Assessment 2

Task 1: Develop learning-based model(s) for Classification

```
In [1]:
         # TensorFlow config to GPU
         import tensorflow as tf
         print(tf.__version__)
         gpus = tf.config.list_physical_devices('GPU')
         if gpus:
             tf.config.set_logical_device_configuration(
                 gpus[0],
                 [tf.config.LogicalDeviceConfiguration(memory_limit=15292)]
             )
         logical_gpus = tf.config.list_logical_devices('GPU')
         print(logical_gpus)
         print(len(gpus), "Physical GPU,", len(logical_gpus), "Logical GPUs")
         from tensorflow.python.client import device_lib
         print(device_lib.list_local_devices())
         print()
         print()
         import tensorflow as tf
         print("Num GPUs Available: ",
         len(tf.config.list_physical_devices('GPU')))
```

```
2.6.0
[LogicalDevice(name='/device:GPU:0', device_type='GPU')]
1 Physical GPU, 1 Logical GPUs
[name: "/device:CPU:0"
    device_type: "CPU"
    memory_limit: 268435456
locality {
}
incarnation: 8707722222192318641
, name: "/device:GPU:0"
    device_type: "GPU"
    memory_limit: 16034824192
locality {
    bus_id: 1
    links {
    }
}
```

```
physical_device_desc: "device: 0, name: NVIDIA GeForce RTX 2060, pci bus id: 0000:0
        1:00.0, compute capability: 7.5"
        Num GPUs Available: 1
In [2]:
         import pandas as pd
         import numpy as np
         import matplotlib.pyplot as plt
         import seaborn as sns
         from sklearn.preprocessing import StandardScaler
         from sklearn.metrics import accuracy_score
         from sklearn.metrics import confusion_matrix
         from sklearn.metrics import f1 score
         from sklearn.model_selection import train_test_split
         import warnings
         warnings.filterwarnings('ignore')
         from sklearn.metrics import precision_recall_curve
         from sklearn.metrics import plot_precision_recall_curve
         from sklearn.metrics import average_precision_score
         from sklearn.metrics import auc
In [3]:
         train_df = pd.read_csv('ECG_dataset/train.csv', header=None)
         test_df = pd.read_csv('ECG_dataset/test.csv', header=None)
         validation_df = pd.read_csv('ECG_dataset/validation.csv', header=None)
In [4]:
         print(train_df.shape)
         print(test_df.shape)
         print(validation df.shape)
        (1081, 141)
        (180, 141)
        (541, 141)
        We will be using train_df as our training data and validation_df as our validation data. We will be
        using test_df as our test data.
In [5]:
         train_df.head()
Out[5]:
                   1
                            2
                                     3
                                                      5
                                                                       7
                                                                                8
```

incarnation: 8088543449005111868

0 1

1 0

2 0

0.024133

0.424380

1.529500

0.016065

0.344420

1.776600

0.044639

0.348130

1.936700

0.340170

1.840200

0.243370

1.800000

0.031001 -0.009473 -0.042663 -0.077283 -0.091508

0.241730

1.724900

0.268780

1.405800

-0.04611

0.35644

0.72472

0.273420

1.008800

	0	1	2	3	4	5	6	7	8	!
3	0	1.286500	1.049900	0.793600	0.473590	0.111730	-0.054857	-0.062095	-0.120750	-0.10301
4	1	-0.175400	-0.121920	-0.053532	-0.024293	0.022917	0.116440	0.187040	0.240710	0.31434

5 rows × 141 columns

Columns: 141 entries, 0 to 140 dtypes: float64(140), int64(1)

memory usage: 1.2 MB

```
In [7]: train_df.shape
```

Out[7]: (1081, 141)

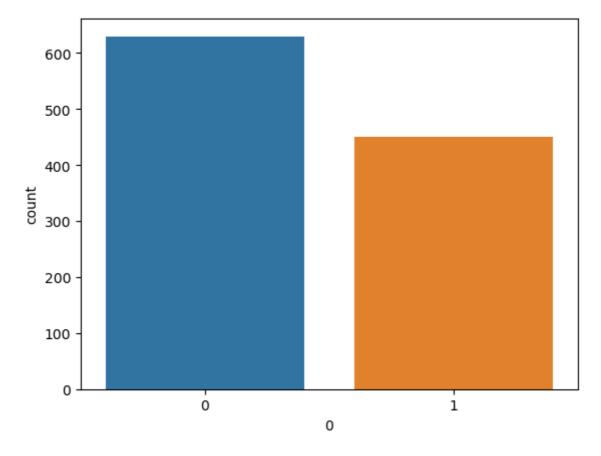
The first column of the data is the label. The remaining columns are the features. The features are the different ECG signals. The label is the class of the ECG signal. The classes are:

- 0 (not having a cardiovascular disease)
- 1 (having a cardiovascular disease)

Aim is to build a model that can predict the class of the ECG signal. This can be formulated as a binary classification problem.

```
In [9]: # plot countplot for first column
sns.countplot(x=0, data=train_df)
```

Out[9]: <AxesSubplot:xlabel='0', ylabel='count'>



```
#Showing the info of train and test data
print('###########Train data#########")
print(train_df.info())
print('###########Test data#########")
print(test_df.info())
#taking a sample of the test data (for reasons of plotting later)
sample = train_df.sample(25)
sampleX= sample.iloc[:,1:]
sampleY=sample.iloc[:,0]
print('###########Sample Info#########")
print(sampleX.info())
```

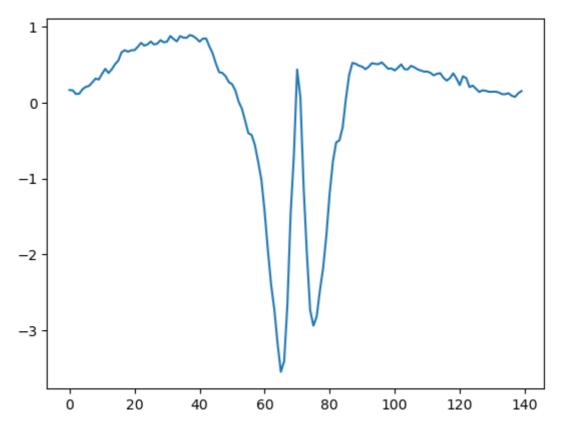
```
############Train data############
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 1081 entries, 0 to 1080
Columns: 141 entries, 0 to 140
dtypes: float64(140), int64(1)
memory usage: 1.2 MB
None
###########Test data###########
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 180 entries, 0 to 179
Columns: 141 entries, 0 to 140
dtypes: float64(140), int64(1)
memory usage: 198.4 KB
None
#############Sample Info############
<class 'pandas.core.frame.DataFrame'>
Int64Index: 25 entries, 10 to 327
Columns: 140 entries, 1 to 140
dtypes: float64(140)
```

memory usage: 27.5 KB

None

```
import matplotlib.pyplot as plt
plt.plot(np.array(range(0,140)),sampleX.iloc[10])
```

Out[11]: [<matplotlib.lines.Line2D at 0x21808f29d30>]

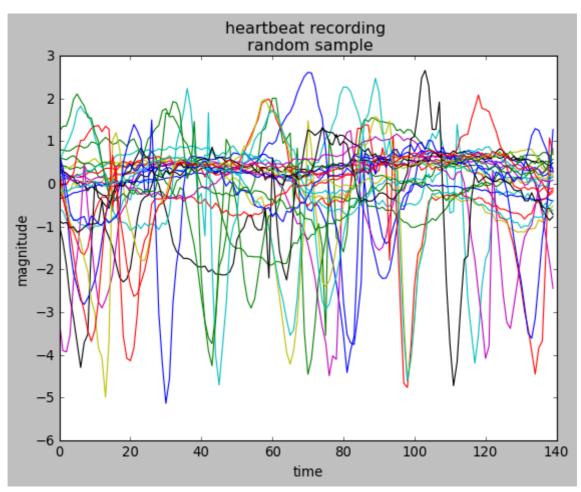


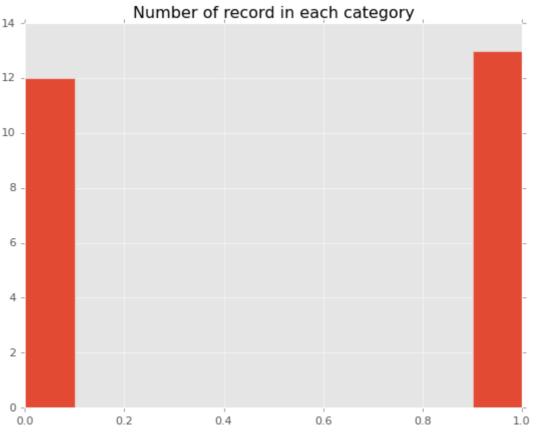
```
#ploting sample data information

plt.style.use('classic')
#ploting the samples
for index, row in sampleX.iterrows():
    plt.plot(np.array(range(0,140)), row)

plt.xlabel('time')
plt.ylabel('magnitude')
plt.title("heartbeat recording \n random sample")
plt.show()

plt.style.use('ggplot')
plt.title("Number of record in each category")
plt.hist(sample.iloc[:,0].transpose())
plt.show()
```





```
#number of labels for train and test
print("Train data")
print("Type\tCount")
print(train_df.iloc[:,0].value_counts())
```

```
print('#########################"')
print("Test data")
print("Type\tCount")
print(test_df.iloc[:,0].value_counts())
```

The models that we will be using are:

- Naive Bayes Classifier
- K-Nearest Neighbors Classifier
- Ensemble Learning Classifier
- Support Vector Machine Classifier

Then we will compare the performance of the models and select the best model for the task.

```
In [8]: # split the data into X_train and y_train, X_test and y_test
X_train = train_df.iloc[:, 1:]
y_train = train_df.iloc[:, 0]

X_valid = validation_df.iloc[:, 1:]
y_valid = validation_df.iloc[:, 0]

# Scale the data
scaler = StandardScaler()
X_train = scaler.fit_transform(X_train)
X_valid = scaler.transform(X_valid)
```

After splitting the data into training and validation sets, we need to scale the data. We will be using StandardScaler for this purpose.

1. Naive Bayes Classifier

```
In [66]: # Naive Bayes
from sklearn.naive_bayes import GaussianNB

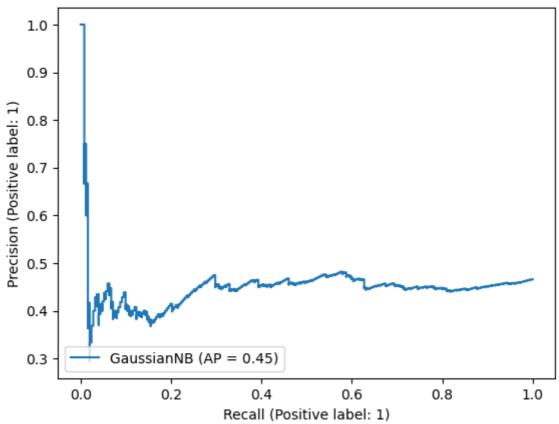
# Fit the model
gnb = GaussianNB()
gnb.fit(X_train, y_train)
```

Accuracy: 0.49353049907578556 F1 score: 0.4922838801304605 Confusion matrix: [[159 130]

[144 108]]

Average precision-recall score: 0.46

2-class Precision-Recall curve: AP=0.46



The naive Bayes classifier is a simple probabilistic classifier based on applying Bayes' theorem with strong (naive) independence assumptions between the features. A Naive Bayes classifier assumes that the presence of a particular feature in a class is unrelated to the presence of any other feature. For example, a fruit may be considered to be an apple if it is red, round, and about 3 inches in diameter. A Naive Bayes classifier considers each of these features to contribute independently to the probability that this fruit is an apple, regardless of any possible correlations between the color, roundness, and diameter features.

Gaussian Naive Bayes is used when the data is normally distributed. It is assumed that the data follows a normal distribution. Inorder to use Gaussian Naive Bayes, we need to convert the data into a normal distribution.

This model gave an accuracy of 0.49 on the validation set which is not good.

2. K-Nearest Neighbors Classifier

```
# KNN
from sklearn.neighbors import KNeighborsClassifier

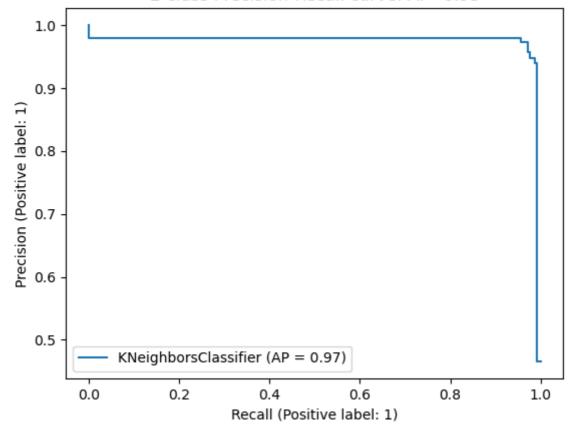
# Fit the model
knn = KNeighborsClassifier(n_neighbors=5)
knn.fit(X_train, y_train)

# Predict the model
y_pred_knn = knn.predict(X_valid)
```

Accuracy: 0.9685767097966729 F1 score: 0.968593948287203 Confusion matrix: [[278 11] [6 246]]

Average precision-recall score: 0.95

2-class Precision-Recall curve: AP=0.95



KNN (K-Nearest Neighbors) is a simple algorithm that stores all available cases and classifies new cases based on a similarity measure (e.g., distance functions). KNN has been used in statistical estimation and pattern recognition already in the beginning of 1970's as a non-parametric technique. The algorithm is among the simplest of all machine learning algorithms.

In KNN, the hyperparameter K is the number of nearest neighbors to be considered. The KNN algorithm assumes that similar things exist in close proximity. In other words, similar things are near to each other.

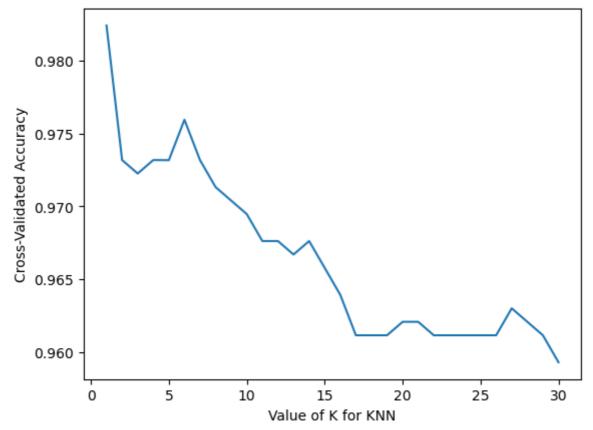
This k value is a hyperparameter that we need to tune. We will be using GridSearchCV to find the best value of k.

```
In [27]:
```

```
# using cross validation to find the best k
from sklearn.model_selection import cross_val_score
# search for an optimal value of K for KNN
k_range = range(1, 31)
k_scores = []
k values = []
for k in k_range:
    knn = KNeighborsClassifier(n neighbors=k)
    scores = cross_val_score(knn, X_train, y_train, cv=10,
scoring='accuracy')
    k_values.append(k)
    k_scores.append(scores.mean())
print(k_scores)
print()
print(max(k_scores))
optimal_k = k_values[k_scores.index(max(k_scores))]
print(optimal_k)
# plot the value of K for KNN (x-axis) versus the cross-validated
accuracy (y-axis)
plt.plot(k_range, k_scores)
plt.xlabel('Value of K for KNN')
plt.ylabel('Cross-Validated Accuracy')
plt.show()
```

[0.9824159021406726, 0.9731736323479444, 0.9722477064220184, 0.9731736323479444, 0.9731651376146789, 0.9759429153924566, 0.9731651376146789, 0.971313285762827, 0.970387 359836901, 0.9694614339109752, 0.9676095820591233, 0.9676095820591233, 0.96668365613 31975, 0.9676095820591234, 0.9657662249405368, 0.963914373088685, 0.961136595310907 3, 0.9611365953109073, 0.9611365953109073, 0.9620625212368331, 0.9620625212368333, 0.9611365953109073, 0.9611365953109073, 0.9611365953109073, 0.9620625212368331, 0.9620625212368331, 0.9629884471627591, 0.9620625212368331, 0.9611365953109072, 0.959284 7434590555]

```
0.9824159021406726
```

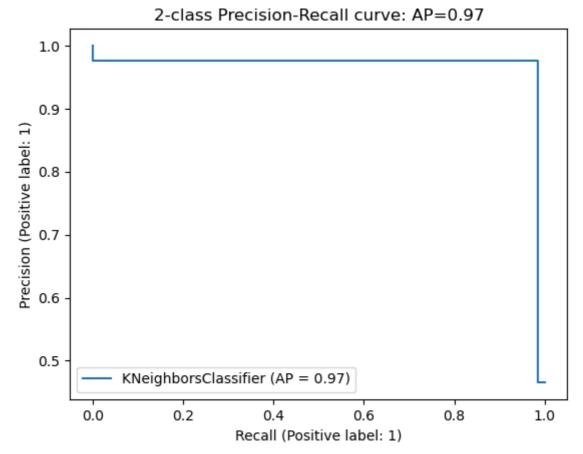


```
In [68]:
          # Create a new KNN model with the optimal k
          knn = KNeighborsClassifier(n_neighbors=optimal_k)
          # Fit the model
          knn.fit(X_train, y_train)
          # Predict the model
          y_pred_knn = knn.predict(X_valid)
          # Evaluate the model
          print("Accuracy: ", accuracy_score(y_valid, y_pred_knn))
          print("F1 score: ", f1_score(y_valid, y_pred_knn, average='weighted'))
          print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_knn))
          # Precision and recall and Area under the curv
          disp = plot_precision_recall_curve(knn, X_valid, y_valid)
          disp.ax_.set_title('2-class Precision-Recall curve: '
                                  'AP=
          {0:0.2f}'.format(average_precision_score(y_valid, y_pred_knn)))
          # Area under the curve
          average_precision = average_precision_score(y_valid, y_pred_knn)
```

Accuracy: 0.9815157116451017 F1 score: 0.9815201510808478 Confusion matrix: [[283 6]

[4 248]]

Average precision-recall score: 0.97



Cross validation helped us to find the best value of k. The best value of k is 1. This model gave an accuracy of 96% when k was set to 5. Now that we have found the best value of k, it gave an accuracy of 98% on the validation set.

3. Ensemble Learning Classifier

Ensemble learning is a machine learning paradigm where multiple models, such as classifiers or experts, are strategically generated and combined to solve a particular computational intelligence problem. Ensemble methods use multiple learning algorithms to obtain better predictive performance than could be obtained from any of the constituent learning algorithms alone.

We will be using the following ensemble learning models:

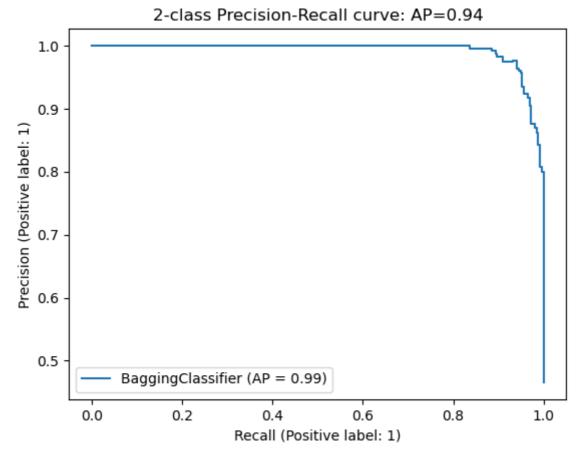
Bagging Classifier: A Bagging classifier is an ensemble meta-estimator that fits base
classifiers each on random subsets of the original dataset and then aggregate their
individual predictions (either by voting or by averaging) to form a final prediction. Such a
meta-estimator can typically be used as a way to reduce the variance of a black-box
estimator (e.g., a decision tree), by introducing randomization into its construction
procedure and then making an ensemble out of it.

- Random Forest Classifier: A random forest is a meta estimator that fits a number of
 decision tree classifiers on various sub-samples of the dataset and uses averaging to
 improve the predictive accuracy and control over-fitting. The sub-sample size is always the
 same as the original input sample size but the samples are drawn with replacement if
 bootstrap=True (default).
- AdaBoost Classifier: AdaBoost (Adaptive Boosting) is a machine learning meta-algorithm
 formulated by Yoav Freund and Robert Schapire. It is a boosting algorithm that can be used
 to convert weak learners to strong ones. It is one of the most popular boosting algorithms.
 It is used in a variety of areas including computer vision and speech recognition.
- Gradient Boosting Classifier: Gradient Boosting is a machine learning technique for regression and classification problems, which produces a prediction model in the form of an ensemble of weak prediction models, typically decision trees. It builds the model in a stagewise fashion like other boosting methods do, and it generalizes them by allowing optimization of an arbitrary differentiable loss function.

```
In [69]:
          # Bagging Classifier
          from sklearn.ensemble import BaggingClassifier
          # fit the model
          bagging = BaggingClassifier(n_estimators=100, max_samples=0.5,
          max_features=0.5)
          bagging.fit(X_train, y_train)
          # Predict the model
          y_pred_bagging = bagging.predict(X_valid)
          # Evaluate the model
          print("Accuracy: ", accuracy_score(y_valid, y_pred_bagging))
          print("F1 score: ", f1_score(y_valid, y_pred_bagging,
          average='weighted'))
          print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_bagging))
          # Precision and recall and Area under the curv
          disp = plot_precision_recall_curve(bagging, X_valid, y_valid)
          disp.ax_.set_title('2-class Precision-Recall curve: '
                                  'AP=
          {0:0.2f}'.format(average_precision_score(y_valid, y_pred_bagging)))
          # Area under the curve
          average_precision = average_precision_score(y_valid, y_pred_bagging)
```

Accuracy: 0.9611829944547134 F1 score: 0.9611276875860413 Confusion matrix: [[283 6] [15 237]]

Average precision-recall score: 0.94



```
# implementing cross validation to find the best n_estimators,
    max_samples and max_features using GridSearchCV
from sklearn.model_selection import GridSearchCV

# Create the parameter grid based on the results of random search
param_grid = {
    'n_estimators': [100, 200, 300, 400, 500],
    'max_samples': [0.5, 0.6, 0.7, 0.8, 0.9],
    'max_features': [0.5, 0.6, 0.7, 0.8, 0.9]
}

# Create a based model
bagging = BaggingClassifier()

# Instantiate the grid search model
grid_search = GridSearchCV(estimator=bagging, param_grid=param_grid, cv=3, n_jobs=-1, verbose=2)
```

```
# Fit the grid search to the data
grid search.fit(X train, y train)
# print the best parameters
print(grid_search.best_params_)
print(grid_search.best_score_)
print(grid_search.best_estimator_)
```

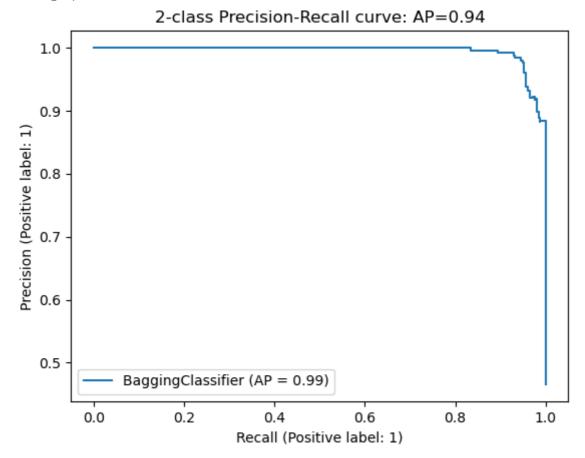
```
Fitting 3 folds for each of 125 candidates, totalling 375 fits
{'max_features': 0.6, 'max_samples': 0.9, 'n_estimators': 300}
0.9491022878834512
BaggingClassifier(max_features=0.6, max_samples=0.9, n estimators=300)
Accuracy: 0.9648798521256932
F1 score: 0.9648545004581948
Confusion matrix: [[282
[ 12 240]]
```

```
In [72]:
```

```
# Create a new Bagging Classifier model with the optimal parameters
 bagging = BaggingClassifier(n_estimators=300,
                             max_samples=0.9,
                             max features=0.6)
 # Fit the model
 bagging.fit(X_train, y_train)
 # Predict the model
 y_pred_bagging = bagging.predict(X_valid)
 # Evaluate the model
 print("Accuracy: ", accuracy_score(y_valid, y_pred_bagging))
 print("F1 score: ", f1_score(y_valid, y_pred_bagging,
 average='weighted'))
 print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_bagging))
 # Precision and recall and Area under the curv
 disp = plot_precision_recall_curve(bagging, X_valid, y_valid)
 disp.ax_.set_title('2-class Precision-Recall curve: '
                         'AP=
 {0:0.2f}'.format(average_precision_score(y_valid, y_pred_bagging)))
 # Area under the curve
 average_precision = average_precision_score(y_valid, y_pred_bagging)
```

Accuracy: 0.9630314232902033 F1 score: 0.9630105887658721 Confusion matrix: [[281 8] [12 240]]

Average precision-recall score: 0.94



No big difference in the accuracy of the models while using cross validation with 96% accuracy.

```
# Random Forest
from sklearn.ensemble import RandomForestClassifier

# fit the model
rf = RandomForestClassifier(n_estimators=100, max_depth=2,
random_state=0)
rf.fit(X_train, y_train)

# Predict the model
y_pred_rf = rf.predict(X_valid)

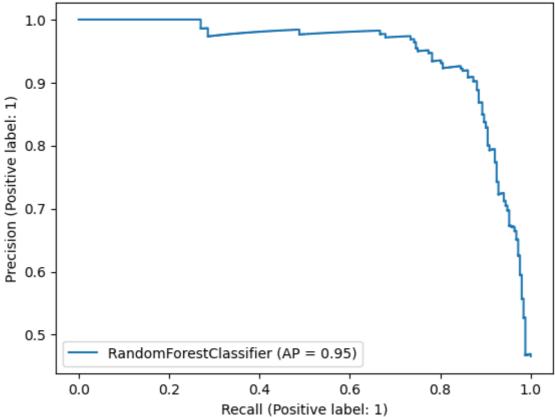
# Evaluate the model
print("Accuracy: ", accuracy_score(y_valid, y_pred_rf))
print("F1 score: ", f1_score(y_valid, y_pred_rf, average='weighted'))
print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_rf))
```

Accuracy: 0.8188539741219963 F1 score: 0.8110658237266908 Confusion matrix: [[286 3]

[95 157]]

Average precision-recall score: 0.79





```
# implementing cross validation to find the best n_estimators, max_depth
and random_state using GridSearchCV
from sklearn.model_selection import GridSearchCV

# Create the parameter grid based on the results of random search
param_grid = {
    'n_estimators': [100, 200, 300, 400, 500],
    'max_depth': [2, 3, 4, 5, 6, 7, 8, 9, 10],
    'random_state': [0, 1, 2, 3, 4, 5, 6, 7, 8, 9]
```

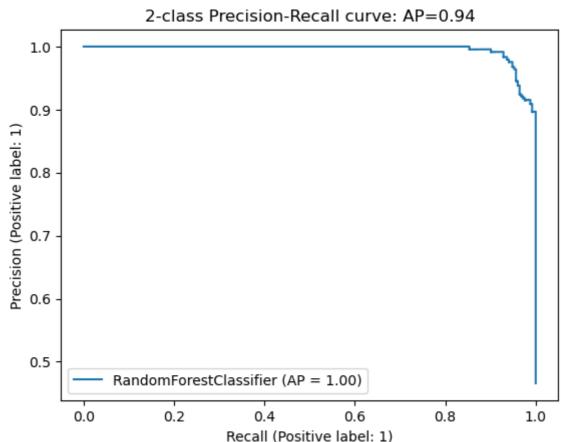
```
# Create a based model
rf = RandomForestClassifier()
 # Instantiate the grid search model
 grid search = GridSearchCV(estimator=rf, param_grid=param_grid, cv=3,
 n_jobs=-1, verbose=2)
# Fit the grid search to the data
 grid_search.fit(X_train, y_train)
# print the best parameters
 print(grid_search.best_params_)
print(grid_search.best_score_)
print(grid_search.best_estimator_)
Fitting 3 folds for each of 450 candidates, totalling 1350 fits
{'max_depth': 10, 'n_estimators': 300, 'random_state': 0}
0.9528008618036319
RandomForestClassifier(max_depth=10, n_estimators=300, random_state=0)
# Create a new Random Forest model with the optimal parameters
 rf = RandomForestClassifier(n_estimators=300,
                              max_depth=10,
                              random_state=0)
```

In [74]: # Fit the model rf.fit(X_train, y_train) # Predict the model y pred rf = rf.predict(X valid) # Evaluate the model print("Accuracy: ", accuracy_score(y_valid, y_pred_rf)) print("F1 score: ", f1_score(y_valid, y_pred_rf, average='weighted')) print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_rf)) # Precision and recall and Area under the curv disp = plot_precision_recall_curve(rf, X_valid, y_valid) disp.ax_.set_title('2-class Precision-Recall curve: ' 'AP= {0:0.2f}'.format(average_precision_score(y_valid, y_pred_rf)))

Accuracy: 0.9630314232902033 F1 score: 0.9630105887658721 Confusion matrix: [[281 8]

[12 240]]

Average precision-recall score: 0.94



Random Forest Classifier is an extension of bagging classifier. It is a very powerful model. It gave an accuracy of 96% on the validation set which is an improvement over the previous model without cross validation.

```
# AdaBoost
from sklearn.ensemble import AdaBoostClassifier

# fit the model
ada = AdaBoostClassifier(n_estimators=100, random_state=0)
ada.fit(X_train, y_train)

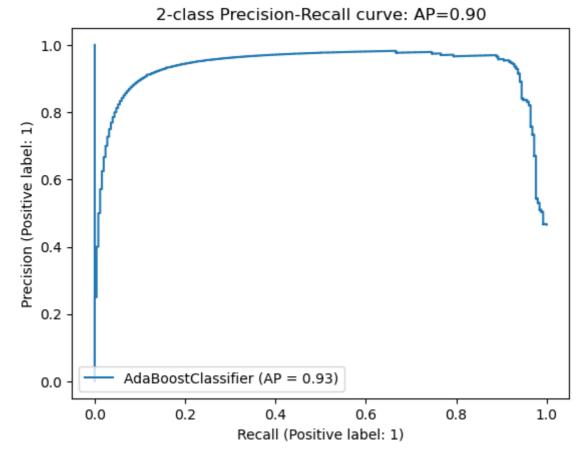
# Predict the model
y_pred_ada = ada.predict(X_valid)

# Evaluate the model
print("Accuracy: ", accuracy_score(y_valid, y_pred_ada))
```

Accuracy: 0.9353049907578558 F1 score: 0.935327626521316 Confusion matrix: [[270 19]

[16 236]]

Average precision-recall score: 0.90



```
# implementing cross validation to find the best n_estimators and
random_state using GridSearchCV
from sklearn.model_selection import GridSearchCV

# Create the parameter grid based on the results of random search
param_grid = {
```

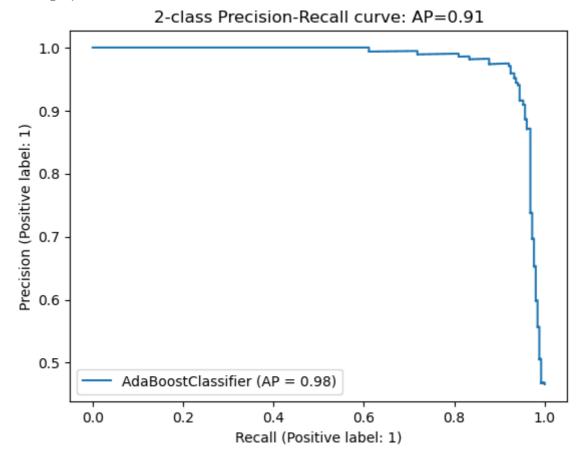
```
'n_estimators': [100, 200, 300, 400, 500],
               'random_state': [0, 1, 2, 3, 4, 5, 6, 7, 8, 9]
          }
          # Create a based model
          ada = AdaBoostClassifier()
          # Instantiate the grid search model
          grid search = GridSearchCV(estimator=ada, param grid=param grid, cv=3,
          n_jobs=-1, verbose=2)
          # Fit the grid search to the data
          grid_search.fit(X_train, y_train)
          # print the best parameters
          print(grid_search.best_params_)
          print(grid_search.best_score_)
          print(grid_search.best_estimator_)
         Fitting 3 folds for each of 50 candidates, totalling 150 fits
         {'n_estimators': 400, 'random_state': 0}
         0.9306094182825485
         AdaBoostClassifier(n_estimators=400, random_state=0)
In [45]:
         print(grid_search.best_estimator_)
         AdaBoostClassifier(n_estimators=400, random_state=0)
         # Create a new AdaBoost model with the optimal parameters
```

```
In [77]:
          ada = AdaBoostClassifier(n_estimators=400,
                                      random_state=0)
          # Fit the model
          ada.fit(X_train, y_train)
          # Predict the model
          y_pred_ada = ada.predict(X_valid)
          # Evaluate the model
          print("Accuracy: ", accuracy score(y valid, y pred ada))
          print("F1 score: ", f1_score(y_valid, y_pred_ada, average='weighted'))
          print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_ada))
          # Precision and recall and Area under the curv
```

Accuracy: 0.944547134935305 F1 score: 0.9445322794487168 Confusion matrix: [[275 14]

[16 236]]

Average precision-recall score: 0.91



AdaBoost Classifier is a boosting classifier. It is a very powerful model. It gave an accuracy of 94% on the validation set.

```
# Gradient Boosting
from sklearn.ensemble import GradientBoostingClassifier

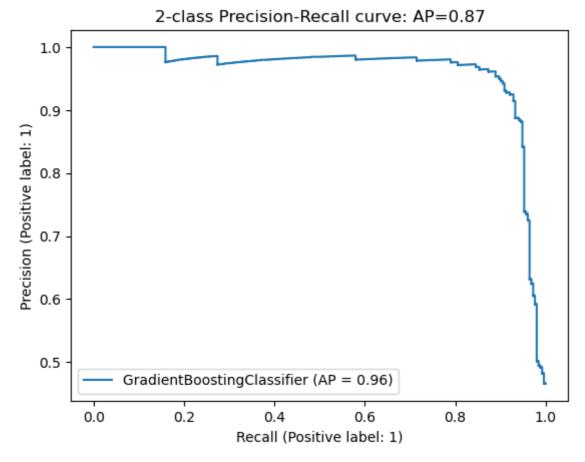
# fit the model
gb = GradientBoostingClassifier(n_estimators=100, learning_rate=1.0,
max_depth=1, random_state=0)
gb.fit(X_train, y_train)

# Predict the model
```

Accuracy: 0.922365988909427 F1 score: 0.9224277045432434 Confusion matrix: [[264 25]

[17 235]]

Average precision-recall score: 0.87



In [48]:

implementing cross validation to find the best n_estimators, learning_rate, max_depth and random_state using GridSearchCV

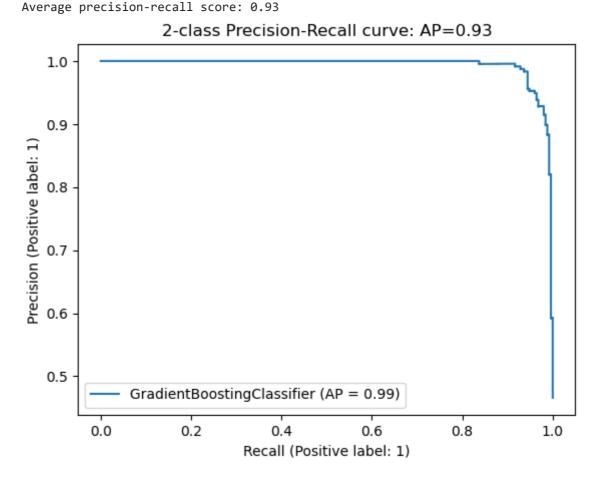
```
from sklearn.model_selection import GridSearchCV
          # Create the parameter grid based on the results of random search
          param_grid = {
              'n_estimators': [100, 200, 300, 400, 500],
               'learning_rate': [0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1.0],
               'max_depth': [1, 2, 3, 4, 5, 6, 7, 8, 9, 10],
               'random_state': [0, 1, 2, 3, 4, 5, 6, 7, 8, 9]
          }
          # Create a based model
          gb = GradientBoostingClassifier()
          # Instantiate the grid search model
          grid_search = GridSearchCV(estimator=gb, param_grid=param_grid, cv=3,
          n_jobs=-1, verbose=2)
          # Fit the grid search to the data
          grid_search.fit(X_train, y_train)
          # print the best parameters
          print(grid_search.best_params_)
          print(grid_search.best_score_)
          print(grid_search.best_estimator_)
         Fitting 3 folds for each of 5000 candidates, totalling 15000 fits
         {'learning_rate': 0.5, 'max_depth': 7, 'n_estimators': 100, 'random_state': 2}
         0.9583538524674259
         GradientBoostingClassifier(learning_rate=0.5, max_depth=7, random_state=2)
In [49]:
          print(grid search.best estimator )
          # GradientBoostingClassifier(learning rate=0.5, max depth=7,
          random state=2)
         GradientBoostingClassifier(learning rate=0.5, max depth=7, random state=2)
In [79]:
          # Create a new Gradient Boosting model with the optimal parameters
          gb = GradientBoostingClassifier(n_estimators=100,
                                           learning rate=0.5,
                                           max depth=7,
                                           random_state=2)
```

Fit the model

gb.fit(X_train, y_train)

```
# Predict the model
y_pred_gb = gb.predict(X_valid)
# Evaluate the model
print("Accuracy: ", accuracy_score(y_valid, y_pred_gb))
print("F1 score: ", f1_score(y_valid, y_pred_gb, average='weighted'))
print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_gb))
# Precision and recall and Area under the curv
disp = plot_precision_recall_curve(gb, X_valid, y_valid)
disp.ax_.set_title('2-class Precision-Recall curve: '
{0:0.2f}'.format(average precision_score(y valid, y pred gb)))
# Area under the curve
average_precision = average_precision_score(y_valid, y_pred_gb)
print('Average precision-recall score: {0:0.2f}'.format(
        average_precision))
```

Accuracy: 0.955637707948244 F1 score: 0.9556258235589734 Confusion matrix: [[278 11] [13 239]]



Gradient Boosting Classifier is a boosting classifier. It is a very powerful model. It gave an

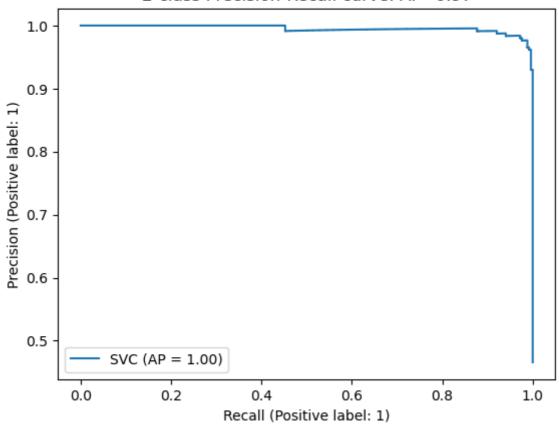
accuracy of 95% on the validation set which is an improvement over the previous model without cross validation (92%).

4. Support Vector Machine Classifier

```
In [80]:
        # SVM Classifier
          from sklearn.svm import SVC
          # fit the model
          svm = SVC(gamma='auto')
          svm.fit(X_train, y_train)
          # Predict the model
          y_pred_svm = svm.predict(X_valid)
          # Evaluate the model
          print("Accuracy: ", accuracy_score(y_valid, y_pred_svm))
          print("F1 score: ", f1_score(y_valid, y_pred_svm, average='weighted'))
          print("Confusion matrix: ", confusion_matrix(y_valid, y_pred_svm))
          # Precision and recall and Area under the curv
          disp = plot_precision_recall_curve(svm, X_valid, y_valid)
          disp.ax_.set_title('2-class Precision-Recall curve: '
                                  'AP=
          {0:0.2f}'.format(average_precision_score(y_valid, y_pred_svm)))
          # Area under the curve
          average_precision = average_precision_score(y_valid, y_pred_svm)
          print('Average precision-recall score: {0:0.2f}'.format(
                  average_precision))
```

```
Accuracy: 0.9815157116451017
F1 score: 0.9815201510808478
Confusion matrix: [[283 6]
[ 4 248]]
Average precision-recall score: 0.97
```

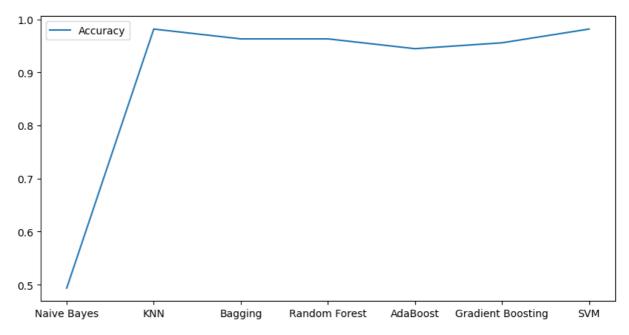
2-class Precision-Recall curve: AP=0.97



SVM (Support Vector Machine) is a supervised machine learning algorithm which can be used for both classification or regression challenges. However, it is mostly used in classification problems. In this algorithm, we plot each data item as a point in n-dimensional space (where n is number of features you have) with the value of each feature being the value of a particular coordinate. Then, we perform classification by finding the hyper-plane that differentiate the two classes very well.

And from this, we can see that this model gave an accuracy of 98% on the validation set.

5. Comparison of the models



Based on the accuracy of the models on validation data, K-Nearest Neighbors Classifier and SVM, both gave an accuracy of 98% which is impressive.

```
In [85]:
         # Testing these models on test data
         X_test = test_df.iloc[:, 1:]
         y_test = test_df.iloc[:, 0]
         # Scale the data
         # scaler = StandardScaler()
         # X_test = scaler.fit_transform(X_test)
         # Naive Bayes
         print("Naive Bayes")
         y_pred_gnb = gnb.predict(X_test)
         # Evaluate the model
         print("Accuracy: ", accuracy_score(y_test, y_pred_gnb))
         print("F1 score: ", f1_score(y_test, y_pred_gnb, average='weighted'))
         print("Confusion matrix: ", confusion_matrix(y_test, y_pred_gnb))
         # KNN
         print("KNN")
```

```
y_pred_knn = knn.predict(X_test)
# Evaluate the model
print("Accuracy: ", accuracy_score(y_test, y_pred_knn))
print("F1 score: ", f1_score(y_test, y_pred_knn, average='weighted'))
print("Confusion matrix: ", confusion_matrix(y_test, y_pred_knn))
print("**************************")
# Bagging
print("Bagging")
y_pred_bagging = bagging.predict(X_test)
# Evaluate the model
print("Accuracy: ", accuracy_score(y_test, y_pred_bagging))
print("F1 score: ", f1_score(y_test, y_pred_bagging,
average='weighted'))
print("Confusion matrix: ", confusion_matrix(y_test, y_pred_bagging))
print("*************************")
# Random Forest
print("Random Forest")
y_pred_rf = rf.predict(X_test)
# Evaluate the model
print("Accuracy: ", accuracy_score(y_test, y_pred_rf))
print("F1 score: ", f1_score(y_test, y_pred_rf, average='weighted'))
print("Confusion matrix: ", confusion_matrix(y_test, y_pred_rf))
# AdaBoost
print("AdaBoost")
y_pred_ada = ada.predict(X_test)
# Evaluate the model
print("Accuracy: ", accuracy_score(y_test, y_pred_ada))
print("F1 score: ", f1_score(y_test, y_pred_ada, average='weighted'))
print("Confusion matrix: ", confusion_matrix(y_test, y_pred_ada))
print("*************************")
# Gradient Boosting
print("Gradient Boosting")
y_pred_gb = gb.predict(X_test)
# Evaluate the model
print("Accuracy: ", accuracy_score(y_test, y_pred_gb))
```

```
Accuracy: 0.5277777777778
F1 score: 0.5265700483091788
Confusion matrix: [[61 41]
[44 34]]
*************
KNN
Accuracy: 0.9833333333333333
F1 score: 0.9833453609864768
Confusion matrix: [[100
[ 1 77]]
*************
Bagging
Accuracy: 0.97777777777777
F1 score: 0.977741352498634
Confusion matrix: [[101
[ 3 75]]
*************
Random Forest
Accuracy: 0.9833333333333333
F1 score: 0.9833202202989771
Confusion matrix: [[101
[ 2 76]]
************
AdaBoost
Accuracy: 0.9611111111111111
F1 score: 0.9611391756351123
Confusion matrix: [[98 4]
[ 3 75]]
***************
Gradient Boosting
Accuracy: 0.988888888888888
F1 score: 0.988888888888888
Confusion matrix: [[101
[ 1 77]]
*************
SVM
Accuracy: 0.97777777777777
F1 score: 0.9778084137527677
Confusion matrix: [[99 3]
[ 1 77]]
************
```

```
accuracy_score(y_test, y_pred_ada), accuracy_score(y_test, y_pred_gb),
           accuracy_score(y_test, y_pred_svm)]
In [87]:
           test_accuracy
         [0.527777777777778,
Out[87]:
           0.98333333333333333333
          0.9777777777777777
          0.98333333333333333333
           0.9611111111111111,
           0.98888888888889,
          0.97777777777777777
In [88]:
           plt.figure(figsize=(10,5))
           plt.plot(models, accuracy, label='Validation Accuracy')
           plt.plot(models, test_accuracy, label='Test Accuracy')
           plt.legend()
           plt.show()
          1.0
          0.9
          0.8
          0.7
          0.6
                                                                             Validation Accuracy
          0.5
                                                                             Test Accuracy
```

But, Gradient Boosting seems to produce better results than K-Nearest Neighbors Classifier and SVM on Testing Data. It gave an accuracy of 98% on the test set.

Bagging

Random Forest

AdaBoost

Gradient Boosting

KNN

Naive Bayes

Its hard to say which model is better. Gradient Boosting Classifier and SVM both gave an accuracy of 98% on the validation set. But, Gradient Boosting Classifier gave an accuracy of 98% on the test set. So, we will be using Gradient Boosting Classifier for the task.

Assessment 2

Task 2: Develop recurrent neural network(s) for sequence-to-sequence classification

```
In [1]:
         # TensorFlow config to GPU
         import tensorflow as tf
         print(tf.__version__)
         gpus = tf.config.list_physical_devices('GPU')
         if gpus:
             tf.config.set_logical_device_configuration(
                 gpus[0],
                 [tf.config.LogicalDeviceConfiguration(memory_limit=15292)]
             )
         logical_gpus = tf.config.list_logical_devices('GPU')
         print(logical_gpus)
         print(len(gpus), "Physical GPU,", len(logical_gpus), "Logical GPUs")
         from tensorflow.python.client import device_lib
         print(device_lib.list_local_devices())
         print()
         print()
         import tensorflow as tf
         print("Num GPUs Available: ",
         len(tf.config.list_physical_devices('GPU')))
```

```
2.6.0
[LogicalDevice(name='/device:GPU:0', device_type='GPU')]
1 Physical GPU, 1 Logical GPUs
[name: "/device:CPU:0"
  device_type: "CPU"
  memory_limit: 268435456
locality {
}
incarnation: 2201299975258163592
, name: "/device:GPU:0"
  device_type: "GPU"
  memory_limit: 16034824192
locality {
    bus_id: 1
    links {
    }
}
```

```
incarnation: 3233997999066038929
physical_device_desc: "device: 0, name: NVIDIA GeForce RTX 2060, pci bus id: 0000:0
1:00.0, compute capability: 7.5"
Num GPUs Available: 1
#loading environment
 import pandas as pd
 import numpy as np
 ''' Data visualisation'''
 import matplotlib.pyplot as plt
 import seaborn as sns
 import warnings
 ''' Scikit-Learn'''
 from sklearn.model_selection import train_test_split
 from sklearn.preprocessing import StandardScaler, RobustScaler
 from sklearn.preprocessing import OneHotEncoder
 from sklearn.model_selection import cross_validate, cross_val_score
 from sklearn.model_selection import GridSearchCV
 from sklearn.model_selection import RandomizedSearchCV
 from sklearn import set_config
 set_config(display='diagram')
 from sklearn.compose import ColumnTransformer
 from sklearn.compose import make_column_selector
 from sklearn.metrics import confusion_matrix
 ''' Imbalanced Classes'''
 import imblearn
 from imblearn.over sampling import SMOTE
 from imblearn.under_sampling import RandomUnderSampler
 from imblearn.pipeline import Pipeline
 ''' Tensorflow Keras'''
 from tensorflow import keras
 from tensorflow.keras import models
 from tensorflow.keras import Sequential, layers
 from tensorflow.keras.callbacks import EarlyStopping
```

```
In [2]:
         from tensorflow.keras import regularizers
         from tensorflow.keras.optimizers import Adam
         from tensorflow.keras.optimizers.schedules import ExponentialDecay
         from keras.utils import np_utils
         from tensorflow.keras.utils import to_categorical
         from keras.preprocessing import image
```

```
from os import listdir
from os.path import isdir, join
import warnings
warnings.filterwarnings('ignore')
```

```
In [46]:
```

```
# Set Pandas options to display more columns
pd.options.display.max_columns=200

# Read in the data csv
df=pd.read_csv('ECG_dataset/trainval.csv', encoding='utf-8', header=None)

# Show a snaphsot of data
df
```

Out[46]:		0	1	2	3	4	5	6	7	8	
	0	1	0.024133	0.016065	0.044639	0.031001	-0.009473	-0.042663	-0.077283	-0.091508	-0.04
	1	0	0.424380	0.344420	0.348130	0.340170	0.243370	0.241730	0.268780	0.273420	0.35
	2	0	1.529500	1.776600	1.936700	1.840200	1.800000	1.724900	1.405800	1.008800	0.72
	3	0	1.286500	1.049900	0.793600	0.473590	0.111730	-0.054857	-0.062095	-0.120750	-0.10
	4	1	-0.175400	-0.121920	-0.053532	-0.024293	0.022917	0.116440	0.187040	0.240710	0.31
	•••										
	1617	1	0.429370	0.603980	0.663720	0.501780	0.323160	0.296740	0.354540	0.537550	0.44
	1618	0	-0.552610	-0.382620	-0.331440	-0.240030	-0.227290	-0.208440	-0.146600	-0.050338	-0.05
	1619	0	1.716600	1.884900	1.808500	1.750200	1.541700	1.352600	1.107600	0.884450	0.53
	1620	0	1.120500	1.227900	1.311500	1.814600	1.954700	1.891200	1.792000	1.437400	1.29

1622 rows × 141 columns

The first column of the data is the label. The remaining columns are the features. The features are the different ECG signals. The label is the class of the ECG signal. The classes are:

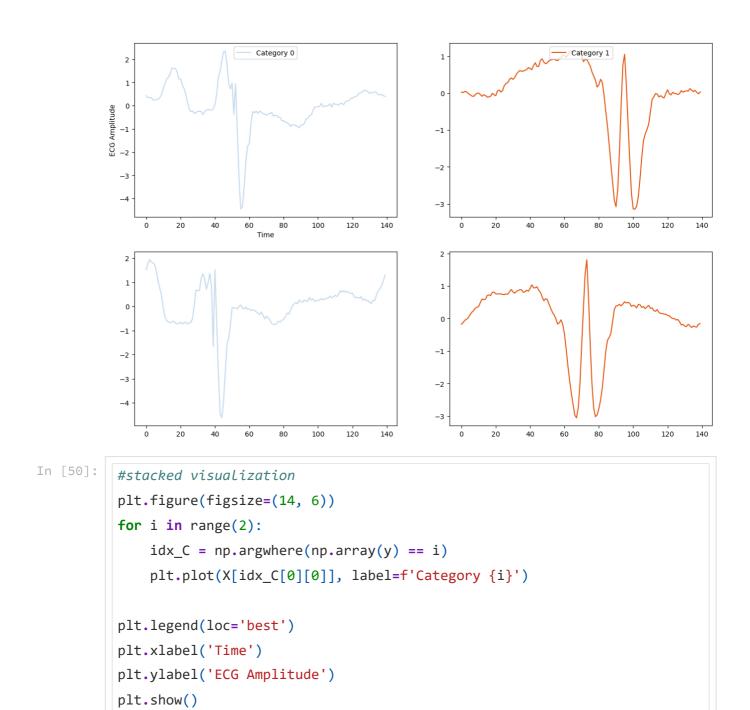
1621 1 1.453000 1.084100 1.098700 1.135100 1.466200 1.482600 0.970650 0.205070 -0.83

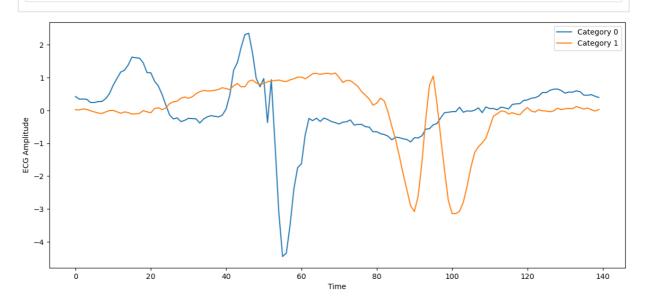
- 0 (not having a cardiovascular disease)
- 1 (having a cardiovascular disease)

Aim is to build a model that can predict the class of the ECG signal using RNNs.

```
Out[47]: 0.024133
In [48]:
          y = df.iloc[:, 0]
          y = y.to_numpy().tolist()
          y[0]
Out[48]: 1
In [49]:
          # visualizing data
          from matplotlib.cm import get_cmap
          name = 'tab20c'
          cmap = get_cmap(name)
          colors_list = cmap.colors
          fix, axs = plt.subplots(2, 2, figsize=(15,10))
          for i in range(2):
              for j in range(2):
                  idx_C = np.argwhere(np.array(y) == j)
                  axs[i,j].plot(X[idx_C[i][0]], label=f'Category {j}',
          c=colors_list[j+3])
                  if i == 0:
                      axs[i,j].legend(loc='upper center')
                      if j ==0:
                          axs[i,j].set_xlabel('Time')
```

axs[i,j].set_ylabel('ECG Amplitude')





```
In [51]: from tensorflow.keras.preprocessing.sequence import pad_sequences
X_pad = pad_sequences(X, dtype='float32', padding='post', value=-1)
```

```
print(X_pad.shape)
          print(X_pad[0][0])
         (1622, 140)
         0.024133
In [52]:
         #reshaping X
          X_pad= np.expand_dims(X_pad, axis=-1)
In [53]:
         #classes imbalance
          np.unique(y, return_counts=True)
Out[53]: (array([0, 1]), array([919, 703], dtype=int64))
In [54]:
          # to categorical
          y_cat = to_categorical(y)
          y_cat.shape
Out[54]: (1622, 2)
In [55]:
          np.unique(y_cat, return_counts=True)
Out[55]: (array([0., 1.], dtype=float32), array([1622, 1622], dtype=int64))
In [56]:
          #Splitting data
          X train, X test, y train, y test= train_test_split(X pad, y_cat,
          test_size=0.2, random_state=42)
In [14]:
          #baseline model
          sum = np.sum(y train, axis=0)
          predicted_category = np.argmax(sum_)
          good_prediction = np.sum(y_test, axis=0)[predicted_category]
          baseline_result = good_prediction/len(y_test)
          print(f'Baseline accuracy: {baseline_result}')
         Baseline accuracy: 0.5692307692307692
In [15]:
          model = Sequential()
          model.add(layers.Masking(mask_value=-1., input_shape=(140,1)))
          model.add(layers.GRU(units=20, activation="tanh", input_shape=(140,1)))
          model.add(layers.Dense(50, activation='relu'))
          model.add(layers.Dropout(0.3))
          model.add(layers.Dense(2, activation='softmax'))
```

model.summary()

Model: "sequential"

Layer (type)	Output Shape	Param #
masking (Masking)	(None, 140, 1)	0
gru (GRU)	(None, 20)	1380
dense (Dense)	(None, 50)	1050
dropout (Dropout)	(None, 50)	0
dense_1 (Dense)	(None, 2)	102
_	============	==========
Total params: 2,532		
Trainable params: 2,532		

Non-trainable params: 0

```
In [16]:
```

```
Epoch 1/100
45 - val loss: 0.6823 - val accuracy: 0.6115
Epoch 2/100
4 - val loss: 0.6751 - val accuracy: 0.6115
Epoch 3/100
1 - val loss: 0.6746 - val accuracy: 0.6115
Epoch 4/100
9 - val loss: 0.6701 - val accuracy: 0.6115
Epoch 5/100
7 - val loss: 0.6691 - val accuracy: 0.6115
Epoch 6/100
2 - val loss: 0.6656 - val accuracy: 0.6115
Epoch 7/100
6 - val loss: 0.6663 - val accuracy: 0.6115
```

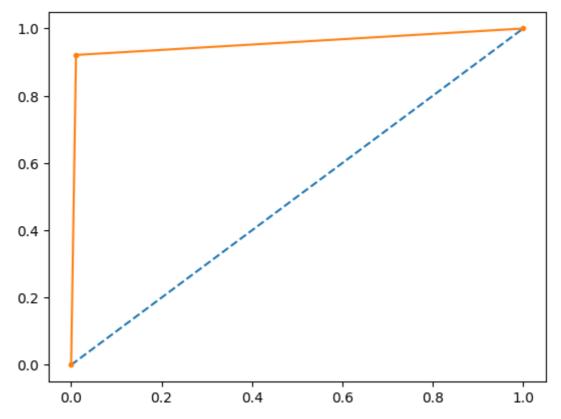
```
Epoch 8/100
6 - val_loss: 0.6652 - val_accuracy: 0.6231
Epoch 9/100
5 - val_loss: 0.6636 - val_accuracy: 0.6038
Epoch 10/100
6 - val_loss: 0.6642 - val_accuracy: 0.5885
Epoch 11/100
9 - val_loss: 0.6639 - val_accuracy: 0.5885
Epoch 12/100
7 - val loss: 0.6595 - val accuracy: 0.5692
Epoch 13/100
3 - val loss: 0.6571 - val accuracy: 0.5846
Epoch 14/100
6 - val_loss: 0.6528 - val_accuracy: 0.6154
Epoch 15/100
1 - val_loss: 0.6532 - val_accuracy: 0.5808
Epoch 16/100
5 - val_loss: 0.6541 - val_accuracy: 0.5769
Epoch 17/100
0 - val_loss: 0.6500 - val_accuracy: 0.5577
Epoch 18/100
3 - val_loss: 0.6506 - val_accuracy: 0.5731
Epoch 19/100
1 - val_loss: 0.6525 - val_accuracy: 0.5731
Epoch 20/100
8 - val_loss: 0.6496 - val_accuracy: 0.5808
Epoch 21/100
1 - val_loss: 0.6478 - val_accuracy: 0.5885
Epoch 22/100
9 - val_loss: 0.6466 - val_accuracy: 0.5769
Epoch 23/100
9 - val loss: 0.6396 - val accuracy: 0.5615
4 - val loss: 0.6402 - val accuracy: 0.5731
Epoch 25/100
2 - val loss: 0.6359 - val accuracy: 0.5769
7 - val loss: 0.6397 - val accuracy: 0.5923
Epoch 27/100
2 - val loss: 0.6402 - val accuracy: 0.5731
6 - val loss: 0.6361 - val accuracy: 0.5962
Epoch 29/100
8 - val_loss: 0.6368 - val_accuracy: 0.6038
5 - val loss: 0.6473 - val accuracy: 0.5846
```

```
Out[16]: [0.6418614387512207, 0.5907692313194275]
In [17]:
        testLoss, testAcc = model.evaluate(X_test, y_test)
         print('Test accuracy:', testAcc)
         print('Test loss:', testLoss)
        Test accuracy: 0.5907692313194275
        Test loss: 0.6418614387512207
In [18]:
        #predict
        y_pred = model.predict(X_test)
In [19]:
        #predicted classes
         pd.DataFrame(y_pred).mean().sort_values()
            0.442444
Out[19]: 1
            0.557556
        dtype: float32
In [20]:
        #actual classes count
         pd.value_counts(y,normalize=True, ascending=True)
Out[20]: 1
            0.433416
            0.566584
        dtype: float64
In [27]:
        # confusion matrix
         y_pred_classes = np.argmax(y_pred, axis=1)
         y_true_classes = np.argmax(y_test, axis=1)
         confusion_matrix(y_true_classes, y_pred_classes)
Out[27]: array([[183,
             [ 11, 129]], dtype=int64)
In [35]:
        # Precision, Recall, F1-score, accuracy, and AUC
         from sklearn.metrics import classification_report
         print(classification_report(y_true_classes, y_pred_classes))
         # AUC
         from sklearn.metrics import roc_auc_score
         from sklearn.metrics import roc_curve
         # calculate AUC
         auc = roc_auc_score(y_true_classes, y_pred_classes)
```

```
print('AUC: %.3f' % auc)
# calculate roc curve
fpr, tpr, thresholds = roc_curve(y_true_classes, y_pred_classes)
# plot no skill
plt.plot([0, 1], [0, 1], linestyle='--')
# plot the roc curve for the model
plt.plot(fpr, tpr, marker='.')
# show the plot
plt.show()
```

	precision	recall	f1-score	support
0	0.94	0.99	0.97	185
1	0.98	0.92	0.95	140
accuracy			0.96	325
macro avg	0.96	0.96	0.96	325
weighted avg	0.96	0.96	0.96	325

AUC: 0.955



```
# plot RNN model
from tensorflow.keras.utils import plot_model
plot_model(model, to_file='model_plot.png', show_shapes=True,
show_layer_names=True)
```

('You must install pydot (`pip install pydot`) and install graphviz (see instruction s at https://graphviz.gitlab.io/download/) ', 'for plot_model_model_to_dot to wor k.')

```
In [21]:
```

```
model = Sequential()
model.add(layers.Masking(mask_value=-1., input_shape=(140,1)))
model.add(layers.LSTM(units=20, activation="tanh", input_shape=(140,1)))
model.add(layers.Dense(50, activation='relu'))
model.add(layers.Dropout(0.3))
model.add(layers.Dense(2, activation='softmax'))
model.summary()
```

Model: "sequential_1"

Layer (type)	Output Shape	Param #
masking_1 (Masking)	(None, 140, 1)	0
lstm (LSTM)	(None, 20)	1760
dense_2 (Dense)	(None, 50)	1050
dropout_1 (Dropout)	(None, 50)	0
dense_3 (Dense)	(None, 2)	102
Total names: 2 012		

Total params: 2,912 Trainable params: 2,912 Non-trainable params: 0

```
1 - val_loss: 0.6619 - val_accuracy: 0.6308
Epoch 5/100
9 - val_loss: 0.6586 - val_accuracy: 0.6538
Epoch 6/100
5 - val_loss: 0.6522 - val_accuracy: 0.6346
Epoch 7/100
6 - val_loss: 0.6420 - val_accuracy: 0.6500
Epoch 8/100
2 - val_loss: 0.6339 - val_accuracy: 0.6462
Epoch 9/100
5 - val loss: 0.6134 - val accuracy: 0.6846
Epoch 10/100
6 - val_loss: 0.5795 - val_accuracy: 0.7154
Epoch 11/100
0 - val_loss: 0.4300 - val_accuracy: 0.8808
Epoch 12/100
7 - val_loss: 0.4164 - val_accuracy: 0.8577
Epoch 13/100
6 - val_loss: 0.2728 - val_accuracy: 0.9385
Epoch 14/100
3 - val_loss: 0.2470 - val_accuracy: 0.9462
Epoch 15/100
7 - val_loss: 0.2191 - val_accuracy: 0.9500
Epoch 16/100
6 - val_loss: 0.3418 - val_accuracy: 0.8538
Epoch 17/100
9 - val_loss: 0.2088 - val_accuracy: 0.9423
Epoch 18/100
4 - val_loss: 0.1850 - val_accuracy: 0.9462
Epoch 19/100
4 - val_loss: 0.1678 - val_accuracy: 0.9500
4 - val loss: 0.1745 - val accuracy: 0.9615
Epoch 21/100
2 - val loss: 0.1573 - val accuracy: 0.9577
Epoch 22/100
3 - val loss: 0.1428 - val accuracy: 0.9654
Epoch 23/100
0 - val loss: 0.1942 - val accuracy: 0.9538
1 - val loss: 0.1299 - val accuracy: 0.9654
Epoch 25/100
3 - val loss: 0.1993 - val accuracy: 0.9462
0 - val_loss: 0.1218 - val_accuracy: 0.9692
Epoch 27/100
```

```
Epoch 28/100
    0 - val_loss: 0.1375 - val_accuracy: 0.9577
    Epoch 29/100
    1 - val_loss: 0.1770 - val_accuracy: 0.9423
    Epoch 30/100
    9 - val_loss: 0.1183 - val_accuracy: 0.9731
    Epoch 31/100
    9 - val_loss: 0.1882 - val_accuracy: 0.9308
    Epoch 32/100
    8 - val loss: 0.1059 - val accuracy: 0.9731
    Epoch 33/100
    7 - val_loss: 0.1426 - val_accuracy: 0.9577
    Epoch 34/100
    7 - val_loss: 0.1014 - val_accuracy: 0.9769
    Epoch 35/100
    6 - val_loss: 0.0989 - val_accuracy: 0.9769
    Epoch 36/100
    6 - val_loss: 0.0880 - val_accuracy: 0.9846
    Epoch 37/100
    7 - val_loss: 0.0852 - val_accuracy: 0.9808
    Epoch 38/100
    6 - val_loss: 0.0857 - val_accuracy: 0.9808
    Epoch 39/100
    6 - val_loss: 0.1168 - val_accuracy: 0.9615
    Epoch 40/100
    6 - val_loss: 0.0873 - val_accuracy: 0.9846
    Epoch 41/100
    4 - val_loss: 0.0857 - val_accuracy: 0.9808
    Epoch 42/100
    6 - val_loss: 0.1746 - val_accuracy: 0.9500
    600
Out[22]: [0.14067725837230682, 0.9599999785423279]
In [23]:
    testLoss, testAcc = model.evaluate(X_test, y_test)
    print('Test accuracy:', testAcc)
     print('Test loss:', testLoss)
    Test accuracy: 0.9599999785423279
    Test loss: 0.14067725837230682
In [26]:
    #predict
    y_pred = model.predict(X_test)
     #predicted classes
```

0 - val_loss: 0.1253 - val_accuracy: 0.9654

```
print(pd.DataFrame(y_pred).mean().sort_values())
          print()
          #actual classes count
          print(pd.value_counts(y,normalize=True, ascending=True))
             0.422169
             0.577831
         dtype: float32
             0.433416
             0.566584
         dtype: float64
In [30]:
         # confusion matrix
          y_pred_classes = np.argmax(y_pred, axis=1)
          y_true_classes = np.argmax(y_test, axis=1)
          confusion_matrix(y_true_classes, y_pred_classes)
Out[30]: array([[183,
                      2],
               [ 11, 129]], dtype=int64)
In [34]:
          # Precision, Recall, F1-score, accuracy, and AUC
          from sklearn.metrics import classification_report
          print(classification_report(y_true_classes, y_pred_classes))
          # AUC
          from sklearn.metrics import roc_auc_score
          from sklearn.metrics import roc_curve
          # calculate AUC
          auc = roc_auc_score(y_true_classes, y_pred_classes)
          print('AUC: %.3f' % auc)
          # calculate roc curve
          fpr, tpr, thresholds = roc_curve(y_true_classes, y_pred_classes)
          # plot no skill
          plt.plot([0, 1], [0, 1], linestyle='--')
          # plot the roc curve for the model
          plt.plot(fpr, tpr, marker='.')
          # show the plot
          plt.show()
                      precision recall f1-score
                                                   support
                          0.94
                                   0.99
                                            0.97
                   0
                                                       185
```

1

accuracy

0.98

0.92

0.95

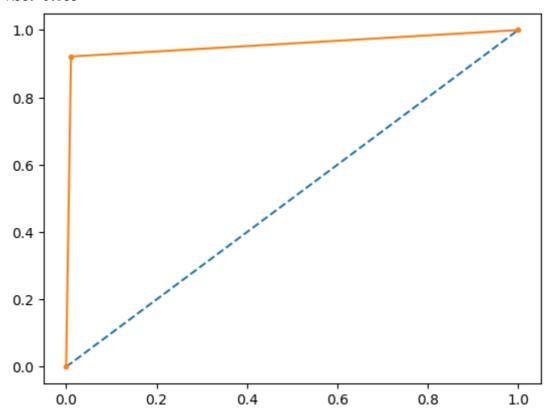
0.96

140

325

macro avg 0.96 0.96 0.96 325 weighted avg 0.96 0.96 0.96 325

AUC: 0.955



```
# plot RNN model
from tensorflow.keras.utils import plot_model
plot_model(model, to_file='model_plot.png', show_shapes=True,
show_layer_names=True)
```

('You must install pydot (`pip install pydot`) and install graphviz (see instruction s at https://graphviz.gitlab.io/download/) ', 'for plot_model/model_to_dot to wor k.')

From both models (LSMT and GRU):

- The accuracy of the model on the test set is higher in the LSTM model than in the GRU model.
- The loss of the model on the test set is lower in the LSTM model than in the GRU model.

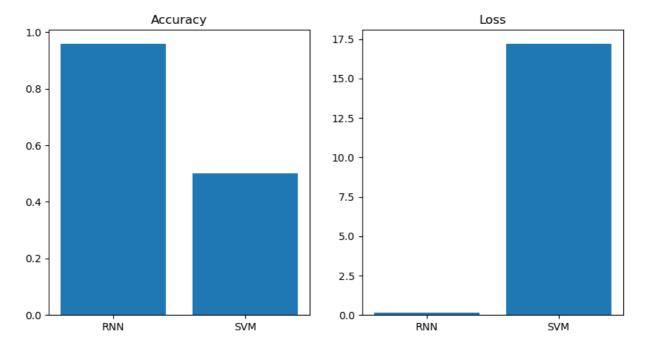
```
In [63]: # Comparing the RNN with SVM
from sklearn.svm import SVC
from sklearn.metrics import accuracy_score

X = df.iloc[:, 1:]
# convert X to numpy array
# X = X.to_numpy().tolist()
# X[0][0]

y = df.iloc[:, 0]
```

```
# y = y.to_numpy().tolist()
# y[0]
# split data into train and test
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2,
random_state=42)
# SVM
svm = SVC(kernel='linear', C=1.0, random_state=0)
svm.fit(X_train, y_train)
y_pred = svm.predict(X_test)
print('Accuracy: %.2f' % accuracy_score(y_test, y_pred))
# calculate loss in SVM
from sklearn.metrics import log_loss
svm_loss = log_loss(y_test, y_pred)
print('SVM loss: %.2f' % svm_loss)
# plotting accuracy and loss between RNN and SVM
plt.figure(figsize=(10,5))
plt.subplot(1,2,1)
plt.bar(['RNN','SVM'],[testAcc,accuracy_score(y_test, y_pred)])
plt.title('Accuracy')
plt.subplot(1,2,2)
plt.bar(['RNN','SVM'],[testLoss,svm_loss])
plt.title('Loss')
plt.show()
```

Accuracy: 0.50 SVM loss: 17.22



From the above graphs, we can see that the RNN is better than the SVM. The RNN has a higher accuracy and lower loss than the SVM. The RNN is also faster than the SVM.