

VISVESVARAYA TECHNOLOGICAL UNIVERSITY

“JnanaSangama”, Belgaum -590014, Karnataka.



LAB REPORT
on

Machine Learning (23CS6PCMAL)

Submitted by

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in partial fulfillment for the award of the degree of

BACHELOR OF ENGINEERING
in
COMPUTER SCIENCE AND ENGINEERING



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CERTIFICATE

This is to certify that the Lab work entitled “Machine Learning (23CS6PCMAL)” carried out by **Arya Himanshu (1BM22CS055)**, who is bonafide student of **B.M.S. College of Engineering**. It is in partial fulfillment for the award of **Bachelor of Engineering in Computer Science and Engineering** of the Visvesvaraya Technological University, Belgaum. The Lab report has been approved as it satisfies the academic requirements in respect of an Machine Learning (23CS6PCMAL) work prescribed for the said degree.

Sarala Assistant Professor Department of CSE, BMSCE	Dr. Kavitha Sooda Professor & HOD Department of CSE, BMSCE
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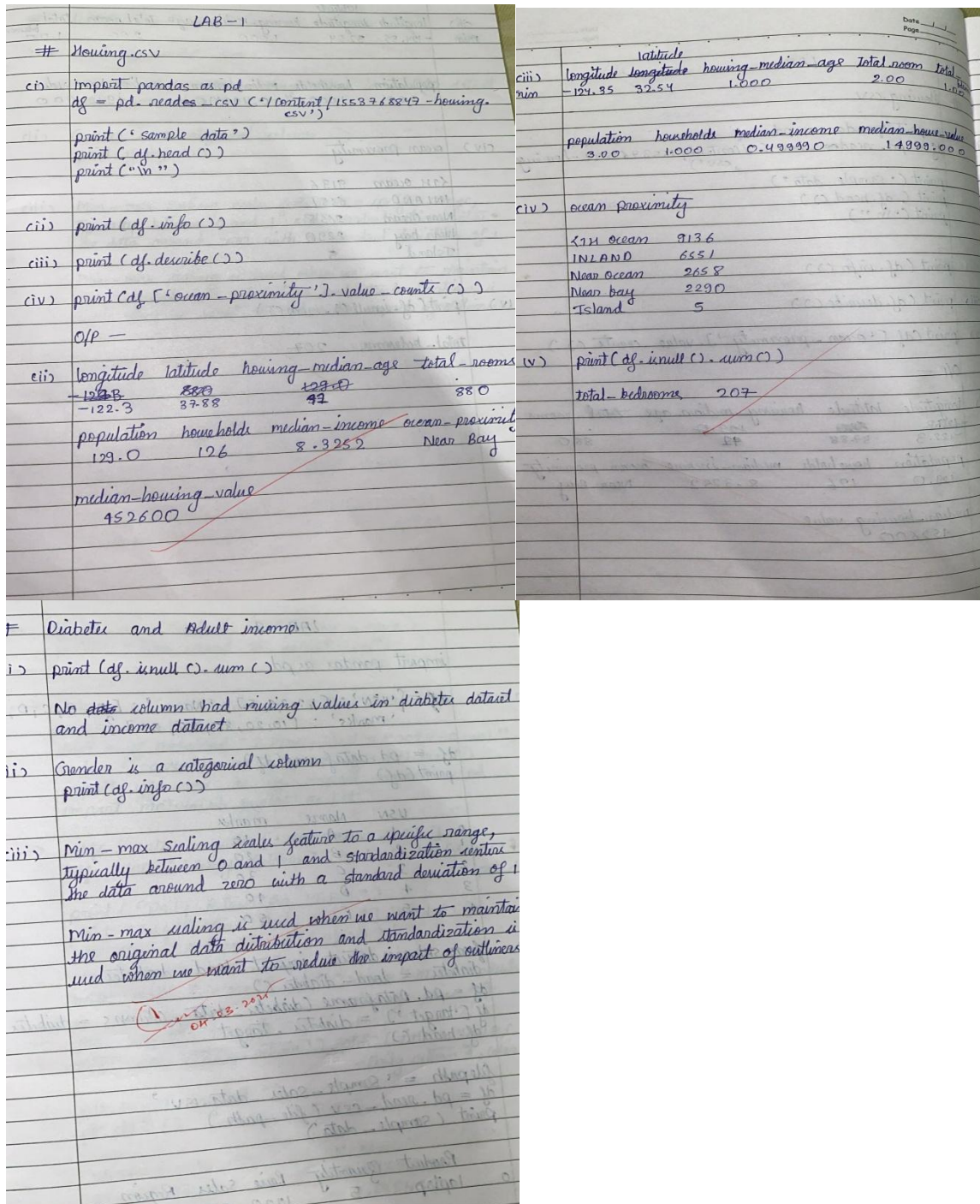
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Program 1

Write a python program to import and export data using Pandas library functions

Screenshot



Code:

```
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
from sklearn.model_selection import train_test_split
from sklearn.impute import SimpleImputer
from sklearn.preprocessing import OrdinalEncoder, OneHotEncoder
from sklearn.preprocessing import StandardScaler, MinMaxScaler
from scipy import stats

***Diabetes Dataset**
df=pd.read_csv('/content/Dataset of Diabetes .csv')
df.head()
df.shape
print(df.info())
# Summary statistics
print(df.describe())
missing_values=df.isnull().sum()
print(missing_values[missing_values > 0])
categorical_cols = df.select_dtypes(include=['object']).columns
print("Categorical columns identified:", categorical_cols)
if len(categorical_cols) > 0:
    df = pd.get_dummies(df, columns=categorical_cols, drop_first=True)
    print("\nDataFrame after one-hot encoding:")
    print(df.head())
else:
    print("\nNo categorical columns found in the dataset.")
from sklearn.preprocessing import MinMaxScaler, StandardScaler
import pandas as pd

numerical_cols = df.select_dtypes(include=['number']).columns

scaler = MinMaxScaler()
df_minmax = df.copy() # Create a copy to avoid modifying the original
df_minmax[numerical_cols] = scaler.fit_transform(df[numerical_cols])

scaler = StandardScaler()
df_standard = df.copy()
df_standard[numerical_cols] = scaler.fit_transform(df[numerical_cols])
print("\nDataFrame after Min-Max Scaling:")
print(df_minmax.head())
print("\nDataFrame after Standardization:")
print(df_standard.head())

***Adult Income Dataset**
df1=pd.read_csv('/content/adult.csv')
df1.head()
df1.shape
```

```

print(df1.info())
# Summary statistics
print(df.describe())
missing_values=df1.isnull().sum()
print(missing_values[missing_values > 0])
categorical_cols = df1.select_dtypes(include=['object']).columns
print("Categorical columns identified:", categorical_cols)
if len(categorical_cols) > 0:
    df1 = pd.get_dummies(df1, columns=categorical_cols, drop_first=True)
    print("\nDataFrame after one-hot encoding:")
    print(df1.head())
else:
    print("\nNo categorical columns found in the dataset.")
from sklearn.preprocessing import MinMaxScaler, StandardScaler
import pandas as pd

numerical_cols = df1.select_dtypes(include=['number']).columns

scaler = MinMaxScaler()
df_minmax = df1.copy() # Create a copy to avoid modifying the original
df_minmax[numerical_cols] = scaler.fit_transform(df1[numerical_cols])

scaler = StandardScaler()
df_standard = df1.copy()
df_standard[numerical_cols] = scaler.fit_transform(df1[numerical_cols])
print("\nDataFrame after Min-Max Scaling:")
print(df_minmax.head())
print("\nDataFrame after Standardization:")
print(df_standard.head())

```

PROGRAM 2 Demonstrate various data pre-processing techniques for a given dataset

Screenshot

LAB-3

```
import pandas as pd
df = {'USN': [1, 2, 3, 4, 5], 'Name': ['A', 'B', 'C', 'D', 'E'],
      'marks': [10, 20, 30, 40, 50]}
df = pd.DataFrame(df)
print(df)
```

	USN	Name	marks
0	1	A	10
1	2	B	20
2	3	C	30
3	4	D	40
4	5	E	50

```
from sklearn.datasets import load_diabetes
diabetes = load_diabetes()
df = pd.DataFrame(diabetes.data, columns = diabetes.feature_names)
df['target'] = diabetes.target
df.head()
```

```
filepath = 'sample-sales-data.csv'
df = pd.read_csv(filepath)
print(sample_data)
```

	Product	Quantity	Price	Sales	Region
0	laptop	5	1000	5000	North

```
df.to_csv('output.csv', index = False)
print("Data saved to output.csv")

-> tickers = ['HDFCBANK.NS', 'ICICIBANK.NS', 'KOTAKBANK.NS']
data = yf.download(tickers, start = "2024-01-01",
                    end = "2024-12-30", group_by = 'tickers')
import yfinance as yf, import pandas as pd
import matplotlib.pyplot as plt

-> hdfc_data = data['HDFCBANK.NS']
hdfc_data['Daily-Return'] = hdfc_data['close'].pct_change()

print("Daily Return for HDFC")
print(hdfc_data['Daily-Return'].head())
plt.figure(figsize = (10, 6))
plt.subplot(2, 1, 1)
hdfc_data['close'].plot(title = "HDFC Industries",
                        label = "closing price")
plt.subplot(2, 1, 2)
hdfc_data['Daily-Return'].plot(title = "HDFC Industries",
                               label = "Daily return", color = 'red')

plt.tight_layout()
plt.show()
```

Date	close	Daily Return
2024-01-01	NAN	
2024-01-02	1542.0	0.015420
2024-01-03	1542.0	-0.015420

Code

```
import pandas as pd

import numpy as np

import seaborn as sns

import matplotlib.pyplot as plt

df=pd.read_csv('housing.csv')

df.head(2)

df.describe()

df.info()

sns.histplot(df['median_income'], kde=True, color='green')

sns.histplot(df['housing_median_age'])

from sklearn.model_selection import train_test_split

X = df.drop("median_house_value", axis=1)

y = df["median_house_value"]

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

X = df.drop("median_house_value", axis=1)

y = df["median_house_value"]

df["income_cat"] = pd.cut(df["median_house_value"],

bins=[0, 100000, 200000, 300000, 400000, np.inf],

labels=[1, 2, 3, 4, 5])

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42,

stratify=df["income_cat"])
```



```

train_set = X_train.copy()

train_set["median_house_value"] = y_train

train_set.plot(kind="scatter", x="longitude", y="latitude", alpha=0.4,s=train_set["population"]/100,
label="population",figsize=(10,7), c="median_house_value", cmap=plt.get_cmap("jet"),

colorbar=True)

plt.legend()

numerical_columns = df.select_dtypes(include=['float64', 'int64'])

correlation_matrix = numerical_columns.corr()

print(correlation_matrix["median_house_value"].sort_values(ascending=False))

df.plot(kind="scatter", x="median_income", y="median_house_value", alpha=0.1)

# Combine 'median_income' and 'households'

df["income_households"] = df["median_income"] * df["households"]


numerical_columns = df.select_dtypes(include=['float64', 'int64'])

correlation_matrix = numerical_columns.corr()

print(correlation_matrix["median_house_value"].sort_values(ascending=False))

df.plot(kind="scatter", x="income_households", y="median_house_value", alpha=0.1)

plt.show()

missing_values = df.isnull().sum()

print(missing_values[missing_values > 0])

h=df

h.dropna(subset=["total_bedrooms"])

from sklearn.preprocessing import OneHotEncoder

df1=pd.read_csv('housing.csv')

hc=df1[["ocean_proximity"]]

```

```

encoder=OneHotEncoder()

hc_encoded=encoder.fit_transform(hc).toarray()

hc_1hot_df = pd.DataFrame(hc_encoded, columns=encoder.get_feature_names_out(hc.columns))

hc_1hot_df.head()

```

Feature scaling is crucial in machine learning for several reasons, particularly when using algorithms that are sensitive to the scale of features. Here's a breakdown of its importance:

1. Improved Performance of Distance-Based Algorithms:

2. Faster Convergence of Gradient Descent:

3. Improved Regularization:

4. Better Interpretation of Coefficients:

5. Numerical Stability:

```

from sklearn.base import BaseEstimator, TransformerMixin

```

```

from sklearn.pipeline import Pipeline

```

```

from sklearn.compose import ColumnTransformer

```

```

from sklearn.preprocessing import StandardScaler

```

```

# Custom transformer to add engineered attributes

```

```

class CombinedAttributesAdder(BaseEstimator, TransformerMixin):

```

```

    def __init__(self, add_bedrooms_per_room=True):

```

```

self.add_bedrooms_per_room = add_bedrooms_per_room

def fit(self, X, y=None):

    return self

def transform(self, X):

    # Assumes X is a NumPy array with the following columns:

    # total_rooms (index 3), total_bedrooms (index 2), population (index 4), households (index 5)

    rooms_per_household = X[:, 3] / X[:, 5]

    population_per_household = X[:, 4] / X[:, 5]

    if self.add_bedrooms_per_room:

        bedrooms_per_room = X[:, 2] / X[:, 3]

        return np.c_[X, rooms_per_household, population_per_household, bedrooms_per_room]

    else:

        return np.c_[X, rooms_per_household, population_per_household]

# Identify numerical and categorical columns

num_attribs = df1.drop("ocean_proximity", axis=1).columns # All numeric columns

cat_attribs = ["ocean_proximity"]

# Build numerical pipeline: impute missing values, add new attributes, then scale

num_pipeline = Pipeline([

    ('imputer', SimpleImputer(strategy="median")),

    ('attribs_adder', CombinedAttributesAdder()),

    ('std_scaler', StandardScaler()),

```

```
# Build the full pipeline combining numerical and categorical processing
```

```
full_pipeline = ColumnTransformer([  
    ("num", num_pipeline, num_attribs),  
    ("cat", OneHotEncoder(), cat_attribs),
```

```
# Process the dataset using the pipeline
```

```
housing_prepared = full_pipeline.fit_transform(housing)  
print("Shape of processed data:", housing_prepared.shape)
```

PROGRAM 3 Implement Linear and Multi-Linear Regression algorithm using appropriate dataset

Screenshot

lab - 4

Linear Regression

week x_i	Sales y_i
1	2
2	4
3	5
4	9

$$x^T = \begin{bmatrix} 1 & 1 & 1 & 1 \\ 1 & 2 & 3 & 4 \end{bmatrix}$$

$$y^T = \begin{bmatrix} 2 & 4 & 5 & 9 \end{bmatrix}$$

$$X = \begin{bmatrix} 1 & 1 \\ 1 & 2 \\ 1 & 3 \\ 1 & 4 \end{bmatrix}, Y = \begin{bmatrix} 2 \\ 4 \\ 5 \\ 9 \end{bmatrix}$$

$$X^T X = \begin{bmatrix} 4 & 10 \\ 10 & 30 \end{bmatrix}$$

$$(X^T X)^{-1} = \begin{bmatrix} 1.5 & -0.5 \\ -0.5 & 0.2 \end{bmatrix}$$

$$(X^T X)^{-1} X^T = \begin{bmatrix} 1.0 & 0.5 & 0 & -0.5 \\ -0.3 & -0.1 & 0.1 & 0.3 \end{bmatrix}$$

$((X^T X)^{-1} X^T) Y = \begin{bmatrix} 1.0 & 0.5 & 0 & -0.5 \\ -0.3 & -0.1 & 0.1 & 0.3 \end{bmatrix} \begin{bmatrix} 2 \\ 4 \\ 5 \\ 9 \end{bmatrix}$

$= \begin{bmatrix} -0.5 \\ 2.2 \end{bmatrix}$

$y = a_0 + a_1 x_1 + a_2 x_2$

for $x = 5$

$y = -0.5 + (2.2) \times 5$

$= 10.5$

* Steps (algo)

- 1) import libraries
- 2) import data - distribution
- 3) Analyse data - distribution
- 4) Distribution plot - visualisation
- 5) Relationship b/w variables
- 6) split the data
- 7) Train the model
- 8) Predict the result
- 9) Visualize prediction
- 10) Values of coefficient and intercept

11.05.2021

Code

```
# -*- coding: utf-8 -*-

import pandas as pd

import numpy as np

from sklearn import linear_model

import matplotlib.pyplot as plt


df = pd.read_csv('/content/housing_area_price.csv')

df

# Commented out IPython magic to ensure Python compatibility.

# %matplotlib inline

plt.xlabel('area')

plt.ylabel('price')

plt.scatter(df.area,df.price,color='red',marker='+')


new_df = df.drop('price',axis='columns')

new_df


price = df.price

price

# Create linear regression object

reg = linear_model.LinearRegression()

reg.fit(new_df,price)
```

```
"""(1) Predict price of a home with area = 3300 sqr ft"""
```

```
reg.predict([[3300]])
```

```
reg.coef_
```

```
reg.intercept_
```

```
"""Y = m * X + b (m is coefficient and b is intercept)"""
```

```
3300*135.78767123 + 180616.43835616432
```

```
"""(1) Predict price of a home with area = 5000 sqr ft"""
```

```
reg.predict([[5000]])
```

```
# -*- coding: utf-8 -*-
```

```
import pandas as pd
```

```
import numpy as np
```

```
from sklearn import linear_model
```

```
df = pd.read_csv('/content/homeprices_Multiple_LR.csv')
```

```
df
```


"""Data Preprocessing: Fill NA values with median value of a column"""

```
df.bedrooms.median()
```

```
df.bedrooms = df.bedrooms.fillna(df.bedrooms.median())
```

```
df
```

```
reg = linear_model.LinearRegression()
```

```
reg.fit(df.drop('price',axis='columns'),df.price)
```

```
reg.coef_
```

```
reg.intercept_
```

"""Find price of home with 3000 sqr ft area, 3 bedrooms, 40 year old"""

```
reg.predict([[3000, 3, 40]])
```

```
112.06244194*3000 + 23388.88007794*3 + -3231.71790863*40 + 221323.00186540384
```

```
import pandas as pd
```

```
from sklearn.linear_model import LinearRegression
```

```
# Load the dataset
```

```
df1 = pd.read_csv('/content/canada_per_capita_income.csv')
```

```

# Prepare the data

X = df1.year.values.reshape(-1, 1) # Features (year)

y = df1['per capita income (US$)'] # Target (per capita income)


# Create and train the linear regression model

model = LinearRegression()

model.fit(X, y)


# Predict per capita income for 2020

year_2020 = [[2020]]

predicted_income = model.predict(year_2020)


print(f"Predicted per capita income for Canada in 2020: {predicted_income[0]:.2f}")


import pandas as pd

from sklearn.linear_model import LinearRegression

import matplotlib.pyplot as plt


# Load the dataset (canada_per_capita_income.csv)

df1 = pd.read_csv('/content/canada_per_capita_income.csv')


# Prepare the data

X = df1.year.values.reshape(-1, 1) # Features (year)

```

```

y = df1['per capita income (US$)'] # Target (per capita income)

# Create and train the linear regression model

model = LinearRegression()

model.fit(X, y)

# Create the plot

plt.figure(figsize=(8, 6))

plt.scatter(X, y, color='blue', label='Data Points') # Now using the correct X and y

plt.plot(X, model.predict(X), color='red', label='Regression Line')

plt.xlabel('Year')

plt.ylabel('Per Capita Income (US$)')

plt.title('Per Capita Income in Canada over Time')

plt.legend()

plt.grid(True)

plt.show()

import pandas as pd

from sklearn.linear_model import LinearRegression

from sklearn.impute import SimpleImputer

# Load the dataset

df = pd.read_csv('/content/salary.csv')

# Prepare the data

```

```

X = df.iloc[:, :-1].values # Features (years of experience)

y = df.iloc[:, 1].values # Target (salary)


# Impute missing values with the mean

imputer = SimpleImputer(strategy='mean') # Create an imputer object with strategy as mean

X = imputer.fit_transform(X) # Fit and transform the imputer on feature data 'X'


# Create and train the linear regression model

model = LinearRegression()

model.fit(X, y)


# Predict salary for 12 years of experience

years_experience = [[12]]

predicted_salary = model.predict(years_experience)


print(f"Predicted salary for 12 years of experience: {predicted_salary[0]:.2f}")

import pandas as pd

from sklearn.linear_model import LinearRegression

from sklearn.impute import SimpleImputer


# Load the dataset

df = pd.read_csv('/content/hiring.csv')


# Handle missing values

```

```

# Convert 'experience' column to numeric, replacing non-numeric with NaN
df['experience'] = pd.to_numeric(df['experience'], errors='coerce')

imputer = SimpleImputer(strategy='mean')

df['experience'] = imputer.fit_transform(df[['experience']])
df['test_score(out of 10)'] = imputer.fit_transform(df[['test_score(out of 10)']])

# Prepare the data
X = df.drop('salary($)', axis='columns')
y = df['salary($)']

# Create and train the linear regression model
model = LinearRegression()
model.fit(X, y)

# Predict salaries for the given candidates
candidate1 = [[2, 9, 6]]
candidate2 = [[12, 10, 10]]

predicted_salary1 = model.predict(candidate1)
predicted_salary2 = model.predict(candidate2)

print(f"Predicted salary for candidate 1: ${predicted_salary1[0]:.2f}")
print(f"Predicted salary for candidate 2: ${predicted_salary2[0]:.2f}")

```

```

import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.linear_model import LinearRegression

from sklearn.preprocessing import LabelEncoder, OneHotEncoder

from sklearn.compose import ColumnTransformer


# Load the dataset

df = pd.read_csv('/content/1000_Companies.csv')


# Separate features (X) and target (y)

X = df.iloc[:, :-1].values

y = df.iloc[:, 4].values


# Encode categorical data (State)

labelencoder = LabelEncoder()

X[:, 3] = labelencoder.fit_transform(X[:, 3])

ct = ColumnTransformer(

    transformers=[('encoder', OneHotEncoder(), [3])],

    remainder='passthrough'

)

X = ct.fit_transform(X)


# Avoid dummy variable trap (remove one encoded column)

```

```
X = X[:, 1:]
```

```
# Split data into training and testing sets
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=0)
```

```
# Create and train the multiple linear regression model
```

```
regressor = LinearRegression()
```

```
regressor.fit(X_train, y_train)
```

```
# Predict profit for the given values
```

```
new_prediction = regressor.predict([[1, 0, 91694.48, 515841.3, 11931.24]])
```

```
print(f"Predicted Profit: {new_prediction[0]:.2f}")
```

PROGRAM 4 Build Logistic Regression Model for a given dataset

Screenshot

lab = 3

i) Given $a_0 = -5$, $a_1 = 0.8$

ii) logistic regression equation

$$p(x) = \frac{1}{1 + e^{-(a_0 + a_1 x)}} = \frac{1}{1 + e^{-(-5 + 0.8x)}}$$

iii) Calc probability that a student who studies 7 hrs will pass

→ $x = 7$ $p(x) = \frac{1}{1 + e^{-(-5 + 0.8(7))}} = 0.6457$

iv) Determine the predicted class (P.F) for this student based on threshold of 0.5

→ $P(x) = 0.6457$
 $P(x) > 0.5$
 Thus $y = 1$ (pass)

2) Consider $z = [2, 1, 0]$ for three classes apply to find probabilities values of 3 classes.

$$\text{Softmax}(z_i) = \frac{e^{z_i}}{\sum_{j=1}^n e^{z_j}}$$

Softmax(z_1) = $\frac{e^2}{e^2 + e^1 + e^0} = 0.645$

Softmax(z_2) = $\frac{e^1}{e^2 + e^1 + e^0} = 0.247$

Softmax(z_3) = $\frac{e^0}{e^2 + e^1 + e^0} = 0.091$

probabilities of the 3 classes are approx 64.5%, 24.7% and 9.1%.

⇒ Binary logistic Regression

import pandas as pd
 from matplotlib import pyplot as plt
 df = pd.read_csv('content/insurance.csv')
 df.head()

fit scatter of age, df.bought, insurance, marked = 1
 color = red

from sklearn.model_selection import train_test_split
 x_train, x_test, y_train, y_test = train_test_split(df[['age', 'bought', 'insurance']], df['marked'], test_size=0.3, random_state=0)

x_train
 from sklearn.linear_model import LogisticRegression
 model = LogisticRegression()
 model.fit(x_train, y_train)
 x_test
 y_test
 x_predicted = model.predict(x_test)
 y_predicted
 model.coef_
 model.intercept_

import math
 def sigmoid(z):
 return 1 / (1 + math.exp(-x))

def prediction_func(age):
 z = 0.12 * age - 4.823
 y = sigmoid(z)
 return(y)

age = 35
 prediction = function(age)
 o/p — 0.3709894

⇒ Multiclass logistic Regression

import pandas as pd
 from sklearn.datasets import load_iris
 from sklearn.model_selection import train_test_split
 from sklearn.linear_model import LogisticRegression
 from sklearn.metrics import accuracy_score

iris = pd.read_csv('content/iris.csv')
 iris.head()
 X = iris.data['species', axis = 'columns']
 y = iris.target
 X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=0)

model = LogisticRegression(multi_class='multinomial')
 model.fit(X_train, y_train)
 y_pred = model.predict(X_test)
 accuracy = accuracy_score(y_test, y_pred)

print('Accuracy on the test set: {accuracy: 27.9%}')
 Confusion matrix = metrics.confusion_matrix(y_test, y_pred)

`cm = display.plot()`
`plt.show()`
 O/P —
 accuracy on the train set : 1.00
 accuracy : 85.2%
 it is a good accuracy, but accuracy alone can
 be misleading due to potential class imbalance.
 o dataset
 performed data processing steps.
 used the animal name column, standardized the features
 using standard scaler to improve model performance.

cii) There were no missing or inconsistent values
 ciii) The confusion matrix showed achieved 95.2% accuracy.
 civ) Reptiles and amphibians were mostly misclassified classes. They share common features. Mammals and birds also as they had overlapping features.

Code

```
import pandas as pd

import numpy as np

df=pd.read_csv("/content/HR_comma_sep.csv")

df.head(3)

print(df.isnull().sum())

print(df.groupby('left').mean(numeric_only=True))

print(df.groupby('salary').mean(numeric_only=True))

import matplotlib.pyplot as plt

pd.crosstab(df.salary,df.left).plot(kind='bar')

plt.title('Employee Retention vs Salary')

plt.xlabel('Salary')

plt.ylabel('Number of Employees')

plt.show()

pd.crosstab(df.Department,df.left).plot(kind='bar')

plt.title('Employee Retention vs Department')

plt.xlabel('Department')

plt.ylabel('Number of Employees')

plt.show()

salary_dummies = pd.get_dummies(df.salary, prefix="salary")

dept_dummies = pd.get_dummies(df.Department, prefix="dept")
```

```
df_with_dummies = pd.concat([df, salary_dummies, dept_dummies], axis=1)
```

```
df_with_dummies = df_with_dummies.drop(['salary', 'Department'], axis=1)
```

```
X_features = ['satisfaction_level', 'last_evaluation', 'number_project', 'average_monthly_hours',  
'time_spend_company', 'Work_accident', 'promotion_last_5years'] + list(salary_dummies.columns) +  
list(dept_dummies.columns)
```

```
X = df_with_dummies[X_features]
```

```
y = df_with_dummies.left
```

```
from sklearn.model_selection import train_test_split
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=0)
```

```
from sklearn.linear_model import LogisticRegression
```

```
model = LogisticRegression()
```

```
model.fit(X_train, y_train)
```

```
from sklearn.metrics import accuracy_score
```

```
y_pred = model.predict(X_test)
```

```
accuracy = accuracy_score(y_test, y_pred)
```

```
print("Accuracy of the model:", accuracy)
```

PROGRAM 5 Use an appropriate data set for building the decision tree (ID3) and apply this knowledge to classify a new sample.

Screenshot

Decision Tree

Instance	a_1	a_2	Classification
1	Hot	High	No
2	Hot	High	No
3	Cool	High	No
4	Hot	High	No
5	Hot	Normal	Yes

Entropy (g) = $-\frac{4}{5} \log_2(\frac{4}{5}) - \frac{1}{5} \log_2(\frac{1}{5})$
 $= 0.7219$

for a_1 :

Split (1+3-2) = $-\frac{4}{4} \log_2(\frac{4}{4}) - \frac{3}{4} \log_2(\frac{3}{4})$
 $S_{cool}(1+3) = 0$

Gain(S, a_1) = $0.7219 - \frac{4}{5} \times 0.813 = 0.0728$

for a_2 :

Split (1+4-3) = $-\frac{4}{4} \log_2(\frac{4}{4}) - \frac{3}{4} \log_2(\frac{3}{4})$
 $S_{high}(1+4) = 0$
 $S_{normal}(1+4) = 0$

Gain(S, a_2) = $0.7219 - 0 - 0 = 0.7219$

Decision Tree Diagram:

```

graph TD
    A((a2)) -- High --> B((1,4,3))
    A -- Normal --> C((5))
    B --> D[No]
    C --> E[Yes]
  
```

Decision Tree

import pandas as pd
 from sklearn.tree import DecisionTreeClassifier
 from sklearn.metrics import accuracy_score
 import matplotlib.pyplot as plt

df = pd.read_csv('Content/drug.csv')
 X = df.drop('Drug', axis=1)
 y = df['Drug']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2)

dtree = DecisionTreeClassifier()
 dtree.fit(X_train, y_train)
 y_pred = dtree.predict(X_test)

accuracy = accuracy_score(y_test, y_pred)
 print("Accuracy: %f" % accuracy)

cm = confusion_matrix(y_test, y_pred)
 plt.figure(figsize=(8, 8))
 plt.show()

O/P —

Accuracy: 1.0

Confusion Matrix:

	Actual	Predicted
Actual	6 0 0 0 0	6 0 0 0 0
0 3 0 0 0	0 0 5 0 0	0 0 0 11 0
0 0 0 0 15		

⇒ O/P —

Mean absolute error: 59.6
 Mean squared error: 4338.2
 Root mean squared error: 39.6128

Q.1. Accuracy was 1.0 for this dataset.

The Confusion matrix only has diagonal elements, the model made zero errors.

Key features like population density and average income are crucial for predicting petrol consumption.

Code

```
from sklearn.datasets import load_iris

from sklearn.model_selection import train_test_split

from sklearn.tree import DecisionTreeClassifier

from sklearn.metrics import accuracy_score, confusion_matrix

from sklearn import tree

import matplotlib.pyplot as plt


iris = load_iris()

X = iris.data

y = iris.target


X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)


clf = DecisionTreeClassifier()


clf.fit(X_train, y_train)


y_pred = clf.predict(X_test)


accuracy = accuracy_score(y_test, y_pred)

conf_matrix = confusion_matrix(y_test, y_pred)


print("Accuracy:", accuracy)
```

```
print("Confusion Matrix:\n", conf_matrix)
```

```
plt.figure(figsize=(12, 8))
```

```
tree.plot_tree(clf, feature_names=iris.feature_names, class_names=iris.target_names, filled=True)
```

```
plt.show()
```

```
from sklearn.datasets import load_iris
```

```
from sklearn.model_selection import train_test_split
```

```
from sklearn.tree import DecisionTreeClassifier
```

```
from sklearn.metrics import accuracy_score, confusion_matrix
```

```
from sklearn import tree
```

```
import matplotlib.pyplot as plt
```

```
iris = load_iris()
```

```
X = iris.data
```

```
y = iris.target
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)
```

```
clf = DecisionTreeClassifier()
```

```
clf.fit(X_train, y_train)
```

```
y_pred = clf.predict(X_test)
```

```
accuracy = accuracy_score(y_test, y_pred)
```

```
conf_matrix = confusion_matrix(y_test, y_pred)
```

```
print("Accuracy:", accuracy)
```

```
print("Confusion Matrix:\n", conf_matrix)
```

```
plt.figure(figsize=(12, 8))
```

```
tree.plot_tree(clf, feature_names=iris.feature_names, class_names=iris.target_names, filled=True)
```

```
plt.show()
```

```
import pandas as pd
```

```
from sklearn.model_selection import train_test_split
```

```
from sklearn.tree import DecisionTreeRegressor
```

```
from sklearn.metrics import mean_absolute_error, mean_squared_error
```

```
import numpy as np # import numpy
```

```
data = pd.read_csv("petrol_consumption.csv")
```

```
X = data[['Petrol_tax', 'Average_income', 'Paved_Highways',
```

```
         'Population_Driver_licence(%)']]
```

```
y = data['Petrol_Consumption']
```

```

X_train, X_test, y_train, y_test = train_test_split(
    X, y, test_size=0.2, random_state=42)

regressor = DecisionTreeRegressor()

regressor.fit(X_train, y_train)

y_pred = regressor.predict(X_test)

mae = mean_absolute_error(y_test, y_pred)
mse = mean_squared_error(y_test, y_pred)
rmse = np.sqrt(mse)

print("Mean Absolute Error:", mae)
print("Mean Squared Error:", mse)
print("Root Mean Squared Error:", rmse)

from sklearn.tree import plot_tree
import matplotlib.pyplot as plt

plt.figure(figsize=(15, 10))

# Assuming 'data' is your original pandas DataFrame
plot_tree(regressor, feature_names=data[['Petrol_tax', 'Average_income', 'Paved_Highways',
'Population_Driver_licence(%)']].columns, filled=True, rounded=True)

plt.show()

```


PROGRAM 6 Build KNN Classification model for a given dataset.

Screenshot

Pearson	Age	Salary	Target	Distance
A	18	50	N	52.8
B	23	55	N	44.59
C	24	70	N	31.95
D	41	60	Y	40.44
E	43	70	Y	31.04
F	38	40	Y	10.07
X	35	100	?	

Distance = $\sqrt{(x_0 - x)^2 + (y_0 - y)^2}$

for $k=3$, Rank in ascending order

Distance	Rank	Target
31.04	1	Y
31.95	2	Y
40.44	3	Y

Since majority is yes, for (35, 100) target will be yes.

Lab-6

```

import pandas as pd
import seaborn as sns
from sklearn.model_selection import train_test_split
from sklearn.neighbors import KNeighborsClassifier

iris_data = pd.read_csv('iris-data.csv')
X = iris_data.iloc[:, :-1]
y = iris_data.iloc[:, -1]
X_train = X, X_test = X, y_train = y, y_test = y
test_size = 0.2, random_state=0

k = 3
knn_classifier = KNeighborsClassifier(n_neighbors=k)
knn_classifier.fit(X_train, y_train)

accuracy = accuracy_score(y_test, y_pred = knn_classifier.predict(X_test))
conf_matrix = confusion_matrix(y_test, y_pred = knn_classifier.predict(X_test))
class_report = classification_report(y_test, y_pred = knn_classifier.predict(X_test))

print(f"Accuracy score: {accuracy}")
print(conf_matrix)
print(class_report)

```

O/p —

Accuracy score: 1.0

Confusion :

10	0	0
0	9	0
0	0	11

Classification Report				
	Precision	Recall	f1-score	Support
Setosa	1.00	1.00	1.00	10
Versicolour	1.00	1.00	1.00	9
Virginica	1.00	1.00	1.00	11
accuracy	1.00	1.00	1.00	30
macro avg	1.00	1.00	1.00	30
weighted avg	1.00	1.00	1.00	30

1. $k=3$ gives 100% accuracy here but generally takes $k=5$, $\text{error rate} = 1 - \text{accuracy} = 0.0$ (no misclassification to find that k , test multiple value & plot the error rate)

It is needed because features have diff ranges (eg glucose vs BMI), standardisation scales ensure equal features contribute by normalising data. Improves KNN performance (accuracy = 70.07% after scaling).

Code

```
import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.neighbors import KNeighborsClassifier

from sklearn.metrics import classification_report, confusion_matrix, accuracy_score

import seaborn as sns

import matplotlib.pyplot as plt


try:

    data = pd.read_csv('/content/iris (1).csv')

except FileNotFoundError:

    print("Error: 'iris.csv' not found. Please upload the file to your Colab environment.")

    exit()


X = data.drop('species', axis=1)

y = data['species']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

knn = KNeighborsClassifier(n_neighbors=3)

knn.fit(X_train, y_train)

y_pred = knn.predict(X_test)

print("Accuracy Score:", accuracy_score(y_test, y_pred))

print("\nConfusion Matrix:")

cm = confusion_matrix(y_test, y_pred)

print(cm)
```

```

plt.figure(figsize=(8, 6))

sns.heatmap(cm, annot=True, fmt='d', cmap='Blues',

            xticklabels=knn.classes_, yticklabels=knn.classes_)

plt.xlabel('Predicted')

plt.ylabel('Actual')

plt.title('Confusion Matrix')

plt.show()


print("\nClassification Report:")

print(classification_report(y_test, y_pred))


import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.neighbors import KNeighborsClassifier

from sklearn.metrics import classification_report, confusion_matrix, accuracy_score

from sklearn.preprocessing import StandardScaler

import seaborn as sns

import matplotlib.pyplot as plt


try:

    diabetes = pd.read_csv('diabetes.csv')

except FileNotFoundError:

    print("Error: 'diabetes.csv' not found. Please ensure the file is in the current directory.")

```

```

exit()

X = diabetes.drop('Outcome', axis=1)

y = diabetes['Outcome']

scaler = StandardScaler()

X = scaler.fit_transform(X)

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

knn = KNeighborsClassifier(n_neighbors=5)

knn.fit(X_train, y_train)

y_pred = knn.predict(X_test)

accuracy = accuracy_score(y_test, y_pred)

print(f"Accuracy: {accuracy}")

cm = confusion_matrix(y_test, y_pred)

print("Confusion Matrix:")

print(cm)

sns.heatmap(cm, annot=True, fmt="d")

plt.title('Confusion Matrix')

plt.xlabel('Predicted')

plt.ylabel('True')

plt.show()

print("Classification Report:")

print(classification_report(y_test, y_pred))

```

```

import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.neighbors import KNeighborsClassifier

from sklearn.metrics import classification_report, confusion_matrix, accuracy_score

import seaborn as sns

import matplotlib.pyplot as plt

from sklearn.preprocessing import StandardScaler

try:

    heart = pd.read_csv('heart.csv')

except FileNotFoundError:

    print("Error: 'heart.csv' not found. Please ensure the file is in the current directory.")

    exit()

X = heart.drop('target', axis=1)

y = heart['target']

scaler = StandardScaler()

X = scaler.fit_transform(X)

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

best_k = 1

best_accuracy = 0

```

```

for k in range(1, 21):

    knn = KNeighborsClassifier(n_neighbors=k)

    knn.fit(X_train, y_train)

    y_pred = knn.predict(X_test)

    accuracy = accuracy_score(y_test, y_pred)

    if accuracy > best_accuracy:

        best_accuracy = accuracy

        best_k = k

print(f"Best k: {best_k} with accuracy {best_accuracy}")

knn = KNeighborsClassifier(n_neighbors=best_k)

knn.fit(X_train, y_train)

y_pred = knn.predict(X_test)

accuracy = accuracy_score(y_test, y_pred)

print(f"Accuracy: {accuracy}")

cm = confusion_matrix(y_test, y_pred)

print("Confusion Matrix:")

print(cm)

sns.heatmap(cm, annot=True, fmt="d")

plt.title('Confusion Matrix')

plt.xlabel('Predicted')

plt.ylabel('True')

```

```
plt.show()
```

```
print("Classification Report:")
```

```
print(classification_report(y_test, y_pred))
```

```
import matplotlib.pyplot as plt
```

```
import seaborn as sns
```

```
from sklearn.metrics import classification_report, confusion_matrix
```

```
cm = confusion_matrix(y_test, y_pred)
```

```
plt.figure(figsize=(8, 6))
```

```
sns.heatmap(cm, annot=True, fmt="d", cmap="Blues")
```

```
plt.title("Confusion Matrix")
```

```
plt.xlabel("Predicted")
```

```
plt.ylabel("Actual")
```

```
plt.show()
```

```
print(classification_report(y_test, y_pred))
```

```
# prompt: For Iris dataset
```

```
# How to choose the k value? Demonstrate using accuracy rate and error
```

```
# rate. Give theory
```

```
import pandas as pd
```

```
from sklearn.model_selection import train_test_split
```

```

from sklearn.neighbors import KNeighborsClassifier

from sklearn.metrics import classification_report, confusion_matrix, accuracy_score

import seaborn as sns

import matplotlib.pyplot as plt

from sklearn.preprocessing import StandardScaler


# Load the Iris dataset

try:

    data = pd.read_csv('/content/iris (1).csv')

except FileNotFoundError:

    print("Error: 'iris (1).csv' not found. Please upload the file to your Colab environment.")

    exit()


# Prepare the data

X = data.drop('species', axis=1)

y = data['species']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)


# Scale the data (important for KNN)

scaler = StandardScaler()

X_train = scaler.fit_transform(X_train)

X_test = scaler.transform(X_test)


# Find the optimal k value

```



```

error_rates = []

for k in range(1, 31): # Test k values from 1 to 30

    knn = KNeighborsClassifier(n_neighbors=k)

    knn.fit(X_train, y_train)

    y_pred = knn.predict(X_test)

    error_rates.append(1 - accuracy_score(y_test, y_pred)) # Error rate = 1 - accuracy


# Plot error rates

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

plt.plot(range(1, 31), error_rates, color='blue', linestyle='dashed', marker='o',
         markerfacecolor='red', markersize=10)

plt.title('Error Rate vs. K Value')

plt.xlabel('K')

plt.ylabel('Error Rate')

plt.show()

# Theory for choosing k:

# The optimal 'k' value minimizes the error rate.

# Very small k (e.g., 1) can lead to overfitting, being too sensitive to noise.

# Very large k (e.g., 30) can lead to underfitting, smoothing out the decision boundaries too much.

# We seek a k that balances these extremes, as shown by the error rate plot.


#Select k based on the minimum error rate observed in the plot

best_k = error_rates.index(min(error_rates)) + 1 #Add 1 as the index starts from 0

# Train and evaluate the model with the best k

```

```

knn = KNeighborsClassifier(n_neighbors=best_k)

knn.fit(X_train, y_train)

y_pred = knn.predict(X_test)


# Evaluate the model

print("Accuracy Score:", accuracy_score(y_test, y_pred))

print("\nConfusion Matrix:")

cm = confusion_matrix(y_test, y_pred)

print(cm)

print("\nClassification Report:")

print(classification_report(y_test, y_pred))


plt.figure(figsize=(8, 6))

sns.heatmap(cm, annot=True, fmt='d', cmap='Blues',

            xticklabels=knn.classes_, yticklabels=knn.classes_)

plt.xlabel('Predicted')

plt.ylabel('Actual')

plt.title('Confusion Matrix')

plt.show()


import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.neighbors import KNeighborsClassifier

from sklearn.metrics import accuracy_score

```

```

import matplotlib.pyplot as plt

# Load data

df = pd.read_csv('/content/iris (1).csv')

X = df.iloc[:, :-1]

y = df.iloc[:, -1]

# Train-test split

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=1)

# Store accuracy and error rate

accuracy = []

error_rate = []

# Try k from 1 to 20

for k in range(1, 21):

    knn = KNeighborsClassifier(n_neighbors=k)

    knn.fit(X_train, y_train)

    preds = knn.predict(X_test)

    acc = accuracy_score(y_test, preds)

    accuracy.append(acc)

    error_rate.append(1 - acc)

# Plot

```

```

plt.figure(figsize=(10,5))

plt.plot(range(1, 21), accuracy, label='Accuracy')

plt.plot(range(1, 21), error_rate, label='Error Rate')

plt.xlabel('K Value')

plt.ylabel('Rate')

plt.title('K vs Accuracy and Error Rate')

plt.legend()

plt.show()


import pandas as pd

from sklearn.preprocessing import StandardScaler


# Load data

df = pd.read_csv('/content/diabetes.csv')

X = df.drop('Outcome', axis=1) # Features

y = df['Outcome']             # Target


# Perform scaling

scaler = StandardScaler()

X_scaled = scaler.fit_transform(X)

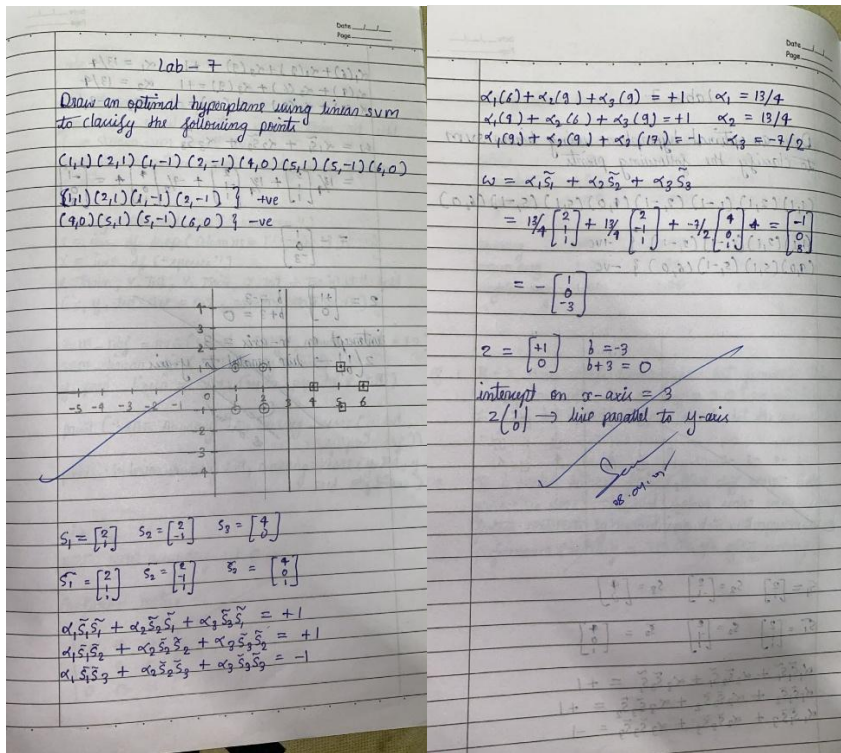

# Convert back to DataFrame (optional)

X_scaled_df = pd.DataFrame(X_scaled, columns=X.columns)

```

PROGRAM 7 Build Support vector machine model for a given dataset

Screenshot



Code SVM -

```
import pandas as pd
from sklearn.model_selection import train_test_split
from sklearn.svm import SVC
from sklearn.metrics import accuracy_score, confusion_matrix

iris_df = pd.read_csv("iris.csv")
X = iris_df.drop(columns=['species'])
y = iris_df['species']
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

svm = svm = SVC(kernel='rbf', random_state=42)
svm.fit(X_train, y_train)
y_pred_linear = svm.predict(X_test)

print("RBF Accuracy:", accuracy_score(y_test, y_pred_linear))
print("Linear Accuracy:", accuracy_score(y_test, y_pred_linear))
```

O/P -

RBF SVM accuracy: 1.0

Confusion matrix -

10	0	0
0	9	0
0	0	11

Linear SVM accuracy: 1.0

Confusion matrix

10	0	0
0	9	0
0	0	11

Questions -

Both RBF and linear gave the same accuracy for the iris dataset.

The AUC score was 95-98%. It shows that the AUC model performed pretty well for the letter-recognition.csv

The AUC score for iris.csv was 1.00 as it was very small and clean dataset with well-separated classes.

Code

```
import numpy as np

import matplotlib.pyplot as plt

positive_class = np.array([[4, 1], [4, -1], [6, 0]])

negative_class = np.array([[1, 0], [0, 1], [0, -1]])

plt.figure(figsize=(8, 6))

plt.scatter(positive_class[:, 0], positive_class[:, 1], color='red', label='Positive Class', s=100,
            edgecolors='black')

plt.scatter(negative_class[:, 0], negative_class[:, 1], color='blue', label='Negative Class', s=100,
            edgecolors='black')

all_points = np.concatenate([positive_class, negative_class])

labels = ["(4,1)", "(4,-1)", "(6,0)", "(1,0)", "(0,1)", "(0,-1)"]

for i, txt in enumerate(labels):

    plt.annotate(txt, (all_points[i][0], all_points[i][1]), textcoords="offset points", xytext=(0,5),
                ha='center', fontsize=10)

x_values = np.linspace(-1, 7, 100)

y_values = np.zeros_like(x_values)

plt.plot(x_values, y_values, color='black', linestyle='--', label='Optimal Hyperplane (y = 0)')
```

```

plt.plot(x_values, y_values + 1, color='gray', linestyle=':', label='Margin at y = 1')
plt.plot(x_values, y_values - 1, color='gray', linestyle=':', label='Margin at y = -1')

plt.title('Optimal Hyperplane for SVM (Visual Approximation)', fontsize=14)
plt.xlabel('x1')
plt.ylabel('x2')
plt.xlim(-1, 7)
plt.ylim(-2, 2)
plt.axhline(0, color='black',linewidth=0.5)
plt.axvline(0, color='black',linewidth=0.5)
plt.legend()

plt.grid(True)
plt.show()

import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.svm import SVC

from sklearn.metrics import accuracy_score, confusion_matrix

import seaborn as sns

import matplotlib.pyplot as plt

data = pd.read_csv('/content/iris (1) (1).csv')

X = data.drop('species', axis=1)

```

```

y = data['species']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

svm_rbf = SVC(kernel='rbf')

svm_rbf.fit(X_train, y_train)

y_pred_rbf = svm_rbf.predict(X_test)

accuracy_rbf = accuracy_score(y_test, y_pred_rbf)

cm_rbf = confusion_matrix(y_test, y_pred_rbf)

print("SVM with RBF Kernel:")

print("Accuracy:", accuracy_rbf)

print("Confusion Matrix:\n", cm_rbf)

plt.figure(figsize=(6, 4))

sns.heatmap(cm_rbf, annot=True, fmt='d', cmap='Blues',

            xticklabels=data['species'].unique(),

            yticklabels=data['species'].unique())

plt.xlabel('Predicted')

plt.ylabel('Actual')

plt.title('Confusion Matrix (RBF Kernel)')

plt.show()

svm_linear = SVC(kernel='linear')

svm_linear.fit(X_train, y_train)

```



```

y_pred_linear = svm_linear.predict(X_test)

accuracy_linear = accuracy_score(y_test, y_pred_linear)

cm_linear = confusion_matrix(y_test, y_pred_linear)


print("\nSVM with Linear Kernel:")

print("Accuracy:", accuracy_linear)

print("Confusion Matrix:\n", cm_linear)


plt.figure(figsize=(6, 4))

sns.heatmap(cm_linear, annot=True, fmt='d', cmap='Blues',

            xticklabels=data['species'].unique(),

            yticklabels=data['species'].unique())

plt.xlabel('Predicted')

plt.ylabel('Actual')

plt.title('Confusion Matrix (Linear Kernel)')

plt.show()

import numpy as np

import matplotlib.pyplot as plt

import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.svm import SVC

from sklearn.metrics import accuracy_score, confusion_matrix, roc_curve, auc

import seaborn as sns

from sklearn.preprocessing import label_binarize

```

```

from sklearn.multiclass import OneVsRestClassifier

data = pd.read_csv('/content/letter-recognition.csv') # Replace with the correct path if necessary

X = data.drop('letter', axis=1)
y = data['letter']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

svm_classifier = SVC(kernel='rbf', probability=True) # probability=True is needed for ROC curve
svm_classifier.fit(X_train, y_train)

y_pred = svm_classifier.predict(X_test)

accuracy = accuracy_score(y_test, y_pred)

cm = confusion_matrix(y_test, y_pred)

print("SVM Classifier:")
print("Accuracy:", accuracy)
print("Confusion Matrix:\n", cm)

plt.figure(figsize=(10, 8))

sns.heatmap(cm, annot=True, fmt="d", cmap="Blues", xticklabels=np.unique(y),
            yticklabels=np.unique(y))

plt.xlabel('Predicted')

```

```

plt.ylabel('Actual')

plt.title('Confusion Matrix')

plt.show()


y_test_bin = label_binarize(y_test, classes=np.unique(y))
n_classes = y_test_bin.shape[1]


classifier = OneVsRestClassifier(SVC(kernel='rbf', probability=True))
classifier.fit(X_train, y_train)
y_score = classifier.predict_proba(X_test)


fpr = dict()
tpr = dict()
roc_auc = dict()
for i in range(n_classes):
    fpr[i], tpr[i], _ = roc_curve(y_test_bin[:, i], y_score[:, i])
    roc_auc[i] = auc(fpr[i], tpr[i])


fpr["micro"], tpr["micro"], _ = roc_curve(y_test_bin.ravel(), y_score.ravel())
roc_auc["micro"] = auc(fpr["micro"], tpr["micro"])

plt.figure(figsize=(8, 6))
plt.plot(fpr["micro"], tpr["micro"],

        label='micro-average ROC curve (area = {0:0.2f})'

        ".format(roc_auc["micro"]))

```

```
plt.plot([0, 1], [0, 1], 'k--')  
plt.xlim([0.0, 1.0])  
plt.ylim([0.0, 1.05])  
plt.xlabel('False Positive Rate')  
plt.ylabel('True Positive Rate')  
plt.title('Micro-averaged ROC Curve')  
plt.legend(loc="lower right")  
plt.show()  
print(f"Micro-averaged AUC: {roc_auc['micro']}")
```

PROGRAM 8 Implement Random forest ensemble method on a given dataset.

Screenshot

lab = 8

→ Difference between Decision tree and Random Forest

DT	RF
1) Model type is single tree	Model type is ensemble of trees
2) Prone to overfitting	Less prone due to averaging
3) accuracy lower on complex data	accuracy higher due to ensemble learning
4) Faster to train and predict	Slower due to multiple trees
5) low bias, high variance	low bias, reduced variance

→ Parameters of RF

- n_estimators — No. of trees in the forest
- class_weight — weight associated with classes
- max_depth — max depth of each tree
- bootstrap — whether bootstrap samples are used when building trees
- n_jobs — no. of jobs to run in parallel
- random_state — seed for reproducibility
- verbose — controls the verbosity of output

→ Algorithm of Random Forest

- 1) Input dataset D with features & class labels
- 2) Set the no. of trees n_estimators
- 3) For each tree i from 1 to n_estimators

Code:

```
import pandas as pd
from sklearn.model_selection import train_test_split
from sklearn.ensemble import RandomForestClassifier
from sklearn.metrics import accuracy_score

iris = pd.read_csv('iris.csv')
X, y = iris[['sepal_length', 'sepal_width', 'petal_length'], iris['species']]
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

rf_default = RandomForestClassifier(n_estimators=10,
                                    random_state=42)
rf_default.fit(X_train, y_train)
y_pred_default = rf_default.predict(X_test)
accuracy_default = accuracy_score(y_test, y_pred_default)
print('Accuracy default: %f' % accuracy_default)

estimators = [10, 50, 100, 200, 500]
scores = []
for n in estimators:
    rf = RandomForestClassifier(n_estimators=n,
                               random_state=42)
    rf.fit(X_train, y_train)
    y_pred = rf.predict(X_test)
    score = accuracy_score(y_test, y_pred)
    scores.append(score)

print(scores)
print('Best n: %d' % estimators[scores.index(max(scores))])
```

o/p - accuracy with n_estimators = 10 : 1.00
 " " " " = 90 : 1
 " " " " = 100 : 1
 " " " " = 500 : 1

But n_estimators : 10
 best accuracy score : 1

Code

```
import pandas as pd

from sklearn.model_selection import train_test_split

from sklearn.ensemble import RandomForestClassifier

from sklearn.metrics import accuracy_score

import matplotlib.pyplot as plt


# Load the dataset

df = pd.read_csv('/content/iris (1).csv')


# Prepare features and target

X = df.drop(columns=['species']) # Assuming 'species' is the target column
y = df['species']


# Split into training and testing sets

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=42)


# Build Random Forest with default n_estimators (10)

rf_default = RandomForestClassifier(n_estimators=10, random_state=42)

rf_default.fit(X_train, y_train)

y_pred_default = rf_default.predict(X_test)


# Measure accuracy

default_score = accuracy_score(y_test, y_pred_default)
```

```

print(f"Default RF accuracy (n_estimators=10): {default_score:.4f}")

# Fine-tune the number of trees

scores = []

n_range = range(1, 101)

for n in n_range:

    rf = RandomForestClassifier(n_estimators=n, random_state=42)

    rf.fit(X_train, y_train)

    y_pred = rf.predict(X_test)

    score = accuracy_score(y_test, y_pred)

    scores.append(score)

# Find the best score and number of trees

best_score = max(scores)

best_n = n_range[scores.index(best_score)]

print(f"Best RF accuracy: {best_score:.4f} with n_estimators={best_n}")

# Optional: Plot accuracy vs number of estimators

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

plt.plot(n_range, scores, marker='o')

plt.title('Random Forest Accuracy vs Number of Trees')

plt.xlabel('Number of Trees (n_estimators)')

plt.ylabel('Accuracy')

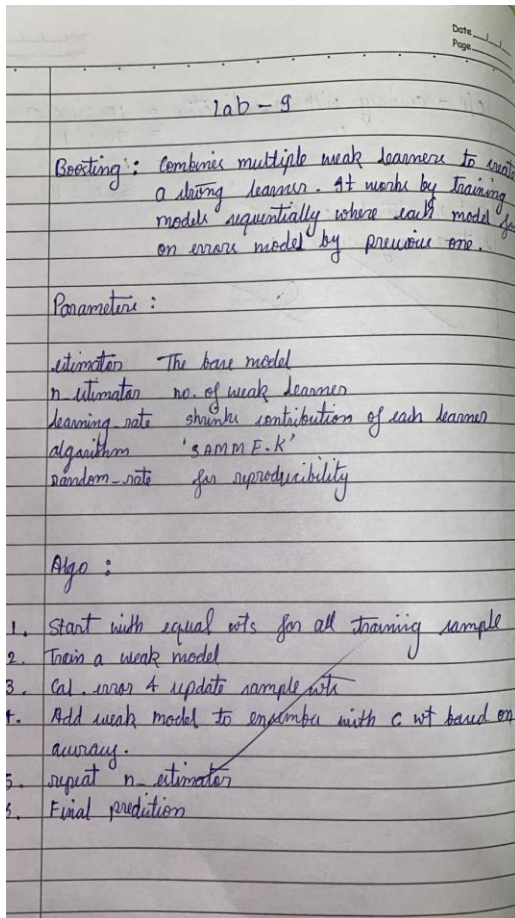
plt.grid(True)

plt.show()

```

PROGRAM 9 Implement Boosting ensemble method on a given dataset.

Screenshot



Code

```
import pandas as pd

import matplotlib.pyplot as plt

from sklearn.model_selection import train_test_split

from sklearn.preprocessing import LabelEncoder

from sklearn.ensemble import AdaBoostClassifier

from sklearn.metrics import accuracy_score

from sklearn.tree import DecisionTreeClassifier
```



```

# Load dataset

df = pd.read_csv("/content/income.csv")

# Drop rows with missing values

df.dropna(inplace=True)

# Encode categorical columns

label_encoders = {}

for column in df.select_dtypes(include=['object']).columns:

    le = LabelEncoder()

    df[column] = le.fit_transform(df[column])

    label_encoders[column] = le

# Separate features and target

X = df.drop(columns=['income_level'], errors='ignore', axis=1)

y = df['income_level']

# Split into training and testing sets

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

# AdaBoost with 10 estimators

model_10 = AdaBoostClassifier(n_estimators=10, random_state=42)

model_10.fit(X_train, y_train)

y_pred_10 = model_10.predict(X_test)

score_10 = accuracy_score(y_test, y_pred_10)

print(f"Accuracy with 10 estimators: {score_10:.4f}")

# Fine-tune number of estimators

best_score = 0

```

```

best_n = 0

estimators_range = list(range(10, 201, 10))

scores = []

for n in estimators_range:

    model = AdaBoostClassifier(n_estimators=n, random_state=42)

    model.fit(X_train, y_train)

    y_pred = model.predict(X_test)

    score = accuracy_score(y_test, y_pred)

    scores.append(score)

    print(f'n_estimators={n}, Accuracy={score:.4f}')

    if score > best_score:

        best_score = score

        best_n = n

print(f"\nBest Accuracy: {best_score:.4f} using {best_n} estimators")

# Plot accuracy vs number of estimators

plt.figure(figsize=(7, 4))

plt.plot(estimators_range, scores, marker='o', linestyle='-', color='blue')

plt.title("Accuracy vs Number of Estimators (AdaBoost)")

plt.xlabel("Number of Estimators (Trees)")

plt.ylabel("Accuracy")

plt.grid(True)

plt.xticks(estimators_range)

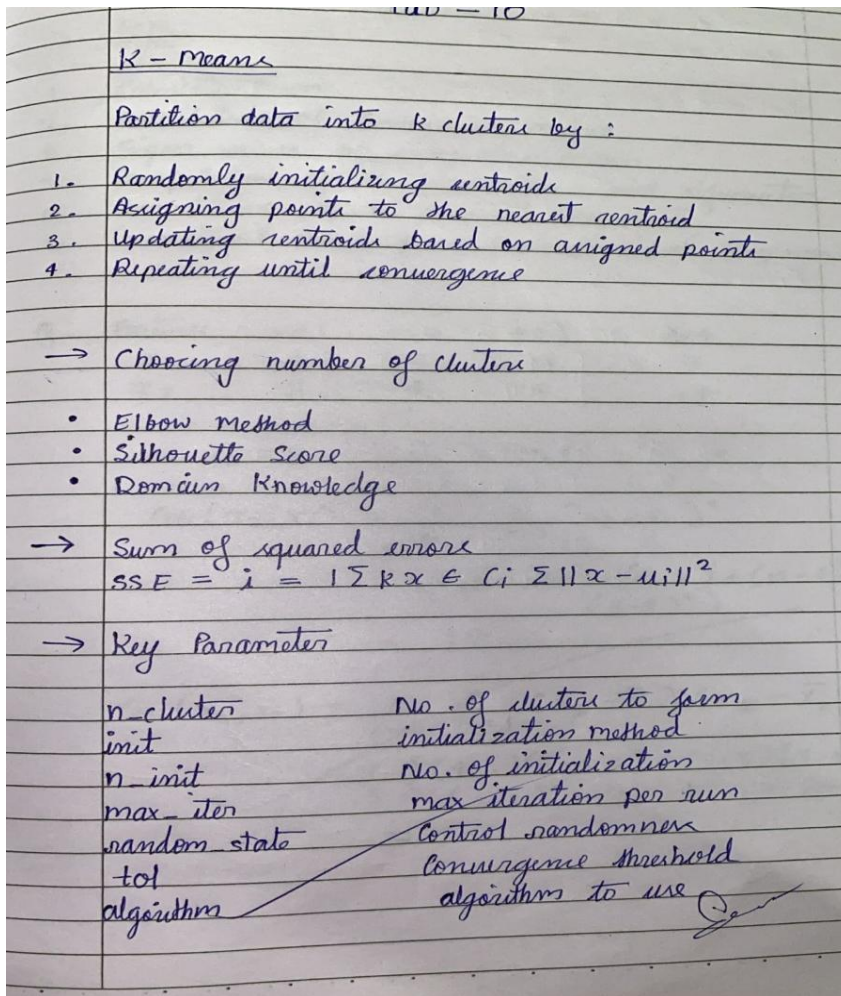
plt.tight_layout()

plt.show()

```

PROGRAM 10 Build k-Means algorithm to cluster a set of data stored in a .CSV file.

Screenshot



Code

```
import pandas as pd

import numpy as np

from sklearn.model_selection import train_test_split

from sklearn.preprocessing import StandardScaler

from sklearn.cluster import KMeans

from sklearn.metrics import accuracy_score
```

```

from scipy.stats import mode

import matplotlib.pyplot as plt


# Step 1: Generate sample data and save to CSV

np.random.seed(42)

names = [f"Person_{i}" for i in range(50)]

ages = np.random.randint(20, 60, 50)

income = np.random.randint(30000, 120000, 50)


df = pd.DataFrame({'Name': names, 'Age': ages, 'Income': income})

df.to_csv("income.csv", index=False)


# Step 2: Load the data

data = pd.read_csv("income.csv")


# Drop 'Name' and extract features

X = data[['Age', 'Income']]


# Step 3: Split the data

X_train, X_test = train_test_split(X, test_size=0.2, random_state=42)


# Step 4: Perform scaling

scaler = StandardScaler()

X_train_scaled = scaler.fit_transform(X_train)

```

```

X_test_scaled = scaler.transform(X_test)

# Step 5: Plot SSE vs number of clusters (Elbow method)

sse = []

k_range = range(1, 11)

for k in k_range:

    kmeans = KMeans(n_clusters=k, random_state=42)

    kmeans.fit(X_train_scaled)

    sse.append(kmeans.inertia_)


plt.figure(figsize=(8, 4))

plt.plot(k_range, sse, marker='o')

plt.xlabel('Number of clusters')

plt.ylabel('SSE (Inertia)')

plt.title('Elbow Method For Optimal k')

plt.grid(True)

plt.show()


# Step 6: Choose optimal number of clusters (say 3) and fit model

optimal_k = 3

kmeans = KMeans(n_clusters=optimal_k, random_state=42)

kmeans.fit(X_train_scaled)


# Predict on test data

```

```

predictions = kmeans.predict(X_test_scaled)

# Note: There's no ground truth labels, but for demonstration,
# we can try assigning true clusters (via KMeans on full data)
# and see if predicted clusters align

# Fit on full data to assign pseudo-labels
full_kmeans = KMeans(n_clusters=optimal_k, random_state=42)
true_clusters = full_kmeans.fit_predict(scaler.fit_transform(X))

# Align predicted clusters using majority voting (only for demonstration)
# Match predicted labels to closest true labels
def map_clusters(true_labels, pred_labels):
    labels = np.zeros_like(pred_labels)

    for i in range(optimal_k):
        mask = (pred_labels == i)

        if np.sum(mask) == 0:
            continue

        labels[mask] = mode(true_labels[mask])[0]

    return labels

mapped_preds = map_clusters(true_clusters[X_test.index], predictions)
accuracy = accuracy_score(true_clusters[X_test.index], mapped_preds)
print(f"Approximate Clustering Accuracy: {accuracy:.2f}")

```

```

import numpy as np

import pandas as pd

import matplotlib.pyplot as plt

from sklearn.datasets import load_iris

from sklearn.cluster import KMeans

from sklearn.preprocessing import StandardScaler

from sklearn.metrics import silhouette_score


# Step 1: Load Iris dataset

iris = load_iris()

df = pd.DataFrame(iris.data, columns=iris.feature_names)

df['target'] = iris.target


# Keep only petal length and petal width

X = df[['petal length (cm)', 'petal width (cm)']].values


# Step 2: Check impact of scaling

# Try without scaling

sse_unscaled = []

for k in range(1, 11):

    kmeans = KMeans(n_clusters=k, random_state=42)

    kmeans.fit(X)

    sse_unscaled.append(kmeans.inertia_)

```

```

# Now scale the features

scaler = StandardScaler()

X_scaled = scaler.fit_transform(X)


sse_scaled = []

for k in range(1, 11):

    kmeans = KMeans(n_clusters=k, random_state=42)

    kmeans.fit(X_scaled)

    sse_scaled.append(kmeans.inertia_)


# Step 3: Plot Elbow Comparison (Scaled vs Unscaled)

plt.figure(figsize=(10, 5))


plt.plot(range(1, 11), sse_unscaled, marker='o', label='Unscaled')
plt.plot(range(1, 11), sse_scaled, marker='s', label='Scaled')

plt.title('Elbow Method (Petal Features Only)')

plt.xlabel('Number of Clusters (k)')

plt.ylabel('SSE (Inertia)')

plt.legend()

plt.grid(True)

plt.show()

```


PROGRAM 11 Implement Dimensionality reduction using Principal Component Analysis (PCA) method.

Screenshot

lab - 11

PCA

1. Calculate Mean
2. Calculation of Covariance matrix
3. Eigen values of covariance matrix
4. Computation of the eigenvector - unit eigenvector
5. Computation of 1st principal component
6. Geometrical meaning of first principal component

Q. Features ex1 ex2 ex3 ex4

x1	4	8	13	7
x2	11	4	5	17

$\bar{x}_1 = 8$ $\bar{x}_2 = 8.5$

$$\text{Cov}(x_1, x_1) = \frac{1}{N-1} \sum_{i=1}^N (x_{1i} - \bar{x}_1)^2$$

$$= \frac{1}{3} ((4-8)^2 + (8-8)^2 + (13-8)^2 + (7-8)^2)$$

$$= 14$$

$$\text{Cov}(x_1, x_2) = \frac{1}{N-1} \sum_{i=1}^N (x_{1i} - \bar{x}_1)(x_{2i} - \bar{x}_2)$$

$$= \frac{1}{3} ((4-8)(11-8.5) + (8-8)(4-8.5) + (13-8)(5-8.5) + (7-8)(17-8.5))$$

$$= -11$$

$$\text{Cov}(x_2, x_2) = \frac{1}{N-1} \sum_{i=1}^N (x_{2i} - \bar{x}_2)^2$$

$$= \frac{1}{3} ((11-8.5)^2 + (4-8.5)^2 + (5-8.5)^2 + (17-8.5)^2)$$

$$= 23$$

Cov mat (S) =
$$\begin{bmatrix} \text{Cov}(x_1, x_1) & \text{Cov}(x_1, x_2) \\ \text{Cov}(x_2, x_1) & \text{Cov}(x_2, x_2) \end{bmatrix}$$

$$= \begin{bmatrix} 14 & -11 \\ -11 & 23 \end{bmatrix}$$

Eigen values of Covariance matrix

$$0 = \det(S - \lambda I)$$

$$= \begin{vmatrix} 14-\lambda & -11 \\ -11 & 23-\lambda \end{vmatrix}$$

$$\Rightarrow \lambda^2 - 37\lambda + 201$$

$$\lambda = \frac{1}{2} (37 \pm \sqrt{565})$$

$$= 30.3849, 6.6151$$

$$= \lambda_1, \lambda_2$$

Computation of eigenvectors

$$\lambda_1 = \begin{bmatrix} 11 \\ 4 \end{bmatrix}$$

$$\begin{bmatrix} 0 \\ 0 \end{bmatrix} = (S - \lambda_1 I) X$$

$$= \begin{bmatrix} 14-21 & -11 \\ -11 & 23-21 \end{bmatrix} X = \begin{bmatrix} -7 & -11 \\ -11 & 2 \end{bmatrix} X$$

$$\Rightarrow \frac{11}{11} = \frac{4}{(14-21)} = \lambda_1$$

$$\mu_1 = 11\hat{i}, \mu_2 = (14-21)\hat{j}$$

$$\mu_1 = \begin{bmatrix} 11 \\ 14-21 \end{bmatrix}$$

To find eigenvector, we compute length of x_1

$$\| \mu_1 \| = \sqrt{11^2 + (14-21)^2}$$

$$= 19.7348$$

$$\therefore \text{unit eigenvector corresponding to } \lambda_1, e_1 = \frac{1}{\| \mu_1 \|} \begin{bmatrix} 11 \\ (14-21) \end{bmatrix}$$

$$e_1 = \begin{bmatrix} 0.5574 \\ 0.5574 \end{bmatrix}$$

Computation of 1st principal component

$$\begin{bmatrix} x_{11} \\ x_{21} \end{bmatrix}$$

$$e_1^T \begin{bmatrix} x_{11} - \bar{x}_1 \\ x_{21} - \bar{x}_2 \end{bmatrix} = \begin{bmatrix} 0.5574 & 0.5574 \end{bmatrix} \begin{bmatrix} x_{11} - \bar{x}_1 \\ x_{21} - \bar{x}_2 \end{bmatrix}$$

$$= 0.5574(x_{11} - \bar{x}_1) + 0.5574(x_{21} - \bar{x}_2)$$

$$= 0.5574(x_{11} + x_{21} - \bar{x}_1 - \bar{x}_2)$$

Code

```
import pandas as pd

from sklearn.preprocessing import LabelEncoder, OneHotEncoder, StandardScaler

from sklearn.compose import ColumnTransformer

from sklearn.pipeline import Pipeline

from sklearn.model_selection import train_test_split

from sklearn.svm import SVC

from sklearn.linear_model import LogisticRegression

from sklearn.ensemble import RandomForestClassifier

from sklearn.decomposition import PCA

from sklearn.metrics import accuracy_score


# 1. Load data

df = pd.read_csv("heart.csv")


# 2. Label-encode binary text columns

le = LabelEncoder()

for col in ["Sex", "ExerciseAngina"]:

    df[col] = le.fit_transform(df[col])


# 3. Separate features and target

X = df.drop("HeartDisease", axis=1)

y = df["HeartDisease"]
```

4. Build preprocessing pipeline:

- One-hot for multi-category columns (using sparse_output=False)

- passthrough the rest

- then scale everything

cat_cols = ["ChestPainType", "RestingECG", "ST_Slope"]

```
preprocessor = Pipeline([
    ("onehot", ColumnTransformer([
        ("ohe", OneHotEncoder(sparse_output=False, drop="first"), cat_cols)
    ], remainder="passthrough")),
    ("scaler", StandardScaler())
])
```

5. Apply preprocessing

X_proc = preprocessor.fit_transform(X)

6. Train/test split

```
X_train, X_test, y_train, y_test = train_test_split(
    X_proc, y, test_size=0.2, random_state=42
)
```

7. Define models

```
models = {
    "SVM": SVC(random_state=42),
    "LogisticRegression": LogisticRegression(max_iter=1000, random_state=42),
```

```

    "RandomForest": RandomForestClassifier(random_state=42)
}

# 8. Train & evaluate before PCA

print("=== Accuracies BEFORE PCA ===")

scores_before = {}

for name, clf in models.items():

    clf.fit(X_train, y_train)

    preds = clf.predict(X_test)

    acc = accuracy_score(y_test, preds)

    scores_before[name] = acc

    print(f"{name:17s}: {acc:.4f}")


# 9. Apply PCA (retain 95% variance)

pca = PCA(n_components=0.95, random_state=42)

X_train_pca = pca.fit_transform(X_train)

X_test_pca = pca.transform(X_test)

print(f"\nPCA retained {pca.n_components_} components, "

      f"explained variance = {pca.explained_variance_ratio_.sum():.4f}\n")

# 10. Train & evaluate after PCA

print("=== Accuracies AFTER PCA ===")

scores_after = {}

for name, clf in models.items():

    clf.fit(X_train_pca, y_train)

```

```
preds = clf.predict(X_test_pca)

acc = accuracy_score(y_test, preds)

scores_after[name] = acc

print(f" {name:17s}: {acc:.4f}")
```