renduLab1

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1 Lab 1 Evaluation

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In the following, we consider the (binarized) Compas dataset that we studied in the Lab

```
[9]: from sklearn import tree from matplotlib import pyplot as plt # for a good visualization of the trees import csv import numpy as np from utils import load_from_csv

from sklearn.model_selection import train_test_split from sklearn.model_selection import cross_val_score from sklearn.metrics import confusion_matrix
```

```
[[1 0 0 ... 0 0 0]

[1 0 1 ... 0 0 0]

[1 0 0 ... 0 0 0]

...

[1 0 0 ... 0 0 0]

[0 1 0 ... 0 0 0]

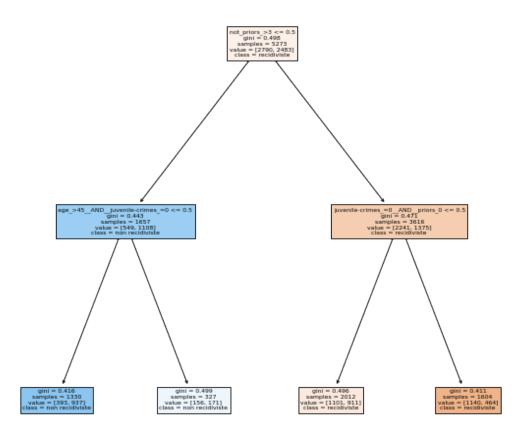
[1 0 0 ... 1 0 1]
```

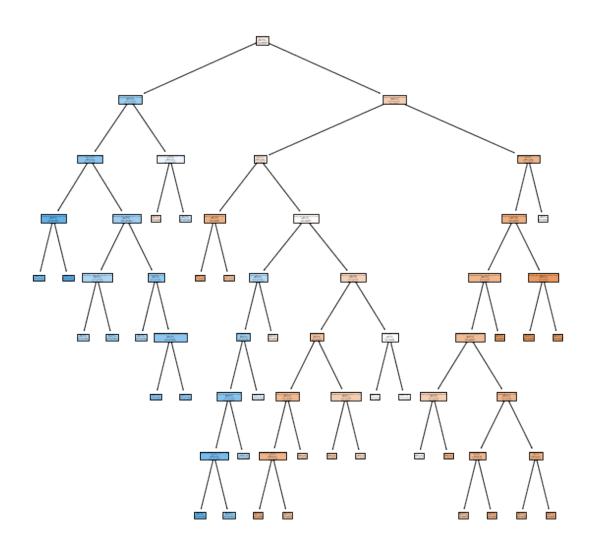
A decision tree configuration is a set of parameters that one can use to build decision trees. Propose 6 configurations that are likely to provide different topologies and caracteristics

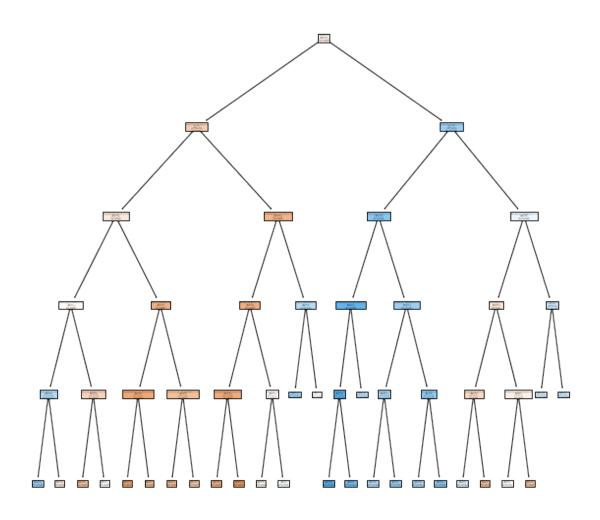
```
[11]: treesParmeters = [
#expected topologie and caracteristic
```

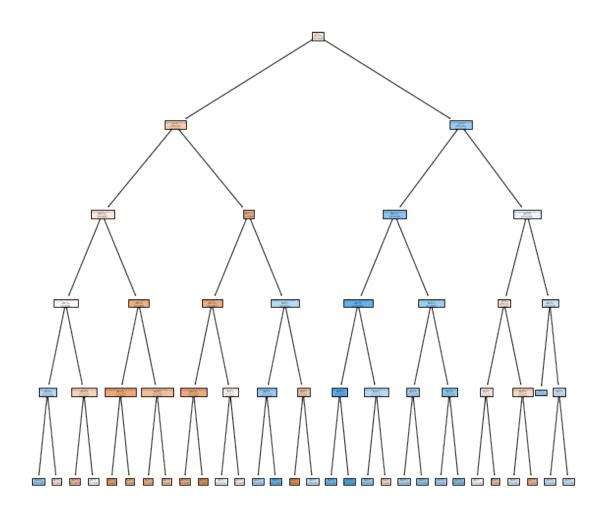
```
{'splitter':'best', 'max_depth': 2, 'min_samples_leaf': 1},
#expected topologie and caracteristic
    {'splitter':'random', 'max_depth': 10, 'min_samples_leaf': 100},
#expected topologie and caracteristic
    {'splitter':'best', 'max_depth': 5, 'min_samples_leaf': 20},
#expected topologie and caracteristic
    {'splitter':'best', 'max_depth': 5, 'min_samples_leaf':5},
#expected topologie and caracteristic
    {'splitter':'best', 'max_depth': 100, 'min_samples_leaf':60},
#expected topologie and caracteristic
    {'splitter':'best', 'max_depth': 100, 'min_samples_leaf':60},
]
nbTree = len(treesParmeters)
```

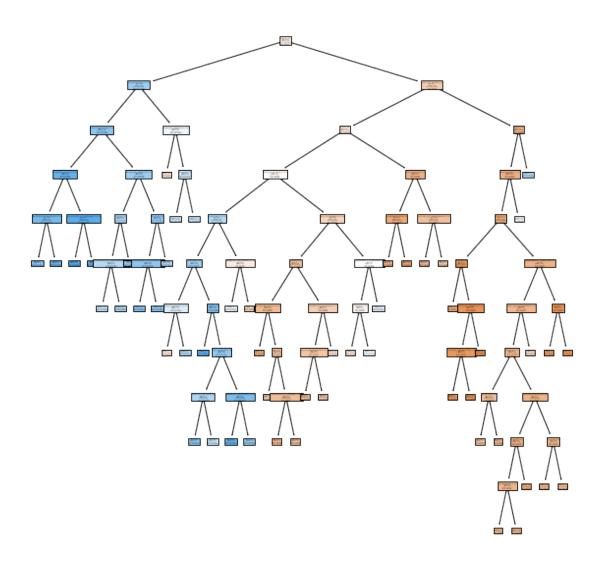
Train a decision tree for each of the previous configurations on the full dataset

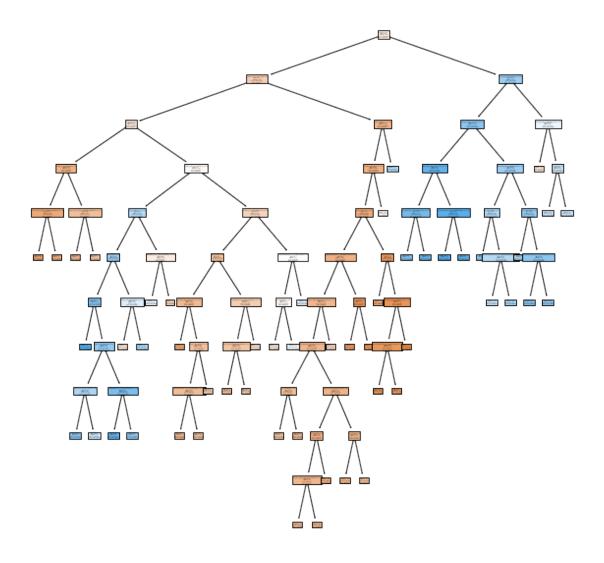










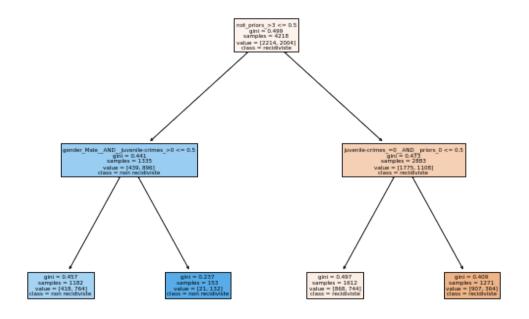


Propose an evaluation in terms of training and testing accuracies using 5-cross validation on two decision trees that have different typologies

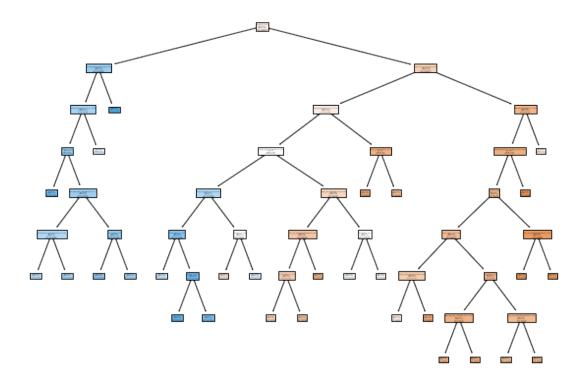
```
[13]: # We only keep the two first topologies
treesParmeters = treesParmeters[:2]
nbTree = len(treesParmeters)
```

```
for i in range (nbTree) :
    parameters= treesParmeters[i]
    clf = tree.DecisionTreeClassifier(
        splitter=parameters['splitter'],
        max_depth =parameters['max_depth'],
        min_samples_leaf =parameters['min_samples_leaf']
)
```

0.66 accuracy with a standard deviation of 0.02

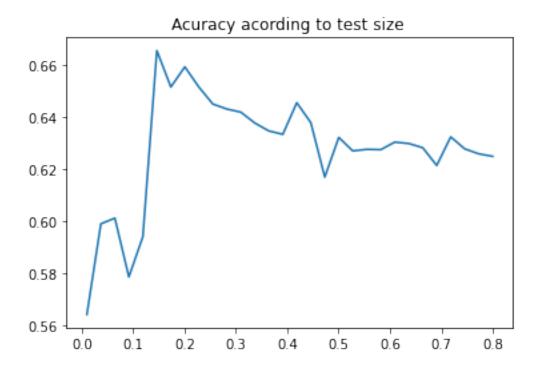


0.64 accuracy with a standard deviation of 0.01



Propose an experimental study that shows the transition phase from underfitting to overfitting

```
[15]: parameters= treesParmeters[0]
      acuracy =[]
      test_sizes=np.linspace(0.01,0.8,30)
      for test_size in test_sizes :
          clf = tree.DecisionTreeClassifier(
              splitter=parameters['splitter'],
              max_depth =parameters['max_depth'],
              min_samples_leaf =parameters['min_samples_leaf']
          X_train, X_test, y_train, y_test = train_test_split(train_examples,__
       strain_labels, test_size=test_size, random_state=42)
          clf = clf.fit(X_train, y_train)
          scores = cross_val_score(clf, X_test, y_test, cv=2)
          acuracy.append(scores.mean())
          \#print(f"for a test size of \{1/i\} we have an accuracy of \{scores.mean()\}_{\sqcup}
       ⇒with a standard deviation of {scores.std()}")
      plt.title('Acuracy acording to test size ')
      plt.plot(test_sizes,acuracy)
      plt.show()
```



1.0.3 Remarks

- With small tests size, we have bad result because our graph is to close of our data
 - this is overfitting
- With high tests size, our results start to be be lower because the graph is constructed with too few data
 - this is underfitting

Construct the confusion matrix on a particular good configuration (after explaining your choice)

Provide an evaluation of the fairness of the model based on the False Positive Rate

```
[]:
```