RANDOM FOREST

Individual decision tree algorithms can be prone to problems, such as bias and over-fitting. A more practical approach is to construct multiple decision trees and combine the results into a single output. This approach is termed **ensemble methods**. The most well-known ensemble method is **bagging** (also known as **bootstrap aggregation**, **bagging=bootstrap+aggregation**). This method was introduced in 1996 by Leo Breiman. In this method, data points in the training set are sampled with replacement (producing a **bootstrap sample**), one-third of which, known as the **out-of-bag (OOB) sample**, is set aside for cross-validation, and the remaining two-thirds of the sample is used to build a decision tree. Decision trees are generated for each bootstrap sample independently from others. The results are then aggregated depending on the type of trees used. If regression trees are trained, the average of predicted values are computed. If classification trees are fitted, the majority of the predictions define the final predicted class. The ensemble method reduces variance and, as the result, yields more accurate predictions than individual decision trees.

The random forest algorithm is an extension of the bagging method as it utilizes both bagging and feature randomness to create an uncorrelated forest of decision trees. Feature randomness (also known as feature bagging or the random subspace method) generates a random subset of variables (also called features), which ensures low correlation among decision trees. This is a key difference between decision trees and random forests. While decision trees consider all the possible variable splits, random forests only select a subset of those variables.

Random forest algorithms have three main hyper-parameters that need to be set before training. These are node size, the number of trees, and the number of variables sampled.

Variable Importance

Random forest makes it easy to evaluate the **variable importance** (or the **contribution** of each splitting variable to the model). It is sometimes termed the **feature importance**. There are two commonly used ways to evaluate variable importance that are characteristically different from each other. One method is termed the **loss reduction** or **Gini increase** or **Gini importance** or **impurity reduction** or **mean decrease in impurity (MDI)**. It is used to measure how much the model's accuracy decreases when a given variable is excluded. For a random forest with regression trees, the loss functions are the **mean squared error** MSE = RSS/n, and the **absolute error**

$$\frac{1}{n}\sum_{i=1}^{n}|y_i-\hat{y}_i|$$
. The second feature importance method is the **permutation importance** (or the

mean decrease accuracy (MDA)), which identifies the average decrease in accuracy by randomly permuting the variable values in OOB samples.

Example. Consider the data in the file "housing data.csv". We construct random forest regres-

sion for this data set.

```
In SAS:
```

```
proc import out=housing
datafile="./housing_data.csv" dbms=csv replace;
run;
/*SPLITTING DATA INTO 80% TRAINING AND 20% TESTING*/
proc surveyselect data=housing rate=0.8 seed=502305
out=housing outall method=srs;
run;
/*BUILDING RANDOM FOREST REGRESSION*/
proc hpforest data=housing seed=829743
maxtrees=60 vars_to_try=4 trainfraction=0.7
maxdepth=50;
target median_house_value/level=interval;
input ocean_proximity/level=nominal;
input housing_median_age total_rooms total_bedrooms
population households median_income/level=interval;
partition rolevar=selected(train='1');
save file='C:/Users/000110888/Desktop/random_forest.bin';
run;
```

Loss Reduction Variable Importance						
Variable	Number of Rules	MSE	OOB MSE	Absolute Error	OOB Absolute Error	
median_income	35554	5.2672E9	4.3027E9	36134	23887	
ocean_proximity	336	1.6836E9	1.7055E9	12553	12796	
total_rooms	7375	4.5494E8	-5.897E7	4665.515799	-423.229467	
housing_median_age	7987	5.285E8	-7.79E7	5698.148519	141.283900	
total_bedrooms	8196	3.22E8	-1.709E8	4098.769334	-795.029912	
households	20105	5.2346E8	-2.364E8	6723.747118	-1030.371514	
population	14896	4.4438E8	-2.694E8	5805.772403	-1414.692176	

```
/*COMPUTING PREDICTED VALUES FOR TESTING DATA*/
data test;
set housing;
if(selected='0');
```

```
run;
proc hp4score data=test;
id median_house_value;
score file='C:/Users/000110888/Desktop/random_forest.bin'
out=predicted;
run;
/*DETERMINING 10%, 15%, AND 20% ACCURACY*/
data accuracy;
set predicted;
if(abs(median_house_value-P_median_house_value)
<0.10*median_house_value)
then ind10=1; else ind10=0;
if(abs(median_house_value-P_median_house_value)
<0.15*median_house_value)
then ind15=1; else ind15=0;
if(abs(median_house_value-P_median_house_value)
<0.20*median_house_value)
then ind20=1; else ind20=0;
run;
proc sql;
 select sum(ind10)/count(*) as accuracy10,
sum(ind15)/count(*) as accuracy15,
 sum(ind20)/count(*) as accuracy20
 from accuracy;
 quit;
```

accuracy10	accuracy15	accuracy20
0.352113	0.485915	0.612676

In R:

```
#install.packages("randomForest")
library(randomForest)
```

```
housing.data<- read.csv(file="./housing data.csv", header=TRUE, sep=",")
#SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS
set.seed(902881)
sample <- sample(c(TRUE, FALSE), nrow(housing.data), replace=TRUE, prob=c(0.8,0.2))
train<- housing.data[sample,]
test<- housing.data[!sample,]
#BUILDING RANDOM FOREST REGRESSION
rf.reg < -randomForest(median house value \sim housing median age + total rooms + total bedrooms
+ population + households + median income + ocean proximity, data=train, ntree=150, mtry=5,
maxnodes=30)
#DISPLAYING FEATURE IMPORTANCE
print(importance(rf.reg,type=2))
                    IncNodePurity
housing_median_age
                     4.241691e+11
total_rooms
                     4.108330e+11
total_bedrooms
                     2.028437e+11
population
                     2.153154e+11
households
                     2.400010e+11
median_income
                     1.245747e+13
ocean_proximity
                     2.914203e+12
#COMPUTING PREDICTION ACCURACY FOR TESTING DATA
P median house value<- predict(rf.reg, newdata=test)
#accuracy within 10%
accuracy10<-
ifelse(abs(test$median house value-P median house value)<0.10*test$median house value,1,0)
print(accuracy10<- sum(accuracy10)/length(accuracy10))
0.2857143
#accuracy within 15%
accuracy15<-
ifelse(abs(test$median house value-P median house value)<0.15*test$median house value,1,0)
print(accuracy15<- sum(accuracy15)/length(accuracy15))
0.4303797
```

```
\label{eq:local_sum} $\#$ accuracy within 20\% $$ accuracy 20 < -$ ifelse (abs(test\$median\_house\_value-P\_median\_house\_value) < 0.20*test\$median\_house\_value,1,0) $$ print(accuracy 20 < -$ sum(accuracy 20)/length(accuracy 20))$$ 0.5334539
```

In Python:

```
import pandas
from sklearn.ensemble import RandomForestRegressor
from sklearn.model_selection import train_test_split
housing=pandas.read_csv('C:/Users/000110888/Desktop/housing data.csv')
coding={'<1H OCEAN': 1, 'INLAND': 2, 'NEAR BAY': 3, 'NEAR OCEAN': 4}
housing['ocean_proximity']=housing['ocean_proximity'].map(coding)
X=housing.iloc[:,0:7].values
y=housing.iloc[:,7].values
#SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS
X train, X test, y train, y test=train test split(X, y, test size=0.20,
random_state=348644)
#FITTING RANDOM FOREST REGRESSION TREE
rf_reg=RandomForestRegressor(n_estimators=100, random_state=323445,
max_depth=50, max_features=4)
rf_reg.fit(X_train, y_train)
#DISPLAYING VARIABLE IMPORTANCE
from sklearn.ensemble import ExtraTreesClassifier
var_names=pandas.DataFrame(['housing_median_age','total_rooms','total_bedrooms','population',
'households','median_income','ocean_proximity'], columns=['var_name'])
loss reduction=pandas.DataFrame(rf_reg.feature_importances_, columns=['loss_reduction'])
var_importance=pandas.concat([var_names, loss_reduction], axis=1)
var importance=var importance.sort values("loss reduction", axis=0, ascending=False)
print(var importance)
```

```
var_name loss_reduction
5 median_income 0.592797
6 ocean_proximity 0.154105
1 total_rooms 0.061499
0 housing_median_age 0.053247
```

```
3 population 0.051611
4 households 0.044158
2 total bedrooms 0.042583
```

```
#COMPUTING PREDICTION ACCURACY FOR TESTING DATA
y_pred=rf_reg.predict(X_test)
ind10=[]
ind15=[]
ind20=[]
for sub1, sub2 in zip(y_pred, y_test):
    ind10.append(1) if abs(sub1-sub2)<0.10*sub2 else ind10.append(0)
    ind15.append(1) if abs(sub1-sub2)<0.15*sub2 else ind15.append(0)</pre>
    ind20.append(1) if abs(sub1-sub2)<0.20*sub2 else ind20.append(0)
#accuracy within 10%
accuracy10=sum(ind10)/len(ind10)
print(accuracy10)
#accuracy within 15%
accuracy15=sum(ind15)/len(ind15)
print(accuracy15)
#accuracy within 20%
accuracy20=sum(ind20)/len(ind20)
print(accuracy20)
```

- 0.36203866432337434
- 0.5202108963093146
- 0.648506151142355

Example. Consider the data in the file "pneumonia_data.csv". We construct a random forest binary classifier for this data set.

In SAS:

```
proc import out=pneumonia datafile="./pneumonia_data.csv" dbms=csv replace;
/*SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS*/
proc surveyselect data=pneumonia rate=0.8 seed=6132208
```

```
out=pneumonia outall method=srs;
run;
/*BUILDING RANDOM FOREST BINARY CLASSIFIER*/
proc hpforest data=pneumonia seed=115607
maxtrees=60 vars_to_try=4 trainfraction=0.7
maxdepth=50;
target pneumonia/level=binary;
input gender tobacco_use/level=nominal;
input age PM2_5/level=interval;
partition rolevar=selected(train='1');
save file='C:/Users/000110888/Desktop/random_forest.bin';
run;
/*COMPUTING PREDICTED VALUES FOR TESTING DATA*/
data test;
set pneumonia;
if(selected='0');
run;
proc hp4score data=test;
id pneumonia;
score file='C:/Users/000110888/Desktop/random_forest.bin'
out=predicted;
run;
```

Loss Reduction Variable Importance						
Variable	Number of Rules	Gini	OOB Gini	Margin	OOB Margin	
gender	84	0.056581	0.05607	0.113161	0.112805	
tobacco_use	263	0.044094	0.04157	0.088188	0.086383	
PM2_5	7647	0.271965	0.00917	0.543929	0.279844	
age	3208	0.085176	-0.05001	0.170351	0.035585	

```
/*COMPUTING PREDICTION ACCURACY FOR TESTING DATA*/
data predicted;
set predicted;
match=(pneumonia=lowcase(I_pneumonia));
proc sql;
select sum(match)/count(*) as accuracy
from predicted;
quit;
                                       ассигасу
                                        0.823188
In R:
pneumonia.data<- read.csv(file="./pneumonia data.csv", header=TRUE, sep=",")
#SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS
set.seed(447558)
sample <- sample(c(TRUE, FALSE), nrow(pneumonia.data), replace=TRUE, prob=c(0.8,0.2))
train <- pneumonia.data[sample,]
test<- pneumonia.data[!sample,]
#BUILDING RANDOM FOREST BINARY CLASSIFIER
library(randomForest)
rf.class < -randomForest(as.factor(pneumonia) \sim age + gender + tobacco use + PM2 5, data=train,
ntree=150, mtry=4, maxnodes=30)
#DISPLAYING FEATURE IMPORTANCE
print(importance(rf.class,type=2))
            MeanDecreaseGini
                     36.39492
age
gender
                     67.32295
tobacco_use
                     56.93959
```

151.83923

PM2_5

In Python:

```
import pandas
from sklearn.model_selection import train_test_split
from sklearn.ensemble import RandomForestClassifier
pneumonia_data=pandas.read_csv('C:/Users/000110888/Desktop/pneumonia_data.csv')
code_gender={ 'M':1, 'F':0}
code_tobacco_use={'yes':1,'no':0}
code_pneumonia={'yes':1,'no':0}
pneumonia_data['gender']=pneumonia_data['gender'].map(code_gender)
pneumonia_data['tobacco_use']=pneumonia_data['tobacco_use'].map(code_tobacco_use)
pneumonia_data['pneumonia']=pneumonia_data['pneumonia'].map(code_pneumonia)
X=pneumonia_data.iloc[:,0:4].values
y=pneumonia_data.iloc[:,4].values
#SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS
X_train, X_test, y_train, y_test=train_test_split(X, y, test_size=0.20,
random_state=786756)
#FITTING RANDOM FOREST BINARY CLASSIFIER
rf class=RandomForestClassifier(n estimators=150, criterion='entropy',
random state=778554, max depth=50, max features=4)
rf_class.fit(X_train, y_train)
#DISPLAYING VARIABLE IMPORTANCE
from sklearn.ensemble import ExtraTreesClassifier
var_names=pandas.DataFrame(['gender','age','tobacco_use','PM2_5'], columns=['var_name'])
loss_reduction=pandas.DataFrame(rf_class.feature_importances_, columns=['loss_reduction'])
var_importance=pandas.concat([var_names, loss_reduction], axis=1)
var_importance=var_importance.sort_values("loss_reduction", axis=0, ascending=False)
print(var importance)
```

```
    var_name
    loss_reduction

    3
    PM2_5
    0.581401

    1
    age
    0.233343

    0
    gender
    0.095434

    2
    tobacco_use
    0.089823
```

```
#COMPUTING PREDICTION ACCURACY FOR TESTING DATA
y_pred=rf_class.predict(X_test)
y_test=pandas.DataFrame(y_test,columns=['pneumonia'])
y_pred=pandas.DataFrame(y_pred,columns=['predicted'])
df=pandas.concat([y_test,y_pred],axis=1)

match=[]
for i in range(len(df)):
    if df['pneumonia'][i]==df['predicted'][i]:
        match.append(1)
    else:
        match.append(0)

accuracy=sum(match)/len(match)
print(accuracy)
```

0.8121387283236994

Example. Consider the data in the file "movie_data.csv". We construct random forest multinomial classifier for this data set.

In SAS:

```
proc import out=movie
datafile="./movie_data.csv" dbms=csv replace;
run;

/*SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS*/
proc surveyselect data=movie rate=0.8 seed=550040
out=movie outall method=srs;
```

```
run;

/*BUILDING RANDOM FOREST MULTINOMIAL CLASSIFIER*/
proc hpforest data=movie seed=454545
maxtrees=150 vars_to_try=4 trainfraction=0.7
maxdepth=10;
target rating/level=ordinal;
input age member/level=nominal;
input age nmovies/level=interval;
partition rolevar=selected(train='1');
```

save file='C:/Users/000110888/Desktop/random_forest.bin';

run;

Loss Reduction Variable Importance						
Variable	Number of Rules	Gini	OOB Gini	Margin	OOB Margin	
member	378	0.004665	-0.00439	0.002648	-0.00475	
age	1234	0.049053	-0.02003	0.038625	-0.01943	
nmovies	3134	0.051944	-0.02365	0.059274	-0.00833	

```
/*COMPUTING PREDICTED VALUES FOR TESTING DATA*/
data test;
set movie;
if(selected='0');
run;

proc hp4score data=test;
id rating;
score file='C:/Users/000110888/Desktop/random_forest.bin'
out=predicted;
run;

/*COMPUTING PREDICTION ACCURACY FOR TESTING DATA*/
```

```
data predicted;
set predicted;
match=(rating=lowcase(I_rating));
run;
proc sql;
select sum(match)/count(*) as accuracy
from predicted;
quit;
                                       accuracy
                                       0.271523
In R:
movie.data<- read.csv(file="./movie_data.csv", header=TRUE, sep=",")
#SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS
set.seed(566222)
sample <- sample(c(TRUE, FALSE), nrow(movie.data),replace=TRUE, prob=c(0.8,0.2))
train <- movie.data[sample,]
test<- movie.data[!sample,]
#BUILDING RANDOM FOREST MULTINOMIAL CLASSIFIER
library(randomForest)
rf.class < -randomForest(as.factor(rating) \sim age + gender + member + nmovies, data=train, ntree=150,
mtry=4, maxnodes=30)
#DISPLAYING FEATURE IMPORTANCE
print(importance(rf.class,type=2))
        MeanDecreaseGini
               64.973483
age
               10.542498
gender
member
                5.531971
nmovies
               25.470534
#COMPUTING PREDICTION ACCURACY FOR TESTING DATA
```

predclass<- predict(rf.class, newdata=test)</pre>

```
\begin{split} & \operatorname{accuracy} <\text{-} \operatorname{c()} \\ & \operatorname{accuracy} <\text{-} \operatorname{c()} \\ & \operatorname{n<-} \operatorname{nrow}(\operatorname{test}) \\ & \operatorname{for} \ (i \ in \ 1:n) \\ & \operatorname{accuracy}[i] <\text{-} \ ifelse(\operatorname{test\$rating}[i] == \operatorname{test\$predclass}[i],1,0) \\ & \operatorname{print}(\operatorname{accuracy} <\text{-} \operatorname{sum}(\operatorname{accuracy}) / \operatorname{length}(\operatorname{accuracy})) \\ & 0.3212435 \end{split}
```

In Python:

```
import pandas
from sklearn.model_selection import train_test_split
from sklearn.ensemble import RandomForestClassifier
movie data=pandas.read csv('C:/Users/000110888/Desktop/movie data.csv')
code_gender={ 'M':1, 'F':0}
code member={'yes':1,'no':0}
code_rating={'very bad':1,'bad':2,'okay':3,'good':4,'very good':5}
movie_data['gender']=movie_data['gender'].map(code_gender)
movie data['member']=movie data['member'].map(code member)
movie_data['rating']=movie_data['rating'].map(code_rating)
X=movie data.iloc[:,0:4].values
y=movie_data.iloc[:,4].values
#SPLITTING DATA INTO 80% TRAINING AND 20% TESTING SETS
X_train, X_test, y_train, y_test=train_test_split(X, y, test_size=0.20, random_state=599555)
#FITTING RANDOM FOREST FOR MULTINOMIAL CLASSIFIER
rf_class=RandomForestClassifier(n_estimators=150, random_state=663474, max_depth=50, max_features=4)
rf_class.fit(X_train, y_train)
#DISPLAYING VARIABLE IMPORTANCE
from sklearn.ensemble import ExtraTreesClassifier
var_names=pandas.DataFrame(['age','gender','member','nmovies'], columns=['var_name'])
loss_reduction=pandas.DataFrame(rf_class.feature_importances_, columns=['loss_reduction'])
var_importance=pandas.concat([var_names, loss_reduction], axis=1)
var importance=var importance.sort values("loss reduction", axis=0, ascending=False)
print(var importance)
```

```
      var_name
      loss_reduction

      0
      age
      0.572954

      3
      nmovies
      0.269890

      2
      member
      0.084660

      1
      gender
      0.072495
```

```
#COMPUTING PREDICTION ACCURACY FOR TESTING DATA
y_pred=rf_class.predict(X_test)
y_test=pandas.DataFrame(y_test,columns=['rating'])
y_pred=pandas.DataFrame(y_pred,columns=['predicted'])
df=pandas.concat([y_test,y_pred],axis=1)

match=[]
for i in range(len(df)):
    if df['rating'][i]==df['predicted'][i]:
        match.append(1)
    else:
        match.append(0)

accuracy=sum(match)/len(match)
print(accuracy)
```

0.3026315789473684