# STA 545 Statistical Data Mining I, Fall 2020 Homework 10

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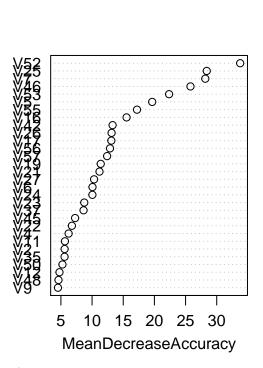
December 9, 2020

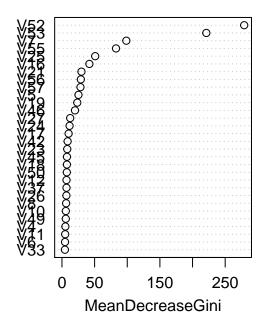
# Problem 1

```
a)
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
spam <- read.table("spam.data.txt")</pre>
#spam$V58 <- as.factor(spam$V58)</pre>
set.seed(2)
train.num <- sample(nrow(spam), 2301)</pre>
spam.train <- spam[train.num , ]</pre>
spam.test <- spam[-train.num ,]</pre>
frac.spam.train <- nrow(subset(spam.train, V58 == 1))/nrow(spam.train)</pre>
frac.spam.test <- nrow(subset(spam.test, V58 == 1))/nrow(spam.test)</pre>
# the fraction of the training data is spam
frac.spam.train
## [1] 0.398957
# the fraction of the testng data is spam
frac.spam.test
## [1] 0.3891304
  b) Bagging
library(randomForest)
## randomForest 4.6-14
## Type rfNews() to see new features/changes/bug fixes.
```

```
##
## Attaching package: 'randomForest'
## The following object is masked from 'package:dplyr':
##
## combine
bag.spam = randomForest(as.factor(V58) ~ .,data = spam.train, mtry = 57, importance = TRUE, ntree = 100
pred.bag.spam <- predict(bag.spam, newdata = spam.test)
bag.error.rate <- mean(pred.bag.spam != spam.test$V58)
#the error rate of the ensemble on the testing data
bag.error.rate
## [1] 0.06347826
varImpPlot(bag.spam)</pre>
```

bag.spam





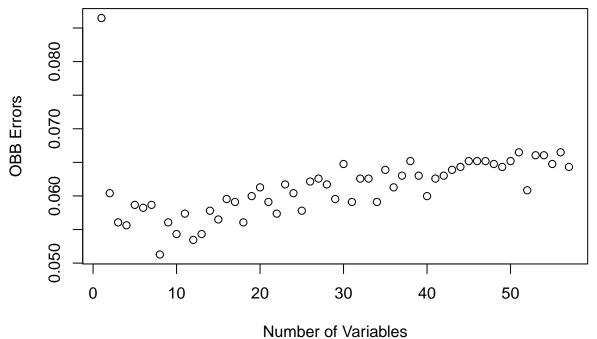
## c)Random Forest

```
m.vec <- c(1:57)

cal.error <- function(vec){
    n = length(vec)
    obb.err.vec = rep(0,n)
    test.err.vec = rep(0,n)

for(i in vec){
    rf <- randomForest(as.factor(V58) ~ .,data = spam.train, mtry = i, ntree = 100)
    pred <- predict(rf, newdata = spam.test)</pre>
```

```
#calculate obb error
obb.err.vec[i] <- mean((rf$predicted != spam.train$V58)^2)
#calculate test error
test.err.vec[i] <- mean((pred != spam.test$V58)^2)
}
return(data.frame(obb = obb.err.vec,test.error = test.err.vec))
}
errors <- cal.error(m.vec)
plot(m.vec, errors$obb, xlab = "Number of Variables", ylab = "OBB Errors")</pre>
```



plot(m.vec,errors\$test.error, xlab = "Number of Variables", ylab = "Test Errors")

```
0
     0.075
Test Errors
     0.065
                                             0
                                  0000
                                                 00
               , 000<sub>0</sub>000,
     0.055
                           0
           0
                      10
                                  20
                                             30
                                                         40
                                                                     50
                                   Number of Variables
```

```
opt.m.test <- m.vec[which.min(errors$test.error)]
opt.m.obb <- m.vec[which.min(errors$obb)]

#optimal parameter m making test error smallest
opt.m.test

## [1] 14

#optimal parameter m making OBB error smallest
opt.m.obb

## [1] 8

# The error rate of the random forest on the testing data
rf.error.rate <- min(errors$test.error)
rf.error.rate

## [1] 0.05565217
d)AdaBoost
library(ada)</pre>
```

## [1] 0.05478261

e)Logistic Regression and Evaluation

```
glm.spam = glm(V58 ~ ., data = spam.train, family = "binomial")

## Warning: glm.fit: fitted probabilities numerically 0 or 1 occurred

pred.glm.spam <- ifelse(predict(glm.spam, newdata = spam.test , "response") > 0.5, 1, 0)
glm.error.rate <- mean(spam.test[,58]!=pred.glm.spam)

# The error rate of the Logistic Regression classifier on the testing data
glm.error.rate</pre>
```

## [1] 0.07608696

To sum up, the error rate on the testing data of Bagging, Random Forest, AdaBoost and Logistic Regression are around 0.06, 0.06, 0.05, 0.08; based on the test error, AdaBoost has the best performance in this case.

## Problem 2

```
# separate data for k-fold cross validation;
cv.group<-function(k,datasize,seed){</pre>
    cvlist<-list()</pre>
    set.seed(seed)
    n<-rep(1:k,ceiling(datasize/k))[1:datasize]</pre>
    temp<-sample(n,datasize)</pre>
    x<-1:k
    dataseq<-1:datasize
  cvlist<-lapply(x,function(x) dataseq[temp==x])</pre>
  return(cvlist)
}
whole.cv.list = cv.group(5,nrow(spam.train),2)
#create formula
names=names(spam.train)
f =as.formula(paste("V58 ~", paste(names[!names %in% "V58"], collapse = " + ")))
#implement neural network model and get the test error one time
library(neuralnet)
## Attaching package: 'neuralnet'
## The following object is masked from 'package:dplyr':
##
##
       compute
cv.test.nn<-function(i,j,cv.list){</pre>
  #i refers to i-th dataset, the testing data
  #j refers to the number of neurons
  data.train <- spam.train[-cv.list[[i]],]</pre>
    data.test<- spam.train[cv.list[[i]],]</pre>
    nn = neuralnet(f, data = data.train, hidden = j, threshold=0.01,
                    algorithm = "rprop+", err.fct="ce",
                    act.fct="logistic",linear.output=FALSE)
```

```
pred.nn = compute(nn,data.test)
    nn.pred<- lapply(pred.nn\net.result, function(x) ifelse(x>0.5,1,0))
  accu <- sum(nn.pred == data.test$V58)/nrow(data.test)</pre>
 return(list(accu,j))
#qet the optimal number of neurons in the single layer for neural network model
i < c(1:5)
j \leftarrow c(10,20,30,40,50) # the number of neurons that we need to test
i.s<-rep(i,times=length(j)) # [1] 1 2 3 4 5 1 2 3 4 5 1 2 3 4 5 1 2 3 4 5 5 5 1 2 3 4 5 5 5 5 5 6 6 7 8 9 9 9 9
i.s<-rep(j,each=5) #[1] 10 10 10 10 10 20 20 20 20 20 30 30 30 30 40 40 40 40 40 50 50 50 50
ijs <- cbind(i.s,j.s)</pre>
i = 0
j = 0
tmp.accus <-NULL
accus <- rep(0,5) # contain accuracy of all numbers of neurons that we need to test
neus \leftarrow \text{rep}(0,5)
index = 1
old neu = 10
for(p in 1:nrow(ijs)){
 if(ijs[p,2] == old_neu){
    i = ijs[p,1]
    j = ijs[p,2]
    accu_neuron = cv.test.nn(i,j,whole.cv.list)
    accu = accu_neuron[[1]]
    neu = accu_neuron[[2]]
    tmp.accus <- c(tmp.accus,accu)</pre>
  accus[index] = mean(tmp.accus)
  neus[index] = j
  index = as.integer(j/10)
  old_neu = as.integer(j+10)
opt.h = neus[which.max(accus)]
#the optimal number of neurons in the single hidden layer
opt.h
## [1] 30
#fit the model to the spam data using opt.h
nn.spam = neuralnet(f, data = spam.train, hidden = opt.h, threshold=0.01,
                     algorithm = "rprop+", err.fct="ce",
                     act.fct="logistic",linear.output=FALSE)
pred.nn.spam = compute(nn.spam,spam.test)
net.prediction.spam<- lapply(pred.nn.spam$net.result, function(x) ifelse(x>0.5,1,0))
nn.error <- mean((net.prediction.spam != spam.test$V58)^2)</pre>
```

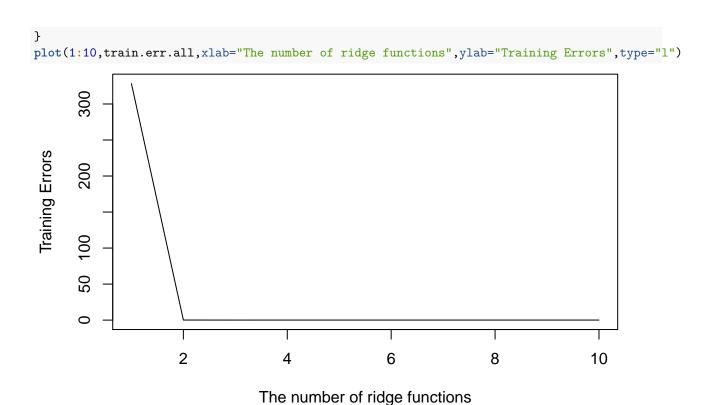
```
nn.error
```

#### ## [1] 0.09391304

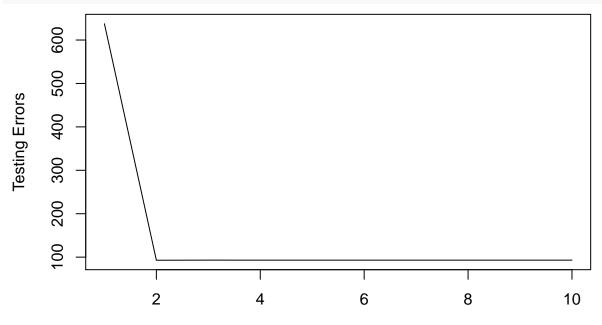
In this case, we separately used 10, 20, 30, 40 and 50 as the number of neurons in our 5-fold cross validation to choose the optimal number of neurons, and based on the accuracy of classification results, we finally chose 10 as the the optimal number of neurons in the single hidden layer.

#### Problem 3

```
a)data preparing
set.seed(50321222)
X1 \leftarrow rnorm(100,0,1)
X2 \leftarrow rnorm(100,0,1)
Z \leftarrow rnorm(100,0,1)
X.train <- as.matrix(rbind(X1,X2))</pre>
a1 <- as.matrix(rbind(3,3))
a2 <- as.matrix(rbind(3,-3))</pre>
v = t(a1) \% X.train
Y.train \leftarrow t(1/(1 + \exp(-v)) + (t(a2)%*%X.train)^2 + 0.3*Z)
X.train <- t(X.train)</pre>
X1 \leftarrow rnorm(1000,0,1)
X2 \leftarrow rnorm(1000,0,1)
Z \leftarrow rnorm(1000,0,1)
X.test <- as.matrix(rbind(X1,X2))</pre>
a1 <- as.matrix(rbind(3,3))
a2 <- as.matrix(rbind(3,-3))
v = t(a1) \% X.test
Y.test <-t(1/(1 + \exp(-v))+(t(a2)%*%X.test)^2+0.3*Z)
X.test <- t(X.test)</pre>
  b)
library(MASS)
##
## Attaching package: 'MASS'
## The following object is masked from 'package:dplyr':
##
##
       select
train.err.all=rep(NA,10)
test.err.all=rep(NA,10)
for(i in 1:10){
  ppr.model=ppr(X.train,Y.train,nterms=i,
                 sm.method="gcvspline")
  predicted.values=predict(ppr.model,newdata=X.test)
  train.err.all[i] = mean((ppr.model$fitted.values-Y.train)^2)
  test.err.all[i]=mean((predicted.values-Y.test)^2)
```



plot(1:10,test.err.all,xlab="The number of ridge functions",ylab="Testing Errors",type="l")



c)
library(MASS)
train.original <- data.frame(X.train, Y.train)
test.original <- data.frame(X.test, Y.test)
maxs=apply(train.original, 2, max)</pre>

The number of ridge functions

```
mins=apply(train.original, 2, min)
train.scaled=as.data.frame(scale(train.original, center = mins, scale = maxs - mins))
test.scaled=as.data.frame(scale(test.original, center = mins, scale = maxs - mins))
train.err.all=rep(NA,10)
test.err.all=rep(NA,10)
for(i in 1:10){
  nn=neuralnet(Y.train~X1+X2,data=train.scaled, hidden=i, threshold=0.01,
              err.fct="sse", act.fct="logistic",linear.output=T)
  nn.scaled.prediction=compute(nn,test.scaled[,1:2])
  nn.scaled.fitted=compute(nn,train.scaled[,1:2])
  nn.prediction=nn.scaled.prediction$net.result*(max(train.original$Y.train)
                                          -min(train.original$Y.train))+min(train.original$Y.train)
  nn.fitted=nn.scaled.fitted$net.result*(max(train.original$Y.train)
                                          -min(train.original$Y.train))+min(train.original$Y.train)
  train.err.all[i] = mean((nn.fitted-train.original[,3])^2)
  test.err.all[i] = mean((nn.prediction-test.original[,3])^2)
plot(1:10,train.err.all,xlab="The number of neurons",ylab="Training Errors",type="l")
      200
      150
Training Errors
      100
      50
      0
                                                    6
                     2
                                    4
                                                                   8
                                                                                  10
```

plot(1:10,test.err.all,xlab="The number of neurons",ylab="Testing Errors",type="l")

The number of neurons

