Cheatsheet.R

```
#####
# R #
#####
# install.packages("<package name>")
# library(<package name>)
# print(<something>)
# Assignment statement: x <- <something>
# Manipulating objects in the workspace:
                            # list all objects in memory
# Ls()
# rm(<o1>, <o2>, <o3>, ...) # remove one or more objects from memory by their names
# rm(list = ls())  # remove all objects from memory (usually not recommended)
# Operators:
# + Add, 2 + 3 = 5
# - Subtract, 5 - 2 = 3
# * Multiply, 2 * 3 = 6
\# / Divide, 6 / 2 = 3
# ^ Exponent, 2 ^ 3 = 8
# %% Modulus operator, 9%%2 = 1
# %/% Integer division, 9 %/% 2 = 4
# < Less than
# > Greater than
# = Equal to
# <= Less than or equal to
# >= Greater than or equal to
# != Not equal to
#! Not
# |
      OR
# &
     And
# Expressions:
\# \langle x \rangle / \langle y \rangle - \langle z \rangle^2 \dots
```

```
# Absolute value:
# abs(<value>)
# Vectors:
# <y> <- c(<something1>, <something2>, <something3>, ...)
# <y> <- rep(<something>, <times>)
# <y> <- <int1>:<int2>
\# \langle y \rangle \langle -seq(from = \langle value1 \rangle, to = \langle value2 \rangle, by = \langle step \rangle)
# Matrices:
\# < m > < - matrix(c(3, 5, 7, 1, 9, 4), nrow = 3, ncol = 2, byrow = TRUE)
# <m>.nrow <- nrow(<m>) # number of rows
# <m>.ncol <- ncol(<m>) # number of columns
# <m> <- t(<m>) # transpose <m>
\# < m > [3,2]
\# < m > [2, ]
# Lists: ordered collections of elements of different types
# * <- list(<e1.name> = <e1>, <e2.name> = <e2>, <e3.name> = <e3>, ...)
# <list>[[<index>]] # accessing list element by index, showing value only (returns a vector)
# <list>[<index>] # accessing list element by index, showing both name and value (returns a list)
# <list>$<element.name> # accessing list element by its name
# is.list(<something>)
                                                          # Is <something> a list?
# <combined.list> <- c(<list1>, <list2>, <list3>, ...) # list concatenation
# names(<list>)
                                           # names of list elements
# <list>[names(<list>) == <element.name>] # all elements of a list having the same name
# unlist(<list>)
                                          # convert list into a named vector
# unlist(<list>, use.names = FALSE) # convert list into a character vector
# append(<list>,
                                          # insert new element into an existing list, after index <n>
         list(<e1.name> = <e>),
                                       # new element must be a list itself, that's why list(<e1.name> = <e>)
                                          # <n> is optional; if omitted, new element is appended at the end
              <n>)
# <list>[[<n>]] <- NULL
                                           # remove <n>th element from <list>
# class(<something>)
                                         # data type
# mode(something), typeof(<something>) # how a data item is internally stored in memory
# Factors:
# b \leftarrow c(1, 2, 2, 2, 3, 1, 1, 4, 5, 4)
# b.as.factor <- as.factor(b)</pre>
# levels(b.as.factor)
# f < -factor(c(1, 2, 3))
```

```
# Dataframes:
# e.g., <dataframe> <- as.data.frame(<matrix>)
# str(<dataframe>)
# Reading a dataset:
# <dataframe> <- read.csv("<filename>", stringsAsFactors = FALSE)
# Saving a dataset (modified or newly created dataset):
# write.csv(x = \langle dataframe \rangle, file = "\langle filename \rangle", row.names = F) # do not include the row names (row numbers) column
# saveRDS(object = <dataframe or another R object>, file = "<filename>") # save R object for the next session
# <dataframe or another R object> <- readRDS(file = "<filename>") # restore R object in the next session
# Examining a dataframe:
# str(<dataframe>)
                               # structure of <dataframe>, all variables/columns
                               # showing dimensions (numbers of rows and columns) of a dataframe
# dim(<dataframe>)
# names(<dataframe>)
                               # showing column names
# head(<dataframe>)
                               # the first few rows
                         # the last few rows
# tail(<dataframe>)
# <dataframe>[ , ] # the entire dataframe
# <dataframe>
                            # the entire dataframe
# <dataframe>[<m>, ]
                     # m-th row
# <dataframe>[ ,<n>]
                             # n-th column
# summary(<dataframe>$<column>) # summarizing a variable/column values
# fix(<dataframe>)
                     # editing a dataframe
# new.df <- edit(<dataframe>) # editing a dataframe and assigning the modified dataframe to another datavrame
# Adding/Removing columns to/from a dataframe:
# <dataframe>$<new column name> <- <default value> # adding a new column (default values)
# <dataframe>$<column name> <- NULL
                                                # removing a column
# Adding a new row to a dataframe - the row must be a 1-line dataframe with the same column names:
# <new row> <- data.frame(<column name 1> = <value 1>, <column name 2> = <value 2>,...)
# <new data frame> <- rbind(<dataframe>, <new row>) # append new row to the end of the existing dataframe
# <new data frame> <- rbind(<dataframe>[1:i, ], # insert new row in the middle
#
                           <new row>.
#
                           <dataframe>[(i + 1):nrow(<dataframe>), ])
# Removing rows from a dataframe
# <dataframe>[-i, ]
                                                  # show dataframe without i-th row
# <dataframe>[-c(i, j, k), ]
                                                  # show dataframe without rows i, j, k
# <dataframe> <- <dataframe>[-i, ]
                                                  # remove i-th row from dataframe
```

```
# <dataframe> <- <dataframe>[-c(i, j, k), ]
                                                     # remove rows i, j, k from dataframe
# <dataframe> <- <dataframe>[-(i:k), ]
                                                      # remove rows i to k from dataframe
# Changing column names:
# colnames(<dataframe>)[i] <- "<new name>"
# Changing row names:
# rownames(<dataframe>)[i] <- "<new name>"
# rownames(<dataframe>) <- c("<new name 1>", "<new name 2>",...)
# rownames(<dataframe>) <- c(1, 2,...)</pre>
# rownames(<dataframe>) <- list("<new name 1>", <numeric 2>,...)
# Slicing and dicing dataframes:
# <selection> <- <dataframe>[<some rows>, <some columns>]
# <selection> <- <dataframe>[i:k, c("<column 1>", "<column 2>",...)]
# <selection> <- <dataframe>[<indexes>, ]
# <selection> <- subset(<dataframe>,
                                                                         # subset() is much like SELECT... FROM... WHERE
                         <logical condition for the rows to return>,
                         <select statement for the columns to return>) # can be omitted; column names not prefixed by
<dataframe>$
# <new dataframe> <- <dataframe>[, c("<col1 name>", "<col2 name>")]
# <new dataframe> <- <dataframe>[, <col1 index>:<col2 index>)]
# Shuffling rows/columns:
# <dataframe> <- <dataframe>[sample(nrow(<dataframe>)), ] # shuffle row-wise
                                                              # shuffle column-wise
# <dataframe> <- <dataframe>[, sample(ncol(<dataframe>))]
# Replacing selected values in a column:
# <selected var name> <- <dataframe>$<column> == <selected value>
# <dataframe>$<column>[<selected var name>] <- <new value>
# Applying functions to all elements in rows/columns of a dataframe:
# apply(\langle dataframe \rangle, \langle 1 \mid 2 \rangle, \langle function(x) \mid \{...\} \rangle) # 1 | 2: apply function(x) by row | column
# sapply(\langle vector \rangle, FUN = function(x) {...}) # function(x): function to be applied to each element of \langle vector \rangle
# Partitioning a dataframe:
# install.packages('caret')
# librarv(caret)
# set.seed(<any specific int>) # allows for repeating the randomization process exactly
# <indexes> <- createDataPartition(<dataframe>$<column>, p = 0.8, list = FALSE)
# <partition 1> <- <dataframe>[<indexes>, ]
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# <partition 2> <- <dataframe>[-<indexes>, ]
# for, if, break, next:
# for (<i> in <int vector>) {
   if (<logical condition>) {
    <i 1>
     ine i2>
#
     . . .
#
     break # break: exit the loop; next: skip the remaining lines in this iteration
#
   e n>
# while, if-else, break, next:
# <i> <- <initial value>
# while (logical condition involving <i>) {
   if (<logical condition>) {
    ine i1>
#
     ine i2>
#
     break
               # break: exit the loop; next: skip the remaining lines in this iteration
  } else {
     ine j1>
#
     ine j2>
#
#
   e n>
   <i> <- <modify <i>>>
# ifelse(<condition>, v1, v2) # can return a vector
```

```
# Data type conversion
# b \leftarrow c(1, 2, 2, 2, 3, 1, 1, 4, 5, 4)
# b.as.factor <- as.factor(b)</pre>
# Levels(b.as.factor)
# e.q., <dataframe> <- as.data.frame(<matrix>)
# str(<dataframe>)
# ...
# Convert numeric to factor:
# <dataframe>$<numeric column with few different values> <-
      factor(<dataframe>$<numeric column with few different values>,
              levels = c(0, 1, ..., k), labels = c("\langle l1 \rangle", "\langle l2 \rangle", ..., "\langle lk \rangle"))
#
# Attributes of R objects (dataframes, matrices, factors, lists, tables...)
# attributes(<dataframe> | <matrix> | <factor> | table | ...)
# Tables
# The table() function:
# table(<var>) # typically a factor or an integer var
# The prop.table() function:
# prop.table(table(<var>))
# round(prop.table(table(<var>>)), digits = <n>)
# Row and column margins:
# table(<var1>, <var2>)
                                                          # <var1>, <var2>: usually factors or integers
# table(<rows title> = <var1>, <columns title> = <var2>) # add common titles for rows/columns
# prop.table(table(<var1>, <var2>), margin = 1)
                                                          # all row margins (sums of values by row) are 1.0
# prop.table(table(<var1>, <var2>), margin = 2)
                                                          # all column margins (sums of values by column) are 1.0
# Vectors
# Differences in initializing vectors and dataframe columns:
# <vector> <- rep(<value>, <times>)
# <vector> <- <value>
# <dataframe>$<column> <- rep(<value>, <times>)
# <dataframe>$<column> <- <value>
# Length of a vector:
# Length(<vector>)
# Counting the number of elements with the values of <x> in a vector:
# 1.  <- table(<vector>)
      ["<x>"], or [names() == <x>]
# 2. sum(<vector> == <x>)
# 3. Length(which(<vector> == <x>)) # which() is like WHERE in SQL
```

```
# Appending an element to a vector:
# <vector> <- c(<vector>, <element>)
                                              # type conversion occurs if <element> is of different type than v[i]
                                             # type conversion occurs if <element> is of different type than v[i]
# <vector> <- append(<vector>, <element>)
# <vector> <- append(<vector>, <element>,
                             after = \langle n \rangle
                                              # insert <=> append at a desired location
# <vector> <- append(<vector>, NA)
# Removing NAs from a vector in NA-sensitive functions:
# <function>(<vector>, na.rm = TRUE)
# Selecting items matching criteria from a numeric vector (added check for NAs and NaNs):
# \langle numeric \ vector \rangle \langle -c(\langle n1 \rangle, \langle n2 \rangle, \langle n3 \rangle, \ldots, NA, \ldots NaN)
# <selected> <- <numeric vector>[<logical criterion> & !is.na(<numeric vector>)] # is.na() is TRUE for both NA and NaN
# is.na() is the only way to test if <something> is NA (<something> == NA does not work)
# Range of a numeric vector:
# range(<vector>)
# Create numeric vector with <length> elements:
# <vector> <- vector(mode = "numeric", length = <length>)
# Number of unique values in a vector:
# unique(<vector>)
# Check if numeric variables follow normal distribution:
# summary(<numeric variable>)
                                   # the mean and the median values similar: probably normal distribution
# plot(density((<numeric variable>)) # visual inspection
# hist(<numeric variable>, breaks = <n>) # visual inspection; <n>: number of bins in the histogram
# ggnorm(<numeric variable>)
                                        # values lie more or less along the diagonal (straight line)
# shapiro.test(<numeric variable>) # good for small sample sizes, e.g. n < ~2000; H0: normal distribution
# Discretizing numeric variables (using bnlearn::discretize()):
# library(bnlearn)
# ?discretize()
# <new dataframe with discretized variables> <-
# discretize(<numeric dataframe>,
                                                    # <original dataframe>[, c(<num. col. 1>, <num. col. 1>, ...]
               method = "quantile" | # use equal-frequency intervals (default)
#
#
               method = "interval",
                                                 # use equal-length intervals
               breaks = c(\langle n1 \rangle, \langle n2 \rangle, \ldots, \langle ncol \rangle)) # no. of discrete intervals for each column
# Discretizing numeric variables (using base::cut())
# <dataset>$<new factor feature> <-
    cut(<dataset>$<numeric feature>,
        breaks = \langle n \rangle,
#
                                         # number of intervals to cut the <numeric feature> into
        labels = c("\langle lab 1 \rangle", "\langle lab 2 \rangle", \ldots, "\langle lab n \rangle")) # factor labels
```

```
# Scatterplot matrices (useful for examining the presence of linear relationship between several pairs of variables):
# pairs(~<x1> + <x2> + ..., data = <dataframe>)
# Data normalization:
# librarv(clusterSim)
# <dataframe with numeric columns> <-</pre>
                                                              # works with vectors and matrices as well
# data.Normalization(<dataframe with numeric columns>,
                        type = "n4",
                                                              # normalization: (x - min(x)) / (max(x) - min(x))
#
                        normalization = "column")
                                                              # normalization by columns
                                                 # correlations between numeric variables in the dataset
# Correlation plots:
# <numeric dataframe> <-</pre>
                                                                  # create all-numeric dataframe,
    data.frame(<num col 1 name> = <dataframe>$<num col 1>,  # leave out all non-numeric columns
               <num col 2 name> = <dataframe>$<num col 2>,  # from the original dataframe
# <correlation matrix> <- cor(<numeric dataframe>) # all-numeric dataframe
# library(corrplot)
# corrplot.mixed(<correlation matrix>, tl.cex = <text font size>, number.cex = <number font size>)
# Quantiles/Percentiles:
# <quantiles> <- quantile(<dataset>$<column name>, # examine the 0th, 2.5th, ..., percentile
                           probs = seq(from = 0.0, to = 0.1, by = 0.025))
# Sorting:
# sort(<numeric vector>) # sort <numeric vector>
# install.packages("knitr")
                                                              # pretty-printing tables etc. in the console
                                                              # (a set of "fancy" reporting tools)
# library(knitr)
# kable(x = <stats>, format = "rst")
# ggplot2
# Bar graphs:
# ggplot(data = <dataframe>,
         aes(x = \langle column \ 1 \rangle, y = \langle column \ 2 \rangle, fill = \langle column \ 1 \rangle)) + # fill = \langle column \ 1 \rangle is optional; no y for counts
    geom bar(stat = "identity") +
                                                                      # "identity" for values, "count" for counts
# xlab("<x-axis label>") + ylab("<y-axis label>") +
    agtitle("<graph title>")
\# qqplot(\langle dataframe \rangle, aes(x = \langle column 1 \rangle, fill = \langle column 2 \rangle)) +
    geom bar(position = "dodge", width = <bin width>) +
                                                                     # "dodge": bar grouping, <bin width>: 0.2-0.6
```

```
labs(x = "\langle x-label \rangle", y = "\langle y-label \rangle", title = "\langle title \rangle") +
   theme bw()
# Line graphs:
# ggplot(data = <dataframe>,
          aes(x = \langle column \ 1 \rangle, y = \langle column \ 2 \rangle, qroup = 1)) + # qroup = 1: one line, all points connected
    geom line(colour = "<colour>", linetype = "<linetype>", size = <line thickness>) +
# geom point(colour="<colour>", size = <point size>, shape = <point shape>, fill = "<point fill colour>") +
# xlab("<x-axis label>") + ylab("<y-axis label>") +
# ggtitle("<graph title>")
# All parameters in geom_line() and in geom_point() are optional.
# The defaults are: colour = "black", linetype = "solid", size = 1, shape = 21 (circle), fill = "black"
# See http://www.cookbook-r.com/Graphs/Colors_(ggplot2)/
# for more information on colors.
# See http://www.cookbook-r.com/Graphs/Shapes and line types/
# for information on shapes and line types.
# Scatterplots:
# qqplot(\langle dataset \rangle, aes(x = \langle num.var.1 \rangle, y = \langle num.var.2 \rangle)) +
    geom point(shape = \langle n \rangle, # \langle n \rangle = 1: hollow circle
                fill = <color 1>, # color of point fill (optional)
#
                color = <color 2>, # color of point line (optional)
#
#
                size = <s>) + # size of point line (optional)
    geom_smooth(method = lm,  # add regression line (optional); if left out, nonlinear best-fit line is shown
                 se=FALSE) # do NOT show 95% confidence region as a shaded area (optional)
# <scatterplot> <-
    ggplot(\langle dataset \rangle, aes(x = \langle num.var.1 \rangle, y = \langle num.var.2 \rangle)) +
#
      geom\_point(shape = \langle n \rangle, \qquad \# \langle n \rangle = 1: hollow circle, no fill; \langle n \rangle = 21: circle that can be filled
                  fill = <color 1>, # color of point fill (optional)
#
                   color = <color 2>, # color of point line (optional)
#
                                       # size of point line (optional)
                   size = \langle s \rangle)
# <scatterplot> <- <scatterplot> + xlab("<x label>")
                                                                                # label/caption on x-axis
# <scatterplot> <- <scatterplot> + ylab("<y label>")
                                                                               # label/caption on x-axis
# <scatterplot> <- <scatterplot> + ggtitle("<scatterplot title>")
                                                                                # scatterplot title
# Boxplots:
# boxplot(<dataset>$<column name>, xlab = "<column name>")
                                                                    # basic boxplot for <column name>
                                                                    # returns the stats used for drawing a boxplot
# boxplot.stats(<dataset>$<column name>)
# ggplot(<dataset>,
                                                                     # ggplot2 boxplots
          aes(x = "", y = \langle column \ name \rangle, fill = "\langle color \rangle")) + # show boxplot of \langle column \ name \rangle
#
                                                                    # boxplot width
          geom\ boxplot(width = 0.5) +
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```
stat boxplot(qeom = 'errorbar', width = 0.15) + # show whiskers, control their width
#
       quides(fill = FALSE) +
                                                     # no legend (it makes no sense here)
       xLab("")
                                                     # no x-axis label (it makes no sense here)
# Histograms:
# gqplot(data = <dataset>, mapping = aes(x = <column name>)) +
   geom histogram(bins = <nbins>,
                fill = "<fill color>",
                color = "<line color>")
#
# Density graphs:
# ggplot(data = <dataset>,
       mapping = aes(x = \langle num. \ var. \rangle, \ fill = \langle fill \ var. \rangle)) +
 geom density(alpha = <value>) +
                                                     # alpha: plot transparency (0-1, optional)
# theme_bw()
######
# ML #
######
# Model building and examination:
# <x> and <y> are numeric variables from <dataset>
           data = <dataset>)
# <model>
                       # show the model
# summary(<model>)
                 # show the model statistics
# library(rpart)
# <model> <- rpart(<output variable> ~
                                                                # build the tree
                data = <train dataset>,
                method = "class",
                                                                # build classification tree
                control = rpart.control(minsplit = \langle n \rangle, cp = \langle q \rangle)) # decrease both for larger tree
# library(rattle)
# library(rpart.plot)
# library(RColorBrewer)
# fancyRpartPlot(<decision tree>)
# <model> <- kmeans(x = <normalized dataframe>,
                 centers = \langle k \rangle,
                                                 \# K = \langle k \rangle
```

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iter.max = \langle i \rangle,
                                                      # max number of iterations allowed, e.g. 20
#
                   nstart = \langle n \rangle)
                                                      # no. of initial configurations, e.g. 1000 (report on the best
one)
# library(e1071)
# library(caret)
# <folds> = trainControl(method = "cv", number = <k>)
# define <k>-fold cross-validation parameters
# <cpGrid> = expand.grid(.cp =
                                                             # specify the range of the cp values to examine
                        seq(from = <start value>, to = <end value>, by = <step>))
# train(<output variable> ~
                                                             # find the optimal value for cp
       data = <train dataset>,
       method = "rpart",
                                                             # use rpart() to build multiple classification trees
       control = rpart.control(minsplit = <min split>),
                                                             # default <min split> is 20
       trControl = <folds>, tuneGrid = <cpGrid>)
                                                             # <folds> and <cpGrid> from above
# <prunned model> <- prune(<model>, cp = <optimal cp value>)
# library(class)
# <model> <- knn(train = <training dataset>, # training data without the output (class) variable
                test = <test dataset>, # test data without the output (class) variable
                cl = <class values for training>, # output (class) variable is specified here
                                                 # <n>: random guess, or obtained from cross-validation
                k = \langle n \rangle
# library(e1071)
# ?naiveBayes
# <model> <- naiveBayes(<output variable> ~ .,
                                                    # include all predictors from the training set
                       data = <training dataset>)
# <model> <- naiveBayes(<output variable> ~
                                                    # include only selected predictors from the training set
                       \langle var 1 \rangle + \langle var 2 \rangle + ...,
                       data = <training dataset>)
# Multicolinearity:
# library(car)
# vif(<model>)
# sqrt(vif(<model>)) # variables with sqrt(vif) > 2 (2.5 - disagreement) are problematic
# Making predictions:
# ctions> <- predict(<model>,
         <test dataframe>,
         interval = "confidence" |
                                       # include the confidence interval for the predictions (optional; used
only in linear regression)
                    "predict")
                                  # include prediction intervals (optional)
# cpredictions> <- predict(object = <decision tree>,
                         newdata = <test dataset>,
                          type = "class") # for classification task
```

```
# constant
#
                        newdata = <test dataset>,
                        type = "raw") # compute probabilities, not classes
                                         # examine some of the predictions
# // ci1>:<ik>//
# cpredictions dataframe> <-</pre>
       data.frame(<observation ID> = <test dataset>$<observation ID column>,
#
                 <another relevant feature> = <test dataset>$<another relevant feature column>,
#
                 <output feature> = <test dataset>$<output variable>,
                 <predictions feature> = <predictions>)
# Diagnostic plots:
                    # set up the plotting panel for 4 graphs
\# par(mfrow = c(2,2))
# plot(<model>) # plot the 4 graphs
# par(mfrow = c(1,1)) # reset the plotting panel
# R-sauared and RMSE:
# Compute R-squared on the test data for a model:
# R-squared = 1 - RSS/TSS, where RSS is the residual sum of squares, and TSS is the total sum of squares
# # ctions RSS> <-</pre>
# sum((<predictions> - <test dataset>$<output variable>)^2)
# # ctions TSS> <-
# sum(mean(<train dataset>$<output variable>) - <test dataset>$<output variable>)^2)
# <R-squared>
# Compute Root Mean Squared Error (RMSE) for a model on the test data
# to see how much error we are making with the predictions:
\# RMSE = sqrt(RSS/n)
# credictions RMSE> <- sqrt(<pre>credictions RSS> / nrow(<test dataset>))
# # ctions RMSE>
# ROC curve (Receiver Operating Characteristic)
# Library(pROC)
# <ROC curve parameters> <-
                                                    # compute ROC curve parameters
   roc(response = <test dataset>$<output variable>,
       predictor = col. no. of the "positive class" (can be the No class!)
# <ROC curve parameters>$auc
                                                    # extract and show AUC
# plot.roc(<ROC curve parameters>,
                                     # computed in the previous step
         print.thres = TRUE,
                                           # show the probability threshold (cut-off point) on the plot
```

```
print.thres.best.method =
#
              "vouden" |
                                # maximize the sum of sensitivity and specificity (the distance to the diag. line)
              "closest.topleft") # minimize the distance to the top-left point of the plot
# <ROC coords> <- coords(<ROC curve parameters>,
                                                                           # computed in the previous step
                        ret = c("accuracy", "spec", "sens", "thr", ...), # ROC curve parameters to return
#
#
                                               # the coordinates to look for:
#
                             "Local maximas" | # Local maximas of the ROC curve
                             "best" | ...) # the point with the best sum of sensitivity and specificity, i.e.
#
#
                                               # the same as the one shown on the ROC curve
# Compare multiple clustering results/schemes:
# install.packages("fpc")
# library(fpc)
# ?cluster.stats
# <comparison criteria> <-
                                      # specify criteria (from cluster.stats()) for comparing
 c("<criterion 1>",
                                      # different clusterings (e.g., "max.diameter", "min.separation",
                                      # "average.between", "average.within", "within.cluster.ss", ...)
      "<criterion 2>", ...)
# <distance matrix> <-
# dist(x = <normalized dataset>)
# <comparison> <- sapply(list(<clustering 1 name> = <clustering 1>, # <clustering 1> computed by kmeans()
                             <clustering 1 name> = <clustering 2>, # <clustering 2> computed by kmeans()
#
                         FUN = function(x)
                         cluster.stats(<distance matrix>, x))[<comparison criteria>, ]
# install.packages("knitr")
                                                           # pretty-printing tables etc. in the console
                                                           # (a set of "fancy" reporting tools)
# library(knitr)
# kable(x = comparison, format = "rst")
```