Loan Prediction

The aim of this project is to predict real-estate prices using the machine learning algorithms: Logistic Regression, Decision tree, Random Forest. The three of them will show different results for the accuracy.

Imports

```
import pandas as pd
import requests
import matplotlib.pyplot as plt
import seaborn as sns
from google.colab import files
from datetime import datetime
import io
import mpl_toolkits
import numpy as np
%matplotlib inline

# Load the data
local_file = files.upload()
train_data = io.BytesIO(local_file['results.csv'])
df = pd.read_csv(train_data)
```

Choose Files | results.csv
 results.csv(application/vnd.ms-excel) - 104433 bytes, last modern

• **results.csv**(application/vnd.ms-excel) - 104433 bytes, last modified: 4/13/2021 - 100% done Saving results.csv to results.csv

Preparing the data for training the models

Encoding to numeric data in order to start the training of the models.

```
#drop the uniques loan id
df.drop('Loan_ID', axis = 1, inplace = True)
df.drop('Unnamed: 0', axis = 1, inplace = True)
```

df.info()

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 981 entries, 0 to 980
Data columns (total 15 columns):
     Column
                        Non-Null Count Dtype
---
    -----
 0
     Gender
                        981 non-null
                                        float64
 1
     Married
                        981 non-null
                                        float64
 2
     Dependents
                        981 non-null
                                        float64
 3
     Education
                        981 non-null
                                        int64
 4
     Self Employed
                        981 non-null
                                        float64
 5
     ApplicantIncome
                        981 non-null
                                        int64
 6
     CoapplicantIncome 981 non-null
                                        float64
 7
     LoanAmount
                        981 non-null
                                        float64
 8
     Loan Amount Term
                        981 non-null
                                        float64
 9
     Credit_History
                        981 non-null
                                        float64
 10
    Property_Area
                        981 non-null
                                        int64
 11 Loan Status
                        981 non-null
                                        float64
 12 LoanAmount log
                        981 non-null
                                        float64
 13
    TotalIncome
                        981 non-null
                                        float64
 14 TotalIncome log
                        981 non-null
                                        float64
dtypes: float64(12), int64(3)
memory usage: 115.1 KB
```

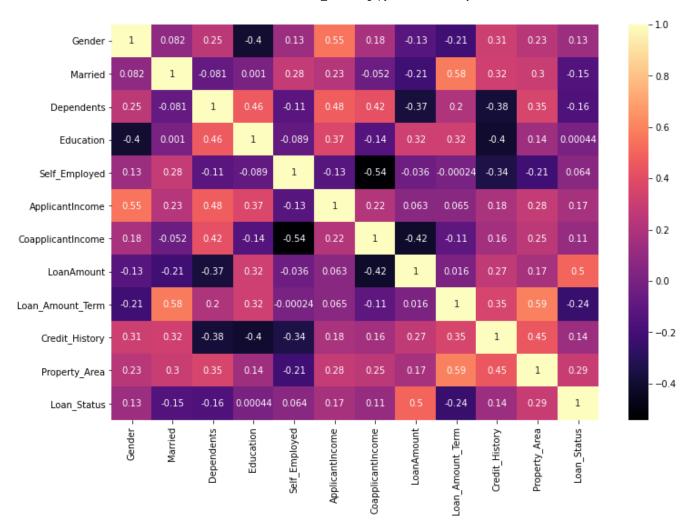
Train-Test Split dataset

Heatmaps are very useful to find relations between two variables in a dataset and this way the user gets a visualisation of the numeric data. No correlations are extremely high. Each square shows the correlation between the variables on each axis.

The correlations between the feautures can be explained:

The close to 1 the correlation is the more positively correlated they are; that is as one increases so does the other and the closer to 1 the stronger this relationship is. It is noticable that the correlation between the ApplicantIncome and LoanAmount is 0.57, which mean that they have a positive correlation, but not strong.

```
from pandas import DataFrame
%matplotlib inline
plt.figure(figsize=(12, 8))
df_temp = df.copy()
Index= ['Gender', 'Married', 'Dependents', 'Education', 'Self_Employed', 'ApplicantIncome'
Cols = ['Gender', 'Married', 'Dependents', 'Education', 'Self_Employed', 'ApplicantIncome'
df_temp = DataFrame(abs(np.random.randn(12, 12)), index=Index, columns=Cols)
sns.heatmap(df_temp.corr(), annot=True, cmap = 'magma')
plt.show()
```



Importing sklearn libraries

```
from sklearn.model_selection import train_test_split
from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import RandomForestClassifier
from sklearn.linear_model import LogisticRegression
from sklearn.metrics import f1_score
```

Splitting into train and test set after choosing the right features X and labels y

```
y = df['Loan_Status']
X = df.drop('Loan_Status', axis = 1)
```

To split the dataset, I will use random sampling with 80/20 train-test split; that is, 80% of the dataset will be used for training and set aside 20% for testing:

Analyzing the numeric features.

4/15/2021

Modeling:

Three models will be built and evaluated by their performances with R-squared metric. Additionally, insights on the features that are strong predictors of house prices, will be analised.

Logistic Regression:

- 1. Creating
- 2. Fitting with train data

Do predictions on a test set. **Testing** the model by testing the test data.

```
y_reg=model.predict(X_test)
```

The F1 score can be interpreted as a weighted average of the precision and recall, where an F1 score reaches its best value at 1 and worst score at 0.

```
evaluation = f1_score(y_test, y_reg)
evaluation
0.8895522388059702
```

Reporting the coefficient value for each feature. Notice that the coefficients are both positive and negative. The positive scores indicate a feature that predicts class 1, whereas the negative scores indicate a feature that predicts class 0.

The importance of a feature is measured by calculating the increase in the model's prediction error after permuting the feature. A feature is "important" if shuffling its values increases the model error, because in this case the model relied on the feature for the prediction.

```
# get importance
importance = model.coef_[0]
# summarize feature importance
for i,v in enumerate(importance):
   print('Feature: %0d, Score: %.5f' % (i,v))
# plot feature importance
plt.bar([x for x in range(len(importance))], importance)
plt.show()
```

```
Feature: 0, Score: -0.04457
Feature: 1, Score: 0.00662
Feature: 2, Score: -0.06491
Feature: 3, Score: -0.00553
Feature: 4, Score: -0.04740
Feature: 5, Score: -0.00002
Feature: 6, Score: 0.00003
Feature: 7, Score: 0.00319
Feature: 8, Score: 0.00202
Feature: 9, Score: -0.12728
Feature: 10, Score: -0.09592
Feature: 11, Score: -0.08757
Feature: 12, Score: 0.00001
Feature: 13, Score: -0.18096
```

This might mean that your model is underfit (not enough iteration and it has not used the feature enough) or that the feature is not good and you can try removing it to improve final quality.

```
-V.123 ]
```

Decision tree:

- 1. Creating classifier
- 2. Fitting classifier with train data

Do predictions on a test set. **Testing** the model by testing the test data.

```
y tree=dtree.predict(X test)
```

The F1 score can be interpreted as a weighted average of the precision and recall, where an F1 score reaches its best value at 1 and worst score at 0.

```
evaluation = f1_score(y_test, y_tree)
evaluation
```

0.872611464968153

Evaluate classifier measures accuracy by using F1 score. The result shows that the model is precise.

Random forests

```
# Feature Scaling
from sklearn.preprocessing import StandardScaler
from sklearn.ensemble import RandomForestRegressor
from sklearn.model_selection import cross_val_score, train_test_split, GridSearchCV
sc = StandardScaler()
X train = sc.fit transform(X train)
X test = sc.transform(X test)
regressor = RandomForestRegressor(n estimators=20, random state=0)
regressor.fit(X_train, y_train)
y pred = regressor.predict(X test)
from sklearn import metrics
print('Mean Absolute Error:', metrics.mean_absolute_error(y_test, y_pred))
print('Mean Squared Error:', metrics.mean squared error(y test, y pred))
print('Root Mean Squared Error:', np.sqrt(metrics.mean_squared_error(y_test, y_pred)))
     Mean Absolute Error: 0.24162436548223348
     Mean Squared Error: 0.12743654822335024
     Root Mean Squared Error: 0.3569825601109251
# Import the model we are using
from sklearn.ensemble import RandomForestRegressor
# Instantiate model with 1000 decision trees
rf = RandomForestRegressor(n estimators = 1000, random state = 42)
# Train the model on training data
rf.fit(X train, y train)
     RandomForestRegressor(bootstrap=True, ccp alpha=0.0, criterion='mse',
                           max depth=None, max features='auto', max leaf nodes=None,
                           max samples=None, min impurity decrease=0.0,
                           min_impurity_split=None, min_samples_leaf=1,
                           min samples split=2, min weight fraction leaf=0.0,
                           n_estimators=1000, n_jobs=None, oob_score=False,
                           random state=42, verbose=0, warm start=False)
```

```
# Calculate the absolute errors
errors = abs(predictions - y test)
# Print out the mean absolute error (mae)
print('Mean Absolute Error:', round(np.mean(errors), 2), 'degrees.')
    Mean Absolute Error: 0.24 degrees.
# Calculate mean absolute percentage error (MAPE)
mape = 100 * (errors /y_test)
# Calculate and display accuracy
accuracy = 100 - np.mean(mape)
print('Accuracy:', round(accuracy, 2), '%.')
    Accuracy: 81.92 %.
#Random forest determined feature importances
rf.feature importances
     array([0.01863798, 0.02443568, 0. , 0.02182857, 0.01956634,
            0.16946051, 0.10084354, 0.1094957, 0.04036644, 0.12931742,
            0.03932196, 0.11745444, 0.1039604, 0.10531102])
forest = RandomForestClassifier()
forest.fit(X train, y train)
     RandomForestClassifier(bootstrap=True, ccp_alpha=0.0, class_weight=None,
                            criterion='gini', max_depth=None, max_features='auto',
                            max leaf nodes=None, max samples=None,
                            min impurity decrease=0.0, min impurity split=None,
                            min samples leaf=1, min samples split=2,
                            min_weight_fraction_leaf=0.0, n_estimators=100,
                            n jobs=None, oob score=False, random state=None,
                            verbose=0, warm start=False)
```

Testing the model by testing the test data.

```
y_forest=forest.predict(X_test)
```

The F1 score can be interpreted as a weighted average of the precision and recall, where an F1 score reaches its best value at 1 and worst score at 0.

Result:

```
evaluation_f= f1_score(y_test, y_forest)
evaluation f
```

0.8982035928143713

After using the F1, it is determined that the model is precised to be used in the deployment.

```
importance = rf.feature_importances_

# map feature importance values to the features
feature_importances = zip(importance, X.columns)

#list(feature_importances)
sorted_feature_importances = sorted(feature_importances, reverse = True)

#print(sorted_feature_importances)
top_15_predictors = sorted_feature_importances[0:15]
values = [value for value, predictors in top_15_predictors]
predictors = [predictors for value, predictors in top_15_predictors]
print(predictors)

values_importances_value, reverse_value_value, reverse_value_value, reverse_value, reverse_value_value, reverse_value, reverse_v
```

Saving the model that I am going to use in the deployment part of the project.

```
# Saving the model
import pickle

filename = 'classifier.pkl'
pickle.dump(forest, open(filename, 'wb'))
```

Conclusion

From the Modelling, it can be concluded:

- 1. The Logistic regression showed accuracy of 88%.
- 2. The Decision tree showed accuracy of 87%.
- 3. Rando forest showed accuracy of 89%.

From the evaluation of the three models, it can be noticed that Random forest performed better than the others.

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