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🚢 Titanic - Machine Learning from Disaster: A Survival Prediction  
using Various Machine Learning

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# 1. | Introduction 🙌



## 🤖 Problem Statement

The sinking of the Titanic is one of the most infamous shipwrecks in history. On **April 15, 1912**, during her maiden voyage, the widely considered “unsinkable” RMS Titanic sank after colliding with an iceberg. Unfortunately, there weren’t enough lifeboats for everyone onboard, resulting in the death of **1502 out of 2224** passengers and crew. While there was some element of luck involved in surviving, it seems some groups of people were more likely to survive than others. In this notebook, we are going to build a **predictive model** that answers the question: “what sorts of people were more likely to survive?” using passenger data (ie name, age, gender, socio-economic class, etc).

## 😬 Dataset Problems

This dataset is taken from the [Kaggle Website](#). This dataset contains **passenger data (ie name, age, gender, socio-economic class, etc)** of whether the passenger has survived. **With the help of Machine learning models** based on the passenger information provided in the dataset. The **variables that most influence** a survival chances will also be **explored more deeply** in this notebook.

## 📌 Notebook Objectives

This notebook **aims** to:

- Perform dataset exploration using various types of data visualization.
- Build machine learning model that can predict survival.
- Export prediction result on test data into files.
- Save/dump the complete machine learning pipeline for later usage.
- Perform prediction on new example data given and export the prediction result.



## Machine Learning Model

The **models** used in this notebook:

1. **Logistic Regression**,
2. **K-Nearest Neighbour (KNN)**,
3. **Support Vector Machine (SVM)**,
4. **Gaussian Naive Bayes**,
5. **Decision Tree**,
6. **Random Forest**,
7. **Extra Tree Classifier**,
8. **Gradient Boosting**, and
9. **AdaBoost**.

## 2. | Installing and Importing Libraries

**Installing and Importing libraries** that will be used in this notebook.

```
In [1]: ## --- Installing Libraries ---
#!pip install ydata-profiling
#!pip install highlight-text
#!pip install Pillow
```

```
In [2]: # --- Importing Libraries ---
from IPython.display import display, HTML, Javascript
import numpy as np
import pandas as pd
import ydata_profiling
import matplotlib.pyplot as plt
import matplotlib.patches as mpatches
%matplotlib inline
import seaborn as sns
import warnings
import os
import yellowbrick
import joblib
```

```

from ydata_profiling import ProfileReport
#from pywaffle import Waffle
from statsmodels.graphics.gofplots import qqplot
from PIL import Image
from highlight_text import fig_text
from sklearn.pipeline import Pipeline
from sklearn.compose import ColumnTransformer
from sklearn.preprocessing import RobustScaler, OneHotEncoder
from sklearn.model_selection import train_test_split, GridSearchCV
from sklearn.linear_model import LogisticRegression
from sklearn.neighbors import KNeighborsClassifier
from sklearn.svm import SVC
from sklearn.naive_bayes import GaussianNB
from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import RandomForestClassifier, GradientBoostingClassifier, Ada
from sklearn.metrics import classification_report, accuracy_score
from yellowbrick.classifier import PrecisionRecallCurve, ROCAUC, ConfusionMatrix
from yellowbrick.model_selection import LearningCurve, FeatureImportances
from yellowbrick.contrib.wrapper import wrap
from yellowbrick.style import set_palette
warnings.filterwarnings("ignore")

```

### 3. | Reading Dataset 🕶

After importing libraries, **the dataset that will be used will be imported.**

```

In [3]: # --- Importing Dataset ---
df = pd.read_csv("titanic.csv")

# --- Reading Train Dataset ---
class Color:
    # Define color codes
    start = '\033[91m'
    end = '\033[0m'
    color = '\033[94m'

# Create an instance of the Color class
clr = Color()

# Reading Train Dataset
print(clr.start + '.: Imported Dataset :.' + clr.end)
print(clr.color + '*' * 23)
styled_df = df.head(10).reset_index(drop=True).style.background_gradient(cmap='Blue')
styled_df

.: Imported Dataset :.
*****

```

Out[3]:

	PassengerId	Pclass	Name	Sex	Age	SibSp	Parch	Ticket	Fare
0	1	3	Braund, Mr. Owen Harris	male	22.000000	1	0	A/5 21171	7.250000
1	2	1	Cumings, Mrs. John Bradley (Florence Briggs Thayer)	female	38.000000	1	0	PC 17599	71.283300
2	3	3	Heikkinen, Miss. Laina	female	26.000000	0	0	STON/O2. 3101282	7.925000
3	4	1	Futrelle, Mrs. Jacques Heath (Lily May Peel)	female	35.000000	1	0	113803	53.100000
4	5	3	Allen, Mr. William Henry	male	35.000000	0	0	373450	8.050000
5	6	3	Moran, Mr. James	male	nan	0	0	330877	8.458300
6	7	1	McCarthy, Mr. Timothy J	male	54.000000	0	0	17463	51.862500
7	8	3	Palsson, Master. Gosta Leonard	male	2.000000	3	1	349909	21.075000
8	9	3	Johnson, Mrs. Oscar W (Elisabeth Vilhelmina Berg)	female	27.000000	0	2	347742	11.133300
9	10	2	Nasser, Mrs. Nicholas (Adele Achem)	female	14.000000	1	0	237736	30.070800

## Dataset Description

The following is the **structure of the dataset**.

Variable Name	Description	Sample Data
PassengerId	Passenger's ID (in years)	63; 37; ...
pclass	Ticket Class	1 = 1st, 2 = 2nd, 3 = 3rd
Name	Name of the Passenger	Braund, Mr. Owen Harris
Sex	Gender of patient 0 = male 1 = female	1; 0; ...
Age	Passengers age (in years)	63; 37; ...
SibSp	No. of siblings / spouses aboard the Titanic	3; 1; 2; ...
Parch	No. of parents / children aboard the Titanic	4; 3; ...
thalach	Maximum heart rate achieved	150; 187; ...
Ticket	Ticket number	2649; 349909; ...
Fare	Passenger fare	31.3875; 263; ...
Cabin	Cabin number	C85;E46;; ...
Survival	Survival of Passenger	0 =No , 1=Yes

## 4. | Initial Dataset Exploration

This section will focused on **initial data exploration on the dataset** with [Pandas Profiling](#) before pre-processing performed. In addition, **variables correlation** will be examined as well.

```
In [4]: # --- Dataset Report ---
ProfileReport(df,
              title="Titanic Survival Prediction",
              minimal=True,
              progress_bar=False,
              samples=None,
              interactions=None,
              explorative=True,
              dark_mode=True,
              notebook={'iframe': {'height': '600px'},
                       'html': {'style': {'primary_color': clr}},
                       'missing_diagrams': {'heatmap': False, 'dendrogram': False}}
              ).to_notebook_iframe()
```

# Overview

## Dataset statistics

Number of variables	12
Number of observations	891
Missing cells	866
Missing cells (%)	8.1%
Total size in memory	315.0 KiB
Average record size in memory	362.1 B

## Variable types

Numeric	7
Text	5

## 4. | Data Preprocessing

Since **Name** and **Ticket** columns are not directly related to the survival we can drop them.

```
In [5]: df = df.drop(["Name", "Ticket"], axis=1)
```

In the **Cabin** columns there are **687 (77.1%) missing values** hence removing the cabin column

```
In [6]: df = df.drop("Cabin", axis=1)
```

```
In [7]: df["Embarked"] = df["Embarked"].fillna(df["Embarked"].mode()[0])
```

```
In [8]: df['Age'] = df.groupby(['Pclass', 'Sex'])['Age'].transform(lambda x: x.fillna(x.med
```

```
In [9]: df = df.drop(["PassengerId"],axis=1)
```

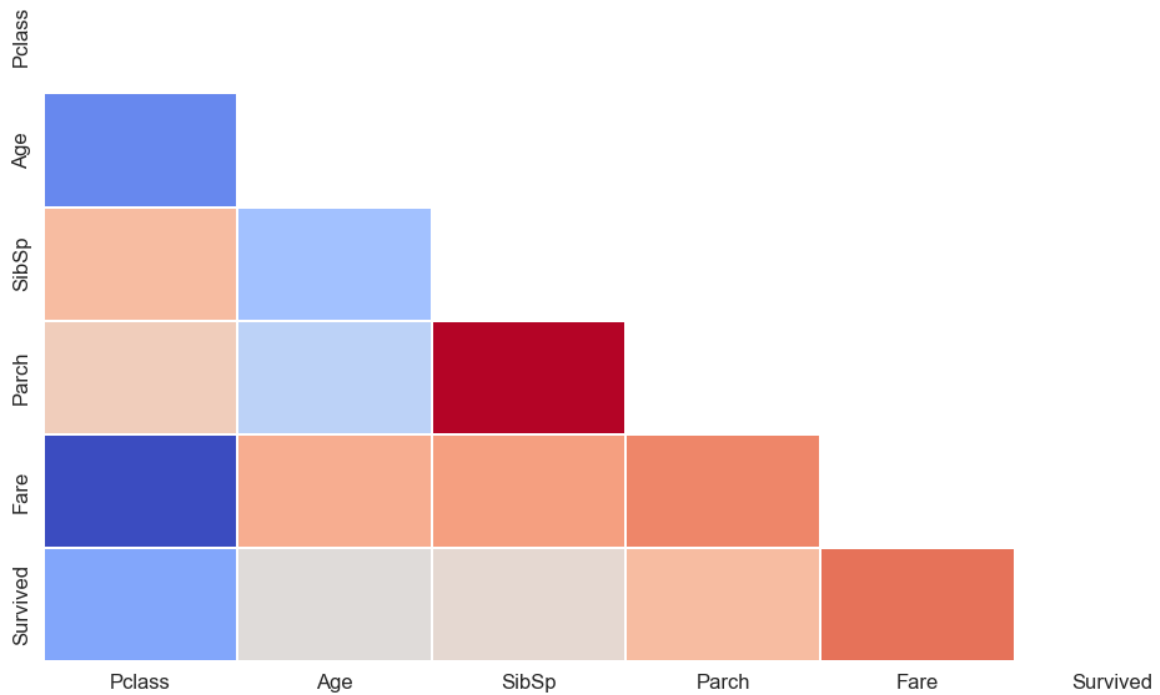
```
In [10]: corr = df.corr(numeric_only=True)
```

```
In [11]: # --- Correlation Map Variables ---
colors = ['red','blue','green']
suptitle = dict(x=0.1, y=1.01, fontsize=20, weight='heavy', ha='left', va='bottom',
title = dict(x=0.1, y=0.98, fontsize=14, weight='normal', ha='left', va='bottom', f
xy_label = dict(size=12)
highlight_textprops = [{'weight':'bold', 'color': colors[0]}, {'weight':'bold', 'co

# --- Correlation Map (Heatmap) ---
mask = np.triu(np.ones_like(corr, dtype=bool))
fig, ax = plt.subplots(figsize=(10, 8))
sns.heatmap(corr, mask=mask, annot=True, cmap='coolwarm', linewidths=0.2, cbar=Fa
yticks, ylabels = plt.yticks()
xticks, xlabels = plt.xticks()
ax.set_xticklabels(xlabels, rotation=0, **xy_label)
ax.set_yticklabels(ylabels, **xy_label)
ax.grid(False)
fig_text(s='Numerical Variables Correlation Map', **suptitle)
fig_text(s='<Chest pain type, max heart rate, and slope> positively correlate with
plt.tight_layout(rect=[0, 0.08, 1, 1.01])
plt.show();
```

## Numerical Variables Correlation Map

Chest pain type, max heart rate, and slope positively correlate with target variables.



From **dataset report** and **correlation matrix**, it can be **concluded** that:

- There are **no missing values** detected in the dataset. In addition, it also can be seen that **the number of categorical columns is more than the numerical columns**.
- As can be seen from the profiling report, **the number of male passengers is greater than female passengers**. In addition, **passenger class 3 is higher than other types**. **Most of the passengers in the dataset had age that was less than 40 years**.
- Furthermore, **passengers with 0 Pclass are more numerous than those with other Pclass**.
- **The Fare column has a left skewed distribution** based on the histogram and skewness value.
- **The mean age of the passengers in the dataset was 29 years old**, with the most senior patient being 80 years old and the youngest being less than a year old.



This section will perform some **EDA** to get more insights about dataset.

```
In [12]: # --- EDA 3 Dataframes ---
df_eda3 = df[['Pclass', 'Survived']]
df_eda3 = pd.DataFrame(df_eda3.groupby(['Pclass', 'Survived']).size().reset_index(name='total'))
df_eda3.loc[len(df_eda3.index)] = [1, 2, 0]
df_eda3_0 = df_eda3.query('Pclass == 1').drop('Pclass', axis=1)
df_eda3_1 = df_eda3.query('Pclass == 2').drop('Pclass', axis=1)
df_eda3_2 = df_eda3.query('Pclass == 3').drop('Pclass', axis=1)

# --- EDA 3 Variables ---
total_list = [df_eda3_0['total'], df_eda3_1['total'], df_eda3_2['total']]
suptitle = dict(x=0.5, y=0.94, fontsize=14, weight='heavy', ha='center', va='center')
exp_text = dict(x=0.5, y=-0.08, fontsize=8, weight='normal', ha='center', va='center')
highlight_explanation = [{'weight': 'bold', 'color': 'red'}, {'weight': 'bold', 'color': 'green'}]
survived_0 = mpatches.Patch(color='red', label='Not_Survived')
survived_1 = mpatches.Patch(color='green', label='Survived')

def my_autopct(pct):
    return f'{pct:.2f}%' if pct != 0 else ''

# --- EDA 3 Functions ---
def display_eda3(subplot_num, pclass_type, total, colors, start_angle):
    centre = plt.Circle((0, 0), 0.85, fc='white', edgecolor='black', linewidth=0.5)
    total_passengers = total.sum()

    plt.subplot(1, 3, subplot_num)
    plt.tight_layout(rect=[0, 0, 1, 1.01])
    plt.pie(total, colors=colors, autopct=my_autopct, pctdistance=0.65, startangle=start_angle)
    plt.text(0, 0.08, f"Class {pclass_type}", weight='bold', ha='center', fontsize=12)
    plt.text(0, -0.08, f"{total_passengers} passengers", ha='center', fontsize=8)
    fig = plt.gcf()
    fig.gca().add_artist(centre)

# --- Display EDA 3 ---
plt.figure(figsize=(9, 4))
for idx, total in enumerate(total_list):
    display_eda3(idx + 1, idx, total, ['red', 'green', 'blue'], 90 * idx)
    if idx == 1:
        plt.legend(handles=[survived_0, survived_1], loc='upper center', bbox_to_anchor=(0.5, 1.1))

plt.suptitle("Survival Distribution by Passenger Class", **suptitle)
plt.title("Explanation: Pclass 3 has the highest survival rate, Class 2 is moderate")
plt.show()
```



## 6. | Data Preprocessing

This section will **prepare the dataset** before building the machine learning models.

### 6.1 | Features Separating and Splitting

In this section, the 'Survived' (dependent) column will be separated from independent columns. Also, the dataset will be splitted into 90:10 ratio (90% training and 10% testing).

```
In [13]: # --- Separating Dependent Features ---
x = df.drop(['Survived'], axis=1)
y = df['Survived']

# --- Splitting Dataset ---
x_train, x_test, y_train, y_test = train_test_split(x, y, test_size=0.1, random_sta
```

### 6.2 | Processing Pipeline

This section will create a preprocessing pipeline for numerical and categorical columns and apply them to the x\_train and x\_test data. Not all columns will go through preprocessing. For all numerical columns, scaling will be carried out using a robust scaler since the dataset used is a **small dataset** where the presence of outliers dramatically affects the performance of a model. While for categorical columns with more than two categories, one-hot encoding will be carried out.

```
In [14]: # --- Numerical Pipeline ---
num_column = ['Age', 'Fare']
num_pipeline = Pipeline([
    ('scaling', RobustScaler())
])

# --- Categorical Pipeline ---
cat_column = ['Pclass', 'Sex', 'Embarked']
cat_pipeline = Pipeline([
    ('onehot', OneHotEncoder(drop='first', sparse=False))
```

```

])

# --- Combine Both Pipelines into Transformer ---
preprocessor = ColumnTransformer([
    ('categorical', cat_pipeline, cat_column)
    , ('numerical', num_pipeline, num_column)]
    , remainder='passthrough')

# --- Apply Transformer to Pipeline ---
process_pipeline = Pipeline([
    ('preprocessor', preprocessor)
])

# --- Apply to Dataframe ---
x_train_process = process_pipeline.fit_transform(x_train)
x_test_process = process_pipeline.fit_transform(x_test)

```

## 7. | Machine Learning Model Implementation

This section will **implement various machine learning models** as mentioned in Introduction section. In addition, explanation for each models also will be discussed.

```

In [15]: # --- Functions: Model Fitting and Performance Evaluation ---
color_yb = sns.color_palette("Paired")
color_line = 'red'
color = 'red'
def fit_ml_models(algo, algo_param, algo_name):

    # --- Algorithm Pipeline ---
    algo = Pipeline(['algo', algo])

    # --- Apply Grid Search ---
    model = GridSearchCV(algo, param_grid=algo_param, cv=10, n_jobs=-1, verbose=1)

    # --- Fitting Model ---
    print clr.start+f"..Fitting {algo_name} .."+clr.end)
    fit_model = model.fit(x_train_process, y_train)

    # --- Model Best Parameters ---
    best_params = model.best_params_
    print("\n>> Best Parameters: "+clr.start+f"{best_params}" +clr.end)

    # --- Best & Final Estimators ---
    best_model = model.best_estimator_
    best_estimator = model.best_estimator_.final_estimator
    best_score = round(model.best_score_, 4)
    print(">> Best Score: "+clr.start+"{:.3f}".format(best_score)+clr.end)

    # --- Create Prediction for Train & Test ---
    y_pred_train = model.predict(x_train_process)
    y_pred_test = model.predict(x_test_process)

```

```

# --- Train & Test Accuracy Score ---
acc_score_train = round(accuracy_score(y_pred_train, y_train)*100, 3)
acc_score_test = round(accuracy_score(y_pred_test, y_test)*100, 3)
print("\n"+clr.start+f"... Train and Test Accuracy Score for {algo_name} ..."+clr.end)
print("\t>> Train Accuracy: "+clr.start+"{:.2f}%".format(acc_score_train)+clr.end)
print("\t>> Test Accuracy: "+clr.start+"{:.2f}%".format(acc_score_test)+clr.end)

# --- Classification Report ---
print("\n"+clr.start+f"... Classification Report for {algo_name} ..."+clr.end)
print(classification_report(y_test, y_pred_test))

# --- Figures Settings ---
xy_label = dict(fontweight='bold', fontsize=12)
grid_style = dict(color='black', linestyle='dotted', zorder=1)
title_style = dict(fontsize=14, fontweight='bold')
tick_params = dict(length=3, width=1, color='red')
bar_style = dict(zorder=3, edgecolor='black', linewidth=0.5, alpha=0.85)
set_palette(color_yb)
fig, ((ax1, ax2), (ax3, ax4)) = plt.subplots(2, 2, figsize=(16, 14))

# --- Confusion Matrix ---
conf_matrix = ConfusionMatrix(best_estimator, ax=ax1, cmap='Reds')
conf_matrix.fit(x_train_process, y_train)
conf_matrix.score(x_test_process, y_test)
conf_matrix.finalize()
conf_matrix.ax.set_title('Confusion Matrix\n', **title_style)
conf_matrix.ax.tick_params(axis='both', labelsz=10, bottom='on', left='on', *
for spine in conf_matrix.ax.spines.values(): spine.set_color(color_line)
conf_matrix.ax.set_xlabel('\nPredicted Class', **xy_label)
conf_matrix.ax.set_ylabel('True Class\n', **xy_label)
conf_matrix.ax.xaxis.set_ticklabels(['False', 'True'], rotation=0)
conf_matrix.ax.yaxis.set_ticklabels(['True', 'False'])

# --- ROC AUC ---
logroauc = ROCAUC(best_estimator, classes=['False', 'True'], ax=ax2, colors=co
logroauc.fit(x_train_process, y_train)
logroauc.score(x_test_process, y_test)
logroauc.finalize()
logroauc.ax.set_title('ROC AUC Curve\n', **title_style)
logroauc.ax.tick_params(axis='both', labelsz=10, bottom='on', left='on', **t
logroauc.ax.grid(axis='both', alpha=0.4, **grid_style)
for spine in logroauc.ax.spines.values(): spine.set_color('None')
for spine in ['bottom', 'left']:
    logroauc.ax.spines[spine].set_visible(True)
    logroauc.ax.spines[spine].set_color(color_line)
logroauc.ax.legend(loc='upper center', bbox_to_anchor=(0.5, -0.12), ncol=2, bo
logroauc.ax.set_xlabel('\nFalse Positive Rate', **xy_label)
logroauc.ax.set_ylabel('True Positive Rate\n', **xy_label)

# --- Learning Curve ---
lcurve = LearningCurve(best_estimator, scoring='f1_weighted', ax=ax3, colors=co
lcurve.fit(x_train_process, y_train)
lcurve.finalize()
lcurve.ax.set_title('Learning Curve\n', **title_style)
lcurve.ax.tick_params(axis='both', labelsz=10, bottom='on', left='on', **tick

```

```

lcurve.ax.grid(axis='both', alpha=0.4, **grid_style)
for spine in lcurve.ax.spines.values(): spine.set_color('None')
for spine in ['bottom', 'left']:
    lcurve.ax.spines[spine].set_visible(True)
    lcurve.ax.spines[spine].set_color(color_line)
lcurve.ax.legend(loc='upper center', bbox_to_anchor=(0.5, -0.12), ncol=2, border=0)
lcurve.ax.set_xlabel('\nTraining Instances', **xy_label)
lcurve.ax.set_ylabel('Scores\n', **xy_label)

# --- Feature Importance or Precision Recall Curve ---
try:
    feat_importance = FeatureImportances(best_estimator, labels=columns_list_on)
    feat_importance.fit(x_train_process, y_train)
    feat_importance.finalize()
    feat_importance.ax.set_title('Feature Importances (Top 5 Features)\n', **title_style)
    feat_importance.ax.tick_params(axis='both', labelsize=10, bottom='on', left='on')
    feat_importance.ax.grid(axis='x', alpha=0.4, **grid_style)
    feat_importance.ax.grid(axis='y', alpha=0, **grid_style)
    for spine in feat_importance.ax.spines.values(): spine.set_color('None')
    for spine in ['bottom']:
        feat_importance.ax.spines[spine].set_visible(True)
        feat_importance.ax.spines[spine].set_color(color_line)
    feat_importance.ax.set_xlabel('\nRelative Importance', **xy_label)
    feat_importance.ax.set_ylabel('Features\n', **xy_label)
except:
    prec_curve = PrecisionRecallCurve(best_estimator, ax=ax4, ap_score=True, is_ordinal=False)
    prec_curve.fit(x_train_process, y_train)
    prec_curve.score(x_test_process, y_test)
    prec_curve.finalize()
    prec_curve.ax.set_title('Precision-Recall Curve\n', **title_style)
    prec_curve.ax.tick_params(axis='both', labelsize=10, bottom='on', left='on')
    for spine in prec_curve.ax.spines.values(): spine.set_color('None')
    for spine in ['bottom', 'left']:
        prec_curve.ax.spines[spine].set_visible(True)
        prec_curve.ax.spines[spine].set_color(color_line)
    prec_curve.ax.legend(loc='upper center', bbox_to_anchor=(0.5, -0.12), ncol=2, border=0)
    prec_curve.ax.set_xlabel('\nRecall', **xy_label)
    prec_curve.ax.set_ylabel('Precision\n', **xy_label)

plt.suptitle(f'\n{algo_name} Performance Evaluation Report\n', fontsize=18, fontweight='bold')
plt.gcf().text(0.88, 0.02, 'kaggle.com/caesarmario', style='italic', fontsize=10)
plt.tight_layout();

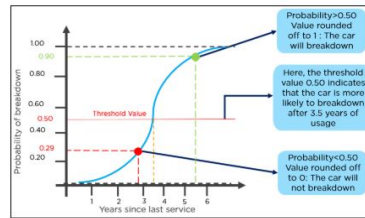
return acc_score_train, acc_score_test, best_score


```

## 7.1 | Logistic Regression

**Logistic regression** is a statistical method that is used for building machine learning models where **the dependent variable is dichotomous: i.e. binary**. Logistic regression is used to describe data and **the relationship between one dependent variable and one or more independent variables**. The independent variables can be nominal, ordinal, or of interval type.

The name "logistic regression" is derived from the concept of the logistic function that it uses. **The logistic function is also known as the sigmoid function.** The value of this logistic function lies between zero and one.



 Logistic Function by Simplilearn

```
In [16]: # --- Logistic Regression Parameters ---
parameter_lr = {"algo_solver": ["lbfgs", "saga", "newton-cg"],
                 , "algo_C": [0.1, 0.2, 0.5, 0.8]}

# --- Logistic Regression Algorithm ---
algo_lr = LogisticRegression(penalty="l2", random_state=42, n_jobs=-1)

# --- Applying Logistic Regression ---
acc_score_train_lr, acc_score_test_lr, best_score_lr = fit_ml_models(algo_lr, param
```

..< Fitting Logistic Regression ..

Fitting 10 folds for each of 12 candidates, totalling 120 fits

>> Best Parameters: {'algo\_C': 0.1, 'algo\_solver': 'lbfgs'}

>> Best Score: 0.808

..< Train and Test Accuracy Score for Logistic Regression ..

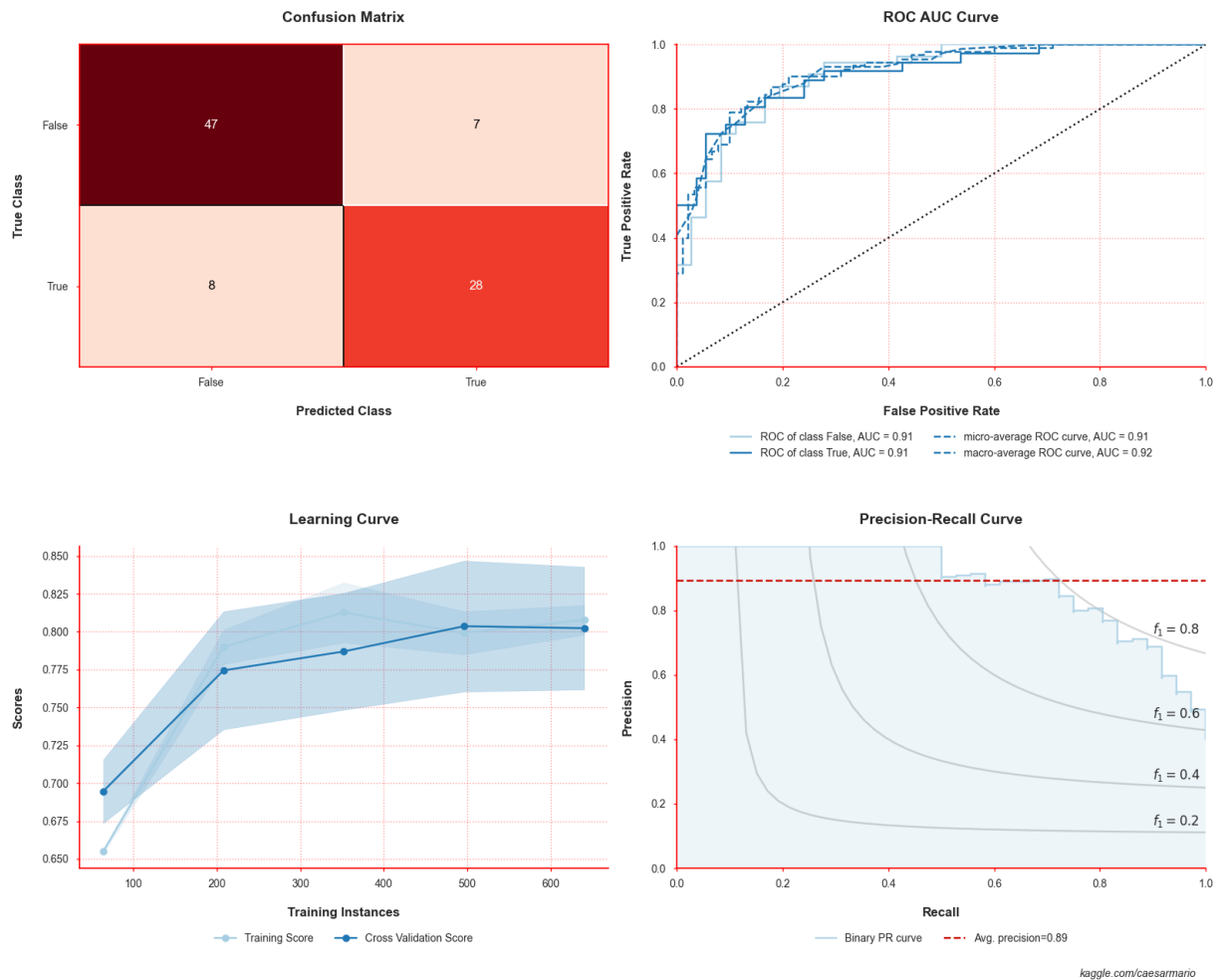
>> Train Accuracy: 81.15%

>> Test Accuracy: 83.33%

..< Classification Report for Logistic Regression ..

	precision	recall	f1-score	support
0	0.85	0.87	0.86	54
1	0.80	0.78	0.79	36
accuracy			0.83	90
macro avg	0.83	0.82	0.83	90
weighted avg	0.83	0.83	0.83	90

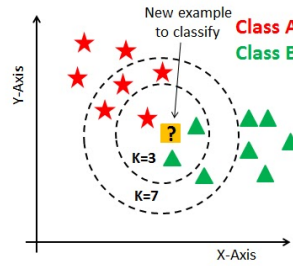
## Logistic Regression Performance Evaluation Report



## 7.2 | K-Nearest Neighbour (KNN)

**The k-nearest neighbors (KNN)** algorithm is a data classification method **for estimating the likelihood that a data point will become a member of one group or another** based on what group the data points nearest to it belong to. The k-nearest neighbor algorithm is a type of supervised machine learning algorithm used **to solve classification and regression problems**.

It's called a **lazy learning algorithm or lazy learner** because it doesn't perform any training when you supply the training data. Instead, it just stores the data during the training time and doesn't perform any calculations. It doesn't build a model until a query is performed on the dataset. This makes KNN ideal for data mining.



 KNN by Kita Informatika

```
In [17]: # --- KNN Parameters ---
parameter_knn = {"algo__n_neighbors": [2, 5, 10, 17]
                  , "algo__leaf_size": [1, 10, 11, 30]}

# --- KNN Algorithm ---
algo_knn = KNeighborsClassifier(n_jobs=-1)

# --- Applying KNN ---
acc_score_train_knn, acc_score_test_knn, best_score_knn = fit_ml_models(algo_knn, p

... Fitting K-Nearest Neighbour (KNN) ...
Fitting 10 folds for each of 16 candidates, totalling 160 fits

>> Best Parameters: {'algo__leaf_size': 10, 'algo__n_neighbors': 10}
>> Best Score: 0.805

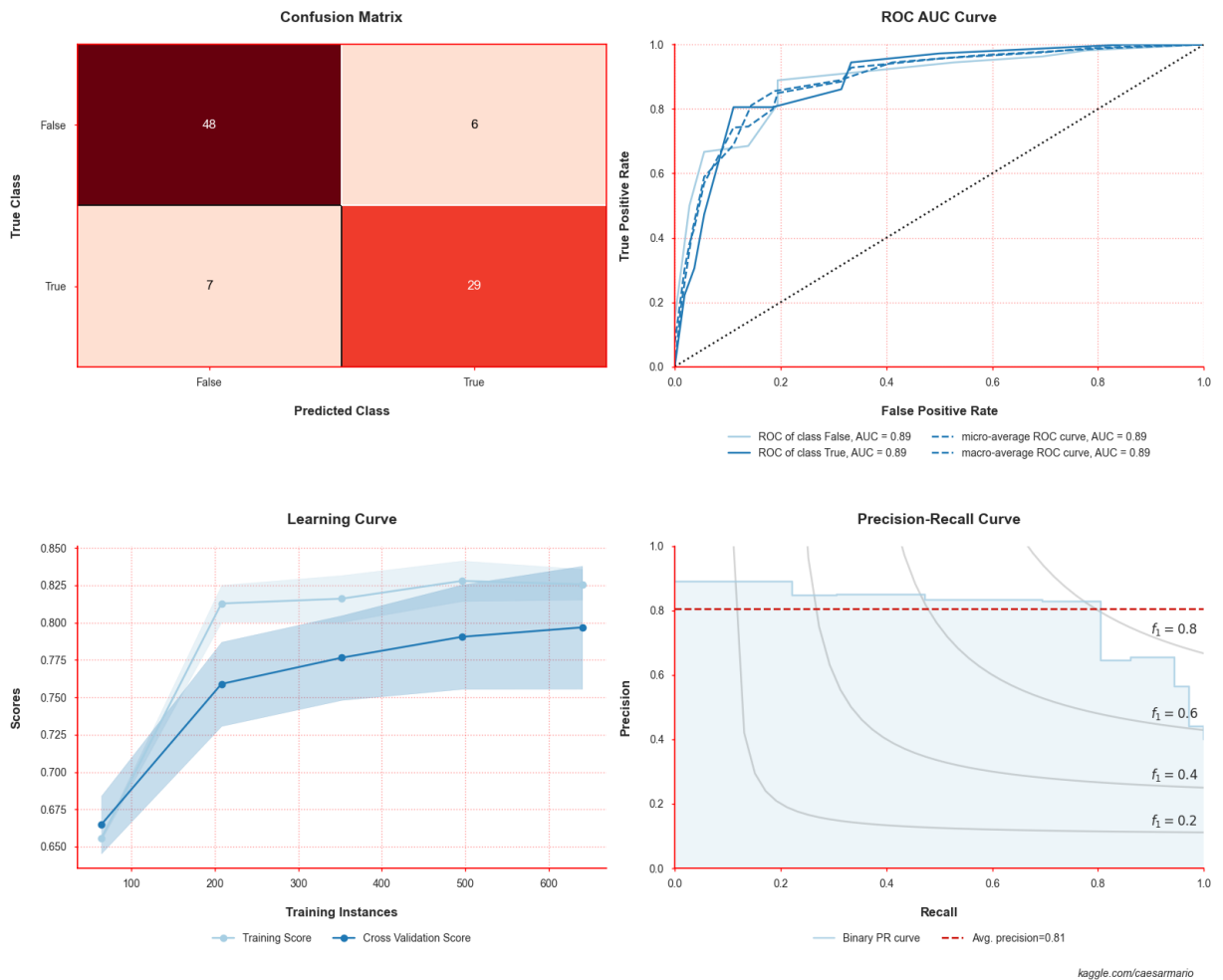
... Train and Test Accuracy Score for K-Nearest Neighbour (KNN) ...
>> Train Accuracy: 82.77%
>> Test Accuracy: 85.56%

... Classification Report for K-Nearest Neighbour (KNN) ...
```

	precision	recall	f1-score	support
0	0.87	0.89	0.88	54
1	0.83	0.81	0.82	36
accuracy			0.86	90
macro avg	0.85	0.85	0.85	90
weighted avg	0.86	0.86	0.86	90



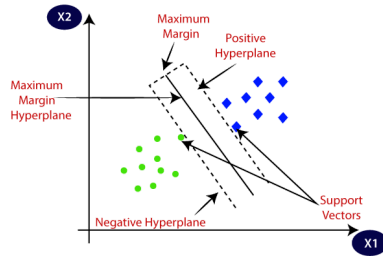
### K-Nearest Neighbour (KNN) Performance Evaluation Report




## 7.3 | Support Vector Machine (SVM)

**Support Vector Machine (SVM)** is one of the most popular Supervised Learning algorithms, which is used for Classification as well as Regression problems. The goal of the SVM algorithm is **to create the best line or decision boundary that can segregate n-dimensional space into classes** so that we can easily put the new data point in the correct category in the future. This best decision boundary is called a hyperplane.

SVM chooses the **extreme points/vectors** that help in creating the hyperplane. These extreme cases are called as support vectors, and hence algorithm is termed as Support Vector Machine.



 SVM by JavaTPoint

```
In [18]: # --- SVM Parameters ---
parameter_svc = [
    {'algo__kernel': ['rbf'], 'algo__gamma': np.arange(0.1, 1, 0.1), 'algo__C': np.
    , {'algo__kernel': ['linear'], 'algo__C': np.arange(0.1, 1, 0.1)}
    , {'algo__kernel': ['poly'], 'algo__degree' : np.arange(1, 10, 1), 'algo__C': n
]

# --- SVM Algorithm ---
algo_svc = SVC(random_state=1, probability=True)

# --- Applying SVM ---
acc_score_train_svc, acc_score_test_svc, best_score_svc = fit_ml_models(algo_svc, p
```

... Fitting Support Vector Machine (SVM) ...

Fitting 10 folds for each of 171 candidates, totalling 1710 fits

>> Best Parameters: {'algo\_\_C': 0.6, 'algo\_\_degree': 2, 'algo\_\_kernel': 'poly'}

>> Best Score: 0.825

... Train and Test Accuracy Score for Support Vector Machine (SVM) ...

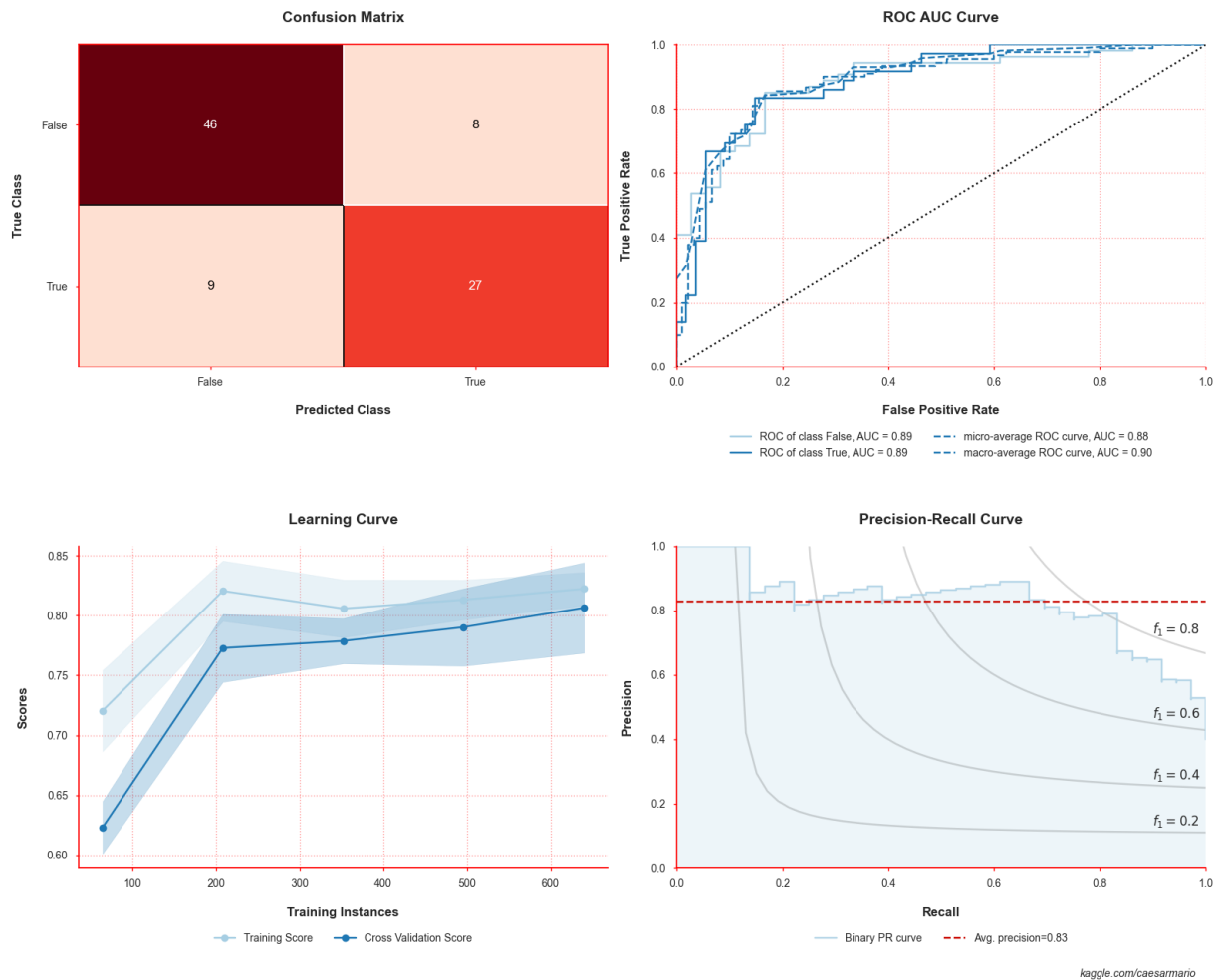
>> Train Accuracy: 83.15%

>> Test Accuracy: 81.11%

... Classification Report for Support Vector Machine (SVM) ...

	precision	recall	f1-score	support
0	0.84	0.85	0.84	54
1	0.77	0.75	0.76	36
accuracy			0.81	90
macro avg	0.80	0.80	0.80	90
weighted avg	0.81	0.81	0.81	90

## Support Vector Machine (SVM) Performance Evaluation Report



## 7.4 | Gaussian Naive Bayes

**Naive Bayes Classifiers** are based on the Bayes Theorem, which **one assumption taken is the strong independence assumptions between the features**. These classifiers assume that the value of a particular feature is independent of the value of any other feature. In a supervised learning situation, Naive Bayes Classifiers are trained very efficiently. Naive Bayes classifiers **need a small training data to estimate the parameters needed for classification**. Naive Bayes Classifiers have simple design and implementation and they can be applied to many real life situations.

**Gaussian Naive Bayes** is a **variant of Naive Bayes that follows Gaussian normal distribution and supports continuous data**. When working with continuous data, an assumption often taken is that the continuous values associated with each class are distributed according to a normal (or Gaussian) distribution.



```
In [19]: # --- Gaussian NB Parameters ---
parameter_gnb = {"algo__var_smoothing": [1e-2, 1e-3, 1e-4, 1e-6]}

# --- Gaussian NB Algorithm ---
algo_gnb = GaussianNB()

# --- Applying Gaussian NB ---
acc_score_train_gnb, acc_score_test_gnb, best_score_gnb = fit_ml_models(algo_gnb, p
```

... Fitting Gaussian Naive Bayes ...

Fitting 10 folds for each of 4 candidates, totalling 40 fits

>> Best Parameters: {'algo\_\_var\_smoothing': 0.01}

>> Best Score: 0.790

... Train and Test Accuracy Score for Gaussian Naive Bayes ...

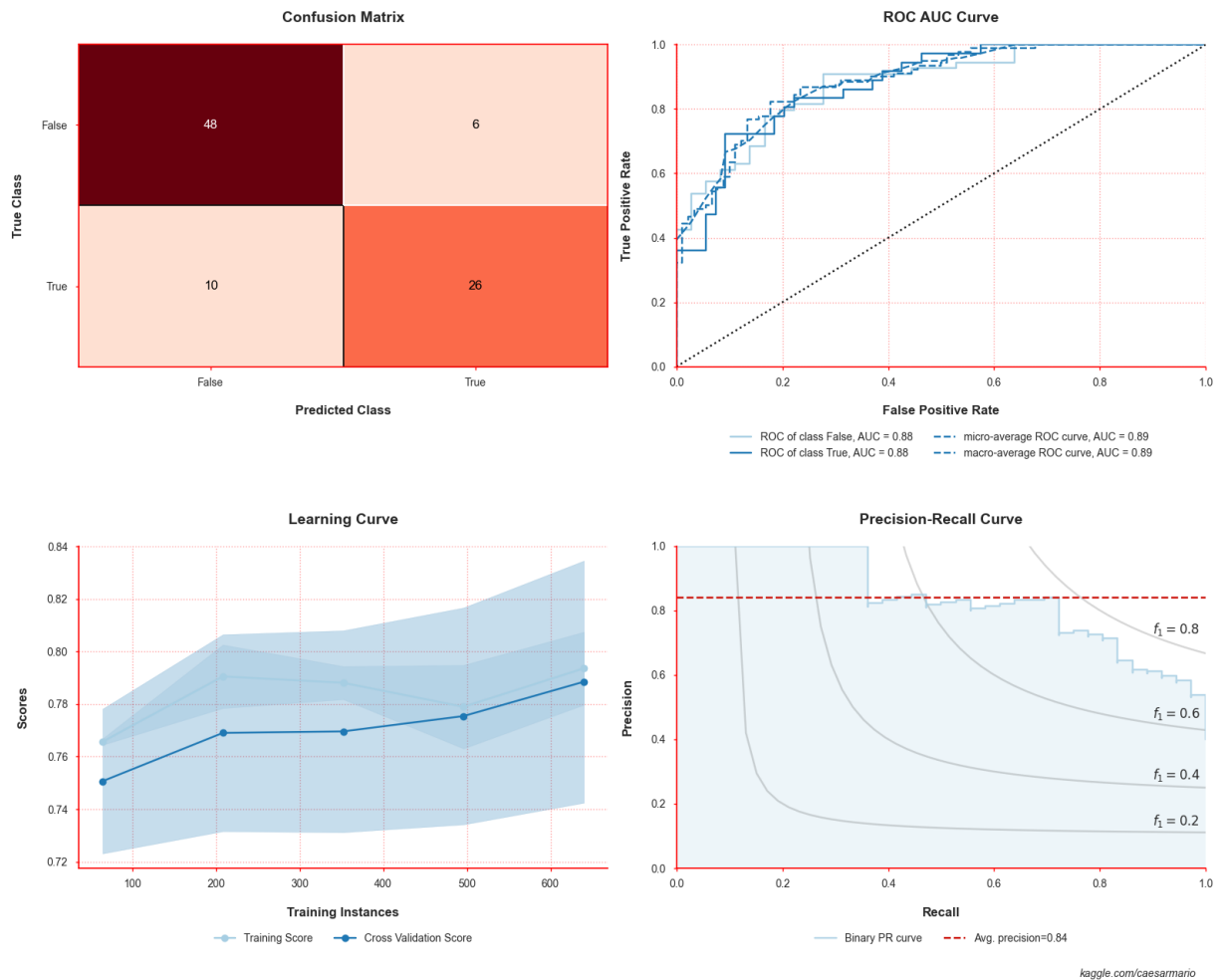
>> Train Accuracy: 80.40%

>> Test Accuracy: 82.22%

... Classification Report for Gaussian Naive Bayes ...

	precision	recall	f1-score	support
0	0.83	0.89	0.86	54
1	0.81	0.72	0.76	36
accuracy			0.82	90
macro avg	0.82	0.81	0.81	90
weighted avg	0.82	0.82	0.82	90

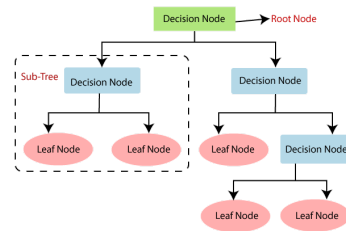
## Gaussian Naive Bayes Performance Evaluation Report



## 7.5 | Decision Tree

**Decision Tree** is a Supervised learning technique that can be used for both classification and Regression problems, but mostly it is preferred for solving Classification problems. It is a tree-structured classifier, where **internal nodes represent the features of a dataset, branches represent the decision rules and each leaf node represents the outcome.**

In a Decision tree, there are **two nodes**, which are the **Decision Node** and **Leaf Node**. Decision nodes are used to make any decision and have multiple branches, whereas Leaf nodes are the output of those decisions and do not contain any further branches.



 Decision Tree by Javatpoint

```
In [20]: # --- Decision Tree Parameters ---
parameter_dt = {"algo__max_depth": [1, 2, 3]}

# --- Decision Tree Algorithm ---
algo_dt = DecisionTreeClassifier(random_state=42)

# --- Applying Decision Tree ---
acc_score_train_dt, acc_score_test_dt, best_score_dt = fit_ml_models(algo_dt, param
```

... Fitting Decision Tree ...

Fitting 10 folds for each of 3 candidates, totalling 30 fits

>> Best Parameters: {'algo\_\_max\_depth': 3}

>> Best Score: 0.815

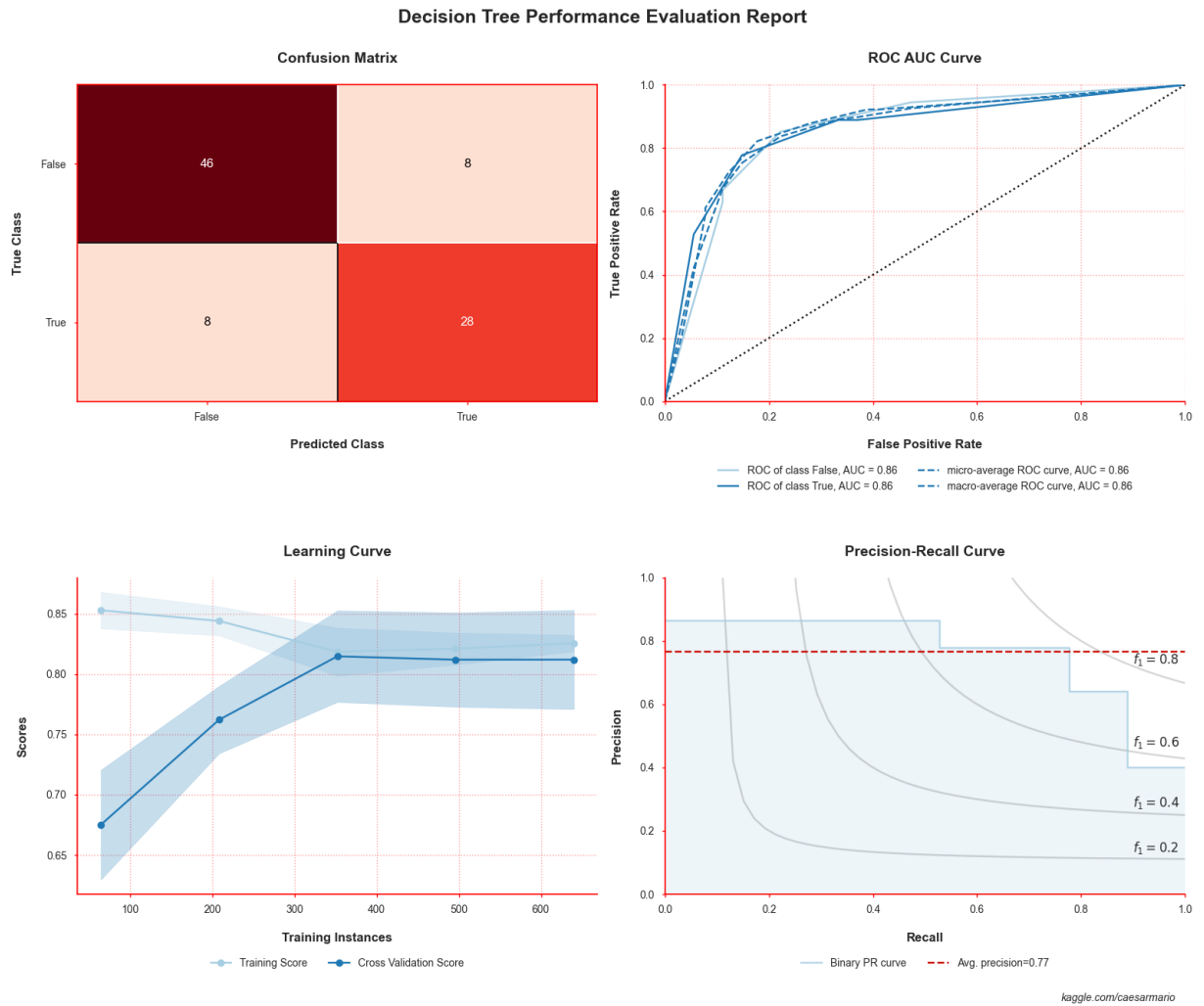
... Train and Test Accuracy Score for Decision Tree ...

>> Train Accuracy: 82.77%

>> Test Accuracy: 82.22%

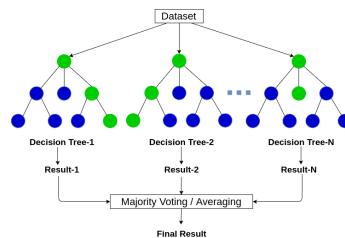
... Classification Report for Decision Tree ...

	precision	recall	f1-score	support
0	0.85	0.85	0.85	54
1	0.78	0.78	0.78	36
accuracy			0.82	90
macro avg	0.81	0.81	0.81	90
weighted avg	0.82	0.82	0.82	90



## 7.6 | Random Forest

**Random Forest** is a tree-based machine learning algorithm that **leverages the power of multiple decision trees for making decisions**. Each individual tree in the random forest spits out a class prediction and the class with the most votes becomes our model's prediction. **A large number of relatively uncorrelated models (trees) operating as a committee will outperform any of the individual constituent models.**



```
In [21]: # --- Random Forest Parameters ---
parameter_rf = {"algo_max_depth": np.arange(1, 10, 1)}

# --- Random Forest Algorithm ---
```

```

algo_rf = RandomForestClassifier(random_state=99, n_jobs=-1)

# --- Applying Random Forest ---
acc_score_train_rf, acc_score_test_rf, best_score_rf = fit_ml_models(algo_rf, param

```

... Fitting Random Forest ...

Fitting 10 folds for each of 9 candidates, totalling 90 fits

>> Best Parameters: {'algo\_max\_depth': 5}

>> Best Score: 0.837

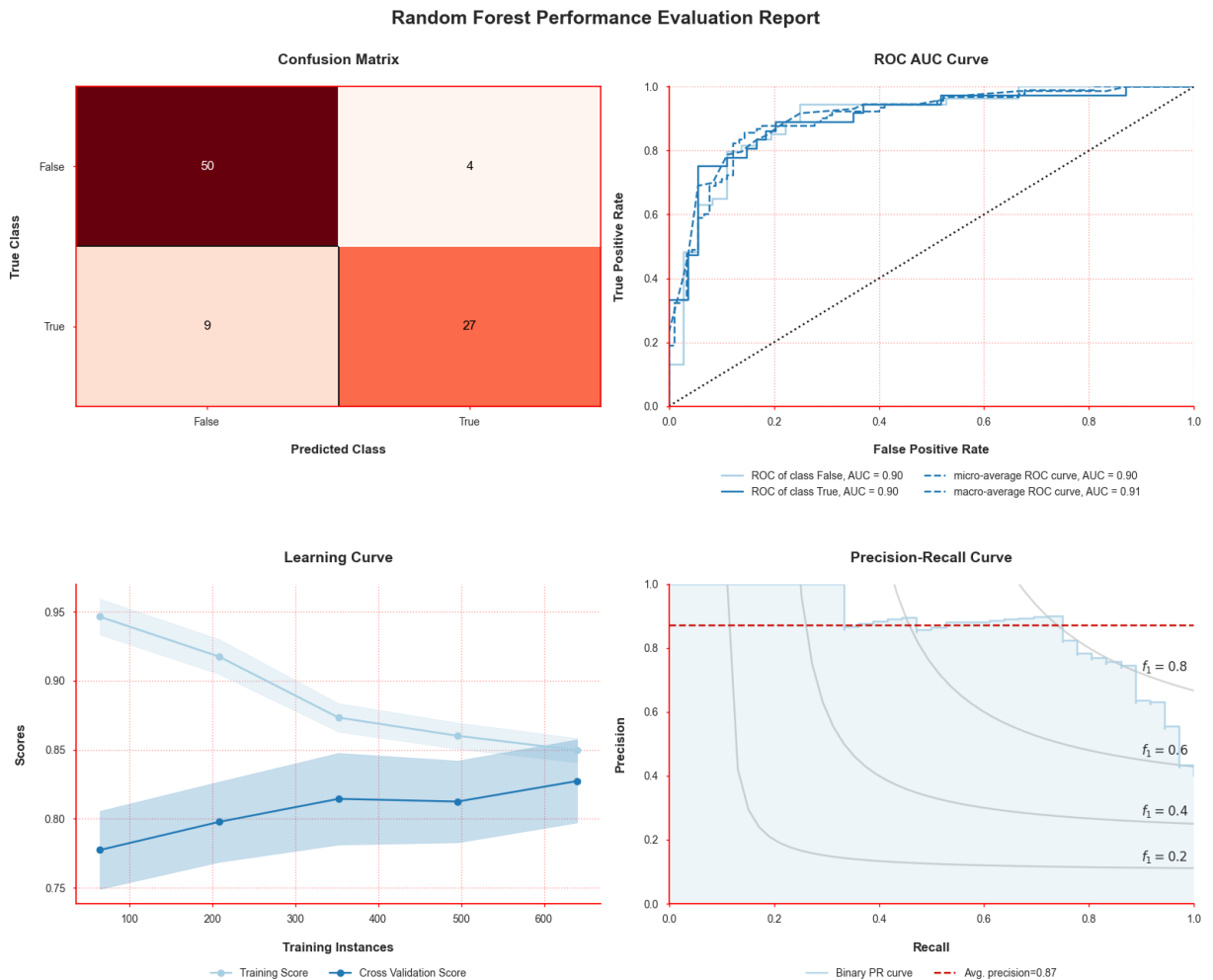
... Train and Test Accuracy Score for Random Forest ...

>> Train Accuracy: 85.52%

>> Test Accuracy: 85.56%

... Classification Report for Random Forest ...

	precision	recall	f1-score	support
0	0.85	0.93	0.88	54
1	0.87	0.75	0.81	36
accuracy			0.86	90
macro avg	0.86	0.84	0.85	90
weighted avg	0.86	0.86	0.85	90





## 7.7 | Extra Tree Classifier

**Extra Trees Classifier** is a type of ensemble learning technique which **aggregates the results of multiple de-correlated decision trees collected in a "forest" to output it's classification result**. In concept, it is very similar to a Random Forest Classifier and only differs from it in the manner of construction of the decision trees in the forest.

Each Decision Tree in the Extra Trees Forest is **constructed from the original training sample**. Then, at each test node, each tree is provided with a **random sample of k features** from the feature-set from which each decision tree must select the best feature to split the data based on some mathematical criteria (typically the Gini Index). This random sample of features leads to the creation of multiple de-correlated decision trees.

```
In [22]: # --- Extra Tree Parameters ---
parameter_et = {"algo__max_depth": [2, 3]
               , "algo__max_leaf_nodes": [3, 5, 7]}

# --- Extra Tree Algorithm ---
algo_et = ExtraTreesClassifier(random_state=42, n_jobs=-1)

# --- Applying Extra Tree ---
acc_score_train_et, acc_score_test_et, best_score_et = fit_ml_models(algo_et, param
```

... Fitting Extra Tree Classifier ...

Fitting 10 folds for each of 6 candidates, totalling 60 fits

>> Best Parameters: {'algo\_\_max\_depth': 3, 'algo\_\_max\_leaf\_nodes': 5}

>> Best Score: 0.809

... Train and Test Accuracy Score for Extra Tree Classifier ...

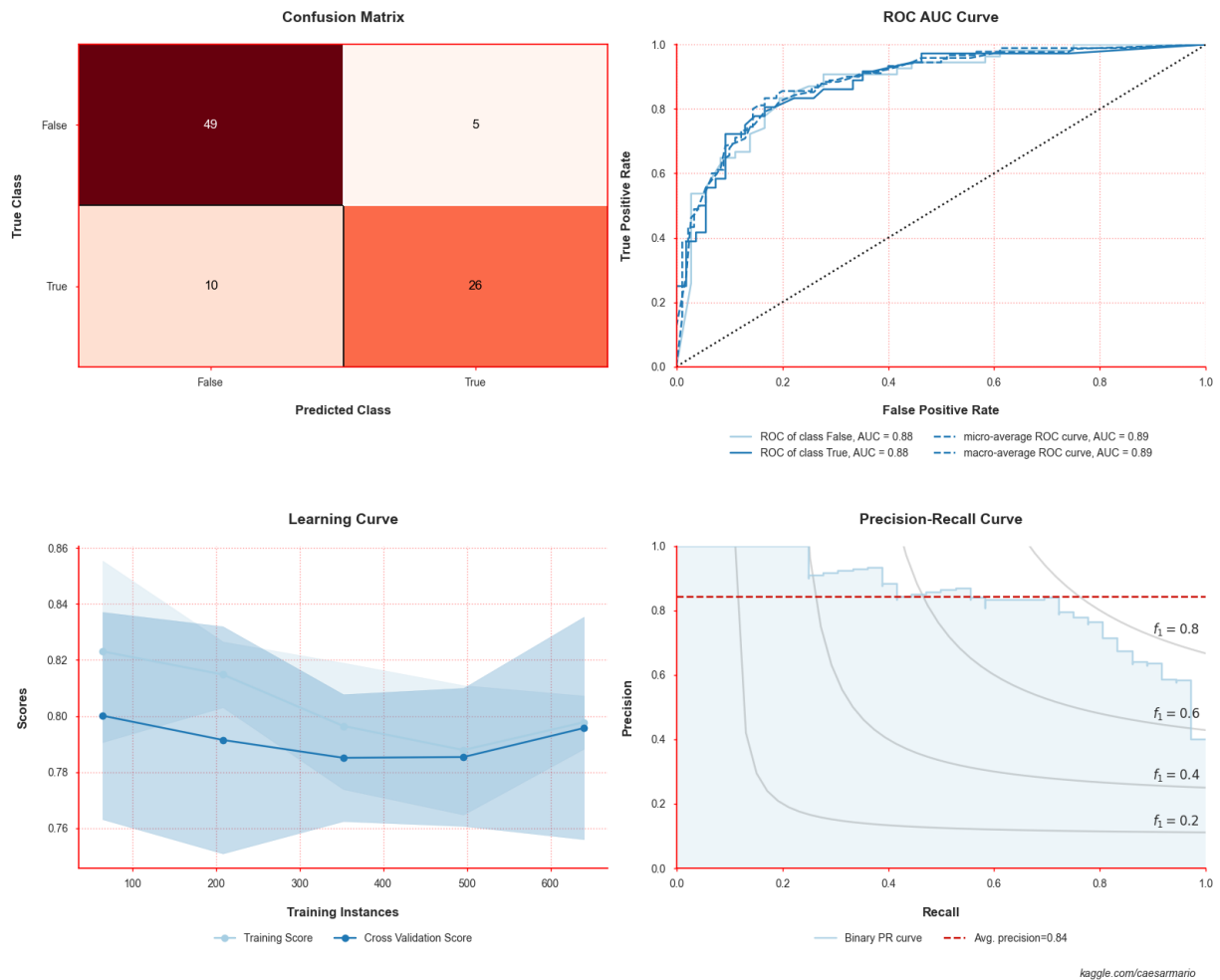
>> Train Accuracy: 80.90%

>> Test Accuracy: 83.33%

... Classification Report for Extra Tree Classifier ...

	precision	recall	f1-score	support
0	0.83	0.91	0.87	54
1	0.84	0.72	0.78	36
accuracy			0.83	90
macro avg	0.83	0.81	0.82	90
weighted avg	0.83	0.83	0.83	90

### Extra Tree Classifier Performance Evaluation Report



## 7.8 | Gradient Boosting

**Boosting** is a method of **converting weak learners into strong learners**. In boosting, **each new tree is a fit on a modified version** of the original data set. It strongly relies on the prediction that the next model will reduce prediction errors when blended with previous ones. The main idea is **to establish target outcomes for this upcoming model to minimize errors**.

**Gradient Boosting** trains many models in a **gradual, additive and sequential manner**. The term gradient boosting emerged because every case's target outcomes are based on the gradient's error with regards to the predictions. Every model reduces prediction errors by taking a step in the correct direction.



Boosting Algorithm by Rui Guo et al.

```
In [23]: # --- Gradient Boosting Parameters ---
parameter_gb = {
```

```

    "algo__learning_rate": [0.1, 0.3, 0.5]
    , "algo__n_estimators": [2, 4, 6]
    , "algo__min_weight_fraction_leaf": [0.1, 0.2, 0.5]
}

# --- Gradient Boosting Algorithm ---
algo_gb = GradientBoostingClassifier(loss="exponential", random_state=2)

# --- Applying Gradient Boosting ---
acc_score_train_gb, acc_score_test_gb, best_score_gb = fit_ml_models(algo_gb, param

```

... Fitting Gradient Boosting ...

Fitting 10 folds for each of 27 candidates, totalling 270 fits

```
>> Best Parameters: {'algo__learning_rate': 0.3, 'algo__min_weight_fraction_leaf':
0.1, 'algo__n_estimators': 6}
```

```
>> Best Score: 0.803
```

... Train and Test Accuracy Score for Gradient Boosting ...

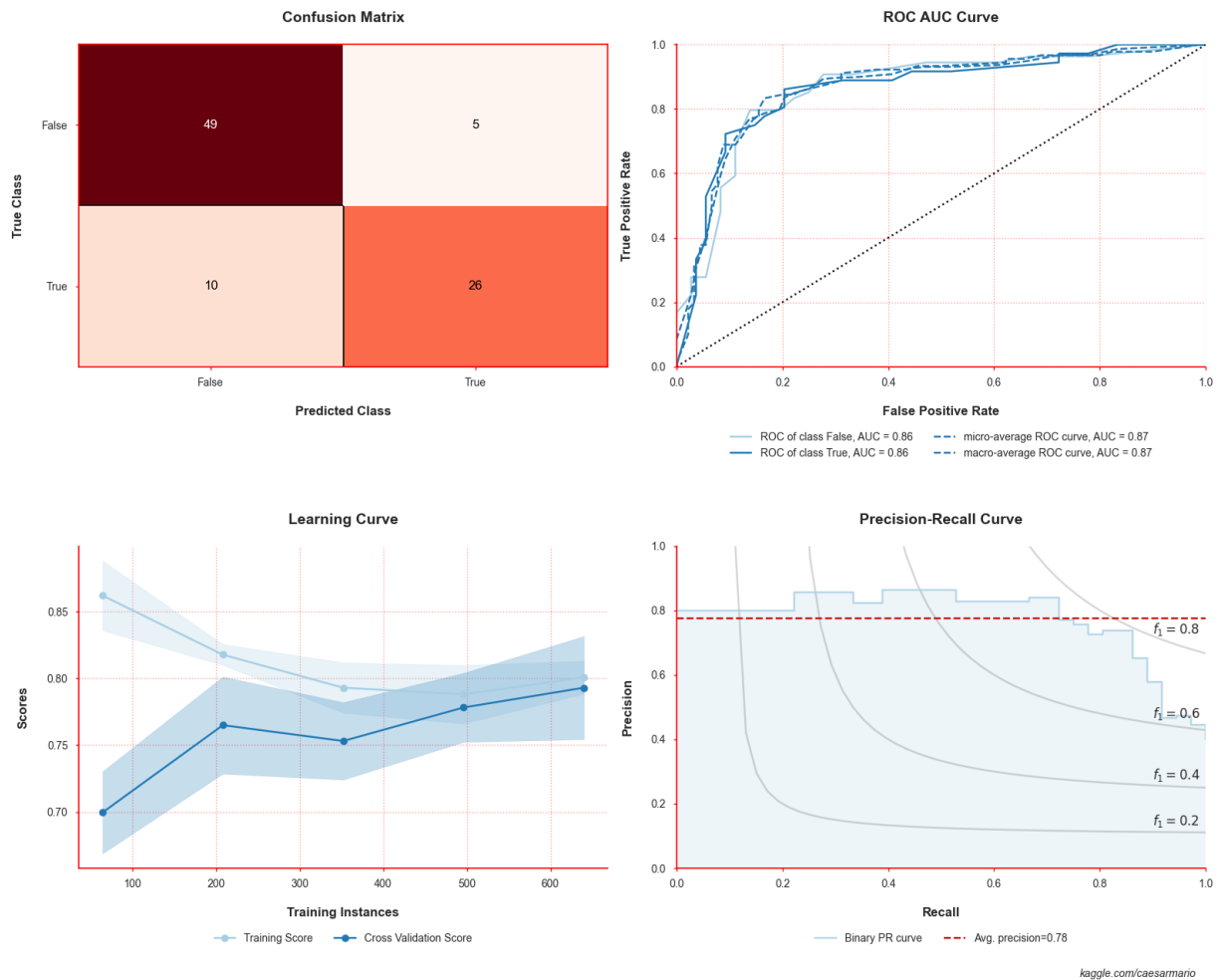
```
>> Train Accuracy: 80.90%
```

```
>> Test Accuracy: 83.33%
```

... Classification Report for Gradient Boosting ...

	precision	recall	f1-score	support
0	0.83	0.91	0.87	54
1	0.84	0.72	0.78	36
accuracy			0.83	90
macro avg	0.83	0.81	0.82	90
weighted avg	0.83	0.83	0.83	90

## Gradient Boosting Performance Evaluation Report



## 7.9 | AdaBoost

**AdaBoost** also called **Adaptive Boosting** is a technique in Machine Learning used as an Ensemble Method. The most common algorithm used with AdaBoost is **decision trees with one level** that means with Decision trees with only 1 split. These trees are also called **Decision Stumps**. **AdaBoost builds a model and gives equal weights to all the data points**. It then assigns higher weights to points that are wrongly classified. Now, all the points which have higher weights are given more importance in the next model. It will keep training models until and unless a low error is received.

```
In [24]: # --- AdaBoost Parameters ---
parameter_ab = {
    "algo_n_estimators": [6, 7, 10]
    , "algo_learning_rate": [0.2, 0.4, 0.8]
}

# --- AdaBoost Algorithm ---
algo_ab = AdaBoostClassifier(random_state=1)
```

```
# --- Applying AdaBoost ---
```

```
acc_score_train_ab, acc_score_test_ab, best_score_ab = fit_ml_models(algo_ab, param
```

```
... Fitting AdaBoost ...
```

```
Fitting 10 folds for each of 9 candidates, totalling 90 fits
```

```
>> Best Parameters: {'algo_learning_rate': 0.8, 'algo_n_estimators': 6}
```

```
>> Best Score: 0.803
```

```
... Train and Test Accuracy Score for AdaBoost ...
```

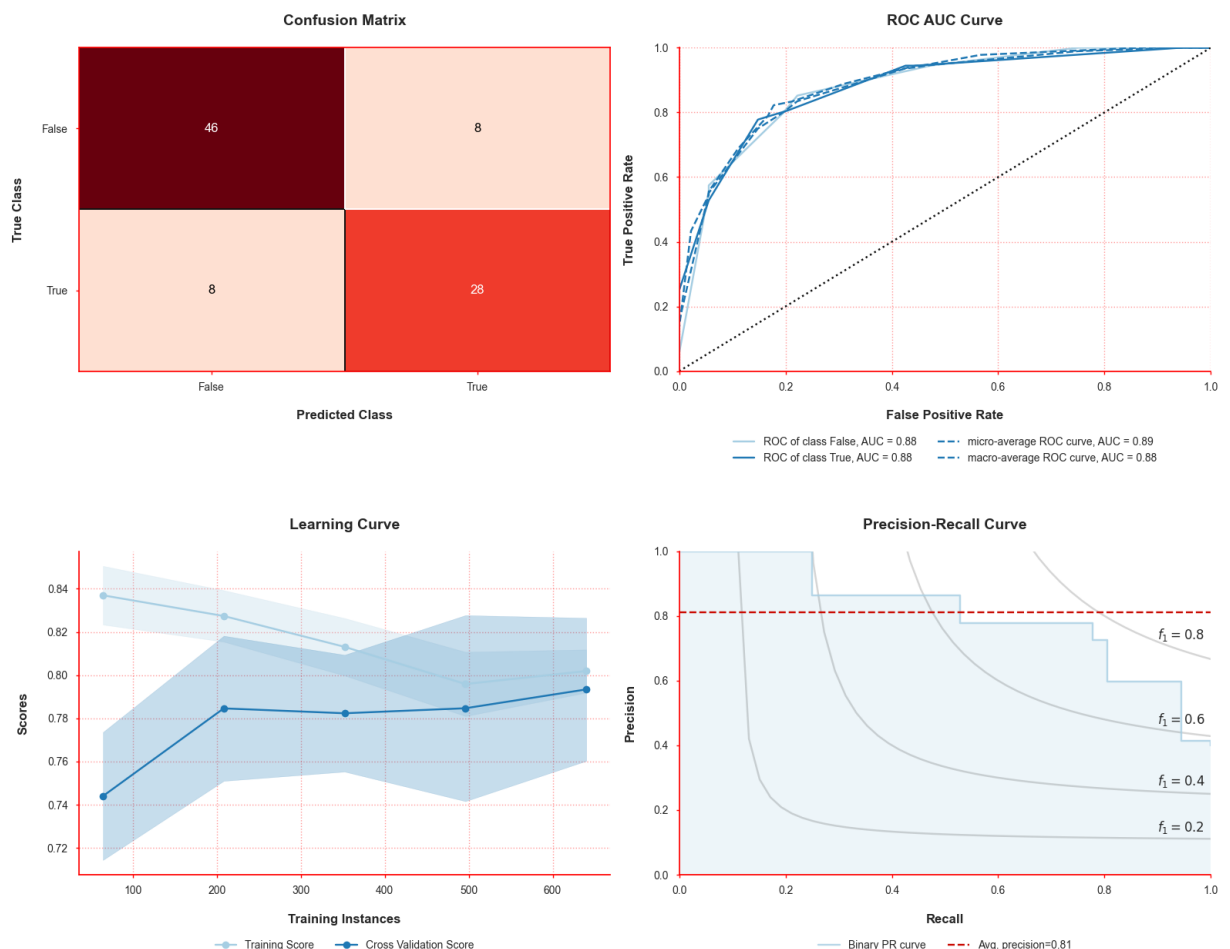
```
>> Train Accuracy: 80.65%
```

```
>> Test Accuracy: 82.22%
```

```
... Classification Report for AdaBoost ...
```

	precision	recall	f1-score	support
0	0.85	0.85	0.85	54
1	0.78	0.78	0.78	36
accuracy			0.82	90
macro avg	0.81	0.81	0.81	90
weighted avg	0.82	0.82	0.82	90

AdaBoost Performance Evaluation Report



## 7.10 | Model Comparison 🧐

After implementing and tuning 9 models, this section will **compare all machine learning models accuracy and best score**.

```
In [25]: # --- Create Accuracy Comparison Table ---
df_compare = pd.DataFrame({'Model': ['Logistic Regression', 'K-Nearest Neighbour',
                                     'Decision Tree', 'Random Forest', 'Extra Tree'],
                           'Accuracy Train': [acc_score_train_lr, acc_score_train_dt, acc_score_train_rf,
                                              acc_score_train_dt, acc_score_train_rf],
                           'Accuracy Test': [acc_score_test_lr, acc_score_test_knn, acc_score_test_dt,
                                              acc_score_test_dt, acc_score_test_rf],
                           'Best Score': [best_score_lr, best_score_knn, best_score_dt,
                                          best_score_et, best_score_gb, best_score_rf])

# --- Create Comparison Table ---
print(clear.start+f"... Models Comparison ..."+clear.end)
print(clear.color+'*' * 26)
df_compare.sort_values(by='Best Score', ascending=False).reset_index(drop=True).style

... Models Comparison ...
*****
```

Out[25]:

	Model	Accuracy Train	Accuracy Test	Best Score
0	Random Forest	85.518000	85.556000	0.836500
1	Support Vector Machine	83.146000	81.111000	0.825200
2	Decision Tree	82.772000	82.222000	0.815300
3	Extra Tree Classifier	80.899000	83.333000	0.809000
4	Logistic Regression	81.149000	83.333000	0.807800
5	K-Nearest Neighbour	82.772000	85.556000	0.805200
6	Gradient Boosting	80.899000	83.333000	0.802800
7	AdaBoost	80.649000	82.222000	0.802800
8	Gaussian NB	80.400000	82.222000	0.790300

- From the results of the **accuracy of the train and test** above, **most models experienced overfitting or underfitting**. However, **several models have a good fit**, where the difference between train and test accuracy or vice versa is a little. These models are **random forest, Support Vector Machine, Decision Tree, and extra tree classifier**.
- As seen in the data frame above, of the four models, **random forest and Support Vector Machine have the highest accuracy compared to the other models**. This is also supported by the ROC AUC curve figure for

random forest and Support Vector Machine, where the AUC value for both models is close to 1, which means that both models can predict well whether passengers survival. The **confusion matrix** shows that the prediction results between the actual target and the predicted target for the random forest and Support Vector Machine models in each class in the test data are better than those of other models.

- Judging from the F1 scores of both models, both models do a very good job differentiating passengers from those who are not (scores above 0.80). If seen from the precision value for Random Forest, 85% of all the passengers that the model predicted have survived. Whereas in the Support Vector Machine precision value, only 82% out of all the passengers that the model predicted have survived, slightly lower than the random forest precision value.
- Furthermore, in the learning curve between Support Vector Machine and random forest, the learning curve for the random forest is more ideal than Support Vector Machine. This is because both training and validation scores of Support Vector Machine stay too close together (indicates low variance and high bias). This will more likely result in poor fit and especially poor generalization of the data (towards the data it has not seen before).
- In a random forest, the validation score constantly improves as the number of training set sizes gets larger (notice the difference in the x-axis scale from the previous two curves). Both the training and validation scores also converge to nearly similar values. This is a model that can generalize very well. From the analysis above, it can be concluded that the random forest model best predicts whether a person has heart disease.

## 8. | Miscellaneous

This section focuses on creating a complete pipeline, starting from data processing to a machine learning pipeline, using the best model concluded in the previous section and exporting it to joblib and pickle (.pkl) files. Besides that, test dataset predicted results would also be exported along with actual results in CSV and JSON files. Moreover, this section will also make predictions on dummy data (data generated using Python functions) and export them to CSV and JSON files.

### 8.1 | Creating Outputs

The complete pipeline will be exported in this section. The pipeline will be stored using the joblib library into joblib and pickle (.pkl) files. This section will also show the test data frame before exporting the predicted results and the actual results to the CSV and JSON files.

```
In [26]: # --- Complete Pipeline: Preprocessor & RF ---
rf_pipeline = Pipeline([
    ('preprocessor', preprocessor)
    , ('algo', RandomForestClassifier(max_depth=3, random_state=99, n_jobs=-1))
])

# --- Save Complete Pipeline (joblib and pickle) ---
directory = r'D:\Data_Science\Machine_Learning\Titanic_Classification'
file_name = 'pipeline_titanic_survival_random_forest_darshanpathak'

# Save pipeline using joblib
for ext in ['joblib', 'pkl']:
    file_path = os.path.join(directory, f'{file_name}.{ext}')
    joblib.dump(rf_pipeline, file_path)

In [27]: # --- Dataframes to Create Test Output Dataframe ---
rf_pipeline.fit(x_train, y_train)
y_pred_rf = rf_pipeline.predict(x_test)
pred_target = pd.DataFrame(y_pred_rf, columns=['pred_target'])

x_test_output = x_test.reset_index()
actual_target = y_test.to_frame(name='actual_target').reset_index()

# --- Combining and Creating Test Output Dataframe ---
df_test_output = pd.concat([x_test_output, actual_target, pred_target], axis=1).dro

# --- Showing Sample Test Output Dataframe ---
print(clear.start+'.: Sample Test Dataframe :.'+clear.end)
print(clear.color+'*' * 28)
df_test_output.sample(n=10, random_state=0).reset_index(drop=True).style.background

.: Sample Test Dataframe :.
*****
```



Out[27]:

	Pclass	Sex	Age	SibSp	Parch	Fare	Embarked	actual_target	pred_targ
0	3	male	20.000000	0	0	7.925000	S	0	
1	3	male	30.000000	0	0	7.225000	C	0	
2	1	male	11.000000	1	2	120.000000	S	1	
3	3	male	29.000000	1	0	7.045800	S	0	
4	2	male	21.000000	0	0	73.500000	S	0	
5	3	male	19.000000	0	0	7.895800	S	0	
6	3	male	16.000000	1	3	34.375000	S	0	
7	3	male	25.000000	0	0	7.750000	Q	0	
8	2	female	38.000000	0	0	13.000000	S	0	
9	3	male	25.000000	0	0	7.250000	S	0	

In [28]:

```
# --- Export to CSV and JSON Files ---
output_name = 'titanic_survival_random_forest_darshanpathak'
df_test_output.to_csv(f'D:\Data_Science\Machine_Learning\Titanic_Classification.csv')
df_test_output.to_json(f'D:\Data_Science\Machine_Learning\Titanic_Classification.js')
```

## 9. | Conclusions and Future Improvements



From the results of dataset analysis and implementation of machine learning models in the previous section, it can be concluded as follows:

- **Random forest is the best model** out of 9 machine-learning models implemented in this notebook. This is because **this model fits well with train and test data**. In addition, **this model also performs better than other models when predicting the test data** (can be seen from the performance evaluation graph and classification report of each model).
- **The prediction results on test data, dummy data, and the complete machine learning pipeline have been successfully exported** for other purposes. In addition, data exploration has also been successfully carried out using the `ydata-profiling`, `seaborn`, and `matplotlib` libraries.
- **Several improvements can be implemented in the following research/notebook**. For example, by carrying out A/B Testing on passengers with the same major vessel number in one group. Another example is performing advanced hyperparameter tuning experiments to obtain higher accuracy (~90%).

**Thank You**