupos de datos

Representación de los resultados

Resolución del problema