

## **ASSIGNMENT - 1**

Importing libraries

Preparing the dataset

- Importing the data\_1 and data\_2
- Checking dimensions of the dataset
- Preview dataset
- Check missing values in variables
- Feature vector and target variable declaration (X and y)
- Check target variable distribution - unbalanced distribution

### **(A) Q1**

**Step1:** Splitting X and y into training and test set with a split ratio of 80-20.

For maintaining equal distribution use stratify=y.

**Step2:** Instantiate the decision tree classifier model `tr_model` with the random state as 0 that will control the shuffling applied before each split.

All the default parameters are taken:

```
criterion='gini', splitter='best', max_depth=None, min_samples_split=2,  
min_samples_leaf=1, min_weight_fraction_leaf=0.0, max_features=None,  
random_state=None, max_leaf_nodes=None, min_impurity_decrease=0.0,  
class_weight=None, ccp_alpha=0.0
```

**Step3:** Fit the model and predict test set results

**Step4:** Precision, recall, accuracy, AUC-ROC curve are reported.

**Step5:** Decision tree DT visualization

Save the Decision tree DT as DT\_A\_1

### **(A) Q2**

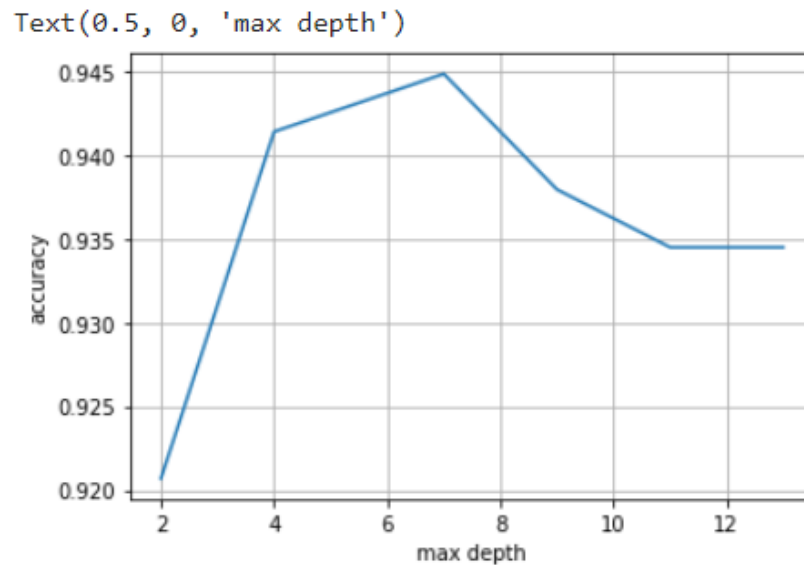
**Step1:** Training the decision tree classifier on 6 different depths

**Step2:** Calculate the accuracy for each depth (ta

- max\_depth=2 accuracy= 0.9206896551724137
- max\_depth=4 accuracy= 0.9413793103448276
- max\_depth=7 accuracy= 0.9448275862068966
- max\_depth=9 accuracy= 0.9379310344827586
- max\_depth=11 accuracy= 0.9344827586206896
- max\_depth=13 accuracy= 0.9344827586206896

**Step3:** Plot the accuracy versus depth graph taking accuracy on the y-axis and max depth in X-axis

From the obtained graph we can observe that we are getting the maximum accuracy for max depth as 7.



### (A) Q3

**Experiment 1:** If we **vary Criterion as entropy** taking other hyperparameters as default, we observe a positive performance score in the precision of classes, recall, and AUC\_ROC score. Though the Gini criterion provides the best features and is faster to compute but here since the class1 labels are more and we aim for the more balanced tree so selecting entropy is a better choice.

precision	recall	f1-score	support		
	1	0.96	0.97	0.97	240
	2	0.79	0.78	0.78	40
	3	0.89	0.80	0.84	10
accuracy				0.94	290
macro avg	0.88	0.85	0.86		290
weighted avg	0.94	0.94	0.94		290

**Experiment 2:** varying **splitter as random** and taking other hyperparameters as default. It is used for choosing the technique to split each node. we observed that it reduces all the accuracy, precision, and recall among others such as random. So we picked **splitter to be best**, which was the default one.

precision	recall	f1-score	support		
	1	0.97	0.96	0.97	240
	2	0.74	0.85	0.79	40
	3	0.71	0.50	0.59	10
accuracy				0.93	290
macro avg	0.81	0.77	0.78		290
weighted avg	0.93	0.93	0.93		290

**Experiment 3:**varying **min sample split as 2** and taking other hyperparameters as default, min sample split is for deciding a minimum number of samples required to split the node. accuracy, precision, recall, and auc\_roc score all increased. So-min sample split as 2 is taken as the best choice.

precision	recall	f1-score	support		
	1	0.96	0.97	0.97	240
	2	0.82	0.78	0.79	40
	3	1.00	0.90	0.95	10
accuracy				0.94	290
macro avg	0.93	0.88	0.90		290
weighted avg	0.94	0.94	0.94		290

**Experiment 4:**varying **max\_depth as 7** and taking other hyperparameters as default. Max depth indicates the maximum depth that the tree is allowed to grow after which we take the majority class label and make it as a leaf(if impure). We observed that among various depths, 7 is giving good results. So it is being chosen.

precision	recall	f1-score	support		
	1	0.97	0.97	0.97	240
	2	0.80	0.80	0.80	40
	3	0.90	0.90	0.90	10
accuracy				0.94	290
macro avg		0.89	0.89	0.89	290
weighted avg		0.94	0.94	0.94	290

**Experiment 5:** varying **min\_samples\_leaf as 2** and taking other hyperparameters as default. The min\_samples\_leaf indicates the minimum number of samples to be needed at the leaf node. By experimenting, we found that min\_samples\_leaf as 2 gives the best results. So it is being chosen.

precision	recall	f1-score	support		
	1	0.96	0.97	0.96	240
	2	0.78	0.78	0.78	40
	3	1.00	0.70	0.82	10
accuracy				0.93	290
macro avg		0.91	0.82	0.85	290
weighted avg		0.93	0.93	0.93	290

**Experiment 6:** varying **max features as sqrt** and taking other hyperparameters as default. It indicates the number of features needed to look at while determining the best split. We have taken the **default that is None** as all the accuracy, precision, recall and auc\_roc score decreased.

precision	recall	f1-score	support		
	1	0.96	0.97	0.97	240
	2	0.83	0.75	0.79	40
	3	0.60	0.60	0.60	10
accuracy				0.93	290
macro avg		0.80	0.78	0.79	290
weighted avg		0.93	0.93	0.93	290

**Experiment 7: varying class weight as balanced** and taking other hyperparameters as default. It is the weight that is associated with the class. **Default** performs better in terms of all scores in comparison to balanced.

precision	recall	f1-score	support	
1	0.97	0.98	0.97	240
2	0.84	0.80	0.82	40
3	0.89	0.80	0.84	10
accuracy			0.95	290
macro avg	0.90	0.86	0.88	290
weighted avg	0.95	0.95	0.95	290

**Experiment 8: varying max-leaf nodes to 20** and taking other hyperparameters as default. Max leaf nodes put the limit on the maximum number of leaves that can be in the decision tree. By experimenting with different values, I found out that 20 gives good results.

precision	recall	f1-score	support	
1	0.97	0.97	0.97	240
2	0.80	0.82	0.81	40
3	1.00	0.80	0.89	10
accuracy			0.95	290
macro avg	0.93	0.87	0.89	290
weighted avg	0.95	0.95	0.95	290

After all the experiments the **best parameters** derived are:

criterion='entropy', splitter='best', max\_depth=7, min\_samples\_split=3, min\_samples\_leaf=5, max\_leaf\_nodes=20, random\_state=0

	precision	recall	f1-score	support
1	0.96	0.99	0.97	240
2	0.86	0.78	0.82	40
3	1.00	0.70	0.82	10
accuracy			0.95	290
macro avg	0.94	0.82	0.87	290
weighted avg	0.95	0.95	0.95	290

## (B) Q1

**Step1:** Take DT\_B\_1 from (A) for the computation

**Step2:** After fitting the model, check the tree structure using id of left and right children  
Total node count=39

**Step3:** Select a random node that needs to be deleted, in our experiment we took node 10. Now, making both the children of node 10 as -1.

We saw the optimization of some performance measures like roc score but the overall accuracy is not changed.

	precision	recall	f1-score	support
1	0.96	0.99	0.97	240
2	0.86	0.78	0.82	40
3	1.00	0.70	0.82	10
accuracy			0.95	290
macro avg	0.94	0.82	0.87	290
weighted avg	0.95	0.95	0.95	290

The optimization of accuracy could be seen if we remove nodes starting from leaf nodes.

These high values of accuracy might be the result of overfitting that can be removed using pruning in further questions.

## (B) Q2

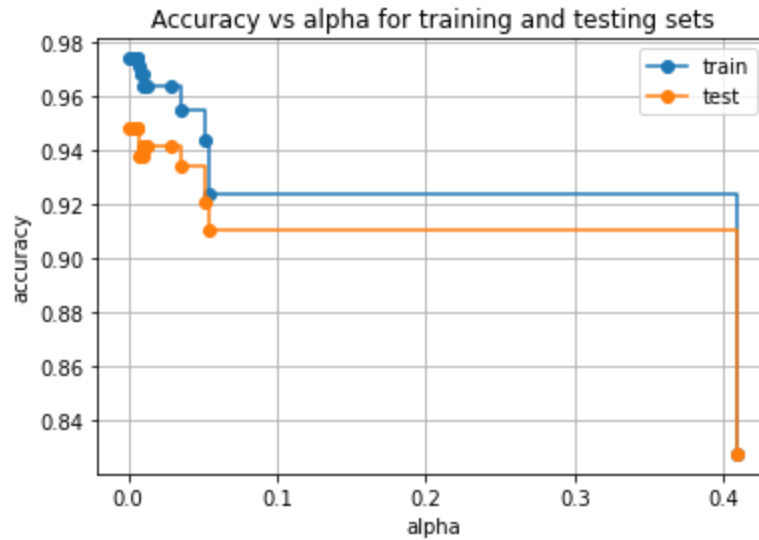
### Cost Complexity pruning technique(Post pruning)

Cost complexity pruning helps to control the size of a tree. Cost complexity pruning technique is parameterized by the cost complexity parameter, `ccp_alpha`. Greater values of `ccp_alpha`, more the number of nodes that are pruned.

We use the DT-A that we got from the end of question (A) and generate all `ccp_alpha` values based on `x_train` and `y_train`.

Then for each `ccp_alpha`, we fit the model and keep on appending them.

Next we plot the accuracy vs `ccp_alpha` graph where we can observe the training and testing accuracy with increasing `ccp_alpha` values.

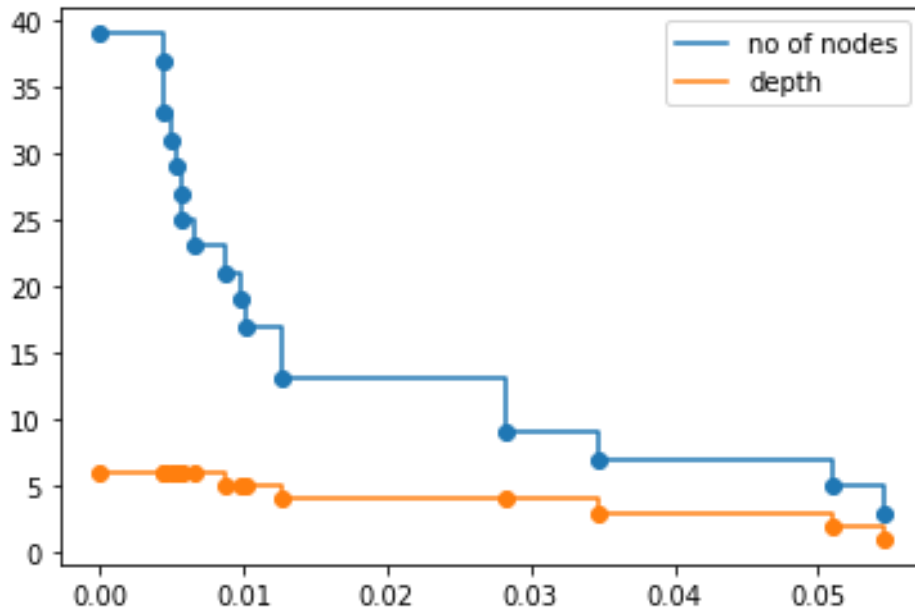


In the above graph, we can observe that at `ccp_alpha` value of 0.025, we get maximum accuracy for both training data and testing data simultaneously. So, we choose `ccp_alpha` value as 0.025.

Then we fit the tree by giving the `ccp_alpha` hyperparameter value as 0.025. The scores after this post pruning technique are observed as follows.

Accuracy score: 0.9413793103448276  
Precision score: 0.9413793103448276  
Recall score: 0.9413793103448276  
AUC\_ROC score: 0.9543027777777778

With increase in value of `alpha`, the number of nodes and depth is decreasing. This can be visualized from the following scatter plot.



### **Other pruning technique - Pre pruning**

We can use pre pruning before the construction of the decision tree using hyper parameter tuning that will help us to overcome the overfitting issue that we observed in the previous question.

Again taking the same `DT_A = tr_model`

For selecting the best parameters for hyperparameters tuning we have used `GrisSearchCV` that results in the best possible parameters in the given range to feed into the classifier.

```
{'criterion': 'entropy', 'max_depth': 7, 'min_samples_leaf': 4,
'min_samples_split': 13, 'splitter': 'best'}
```

Now, using this result in the pre pruning operation.

Test results after pre pruning:

```
Accuracy score: 0.9448275862068966
Precision score: 0.9448275862068966
Recall score: 0.9448275862068966
AUC_ROC score: 0.94145
```



	precision	recall	f1-score	support
1	0.96	0.98	0.97	240
2	0.82	0.80	0.81	40
3	1.00	0.70	0.82	10
accuracy			0.94	290
macro avg	0.93	0.83	0.87	290
weighted avg	0.94	0.94	0.94	290

Observations:

Though pre pruning computes faster and helps in trade-off but here the overall accuracy decreased.

The decision tree that we got is definitely more interpretable as compared to part A.

Also, using pruning we are able to limit overfitting as best parameters are selected so the pruned tree fits the data well.

## (B) Q3

## (C) Q1

Method1:

Step1: Split both the datasets 80-20 ratio.

Step2 Train both the models

Step3: Compare the performances

Method2:

Use VFDT using Hoeffding Tree

Define the dataset complete and another for streaming

Stream the data one after another

Plot the performances after each sample added to the dataset.

So, the performance is increasing with each data sample added to the sample. If more streaming data is added, we will get better performance.

## (C) Q2

### **Decision surface boundaries**

Here we train the decision tree trained on pairs of features. For each possible pair, the decision tree learns the decision boundaries. We have plotted surface boundaries by taking two features at a time.

**Learnings:** We have learnt the following.

- How to fit the decision tree using the decision tree classifier present in sklearn.
- Calculate different metrics such as precision, recall, auc\_roc curve and many more.
- Equal distribution of classes in testing and training data sets.
- Pre pruning and post pruning techniques such as cost complexity pruning technique.
- Different hyperparameters for decision tree classifiers.
- Implementation of decision surface boundaries.

### **References:**

- Scikit-learn documentation
- <https://www.kaggle.com/prashant111/decision-tree-classifier-tutorial>
- Towards data science - Medium
- Analytics Vidya