### **HW 6**

This assignment covers Comparision of Decision Trees and Support Vector Machine. **DO NOT ERASE MARKDOWN CELLS AND INSTRUCTIONS IN YOUR HW submission** 

- Q QUESTION
- A Where to input your answer

#### Instructions

Keep the following in mind for all notebooks you develop:

- Structure your notebook.
- Use headings with meaningful levels in Markdown cells, and explain the questions each piece of code is to answer or the reason it is there.
- Make sure your notebook can always be rerun from top to bottom.
- Objective of this assignment is to help you master python and scikit-learn package.
- See <u>README.md</u> (<u>README.md</u>) for homework submission instructions

#### **Related Tutorials**

- <u>Decision Tree with KFold Cross Validation (https://scikit-learn.org/stable/modules/generated/sklearn.model\_selection.cross\_val\_score.html)</u>
- <u>Decision Tree with Bagging (https://scikit-learn.org/stable/modules/generated</u> /sklearn.ensemble.BaggingRegressor.html#sklearn.ensemble.BaggingRegressor)
- <u>Support Vector Machine (https://towardsdatascience.com/support-vector-machine-introduction-to-machine-learning-algorithms-934a444fca47)</u>

# **Data Processing**

Q1 Get training data from the dataframe

- 1. Load ../data/HW6\_data.csv into a data frame
- 2. Check if there is any NaN in the dataset
- 3. Clean the dataset
- 4. Assign values of target class column to y, note you have to use .values method
- 5. Drop target class column from data frame,
- 6. Assign df values to x
- 7. Split dataset into train and test data use train\_test\_split with test\_size = 0.25, stratify y and random state = 1238

A1 Replace ??? with code in the code cell below

```
In [96]: | import numpy as np
            import pandas as pd
            from sklearn.model selection import train test split
            from sklearn.linear model import LogisticRegression
             #Read the data file using the prropriate separator as input to read cs
            df = pd.read csv('HW6 data.csv')
            df.head(10)
             # check if there is NaN in the dataset
            df.isnull()
             #Drop NaNs if there is any
            df.dropna(inplace = True)
             # Count number of entries for different class target
   Out[96]: 9273
In [97]: ▶ # Assign values of ```target class``` column to y, note you have to us
            y = df.target class.values
            # Drop 'target class' column from data frame,
            df.drop(columns = ['target class'], inplace = True)
            # Assign df values to x
            x = df.values
             # View shape of x and y
            x.shape
            y.shape
            xtrain, xtest, ytrain, ytest = train test split(x,y), test size = 0.25,
```

## **Decision Tree**

## **Decision Tree with different depth**

Q2 Train DecisionTreeClassifier Model at different depths

- 1. Create four <u>DecisionTreeClassifier (https://scikit-learn.org/stable/modules/generated</u>/sklearn.tree.DecisionTreeClassifier.html) models with different parameters
- 2. Create four decision tree classifier models, use max\_depth size = 1, 2, 5, 25 & max\_leaf\_nodes=5, 10, 15, 25 respectively
- 3. Use random state=30 & criterion='entropy' for all models
- 4. Fit the model with the train data. You will have 4 different models for different depth sizes.
- 5. Predict the values with the test data
- 6. Calculate the mean squared error of each model's prediction
- 7. Print precision recall curve for the model with the best MSE value

A2 Replace ??? with code in the code cell below

```
In [98]: | from sklearn.tree import DecisionTreeClassifier
```

```
from sklearn.metrics import mean squared error
from IPython.display import display
#create decision tree classifier
clf 1 = DecisionTreeClassifier(max depth = 1, max leaf nodes = 5, rand
clf 2 = DecisionTreeClassifier(max_depth = 2, max_leaf_nodes = 10, ran
clf 3 = DecisionTreeClassifier(max depth = 5, max leaf nodes = 15, ran
clf 4 = DecisionTreeClassifier(max depth = 25, max leaf nodes = 25, ra
#fit classifier model
clf 1.fit(xtrain, ytrain)
clf 2.fit(xtrain, ytrain)
clf 3.fit(xtrain, ytrain)
clf 4.fit(xtrain, ytrain)
#predict
y pred clf1 = clf 1.predict(xtest)
y pred clf2 = clf 2.predict(xtest)
y pred clf3 = clf 3.predict(xtest)
y pred clf4 = clf 4.predict(xtest)
#calculate mean squared error
mse clf1 = mean squared error(ytest, y pred clf1)
mse clf2 = mean squared error(ytest, y pred clf2)
mse_clf3 = mean_squared_error(ytest, y_pred_clf3)
mse clf4 = mean squared error(ytest, y pred clf4)
#printing mse values
display(mse clf1)
display(mse clf2)
display(mse clf3)
display (mse clf4)
0.0258732212160414
0.0258732212160414
0.02501078050884002
0.0258732212160414
```

#### **Precision-Recall Curve for Best Above**

Important Note: If from\_estimator() function gives Attribute error then it means your sklearn is not updated.

· If you are using conda, you can upgrade with

conda upgrade -c conda-forge scikit-learn

· or, with pip,

python -m pip install scikit-learn --upgrade

```
In [99]:
           from sklearn.metrics import precision recall curve
               from sklearn.metrics import PrecisionRecallDisplay
               import matplotlib.pyplot as plt
               precision recall curve (ytest, y pred clf3)
               PrecisionRecallDisplay.from estimator(clf 3, xtest, ytest)
                  1.0
                Precision (Positive label: 1.0)
                  0.8
                  0.6
                  0.4
                  0.2
                           DecisionTreeClassifier (AP = 0.87)
                       0.0
                                0.2
                                                          0.8
                                                 0.6
                                                                   1.0
```

#### **Decision Tree with K-fold cross validation**

Q3 Use Kfold on the testing dataset, and evaluate the best model

Recall (Positive label: 1.0)

- 1. Finally pick the best max\_depth you got
- 2. Use this max\_depth & max\_leaf\_node. Also, use cross\_val\_score and fit your model with k = 5 fold size
- 3. Calculate average scores in kfold

A3 Replace ??? with code in the code cell below

```
In [100]: | from sklearn.model_selection import KFold, cross_val_score

scores = cross_val_score(clf_3, xtrain, ytrain, cv = KFold(n_splits = print("Cross-validation scores: {}".format(scores))

Cross-validation scores: [0.9798706 0.97843278 0.97699497 0.98634076 0.97482014]

Average cross-validation score: 0.98
```

## **Decision Tree with Bagging**

Q4 Now we will use Bagging technique on the our previous best model, and evaluate it

Part 1:

1. Now, Create a Bagged Model passing model = previous\_best, n\_estimators =

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```
10 & random state=1 to BaggingClassifier()
```

- 2. Fit the model with the train data
- 3. Predict the values with the test data
- Calculate the test MSE
- 5. Plot Precision-Recall Curve from the true & predicted test data

#### A4 Replace ??? with code in the code cell below

```
In [101]:
            from sklearn.ensemble import BaggingClassifier
               # Use BaggingRegressor to fit the training data
               # Calculate the mean squared error
               #load BaggingRegressor model and pass n estimators=10, random state=1
               bagged clf = BaggingClassifier(n estimators = 10, random state = 1)
               bagged clf.fit(xtrain, ytrain)
               pred = bagged clf.predict(xtest)
In [102]:
               #pass necessary parameters to PrecisionRecallDisplay.from estimator()
               PrecisionRecallDisplay.from estimator(bagged clf, xtest, ytest)
                  1.0
                Precision (Positive label: 1.0)
                  0.8
                  0.6
                  0.4
                  0.2
                           BaggingClassifier (AP = 0.84)
                               0.2
                                               0.6
                                       0.4
                                                        0.8
                                                                1.0
                      0.0
                                    Recall (Positive label: 1.0)
```

Part 2:

1. Why BaggingClassifier is called an ensembled technique? why it works better most of the time than the single model classifiers?

The BaggingClassifier is called an ensembled technique due to the nature of using multiple random data points in different models and each getting there own value which is then average or voted on. This a better technique as it creates more accurate predictions due to having multiple models evaluated. The concept is of ensembled technique is that of group thinking and how it's better to have multiple people looking at the data then just one expert.

What is the disadvantage of incresing the number of estimators while using BaggingClassifier? Explain with an appropriate example.

# Increasing the number of estimators provides better accuracy but unfortantly takes much longer to compute the higher you go

# Support Vector Machine(SVM)

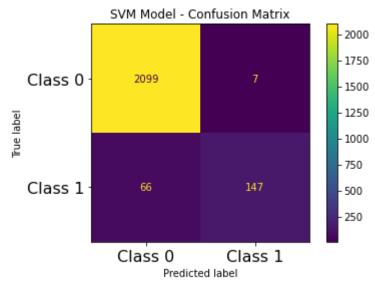
Q5 Create SVM Model on the training set, and evaluate

Part:1

- 1. Now, Create a SVM Model with default parameters
- 2. Fit the model with the train data
- 3. Predict the values with the test data
- 4. Calculate the test MSE
- 5. Plot confusion matrix on the test data

A5 Replace ??? with code in the code cell below

Model accuracy score with default hyperparameters: 0.9685



#### Part2:

1. From the above Confusion Matrix we can see that high number of Class 1 is predicted as Class 0 from the model. What is your reasoning behind this situation?

One of the possible reasons behind this issues is due the training data. If our training data accuracy is low it can give us a lot of erros when making the predictions.

2. What can be done in order to resolve this issue?

If we look into the cleaning up the training data we can get better results.

## **SVM** with high margin

**Q6** Create SVM Model on the training set, and evaluate

Note:

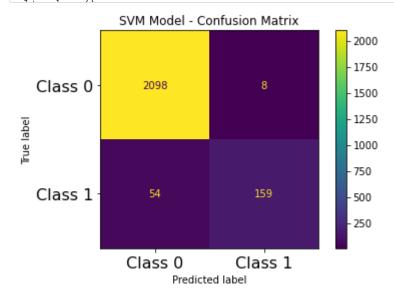
- 1. If we analyze our dataset using df.describe() function, we will see that there are many outliers in the dataset.
- 2. So, we need to increase our margin with <code>HIGH C</code> values so that the SVM model get better generalization

Task:

- 1. Now, Create a SVM Model with rbf kernel and C=100
- 2. Fit the model with the train data
- 3. Predict the values with the test data
- 4. Calculate the test MSE
- 5. Plot Confusion Matrix from the true & predicted test data

#### A6 Replace ??? with code in the code cell below

Model accuracy score with rbf kernel and C=100.0: 0.9733



#### **SVM** with linear kernel

Q7 Create SVM Model on the training set, and evaluate

Task:

- 1. Now, Create a SVM Model with linear kernel and C=1.0
- 2. Fit the model with the train data
- 3. Predict the values with the test data
- 4. Calculate the test MSE
- 5. Plot Confusion Matrix from the true & predicted test data

A7 Replace ??? with code in the code cell below

```
In [107]: | # instantiate classifier with linear kernel and C=1.0
linear_svc= SVC(C = 1.0, kernel = 'linear')

# fit classifier to training set
linear_svc.fit(xtrain, ytrain)

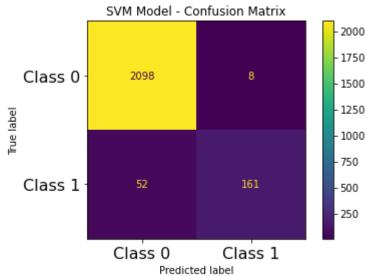
# make predictions on test set
pred_linear_svc = linear_svc.predict(xtest)

# calculate MSE

mse_linear_svc = mean_squared_error(ytest, pred_linear_svc)

# compute and print accuracy score
print('Model accuracy score with linear kernel and C=1.0 : {0:0.4f}'.
```

Model accuracy score with linear kernel and C=1.0: 0.9741

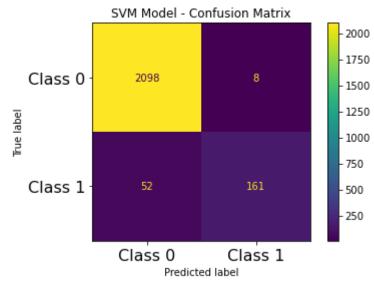


Q8 Create a Grid Search for finetuning the value of C in SVM Model on the training set,

Task:

- 1. Now, Create a SVM Model with linear kernel and evaluate the model for different values of C. Use 'C': [0.01, 0.1, 5, 10, 100]
- 2. Use the sklearn GridSearchCV method for finetuning the linear SVM.
- 3. Use 3 as the number of Cross Validation
- 4. Use accuracy as the scoring technique
- 5. Use clf.cv\_results\_ & clf.best\_params\_ for getting the results from the Cross Validation run.
- 6. Now, Plot the Confusion Matrix for test data, using the best value of C we found from our finetune.

A8 Replace ??? with code in the code cell below



We can see that after using the Best Value of  $\, \, \mathbb{C} \,$ , we have less amount of false positive in our test data prediction.