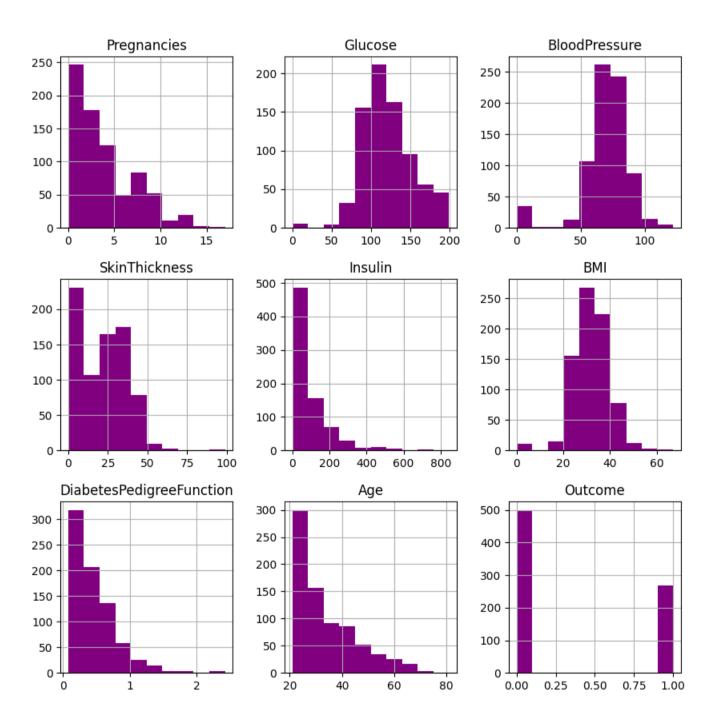
```
Diabetes Prediction using Machine Learning
Diabetes refers to a collection of metabolic disorders characterized by elevated blood sugar levels persisting over an extended period.
Symptoms associated with high blood sugar levels include frequent urination, heightened thirst, and increased appetite. Without proper
treatment, diabetes can lead to various complications. Acute complications may involve conditions such as diabetic ketoacidosis,
hyperosmolar hyperglycemic state, or even mortality. Furthermore, severe long-term consequences encompass cardiovascular ailments,
 strokes, chronic kidney ailments, foot ulcers, and vision impairment.
We will try to build a machine learning model to accurately predict whether or not the patients in the dataset have diabetes or not?
Number of Observation Units: 768
Variable Number: 9
import matplotlib.pyplot as plt
import seaborn as sns
import numpy as np
import pandas as pd
from google.colab import files
uploaded = files.upload()
Choose Files No file chosen
                                    Upload widget is only available when the cell has been executed in the current browser session. Please rerun this cell to enable.
     Saving diabetes.csv to diabetes.csv
Reading the dataset which is the CSV format
df = pd.read_csv('diabetes.csv')
df.head()
        Pregnancies Glucose BloodPressure SkinThickness Insulin BMI DiabetesPedigreeFunction Age Outcome
                                                                                        0.627 50
                                                            0 33.6
                                                            0 26.6
                                                                                       0.351 31
                        183
                                                            0 23.3
                                                                                       0.672 32
                         89
                                                    23
                                                           94 28.1
                                                                                       0.167 21
                        137
                                                           168 43.1
                                                                                       2.288 33
Exploratory Data Analysis (EDA)
df.shape
    (768, 9)
df.columns
     Index(['Pregnancies', 'Glucose', 'BloodPressure', 'SkinThickness', 'Insulin',
            'BMI', 'DiabetesPedigreeFunction', 'Age', 'Outcome'],
           dtype='object')
df.info()
    <class 'pandas.core.frame.DataFrame'>
    RangeIndex: 768 entries, 0 to 767
    Data columns (total 9 columns):
                                 Non-Null Count Dtype
     # Column
     0 Pregnancies
                                  768 non-null int64
      1 Glucose
                                  768 non-null int64
      2 BloodPressure
                                  768 non-null int64
      3 SkinThickness
                                  768 non-null
                                                int64
      4 Insulin
                                  768 non-null int64
                                  768 non-null
                                                float64
      6 DiabetesPedigreeFunction 768 non-null
      7 Age
                                 768 non-null
                                                int64
                                  768 non-null int64
      8 Outcome
     dtypes: float64(2), int64(7)
     memory usage: 54.1 KB
df.describe()
            Pregnancies Glucose BloodPressure SkinThickness Insulin
                                                                              BMI DiabetesPedigreeFunction
                                                                                                                 Age Outcome

        count
        768.000000
        768.000000
        768.000000
        768.000000
        768.000000
        768.000000

                                                                                                 768.000000 768.000000 768.000000
                                                    20.536458 79.799479 31.992578
                                                                                                   0.471876 33.240885 0.348958
                3.845052 120.894531
                                      69.105469
               3.369578 31.972618
                                      19.355807
                                                    15.952218 115.244002 7.884160
                                                                                                   0.331329 11.760232 0.476951
                                                                                                   0.078000 21.000000 0.000000
                         0.000000
                                        0.000000
                                                     0.000000 0.000000 0.000000
                                      62.000000
                                                     0.000000 0.000000 27.300000
                                                                                                   0.243750 24.000000 0.000000
               1.000000 99.000000
                                      72.000000
                                                    23.000000 30.500000 32.000000
                                                                                                   0.372500 29.000000 0.000000
                3.000000 117.000000
              6.000000 140.250000
                                      80.000000
                                                    32.000000 127.250000 36.600000
                                                                                                   0.626250 41.000000 1.000000
              17.000000 199.000000 122.000000
                                                    99.000000 846.000000 67.100000
                                                                                                   2.420000 81.000000 1.000000
df.describe().T
                                                   std min 25% 50% 75% max
            Pregnancies
                             768.0 3.845052 3.369578 0.000 1.00000 3.0000 6.00000 17.00
             Glucose
                              768.0 120.894531 31.972618 0.000 99.00000 117.0000 140.25000 199.00
                             768.0 69.105469 19.355807 0.000 62.00000 72.0000 80.00000 122.00
                             768.0 20.536458 15.952218 0.000 0.00000 23.0000 32.00000 99.00
           SkinThickness
                              768.0 79.799479 115.244002 0.000 0.00000 30.5000 127.25000 846.00
                              768.0 31.992578 7.884160 0.000 27.30000 32.0000 36.60000 67.10
      DiabetesPedigreeFunction 768.0 0.471876 0.331329 0.078 0.24375 0.3725 0.62625 2.42
                              768.0 33.240885 11.760232 21.000 24.00000 29.0000 41.00000 81.00
             Outcome
                             768.0 0.348958 0.476951 0.000 0.00000 0.0000 1.00000 1.00
let's check that if our dataset have null values or not
df.isnull().head(6)
        Pregnancies Glucose BloodPressure SkinThickness Insulin BMI DiabetesPedigreeFunction Age Outcome
                                                  False False False
                                                                                        False False
                                                         False False
                                                                                        False False False
              False
                                    False
                                                  False False False
                                                                                        False False False
              False
                                    False
                                                  False False False
                                                                                        False False False
                                    False
              False
                                                  False False False
                                                                                        False False False
                                                  False False False
                                                                                        False False False
              False
df.isnull().sum()
     Pregnancies
     Glucose
    BloodPressure
     SkinThickness
     Insulin
    DiabetesPedigreeFunction
     Outcome
    dtype: int64
df_copy = df.copy(deep = True)
df_copy[['Glucose','BloodPressure','SkinThickness','Insulin','BMI']] = df_copy[['Glucose','BloodPressure','SkinThickness','Insulin','BMI']].replace(0,np.NaN)
# Showing the Count of NANs
print(df_copy.isnull().sum())
     Pregnancies
     Glucose
    BloodPressure
    SkinThickness
     Insulin
     DiabetesPedigreeFunction
    Age
    Outcome
    dtype: int64
 As mentioned above that now we will be replacing the zeros with the NAN values so that we can impute it later to maintain the authenticity of
the dataset as well as trying to have a better Imputation approach i.e to apply mean values of each column to the null values of the respective
 Data Visalization
```

df.hist(figsize=(10, 10), color="purple")

# Show plot plt.show()



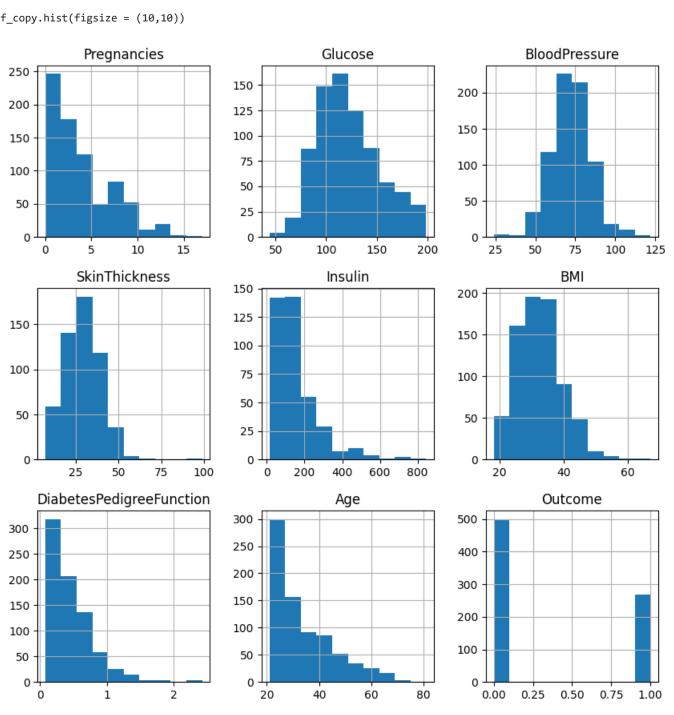
The best way to start the analysis of the dataset is to shows the occurrence of every kind of value in the graphical structure which in turn lets us know the range of the data.

Now we will be imputing the mean value of the column to each missing value of that particular column.

Plotting the distributions after removing the NAN values.

df\_copy['Glucose'].fillna(df\_copy['Glucose'].mean(), inplace = True) df\_copy['BloodPressure'].fillna(df\_copy['BloodPressure'].mean(), inplace = True) df\_copy['SkinThickness'].fillna(df\_copy['SkinThickness'].median(), inplace = True) df\_copy['Insulin'].fillna(df\_copy['Insulin'].median(), inplace = True) df\_copy['BMI'].fillna(df\_copy['BMI'].median(), inplace = True)

p = df\_copy.hist(figsize = (10,10))



import missingno as msno p = msno.bar(df) 768 1.0 0.8 614 0.6

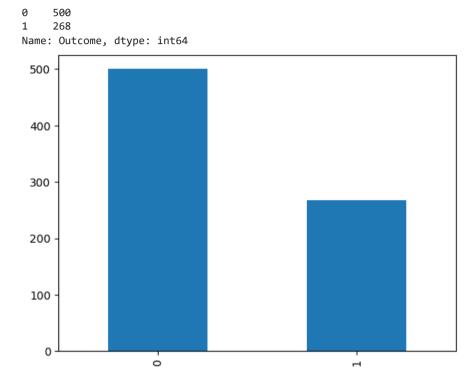
307

Now, let's check that how well our outcome column is balanced

0.4

Now we can clearly see that there are no null values in the dataset.

color\_wheel = {1: "#0392cf", 2: "#7bc043"} colors = df["Outcome"].map(lambda x: color\_wheel.get(x + 1)) print(df.Outcome.value\_counts()) p=df.Outcome.value\_counts().plot(kind="bar")



According above visualization it is clearly visible that our dataset is completely imbalanced in fact the number of patients who are diabetic is half of the patients who are non-diabetic.

plt.subplot(121), sns.distplot(df['Insulin']) plt.subplot(122), df['Insulin'].plot.box(figsize=(16,5))

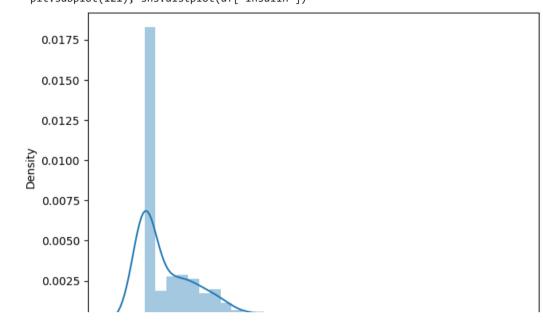
plt.show()

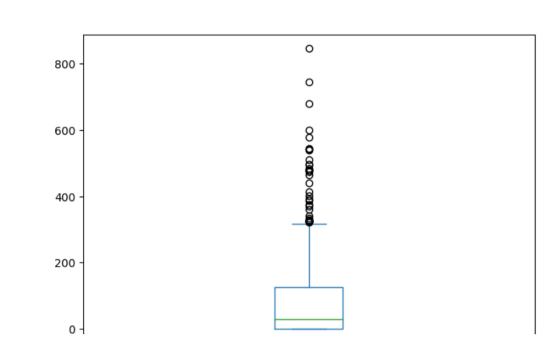
<ipython-input-21-1d062345beba>:1: UserWarning: `distplot` is a deprecated function and will be removed in seaborn v0.14.0.

Please adapt your code to use either `displot` (a figure-level function with similar flexibility) or `histplot` (an axes-level function for histograms).

For a guide to updating your code to use the new functions, please see https://gist.github.com/mwaskom/de44147ed2974457ad6372750bbe5751

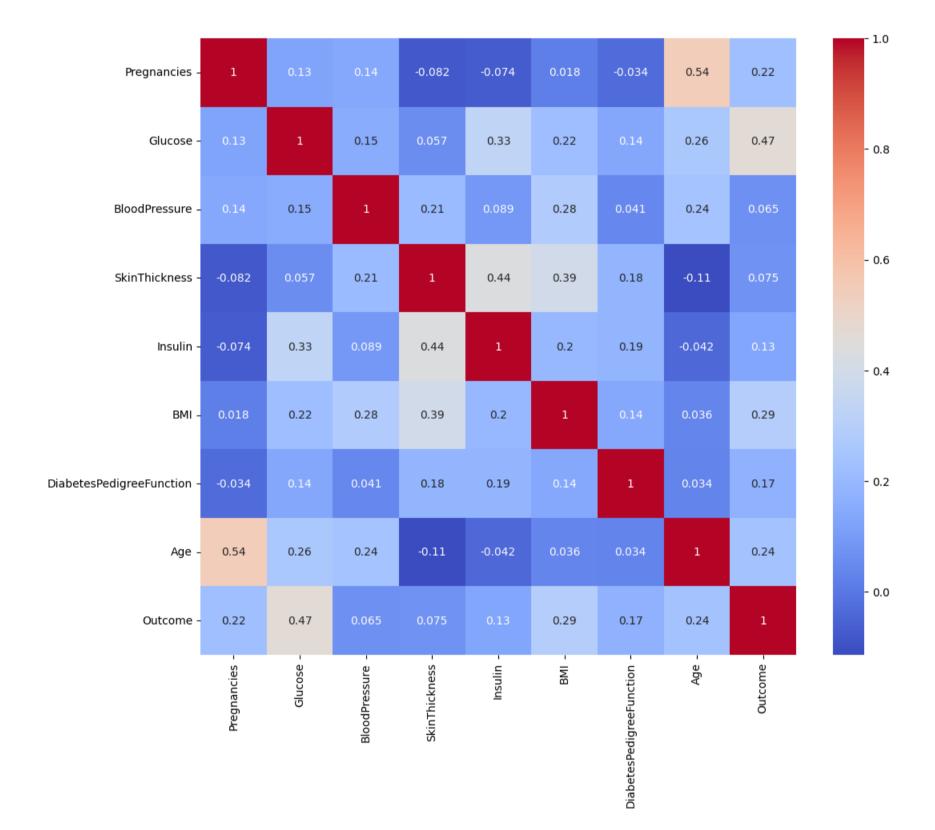
plt.subplot(121), sns.distplot(df['Insulin'])





Correlation between all the features before cleaning

plt.figure(figsize=(12, 10)) # Specify the colormap using the cmap parameter p = sns.heatmap(df.corr(), annot=True, cmap='coolwarm') plt.show()



## **Model Building** Splitting the dataset

x = df.drop('Outcome', axis = 1)y = df['Outcome']

Pregnancies Glucose BloodPressure SkinThickness Insulin BMI DiabetesPedigreeFunction Age 35 0 33.6 0.627 50 29 0 26.6 0.351 31 8 183 0 0 23.3 0.672 32 0.167 21 23 94 28.1 137 35 168 43.1 2.288 33 0.171 63 101 48 180 32.9 27 0 36.8 0.340 27 122 121 23 112 26.2 0.245 30 1 126 0 0 30.1 0.349 47 1 93 31 0 30.4 0.315 23 768 rows × 8 columns

763 0 764 0 765 0 766 1 767 0 Name: Outcome, Length: 768, dtype: int64

x = np.array(x)

y = np.array(y)

...,
[ 5. , 121. , 72. , ..., 26.2 , 0.245, 30. ],
[ 1. , 126. , 60. , ..., 30.1 , 0.349, 47. ],
[ 1. , 93. , 70. , ..., 30.4 , 0.315, 23. ]])

array([1, 0, 1, 0, 1, 0, 1, 0, 1, 1, 0, 1, 0, 1, 1, 1, 1, 1, 0, 1, 0, 0, 1, 0, 0, 0, 1, 0, 1, 0, 0, 0, 0, 1, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 0, 0, 0, 1, 0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0, 0, 0, 1, 1, 1, 0, 0, 1, 1, 1, 0, 0, 0, 1, 0, 0, 0, 1, 1, 0, 0, 1, 1, 1, 1, 1, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 1, 1, 0, 0, 0, 1, 0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 1, 1, 0, 0, 0, 1, 0, 1, 0, 1, 0, 0, 0, 0, 1, 1, 1, 1, 1, 0, 0, 1, 1, 0, 1, 0, 1, 1, 1, 0, 0, 0, 0, 0, 0, 1, 1, 0, 1, 0, 0, 0, 1, 1, 1, 1, 0, 1, 1, 1, 1, 0, 0, 0, 0, 1, 0, 0, 1, 1, 0, 0, 0, 1, 1, 1, 1, 0, 0, 0, 1, 1, 0, 1, 0, 0, 0, 0, 0, 0, 0, 1, 1, 0, 0, 0, 1, 0, 1, 0, 0, 1, 0, 1, 0, 0, 1, 1, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 0, 1, 1, 0, 0, 1, 0, 0, 0, 1, 1, 1, 0, 0, 1, 0, 1, 0, 1, 1, 0, 1, 0, 0, 1, 0, 1, 1, 0, 0, 1, 0, 1, 0, 0, 1, 0, 1, 0, 1, 1, 1, 0, 0, 1, 0, 1, 0, 0, 0, 1, 0, 0, 0, 0, 1, 1, 1, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 1, 0, 1, 1, 0, 1, 0, 1, 0, 1, 1, 0, 0, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 0, 0, 1, 1, 0, 1, 0, 1, 0, 0, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 1, 1, 1, 0, 0, 1, 0, 0, 1, 0, 0, 0, 1, 0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 1, 1, 1, 1, 1, 0, 0, 1, 1, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 1, 0, 1, 1, 0, 0, 0, 1, 0, 1, 0, 1, 0, 1, 0, 1, 0, 0, 1, 0, 0, 1, 0, 0, 0, 0, 1, 1, 0, 1, 0, 0, 0, 0, 1, 1, 0, 1, 0, 0, 0, 1, 1, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 1, 1, 0, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 1, 1, 1, 1, 0, 1, 1, 0, 0, 0, 0, 0, 0, 0, 1, 1, 0, 1, 0, 0, 1, 0, 1, 0, 0, 0, 0, 0, 1, 0, 1, 0, 1, 0, 1, 1, 0, 0, 0, 0, 1, 1, 0, 0, 0, 1, 0, 1, 1, 0, 0, 1, 0, 0, 1, 1, 0, 0, 1, 0, 0, 1, 0, 0, 0, 0, 0, 0, 0, 1, 1, 1, 0, 0, 0, 0, 0, 1, 1, 0, 0, 1, 0, 0, 1, 0,

1, 1, 1, 0, 0, 1, 1, 1, 0, 1, 0, 1, 0, 1, 0, 0, 0, 0, 1, 0])

Normalize

x = scaler.fit\_transform(x)

from sklearn.preprocessing import StandardScaler scaler = StandardScaler()

array([[ 0.63994726, 0.84832379, 0.14964075, ..., 0.20401277, 0.46849198, 1.4259954], [-0.84488505, -1.12339636, -0.16054575, ..., -0.68442195, -0.36506078, -0.19067191], [ 1.23388019, 1.94372388, -0.26394125, ..., -1.10325546, 0.60439732, -0.10558415], ..., [ 0.3429808 , 0.00330087, 0.14964075, ..., -0.73518964,

-0.68519336, -0.27575966], [-0.84488505, 0.1597866, -0.47073225, ..., -0.24020459, -0.37110101, 1.17073215], [-0.84488505, -0.8730192 , 0.04624525, ..., -0.20212881, -0.47378505, -0.87137393]])

Train / Test

from sklearn.model\_selection import train\_test\_split x\_train, x\_test, y\_train, y\_test = train\_test\_split(x, y, test\_size = 0.2)

x\_train.shape, y\_train.shape ((614, 8), (614,))

x\_test.shape, y\_test.shape ((154, 8), (154,))

Naive Bayes Building the model using Naive Bayes

from sklearn.naive\_bayes import GaussianNB

model = GaussianNB() model.fit(x\_train, y\_train)

> ▼ GaussianNB GaussianNB()

y\_pred\_train = model.predict(x\_train) y\_pred\_test = model.predict(x\_test)

from sklearn.metrics import accuracy\_score acc\_train = accuracy\_score(y\_true = y\_train, y\_pred = y\_pred\_train) acc\_test = accuracy\_score(y\_true = y\_test, y\_pred = y\_pred\_test)

acc\_train, acc\_test

(0.7687296416938111, 0.7467532467532467)

from sklearn.metrics import confusion\_matrix, precision\_score, recall\_score

confusion\_matrix(y\_test, y\_pred\_test)

array([[83, 17], [22, 32]])

precision\_score(y\_test, y\_pred\_test) 0.6530612244897959

precision\_score(y\_train, y\_pred\_train)

0.6855670103092784

recall\_score(y\_test, y\_pred\_test)

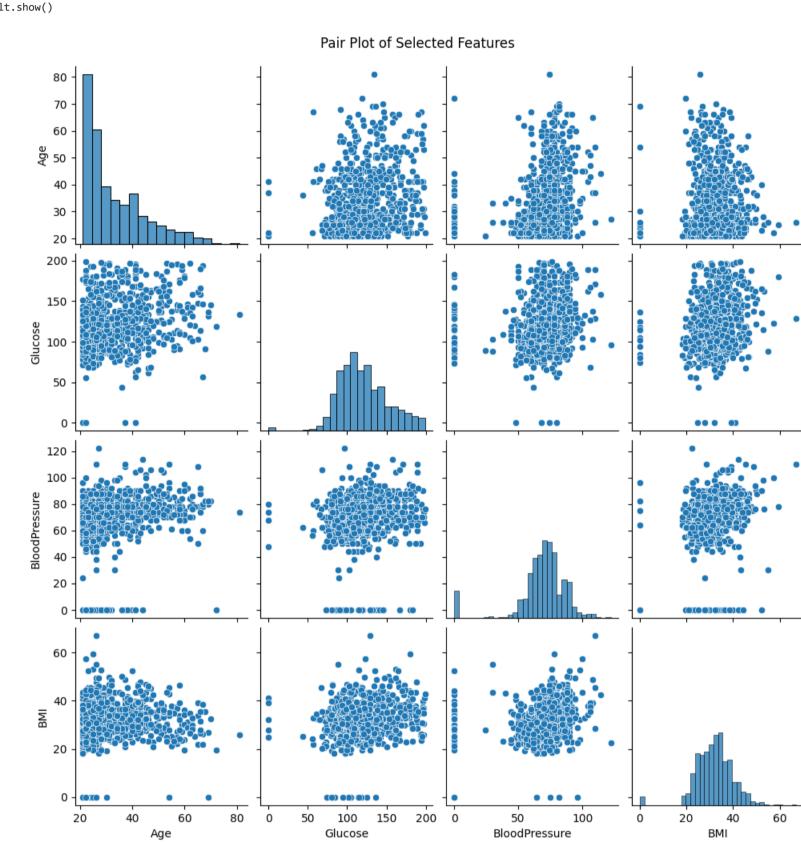
0.5925925925925926

recall\_score(y\_train, y\_pred\_train)

0.6214953271028038 selected\_features = ['Age', 'Glucose', 'BloodPressure', 'BMI']

# Create a pair plot sns.pairplot(df[selected\_features])

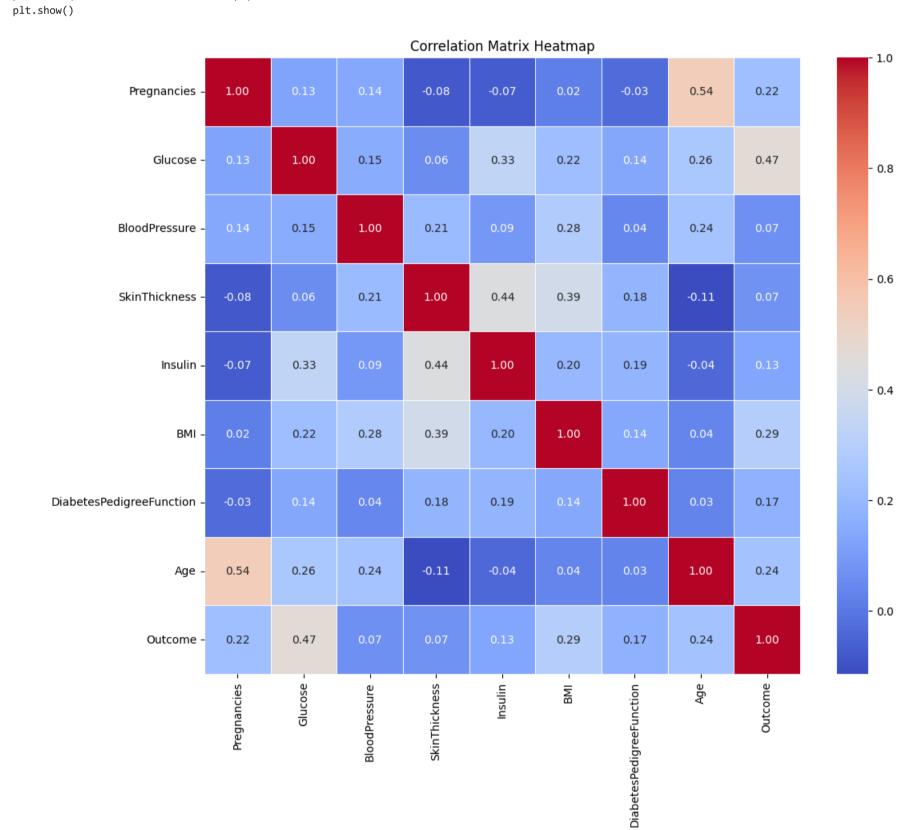
plt.suptitle("Pair Plot of Selected Features", y=1.02) plt.show()



	Pregnancies	Glucose	BloodPressure	SkinThickness	Insulin	BMI	DiabetesPedigreeFunction	Age	Outcome
Pregnancies	1.000000	0.129459	0.141282	-0.081672	-0.073535	0.017683	-0.033523	0.544341	0.221898
Glucose	0.129459	1.000000	0.152590	0.057328	0.331357	0.221071	0.137337	0.263514	0.466581
BloodPressure	0.141282	0.152590	1.000000	0.207371	0.088933	0.281805	0.041265	0.239528	0.065068
SkinThickness	-0.081672	0.057328	0.207371	1.000000	0.436783	0.392573	0.183928	-0.113970	0.074752
Insulin	-0.073535	0.331357	0.088933	0.436783	1.000000	0.197859	0.185071	-0.042163	0.130548
ВМІ	0.017683	0.221071	0.281805	0.392573	0.197859	1.000000	0.140647	0.036242	0.292695
DiabetesPedigreeFunction	-0.033523	0.137337	0.041265	0.183928	0.185071	0.140647	1.000000	0.033561	0.173844
Age	0.544341	0.263514	0.239528	-0.113970	-0.042163	0.036242	0.033561	1.000000	0.238356
Outcome	0.221898	0.466581	0.065068	0.074752	0.130548	0.292695	0.173844	0.238356	1.000000

correlation\_matrix = df.corr()

# Create a heatmap plt.figure(figsize=(12, 10)) sns.heatmap(correlation\_matrix, annot=True, cmap='coolwarm', fmt=".2f", linewidths=.5) plt.title("Correlation Matrix Heatmap")



from sklearn.neighbors import KNeighborsClassifier

knn = KNeighborsClassifier(n\_neighbors=4) knn.fit(x\_train, y\_train)

> KNeighborsClassifier KNeighborsClassifier(n\_neighbors=4)

y\_pred\_train = knn.predict(x\_train) y\_pred\_test = knn.predict(x\_test)

from sklearn.metrics import accuracy\_score acc\_train = accuracy\_score(y\_train, y\_pred\_train) acc\_test = accuracy\_score(y\_test, y\_pred\_test)

acc\_train, acc\_test (0.8094462540716613, 0.7532467532467533)

from sklearn.metrics import recall\_score, precision\_score, confusion\_matrix confusion\_matrix(y\_test, y\_pred\_test)

array([[92, 8], [30, 24]])

p = precision\_score(y\_test, y\_pred\_test)

r = recall\_score(y\_test, y\_pred\_test)

0.75

0.4444444444444444

As evident from the results, the recall\_score is unsatisfactory. By augmenting the value of k, we observe an improvement in the recall score. Therefore, increasing the value of k can enhance the recall performance for this dataset.

from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors= 8)

knn.fit(x\_train, y\_train) y\_pred\_train = knn.predict(x\_train) y\_pred\_test = knn.predict(x\_test) from sklearn.metrics import accuracy\_score acc\_train = accuracy\_score(y\_train, y\_pred\_train) acc\_test = accuracy\_score(y\_test, y\_pred\_test) acc\_train, acc\_test

(0.7931596091205212, 0.7532467532467533)

from sklearn.metrics import recall\_score, precision\_score, confusion\_matrix confusion\_matrix(y\_test, y\_pred\_test)

array([[91, 9], [29, 25]])

p = precision\_score(y\_test, y\_pred\_test)

0.7352941176470589

```
naive_bayes = GaussianNB()
knn = KNeighborsClassifier()
naive_bayes.fit(x_train, y_train)
knn.fit(x_train, y_train)
# Predictions for Naive Bayes
y_pred_nb = naive_bayes.predict(x_test)
y_pred_knn = knn.predict(x_test)
recall_nb = recall_score(y_test, y_pred_nb)
recall_knn = recall_score(y_test, y_pred_knn)
plt.figure(figsize=(8, 6))
plt.bar(['Naive Bayes', 'KNN'], [recall_nb, recall_knn], color=['blue', 'green'])
plt.title('Comparison of Recall Scores')
plt.xlabel('Classifier')
plt.ylabel('Recall Score')
plt.ylim(0, 1) # Limiting y-axis from 0 to 1 for better visualization
plt.grid(True)
plt.show()
                                    Comparison of Recall Scores
        0.8 -
        0.2
                          Naive Bayes
                                                                    KNN
                                               Classifier
Decision Tree
Building the model using DecisionTree
from sklearn.tree import DecisionTreeClassifier
dt = DecisionTreeClassifier(max_depth=8, min_samples_split=4, min_samples_leaf=2)
dt.fit(x_train, y_train)
                                DecisionTreeClassifier
     DecisionTreeClassifier(max_depth=8, min_samples_leaf=2, min_samples_split=4)
 Random Forest
Building the model using Random Forest
from sklearn.ensemble import RandomForestClassifier
rf = RandomForestClassifier(n_estimators=100)
rf.fit(x_train, y_train)
     ▼ RandomForestClassifier
     RandomForestClassifier()
Getting the accuracy score for Decision Tree and Random Forest
from sklearn.metrics import accuracy_score
y_pred_train_dt = dt.predict(x_train)
y_pred_train_rf = rf.predict(x_train)
acc_train_dt = accuracy_score(y_train, y_pred_train_dt)
acc_train_rf = accuracy_score(y_train, y_pred_train_rf)
acc_train_dt, acc_train_rf
    (0.9071661237785016, 1.0)
y_pred_test_dt = dt.predict(x_test)
y_pred_test_rf = rf.predict(x_test)
acc_test_dt = accuracy_score(y_test, y_pred_test_dt)
acc_test_rf = accuracy_score(y_test, y_pred_test_rf)
acc_test_dt, acc_test_rf
    (0.7662337662337663, 0.7402597402597403)
Based on the information provided, it's observed that while the Random Forest model achieves a perfect accuracy of 100% on the training data,
its accuracy drops to 0.7467 on the test data. This indicates the model's tendency to overfit the training data. By introducing a maximum depth
parameter, we anticipate observing alterations in the model's performance.
from sklearn.ensemble import RandomForestClassifier
rf = RandomForestClassifier(n_estimators=100, max_depth=8)
rf.fit(x_train, y_train)
from sklearn.metrics import accuracy_score
y_pred_train_dt = dt.predict(x_train)
y_pred_train_rf = rf.predict(x_train)
acc_train_dt = accuracy_score(y_train, y_pred_train_dt)
acc_train_rf = accuracy_score(y_train, y_pred_train_rf)
acc_train_dt, acc_train_rf
    (0.9071661237785016, 0.9625407166123778)
y_pred_test_dt = dt.predict(x_test)
y_pred_test_rf = rf.predict(x_test)
acc_test_dt = accuracy_score(y_test, y_pred_test_dt)
acc_test_rf = accuracy_score(y_test, y_pred_test_rf)
acc_test_dt, acc_test_rf
    (0.7662337662337663, 0.7272727272727273)
from sklearn.metrics import recall_score
r_dt = recall_score(y_test, y_pred_test_dt)
r_rf = recall_score(y_test, y_pred_test_rf)
r_dt, r_rf
    (0.5740740740740741, 0.5370370370370371)
train_accuracies = [acc_train_dt, acc_train_rf]
test_accuracies = [acc_test_dt, acc_test_rf]
plt.figure(figsize=(10, 6))
plt.plot(['Decision Tree', 'Random Forest'], train_accuracies, marker='o', label='Training Accuracy')
plt.plot(['Decision Tree', 'Random Forest'], test_accuracies, marker='o', label='Testing Accuracy')
plt.title('Training and Testing Accuracies of Decision Tree and Random Forest Classifiers')
plt.xlabel('Classifier')
plt.ylabel('Accuracy')
plt.ylim(0, 1)
plt.legend()
plt.grid(True)
plt.show()
                    Training and Testing Accuracies of Decision Tree and Random Forest Classifiers
        0.8 -
        0.6
              --- Training Accuracy
              Testing Accuracy
        0.0
           Decision Tree
                                                                                                  Random Forest
                                                        Classifier
# Recall scores for decision tree and random forest classifiers
train_recalls = [recall_score(y_train, dt.predict(x_train)), recall_score(y_train, rf.predict(x_train))]
test_recalls = [recall_score(y_test, dt.predict(x_test)), recall_score(y_test, rf.predict(x_test))]
plt.figure(figsize=(10, 6))
plt.plot(['Decision Tree', 'Random Forest'], train_recalls, marker='o', label='Training Recall')
plt.plot(['Decision Tree', 'Random Forest'], test_recalls, marker='o', label='Testing Recall')
plt.title('Training and Testing Recall Scores of Decision Tree and Random Forest Classifiers')
plt.xlabel('Classifier')
plt.ylabel('Recall Score')
plt.ylim(0, 1)
plt.legend()
# Show plot
plt.grid(True)
plt.show()
                  Training and Testing Recall Scores of Decision Tree and Random Forest Classifiers
```

r = recall\_score(y\_test, y\_pred\_test)

Bayes model outperforms KNN on this dataset.

from sklearn.naive\_bayes import GaussianNB

from sklearn.metrics import recall\_score

from sklearn.neighbors import KNeighborsClassifier

When considering this code, if we contrast the recall metrics between the KNN and Naive Bayes models, it becomes apparent that the Naive

0.46296296296296297

import matplotlib.pyplot as plt

→ Training Recall --- Testing Recall 0.2 Decision Tree Random Forest Classifier

from sklearn.metrics import precision\_score p\_dt = precision\_score(y\_test, y\_pred\_test\_dt)

p\_rf = precision\_score(y\_test, y\_pred\_test\_rf)

p\_dt, p\_rf

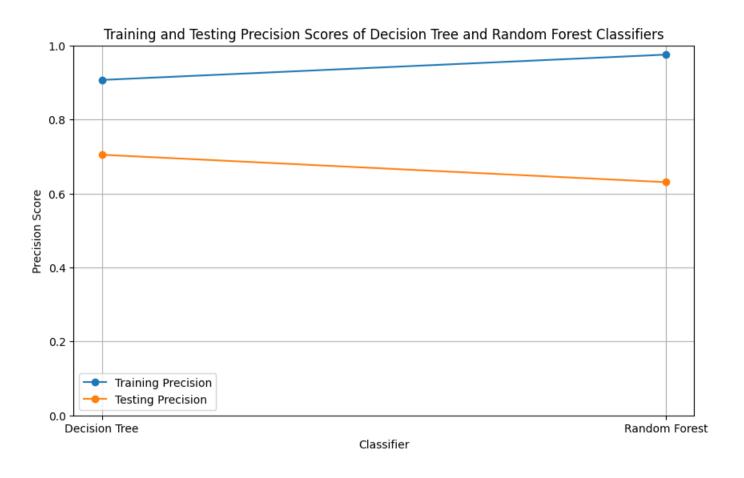
(0.7045454545454546, 0.6304347826086957)

```
train_precisions = [precision_score(y_train, dt.predict(x_train)), precision_score(y_train, rf.predict(x_train))]
test_precisions = [precision_score(y_test, dt.predict(x_test)), precision_score(y_test, rf.predict(x_test))]
```

plt.figure(figsize=(10, 6))

plt.plot(['Decision Tree', 'Random Forest'], train\_precisions, marker='o', label='Training Precision')
plt.plot(['Decision Tree', 'Random Forest'], test\_precisions, marker='o', label='Testing Precision')
plt.title('Training and Testing Precision Scores of Decision Tree and Random Forest Classifiers')
plt.xlabel('Classifier')
plt.ylabel('Precision Score')
plt.ylim(0, 1)
plt.legend()

plt.grid(True)
plt.show()



Based on the information provided earlier, it is evident that Random Forest outperforms the Decision Tree model on this dataset.

Support Vector Machine (SVM)

Building the model using Support Vector Machine (SVM) with linear kernel

from sklearn import svm
model = svm.SVC(kernel = 'linear')
model.fit(x\_train, y\_train)

▼ SVC

SVC(kernel='linear')

Prediction from support vector machine model on the traing and testing data

y\_pred\_train = model.predict(x\_train)
y\_pred\_test = model.predict(x\_test)

Accuracy score , recall\_score, and precision\_score for SVM

from sklearn.metrics import accuracy\_score, recall\_score, precision\_score

acc\_train = accuracy\_score(y\_train, y\_pred = y\_pred\_train)
acc\_test = accuracy\_score(y\_test, y\_pred = y\_pred\_test)

acc\_train, acc\_test

(0.7785016286644951, 0.7727272727272727)

p = precision\_score(y\_test, y\_pred\_test)
r = recall\_score(y\_test, y\_pred\_test)

p, r

(0.711111111111111, 0.5925925925925926)

Building the model using Support Vector Machine (SVM) with linear rbf

from sklearn import svm
model = svm.SVC(kernel = 'rbf')
model fit(x train, x train)

model.fit(x\_train, y\_train)
y\_pred\_train = model.predict(x\_train)

y\_pred\_train = model.predict(x\_train)
y\_pred\_test = model.predict(x\_test)
from sklearn.metrics import accuracy\_score, recall\_score, precision\_score

acc\_train = accuracy\_score(y\_train, y\_pred = y\_pred\_train)
acc\_test = accuracy\_score(y\_test, y\_pred = y\_pred\_test)

acc\_train, acc\_test

(0.8306188925081434, 0.7597402597402597)

p = precision\_score(y\_test, y\_pred\_test)
r = recall\_score(y\_test, y\_pred\_test)

p, r

(0.6976744186046512, 0.555555555555556)

from sklearn.linear\_model import LogisticRegression
model = LogisticRegression()

model.fit(x\_train, y\_train)

▼ LogisticRegression

LogisticRegression()

Logistic Regression

y\_pred\_train = model.predict(x\_train)
y\_pred\_test = model.predict(x\_test)

from sklearn.metrics import accuracy\_score, recall\_score, precision\_score

acc\_train = accuracy\_score(y\_train, y\_pred\_train)
acc\_test = accuracy\_score(y\_test, y\_pred\_test)

(0.8306188925081434, 0.7597402597402597)

acc\_train, acc\_test

p = precision\_score(y\_test, y\_pred\_test)
r = recall score(y test, y pred test)

r = recall\_score(y\_test, y\_pred\_test)
p, r

(0.6976744186046512, 0.555555555555556)

Comparing used models

import matplotlib.pyplot as plt
from sklearn.model\_selection import cross\_val\_score
from sklearn.naive\_bayes import GaussianNB
from sklearn.neighbors import KNeighborsClassifier
from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import RandomForestClassifier

from sklearn.linear\_model import LogisticRegression
# Initialize models
models = {

plt.grid(axis='y')

plt.show()

models = {
 'Naive Bayes': GaussianNB(),
 'KNN': KNeighborsClassifier(),
 'Parising Target Design Transferred

'Decision Tree': DecisionTreeClassifier(),
'Random Forest': RandomForestClassifier(),
'Logistic Regression': LogisticRegression()

'Logistic Regression': LogisticRegression()

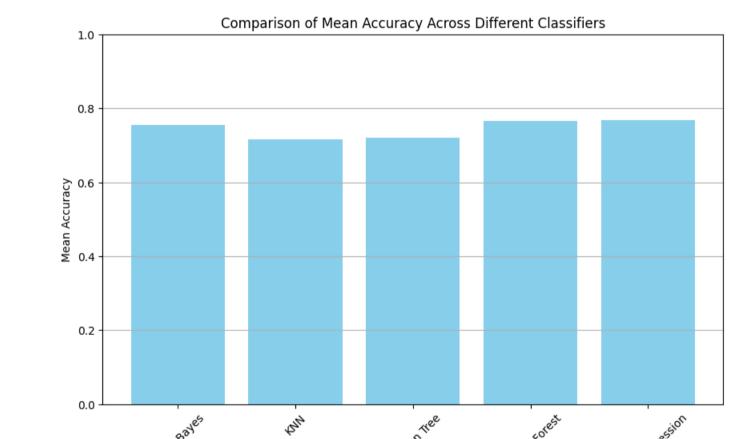
# Cross-validate each model and get accuracy scores
accuracy\_scores = {}
for model\_name, model in models.items():
 scores = cross\_val\_score(model, x\_train, y\_train, cv=5, scoring='accuracy')

accuracy\_scores[model\_name] = scores.mean()

# Plotting
plt.figure(figsize=(10, 6))

plt.bar(accuracy\_scores.keys(), accuracy\_scores.values(), color='skyblue')
plt.xlabel('Classifier')
plt.ylabel('Mean Accuracy')
plt.ylabel('Companies of Mean Accuracy Across Different Classifiers')

plt.title('Comparison of Mean Accuracy Across Different Classifiers')
plt.xticks(rotation=45)
plt.ylim(0, 1) # Limiting y-axis from 0 to 1 for better visualization



Conclusion

The aim of this study was to create classification models for the diabetes data set and to predict whether a person is sick by establishing models and to obtain maximum accurancy scores in the established models. Having utilized the entirety of the patient records, we've successfully constructed a machine learning model, particularly a Logestic Regrassion model, which effectively predicts the presence or absence of diabetes among the patients in the dataset. Additionally, through comprehensive data analysis and visualization, we've gleaned valuable insights from the data.

Classifier