# Homework week 4

2022-06-11

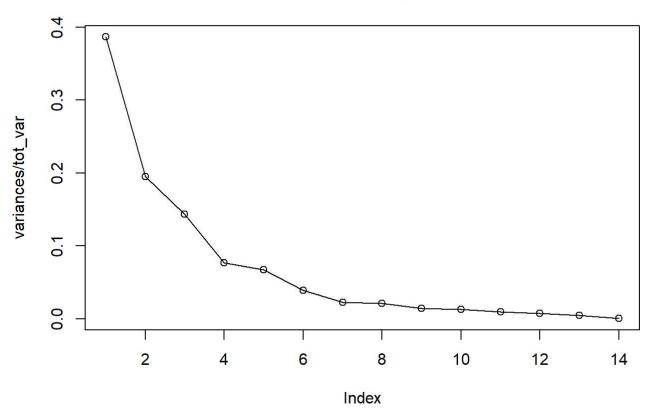
#### Question 9.1

Answer:

First, load the data and perform PCA. We need to exclude the response variable (column 16), the binary variable (column 2) and scale the data before implementing PCA.

Then we display the % of variance explained by each component in a graph.

# % of variance explained



From the line chart above we can see that the first 6 principle components explain most the variances in the data set. Therefore, we will use the first 6 principle components in our model.

```
# Load the first 6 principle components and the response variable (Crime)
# into one data frame
crime_pca_df <- cbind(crime_pca$x[,1:6], data["Crime"])
head(crime_pca_df)</pre>
```

```
##
       PC1
               PC2
                        PC3
                                PC4
                                       PC5
                                               PC6 Crime
## 1 -3.893446 -1.29197714 -1.10138991 0.85371781 -0.2582188 -0.2220763
                                                   791
1635
578
    3.951699 -2.29488126 0.22720214 -0.04741697 -1.7050714 -0.6698441 1969
## 5 1.647039 1.44338543 1.25954528 0.64890563 -0.1162943 0.4327958 1234
## 6 2.861367 -0.11765453 0.51823342 1.44813377 1.0732058 0.6906828
                                                   682
```

Next, we will perform a linear regression using the data frame generated in last step.

```
model <- lm(Crime~PC1+PC2+PC3+PC4+PC5+PC6, data = crime_pca_df)
summary(model)</pre>
```

```
##
## Call:
## lm(formula = Crime ~ PC1 + PC2 + PC3 + PC4 + PC5 + PC6, data = crime pca df)
##
## Residuals:
      Min
##
                1Q Median
                                3Q
                                      Max
## -399.15 -166.78 15.28 150.91 452.53
##
## Coefficients:
               Estimate Std. Error t value Pr(>|t|)
##
                            36.058 25.101 < 2e-16 ***
## (Intercept) 905.085
                                    4.898 1.64e-05 ***
## PC1
                76.750
                            15.668
                -57.648
## PC2
                            22.072 -2.612
                                            0.0126 *
## PC3
                24.313
                            25.744
                                    0.944
                                            0.3506
                            35.157
## PC4
                  3.786
                                    0.108
                                            0.9148
## PC5
               -235.831
                            37.674 -6.260 2.04e-07 ***
                64.174
## PC6
                            49.221
                                    1.304
                                            0.1998
## ---
                  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
## Signif. codes:
##
## Residual standard error: 247.2 on 40 degrees of freedom
## Multiple R-squared: 0.6448, Adjusted R-squared: 0.5915
## F-statistic: 12.1 on 6 and 40 DF, p-value: 1.036e-07
```

Then we need to find the coefficients for each original variable using the results above from both PCA and linear regression model.

```
#Find coefficients for orginal variables(scaled) by multiplying
#coeffients of principle components with eighenvectors
beta_vec <- model$coefficients[2:7]
beta_intercept <- model$coefficients[1]

alpha_vec <- c((crime_pca$rotation[,1:6] %*% beta_vec)[,1])

#display coefficients for original factors, without removing effect of scaling
as.data.frame(alpha_vec)</pre>
```

```
##
          alpha_vec
          95.400235
## M
## Ed
          13.470475
## Po1
         124.092698
## Po2
         119.580606
## LF
         43.199820
         107.629381
## M.F
## Pop
          28.786137
## NW
        103.806024
## U1
          2.756535
## U2
          29.061629
## Wealth 38.505494
## Ineq
           8.489242
## Prob
         -43.150501
          35.340376
## Time
```

Considering that the data has been scaled before applying PCA to it, we need to execute following steps to remove the effect of scaling:

```
#Calculate the means and standard deviations of the variables in the orginal
#data, excluding response variable (Crime) and binary variable (So)
means <- colMeans(data[,c(-2,-16)])
stdev <- apply(data[,c(-2,-16)],2,sd)

#remove effect of scaling
origin_vec <- alpha_vec/stdev
as.data.frame(origin_vec)</pre>
```

```
##
             origin_vec
           7.590946e+01
## M
## Ed
           1.204119e+01
## Po1
           4.175538e+01
## Po2
           4.276644e+01
## LF
           1.068990e+03
## M.F
           3.652494e+01
## Pop
           7.561134e-01
           1.009503e+01
## NW
## U1
           1.528963e+02
## U2
           3.441099e+01
## Wealth 3.990581e-02
           2.127840e+00
## Ineq
## Prob
          -1.897812e+03
## Time
           4.986722e+00
```

```
origin_intercept <- beta_intercept - sum(alpha_vec * means / stdev)
#display coefficients for original factors, removing effect of scaling
as.data.frame(origin_intercept)</pre>
```

```
## origin_intercept
## (Intercept) -5717.968
```

The results above shows the coefficients and intercepts for the original factors. Next, we are going to using them to estimate the crime rate in each row of the original data set (without binary variable, column 2)

Then we are going to compare our estimated result with real results, and calculate adjusted R squared.

```
input <- as.matrix(data[, c(-2,-16)])
estimate <- input %*% origin_vec + origin_intercept

RSS <- sum((data[,16]-estimate)^2)
TSS <- sum((data[,16]-mean(data[,16]))^2)

R2 <- 1 - RSS/TSS
R2</pre>
```

```
## [1] 0.644773
```

```
Adj_R2 = 1- (1-R2)*nrow(data)/(nrow(data)-6-1)
Adj_R2
```

```
## [1] 0.5826083
```

As the last step of this question, we are going to make a prediction with the data of the new city.

```
input <- as.matrix(data.frame("M" = 14.0,</pre>
                                "Ed" = 10.0,
                                "Po1" = 12.0,
                                "Po2" = 15.5,
                                "LF" = 0.64,
                                "M.F" = 94.0,
                                "Pop" = 150,
                                "NW" = 1.1,
                                "U1" = 0.12,
                                "U2" = 3.6,
                                "Wealth" = 3200,
                                 "Ineq" = 20.1,
                                 "Prob" = 0.04,
                                "Time" = 39.0))
prediction <- input %*% origin_vec + origin_intercept</pre>
prediction
```

```
## [,1]
## [1,] 1302.405
```

## Question 10.1

#### Answer:

First, we read the data and run a regression tree model by calling tree() function. Since later we are going to use cross validation to prune the tree model, we need to split the data into 2 parts, 70% for training and cross validation, the rest 30% for testing.

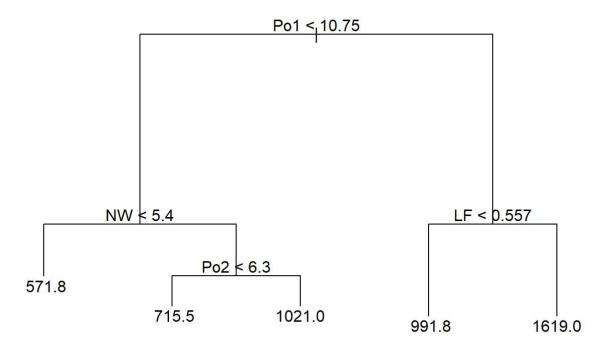
We are going to show the nodes, and branching criteria in the graph below.

```
## node), split, n, deviance, yval
        * denotes terminal node
##
##
   1) root 33 5601000 929.1
##
##
      2) Po1 < 10.75 23 1539000 765.5
##
        4) NW < 5.4 9 318300 571.8 *
##
        5) NW > 5.4 14 666100 890.0
        10) Po2 < 6.3 6 140900 715.5 *
##
        11) Po2 > 6.3 8 205400 1021.0 *
##
##
     3) Po1 > 10.75 10 2031000 1305.0
##
       6) LF < 0.557 5 478200 991.8 *
##
        7) LF > 0.557 5 568900 1619.0 *
```

## summary(tree)

```
##
## Regression tree:
## tree(formula = Crime ~ ., data = training)
## Variables actually used in tree construction:
## [1] "Po1" "NW" "Po2" "LF"
## Number of terminal nodes: 5
## Residual mean deviance: 61130 = 1712000 / 28
## Distribution of residuals:
## Min. 1st Qu. Median Mean 3rd Qu. Max.
## -394.00 -171.90 -29.78  0.00 133.20 563.20
```

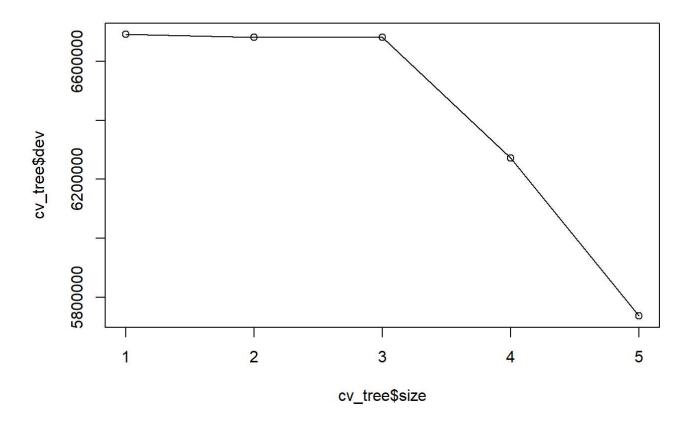
```
plot(tree)
text(tree, pretty = 0)
```



Next, we need to prune the tree and choose the best number of nodes. Here we use cv.tree() to perform the pruning and cross validation process at the same time. We will use deviance to measure the quality of each model.

We will do a 3-fold cross validation and show the deviance of the model under each number of nodes in a line chart.

```
set.seed(10)
#prune the tree & cross validation
cv_tree <- cv.tree(tree,K=3)
plot(cv_tree$size, cv_tree$dev, type = "o")</pre>
```



From the chart above, we can see that the model has lowest deviation when there are 5 nodes in total. Therefore, we will choose have 5 nodes in our tree. We will show this model by calling prune.tree() function with parameter best = 5.

The structure of the tree will be displayed in a graph.

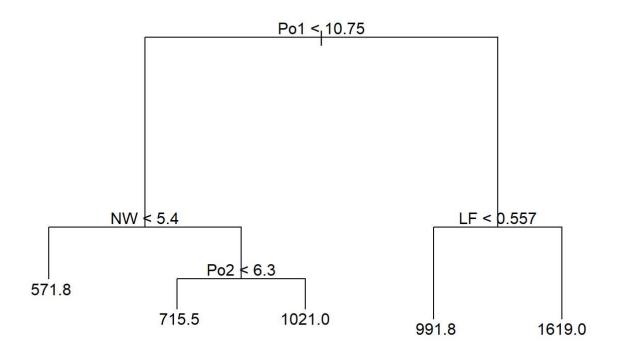
```
tree_pruned <- prune.tree(tree, best = 5)
tree_pruned</pre>
```

```
## node), split, n, deviance, yval
##
         * denotes terminal node
##
    1) root 33 5601000 929.1
##
##
      2) Po1 < 10.75 23 1539000 765.5
##
        4) NW < 5.4 9 318300 571.8 *
        5) NW > 5.4 14 666100 890.0
##
##
         10) Po2 < 6.3 6 140900 715.5 *
##
         11) Po2 > 6.3 8 205400 1021.0 *
##
      3) Po1 > 10.75 10 2031000 1305.0
        6) LF < 0.557 5 478200 991.8 *
##
        7) LF > 0.557 5 568900 1619.0 *
##
```

```
summary(tree_pruned)
```

```
##
## Regression tree:
## tree(formula = Crime ~ ., data = training)
## Variables actually used in tree construction:
## [1] "Po1" "NW" "Po2" "LF"
## Number of terminal nodes: 5
## Residual mean deviance: 61130 = 1712000 / 28
## Distribution of residuals:
## Min. 1st Qu. Median Mean 3rd Qu. Max.
## -394.00 -171.90 -29.78  0.00 133.20 563.20
```

```
plot(tree_pruned)
text(tree_pruned)
```



```
# Calculate R squared
prediction <- predict(tree_pruned, testing[,1:15])
RSS <- sum((testing[,16] - prediction)^2)
TSS <- sum((testing[,16] - mean(testing[,16]))^2)
R2 <- 1- RSS/TSS
R2</pre>
```

From the charts and tables in previous steps, we can have following findings:

- 1. "Po1" is probably the most important factor in the regression tree model. It is the node at the top of the tree. When Po1 < 10.75, the overall estimation of crime rate is much lower comparing to when Po1>=10.75.
- 2. Comparing to the deviation we see in cross validation (5736093), the deviation of the model shown in the last step (1712000) is much lower. That means our model may have a issue of overfitting to training data. The R squared value shown in the testing result also implies this issue. It is as low as only 4.58%.

In the next part of this questions. We are go to run a random forest model using the same training / testing data. To select the best model, we are going to change the values of mytry and nodesize parameters. We are going to select the combination of these two parameters that generates the best R squared value.

In the following code, we are going to try all the combinations of mytry between 1 to 14, and nodesize between 2 and 10.

```
library(randomForest)
```

```
## randomForest 4.7-1.1
```

## Type rfNews() to see new features/changes/bug fixes.

```
set.seed(19)
mtry_vec <- c()</pre>
nodesize vec <- c()</pre>
R2 vec <- c()
#try all possible combinations of mytry and nodesize
for (m in 1:14){
  for (n in 2:10){
    rf <- randomForest(Crime~., data = training, importance = TRUE, mtry = m,</pre>
                        nodeszie = n)
    #calculate R squared value
    RSS <- sum((training[,16] - rf$predicted)^2)</pre>
    TSS <- sum((training[,16] - mean(training[,16]))^2)
    R2 <- 1- RSS/TSS
    mtry_vec <- append(mtry_vec,m)</pre>
    nodesize vec <- append(nodesize vec, n)</pre>
    R2_vec = append(R2_vec, R2)
  }
}
training_result_df <- data.frame("mytry" = mtry_vec,</pre>
                          "node size" = nodesize vec,
                          "R2" = R2 vec)
#display part of the training results
head(training_result_df)
```

```
##
     mytry node.size
                              R2
## 1
         1
                    2 0.3027683
## 2
         1
                    3 0.3569474
## 3
         1
                    4 0.3645463
## 4
         1
                    5 0.3255773
## 5
         1
                    6 0.3370738
## 6
                    7 0.3669648
```

After reviewing all the results, we will find that the combination mytry=3 and nodesize=6 generates the best R squared value. (41.05%) using training data.

Therefore, we will choose mytry = 3 and nodesize = 6, and display the model in more details, running with testing data.

```
##
## Call:
    randomForest(formula = Crime ~ ., data = testing, importance = TRUE,
##
                                                                              mtry = 3, nodeszie
= 6)
##
                  Type of random forest: regression
                        Number of trees: 500
##
## No. of variables tried at each split: 3
##
##
             Mean of squared residuals: 76002.87
                       % Var explained: 12.47
##
```

# rf\$importance

```
%IncMSE IncNodePurity
##
## M
            732.10650
                            38708.27
## So
             23.87658
                             5132.08
## Ed
            363.16270
                            37132.30
## Po1
          10483.79150
                           197348.74
## Po2
          13726.27332
                           157622.28
## LF
           -657.93964
                            29230.19
## M.F
          -1144.77825
                            39535.88
## Pop
          -1156.57353
                            39786.06
## NW
           -381.29790
                            47953.62
## U1
                            35077.22
            472.13250
## U2
           2054.21812
                            57467.56
## Wealth 7482.76063
                            83363.60
## Ineq
          -1950.20277
                            37966.50
           5483.90492
## Prob
                            95448.39
## Time
            477.94403
                            52282.65
```

```
#Calculate R squared using testing data
RSS <- sum((testing[,16] - rf$predicted)^2)
TSS <- sum((testing[,16] - mean(testing[,16]))^2)
R2 <- 1- RSS/TSS
R2</pre>
```

```
## [1] 0.1247075
```

Reviewing the table above and compare the random forest tree model with the regression tree mode, we can have following findings:

- 1. The random forest model has better quality of fit than regression tree model since it has a higher R squared value running on testing data (12.47%)
- 2. In the random forest model, factors "Po1" and "Po2" are two most important factors, since they have the highest value of "%IncMSE"

# Question 10.2

#### Answer:

I am recently applying for a new credit card. I think a logistic model can be used to predict the probability of whether my application will be approved.

The predictors my include: my gross annual income, monthly housing rent payment, credit score, years of longest credit history, etc.

Question 10.3

Question 10.3.1

Answer:

First, we need to read the data. Since the last column (response variable) has values of 1 and 2, we need to convert them to 0 and 1.

To measure the quality of fit, we also need to split the data into training and testing. Here we will use 70% of data for training and the rest 30% for testing.

```
V5 V6 V7 V8 V9 V10 V11 V12 V13 V14 V15 V16 V17 V18
##
     V1 V2 V3 V4
## 1 A11 6 A34 A43 1169 A65 A75 4 A93 A101
                                             4 A121
                                                    67 A143 A152
                                                                    2 A173
## 2 A12 48 A32 A43 5951 A61 A73 2 A92 A101
                                             2 A121 22 A143 A152
                                                                    1 A173
                                                                             1
## 3 A14 12 A34 A46 2096 A61 A74 2 A93 A101
                                             3 A121 49 A143 A152
                                                                    1 A172
                                                                             2
                                             4 A122 45 A143 A153
## 4 A11 42 A32 A42 7882 A61 A74 2 A93 A103
                                                                             2
                                                                    1 A173
## 5 A11 24 A33 A40 4870 A61 A73 3 A93 A101
                                             4 A124 53 A143 A153
                                                                             2
                                                                    2 A173
## 6 A14 36 A32 A46 9055 A65 A73 2 A93 A101
                                             4 A124 35 A143 A153
                                                                    1 A172
                                                                             2
##
     V19 V20 V21
## 1 A192 A201
## 2 A191 A201
## 3 A191 A201
## 4 A191 A201
## 5 A191 A201
## 6 A192 A201
```

```
#split data into training and testing
training_index <- sample(1:nrow(data), size = round(nrow(data)*0.7))
training <- data[training_index,]
testing <- data[-training_index,]</pre>
```

Next, we will run the logistic regression model using all the variables and see the result. We will test the model use testing data set and compare the prediction results and the real response factor in a confusion matrix. Right now, we don't have a probably threshold for the prediction. Therefore we will just round the predicted probabilities to zero and one.

```
model <- glm(V21~., data = training, family=binomial(link="logit"))
summary(model)</pre>
```

```
##
## Call:
### glm(formula = V21 ~ ., family = binomial(link = "logit"), data = training)
##
## Deviance Residuals:
##
      Min
                1Q
                     Median
                                  3Q
                                          Max
## -2.1933 -0.7031 -0.3556 0.7312
                                       2.4779
##
## Coefficients:
##
                Estimate Std. Error z value Pr(>|z|)
## (Intercept) 1.859e-01 1.297e+00
                                      0.143 0.886093
## V1A12
              -5.424e-01 2.674e-01 -2.029 0.042465 *
## V1A13
              -1.424e+00 4.435e-01 -3.210 0.001326 **
## V1A14
              -1.845e+00 2.836e-01 -6.504 7.84e-11 ***
## V2
               2.054e-02 1.132e-02 1.815 0.069525 .
## V3A31
              -4.140e-01 6.402e-01 -0.647 0.517836
## V3A32
              -3.680e-01 5.167e-01 -0.712 0.476258
## V3A33
              -1.217e+00 5.858e-01 -2.077 0.037791 *
## V3A34
              -1.733e+00 5.378e-01 -3.222 0.001271 **
              -1.759e+00 4.531e-01 -3.882 0.000104 ***
## V4A41
              -1.383e+00 1.179e+00 -1.173 0.240744
## V4A410
## V4A42
              -7.525e-01 3.201e-01 -2.351 0.018736 *
## V4A43
              -4.869e-01 3.001e-01 -1.623 0.104668
## V4A44
               -5.411e-01 8.941e-01 -0.605 0.545061
## V4A45
              -2.986e-01 7.163e-01 -0.417 0.676776
## V4A46
               4.326e-01 4.913e-01 0.880 0.378660
## V4A48
              -1.345e+00 1.216e+00 -1.106 0.268696
## V4A49
              -5.121e-01 3.949e-01 -1.297 0.194731
## V5
               1.852e-04 5.651e-05
                                     3.276 0.001051 **
              -3.102e-01 3.523e-01 -0.881 0.378503
## V6A62
## V6A63
              -3.862e-01 4.641e-01 -0.832 0.405241
## V6A64
              -9.202e-01 6.311e-01 -1.458 0.144799
              -8.523e-01 3.293e-01 -2.588 0.009645 **
## V6A65
## V7A72
               5.377e-01 5.183e-01
                                      1.037 0.299572
## V7A73
               3.398e-01 5.035e-01
                                      0.675 0.499767
              -7.879e-01 5.523e-01 -1.427 0.153716
## V7A74
## V7A75
              -5.283e-02 5.055e-01 -0.105 0.916759
## V8
               3.682e-01 1.069e-01
                                      3.446 0.000569 ***
## V9A92
              -6.305e-01 4.577e-01 -1.378 0.168341
## V9A93
              -8.454e-01 4.480e-01 -1.887 0.059181 .
## V9A94
              -5.725e-01 5.269e-01 -1.087 0.277166
## V10A102
               1.680e-01 5.264e-01
                                      0.319 0.749603
## V10A103
              -9.469e-01 4.890e-01 -1.936 0.052819 .
## V11
               1.023e-01 1.035e-01
                                      0.989 0.322717
## V12A122
              -2.861e-02 3.068e-01 -0.093 0.925695
## V12A123
               7.580e-02 2.783e-01
                                      0.272 0.785333
               3.433e-01 5.225e-01
## V12A124
                                      0.657 0.511088
## V13
              -1.834e-02 1.113e-02 -1.647 0.099466 .
## V14A142
               9.205e-02 4.883e-01
                                      0.189 0.850468
## V14A143
              -6.556e-01 2.903e-01 -2.258 0.023923 *
## V15A152
              -5.805e-01 2.756e-01 -2.107 0.035139 *
## V15A153
              -2.242e-01 5.784e-01 -0.388 0.698257
```

```
## V16
               4.997e-01 2.354e-01 2.123 0.033782 *
## V17A172
              -1.320e-01 7.689e-01 -0.172 0.863697
## V17A173
              -5.099e-02 7.360e-01 -0.069 0.944769
## V17A174
               6.705e-02 7.500e-01 0.089 0.928760
## V18
               3.714e-01 3.082e-01 1.205 0.228106
              -2.753e-01 2.428e-01 -1.134 0.256812
## V19A192
## V20A202
              -1.406e+00 7.724e-01 -1.820 0.068728 .
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## (Dispersion parameter for binomial family taken to be 1)
##
      Null deviance: 877.62 on 699 degrees of freedom
##
## Residual deviance: 630.06 on 651 degrees of freedom
## AIC: 728.06
##
## Number of Fisher Scoring iterations: 5
```

```
#test the model use testing data
predicttion <- predict(model, newdata = testing, type = "response")
table(testing[,21], round(predicttion))</pre>
```

```
##
## 0 1
## 0 195 29
## 1 44 32
```

From the results above, the model correctly predicted 227 data points out of 300. But we can see from the summary table of the model that many predictors do not have a significant coefficients.

Therefore, we will run the model again after excluding all the predictors with p value bigger than 0.01.

For each column in original data, we will only keep the most significant coefficient, since the coefficients of the same column may be correlated with each other. Here's the results:

```
#Manually add significtant predictors to the training / testing data sets
add_coefficients <- function(dataset){</pre>
  dataset$V1A14[dataset$V1 == "A14"] <- 1</pre>
  dataset$V1A14[dataset$V1 != "A14"] <- 0</pre>
  dataset$V3A34[dataset$V3 == "A34"] <- 1</pre>
  dataset$V3A34[dataset$V3 != "A34"] <- 0</pre>
  dataset$V4A41[dataset$V4 == "A41"] <- 1</pre>
  dataset$V4A41[dataset$V4 != "A41"] <- 0</pre>
  dataset$V6A65[dataset$V6 == "A65"] <- 1</pre>
  dataset$V6A65[dataset$V6 != "A65"] <- 0</pre>
  return (dataset)
}
training_2 <- add_coefficients(training)</pre>
testing_2 <- add_coefficients(testing)</pre>
#Rerun the model after excluding insignificant coefficients
model 2 <- glm(V21\sim
                V1A14+V3A34+V4A41+V5+V6A65+V8,
              data = training_2, family=binomial(link="logit"))
summary(model 2)
```

```
##
## Call:
## glm(formula = V21 \sim V1A14 + V3A34 + V4A41 + V5 + V6A65 + V8,
##
       family = binomial(link = "logit"), data = training_2)
##
## Deviance Residuals:
##
      Min
                1Q Median
                                  3Q
                                          Max
## -2.3023 -0.8456 -0.4748 1.0095
                                       2.3581
##
## Coefficients:
##
                Estimate Std. Error z value Pr(>|z|)
## (Intercept) -1.691e+00 3.478e-01 -4.864 1.15e-06 ***
             -1.516e+00 2.201e-01 -6.887 5.69e-12 ***
## V1A14
## V3A34
              -8.850e-01 2.262e-01 -3.912 9.14e-05 ***
## V4A41
              -1.073e+00 3.658e-01 -2.934 0.003348 **
               2.259e-04 3.784e-05 5.970 2.38e-09 ***
## V5
## V6A65
              -8.271e-01 2.900e-01 -2.852 0.004346 **
## V8
               3.434e-01 9.014e-02 3.810 0.000139 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## (Dispersion parameter for binomial family taken to be 1)
##
       Null deviance: 877.62 on 699 degrees of freedom
##
## Residual deviance: 724.90 on 693 degrees of freedom
## AIC: 738.9
##
## Number of Fisher Scoring iterations: 5
```

```
#test the model use testing data
predicttion <- predict(model_2, newdata = testing_2, type = "response")
table(testing_2[,21], round(predicttion))</pre>
```

From the results above, even thought the prediction accuracy didn't improve (225 out of 300), there are no longer insignificant coefficients. That means we have reduced the likely hood of overfitting problem. Therefore, we are going to choose the second model.

# Question 10.3.2

# Answer:

In our prediction above, 0 stands for good and 1 stands for bad.

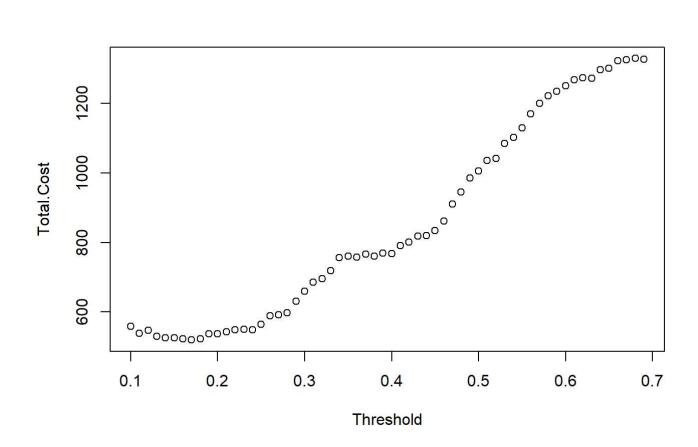
Assume that the cost of incorrectly identifying a bad customer as good is 5. and the cost of incorrectly classifying a good customer as bad is 1. If we do not change the probability threshold, the total cost in the last scenario equals 52 \* 5 + 23 \* 1 = 283

To find the best probability threshold and minimize the total cost, we will try every possible probability threshold

starting from 10% to 70%. If the probability prediction result for one customer is bigger or equal to the threshold, we will identify this customer as "bad" (1). Each time, we will raise the threshold by 1% and calculate the total cost. The probability threshold that minimize the total cost will be our answer.

Let's run second model on the full data set to make the prediction.

```
#add necessary predictors to orginal dataset and run the model
data_2 <- add_coefficients(data)</pre>
pred_prob <- predict(model_2, newdata = data_2, type = "response")</pre>
#try different threshold
p = 0.1
threshold <- c()
cost <- c()
while (p<=0.7){
    prediction = pred_prob
    for (i in 1:length(prediction)){
      if (prediction[i] >= p){
        prediction[i] <- 1</pre>
      }
      else {prediction[i] <- 0}</pre>
    }
    con mat <- table(data 2[,21], prediction)</pre>
    total_cost <- con_mat[1,2] * 1 + con_mat[2,1] * 5
    threshold <- append(threshold, p)</pre>
    cost <- append(cost, total_cost)</pre>
    p = p + 0.01
}
results <- data.frame("Threshold" = threshold, "Total Cost" = cost)</pre>
plot(results)
```



results

##	Threshold	Total.Cost
## 1	0.10	558
## 2	0.11	538
## 3	0.12	547
## 4	0.13	529
## 5	0.14	525
## 6	0.15	525
## 7	0.16	522
## 8	0.17	5 <b>1</b> 9
## 9	0.18	522
## 10	0.19	537
## 11	0.20	537
## 12	0.21	542
## 13	0.22	548
## 14	0.23	550
## 15	0.24	548
## 16	0.25	564
## 17	0.26	589
## 18	0.27	592
## 19	0.28	597
## 20	0.29	631
## 21	0.30	659
## 22	0.31	685
## 23	0.32	696
## 24	0.33	719
## 25	0.34	756
## 26	0.35	760
## 27	0.36	757
## 28	0.37	766
## 29	0.38	760
## 30	0.39	769
## 31	0.40	767
## 32	0.41	791
## 33	0.42	801
## 34	0.43	818
## 35	0.44	820
## 36	0.45	834
## 37	0.46	862
## 38	0.47	911
## 39	0.48	945
## 40	0.49	986
## 41	0.50	1006
## 42	0.51	1036
## 43	0.52	1042
## 44	0.53	1085
## 45	0.54	1102
## 46	0.55	1130
## 47	0.56	1171
## 48	0.57	1201
## 49	0.58	1223
## 50	0.59	1236
## 51	0.60	1251

##	52	0.61	1269
##	53	0.62	1274
##	54	0.63	1273
##	55	0.64	1297
##	56	0.65	1301
##	57	0.66	1324
##	58	0.67	1326
##	59	0.68	1330
##	60	0.69	1328

From the results above, we can easily observe that the total cost will be minimized (equals 519) when the threshold is 17%. Therefore, we would choose the threshold to be 17%