1 Getting data

First of all, we load all the packages we are going to use and we look at the data

There are three types of NAs in the data sets; "NA", "#DIV/0!" and an empty space "". We classify all of those as "NA".

Moreover, the first seven columns of the data sets are just information about users and time/dates, which are not important for the prediction of classe, therefore we do not need those.

2 Cleaning data

Firstly, we get rid of the columns that with more than 40% NA.

```
> data_NA <- apply(is.na(trainData), 2, sum) > 0.4*nrow(trainData)
> train <- (trainData[,!data_NA])
> test <- (testData[,!data_NA])</pre>
```

Secondly, we check the correlation between the predictors. We are interested in pairs of predictors with correlation over 0.80.

```
> train_matrix <- as.matrix(train[,-53])
> correlation <- pcor(train_matrix, method="pearson")$estim
> diag(correlation) <- 0
> which((abs(correlation))>0.80, arr.ind=TRUE)
```

```
    gyros_arm_y
    19
    18

    gyros_arm_x
    18
    19

    magnet_arm_z
    26
    23

    accel_arm_z
    23
    26

    gyros_dumbbell_z
    33
    31

    gyros_dumbbell_x
    31
    33

    magnet_forearm_x
    50
    47

    accel_forearm_x
    47
    50
```

Since we work out ourselves and have read the paper Wearable Qualitative Activity Recognition of Weight Lifting Exercises as well as about methods and devices used for the measurements, we decide to eliminate following predictors: " $gyros_arm_x$ ", " $magnet_arm_z$ ", " $gyros_dumbbell_x$ " and " $magnet_forearm_x$ ".

```
> ## we eliminate gyros_arm_x, magnet_arm_z, gyros_dumbbell_x and magnet_forearm_x
> train_noncorr <- train[,-c(18,26,31,47)]</pre>
```

We use this dataset with non-correlated columns to build our model.

3 Building the model

For the cross validation we use random subsampling approach, i.e. splitting our "train" data into two subsets: 60% used for the model building and 40% for the model testing. (the "test" data set will be used later for the prediction). Also, we change "classe" to factor variable.

```
> set.seed(123)
> inTrain <- createDataPartition(train_noncorr$classe, p=0.6, list=F)
> training <- train_noncorr[inTrain,]
> testing <- train_noncorr[-inTrain,]
> dim(training); dim(testing)

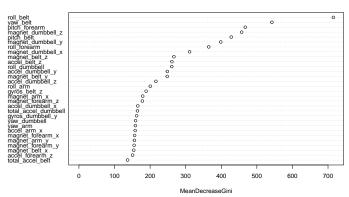
[1] 11776     49

[1] 7846     49
> training$classe <- as.factor(training$classe)</pre>
```

We decided to use random forest approach since it produces the most accurate results.

After eliminating NAs and correlated predictors we end up with 49 predictors, which is still quite a lot. Therefore we order them according to their importance and decide whether we can simplify the model.

```
> set.seed(123)
> RF <- randomForest(classe~., data=training, ntree=500)
> varImpPlot(RF)
```



There is a sig-

nificant break after the 13th predictor, thus we use the first 13 of them. We create the final dataset with the predictors and "classe".

```
> df <- as.data.frame(RF$importance)
> TRAINING <- training[,c(order(df$MeanDecreaseGini, decreasing = T)[1:13], 49)]</pre>
```

Finally, we build the model. In order to improve the performance of random forest approach, we have decided to use bagging.

```
> ## bootstrapped sample size 1/p of the original number of observations,
> ## "iter" is number of iterations
> bagg_RF <- function(training_data,testing_data,p,iter)</pre>
    predictions <- foreach(m=1:iter,.combine=cbind) %do% {</pre>
      sampled_positions <- sample(nrow(training_data), size=floor((nrow(training_data)/p)))</pre>
      training_positions <- 1:nrow(training_data) %in% sampled_positions</pre>
      set.seed(123)
      ## we have chosen randomForest instead of train(..., method="rf",...)
      ## because it is quicker and we get slightly better results
      RF_fit <- randomForest(classe~., data=training_data[training_positions,], ntree=500)
      predict(RF_fit,newdata=testing_data)
    Predictions_num <- round(rowMeans(predictions))</pre>
    Predictions <- as.vector(length(Predictions_num))</pre>
    for(i in 1:length(Predictions_num)){
      if(Predictions_num[i]==1) Predictions[i]="A"
      if(Predictions_num[i]==2) Predictions[i]="B"
      if(Predictions_num[i]==3) Predictions[i]="C"
      if(Predictions_num[i]==4) Predictions[i]="D"
      if(Predictions_num[i]==5) Predictions[i]="E"
    Predictions
```

We use our model to predict on the "testing" dataset for accuracy.

- > testing_pred <- bagg_RF(TRAINING,testing,p=4,iter=100)</pre>
- > confusionMatrix(testing\$classe, testing_pred)

Confusion Matrix and Statistics

Reference Prediction A B C D E A 2184 44 4 0 0 B 41 1386 70 19 2 C 0 52 1290 26 0 D 4 3 48 1231 0

E 2 2 13 33 1392

Overall Statistics

Accuracy : 0.9537

95% CI : (0.9489, 0.9583)

No Information Rate : 0.2843 P-Value [Acc > NIR] : < 2.2e-16

Kappa: 0.9415 Mcnemar's Test P-Value: 1.993e-12

Statistics by Class:

	Class A	Class B	Class: C	Class D	Class E
Sensitivity	0.9789	0.9321	0.9053	0.9404	0.9986
Specificity	0.9915	0.9792	0.9879	0.9916	0.9923
Pos Pred Value	0.9785	0.9130	0.9430	0.9572	0.9653
Neg Pred Value	0.9916	0.9840	0.9792	0.9881	0.9997
Prevalence	0.2843	0.1895	0.1816	0.1668	0.1777
Detection Rate	0.2784	0.1767	0.1644	0.1569	0.1774
Detection Prevalence	0.2845	0.1935	0.1744	0.1639	0.1838
Balanced Accuracy	0.9852	0.9557	0.9466	0.9660	0.9954

We get the accuracy 95.37% on the "testing" set, which is satisfactory. The very last thing to do on this project is to predict "classe" for the "testData".

> bagg_RF(TRAINING,testData,p=4,iter=100)

[1] "B" "A" "B" "A" "A" "E" "D" "B" "A" "A" "B" "C" "B" "A" "E" "E" "A" "B" "B" [20] "B"