# Machine Learning Prediction Project

August 16, 2017

# Summary

We have investigated two models or methods, classification tree and random forest tree, to find out that which method has a better accuracy to predict outcomes. To demonstrate this, both modes were tested on 30% data of training set for cross validation of predictions. It has been observed that random forest tree provides bettre accuracy 99% (0.99) than classification tree method with 49% (0.49) accuracy. Random forest tree model was used to prdict the outcomes of test data set.

# Loading libraries and data sets

```
library(caret); library(rpart); library(randomForest)
training <- read.csv("pml-training.csv", na.strings = c("NA", ""))
testing <- read.csv("pml-testing.csv", na.strings = c("NA", ""))</pre>
```

# Data cleaning

In this step, columns with missing values have been removed from both the test and train data sets, and also variables those are not important or required for the model calculations have been delted form the data sets. This process reduces the calculation time for model and make it computationally efficient.

```
training <- training[ ,-c(1:7)]
new.training <- training[, colSums(is.na(training)) == 0]
testing <- testing[ ,-c(1:7)]
testing <- testing[, colSums(is.na(testing)) == 0]
new.testing <- testing2[ ,-53]</pre>
```

# Data partitioned

After cleaning the data, the training data set is divided into two data sets, (70% training.data and 30% valid.training.data) for training and prediction with models used in this data analysis.

```
inTrain <- createDataPartition(new.training$classe, p = 0.7, list =FALSE)
training.data <- new.training[inTrain,]
Valid.training.data <- new.training[-inTrain,]</pre>
```

# 1. Prediction with clssification tress using validation data set

We used trainControl function with K=5 fold and rpart method for cross validation of predictions

```
Control<- trainControl(method = "cv", number = 5)
Fit.Model <- train(classe ~.,method = "rpart", data =training.data,trControl = Control)
print(Fit.Model, digit=4)</pre>
```

CART

#### 13737 samples

52 predictor

5 classes: 'A', 'B', 'C', 'D', 'E'

No pre-processing

Resampling: Cross-Validated (5 fold)

Summary of sample sizes: 10989, 10989, 10991, 10989, 10990

Resampling results across tuning parameters:

ср	Accuracy	Kappa
0.03428	0.5097	0.3601
0.06076	0.4155	0.2077
0.11647	0.3328	0.0738

Accuracy was used to select the optimal model using the largest value. The final value used for the model was cp = 0.03428.

prediction <- predict(Fit.Model, Valid.training.data)
classification <- confusionMatrix(prediction, Valid.training.data\$classe)
classification</pre>

Confusion Matrix and Statistics

#### Reference

Prediction	Α	В	C	D	Ε
A	1509	454	460	448	155
В	41	383	32	163	147
C	118	302	534	353	302
D	0	0	0	0	0
F.	6	0	0	0	478

Overall Statistics

Accuracy : 0.4935

95% CI : (0.4806, 0.5063)

No Information Rate : 0.2845 P-Value [Acc > NIR] : < 2.2e-16

Kappa : 0.3385

Mcnemar's Test P-Value : NA

Statistics by Class:

	Class:A	Class:B	Class:C	Class:D	Class:E
Sensitivity	0.9014	0.33626	0.52047	0.0000	0.44177
Specificity	0.6398	0.91930	0.77876	1.0000	0.99875
Pos Pred Value	0.4987	0.50000	0.33188	NaN	0.98760
Neg Pred Value	0.9423	0.85231	0.88494	0.8362	0.88817
Prevalence	0.2845	0.19354	0.17434	0.1638	0.18386
Detection Rate	0.2564	0.06508	0.09074	0.0000	0.08122
Detection Prevalence	0.5142	0.13016	0.27341	0.0000	0.08224
Balanced Accuracy	0.7706	0.62778	0.64961	0.5000	0.72026

This method has a 50 % acuracy and out of sample error is 0.50 (50 %). This classification tree doesn't do good job in prediction of outcomes. Next, we will experiment with random forest tree to obtain better accuracy in prediction than calssification tree method.

# 2. Prediction with random forest tree using validation data set

```
Fit.Model2 <- train(classe ~.,method = "rf", data =training.data,trControl = Control)
prediction2 <- predict(Fit.Model2, Valid.training.data)
print(Fit.Model2, digit=4)</pre>
```

Random Forest

13737 samples 52 predictor

5 classes: 'A', 'B', 'C', 'D', 'E'

No pre-processing

Resampling: Cross-Validated (5 fold)

Summary of sample sizes: 10990, 10990, 10989, 10989, 10990

Resampling results across tuning parameters:

$\mathtt{mtry}$	Accuracy	Kappa
2	0.9903	0.9878
27	0.9909	0.9885
52	0.9815	0.9766

Accuracy was used to select the optimal model using the largest value.

The final value used for the model was mtry = 27.

classification2 <- confusionMatrix(prediction2, Valid.training.data\$classe)
classification2</pre>

Confusion Matrix and Statistics

### ${\tt Reference}$

${\tt Prediction}$	Α	В	C	D	Ε
A	1672	1	0	0	0
В	2	1135	3	0	0
C	0	3	1021	13	1
D	0	0	2	950	6
Е	0	0	0	1	1075

Overall Statistics

Accuracy : 0.9946

95% CI: (0.9923, 0.9963)

No Information Rate : 0.2845 P-Value [Acc > NIR] : < 2.2e-16

Kappa: 0.9931

Mcnemar's Test P-Value : NA

#### Statistics by Class:

	Class: A	Class: B	Class: C	Class: D	Class: E
Sensitivity	0.9988	0.9965	0.9951	0.9855	0.9935
Specificity	0.9998	0.9989	0.9965	0.9984	0.9998
Pos Pred Value	0.9994	0.9956	0.9836	0.9916	0.9991
Neg Pred Value	0.9995	0.9992	0.9990	0.9972	0.9985
Prevalence	0.2845	0.1935	0.1743	0.1638	0.1839
Detection Rate	0.2841	0.1929	0.1735	0.1614	0.1827
Detection Prevalence	0.2843	0.1937	0.1764	0.1628	0.1828
Balanced Accuracy	0.9993	0.9977	0.9958	0.9919	0.9967

Random forest tree model has a accuracy of 0.99 and out of sampleor err 0.01 that is far better than classification tree method. We will use this model or method to predict outcomes of 20 test data sets.

# 3. Prediction with random forest tree using test data set

prediction3<- predict(Fit.Model2, new.testing)
prediction3</pre>