Prediction

Autor: "Ali Abdorrahimi"

Studiengang: "Digital Engineering"

Fach: "Data Mining"

Semester : "Wintersemester 2022"

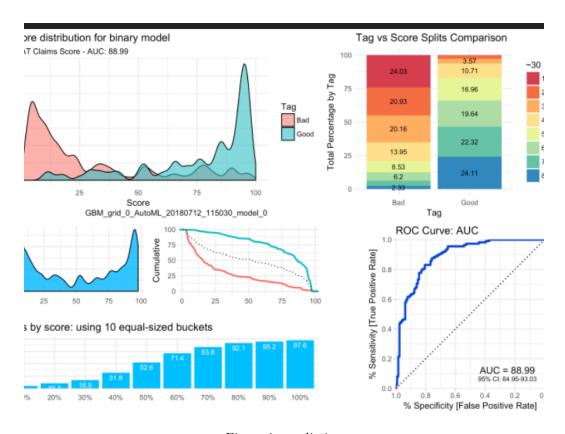


Figure 1: prediction

Introduction

In der R-Programmierung sind Prognosemodelle äußerst nützlich für die Vorhersage zukünftiger Ergebnisse

Mit Hilfe von Modellen können Sie zukünftiges Verhalten auf der Grundlage von vergangenem Verhalten vorhersagen. Nachdem Sie ein Modell aufgebaut haben, verwenden Sie es, um neue Daten zu bewerten, d. h. um Vorhersagen zu treffen.

Mit R können Sie viele Arten von Modellen erstellen.

these R models:

glm : Generalized linear models lda : Linear Discriminant Analysis knn : K-nearest Neighbors Algorithm

rpart: Recursive Partitioning and Regression Trees

kmeans: k-Means clustering

rf: Random Forest

Load Libraries

```
library(magrittr)
library(plyr)
library(dplyr)
library(ggplot2)
library(grid)
library(gridExtra)
library(stringr)
library(caret)
library(mvtnorm)
library(ggplot2)
library(randomForest)
library(caret)
library(pROC)
library(klaR)
library(psych)
library(MASS)
library(devtools)
library(ROCR)
```

Load Data

```
workPath ="D:\\DataMining\\"
cardio <- readRDS(paste (workPath , "cardio.rds", sep =""))</pre>
```

Train test split

Wenn wir Modelle für maschinelles Lernen erstellen, müssen wir unser Modell auf einer Teilmenge der verfügbaren Daten trainieren und die Genauigkeit des Modells auf einer Teilmenge der Daten testen.

verwenden wir die Funktion complete.cases() in R, um (Missing values) in einem Vektor, einer Matrix oder einem Datenrahmen zu entfernen.

```
cardio_complete <- cardio[(complete.cases(cardio)),]
nrow(cardio_complete)</pre>
```

[1] 2927

Nach der Entfernung der Missing Valuses haben wir 2927 Datensätze, von denen wir 1500 Datensätze für den Test verwenden werden.

```
set.seed(1001)

n_test <- 1500

idx_test <- sample(1:nrow(cardio_complete), n_test)

cardio_test <- cardio_complete[ idx_test,]
cardio_train <- cardio_complete[ -idx_test,]</pre>
```

Train first models

mod.cart\$results

Accuarcy ist ein Maßstab für die Bewertung von Klassifizierungsmodellen und Formal ist die accuracy wie folgt definiert:

$$Accuracy = \frac{Number of Correct Prediction}{Total Number of Predicted}$$

```
control <- trainControl(method="cv", number=5)
metric <- "Accuracy"</pre>
```

The train function can be used to:

- evaluate, using resampling, the effect of model tuning parameters on performance
- choose the "optimal" model across these parameters
- estimate model performance from a training set

```
mod.glm <- train(target~., data=cardio_train, method = "glm",family = "binomial",metric=metric,trContro</pre>
```

The Kappa statistic (or value) is a metric that compares an Observed Accuracy with an Expected Accuracy (random chance). The kappa statistic is used not only to evaluate a single classifier, but also to evaluate classifiers amongst themselves. In addition, it takes into account random chance (agreement with a random classifier), which generally means it is less misleading than simply using accuracy as a metric (an Observed Accuracy of 80% is a lot less impressive with an Expected Accuracy of 75% versus an Expected Accuracy of 50%). Computation of Observed Accuracy and Expected Accuracy is integral to comprehension of the kappa statistic, and is most easily illustrated through use of a confusion matrix

```
mod.lda <- train(target~., data=cardio_train, method="lda",metric=metric,trControl=control)
mod.lda$results

## parameter Accuracy Kappa AccuracySD KappaSD
## 1 none 0.8507324 0.128926 0.01813854 0.107302

mod.cart <- train(target~., data=cardio_train, method="rpart",metric=metric,trControl=control)</pre>
```

```
cp Accuracy Kappa AccuracySD
## 1 0.007898894 0.8437370 0.15153304 0.007838914 0.08565146
## 2 0.009478673 0.8437370 0.11206497 0.008588368 0.06216736
## 3 0.026066351 0.8528401 0.01988439 0.002304735 0.04446285
mod.knn <- train(target~., data=cardio_train, method="knn", metric=metric, trControl=control)</pre>
mod.knn$results
                     Kappa AccuracySD
## k Accuracy
## 1 5 0.8374310 0.09635543 0.008192540 0.03566687
## 2 7 0.8437419 0.07216125 0.010423923 0.05015822
## 3 9 0.8437370 0.02081593 0.007015914 0.04194563
mod.svm.radial <- train(target~., data=cardio_train, method="svmRadial", metric=metric, trControl=contr</pre>
mod.svm.radial$results
                                  Kappa AccuracySD
         sigma C Accuracy
## 1 0.0461617 0.25 0.8521401 0.00000000 0.001098966 0.0000000
## 2 0.0461617 0.50 0.8521401 0.00000000 0.001098966 0.0000000
## 3 0.0461617 1.00 0.8521352 0.01290007 0.001741712 0.0211263
mod.svm.linear <- train(target~., data=cardio_train, method="svmLinear", metric=metric, trControl=contr
mod.svm.linear$results
## C Accuracy
                       Kappa AccuracySD
                                             KappaSD
## 1 1 0.8507349 -0.002760661 0.001986759 0.00378019
mod.rf <- train(target~., data=cardio_train, method="rf", metric=metric, trControl=control)</pre>
mod.rf$results
    mtry Accuracy
                        Kappa AccuracySD
                                             KappaSD
## 1 2 0.8514366 0.01155700 0.001892435 0.01602853
## 2 9 0.8479328 0.07532032 0.003120357 0.04284008
     17 0.8444265 0.10546684 0.016218403 0.08464020
## 3
```

summarize all models

##

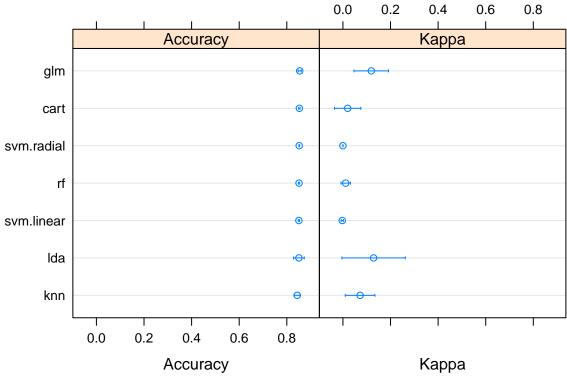
```
results <- resamples(list(glm = mod.glm , lda= mod.lda, cart=mod.cart, knn=mod.knn, svm.radial=mod.svm.summary(results)

##
## Call:
## summary.resamples(object = results)
##
## Models: glm, lda, cart, knn, svm.radial, svm.linear, rf
## Number of resamples: 5</pre>
```

```
## Accuracy
##
                    1st Qu.
                             Median
                                       Mean
                                            3rd Qu.
                                                       Max. NA's
               Min.
           0.8491228 0.8526316 0.8526316 0.8542405 0.8536585 0.8631579
## glm
           0.8286713 0.8421053 0.8491228 0.8507324 0.8561404 0.8776224
## lda
## cart
           0.8496503 0.8526316 0.8526316 0.8528401 0.8531469 0.8561404
           0.8321678 0.8356643 0.8421053 0.8437419 0.8526316 0.8561404
## knn
## svm.radial 0.8501742 0.8526316 0.8526316 0.8521401 0.8526316 0.8526316
## svm.linear 0.8491228 0.8491228 0.8496503 0.8507349 0.8526316 0.8531469
## rf
           0.8491228 0.8496503 0.8526316 0.8514366 0.8526316 0.8531469
##
## Kappa
##
                 Min.
                         1st Qu.
                                   Median
                                                     3rd Qu.
                                               Mean
           ## glm
           ## lda
## cart
           0.00000000 0.00000000 0.00000000 0.019884393 0.00000000
           ## knn
## svm.linear -0.006901651 -0.006901651 0.00000000 -0.002760661 0.00000000
           0.00000000 0.00000000 0.00000000 0.011557004 0.02529229
##
                Max. NA's
## glm
           0.21502996
## lda
           0.29675425
## cart
           0.09942197
## knn
           0.14414414
## svm.radial 0.00000000
## svm.linear 0.00000000
           0.03249273
saveRDS(results,file =paste (workPath , "results.rds", sep =""))
```

plot accuracies

```
dotplot(results)
```



Confidence Level: 0.95

Analyse der ersten Modelle

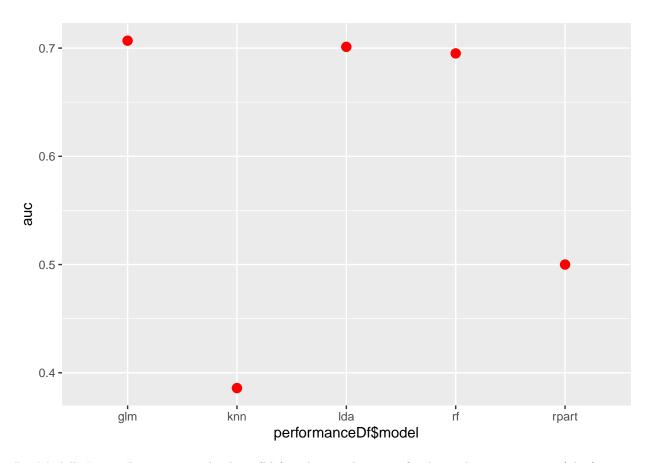
AUC (Area Under The Curve) Die ROC-Kurve (Receiver Operating Characteristics) ist eines der wichtigsten Bewertungsmaße für die Beurteilung der Leistung von binären Klassifikationsproblemen. ROAC ist eine Wahrscheinlichkeitskurve, die die TPR (True Positive Rate) gegen die FPR (False Positive Rate) aufträgt. AUC ist das Maß für die Trennbarkeit und zeigt an, wie gut unser Modell in der Lage ist, zwischen den Klassen zu unterscheiden. Die AUC gibt an, wie gut das Modell zwischen positiven und negativen Klassen unterscheidet. Je größer der AUC, desto besser.

```
computeData <- F
models <- c("rf", "lda", "rpart", "knn", "glm")
workPath = "D:\\DataMining\\Seafile\\01_cardio\\Projekt\\"

if (computeData) {
    performanceDf <- data.frame(model = models)
    performanceDf$accuracy <- NA
    performanceDf$sensitivity <- NA
    performanceDf$specificity <- NA
    performanceDf$specificity <- NA
    for (ir in 1:nrow (performanceDf)) {
        mod <-
            train(
                target ~ .,
                data = cardio_train,
                method = performanceDf$model[ir]</pre>
```

```
metric = metric,
       trControl = control
     )
   y_pred_prob <- predict(mod , cardio_test, type = "prob")</pre>
   y_pred <- predict(mod , cardio_test)</pre>
   table(y_pred, cardio_test$target)
   accuracy = table(y_pred == cardio_test$target)["TRUE"] / length (y_pred)
   performanceDf$accuracy[ir] = accuracy
   sensitivity = sensitivity(factor(y_pred), factor(cardio_test$target))
   performanceDf$sensitivity[ir] = sensitivity
   specificity = specificity(factor(y_pred), factor(cardio_test$target))
   performanceDf$specificity[ir] = specificity
   auc <- roc(cardio_test$target , y_pred_prob$CHD)$auc</pre>
   performanceDf$auc[ir] = auc
 saveRDS(performanceDf, file = paste (workPath , "performanceDf.rds", sep =
                                        ""))
performanceDf <-</pre>
 readRDS(paste (workPath , "performanceDf.rds", sep = ""))
performanceDf
   model accuracy sensitivity specificity
##
## 1 rf 0.8453333 0.01287554 0.9984215 0.6951062
## 3 rpart 0.8446667 0.00000000 1.0000000 0.5000000
## 4 knn 0.8420000 0.04291845 0.9889503 0.3858732
      glm 0.8453333 0.09871245 0.9826361 0.7068537
## 5
ggplot(performanceDf, aes(x = performanceDf$model , y = auc )) +
  geom_point(colour = "red", size = 3)
```

Warning: Use of 'performanceDf\$model' is discouraged. Use 'model' instead.



Die Modelle Linear Discriminant Analysis (lda) und generalization of ordinary linear regression (glm)zeigten das beste Ergebnis mit einem auc von 0.7011527 und 0.7068537.

as Modell k nearest neighbor hat die schlechteste Performance mit einem auc von 0,3858732.