1.Implement A* Search algorithm.

```
from collections import deque
class Graph:
    def __init__(self, adjac_lis):
        \overline{\text{self.adjac}} ac_lis = adjac_lis
    def get neighbors(self, v):
        return self.adjac lis[v]
    # This is heuristic function which is having equal values for all nodes
    def h(self, n):
        H = {
            'A': 1,
            'B': 1,
            'C': 1,
            'D': 1
        return H[n]
    def a star algorithm(self, start, stop):
        # In this open 1st is a lisy of nodes which have been visited, but
        # neighbours haven't all been always inspected, It starts off with
        the start
  #node
        # And closed 1st is a list of nodes which have been visited
        # and who's neighbors have been always inspected
        open lst = set([start])
        closed_lst = set([])
        # poo has present distances from start to all other nodes
        # the default value is +infinity
        poo = {}
        poo[start] = 0
        # par contains an adjac mapping of all nodes
        par = {}
        par[start] = start
        while len(open_lst) > 0:
            n = None
            # it will find a node with the lowest value of f() -
            for v in open lst:
                if n == None or poo[v] + self.h(v) < poo[n] + self.h(n):
                    #print(poo[v] + self.h(v))
                     #print(poo[n] + self.h(n))
                    n = v;
                    #print(v,n)
            if n == None:
                print('Path does not exist!')
                return None
            # if the current node is the stop
            # then we start again from start
            if n == stop:
                reconst_path = []
```

```
while par[n] != n:
                    reconst path.append(n)
                    n = par[n]
                reconst path.append(start)
                reconst path.reverse()
                print('Path found: {}'.format(reconst path))
                return reconst path
            # for all the neighbors of the current node do
            for (m, weight) in self.get neighbors(n):
                #print(m)
              # if the current node is not presentin both open 1st and
        closed 1st
                # add it to open 1st and note n as it's par
                if m not in open 1st and m not in closed 1st:
                    open lst.add(m)
                    par[m] = n
                    poo[m] = poo[n] + weight
                # otherwise, check if it's quicker to first visit n, then m
                # and if it is, update par data and poo data
                # and if the node was in the closed lst, move it to
        open 1st
                else:
                    if poo[m] > poo[n] + weight:
                        poo[m] = poo[n] + weight
                        par[m] = n
                        if m in closed_lst:
                            closed lst.remove(m)
                            open lst.add(m)
            # remove n from the open lst, and add it to closed lst
            # because all of his neighbors were inspected
            open lst.remove(n)
            closed lst.add(n)
        print('Path does not exist!')
        return None
adjac lis ={
    'A': [('B', 1), ('C', 3), ('D', 7)],
    'B': [('D', 5)],
    'C': [('D', 12)]
}
graph1 =Graph(adjac lis)
graph1.a star algorithm('A', 'D')
Output:
Path found: ['A', 'B', 'D']
['A', 'B', 'D']
```

2. Implement AO* Search algorithm.

```
def recAOStar(n):
  global finalPath
  print("Expanding Node : ", n)
  and nodes = []
  or nodes = []
#Segregation of AND and OR nodes
  if (n in allNodes):
    if 'AND' in allNodes[n]:
       and nodes = allNodes[n]['AND']
    if 'OR' in allNodes[n]:
       or nodes = allNodes[n]['OR']
# If leaf node then return
  if len(and nodes) == 0 and len(or nodes) == 0:
    return
  solvable = False
  marked = \{\}
  while not solvable:
# If all the child nodes are visited and expanded, take the least cost of all the child nodes
    if len(marked) == len(and nodes) + len(or nodes):
       min cost least, min cost group least = least cost group(and nodes, or nodes, {})
       solvable = True
       change heuristic(n, min cost least)
       optimal child group[n] = min cost group least
       continue
# Least cost of the unmarked child nodes
  min cost, min cost group = least cost group(and nodes, or nodes, marked)
  is expanded = False
  # If the child nodes have sub trees then recursively visit them to recalculate the heuristic of
the child node
  if len(min cost group) > 1:
    if (min cost group[0] in allNodes):
       is expanded = True
       recAOStar(min cost group[0])
```

```
if (min cost group[1] in allNodes):
       is expanded = True
       recAOStar(min cost group[1])
    else:
       if (min cost group in allNodes):
         is expanded = True
         recAOStar(min cost group)
         # If the child node had any subtree and expanded, verify if the new heuristic value
is still the least among all nodes
    if is expanded:
       min cost verify, min cost group verify = least cost group(and nodes, or nodes,
{})
  if min cost group == min cost group verify:
    solvable = True
    change heuristic(n, min cost verify)
    optimal child group[n] = min cost group
    # If the child node does not have any subtrees then no change in heuristic, so update the
min cost of the current node
  else:
    solvable = True
    change heuristic(n, min cost)
    optimal child group[n] = min cost group
    #Mark the child node which was expanded
  marked[min cost group] = 1
  return heuristic(n)
# Function to calculate the min cost among all the child nodes
def least cost group(and nodes, or nodes, marked):
  node wise cost = \{\}
  for node pair in and nodes:
    if not node pair[0] + node pair[1] in marked:
       cost = 0
       cost = cost + heuristic(node pair[0]) + heuristic(node pair[1]) + 2
       node wise cost[node pair[0] + node pair[1]] = cost
  for node in or nodes:
```

```
if not node in marked:
       cost = 0
       cost = cost + heuristic(node) + 1
       node wise cost[node] = cost
  min cost = 999999
  min cost group = None
  # Calculates the min heuristic
  for costKey in node wise cost:
    if node wise cost[costKey] < min cost:
       min_cost = node_wise_cost[costKey]
       min cost group = costKey
  return [min cost, min cost group]
  # Returns heuristic of a node
def heuristic(n):
  return H_dist[n]
  # Updates the heuristic of a node
def change heuristic(n, cost):
  H dist[n] = cost
  return
  # Function to print the optimal cost nodes
def print_path(node):
  print(optimal child group[node], end="")
  node = optimal_child_group[node]
  if len(node) > 1:
    if node[0] in optimal child group:
       print("->", end="")
       print path(node[0])
    if node[1] in optimal child group:
       print("->", end="")
       print path(node[1])
  else:
    if node in optimal child group:
       print("->", end="")
       print path(node)
```

Output:

Expanding Node : A Expanding Node : B Expanding Node : C Expanding Node : D

Nodes which gives optimal cost are

CD->HI->J

Optimal Cost is :: 5

3.For a given set of training data examples stored in a .CSV file, implement and demonstrate the Candidate-Elimination algorithm to output a description of the set of all hypotheses consistent with the training examples.

```
import numpy as np
import pandas as pd
data = pd.read csv('data4.csv')
concepts = np.array(data.iloc[:.0:-1])
target = np.array(data.iloc[:,-1])
def learn(concepts,target):
  count = 0
  first = ['?','?','?','?','?']
  for i, val in enumerate(target):
     if val == 'Yes':
       #print(specific h)
       break
  specific h = concepts[i].copy()
  generic_h = [["?" for i in range(len(specific_h))] for i in range (len(specific_h))]
  for i,h in enumerate(concepts):
     if target[i] == "Yes":
       for x in range(len(specific h)):
          if h[x] != specific h[x]:
             specific h[x] = "?"
             generic h[x][x] = "?"
     if target[i] == "No":
       for x in range(len(specific h)):
          if h[x] != specific h[x]:
             generic h[x][x] = \text{specific } h[x]
          else:
             generic h[x][x] == "?"
     if generic h[x][x] == "?":
       print(f'S{count} : {specific h}')
       print(f'G{count} : {first}')
       count+=1
     else:
       print(f'S{count} : {specific h}')
       print(f'G{count} : {generic h}')
       for x in range(len(generic h)):
          first[x] = generic h[x]
       count += 1
  indices=[i for i, val in enumerate(generic h)if val == ['?','?','?','?','?','?']]
  for i in indices:
     generic h.remove(['?','?','?','?','?'])
  return specific h, generic h
s final,g final = learn(concepts,target)
print("final s:",s final,sep="\n")
print("final g:",g final,sep="\n")
```

dataset:data4.csv

sunny,warm,normal,strong,warm,same,yes sunny,warm,high,strong,warm,same,yes rainy,cold,high,strong,warm,change,no sunny,warm,high,strong,cool,change,yes

Output

```
So: ['sunny' 'warm' 'high' 'strong' 'warm' 'same']
Go: ['?', '?', '?', '?', '?', '?']
S1: ['sunny' 'warm' 'high' 'strong' 'warm' 'same']
G1: [['sunny', '?', '?', '?', '?', '?'], ['?', 'warm', '?', '?', '?'], ['?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?'], ['?', '?', '?', '?']]
final s:
['sunny' 'warm' 'high' 'strong' '?' '?']
final g:
[['sunny', '?', '?', '?', '?', '?'], ['?', 'warm', '?', '?', '?', '?']]
```

4. Write a program to demonstrate the working of the decision tree based ID3 algorithm. Use an appropriate data set for building the decision tree and apply this knowledge toclassify a new sample.

```
import pandas as pd
import math
import numpy as np
data = pd.read csv("play.csv")
features = [feat for feat in data]
features.remove("classification")
class Node:
  def init__(self):
     self.children = []
     self.value = ""
     self.isLeaf = False
     self.pred = ""
def entropy(examples):
  pos = 0.0
  neg = 0.0
  for , row in examples.iterrows():
     if row["classification"] == "Yes":
       pos += 1
     else:
       neg += 1
  if pos == 0.0 or neg == 0.0:
     return 0.0
  else:
     p = pos / (pos + neg)
     n = neg / (pos + neg)
     return -(p * math.log(p, 2) + n * math.log(n, 2))
def info gain(examples, attr):
  uniq = np.unique(examples[attr])
  #print ("\n",uniq)
  gain = entropy(examples)
  #print ("\n",gain)
  for u in uniq:
     subdata = examples[examples[attr] == u]
     #print ("\n",subdata)
     sub e = entropy(subdata)
     gain -= (float(len(subdata)) / float(len(examples))) * sub e
     #print ("\n",gain)
  return gain
def ID3(examples, attrs):
  root = Node()
  \max gain = 0
```

```
max feat = ""
  for feature in attrs:
    #print ("\n",examples)
    gain = info gain(examples, feature)
    if gain > max gain:
       max gain = gain
       max feat = feature
  root.value = max feat
  #print ("\nMax feature attr",max feat)
  uniq = np.unique(examples[max feat])
  #print ("\n",uniq)
  for u in uniq:
    #print ("\n",u)
    subdata = examples[examples[max feat] == u]
    #print ("\n",subdata)
    if entropy(subdata) == 0.0:
       newNode = Node()
       newNode.isLeaf = True
       newNode.value = u
       newNode.pred = np.unique(subdata["classification"])
       root.children.append(newNode)
    else:
       dummyNode = Node()
       dummyNode.value = u
       new attrs = attrs.copy()
       new attrs.remove(max feat)
       child = ID3(subdata, new attrs)
       dummyNode.children.append(child)
       root.children.append(dummyNode)
  return root
def printTree(root: Node, depth=0):
  for i in range(depth):
    print("\t", end="")
  print(root.value, end="")
  if root.isLeaf:
    print(" -> ", root.pred)
  print()
  for child in root.children:
    printTree(child, depth + 1)
root = ID3(data, features)
printTree(root)
```

dataset:

play.csv

A1,A2,A3,classification True,Hot,High,No True,Hot,High,No False,Hot,High,Yes False,Cool,Normal,Yes False,Cool,Normal,Yes True,Cool,High,No True,Hot,High,No True,Hot,Normal,Yes False,Cool,Normal,Yes False,Cool,High,No

output

```
A3

High

A1

False

A2

Cool -> ['No']

Hot -> ['Yes']

Normal -> ['Yes']
```

5. Build an Artificial Neural Network by implementing the Backpropagation Algorithm and test the same using appropriate data sets.

```
import numpy as np
X = \text{np.array}(([2, 9], [1, 5], [3, 6]), \text{dtype=float})
y = np.array(([92], [86], [89]), dtype=float)
X = X/np.amax(X,axis=0) \# maximum of X array longitudinally <math>y = y/100
#Sigmoid Function
def sigmoid (x):
  return (1/(1 + np.exp(-x)))
#Derivative of Sigmoid Function
def derivatives sigmoid(x):
  return x * (1 - x)
                                            #Variable initialization
epoch=7000
                                    #Setting training iterations
1r=0.1
                                    #Setting learning rate
inputlayer neurons = 2
                                    #number of features in data set
hiddenlayer neurons = 3
                                     #number of hidden layers neurons
output neurons = 1
                                     #number of neurons at output layer
                                            #weight and bias initialization
wh=np.random.uniform(size=(inputlayer neurons,hiddenlayer neurons))
bh=np.random.uniform(size=(1,hiddenlayer neurons))
wout=np.random.uniform(size=(hiddenlayer neurons,output neurons))
bout=np.random.uniform(size=(1,output neurons))
       # draws a random range of numbers uniformly of dim x*y
#Forward Propagation
for i in range(epoch):
  hinp1=np.dot(X,wh)
  hinp=hinp1 + bh
  hlayer act = sigmoid(hinp)
  outinp1=np.dot(hlayer act,wout)
  outinp= outinp1+ bout
  output = sigmoid(outinp)
#Backpropagation
  EO = y-output
  outgrad = derivatives sigmoid(output)
  d output = EO* outgrad
  EH = d output.dot(wout.T)
  hiddengrad = derivatives sigmoid(hlayer act)
```

[[0.89559591] [0.88142069] [0.8928407]]

output

```
#how much hidden layer wts contributed to error
  d hiddenlayer = EH * hiddengrad
  wout += hlayer act.T.dot(d output) *lr
# dotproduct of nextlayererror and currentlayerop
  bout += np.sum(d output, axis=0,keepdims=True) *lr
  wh += X.T.dot(d_hiddenlayer) *lr
#bh += np.sum(d hiddenlayer, axis=0,keepdims=True) *lr
print("Input: \n" + str(X))
print("Actual Output: n'' + str(y))
print("Predicted Output: \n" ,output)
Input:
[[ 0.66666667 1.
[ 0.33333333 0.55555556]
[ 1.
         0.66666667]]
Actual Output:
[[ 0.92]
[0.86]
[0.89]]
Predicted Output:
```

6. Write a program to implement the naïve Bayesian classifier for a sample training data set stored as a .CSV file. Compute the accuracy of the classifier, considering few test data sets.

```
import csv
import random
import math
def loadCsv(filename):
 lines = csv.reader(open(filename, "r"));
 dataset = list(lines)
 for i in range(len(dataset)):
                                       #converting strings into numbers for processing
       dataset[i] = [float(x) for x in dataset[i]]
 return dataset
def splitDataset(dataset, splitRatio):
                                                              #67% training size
 trainSize = int(len(dataset) * splitRatio);
 trainSet = []
 copy = list(dataset);
 while len(trainSet) < trainSize:
          #generate indices for the dataset list randomly to pick ele for
                                                                training
       data index = random.randrange(len(copy));
       trainSet.append(copy.pop(index))
 return [trainSet, copy]
def separateByClass(dataset):
 separated = \{\}
#creates a dictionary of classes 1 and 0 where the values are the instances belonging
to # each class
 for i in range(len(dataset)):
       vector = dataset[i]
       if (vector[-1] not in separated):
               separated[vector[-1]] = []
       separated[vector[-1]].append(vector)
 return separated
def mean(numbers):
 return sum(numbers)/float(len(numbers))
def stdev(numbers):
 avg = mean(numbers)
 variance = sum([pow(x-avg,2) for x in numbers])/float(len(numbers)-1)
 return math.sqrt(variance)
def summarize(dataset):
 summaries = [(mean(attribute), stdev(attribute)) for attribute in zip(*dataset)];
```

```
del summaries[-1]
 return summaries
def summarizeByClass(dataset):
 separated = separateByClass(dataset);
 summaries = \{\}
                          #summaries is a dic of tuples(mean,std) for each class value
 for classValue, instances in separated.items():
       summaries[classValue] = summarize(instances)
 return summaries
def calculateProbability(x, mean, stdev):
 exponent = math.exp(-(math.pow(x-mean,2)/(2*math.pow(stdev,2))))
 return (1 / (math.sqrt(2*math.pi) * stdev)) * exponent
def calculateClassProbabilities(summaries, inputVector):
probabilities = {}
                                     #class and attribute information as mean and sd
 for classValue, classSummaries in summaries.items():
       probabilities[classValue] = 1
       for i in range(len(classSummaries)):
              mean, stdev = classSummaries[i]
                                                #take mean and sd of every attribute
       for class 0 and 1 seperaely
              x = inputVector[i]
                                                              #testvector's first
              attribute
              probabilities[classValue] *= calculateProbability(x, mean, stdev);
                                                                      #use normal dist
   return probabilities
def predict(summaries, inputVector):
 probabilities = calculateClassProbabilities(summaries, inputVector)
 bestLabel, bestProb = None, -1
 for classValue, probability in probabilities.items():
                                        #assigns that class which has he highest prob
       if bestLabel is None or probability > bestProb:
              bestProb = probability
              bestLabel = classValue
 return bestLabel
def getPredictions(summaries, testSet):
predictions = []
 for i in range(len(testSet)):
       result = predict(summaries, testSet[i])
       predictions.append(result)
return predictions
```

```
def getAccuracy(testSet, predictions):
 correct = 0
 for i in range(len(testSet)):
       if testSet[i][-1] == predictions[i]:
               correct += 1
 return (correct/float(len(testSet))) * 100.0
def main():
 filename = 'diabetesdata.csv'
 splitRatio = 0.67
 dataset = loadCsv(filename);
 trainingSet, testSet = splitDataset(dataset, splitRatio)
 print('Split {0} rows into train={1} and test={2} rows'.format(len(dataset),
 len(trainingSet), len(testSet)))
                                                                    # prepare model
 summaries = summarizeByClass(trainingSet);
                                                                    # test model
 predictions = getPredictions(summaries, testSet)
 accuracy = getAccuracy(testSet, predictions)
 print('Accuracy of the classifier is: {0}%'.format(accuracy))
main()
Output
       confusion matrix is as follows
       [[170 0]
        [0170]
        [0\ 0\ 11]]
       Accuracy metrics
               precision
                           recall f1-score support
           0
                   1.00
                           1.00
                                   1.00
                                             17
            1
                   1.00
                           1.00
                                   1.00
                                             17
           2
                                             11
                  1.00
                           1.00
                                   1.00
       avg / total 1.00
                              1.00 1.00
                                                45
```

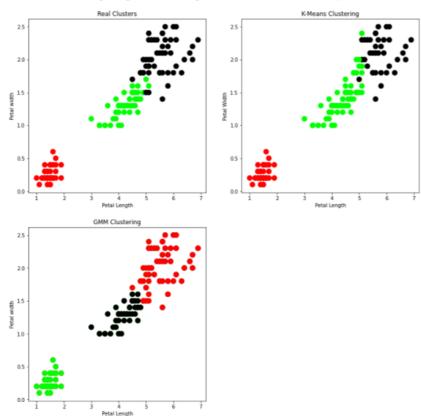
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7. Apply EM algorithm to cluster a set of data stored in a .CSV file. Use the same data set for clustering using k-Means algorithm. Compare the results of these two algorithms and comment on the quality of clustering. You can add Java/Python ML library classes/API in the program.

```
import matplotlib.pyplot as plt
from sklearn import datasets
from sklearn.cluster import KMeans
import pandas as pd
import numpy as np
iris=datasets.load iris()
X=pd.DataFrame(iris.data)
X.columns=['Sepal Length', 'Sepal Width', 'Petal Length', 'Petal Width']
y=pd.DataFrame(iris.target)
y.columns=['Targets']
model=KMeans(n clusters=3)
model.fit(X)
plt.figure(figsize=(14,14))
colormap=np.array(['red','lime','black'])
plt.subplot(2,2,1)
plt.scatter(X.Petal Length, X.Petal Width, c=colormap[y.Targets], s=40)
plt.title('Real Clusters')
plt.xlabel('Petal Length')
plt.ylabel('Petal width')
plt.subplot(2,2,2)
plt.scatter(X.Petal Length,X.Petal Width,c=colormap[model.labels_],s=40)
plt.title('K-Means Clustering')
plt.xlabel('Petal Length')
plt.ylabel('Petal Width')
plt.subplot(2,2,2)
plt.scatter(X.Petal Length, X.Petal Width, c=colormap[model.labels], s=40)
plt.title('K-Means Clustering')
plt.ylabel('Petal Width')
from sklearn import preprocessing
scaler=preprocessing.StandardScaler()
scaler.fit(X)
xsa=scaler.transform(X)
xs=pd.DataFrame(xsa,columns=X.columns)
from sklearn.mixture import GaussianMixture
gmm=GaussianMixture(n components=3)
gmm.fit(xs)
gmm y=gmm.predict(xs)
plt.subplot(2,2,3)
plt.scatter(X.Petal Length, X.Petal Width, c=colormap[gmm y], s=40)
plt.title('GMM Clustering')
plt.xlabel('Petal Length')
plt.ylabel('Petal width')
print('Observation:The GMM using EM algo based clustering matched the true labels more
closely than KMeans.')
```

<u>output</u>

Observation:The GMM using EM algo based clustering matched the true labels more closely than KMeans.



8. Write a program to implement k-Nearest Neighbour algorithm to classify the iris data set. Print both correct and wrong predictions. Java/Python ML library classes can be used for this problem.

```
from sklearn.datasets import load iris
from sklearn.neighbors import KNeighborsClassifier
import numpy as np
from sklearn.model selection import train test split
iris dataset = load iris()
#print(iris dataset)
targets = iris dataset.target names
print("Class : number")
for i in range(len(targets)):
  print(targets[i], ':', i)
X_train, X_test, y_train, y_test = train_test_split(iris_dataset["data"], iris_dataset["target"])
kn = KNeighborsClassifier(1)
kn.fit(X train, y train)
for i in range(len(X test)):
  x_new = np.array([X_test[i]])
  prediction = kn.predict(x new)
  print("Actual:[{0}] [{1}],Predicted:{2} {3}".format(y_test[i], targets[y_test[i]], prediction,
targets[prediction]))
print("\nAccuracy:",kn.score(X test,y test))
```

Output:

```
Class : number
setosa : 0
versicolor : 1
virginica : 2
Actual:[0] [setosa],Predicted:[0] ['setosa']
Actual:[1] [versicolor],Predicted:[1] ['versicolor']
Actual:[1] [versicolor],Predicted:[1] ['versicolor']
Actual:[2] [virginica],Predicted:[2] ['virginica']
Actual:[2] [virginica],Predicted:[2] ['virginica']
Actual:[1] [versicolor],Predicted:[2] ['virginica']
Actual:[1] [versicolor],Predicted:[2] ['virginica']
Actual:[1] [versicolor],Predicted:[2] ['virginica']
Actual:[2] [virginica],Predicted:[2] ['virginica']
Actual:[1] [versicolor],Predicted:[1] ['versicolor']
Actual:[1] [versicolor],Predicted:[1] ['versicolor']
Actual:[2] [virginica],Predicted:[2] ['virginica']
Actual:[0] [setosa],Predicted:[2] ['virginica']
Actual:[0] [setosa],Predicted:[0] ['setosa']
Actual:[0] [setosa],Predicted:[0] ['setosa']
Actual:[0] [setosa],Predicted:[0] ['virginica']
Actual:[0] [virginica],Predicted:[0] ['virginica']
Actual:[0] [virginica],Predicted:[0] ['virginica']
Actual:[0] [virginica],Predicted:[0] ['virginica']
Actual:[0] [virginica],Predicted:[0] ['virginica']
Actual
```

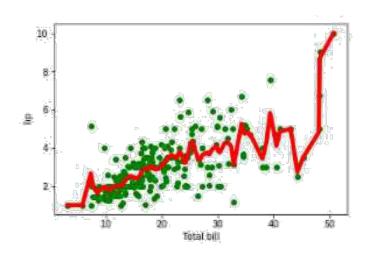
9.Implement the non-parametric Locally Weighted Regressionalgorithm in order to fit data points. Select appropriate data set for your experiment and draw graphs

```
from numpy import *
import operator
from os import listdir
import matplotlib
import matplotlib.pyplot as plt
import pandas as pd
import numpy as np1
import numpy.linalg as np
from scipy.stats.stats import pearsonr
def kernel(point,xmat, k):
  m,n = np1.shape(xmat)
  weights = np1.mat(np1.eye((m)))
  for j in range(m):
    diff = point - X[i]
    weights[j,j] = np1.exp(diff*diff.T/(-2.0*k**2))
  return weights
def localWeight(point,xmat,ymat,k):
  wei = kernel(point,xmat,k)
  W=(X.T*(wei*X)).I*(X.T*(wei*ymat.T))
  return W
def localWeightRegression(xmat,ymat,k):
  m,n = np1.shape(xmat)
  ypred = np1.zeros(m)
  for i in range(m):
    ypred[i] = xmat[i]*localWeight(xmat[i],xmat,ymat,k)
  return ypred
                                                     # load data points
data = pd.read csv('data10.csv')
bill = np1.array(data.total bill)
tip = np1.array(data.tip)
                                            #preparing and add 1 in bill
mbill = np1.mat(bill)
mtip = np1.mat(tip)
m = np1.shape(mbill)[1]
```

one = np1.mat(np1.ones(m))
X= np1.hstack((one.T,mbill.T))
ypred = localWeightRegression(X,mtip,2)
SortIndex = X[:,1].argsort(0)
xsort = X[SortIndex][:,0]

#set k here

Output



Questions

- 1. What is machine learning?
- 2. Define supervised learning
- 3. Define unsupervised learning
- 4. Define semi supervised learning
- 5. Define reinforcement learning
- 6. What do you mean by hypotheses?
- 7. What is classification?
- 8. What is clustering?
- 9. Define precision, accuracy and recall
- 10. Define entropy
- 11. Define regression
- 12. How Knn is different from k-means clustering?
- 13. What is concept learning
- 14. Define specific boundary and general boundary
- 15. Define target function
- 16. Define decision tree
- 17. What is ANN
- 18. Explain gradient descent approximation
- 19. State Bayes theorem
- 20. Define Bayesian belief network
- 21. Differentiate hard and soft clustering
- 22. Define variance
- 23. What is inductive machine learning?
- 24. Why K nearest neighbour algorithm is lazy learning algorithm?
- 25. Why naïve Bayes is naïve?
- 26. Mention classification algorithms
- 27. Define pruning
- 28. Differentiate Clustering and classification
- 29. Mention clustering algorithms
- 30. Define Bias