PART A

1 #FIND S PROG:

```
import pandas as pd
import numpy as np
       #to read the data in the csv file
data = pd.read csv("data.csv")
print(data,"n")
       #making an array of all the attributes
d = np.array(data)[:,:-1]
print("n The attributes are: ",d)
       #segragating the target that has positive and negative examples
target = np.array(data)[:,-1]
print("n The target is: ",target)
       #training function to implement find-s algorithm
def train(c,t):
  for i, val in enumerate(t):
     if val == "Yes":
       specific hypothesis = c[i].copy()
       break
  for i, val in enumerate(c):
     if t[i] == "Yes":
       for x in range(len(specific hypothesis)):
          if val[x] != specific hypothesis[x]:
             specific hypothesis[x] = '?'
          else:
             pass
  return specific hypothesis
       #obtaining the final hypothesis
print("n The final hypothesis is:",train(d,target))
```

2 #CANDIDATE ELIMINATION PROG:

```
import csv
with open("data.csv") as f:
  csv file=csv.reader(f)
  data=list(csv file)
  s=data[1][:-1] #1st record is taken as specific hypothesis
  g=[['?' for i in range(len(s))] for j in range(len(s))] #? is put for all the records
  for i in data:
     if i[-1]=="Yes": # if target attribute is yes, then it should match specific hypothesis
       for j in range(len(s)):
          if i[j]!=s[j]: # if record doesn't match specific hypothesis, put ? in specific columns
             s[i]='?'
             g[j][j]='?'
     elif i[-1]=="No": # if target attribute is no, then specific hypothesis should not match
       for j in range(len(s)):
          if i[j]!=s[j]:
             g[j][j]=s[j]
          else:
             g[j][j]="?"
     print("\nSteps of Candidate Elimination Algorithm",data.index(i)+1)
     print(s)
    print(g)
  gh=[]
  for i in g:
     for j in i:
       if i!='?':
          gh.append(i)
          break
  print("\nFinal specific hypothesis:\n",s)
  print("\nFinal general hypothesis:\n",gh)
```

3 #ID3 / DECISION TREE

```
from sklearn.tree import DecisionTreeClassifier
import pandas as pd
from sklearn.metrics import confusion matrix
dataset=pd.read csv('pima-indians-Diabetes.csv')
print(dataset.head())
train features=dataset.iloc[:80,:-1]
test features=dataset.iloc[80:,:-1]
train targets=dataset.iloc[:80,-1]
test targets=dataset.iloc[80:,-1]
tree1=DecisionTreeClassifier(criterion='entropy').fit(train features,train targets)
prediction=tree1.predict(test features)
cm=confusion matrix(test targets,prediction)
print('Confusion Matrix:\n',cm)
TP = cm[0][0]
FP=cm[0][1]
FN=cm[1][0]
TN=cm[1][1]
print('False Positive \n {}'.format(FP))
print('False Negative \n {}'.format(FN))
print('True Positive \n {}'.format(TP))
print('True Negative \n {}'.format(TN))
TPR=TP/(TP+FN)
print('Senstivity \n {}'.format(TPR))
TNR=TN/(TN+FP)
print('Specificity \n {}'.format(TNR))
Precision=TP/(TP+FP)
print('Precision \n {}'.format(Precision))
```

```
Recall=TP/(TP+FN)

print('Recall \n {}'.format(Recall))

Acc=(TP+TN)/(TP+TN+FP+FN)

print('Accuracy \n {}'.format(Acc))

Fscore=2*(Precision*Recall)/(Precision+Recall)

print('FScore \n {}'.format(Fscore))
```

4.A #BINARIZATION

```
# Python code explaining how to Binarize feature values
#Importing Libraries
import pandas as pd
       #Importing Data
data set = pd.read csv('data.csv')
print(data set.head())
age = data_set.iloc[:, 1].values
salary = data set.iloc[:, 2].values
print ("\nOriginal age data values : \n", age)
print ("\nOriginal salary data values : \n", salary)
       # Binarizing values
from sklearn.preprocessing import Binarizer
x = age
x = x.reshape(1, -1)
y = salary
y = y.reshape(1, -1)
       # For age, let threshold be 35, For salary, let threshold be 61000
binarizer 1 = Binarizer(35) \# Age below 35 is binarized to 0
binarizer_2 = Binarizer(61000) # Salary below 61000 is binarized to 0
       # Transformed features
print ("\nBinarized age : \n", binarizer 1.fit transform(x))
print ("\nBinarized salary : \n", binarizer 2.fit transform(y))
```

4.B #BINNING

```
import pandas as pd
import numpy as np
       # Generate 20 random integers uniformly between 0 and 99
small counts = np.random.randint(0, 100, 20)
print(small counts)
       # Map to evenly spaced bins 0-9 by division
print(np.floor divide(small counts, 10))
large counts = [296, 8286, 64011, 80, 3, 725, 867, 2215, 7689, 11495, 91897, 44, 28, 7971,
926, 12]
       # print(np.floor(np.log10(large counts)))
       #Map the counts to quartiles into 4 bins (quartile)
print(pd.qcut(large counts, 4, labels=False))
       #convert large counts into series data
large counts series = pd.Series(large counts)
       # print(large counts series)
print(large counts series.quantile([0.25, 0.5, 0.75]))
4.C #LOG TRANSFORMATION
import pandas as pd
import numpy as np
       #Log Transform Example
data = pd.DataFrame({'value':[2,45, -23, 85, 28, 2, 35, -12]})
data['log+1'] = (data['value']+1).transform(np.log)
print(data)
       #Negative Values Handling. Note that the values are different
data['log'] = (data['value']-data['value'].min()+1) .transform(np.log)
print(data)
```

4.D #FEATURE SCALING

```
#Importing Libraries
import pandas as pd
       # Sklearn library
from sklearn import preprocessing
       #Import Data
data set = pd.read csv('data.csv')
print(data_set.head())
       # here Features - Age and Salary columns
       # are taken using slicing
       # to handle values with varying magnitude
x = data set.iloc[:, 1:3].values
print ("\nOriginal data values : \n", x)
       #MIN MAX SCALER
min max scaler = preprocessing.MinMaxScaler(feature range =(0, 1))
       # Scaled feature
x after min max scaler = min max scaler.fit transform(x)
print ("\nAfter min max Scaling : \n", x after min max scaler)
       # STANDARDIZATION
Standardisation = preprocessing.StandardScaler()
# Scaled feature
x after Standardisation = Standardisation.fit transform(x)
print ("\nAfter Standardisation : \n", x after Standardisation)
```

PART B

1 #PCA

```
# https://towardsdatascience.com/pca-using-python-scikit-learn-e653f8989e60
import pandas as pd
url = "https://archive.ics.uci.edu/ml/machine-learning-databases/iris/iris.data"
       # load dataset into Pandas DataFrame
       #data = pd.read csv("data5.csv")
       #df = pd.read csv("Iris.csv", names=['sepal length', 'sepal width', 'petal length', 'petal
       width', 'target'])
df = pd.read csv(url, names=['sepal length', 'sepal width', 'petal length', 'petal width', 'target'])
print(df.head)
from sklearn.preprocessing import StandardScaler
features = ['sepal length', 'sepal width', 'petal length', 'petal width']
       # Separating out the features
x = df.loc[:, features].values
       # Separating out the target
y = df.loc[:,['target']].values
       # Standardizing the features
x = StandardScaler().fit transform(x)
print("Standardized features ", x)
from sklearn.decomposition import PCA
pca = PCA(n components=2)
principalComponents = pca.fit transform(x) #Apply PCA
principalDf = pd.DataFrame(data = principalComponents,
columns = ['principal component 1', 'principal component 2'])
finalDf = pd.concat([principalDf, df[['target']]], axis = 1)
print(pca.explained variance ratio )
```

2 #K MEANS

```
#Import libraries
import pandas as pd
from sklearn.cluster import KMeans
       #import the dataset
df = pd.read csv('iris.csv')
print(df.head(10))
x = df.iloc[:, [0,1,2,3]].values # Load Input Attributes
kmeans5 = KMeans(n_clusters=5) # Create 5 Clusters
y kmeans5 = kmeans5.fit predict(x)
print(y kmeans5) #Prints the clusters for each record
print(kmeans5.cluster centers )
Error =[]
for i in range(1, 11):
  kmeans = KMeans(n clusters = i).fit(x)
  kmeans.fit(x)
  Error.append(kmeans.inertia)
import matplotlib.pyplot as plt
plt.plot(range(1, 11), Error)
plt.title('Elbow method')
plt.xlabel('No of clusters')
plt.ylabel('Error')
plt.show()
```

3 #NAIVE BAYES

```
from sklearn import datasets
iris = datasets.load iris()
print("Features: ", iris.feature names)
print("Labels: ", iris.target_names)
from sklearn.model_selection import train_test_split
x_train, x_test, y_train, y_test = train_test_split(iris.data, iris.target, test_size=0.3,
random state=109)
from sklearn.naive_bayes import GaussianNB
gnb = GaussianNB()
gnb.fit(x_train, y_train)
y_pred = gnb.predict(x_test)
from sklearn import metrics
print("Accuracy: " ,metrics.accuracy score(y test, y pred))
```

4 #ADA BOOST CLAASIFIER

```
#https://www.datacamp.com/community/tutorials/adaboost-classifier-python
       # Load libraries
from sklearn.ensemble import AdaBoostClassifier
from sklearn import datasets
# Import train test split function
from sklearn.model selection import train test split
#Import scikit-learn metrics module for accuracy calculation
from sklearn import metrics
iris = datasets.load iris()
X = iris.data
y = iris.target
       # Split dataset into training set and test set
X train, X test, y train, y test = train test split(X, y, test size=0.3)
       # Create adaboost classifer object
abc = AdaBoostClassifier(n estimators=50,
               learning rate=1)
       # Train Adaboost Classifer
model = abc.fit(X_train, y_train)
       #Predict the response for test dataset
y pred = model.predict(X test)
print("Accuracy:",metrics.accuracy score(y test, y pred))
       #Using Different Base Learners
       # Import Support Vector Classifier
from sklearn.svm import SVC
svc=SVC(probability=True, kernel='linear')
```

```
# Create adaboost classifer object
abc =AdaBoostClassifier(n_estimators=50, base_estimator=svc,learning_rate=1)
    # Train Adaboost Classifer
model = abc.fit(X_train, y_train)
    #Predict the response for test dataset
y_pred = model.predict(X_test)
    # Model Accuracy, how often is the classifier correct?
print("Accuracy:",metrics.accuracy_score(y_test, y_pred))
```

5 #RANDOM FOREST

```
import pandas as pd
import numpy as np
dataset = pd.read csv('petrol consumption.csv')
print(dataset.head())
X = dataset.iloc[:, 0:4].values
y = dataset.iloc[:, 4].values
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, random state=0)
       # Feature Scaling
from sklearn.preprocessing import StandardScaler
sc = StandardScaler()
X train = sc.fit transform(X train)
X \text{ test} = \text{sc.transform}(X \text{ test})
from sklearn.ensemble import RandomForestRegressor
regressor = RandomForestRegressor(n estimators=20, random state=0)
regressor.fit(X train, y train)
y pred = regressor.predict(X test)
from sklearn import metrics
print('Mean Absolute Error:', metrics.mean absolute error(y test, y pred))
print('Mean Squared Error:', metrics.mean squared error(y test, y pred))
print('Root Mean Squared Error:', np.sqrt(metrics.mean squared error(y test, y pred)))
```