# Homework 2

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```
get accuracy <- function(predicted, actual){</pre>
  confusion_table = table(predicted, actual)
  TP = confusion_table[2,2]
  TN = confusion_table[1,1]
  FN = confusion_table[1,2]
  FP = confusion_table[2,1]
  accuracy = round((TP + TN) / sum(TP,FP,TN,FN), 2)
  return(accuracy)
get_classification_error_rate <- function(predicted, actual){</pre>
  confusion_table = table(predicted, actual)
  TP = confusion table[2,2]
  TN = confusion_table[1,1]
  FN = confusion_table[1,2]
  FP = confusion_table[2,1]
  classification_error_rate = round((FP + FN) / sum(TP,FP,TN,FN),2)
  return(classification_error_rate)
get_precision <- function(predicted, actual){</pre>
  confusion_table = table(predicted, actual)
  TP = confusion_table[2,2]
  TN = confusion_table[1,1]
  FN = confusion_table[1,2]
  FP = confusion_table[2,1]
  precision = round(TP / (TP + FP), 2)
  return(precision)
}
get_sensitivity <- function(predicted, actual){</pre>
  confusion_table = table(predicted, actual)
  TP = confusion_table[2,2]
  TN = confusion_table[1,1]
  FN = confusion_table[1,2]
  FP = confusion_table[2,1]
  sensitivity = round(TP / (TP + FN), 2)
  return(sensitivity)
get_specificity <- function(predicted, actual){</pre>
```

```
confusion_table = table(predicted, actual)
  TP = confusion table[2,2]
  TN = confusion_table[1,1]
  FN = confusion_table[1,2]
  FP = confusion_table[2,1]
  specificity = round(TN / (TN + FP), 2)
 return(specificity)
get_f1_score <- function(predicted, actual){</pre>
  confusion_table = table(predicted, actual)
  TP = confusion_table[2,2]
  TN = confusion_table[1,1]
  FN = confusion_table[1,2]
  FP = confusion_table[2,1]
  precision = round(TP / (TP + FP), 2)
  sensitivity = round(TP / (TP + FN), 2)
  f1_score = round((2 * precision * sensitivity) / (precision + sensitivity), 2)
  return(f1_score)
get_false_positive_rate <- function(predicted, actual){</pre>
  confusion_table = table(predicted, actual)
  TP = confusion table[2,2]
 TN = confusion_table[1,1]
 FN = confusion_table[1,2]
 FP = confusion_table[2,1]
 fpr = round(FP / (FP + TN), 2)
 return(fpr)
get_false_negative_rate <- function(predicted, actual){</pre>
  confusion_table = table(predicted, actual)
  TP = confusion_table[2,2]
 TN = confusion_table[1,1]
 FN = confusion_table[1,2]
  FP = confusion_table[2,1]
 fnr = round(FN / (FN + TP), 2)
  return(fnr)
}
```

## Section 1

#### Question 1.1

Create a training set of 800 observations, and a test set containing the rest using the following code:

```
# ``{r pressure, echo=FALSE}
n <- nrow(OJ)
set.seed(1234)
id.train=sample(1:n,size=800)</pre>
```

```
id.test=setdiff(1:n,id.train)
OJ$Purchase = as.factor(OJ$Purchase)
OJ$Store7 = as.factor(OJ$Store7)
as.data.frame(sapply(OJ,class))
##
                  sapply(OJ, class)
## Purchase
                              factor
## WeekofPurchase
                             numeric
## StoreID
                             numeric
## PriceCH
                             numeric
## PriceMM
                             numeric
## DiscCH
                             numeric
## DiscMM
                             numeric
## SpecialCH
                             numeric
## SpecialMM
                             numeric
## LoyalCH
                             numeric
## SalePriceMM
                             numeric
## SalePriceCH
                             numeric
## PriceDiff
                             numeric
## Store7
                              factor
## PctDiscMM
                             numeric
## PctDiscCH
                             numeric
## ListPriceDiff
                             numeric
## STORE
                             numeric
OJ.train=OJ[id.train,-3]
OJ.test=OJ[id.test,-3]
```

#### Question 1.2

Construct an unpruned classification tree to predict the variable purchase using the available predictors. Calculate the false positive rate, false negative rate and overall error rate of this tree on the test data (note: you can use your code from the previous assignment directly for this question).

```
# ```{r pressure, echo=FALSE}
tree <- rpart(OJ.train$Purchase ~ . ,data = OJ.train, method = 'class')
#summary(tree)

pred = predict(tree,newdata = OJ.test,type = c("class"))

print(pasteO("False Positive Rate: ", get_false_positive_rate(OJ.test$Purchase,pred)))

## [1] "False Positive Rate: 0.14"

print(pasteO("Error Rate: ", get_classification_error_rate(OJ.test$Purchase,pred)))

## [1] "Error Rate: 0.19"

print(pasteO("False Negative Rate: ", get_false_negative_rate(OJ.test$Purchase,pred)))

## [1] "False Negative Rate: 0.27"</pre>
```

#### Question 1.3

Using all the data (training and test put together), use the bagging approach to do the same analysis again and compare the results of the different error rates.

#### Question 1.4

Using all the data (training and test put together), use the random forest approach to redo the same analysis and compare the results of the different error rates.

#### Question 1.5

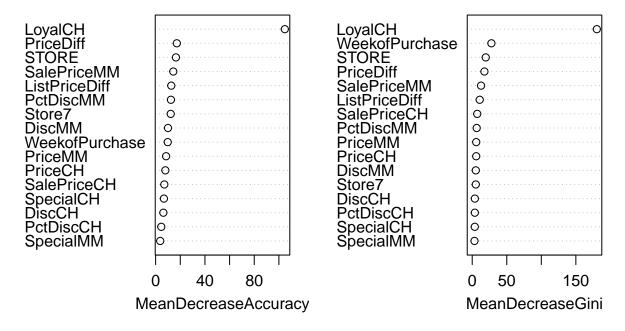
Calculate the importance of the variables in the classification tree constructed in 2) and the forest constructed in 4). Compare.

```
set.seed(1)
as.data.frame(importance(rf.OJ))
```

```
## CH MM MeanDecreaseAccuracy MeanDecreaseGini ## WeekofPurchase 6.2370043 6.263470 9.821831 27.774264
```

##	PriceCH	6.4129690	3.347700	7.894530	5.809040	
##	PriceMM	3.6162856	7.199457	8.496137	5.816289	
##	DiscCH	3.2356695	5.311119	6.259811	3.814248	
##	DiscMM	5.1714256	6.529393	10.040448	5.196968	
##	SpecialCH	2.1385953	5.812328	6.715932	3.727776	
##	SpecialMM	-0.4787197	4.741747	3.719254	3.312229	
##	LoyalCH	73.7608676	98.090066	104.804821	180.327296	
##	${\tt SalePriceMM}$	6.9341087	11.630595	14.330011	12.811005	
##	SalePriceCH	5.3705321	3.347173	7.054213	7.263904	
##	PriceDiff	9.6259636	12.248643	17.187297	17.667201	
##	Store7	4.4216587	12.366033	12.239986	5.174581	
##	PctDiscMM	7.1044541	8.672862	12.428228	6.497433	
##	PctDiscCH	2.3521748	4.010164	4.593497	3.812999	
##	ListPriceDiff	8.0987352	8.405751	12.687143	10.860433	
##	STORE	7.2910690	15.037911	16.467642	19.595198	
<pre>varImpPlot(rf.OJ)</pre>						

# rf.OJ



```
set.seed(1)
as.data.frame(tree$variable.importance)
```

##		tree\$variable.importance
##	LoyalCH	170.416755
##	PriceDiff	21.478485
##	${\tt SalePriceMM}$	20.474434
##	ListPriceDiff	15.244049

```
## PctDiscMM
                                  13.723186
## DiscMM
                                  12.957004
## PriceMM
                                   7.244905
## SpecialMM
                                   7.151039
## WeekofPurchase
                                   5.708685
## SpecialCH
                                   5.376632
## STORE
                                   5.077294
## PriceCH
                                   2.477877
## PctDiscCH
                                   1.722996
## DiscCH
                                   1.656462
## SalePriceCH
                                   1.369650
#varImpPlot(tree)
```

#### Question 1.6

Using all the data (training and test put together), use the boosting approach to do the same analysis again and compare the results of the error rates between all tested methods.

## Section 2

### Question 2.1

Create a training and test dataset of size 300 and 100 respectively using the following code:

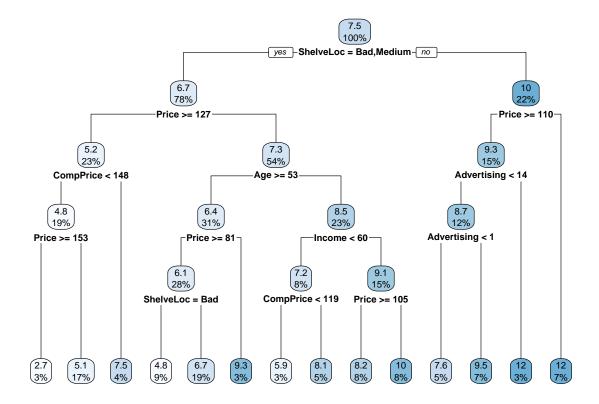
```
set.seed(123456)
n=nrow(Carseats)
id.train=sample(1:n,size=300)
id.test=setdiff(1:n,id.train)
Carseat.train=Carseats[id.train,]
Carseat.test=Carseats[id.test,]
```

### Question 2.2

Construct a regression tree predicting the variable Sales from the other variables available in the training sample. Graph the tree and interpret the results.

# ANSWER: \*\*\* \*\*\*

```
tree.reg <- rpart(Sales ~ . ,data = Carseat.train, method = 'anova')
rpart.plot(tree.reg)</pre>
```



## Question 2.3

Calculate the MSE on the test data.

```
pred.reg.tree = predict(tree.reg, newdata = Carseat.test)
mean((pred.reg.tree - Carseat.test$Sales)^2)
```

## [1] 4.297449

#### Question 2.4

Use cross-validation to determine the optimal level of complexity parameter to prune the tree. Does pruning improve the MSE?

```
set.seed(400)
data_ctrl = trainControl(method = "cv", number = 10)
mytree.caret = train(Sales~.,data=Carseat.train, method = "rpart", trControl=data_ctrl)

## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.

#mytree.caret$resample
#mytree.caret$results
```

#### Question 2.5

## [1] 4.849398

Repeat the previous question several times (with a different seed) - and look at the differences between the results obtained. What do you conclude?

```
#set.seed(0)
seed.list = sample.int(100, 10)
df = data.frame()
for (seed in seed.list){
  set.seed(seed)
  data_ctrl = trainControl(method = "cv", number = 10)
  mytree.caret = train(Sales~., data=Carseat.train, method = "rpart", trControl=data_ctrl)
  #mytree.caret$resample
  #mytree.caret$results
  mytree.caret$bestTune
  pred.cv.tree = predict(mytree.caret, newdata = Carseat.test)
  mse = mean((pred.cv.tree - Carseat.test$Sales)^2)
  output = c(seed,mse)
  df = rbind(df, output)
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
```

```
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
## Warning in nominalTrainWorkflow(x = x, y = y, wts = weights, info = trainInfo, :
## There were missing values in resampled performance measures.
colnames(df)<-c("seed", "mse")</pre>
df
##
      seed
                mse
## 1
        19 4.849398
## 2
        86 4.849398
## 3
        29 4.849398
        52 4.849398
## 4
        25 4.849398
## 5
## 6
        73 4.849398
## 7
        72 4.849398
## 8
        26 4.849398
## 9
        81 4.849398
## 10
        34 4.849398
```

#### Question 2.6

Repeat the previous question several times (with a different seed) - and look at the differences between the results obtained. What do you conclude?

## [1] 2.004661

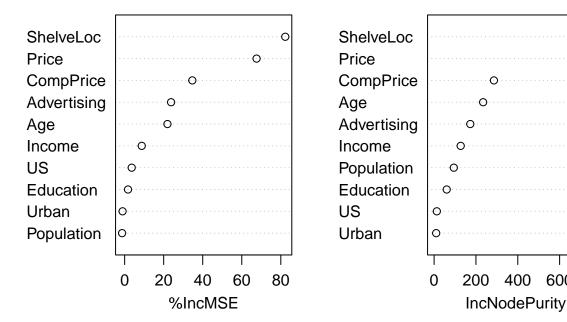
#### Question 2.7

Use the *importance* function to evaluate the importance of variables. Compare with the tree constructed at the beginning of the exercise.

```
as.data.frame(importance(bag.reg))
```

```
%IncMSE IncNodePurity
##
## CompPrice
              34.671069
                            285.920479
## Income
                8.766851
                            127.099848
## Advertising 23.797787
                            172.906967
## Population -1.299732
                            94.290903
## Price
              67.591794
                            668.723626
## ShelveLoc
              82.286261
                            779.935615
## Age
              21.926814
                            234.454410
## Education
               1.736532
                             60.274706
## Urban
              -1.056307
                              9.904163
## US
               3.621940
                             13.064970
```

# bag.reg



```
### Question 2.8 Repeat the 2 previous questions using the random forest algorithm. Comment.**
set.seed(1)
n_pred <- ncol(Carseat.train) - 1</pre>
rf.reg <- randomForest(Carseat.train$Sales ~ .,</pre>
                             data = Carseat.train,
                             #mtry=n_pred,
                             importance=TRUE)
rf.reg.pred = predict(rf.reg, newdata = Carseat.test)#, type = c("anova"))
mean((rf.reg.pred - Carseat.test$Sales)^2)
```

600

800

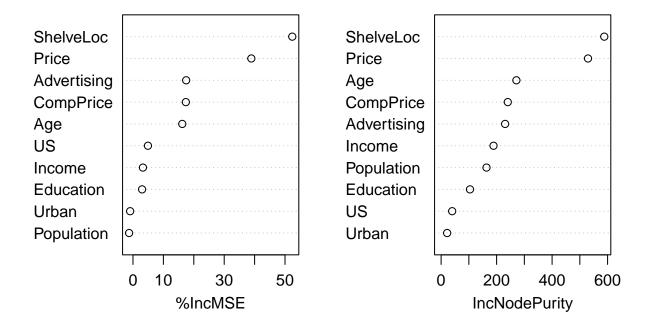
# ## [1] 2.084574

as.data.frame(importance(rf.reg))

```
##
                  %IncMSE IncNodePurity
## CompPrice
               17.3901890
                              240.34938
## Income
                3.2732690
                              188.93112
## Advertising 17.4952006
                              230.71428
## Population -1.2745196
                              163.59608
## Price
                              529.45620
               38.9194656
## ShelveLoc
               52.3866656
                              588.06611
## Age
               16.2092317
                              271.36590
## Education
                2.9971029
                              104.13966
```

```
## Urban -0.9348852 21.80153
## US 4.9195707 39.92127
varImpPlot(rf.reg)
```

# rf.reg



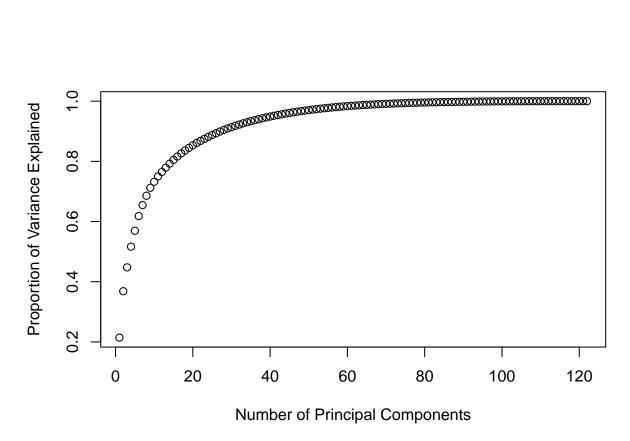
### Section 3

#### Question 3.1

Take the data and example from the principal component regression tutorial. Using the graph of the number of PCs versus the cumulative proportion of variance explained, identify a number of PCs at which the variance explained seems to increase only marginally (note: I don't expect you all to have the same answer). Run a principal component regression with this number of PCs and compare your results with those in the tutorial (MSE of the validation set).

# ANSWER: I would save the first 50 components

```
set.seed(60603)
id=sample(1:nrow(crimeData))
crimeData.pca = crimeData[id[1:1450],][,-123]
crimeData.train=data.frame(apply(crimeData[id[1:1450],],2,scale))
crimeData.valid=data.frame(apply(crimeData[id[1451:nrow(crimeData)],],2,scale))
pca.res = prcomp(crimeData.pca, center = TRUE, scale. = TRUE)
#summary(pca.res)$importance
#round(pca.res$rotation,2)
#data.frame(pca.res$x)
plot(summary(pca.res)$importance[3,], xlab = "Number of Principal Components", ylab = "Proportion of Value")
plot(summary(pca.res)$importance[3,], xlab = "Number of Principal Components", ylab = "Proportion of Value")
```



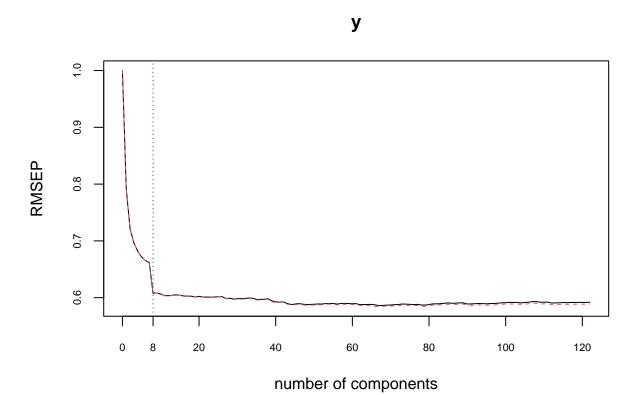
## Question 3.2

It is also possible to choose the number of optimal PCs by using other criteria than the explained variance, such as the MSEP (mean squared error of prediction) or the R2 calculated by cross validation. You can obtain the graph of the number of PCs versus one of these measures with the following command (for the MSEP):

## ANSWER: I would save the first 50 components

```
set.seed(60603)
id=sample(1:nrow(crimeData))
crimeData.pca = crimeData[id[1:1450],][,-123]
crimeData.train=data.frame(apply(crimeData[id[1:1450],],2,scale))
crimeData.valid=data.frame(apply(crimeData[id[1451:nrow(crimeData)],],2,scale))
pca.res = prcomp(crimeData.pca, center = TRUE, scale. = TRUE)

model.pcr=pcr(y~., data = crimeData.train, validation="CV")
validationplot(model.pcr, val.type="RMSEP", cex.axis=0.7)
axis(side = 1, at = c(8), cex.axis=0.7)
abline(v = 8, col = "blue", lty = 3)
```



#### Question 3.3

Repeat the previous question using the R2, with the option val.type="R2" ANSWER: I would save the first 50 components

#"'{r} set.seed(60603) id=sample(1:nrow(crimeData)) crimeData.pca = crimeData[id[1:1450],][,-123] crimeData.train=data.frame(apply(crimeData[id[1:1450],],2,scale)) crimeData.valid=data.frame(apply(crimeData[id[1451:nrow(crimeData.pca, center = TRUE, scale. = TRUE))

 $model.pcr=pcr(y\sim.,\ data=crimeData.train,\ validation="CV")\ validationplot(model.pcr,\ val.type="R2",\ cex.axis=0.7)\ axis(side=1,\ at=c(8),\ cex.axis=0.7)\ abline(v=8,\ col="blue",\ lty=3)\ \#"`$