

VISVESVARAYA TECHNOLOGICAL UNIVERSITY

JNANA SANGAMA, BELAGAVI-590018



3-Day FDP on Machine Learning July 19-21, 2018

Jointly Organized by

DEPARTMENTS OF CSE & ISE

RNS INSTITUTE OF TECHNOLOGY

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Laboratory Manual of Machine Learning [15CSL76]

For VII Sem CSE/ISE

MACHINE LEARNING LABORATORY

[As per Choice Based Credit System (CBCS) scheme]

SEMESTER – VII

Subject Code 15CSL76

Course objectives:

This course will enable students to

1. Make use of Data sets in implementing the machine learning algorithms
2. Implement the machine learning concepts and algorithms in any suitable language of choice

Description (If any):

1. The programs can be implemented in either JAVA or Python.
2. For Problems 1 to 6 and 10, programs are to be developed without using the built-in classes or APIs of Java/Python.
3. Data sets can be taken from standard repositories (<https://archive.ics.uci.edu/ml/datasets.html>) or constructed by the students.

Lab Experiments:

1. Implement and demonstrate the FIND-S algorithm for finding the most specific hypothesis based on a given set of training data samples. Read the training data from a .CSV file.
2. 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.
3. 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 to classify a new sample.
4. Build an Artificial Neural Network by implementing the Back propagation algorithm and test the same using appropriate data sets.
5. 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.
6. Assuming a set of documents that need to be classified, use the naïve Bayesian Classifier model to perform this task. Built-in Java classes/API can be used to write the program. Calculate the accuracy, precision, and recall for your data set.

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7. Write a program to construct a Bayesian network considering medical data. Use this model to demonstrate the diagnosis of heart patients using standard Heart Disease Data Set. You can use Java/Python ML library classes/API.
8. 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.
9. 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.
10. Implement the non-parametric Locally Weighted Regression algorithm in order to fit data points. Select appropriate data set for your experiment and draw graphs.

Course outcomes:

The students should be able to:

1. Understand the implementation procedures for the machine learning algorithms.
2. Design Java/Python programs for various Learning algorithms.
3. Apply appropriate data sets to the Machine Learning algorithms.
4. Identify and apply Machine Learning algorithms to solve real world problems.

Conduction of Practical Examination:

- All laboratory experiments are to be included for practical examination.
- Students are allowed to pick one experiment from the lot.
- Strictly follow the instructions as printed on the cover page of answer script
- Marks distribution: Procedure + Conduction + Viva:20 + 50 +10 (80)

Program 1

Implement and demonstrate the FIND-S algorithm for finding the most specific hypothesis based on a given set of training data samples. Read the training data from a .CSV file.

Task: Find-S algorithm is used to find a maximally specific hypothesis.

Positive and negative training examples for the target concept EnjoySport.

Example	Sky	AirTemp	Humidity	Wind	Water	Forecast	EnjoySport
1	Sunny	Warm	Normal	Strong	Warm	Same	Yes
2	Sunny	Warm	High	Strong	Warm	Same	Yes
3	Rainy	Cold	High	Strong	Warm	Change	No
4	Sunny	Warm	High	Strong	Cool	Change	Yes

Given:

- **Instances X:** Possible days, each described by the attributes
 - ❖ **Sky** (with possible values Sunny, Cloudy, and Rainy),
 - ❖ **AirTemp** (with values Warm and Cold),
 - ❖ **Humidity** (with values Normal and High),
 - ❖ **Wind** (with values Strong and Weak),
 - ❖ **Water** (with values Warm and Cool), and
 - ❖ **Forecast** (with values Same and Change).
- **Hypotheses H:** Each hypothesis is described by a conjunction of constraints on the attributes Sky, AirTemp, Humidity, Wind, Water, and Forecast. The constraints may be "?" (any value is acceptable), "0" (no value is acceptable), or a specific value.
- **Target concept c:** EnjoySport : $X \rightarrow (0,1)$
- **Training examples D:** Positive and negative examples of the target function

The steps of FINDS are given below:

Step1: The first step of FIND-S is to initialize h to the most specific hypothesis in H

$$h_0 \leftarrow \{ \phi, \phi, \phi, \phi, \phi, \phi \}$$

Ex.	Sky	Temp	Humid	Wind	Water	Forecast	Enjoy Sport?	Hypothesis
0								$\langle \phi, \phi, \phi, \phi, \phi, \phi \rangle$

Step 2: Positive training example $x_1 = \langle \text{Sunny, Warm, Normal, Strong, Warm, Same} \rangle$

$$h_1 \leftarrow \langle \text{Sunny, Warm, Normal, Strong, Warm, Same} \rangle$$

Ex.	Sky	Temp	Humid	Wind	Water	Forecast	Enjoy Sport?	Hypothesis
0								$\langle \emptyset, \emptyset, \emptyset, \emptyset, \emptyset \rangle$
1	Sunny	Warm	Normal	Strong	Warm	Same	Yes	$\langle \text{Sunny, Warm, Normal, Strong, Warm, Same} \rangle$

Step 3: Positive training instance $x_2 = \langle \text{Sunny, Warm, High, Strong, Warm, Same} \rangle$

$h_2 \leftarrow \langle \text{Sunny, Warm, ?, Strong, Warm, Same} \rangle$

Ex.	Sky	Temp	Humid	Wind	Water	Forecast	Enjoy Sport?	Hypothesis
0								$\langle \emptyset, \emptyset, \emptyset, \emptyset, \emptyset \rangle$
1	Sunny	Warm	Normal	Strong	Warm	Same	Yes	$\langle \text{Sunny, Warm, Normal, Strong, Warm, Same} \rangle$
2	Sunny	Warm	High	String	Warm	Same	Yes	$\langle \text{Sunny, Warm, ?, Strong, Warm, Same} \rangle$

Step 4: Negative training example $x_3 = \langle \text{Rainy, Cold, High, Strong, Warm, Change} \rangle$

$h_3 \leftarrow \langle \text{Sunny, Warm, ?, Strong, Warm, Same} \rangle$

Ex.	Sky	Temp	Humid	Wind	Water	Forecast	Enjoy Sport?	Hypothesis
0								$\langle \emptyset, \emptyset, \emptyset, \emptyset, \emptyset \rangle$
1	Sunny	Warm	Normal	Strong	Warm	Same	Yes	$\langle \text{Sunny, Warm, Normal, Strong, Warm, Same} \rangle$
2	Sunny	Warm	High	String	Warm	Same	Yes	$\langle \text{Sunny, Warm, ?, Strong, Warm, Same} \rangle$
3	Rainy	Cold	High	Strong	Warm	Change	No	$\langle \text{Sunny, Warm, ?, Strong, Warm, Same} \rangle$

Step 5: Positive training example $x_4 = \langle \text{Sunny, Warm, High, Strong, Cool, change} \rangle$

$h_4 \leftarrow \langle \text{Sunny, Warm, ?, Strong, ?, ?} \rangle$

Ex.	Sky	Temp	Humid	Wind	Water	Forecast	Enjoy Sport?	Hypothesis
0								$\langle \emptyset, \emptyset, \emptyset, \emptyset, \emptyset, \emptyset \rangle$
1	Sunny	Warm	Normal	Strong	Warm	Same	Yes	$\langle \text{Sunny, Warm, Normal, Strong, Warm, Same} \rangle$
2	Sunny	Warm	High	String	Warm	Same	Yes	$\langle \text{Sunny, Warm, ?, Strong, Warm, Same} \rangle$
3	Rainy	Cold	High	Strong	Warm	Change	No	$\langle \text{Sunny, Warm, ?, Strong, Warm, Same} \rangle$
4	Sunny	Warm	High	Strong	Cool	Change	Yes	$\langle \text{Sunny, Warm, ?, Strong, ?, ?} \rangle$

Find-s Algorithm:

Step 1: Initialize h to the most specific hypothesis in H.

Step 2: For each positive training instance x.

- For each attribute constraint a_i in h
 - If the constraint a_i is satisfied by x
Then do nothing
 - Else replace a_i in h by the next more general constraint that is satisfied by x.

Step 3: Output hypothesis h.

Find-s in Python:

```
import csv
#!/usr/bin/python
#list creatin
hypo=['%','%','%','%','%','%'];

with open('Training_examples.csv') as csv_file:
    readcsv = csv.reader(csv_file, delimiter=',')
    print(readcsv)
    data = []
    print("\n\nThe given training examples are:")
    for row in readcsv:
        print(row)
        if row[len(row)-1].upper() == "YES":
            data.append(row)

print("\n\nThe positive examples are:");
for x in data:
    print(x);
print("\n\n");
```

```
TotalExamples = len(data);
i=0;
j=0;
k=0;
print("The steps of the Find-s algorithm are\n",hypo);
list = [];
p=0;
d=len(data[p])-1;
for j in range(d):
    list.append(data[i][j]);
hypo=list;
i=1;
for i in range(TotalExamples):
    for k in range(d):
        if hypo[k]!=data[i][k]:
            hypo[k]='?';
            k=k+1;
        else:
            hypo[k];
    print(hypo);
i=i+1;
print("\n\nThe maximally specific Find-s hypothesis for the given training examples is");
list=[];
for i in range(d):
    list.append(hypo[i]);
print(list);
```

Input csv file:

Training_examples.csv

Output:

The given training examples are:

```
['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']
```

The positive examples are:

```
['Sunny', 'Warm', 'Normal', 'Strong', 'Warm', 'Same', 'Yes']
['Sunny', 'Warm', 'High', 'Strong', 'Warm', 'Same', 'Yes']
['Sunny', 'Warm', 'High', 'Strong', 'Cool', 'Change', 'Yes']
```

The steps of the Find-s algorithm are

['%', '%', '%', '%', '%', '%']

['Sunny', 'Warm', 'Normal', 'Strong', 'Warm', 'Same']

['Sunny', 'Warm', '?', 'Strong', 'Warm', 'Same']

['Sunny', 'Warm', '?', 'Strong', '?', '?']

The maximally specific Find-s hypothesis for the given training examples is

['Sunny', 'Warm', '?', 'Strong', '?', '?']

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Program 2

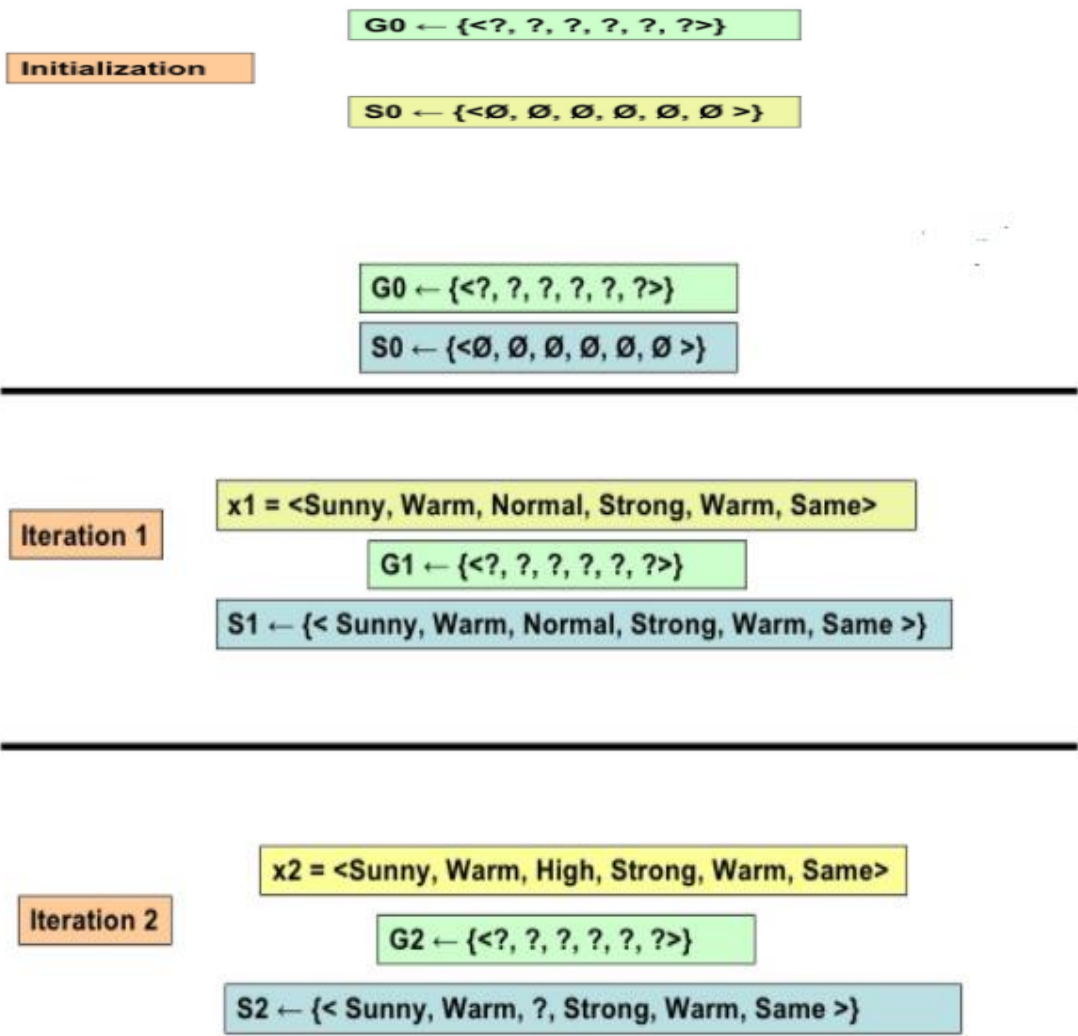
For a given set of training data examples stored in a .CSV file, implement and demonstrate the Candidate-Elimination algorithm to output a description of the set of all hypotheses consistent with the training examples.

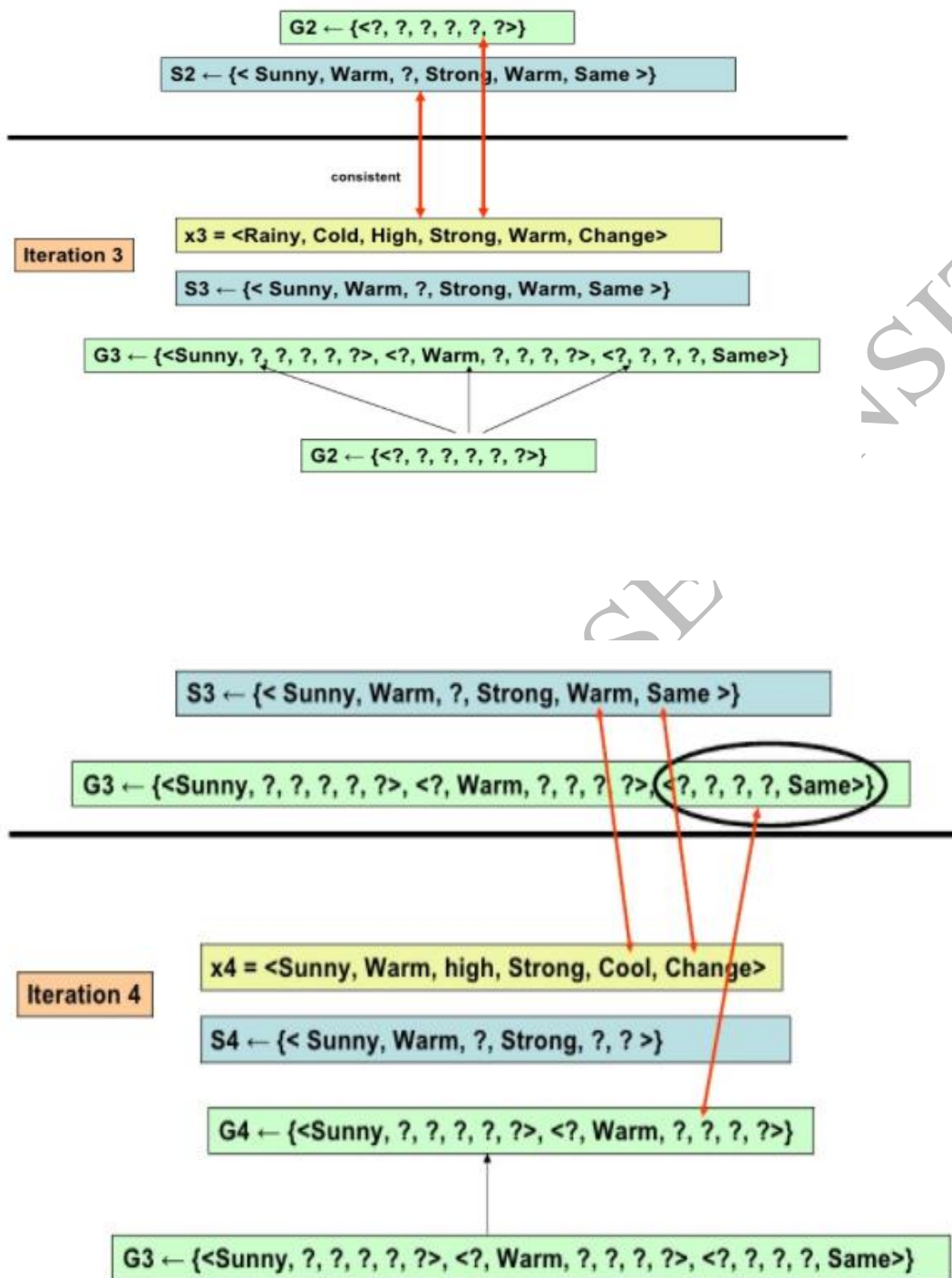
Task: The CANDIDATE-ELIMINATION algorithm computes the version space containing all hypotheses from H that are consistent with an observed sequence of training examples.

Dataset: EnjoySport training examples

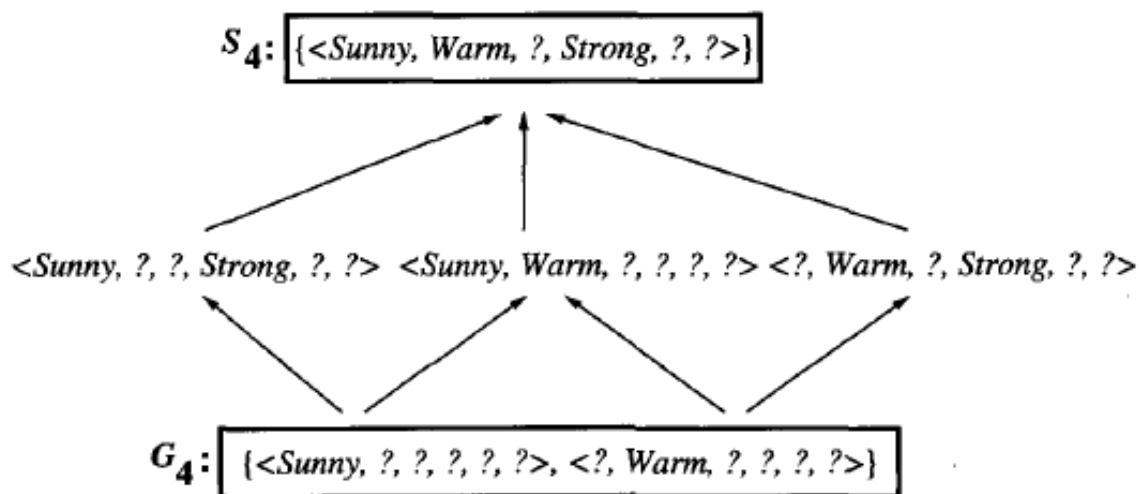
Example

Example	Sky	AirTemp	Humidity	Wind	Water	Forecast	EnjoySport
1	Sunny	Warm	Normal	Strong	Warm	Same	Yes
2	Sunny	Warm	High	Strong	Warm	Same	Yes
3	Rainy	Cold	High	Strong	Warm	Change	No
4	Sunny	Warm	High	Strong	Cool	Change	Yes





Final version space



Candidate-Elimination Algorithm:

Initialize G to the set of maximally general hypotheses in H

Initialize S to the set of maximally specific hypotheses in H

For each training example d , do

- If d is a positive example
 - Remove from G any hypothesis inconsistent with d
 - For each hypothesis s in S that is not consistent with d
 - Remove s from S
 - Add to S all minimal generalizations h of s such that
 - h is consistent with d , and some member of G is more general than h
 - Remove from S any hypothesis that is more general than another hypothesis in S
- If d is a negative example
 - Remove from S any hypothesis inconsistent with d
 - For each hypothesis g in G that is not consistent with d
 - Remove g from G
 - Add to G all minimal specializations h of g such that
 - h is consistent with d , and some member of S is more specific than h
 - Remove from G any hypothesis that is less general than another hypothesis in G

Candidate-Elimiation in Python:

```
class Holder:

    factors={} #Initialize an empty dictionary

    attributes = () #declaration of dictionaries parameters with an arbitrary length

    """
    Constructor of class Holder holding two parameters,
    self refers to the instance of the class
    """

    def __init__(self,attr): #
        self.attributes = attr

    for i in attr:
        self.factors[i]=[]

    def add_values(self,factor,values):
        self.factors[factor]=values

class CandidateElimination:

    Positive={} #Initialize positive empty dictionary

    Negative={} #Initialize negative empty dictionary

    def __init__(self,data,fact):
        self.num_factors = len(data[0][0])
        self.factors = fact.factors

        self.attr = fact.attributes

        self.dataset = data

        #print self.attr
```

```
def run_algorithm(self):
    # print self.dataset
    """
    Initialize the specific and general boundaries, and loop the dataset against the algorithm
    """
    G = self.initializeG()
    S = self.initializeS()
    print("initialize General",G)
    print("initialize Specific",S)

    """
    Programmatically populate list in the iterating variable trial_set
    """
    #count=0
    for trial_set in self.dataset:
        if self.is_positive(trial_set): #if trial set/example consists of positive examples
            G = self.remove_inconsistent_G(G,trial_set[0]) #remove inconsitent data from the general
            boundary
            S_new = S[:] #initialize the dictionary with no key-value pair
            print(S_new)
            for s in S:
                if not self.consistent(s,trial_set[0]):
                    S_new.remove(s)
            generalization = self.generalize_inconsistent_S(s,trial_set[0])
            generalization = self.get_general(generalization,G)
            if generalization:
                S_new.append(generalization)
```

```
S = S_new[:]  
  
S = self.remove_more_general(S)  
  
print(S)  
  
else:#if it is negative  
  
    S = self.remove_inconsistent_S(S,trial_set[0]) #remove inconsitent data from the specific  
boundary  
  
    G_new = G[:] #initialize the dictionary with no key-value pair (dataset can take any value)  
  
print(G_new)  
  
for g in G:  
  
if self.consistent(g,trial_set[0]):  
  
    G_new.remove(g)  
  
specializations = self.specialize_inconsistent_G(g,trial_set[0])  
specializationss = self.get_specific(specializations,S)  
  
if specializations != []:  
  
    G_new += specializations  
  
    G = G_new[:]  
  
print(G)  
  
    G = self.remove_more_specific(G)  
  
print(S)  
  
print(G)  
  
def initializeS(self):  
  
    """ Initialize the specific boundary """  
  
    S = tuple(['-' for factor in range(self.num_factors)]) #6 constraints in the vector  
  
return [S]  
  
def initializeG(self):  
  
    """ Initialize the general boundary """
```

```
G = tuple(['?' for factor in range(self.num_factors)]) # 6 constraints in the vector
return [G]

def is_positive(self,trial_set):
    """ Check if a given training trial_set is positive """
    if trial_set[1] == 'Y':
        return True
    elif trial_set[1] == 'N':
        return False
    else:
        raise TypeError("invalid target value")

def is_negative(self,trial_set):
    """ Check if a given training trial_set is negative """
    if trial_set[1] == 'N':
        return False
    elif trial_set[1] == 'Y':
        return True
    else:
        raise TypeError("invalid target value")

def match_factor(self,value1,value2):
    """ Check for the factors values match,
    necessary while checking the consistency of
    training trial_set with the hypothesis """
    if value1 == '?' or value2 == '?':
        return True
```

```
elif value1 == value2 :

return True

return False

def consistent(self,hypothesis,instance):

    """ Check whether the instance is part of the hypothesis """

for i,factor in enumerate(hypothesis):

if not self.match_factor(factor,instance[i]):

return False

return True

def remove_inconsistent_G(self,hypotheses,instance):

    """ For a positive trial_set, the hypotheses in G

inconsistent with it should be removed """

    G_new = hypotheses[:]

for g in hypotheses:

if not self.consistent(g,instance):

    G_new.remove(g)

return G_new

def remove_inconsistent_S(self,hypotheses,instance):

    """ For a negative trial_set, the hypotheses in S

inconsistent with it should be removed """

    S_new = hypotheses[:]

for s in hypotheses:

if self.consistent(s,instance):

    S_new.remove(s)
```



```
return S_new
```

```
def remove_more_general(self,hypotheses):
```

```
    """ After generalizing S for a positive trial_set, the hypothesis in S
    general than others in S should be removed """
```

```
        S_new = hypotheses[:]
```

```
    for old in hypotheses:
```

```
        for new in S_new:
```

```
            if old!=new and self.more_general(new,old):
```

```
                S_new.remove[new]
```

```
    return S_new
```

```
def remove_more_specific(self,hypotheses):
```

```
    """ After specializing G for a negative trial_set, the hypothesis in G
    specific than others in G should be removed """
```

```
        G_new = hypotheses[:]
```

```
    for old in hypotheses:
```

```
        for new in G_new:
```

```
            if old!=new and self.more_specific(new,old):
```

```
                G_new.remove[new]
```

```
    return G_new
```

```
def generalize_inconsistent_S(self,hypothesis,instance):
```

```
    """ When a inconsistent hypothesis for positive trial_set is seen in the specific boundary S,
    it should be generalized to be consistent with the trial_set ... we will get one hypothesis"""
```

```
    hypo = list(hypothesis) # convert tuple to list for mutability
```

```
for i,factor in enumerate(hypo):
    if factor == '-':
        hypo[i] = instance[i]
    elif not self.match_factor(factor,instance[i]):
        hypo[i] = '?'
    generalization = tuple(hypo) # convert list back to tuple for immutability
    return generalization

def specialize_inconsistent_G(self,hypothesis,instance):
    """ When a inconsistent hypothesis for negative trial_set is seen in the general boundary G
    should be specialized to be consistent with the trial_set.. we will get a set of hypotheses """
    specializations = []
    hypo = list(hypothesis) # convert tuple to list for mutability
    for i,factor in enumerate(hypo):
        if factor == '?':
            values = self.factors[self.attr[i]]
            for j in values:
                if instance[i] != j:
                    hyp=hypo[:]
                    hyp[i]=j
                    hyp=tuple(hyp) # convert list back to tuple for immutability
                    specializations.append(hyp)
    return specializations

def get_general(self,generalization,G):
    """ Checks if there is more general hypothesis in G
```

for a generalization of inconsistent hypothesis in S

in case of positive trial_set and returns valid generalization ""

for g in G:

if self.more_general(g,generalization):

return generalization

return None

def get_specific(self,specializations,S):

"" Checks if there is more specific hypothesis in S

for each of hypothesis in specializations of an

inconsistent hypothesis in G in case of negative trial_set

and return the valid specializations""

valid_specializations = []

for hypo in specializations:

for s in S:

if self.more_specific(s,hypo) or s==self.initializeS()[0]:

valid_specializations.append(hypo)

return valid_specializations

def exists_general(self,hypothesis,G):

""Used to check if there exists a more general hypothesis in
general boundary for version space""

for g in G:

if self.more_general(g,hypothesis):

```
return True
```

```
return False
```

```
def exists_specific(self,hypothesis,S):
```

```
    """Used to check if there exists a more specific hypothesis in  
    general boundary for version space"""
```

```
    for s in S:
```

```
        if self.more_specific(s,hypothesis):
```

```
            return True
```

```
    return False
```

```
def get_version_space(self,specific,general):
```

```
    """ Given the specific and the general boundary of the  
    version space, evaluate the version space in between """
```

```
    while get_order(VS):
```

```
        for hypothesis in VS[:]:
```

```
            hypo = list(hypothesis) # convert tuple to list for mutability
```

```
            for i,factor in enumerate(hypo):
```

```
                if factor != '?':
```

```
                    hyp=hypo[:i]
```

```
                    hyp[i]='?'
```

```
            if self.exists_general(hyp,general)and self.exists_specific(hyp,specific):
```

```
                VS.append(tuple(hyp))
```

```
    return VS
```

```
def get_order(self,hypothesis):
```

```
    pass
```

```
def more_general(self,hyp1,hyp2):
```

```
    """ Check whether hyp1 is more general than hyp2 """
```

```
    hyp = list(zip(hyp1,hyp2))
```

```
    for i,j in hyp:
```

```
        if i == '?':
```

```
            continue
```

```
        elif j == '?':
```

```
            if i != '?':
```

```
                return False
```

```
            elif i != j:
```

```
                return False
```

```
        else:
```

```
            continue
```

```
    return True
```

```
def more_specific(self,hyp1,hyp2):
```

```
    """ hyp1 more specific than hyp2 is
```

```
    equivalent to hyp2 being more general than hyp1 """
```

```
    return self.more_general(hyp2,hyp1)
```

```
dataset=[(('sunny','warm','normal','strong','warm','same'),'Y'),(('sunny','warm','high','strong','warm','same'),  
'Y'),(('rainy','cold','high','strong','warm','change'),'N'),(('sunny','warm','high','strong','cool','change'),'Y')]
```

```
#print((dataset[0][1]))

attributes =('Sky','Temp','Humidity','Wind','Water','Forecast')

f = Holder(attributes)

f.add_values('Sky',('sunny','rainy','cloudy')) #sky can be sunny rainy or cloudy
f.add_values('Temp',('cold','warm')) #Temp can be sunny cold or warm
f.add_values('Humidity',('normal','high')) #Humidity can be normal or high
f.add_values('Wind',('weak','strong')) #wind can be weak or strong
f.add_values('Water',('warm','cold')) #water can be warm or cold
f.add_values('Forecast',('same','change')) #Forecast can be same or change

a = CandidateElimination(dataset,f)

#pass the dataset to the algorithm class and call the run algorithm method

a.run_algorithm()
```

Output

```
['Sunny' 'Warm' 'Normal' 'Strong' 'Warm' 'Same']

[['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?']]

0

['Sunny' 'Warm' 'Normal' 'Strong' 'Warm' 'Same']

[['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?']]

1

['Sunny' 'Warm' '?' 'Strong' 'Warm' 'Same']

[['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?']]

2

['Sunny' 'Warm' '?' 'Strong' 'Warm' 'Same']
```

```
[[Sunny, '?', '?', '?', '?', '?'], ['?', Warm, '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', 'Same']]
```

3

```
['Sunny' Warm '?' Strong' '?' '?']
```

```
[[Sunny, '?', '?', '?', '?', '?'], ['?', Warm, '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?'], ['?', '?', '?', '?', '?', '?']]
```

Final S:

```
['Sunny' Warm '?' Strong' '?' '?']
```

Final G:

```
[[Sunny, '?', '?', '?', '?', '?'], ['?', Warm, '?', '?', '?', '?']]
```

Process finished with exit code 0

Program 3

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 to classify a new sample.

Task: ID3 determines the information gain for each candidate attribute (i.e., Outlook, Temperature, Humidity, and Wind), then selects the one with highest information gain as the root node of the tree. The information gain values for all four attributes are calculated using the following formula:

$$\text{Entropy}(S) = \sum - P(I) \cdot \log_2 P(I)$$

$$\text{Gain}(S,A) = \text{Entropy}(S) - \sum [P(S/A) \cdot \text{Entropy}(S/A)]$$

Dataset:

Table: Training examples for the target concept PlayTennis.

outlook	temperature	humidity	windy	play
sunny	hot	high	false	no
sunny	hot	high	true	no
overcast	hot	high	false	yes
rainy	mild	high	false	yes
rainy	cool	normal	false	yes
rainy	cool	normal	true	no
overcast	cool	normal	true	yes
sunny	mild	high	false	no
sunny	cool	normal	false	yes
rainy	mild	normal	false	yes
sunny	mild	normal	true	yes
overcast	mild	high	true	yes
overcast	hot	normal	false	yes
rainy	mild	high	true	no

Calculation:

Decision/play column consists of 14 instances and includes two labels: yes and no.

There are 9 decisions labeled yes and 5 decisions labeled no.

$$\begin{aligned} \text{Entropy}[\text{Decision}] &= -P(\text{yes}) \cdot \log_2 P(\text{yes}) - P(\text{no}) \cdot \log_2 P(\text{no}) \\ &= - (9/14) \cdot \log_2 (9/14) - (5/14) \cdot \log_2 (5/14) \end{aligned}$$

$$= 0.940$$

Now, we need to find out the most dominant factor for decision.

1) Wind factor on decision:

$$\text{Gain(Decision,wind)} = \text{Entropy(Decision)} - \sum [P(\text{Decision/Wind}). \text{Entropy(Decision/Wind)}]$$

Wind attribute has two labels : Weak and Strong

$$\text{Gain(Decision,Wind)} = \text{Entropy(Decision)} - [P(\text{Decision/Wind=Weak}).$$

$$\text{Entropy(Decision/Wind=Weak)}] - [P(\text{Decision/Wind=strong}).$$

$$\text{Entropy(Decision/Wind=strong)}]$$

There are 8 instances for weak. In that decision of 2 items are no and 6 items are yes.

- $$\begin{aligned} \text{Entropy[Decision/Wind= Weak]} &= -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes}) \\ &= -[2/8].\log_2(2/8) - [6/8].\log_2(6/8) \\ &= 0.811 \end{aligned}$$

- $$\begin{aligned} \text{Entropy[Decision/Wind= Strong]} &= -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes}) \\ &= -[3/6].\log_2(3/6) - [3/6].\log_2(3/6) \\ &= 1 \end{aligned}$$

Note: There are 6 instances for strong. In that decision of 3 items are yes and 3 items are no.

- $$\begin{aligned} \text{Gain(Decision,Wind)} &= 0.940 - [(8/14).0.811] - [(6/14).1] \\ &= 0.048 \end{aligned}$$

Similarly calculate gain for other factors:

2) Outlook factor on decision:

1)
$$\text{Gain(Decision,outlook)} = \text{Entropy (decision)} - \sum [P (\text{Decision /Outlook}).$$

$$\text{Entropy(Decision/Outlook)}]$$

outlook has three parameters : Sunny, Overcast, and rain

$$\text{Gain(Decision,outlook)} = \text{Entropy(decision)} -$$

$$[P(\text{Decision/Outlook=Sunny}).\text{Entropy(Decision/Outlook=Sunny)}] - [$$

$$P(\text{Decision/Outlook=overcast}).\text{Entropy(Decision/Outlook=overcast)}] -$$

$$[P(P(\text{Decision/Outlook=Rain}).\text{Entropy(Decision/Outlook=Rain)})]$$

<u>Sunny</u>	<u>Overcast:</u>	<u>Rain:</u>
Instances: 5	Instances:4	Instances:5
yes:2	yes:4	yes:3
No:3	No: -	No: 2

$$1) \text{Entropy[Decision/Outlook= Sunny]} = -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes}) \\ = 0.97094$$

$$2) \text{Entropy[Decision/Outlook= Overcast]} = -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes}) \\ = 0$$

$$3) \text{Entropy[Decision/Wind= Rain]} = -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes}) \\ = 0.9708$$

$$\text{Gain(Decision,Outlook)} = 0.940 - (5/14)(0.9709) - (4/14)(0) - (5/14)(0.9708) \\ = 0.2473$$

3) **Temperature factor on decision:**

$$\text{Gain(Decision, Temperature)} = \text{Entropy(decision)} - \\ \sum [P(\text{Decision/Temperature}).\text{Entropy(Decision/Temperature)}]$$

Temperature has 3 parameters: hot, mild, cool

$$\text{Gain(Decision, Temp)} = \text{Entropy(decision)} - \\ [P(\text{Decision/Temp=hot}).\text{Entropy(Decision/Temp=hot)}] - \\ [P(\text{Decision/Temp=mild}).\text{Entropy(Decision/Temp=mild)}] - \\ [P(\text{Decision/Temp=cool}).\text{Entropy(Decision/Temp=cool)}]$$

<u>Hot</u>	<u>Mild:</u>	<u>cool:</u>
Instances:4	Instances:6	Instances:4
Yes:2	Yes:	Yes:3
No:2	No: 2	No:1

$$\text{Entropy[Decision/Temp= hot]} = -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes}) \\ = 1$$

$$\text{Entropy[Decision/Temp= mild]} = -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes}) \\ = 0.9182$$

$$\text{Entropy[Decision/Temp= cool]} = -P[\text{no}].\log_2P(\text{no}) - p(\text{yes}).\log_2P(\text{yes})$$

$$=0.8112$$

$$\begin{aligned} \text{Gain}(\text{Decision}, \text{Temp}) &= 0.940 - (4/14)(1) - (6/14)(0.9182) - (4/14)(0.8112) \\ &= 0.0291 \end{aligned}$$

4) Humidity factor on decision:

$$\begin{aligned} \text{Gain}(\text{Decision}, \text{Humidity}) &= \text{Entropy}(\text{decision}) - \\ &\sum [P(\text{Decision}/\text{Humidity}) \cdot \text{Entropy}(\text{Decision}/\text{Humidity})] \end{aligned}$$

Humidity has 2 factors: high and normal

$$\begin{aligned} \text{Gain}(\text{Decision}, \text{humidity}) &= \text{Entropy}(\text{decision}) - \\ &[P(\text{Decision}/\text{humidity}=\text{high}) \cdot \text{Entropy}(\text{Decision}/\text{humidity}=\text{high}) - \\ &P(\text{Decision}/\text{humidity}=\text{normal}) \cdot \text{Entropy}(\text{Decision}/\text{humidity}=\text{normal})] \end{aligned}$$

<u>High</u>	<u>Normal:</u>
Instances:7	Instances:7
Yes:3	Yes:6
No:4	No: 1

$$\begin{aligned} \text{Entropy}[\text{Decision}/ \text{humidity}= \text{high}] &= -P[\text{no}] \cdot \log_2 P(\text{no}) - p(\text{yes}) \cdot \log_2 P(\text{yes}) \\ &= 0.9851 \end{aligned}$$

$$\begin{aligned} \text{Entropy}[\text{Decision}/\text{humidity}= \text{normal}] &= -P[\text{no}] \cdot \log_2 P(\text{no}) - p(\text{yes}) \cdot \log_2 P(\text{yes}) \\ &= 0.5916 \end{aligned}$$

$$\begin{aligned} \text{Gain}(\text{Decision}, \text{Humidity}) &= 0.940 - (7/14)(0.9851) - (7/14)(0.5916) \\ &= 0.1517 \end{aligned}$$

Thus the outlook factor on decision produces the highest score. That's why outlook decision will appear in the root node of the tree. Since Outlook has three possible values, the root node has three branches (sunny, overcast, rain). The next question is "what attribute should be tested at the Sunny branch node?" Since we have used Outlook at the root, we only decide on the remaining three attributes: Humidity, Temperature, or Wind.

Now calculate sunny outlook on decision, overcast outlook on decision, and rain outlook on decision to generate the decision tree.

Sunny outlook on decision:

5 instances of sunny : In that 3 instances are NO and 2 instances are YES

$$\text{Gain(Outlook = Sunny/Temp)} = 0.570$$

$$\text{Gain (Outlook = Sunny/Humidity)} = 0.970$$

$$\text{Gain (Outlook = Sunny/ Wind)} = 0.019$$

Since humidity produces the highest score, if outlook were Sunny.

Overcast outlook on decision:

Decision will always be yes, if outlook were overcast.

Rain outlook on decision:

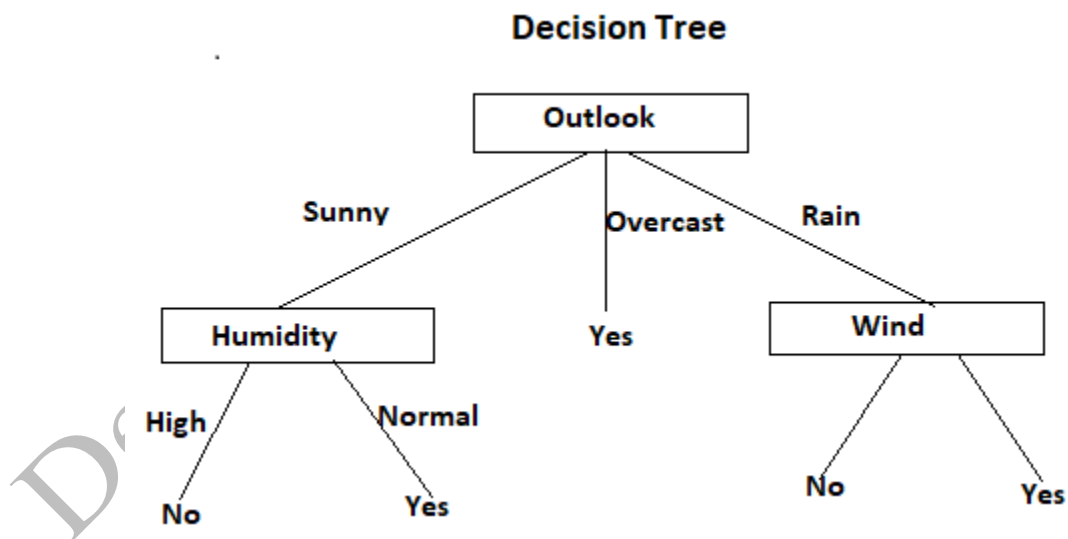
5 instances of rain : In that 3 instances are YES and 2 instances are NO.

$$\text{Gain(Outlook= Rain/Temp)}$$

$$\text{Gain(Outlook= Rain/Humidity)}$$

$$\text{Gain(Outlook= Rain/Wind)}$$

Here, wind produces the highest score . And wind has two attributes namely strong and weak.



ID3 Algorithm:

$$\text{ID3(Examples, Target_attribute, Attributes)}$$

Examples are the training examples. *Target_attribute* is the attribute whose value is to be predicted by the tree. *Attributes* is a list of other attributes that may be tested by the learned decision tree. Returns a decision tree that correctly classifies the given *Examples*.

- Create a *Root* node for the tree
 - If all *Examples* are positive, Return the single-node tree *Root*, with label = +
 - If all *Examples* are negative, Return the single-node tree *Root*, with label = -
 - If *Attributes* is empty, Return the single-node tree *Root*, with label = most common value of *Target_attribute* in *Examples*
 - Otherwise Begin
 - ❖ $A \leftarrow$ the attribute from *Attributes* that best* classifies *Examples*
 - ❖ The decision attribute for *Root* $\leftarrow A$
 - ❖ For each possible value, v_i , of A ,
 - Add a new tree branch below *Root*, corresponding to the test $A = v_i$
 - Let $Examples_{v_i}$, be the subset of *Examples* that have value v_i for A
 - If $Examples_{v_i}$, is empty
 - ❖ Then below this new branch add a leaf node with label=most common value of *Target_attribute* in *Examples*
 - ❖ Else below this new branch add the subtree
- ID3($Examples_{v_i}, Target_attribute, Attributes - \{A\}$)
- End
- Return *Root*

ID3 in Python:

```
import ast
import csv
import sys
import math
import os
def load_csv_to_header_data(filename):
    path = os.path.normpath(os.getcwd() + filename)
    fs = csv.reader(open(path))
```

```
all_row = []
for r in fs:
    all_row.append(r)

headers = all_row[0]
idx_to_name, name_to_idx = get_header_name_to_idx_maps(headers)

data = {
    'header': headers,
    'rows': all_row[1:],
    'name_to_idx': name_to_idx,
    'idx_to_name': idx_to_name
}
return data
```

```
def get_header_name_to_idx_maps(headers):
    name_to_idx = {}
    idx_to_name = {}
    for i in range(0, len(headers)):
        name_to_idx[headers[i]] = i
        idx_to_name[i] = headers[i]
    return idx_to_name, name_to_idx
```

```
def project_columns(data, columns_to_project):
    data_h = list(data['header'])
    data_r = list(data['rows'])

    all_cols = list(range(0, len(data_h)))

    columns_to_project_ix = [data['name_to_idx'][name] for name in
columns_to_project]
    columns_to_remove = [cidx for cidx in all_cols if cidx not in
columns_to_project_ix]

    for delc in sorted(columns_to_remove, reverse=True):
        del data_h[delc]
        for r in data_r:
            del r[delc]
```

```
idx_to_name, name_to_idx = get_header_name_to_idx_maps(data_h)
```

```
return {'header': data_h, 'rows': data_r,  
        'name_to_idx': name_to_idx,  
        'idx_to_name': idx_to_name}
```

```
def get_uniq_values(data):  
    idx_to_name = data['idx_to_name']  
    idxs = idx_to_name.keys()  
  
    val_map = {}  
    for idx in iter(idxs):  
        val_map[idx_to_name[idx]] = set()  
  
    for data_row in data['rows']:  
        for idx in idx_to_name.keys():  
            att_name = idx_to_name[idx]  
            val = data_row[idx]  
            if val not in val_map.keys():  
                val_map[att_name].add(val)  
    return val_map
```

```
def get_class_labels(data, target_attribute):  
    rows = data['rows']  
    col_idx = data['name_to_idx'][target_attribute]  
    labels = {}  
    for r in rows:  
        val = r[col_idx]  
        if val in labels:  
            labels[val] = labels[val] + 1  
        else:  
            labels[val] = 1  
    return labels
```

```
def entropy(n, labels):  
    ent = 0  
    for label in labels.keys():  
        p_x = labels[label] / n
```

```
    ent += - p_x * math.log(p_x, 2)
return ent
```

```
def partition_data(data, group_att):
    partitions = {}
    data_rows = data['rows']
    partition_att_idx = data['name_to_idx'][group_att]
    for row in data_rows:
        row_val = row[partition_att_idx]
        if row_val not in partitions.keys():
            partitions[row_val] = {
                'name_to_idx': data['name_to_idx'],
                'idx_to_name': data['idx_to_name'],
                'rows': list()
            }
        partitions[row_val]['rows'].append(row)
    return partitions
```

```
def avg_entropy_w_partitions(data, splitting_att, target_attribute):
    # find uniq values of splitting att
    data_rows = data['rows']
    n = len(data_rows)
    partitions = partition_data(data, splitting_att)

    avg_ent = 0

    for partition_key in partitions.keys():
        partitioned_data = partitions[partition_key]
        partition_n = len(partitioned_data['rows'])
        partition_labels = get_class_labels(partitioned_data, target_attribute)
        partition_entropy = entropy(partition_n, partition_labels)
        avg_ent += partition_n / n * partition_entropy

    return avg_ent, partitions
```

```
def most_common_label(labels):
    mcl = max(labels, key=lambda k: labels[k])
    return mcl
```



```
def id3(data, uniqs, remaining_atts, target_attribute):
    labels = get_class_labels(data, target_attribute)

    node = {}

    if len(labels.keys()) == 1:
        node['label'] = next(iter(labels.keys()))
        return node

    if len(remaining_atts) == 0:
        node['label'] = most_common_label(labels)
        return node

    n = len(data['rows'])
    ent = entropy(n, labels)

    max_info_gain = None
    max_info_gain_att = None
    max_info_gain_partitions = None

    for remaining_att in remaining_atts:
        avg_ent, partitions = avg_entropy_w_partitions(data, remaining_att,
target_attribute)
        info_gain = ent - avg_ent
        if max_info_gain is None or info_gain > max_info_gain:
            max_info_gain = info_gain
            max_info_gain_att = remaining_att
            max_info_gain_partitions = partitions

    if max_info_gain is None:
        node['label'] = most_common_label(labels)
        return node

    node['attribute'] = max_info_gain_att
    node['nodes'] = {}

    remaining_atts_for_subtrees = set(remaining_atts)
    remaining_atts_for_subtrees.discard(max_info_gain_att)

    uniq_att_values = uniqs[max_info_gain_att]
```

```
for att_value in uniq_att_values:
    if att_value not in max_info_gain_partitions.keys():
        node['nodes'][att_value] = {'label': most_common_label(labels)}
        continue
    partition = max_info_gain_partitions[att_value]
    node['nodes'][att_value] = id3(partition, uniqs, remaining_atts_for_subtrees,
target_attribute)
```

```
return node
```

```
def load_config(config_file):
    with open(config_file, 'r') as myfile:
        data = myfile.read().replace('\n', '')
    return ast.literal_eval(data)
```

```
def pretty_print_tree(root):
    stack = []
    rules = set()
```

```
def traverse(node, stack, rules):
    if 'label' in node:
        stack.append(' THEN ' + node['label'])
        rules.add("".join(stack))
        stack.pop()
    elif 'attribute' in node:
        ifnd = 'IF ' if not stack else ' AND '
        stack.append(ifnd + node['attribute'] + ' EQUALS ')
        for subnode_key in node['nodes']:
            stack.append(subnode_key)
            traverse(node['nodes'][subnode_key], stack, rules)
            stack.pop()
        stack.pop()
```

```
traverse(root, stack, rules)
print(os.linesep.join(rules))
```

```
def main():
    argv = 'tennis.cfg'
```

```
print("Command line args are {}:".format(argv))

config = load_config(argv)

data = load_csv_to_header_data(config['data_file'])
data = project_columns(data, config['data_project_columns'])

target_attribute = config['target_attribute']
remaining_attributes = set(data['header'])
remaining_attributes.remove(target_attribute)

uniqs = get_uniq_values(data)

root = id3(data, uniqs, remaining_attributes, target_attribute)

pretty_print_tree(root)

if __name__ == "__main__": main()
```

Output:

```
IF Outlook EQUALS Rainy AND Windy EQUALS True THEN No
IF Outlook EQUALS Sunny AND Humidity EQUALS Normal THEN Yes
IF Outlook EQUALS Overcast THEN Yes
IF Outlook EQUALS Rainy AND Windy EQUALS False THEN Yes
IF Outlook EQUALS Sunny AND Humidity EQUALS High THEN No
```

Program 4:

Build an Artificial Neural Network by implementing the Back propagation algorithm and test the same using appropriate data set

Back propagation Algorithm

The Backpropagation algorithm is a supervised learning method for multilayer feed-forward networks from the field of Artificial Neural Networks, the backpropagation algorithm is a method for training the weights in a multilayer feed-forward neural network. As such, it requires a network structure to be defined of one or more layers where one layer is fully connected to the next layer. A standard network structure is one input layer, one hidden layer, and one output layer. Backpropagation can be used for both classification and regression problems.

Wheat Seeds Dataset

The seeds dataset involves the prediction of species given measurements seeds from different varieties of wheat. There are 201 records and 7 numerical input variables. It is a classification problem with 3 output classes. The scale for each numeric input value vary, so some data normalization may be required for use with algorithms that weight inputs like the backpropagation algorithm.

Below is a sample of the first 5 rows of the dataset

```
1 15.26,14.84,0.871,5.763,3.312,2.221,5.22,1
2 14.88,14.57,0.8811,5.554,3.333,1.018,4.956,1
3 14.29,14.09,0.905,5.291,3.337,2.699,4.825,1
4 13.84,13.94,0.8955,5.324,3.379,2.259,4.805,1
5 16.14,14.99,0.9034,5.658,3.562,1.355,5.175,1
```

You can download the seeds dataset from the UCI Machine Learning Repository.

Download the seeds dataset and place it into your current working directory with the filename **seeds_dataset.csv**.

The dataset is in tab-separated format, so you must convert it to CSV using a text editor or a spreadsheet program.

Working of algorithm:

1. Initialize Network.
2. Forward Propagate.
3. Back Propagate Error.
4. Train the network

1. Initialize Network

Initialize the network weights to small random numbers in the range of 0 to 1. Below is a function named `initialize_network()` that creates a new neural network ready for training. It accepts three parameters, the number of inputs, the number of neurons to have in the hidden layer and the number of outputs.

The hidden layer we create `n_hidden` neurons and each neuron in the hidden layer has `n_inputs + 1` weights, one for each input column in a dataset and an additional one for the bias. The output layer that connects to the hidden layer has `n_outputs` neurons, each with `n_hidden + 1` weights. This means that each neuron in the output layer connects to (has a weight for) each neuron in the hidden layer.

Let's test out this function. Below is a complete example that creates a small network.

```
1 from random import seed
2 from random import random
3
4 # Initialize a network
5 def initialize_network(n_inputs, n_hidden, n_outputs):
6     network = list()
7     hidden_layer = [{'weights':[random() for i in range(n_inputs + 1)]] for i in
8     range(n_hidden)]
9     network.append(hidden_layer)
10    output_layer = [{'weights':[random() for i in range(n_hidden + 1)]] for i in
11    range(n_outputs)]
12    network.append(output_layer)
13    return network
14
15    seed(1)
16    network = initialize_network(2, 1, 2)
    for layer in network:
        print(layer)
```

Running the example, you can see that the code prints out each layer one by one. You can see the hidden layer has one neuron with 2 input weights plus the bias. The output layer has 2 neurons, each with 1 weight plus the bias.

```
1 [{"weights": [0.13436424411240122, 0.8474337369372327, 0.763774618976614]}]
2 [{"weights": [0.2550690257394217, 0.49543508709194095]}, {"weights":
[0.4494910647887381, 0.651592972722763]}]
```

Forward Propagate

We can calculate an output from a neural network by propagating an input signal through each layer until the output layer outputs its values. We call this forward-propagation.

We can break forward propagation down into three parts:

- Neuron Activation.
- Neuron Transfer.
- Forward Propagation.

Neuron Activation

The first step is to calculate the activation of one neuron given an input.

The input could be a row from our training dataset, as in the case of the hidden layer. It may also be the outputs from each neuron in the hidden layer, in the case of the output layer. Neuron activation is calculated as the weighted sum of the inputs. Much like linear regression.

$$\text{activation} = \text{sum}(\text{weight}_i * \text{input}_i) + \text{bias}$$

$$\text{net}_{h1} = w_1 * i_1 + w_2 * i_2 + b_1 * 1$$

Where weight is a network weight, input is an input, i is the index of a weight or an input and bias is a special weight that has no input to multiply with (or you can think of the input as always being 1.0). Below is an implementation of this in a function named `activate()`. You can see that the function assumes that the bias is the last weight in the list of weights. This helps here and later to make the code easier to read.

```
1 # Calculate neuron activation for an input
2 def activate(weights, inputs):
3     activation = weights[-1]
4     for i in range(len(weights)-1):
5         activation += weights[i] * inputs[i]
6     return activation
```

Neuron Transfer

Once a neuron is activated, we need to transfer the activation to see what the neuron output actually is. Different transfer functions can be used. It is traditional to use the sigmoid activation function. It is also a function of which we can easily calculate the derivative (slope) that we will need later when backpropagating error. We can transfer an activation function using the sigmoid function as follows:

$$\text{output} = 1 / (1 + e^{-(\text{activation})})$$

Where e is the base of the natural logarithms (Euler's number).

Below is a function named `transfer()` that implements the sigmoid equation.

```
1 # Transfer neuron activation
2 def transfer(activation):
3     return 1.0 / (1.0 + exp(-activation))
```

Forward Propagation

Forward propagating an input is straightforward. All of the outputs from one layer become inputs to the neurons on the next layer. Below is a function named `forward_propagate()` that implements the forward propagation for a row of data from our dataset with our neural network.

The function returns the outputs from the last layer also called the output layer.

```
1 # Forward propagate input to a network output
2 def forward_propagate(network, row):
3     inputs = row
4     for layer in network:
5         new_inputs = []
6         for neuron in layer:
7             activation = activate(neuron['weights'], inputs)
8             neuron['output'] = transfer(activation)
9             new_inputs.append(neuron['output'])
10    inputs = new_inputs
11    return inputs
```

Running the example propagates the input pattern `[1, 0]` and produces an output value that is printed. Because the output layer has two neurons, we get a list of two numbers as output.

```
[0.6629970129852887, 0.7253160725279748]
```

Back Propagate Error

The backpropagation algorithm is named for the way in which weights are trained. Error is calculated between the expected outputs and the outputs forward propagated from the network. These errors are then propagated backward through the network from the output layer to the hidden layer, assigning blame for the error and updating weights as they go. The math for backpropagating error is rooted in calculus, but we will remain high level in this section and focus on what is calculated and how rather than why the calculations take this particular form.

This part is broken down into two sections.

Transfer Derivative.

Error Backpropagation.

Transfer Derivative

Given an output value from a neuron, we need to calculate its slope. We are using the sigmoid transfer function, the derivative of which can be calculated as follows:

```
1 derivative = output * (1.0 - output)
```

Below is a function named `transfer_derivative()` that implements this equation.

```
1 # Calculate the derivative of an neuron output
2 def transfer_derivative(output):
3     return output * (1.0 - output)
```

Now, let's see how this can be used.

Error Backpropagation

The first step is to calculate the error for each output neuron, this will give us our error signal (input) to propagate backwards through the network.

The error for a given neuron can be calculated as follows:

$$\text{error} = (\text{expected} - \text{output}) * \text{transfer_derivative}(\text{output})$$

Where expected is the expected output value for the neuron, output is the output value for the neuron and transfer_derivative() calculates the slope of the neuron's output value, as shown above.

The error signal for a neuron in the hidden layer is calculated as the weighted error of each neuron in the output layer. Think of the error traveling back along the weights of the output layer to the neurons in the hidden layer.

The back-propagated error signal is accumulated and then used to determine the error for the neuron in the hidden layer, as follows:

$$\text{error} = (\text{weight_k} * \text{error_j}) * \text{transfer_derivative}(\text{output})$$

Where error_j is the error signal from the jth neuron in the output layer, weight_k is the weight that connects the kth neuron to the current neuron and output is the output for the current neuron.

Below is a function named backward_propagate_error() that implements this procedure. You can see that the error signal calculated for each neuron is stored with the name 'delta'. You can see that the layers of the network are iterated in reverse order, starting at the output and working backwards. This ensures that the neurons in the output layer have 'delta' values calculated first that neurons in the hidden layer can use in the subsequent iteration. I chose the name 'delta' to reflect the change the error implies on the neuron (e.g. the weight delta). You can see that the error signal for neurons in the hidden layer is accumulated from neurons in the output layer where the hidden neuron number j is also the index of the neuron's weight in the output layer neuron['weights'][j].

```
1 # Backpropagate error and store in neurons
2 def backward_propagate_error(network, expected):
3     for i in reversed(range(len(network))):
4         layer = network[i]
5         errors = list()
6         if i != len(network)-1:
7             for j in range(len(layer)):
8                 error = 0.0
9                 for neuron in network[i + 1]:
10                    error += (neuron['weights'][j] * neuron['delta'])
11                errors.append(error)
12            else:
13                for j in range(len(layer)):
14                    neuron = layer[j]
15                    errors.append(expected[j] - neuron['output'])
16                for j in range(len(layer)):
```



```
17 neuron = layer[j]
18 neuron['delta'] = errors[j] * transfer_derivative(neuron['output'])
```

Running the example prints the network after the backpropagation of error is complete. You can see that error values are calculated and stored in the neurons for the output layer and the hidden layer.

```
1 [{ 'output': 0.7105668883115941, 'weights': [0.13436424411240122, 0.8474337369372327,
2 0.763774618976614], 'delta': -0.0005348048046610517}]
  [{ 'output': 0.6213859615555266, 'weights': [0.2550690257394217, 0.49543508709194095],
  'delta': -0.14619064683582808}, { 'output': 0.6573693455986976, 'weights':
  [0.4494910647887381, 0.651592972722763], 'delta': 0.0771723774346327}]
```

Now let's use the backpropagation of error to train the network.

Train Network

The network is trained using stochastic gradient descent. This involves multiple iterations of exposing a training dataset to the network and for each row of data forward propagating the inputs, backpropagating the error and updating the network weights.

This part is broken down into two sections:

- Update Weights.
- Train Network.

Update Weights

Once errors are calculated for each neuron in the network via the back propagation method above, they can be used to update weights. Network weights are updated as follows:

```
weight = weight + learning_rate * error * input
```

Where weight is a given weight, learning_rate is a parameter that you must specify, error is the error calculated by the backpropagation procedure for the neuron and input is the input value that caused the error.

Learning rate controls how much to change the weight to correct for the error. Below is a function named update_weights() that updates the weights for a network given an input row of data, a learning rate and assume that a forward and backward propagation have already been performed. Remember that the input for the output layer is a collection of outputs from the hidden layer.

```
1 # Update network weights with error
2 def update_weights(network, row, l_rate):
3     for i in range(len(network)):
4         inputs = row[:-1]
5         if i != 0:
6             inputs = [neuron['output'] for neuron in network[i - 1]]
7         for neuron in network[i]:
8             for j in range(len(inputs)):
9                 neuron['weights'][j] += l_rate * neuron['delta'] * inputs[j]
```

```
10 neuron['weights'][-1] += l_rate * neuron['delta']
```

Now we know how to update network weights, let's see how we can do it repeatedly.

Train Network

As mentioned, the network is updated using stochastic gradient descent. This involves first looping for a fixed number of epochs and within each epoch updating the network for each row in the training dataset. Below is a function that implements the training of an already initialized neural network with a given training dataset, learning rate, fixed number of epochs and an expected number of output values. The sum squared error between the expected output and the network output is accumulated each epoch and printed.

```
1 # Train a network for a fixed number of epochs
2 def train_network(network, train, l_rate, n_epoch, n_outputs):
3     for epoch in range(n_epoch):
4         sum_error = 0
5         for row in train:
6             outputs = forward_propagate(network, row)
7             expected = [0 for i in range(n_outputs)]
8             expected[row[-1]] = 1
9             sum_error += sum([(expected[i]-outputs[i])**2 for i in range(len(expected))])
10            backward_propagate_error(network, expected)
11            update_weights(network, row, l_rate)
12            print('>epoch=%d, lrate=%.3f, error=%.3f' % (epoch, l_rate, sum_error))
```

ANN in python:

```
from math import exp
from random import seed
from random import random
```

```
# Initialize a network
```

```
def initialize_network(n_inputs, n_hidden, n_outputs):
    network = list()
    hidden_layer = [{'weights':[random() for i in range(n_inputs + 1)]] for i in range(n_hidden)]
    network.append(hidden_layer)
    output_layer = [{'weights':[random() for i in range(n_hidden + 1)]] for i in range(n_outputs)]
```

```
network.append(output_layer)
return network

# Calculate neuron activation for an input
def activate(weights, inputs):
    activation = weights[-1]
    for i in range(len(weights)-1):
        activation += weights[i] * inputs[i]
    return activation

# Transfer neuron activation
def transfer(activation):
    return 1.0 / (1.0 + exp(-activation))

# Forward propagate input to a network output
def forward_propagate(network, row):
    inputs = row
    for layer in network:
        new_inputs = []
        for neuron in layer:
            activation = activate(neuron['weights'], inputs)
            neuron['output'] = transfer(activation)
            new_inputs.append(neuron['output'])
        inputs = new_inputs
    return inputs

# Calculate the derivative of an neuron output
def transfer_derivative(output):
    return output * (1.0 - output)

# Backpropagate error and store in neurons
def backward_propagate_error(network, expected):
    for i in reversed(range(len(network))):
        layer = network[i]
        errors = list()
        if i != len(network)-1:
            for j in range(len(layer)):
                error = 0.0
                for neuron in network[i + 1]:
                    error += (neuron['weights'][j] * neuron['delta'])
                errors.append(error)
        else:
            for j in range(len(layer)):
                neuron = layer[j]
```

```
        errors.append(expected[j] - neuron['output'])
    for j in range(len(layer)):
        neuron = layer[j]
        neuron['delta'] = errors[j] * transfer_derivative(neuron['output'])

# Update network weights with error
def update_weights(network, row, l_rate):
    for i in range(len(network)):
        inputs = row[:-1]
        if i != 0:
            inputs = [neuron['output'] for neuron in network[i - 1]]
        for neuron in network[i]:
            for j in range(len(inputs)):
                neuron['weights'][j] += l_rate * neuron['delta'] * inputs[j]
            neuron['weights'][-1] += l_rate * neuron['delta']

# Train a network for a fixed number of epochs
def train_network(network, train, l_rate, n_epoch, n_outputs):
    for epoch in range(n_epoch):
        sum_error = 0
        for row in train:
            outputs = forward_propagate(network, row)
            expected = [0 for i in range(n_outputs)]
            expected[row[-1]] = 1
            sum_error += sum([(expected[i]-outputs[i])**2 for i in
range(len(expected))])

            backward_propagate_error(network, expected)
            update_weights(network, row, l_rate)
        print('>epoch=%d, lrate=%.3f, error=%.3f' % (epoch, l_rate, sum_error))

# Test training backprop algorithm
seed(1)
dataset = [[2.7810836,2.550537003,0],
[1.465489372,2.362125076,0],
[3.396561688,4.400293529,0],
[1.38807019,1.850220317,0],
[3.06407232,3.005305973,0],
[7.627531214,2.759262235,1],
[5.332441248,2.088626775,1],
[6.922596716,1.77106367,1],
[8.675418651,-0.242068655,1],
[7.673756466,3.508563011,1]]
n_inputs = len(dataset[0]) - 1
n_outputs = len(set([row[-1] for row in dataset]))
```

```
network = initialize_network(n_inputs, 2, n_outputs)
train_network(network, dataset, 0.5, 20, n_outputs)
for layer in network:
    print(layer)
```

Output

```
>epoch=0, lrate=0.500, error=6.350
>epoch=1, lrate=0.500, error=5.531
>epoch=2, lrate=0.500, error=5.22147
>epoch=3, lrate=0.500, error=4.951
>epoch=4, lrate=0.500, error=4.519
>epoch=5, lrate=0.500, error=4.173
>epoch=6, lrate=0.500, error=3.835
>epoch=7, lrate=0.500, error=3.506
>epoch=8, lrate=0.500, error=3.192
>epoch=9, lrate=0.500, error=2.898
>epoch=10, lrate=0.500, error=2.626
>epoch=11, lrate=0.500, error=2.377
>epoch=12, lrate=0.500, error=2.153
>epoch=13, lrate=0.500, error=1.953
>epoch=14, lrate=0.500, error=1.774
>epoch=15, lrate=0.500, error=1.614
>epoch=16, lrate=0.500, error=1.472
>epoch=17, lrate=0.500, error=1.346
>epoch=18, lrate=0.500, error=1.233
>epoch=19, lrate=0.500, error=1.132
```

```
[{'weights': [-1.4688375095432327, 1.850887325439514, 1.0858178629550297], 'output':
0.029980305604426185, 'delta': -0.0059546604162323625}, {'weights': [0.37711098142462157, -
```

```
0.0625909894552989, 0.2765123702642716], 'output': 0.9456229000211323, 'delta':  
0.0026279652850863837}]
```

```
[{'weights': [2.515394649397849, -0.3391927502445985, -0.9671565426390275], 'output':  
0.23648794202357587, 'delta': -0.04270059278364587}, {'weights': [-2.5584149848484263,  
1.0036422106209202, 0.42383086467582715], 'output': 0.7790535202438367, 'delta':  
0.03803132596437354}]
```

```
>>>
```



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Program 5

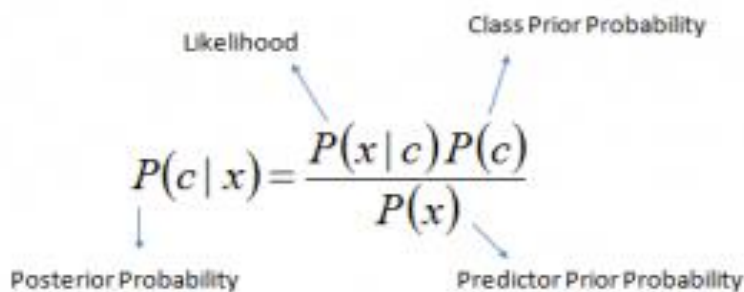
Write a program to implement the naïve Bayesian classifier for a sample training data set stored as a .CSV file. Compute the accuracy of the classifier, considering few test data sets.

Task: It is a classification technique based on Bayes' Theorem with an assumption of independence among predictors. In simple terms, a Naive Bayes classifier assumes that the presence of a particular feature in a class is unrelated to the presence of any other feature. For example, a fruit may be considered to be an apple if it is red, round, and about 3 inches in diameter. Even if these features depend on each other or upon the existence of the other features, all of these properties independently contribute to the probability that this fruit is an apple and that is why it is known as 'Naive'.

Data Set : PlayTennis example

Day	Outlook	Temperature	Humidity	Wind	PlayTennis
D1	Sunny	Hot	High	Weak	No
D2	Sunny	Hot	High	Strong	No
D3	Overcast	Hot	High	Weak	Yes
D4	Rain	Mild	High	Weak	Yes
D5	Rain	Cool	Normal	Weak	Yes
D6	Rain	Cool	Normal	Strong	No
D7	Overcast	Cool	Normal	Strong	Yes
D8	Sunny	Mild	High	Weak	No
D9	Sunny	Cool	Normal	Weak	Yes
D10	Rain	Mild	Normal	Weak	Yes
D11	Sunny	Mild	Normal	Strong	Yes
D12	Overcast	Mild	High	Strong	Yes
D13	Overcast	Hot	Normal	Weak	Yes
D14	Rain	Mild	High	Strong	No

Bayes theorem provides a way of calculating posterior probability $P(c|x)$ from $P(c)$, $P(x)$ and $P(x|c)$.



$$P(c | X) = P(x_1 | c) \times P(x_2 | c) \times \dots \times P(x_n | c) \times P(c)$$

Above,

- $P(c|x)$ is the posterior probability of class (c , target) given predictor (x , attributes).
- $P(c)$ is the prior probability of class.
- $P(x|c)$ is the likelihood which is the probability of predictor given class.
- $P(x)$ is the prior probability of predictor.

So here we have 4 attributes. What we need to do is to create “look-up tables” for each of these attributes, and write in the probability that a game of tennis will be played based on this attribute. In these tables we have to note that there are 5 cases of not being able to play a game, and 9 cases of being able to play a game.

We also must note the following probabilities for $P(C)$:

$$P(\text{Play}=\text{Yes}) = 9/14$$

$$P(\text{Play}=\text{No}) = 5/14$$

OUTLOOK	Play = Yes	Play = No	Total
Sunny	2/9	3/5	5/14
Overcast	4/9	0/5	4/14
Rain	3/9	2/5	5/14

TEMPERATURE	Play = Yes	Play = No	Total
Hot	2/9	2/5	4/14
Mild	4/9	2/5	6/14
Cool	3/9	1/5	4/14

HUMIDITY	Play = Yes	Play = No	Total
High	3/9	4/5	7/14
Normal	6/9	1/5	7/14

WIND	Play = Yes	Play = No	Total
------	------------	-----------	-------

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Strong	3/9	3/5	6/14
Weak	6/9	2/5	8/14

Testing

For this, say we were given a new instance, and we want to know if we can play a game or not, then we need to lookup the results from the tables above. So, this new instance is:

$X = (\text{Outlook}=\text{Sunny}, \text{Temperature}=\text{Cool}, \text{Humidity}=\text{High}, \text{Wind}=\text{Strong})$

Firstly we look at the probability that we can play the game, so we use the lookup tables to get:

$$\begin{aligned}P(\text{Outlook}=\text{Sunny} \mid \text{Play}=\text{Yes}) &= 2/9 \\P(\text{Temperature}=\text{Cool} \mid \text{Play}=\text{Yes}) &= 3/9 \\P(\text{Humidity}=\text{High} \mid \text{Play}=\text{Yes}) &= 3/9 \\P(\text{Wind}=\text{Strong} \mid \text{Play}=\text{Yes}) &= 3/9 \\P(\text{Play}=\text{Yes}) &= 9/14\end{aligned}$$

Next we consider the fact that we cannot play a game:

$$\begin{aligned}P(\text{Outlook}=\text{Sunny} \mid \text{Play}=\text{No}) &= 3/5 \\P(\text{Temperature}=\text{Cool} \mid \text{Play}=\text{No}) &= 1/5 \\P(\text{Humidity}=\text{High} \mid \text{Play}=\text{No}) &= 4/5 \\P(\text{Wind}=\text{Strong} \mid \text{Play}=\text{No}) &= 3/5 \\P(\text{Play}=\text{No}) &= 5/14\end{aligned}$$

Then, using those results, you have to multiple the whole lot together. So you multiple all the probabilities for $\text{Play}=\text{Yes}$ such as:

$$P(X \mid \text{Play}=\text{Yes})P(\text{Play}=\text{Yes}) = (2/9) * (3/9) * (3/9) * (3/9) * (9/14) = 0.0053$$

And this gives us a value that represents ' $P(X|C)P(C)$ ', or in this case ' $P(X|\text{Play}=\text{Yes})P(\text{Play}=\text{Yes})$ '.

We also have to do the same thing for $\text{Play}=\text{No}$:

$$P(X \mid \text{Play}=\text{No})P(\text{Play}=\text{No}) = (3/5) * (1/5) * (4/5) * (3/5) * (5/14) = 0.0206$$

Finally, we have to divide both results by the evidence, or ' $P(X)$ '. The evidence for both equations is the same, and we can find the values we need within the 'Total' columns of the look-up tables. Therefore:

$$\begin{aligned}P(X) &= P(\text{Outlook}=\text{Sunny}) * P(\text{Temperature}=\text{Cool}) * P(\text{Humidity}=\text{High}) * P(\text{Wind}=\text{Strong}) \\P(X) &= (5/14) * (4/14) * (7/14) * (6/14) \\P(X) &= 0.02186\end{aligned}$$

Then, dividing the results by this value:

$$\begin{aligned}P(\text{Play}=\text{Yes} \mid X) &= 0.0053/0.02186 = 0.2424 \\P(\text{Play}=\text{No} \mid X) &= 0.0206/0.02186 = 0.9421\end{aligned}$$

So, given the probabilities, can we play a game or not? To do this, we look at both probabilities and see which one has the highest value, and that is our answer. Therefore:

$$P(\text{Play}=\text{Yes} \mid X) = 0.2424$$

$$P(\text{Play}=\text{No} \mid X) = 0.9421$$

Since 0.9421 is greater than 0.2424 then the answer is 'no', we cannot play a game of tennis today.

Let's understand it using an example. Below I have a training data set of weather and corresponding target variable 'Play' (suggesting possibilities of playing). Now, we need to classify whether players will play or not based on weather condition. Let's follow the below steps to perform it.

Step 1: Convert the data set into a frequency table

Step 2: Create Likelihood table by finding the probabilities like Overcast probability = 0.29 and probability of playing is 0.64.

Step 3: Now, use Naive Bayesian equation to calculate the posterior probability for each class. The class with the highest posterior probability is the outcome of prediction.

Weather	Play
Sunny	No
Overcast	Yes
Rainy	Yes
Sunny	Yes
Sunny	Yes
Overcast	Yes
Rainy	No
Rainy	No
Sunny	Yes
Rainy	Yes
Sunny	No
Overcast	Yes
Overcast	Yes
Rainy	No

Frequency Table		
Weather	No	Yes
Overcast		4
Rainy	3	2
Sunny	2	3
Grand Total	5	9

Likelihood table				
Weather	No	Yes		
Overcast		4	=4/14	0.29
Rainy	3	2	=5/14	0.36
Sunny	2	3	=5/14	0.36
All	5	9		
	=5/14	=9/14		
	0.36	0.64		

Naïve Bayes for PlayTennis dataset in python:

```
print("\nNaive Bayes Classifier for concept learning problem")
import csv
import random
import math
import operator
def safe_div(x,y):
if y == 0:
```

```
return 0
return x / y

def loadCsv(filename):
    lines = csv.reader(open(filename))
    dataset = list(lines)
    for i in range(len(dataset)):
        dataset[i] = [float(x) for x in dataset[i]]
    return dataset

def splitDataset(dataset, splitRatio):
    trainSize = int(len(dataset) * splitRatio)
    trainSet = []
    copy = list(dataset)
    i=0
    while len(trainSet) < trainSize:
        #index = random.randrange(len(copy))

        trainSet.append(copy.pop(i))
    return [trainSet, copy]

def separateByClass(dataset):
    separated = {}
    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 safe_div(sum(numbers),float(len(numbers)))

def stdev(numbers):
    avg = mean(numbers)
    variance = safe_div(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 = {}
for classValue, instances in separated.items():
    summaries[classValue] = summarize(instances)
return summaries

def calculateProbability(x, mean, stdev):
    exponent = math.exp(-safe_div(math.pow(x-mean,2),(2*math.pow(stdev,2))))
    final = safe_div(1, (math.sqrt(2*math.pi) * stdev)) * exponent
    return final

def calculateClassProbabilities(summaries, inputVector):
    probabilities = {}
    for classValue, classSummaries in summaries.items():
        probabilities[classValue] = 1
        for i in range(len(classSummaries)):
            mean, stdev = classSummaries[i]
            x = inputVector[i]
            probabilities[classValue] *= calculateProbability(x, mean, stdev)
    return probabilities

def predict(summaries, inputVector):
    probabilities = calculateClassProbabilities(summaries, inputVector)
    bestLabel, bestProb = None, -1
    for classValue, probability in probabilities.items():
        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
    accuracy = safe_div(correct,float(len(testSet))) * 100.0
    return accuracy
```

```
def main():
    filename = 'ConceptLearning.csv'
    splitRatio = 0.9
    dataset = loadCsv(filename)
    trainingSet, testSet = splitDataset(dataset, splitRatio)
    print('Split {0} rows into'.format(len(dataset)))
    print('Number of Training data: ' + (repr(len(trainingSet))))
    print('Number of Test Data: ' + (repr(len(testSet))))
    print("\nThe values assumed for the concept learning attributes are\n")
    print("OUTLOOK=> Sunny=1 Overcast=2 Rain=3\nTEMPERATURE=> Hot=1 Mild=2
Cool=3\nHUMIDITY=> High=1 Normal=2\nWIND=> Weak=1 Strong=2")
    print("TARGET CONCEPT:PLAY TENNIS=> Yes=10 No=5")
    print("\nThe Training set are:")
    for x in trainingSet:
        print(x)
    print("\nThe Test data set are:")
    for x in testSet:
        print(x)
    print("\n")
    # prepare model
    summaries = summarizeByClass(trainingSet)
    # test model
    predictions = getPredictions(summaries, testSet)
    actual = []
    for i in range(len(testSet)):
        vector = testSet[i]
        actual.append(vector[-1])
# Since there are five attribute values, each attribute constitutes to 20% accuracy. So if all attributes
#match with predictions then 100% accuracy
    print('Actual values: {0}%'.format(actual))
    print('Predictions: {0}%'.format(predictions))
    accuracy = getAccuracy(testSet, predictions)
    print('Accuracy: {0}%'.format(accuracy))
```

main()

The input CSV file is as below



ConceptLearning.csv

Output:

Naive Bayes Classifier for concept learning problem

Split 16 rows into

Number of Training data: 14

Number of Test Data: 2

The values assumed for the concept learning attributes are

OUTLOOK=> Sunny=1 Overcast=2 Rain=3

TEMPERATURE=> Hot=1 Mild=2 Cool=3

HUMIDITY=> High=1 Normal=2

WIND=> Weak=1 Strong=2

TARGET CONCEPT:PLAY TENNIS=> Yes=10 No=5

The Training set are:

[1.0, 1.0, 1.0, 1.0, 5.0]

[1.0, 1.0, 1.0, 2.0, 5.0]

[2.0, 1.0, 1.0, 2.0, 10.0]

[3.0, 2.0, 1.0, 1.0, 10.0]

[3.0, 3.0, 2.0, 1.0, 10.0]

[3.0, 3.0, 2.0, 2.0, 5.0]

[2.0, 3.0, 2.0, 2.0, 10.0]

[1.0, 2.0, 1.0, 1.0, 5.0]

[1.0, 3.0, 2.0, 1.0, 10.0]

[3.0, 2.0, 2.0, 2.0, 10.0]

[1.0, 2.0, 2.0, 2.0, 10.0]

[2.0, 2.0, 1.0, 2.0, 10.0]

[2.0, 1.0, 2.0, 1.0, 10.0]

[3.0, 2.0, 1.0, 2.0, 5.0]

The Test data set are:

[1.0, 2.0, 1.0, 2.0, 5.0]

[1.0, 2.0, 1.0, 2.0, 5.0]

Actual values: [5.0, 5.0]%

Predictions: [5.0, 5.0]%

Accuracy: 100.0%

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Program 6

Assuming a set of documents that need to be classified, use the naïve Bayesian Classifier model to perform this task. Built-in Java classes/API can be used to write the program. Calculate the accuracy, precision, and recall for your data set.

Multinomial Naïve Bayes Classifier

LEARN_NAIVE_BAYES_TEXT(*Examples*, *V*)

Examples is a set of text documents along with their target values. *V* is the set of all possible target values. This function learns the probability terms $P(w_k|v_j)$, describing the probability that a randomly drawn word from a document in class v_j will be the English word w_k . It also learns the class prior probabilities $P(v_j)$.

1. collect all words, punctuation, and other tokens that occur in *Examples*
 - *Vocabulary* ← the set of all distinct words and other tokens occurring in any text document from *Examples*
2. calculate the required $P(v_j)$ and $P(w_k|v_j)$ probability terms
 - For each target value v_j in *V* do
 - $docs_j$ ← the subset of documents from *Examples* for which the target value is v_j
 - $P(v_j) \leftarrow \frac{|docs_j|}{|Examples|}$
 - $Text_j$ ← a single document created by concatenating all members of $docs_j$
 - $n \leftarrow$ total number of distinct word positions in $Text_j$
 - for each word w_k in *Vocabulary*
 - $n_k \leftarrow$ number of times word w_k occurs in $Text_j$
 - $P(w_k|v_j) \leftarrow \frac{n_k+1}{n+|Vocabulary|}$

CLASSIFY_NAIVE_BAYES_TEXT(*Doc*)

Return the estimated target value for the document *Doc*. a_i denotes the word found in the *i*th position within *Doc*.

- *positions* ← all word positions in *Doc* that contain tokens found in *Vocabulary*
- Return v_{NB} , where

$$v_{NB} = \operatorname{argmax}_{v_j \in V} P(v_j) \prod_{i \in \text{positions}} P(a_i|v_j)$$

The 20newsgroups dataset

comp.graphics comp.os.ms-windows.misc comp.sys.ibm.pc.hardware comp.sys.mac.hardware comp.windows.x	rec.autos rec.motorcycles rec.sport.baseball rec.sport.hockey	sci.crypt sci.electronics sci.med sci.space
misc.forsale	talk.politics.misc talk.politics.guns talk.politics.mideast	talk.religion.misc alt.atheism soc.religion.christian

The data set contained two primary folders named as ‘train’ and ‘test.’ Each folder further contained 20 folders, one for each category of documents. Inside these folders were the actual

documents. Each category contained around 600 train documents and around 400 test documents.

Python Code

```
from sklearn.datasets import fetch_20newsgroups
from sklearn.metrics import confusion_matrix
from sklearn.metrics import classification_report
import numpy as np
categories = ['alt.atheism', 'soc.religion.christian','comp.graphics', 'sci.med']
twenty_train = fetch_20newsgroups(subset='train',categories=categories,shuffle=True)
twenty_test = fetch_20newsgroups(subset='test',categories=categories,shuffle=True)
print(len(twenty_train.data))
print(len(twenty_test.data))
print(twenty_train.target_names)
print("\n".join(twenty_train.data[0].split("\n")))
print(twenty_train.target[0])
from sklearn.feature_extraction.text import CountVectorizer
count_vect = CountVectorizer()
X_train_tf = count_vect.fit_transform(twenty_train.data)
from sklearn.feature_extraction.text import TfidfTransformer
tfidf_transformer = TfidfTransformer()
X_train_tfidf = tfidf_transformer.fit_transform(X_train_tf)
X_train_tfidf.shape
from sklearn.naive_bayes import MultinomialNB
from sklearn.metrics import accuracy_score
from sklearn import metrics
mod = MultinomialNB()
mod.fit(X_train_tfidf, twenty_train.target)
X_test_tf = count_vect.transform(twenty_test.data)
X_test_tfidf = tfidf_transformer.transform(X_test_tf)
predicted = mod.predict(X_test_tfidf)
print("Accuracy:", accuracy_score(twenty_test.target, predicted))
print(classification_report(twenty_test.target,predicted,target_names=twenty_test.target_names))

print("confusion matrix is \n",metrics.confusion_matrix(twenty_test.target, predicted))
```

Output

2257

1502

['alt.atheism', 'comp.graphics', 'sci.med', 'soc.religion.christian']

From: sd345@city.ac.uk (Michael Collier)

Subject: Converting images to HP LaserJet III?

Nntp-Posting-Host: hampton

Organization: The City University

Lines: 14

Does anyone know of a good way (standard PC application/PD utility) to convert tif/img/tga files into LaserJet III format. We would also like to do the same, converting to HPGL (HP plotter) files.

Please email any response.

Is this the correct group?

Thanks in advance. Michael.

--

Michael Collier (Programmer)

Email: M.P.Collier@uk.ac.city

Tel: 071 477-8000 x3769

Fax: 071 477-8565

The Computer Unit,

The City University,

London,

EC1V 0HB.

1

Accuracy: 0.8348868175765646

	precision	recall	f1-score	support
alt.atheism	0.97	0.60	0.74	319
comp.graphics	0.96	0.89	0.92	389
sci.med	0.97	0.81	0.88	396
soc.religion.christian	0.65	0.99	0.78	398
avg / total	0.88	0.83	0.84	1502

confusion matrix is

[[192 2 6 119]

[2 347 4 36]

[2 11 322 61]

[2 2 1 393]]

>>>

Testing and Evaluation: The effectiveness of the model is judged by employing the metrics described below. All the metrics obtained range from 0 to 1 where 1 being the ideal and 0 being the worst.

Confusion Matrix: A confusion matrix is a technique for summarizing the performance of a classification algorithm. Positive (P) the number of real positive cases in the data. Negative (N) the number of real negative cases in the data

		Classifier Prediction	
		Positive	Negative
Actual Value	Positive	True Positive	False Negative
	Negative	False Positive	True Negative

The classification **Accuracy** is defined as:

$$\text{Accuracy} = \frac{TP+TN}{(TP+FP+TN+FN)}$$

Precision (P) can be defined as the number of **True positives** (Tp) over the number of **False positives** (Fp) plus the number of True positives (Tp).

$$\text{Precision} = \frac{TP}{(TP + FP)}$$

Recall (R) is defined as the number of **True Positives** (Tp) over the number of **False Negatives** (Fp) plus the number of True Positives (Tp).

$$\text{Recall} = \frac{TP}{(TP+FN)}$$

The **F₁ score** is the harmonic average of the precision and recall, where an F₁ score reaches its best value at 1 (perfect precision and recall) and worst at 0.

$$F1 = 2 \times \frac{(\text{precision} \times \text{recall})}{(\text{precision} + \text{recall})}$$

Consider alt.atheism newsgroup using confusion matrix calculate

$$\text{Accuracy} = \frac{192+1117}{192+127+6+11170}$$

$$\text{Precision} = \frac{192}{192+6} = 0.97$$

$$\text{Recall} = \frac{192}{192+127} = 0.60$$

$$F1\text{score} = 2 \times \frac{(0.97 \times 0.60)}{(0.97+0.60)} = 0.74$$

Program 7

Write a program to construct a Bayesian network considering medical data. Use this model to demonstrate the diagnosis of heart patients using standard Heart Disease Data Set. You can use Java/Python ML library classes/API.

A Bayesian belief network describes the probability distribution over a set of variables.

Probability

$P(A)$ is used to denote the probability of A . For example if A is discrete with states $\{\text{True}, \text{False}\}$ then $P(A)$ might equal $[0.2, 0.8]$. I.e. 20% chance of being True, 80% chance of being False.

Joint probability

A joint probability refers to the probability of more than one variable occurring together, such as the probability of A and B , denoted $P(A,B)$.

Conditional probability

Conditional probability is the probability of a variable (or set of variables) given another variable (or set of variables), denoted $P(A|B)$. For example, the probability of Windy being True, given that Raining is True might equal 50%. This would be denoted $P(\text{Windy} = \text{True} | \text{Raining} = \text{True}) = 50\%$.

Once the structure has been defined (i.e. nodes and links), a Bayesian network requires a probability distribution to be assigned to each node. Each node X in a Bayesian network requires a probability distribution $P(X | \text{pa}(X))$. Note that if a node X has no parents $\text{pa}(X)$ is empty, and the required distribution is just $P(X)$ sometimes referred to as the prior. This is the probability of itself given its parent nodes.

If $U = \{A_1, \dots, A_n\}$ is the universe of variables (all the variables) in a Bayesian network, and $\text{pa}(A_i)$ are the parents of A_i then the joint probability distribution $P(U)$ is simply the product of all the probability distributions (prior and conditional) in the network, as shown in the equation below. This equation is known as the chain rule.

$$P(\mathbf{X}, \mathbf{e}) = \sum_{\mathbf{U} \setminus \mathbf{X}} P(\mathbf{U}, \mathbf{e}) = \sum_{\mathbf{U} \setminus \mathbf{X}} \prod_i P(U_i | \text{pa}(U_i)) \mathbf{e}$$

From the joint distribution over U we can in turn calculate any query we are interested in (with or without evidence set).

Suppose that there are two events which could cause grass to be wet: either the sprinkler is on or it's raining. Also, suppose that the rain has a direct effect on the use of the sprinkler (namely that when it

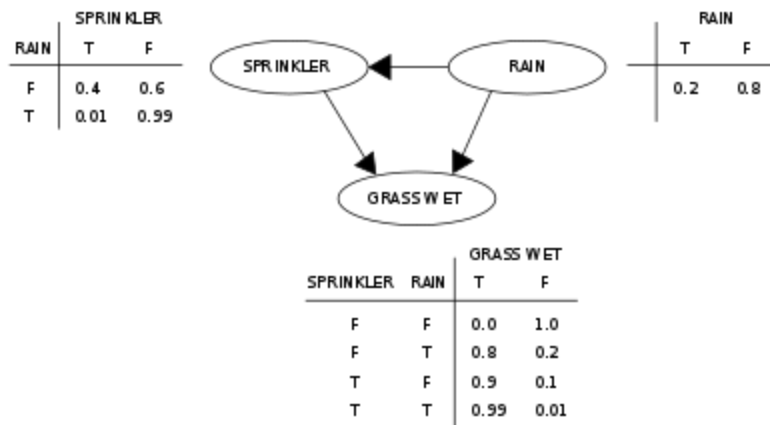
rains, the sprinkler is usually not turned on). Then the situation can be modeled with a Bayesian network (shown to the right). All three variables have two possible values, T (for true) and F (for false).

The joint probability function is:

$$\Pr(G, S, R) = \Pr(G|S, R) \Pr(S|R) \Pr(R)$$

The model can answer questions like "What is the probability that it is raining, given the grass is wet?"

by using the conditional probability formula and summing over all nuisance variables:



$$\Pr(R = T|G = T) = \frac{\Pr(G = T, R = T)}{\Pr(G = T)} = \frac{\sum_{S \in \{T, F\}} \Pr(G = T, S, R = T)}{\sum_{S, R \in \{T, F\}} \Pr(G = T, S, R)}$$

$$\begin{aligned} \Pr(G = T, S = T, R = T) &= \Pr(G = T|S = T, R = T) \Pr(S = T|R = T) \Pr(R = T) \\ &= 0.99 \times 0.01 \times 0.2 \\ &= 0.00198. \end{aligned}$$

BBn in python:

```
import bayespy as bp
import numpy as np
import csv
from colorama import init
from colorama import Fore, Back, Style
init()

# Define Parameter Enum values
```

```
#Age
ageEnum = {'SuperSeniorCitizen':0, 'SeniorCitizen':1, 'MiddleAged':2, 'Youth':3, 'Teen':4}
# Gender
genderEnum = {'Male':0, 'Female':1}
# FamilyHistory
familyHistoryEnum = {'Yes':0, 'No':1}
# Diet(Calorie Intake)
dietEnum = {'High':0, 'Medium':1, 'Low':2}
# LifeStyle
lifeStyleEnum = {'Athlete':0, 'Active':1, 'Moderate':2, 'Sedetary':3}
# Cholesterol
cholesterolEnum = {'High':0, 'BorderLine':1, 'Normal':2}
# HeartDisease
heartDiseaseEnum = {'Yes':0, 'No':1}
#heart_disease_data.csv
with open('heart_disease_data.csv') as csvfile:
    lines = csv.reader(csvfile)
    dataset = list(lines)
    data = []
    for x in dataset:

data.append([ageEnum[x[0]],genderEnum[x[1]],familyHistoryEnum[x[2]],dietEnum[x[3]],lifeStyle
eEnum[x[4]],cholesterolEnum[x[5]],heartDiseaseEnum[x[6]])
# Training data for machine learning todo: should import from csv
data = np.array(data)
N = len(data)

# Input data column assignment
p_age = bp.nodes.Dirichlet(1.0*np.ones(5))
age = bp.nodes.Categorical(p_age, plates=(N,))
age.observe(data[:,0])

p_gender = bp.nodes.Dirichlet(1.0*np.ones(2))
gender = bp.nodes.Categorical(p_gender, plates=(N,))
gender.observe(data[:,1])

p_familyhistory = bp.nodes.Dirichlet(1.0*np.ones(2))
familyhistory = bp.nodes.Categorical(p_familyhistory, plates=(N,))
familyhistory.observe(data[:,2])

p_diet = bp.nodes.Dirichlet(1.0*np.ones(3))
diet = bp.nodes.Categorical(p_diet, plates=(N,))
diet.observe(data[:,3])
```

```
p_lifestyle = bp.nodes.Dirichlet(1.0*np.ones(4))
lifestyle = bp.nodes.Categorical(p_lifestyle, plates=(N,))
lifestyle.observe(data[:,4])

p_cholesterol = bp.nodes.Dirichlet(1.0*np.ones(3))
cholesterol = bp.nodes.Categorical(p_cholesterol, plates=(N,))
cholesterol.observe(data[:,5])

# Prepare nodes and establish edges
# np.ones(2) -> HeartDisease has 2 options Yes/No
# plates(5, 2, 2, 3, 4, 3) -> corresponds to options present for domain values
p_heartdisease = bp.nodes.Dirichlet(np.ones(2), plates=(5, 2, 2, 3, 4, 3))
heartdisease = bp.nodes.MultiMixture([age, gender, familyhistory, diet, lifestyle, cholesterol],
bp.nodes.Categorical, p_heartdisease)
heartdisease.observe(data[:,6])
p_heartdisease.update()

# Sample Test with hardcoded values
#print("Sample Probability")
#print("Probability(HeartDisease|Age=SuperSeniorCitizen, Gender=Female, FamilyHistory=Yes,
DietIntake=Medium, LifeStyle=Setetary, Cholesterol=High)")
#print(bp.nodes.MultiMixture([ageEnum['SuperSeniorCitizen'], genderEnum['Female'],
familyHistoryEnum['Yes'], dietEnum['Medium'], lifeStyleEnum['Setetary'],
cholesterolEnum['High']], bp.nodes.Categorical,
p_heartdisease).get_moments()[0][heartDiseaseEnum['Yes']])

# Interactive Test
m = 0
while m == 0:
    print("\n")
    res = bp.nodes.MultiMixture([int(input('Enter Age: ' + str(ageEnum))), int(input('Enter Gender: '
+ str(genderEnum))), int(input('Enter FamilyHistory: ' + str(familyHistoryEnum))),
int(input('Enter dietEnum: ' + str(dietEnum))), int(input('Enter LifeStyle: ' + str(lifeStyleEnum))),
int(input('Enter Cholesterol: ' + str(cholesterolEnum))), bp.nodes.Categorical,
p_heartdisease).get_moments()[0][heartDiseaseEnum['Yes']]
    print("Probability(HeartDisease) = " + str(res))
    #print(Style.RESET_ALL)
    m = int(input("Enter for Continue:0, Exit :1 "))
```

Output:

```
Enter Age: {'SuperSeniorCitizen': 0, 'SeniorCitizen': 1, 'MiddleAged': 2, 'Youth': 3, 'Teen':
4} 1
Enter Gender: {'Male': 0, 'Female': 1} 1
```

Enter FamilyHistory: {'Yes': 0, 'No': 1}1
Enter dietEnum: {'High': 0, 'Medium': 1, 'Low': 2}2
Enter LifeStyle: {'Athlete': 0, 'Active': 1, 'Moderate': 2, 'Sedetary': 3}2
Enter Cholesterol: {'High': 0, 'BorderLine': 1, 'Normal': 2}1
Probability(HeartDisease) = 0.5
Enter for Continue:0, Exit :1 0

Enter Age: {'SuperSeniorCitizen': 0, 'SeniorCitizen': 1, 'MiddleAged': 2, 'Youth': 3, 'Teen': 4}0
Enter Gender: {'Male': 0, 'Female': 1}0
Enter FamilyHistory: {'Yes': 0, 'No': 1}0
Enter dietEnum: {'High': 0, 'Medium': 1, 'Low': 2}0
Enter LifeStyle: {'Athlete': 0, 'Active': 1, 'Moderate': 2, 'Sedetary': 3}3
Enter Cholesterol: {'High': 0, 'BorderLine': 1, 'Normal': 2}0
Probability(HeartDisease) = 0.5
Enter for Continue:0, Exit :1

Program 8

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.

Introduction to Expectation-Maximization (EM)

The EM algorithm tends to get stuck less than K-means algorithm. The idea is to assign data points partially to different clusters instead of assigning to only one cluster. To do this partial assignment, we model each cluster using a probabilistic distribution So a data point associates with a cluster with certain probability and it belongs to the cluster with the highest probability in the final assignment.

Expectation-Maximization (EM) algorithm

Step 1: An initial guess is made for the model's parameters and a probability distribution is created. This is sometimes called the "E-Step" for the "Expected" distribution.

Step 2: Newly observed data is fed into the model.

Step 3: The probability distribution from the E-step is drawn to include the new data. This is sometimes called the "M-step."

Step 4: Steps 2 through 4 are repeated until stability.

Data set:

1	1
3	3
5	5
8	8
2	2
11	11
14	14
18	18

Expectation-Maximization in Python:

```
import csv
import math
import copy
k=2
```

```
class cluster:
    def __init__(self,cluster_head_data):
        self.head=cluster_head_data
        self.l=[]
        self.mean=cluster_head_data[0]
        self.variance=1.0
    def display_head(self):
        print("m=",self.mean)
        print("v=",self.variance)
    def add_ele_cluster(self,data):
        self.l.append(data)
        print(self.l)
    def display_ele(self):
        print('list contains',self.l)

def find_insert_individual_cluster(cluster,element,n):
    ele=float(element[0])
    prob=[]

    for i in range(len(cluster)):
        pxb= 1/math.sqrt(2*3.142*cluster[i].variance)* math.exp(-1* ( ele - float(cluster[i].mean)
)**2 / (2*float(cluster[i].variance)))
        print('pxb exact==',pxb)
        prob.append(pxb)
    print('prob elem',prob)
    bi_den=0
    for i in range(len(prob)):
        bi_den=bi_den+prob[i]*1/n
        print('bi den', bi_den)
    #insert ele in to the cluster-- ele+pxb+bi
    for i in range(len(cluster)):
        clust_data=[]
        clust_data.append(ele)
        bi=(prob[i]*1/n)/bi_den
        clust_data.append(bi)
        #add the contents on to the cluster
        cluster[i].add_ele_cluster(clust_data)

def recalculate_cluster_mean_variance(cluster):
    l1=cluster.l
```

```
print('list enteries',l1)
#recalculating mean
mean_num=0.0
mean_den=0.0
var_num=0.0
var_den=0.0
for i in range(len(l1)):
    mean_num=mean_num+l1[i][0]*l1[i][1]
    mean_den=mean_den+l1[i][1]
mean=mean_num/mean_den
cluster.mean=mean
#recalculating varaiance
for i in range(len(l1)):
    var_num=var_num+l1[i][1]*(l1[i][0]-mean)**2
    var_den=var_den+l1[i][1]
variance=var_num/var_den
cluster.variance=mean

def find_nearest(cluster,ele):
    ele=float(ele[0])
    prob=[]
    nearest_prob=None
    index=1
    for i in range(len(cluster)):
        pxb= 1/math.sqrt(2*3.142*cluster[i].variance)* math.exp(-1* ( ele - float(cluster[i].mean)
)**2 / (2*float(cluster[i].variance)))
        print('pxb for cluster i,i','=',pxb)
        if nearest_prob is None:
            nearest_prob=pxb
            index=i
        else:
            if nearest_prob < pxb:
                nearest_prob=pxb
                index=i
    print('index',index,'nearest_prob','=',nearest_prob)
    cluster[index].l.append(ele)

#read the contents of CSV file
with open('cluster.csv') as csvfile:
    spamreader = csv.reader(csvfile, delimiter=',')
    #inserting ellements in to the list
```

```
db=[]
for row in spamreader:
    db.append(row)
#creating individual cluster heads
#displaying elements of the list
print('Db entries')
print(db)

#initialize the cluster
c=[]

#initialize the cluster heads
for i in range(k):
    new_clust=cluster(db[i])
    c.append(new_clust)
print('initial cluster Mean Variance')
#print cluster mean and variance
for i in range(k):
    print('----cluster',i,'-----')
    c[i].display_head()

error_ratio=1
# Iteration and including elements in the cluster
while error_ratio>0:
    error_ratio=0
    prevc=copy.deepcopy(c)
    #estimation
    for i in range(len(db)):
        find_insert_individual_cluster(c,db[i],len(db))
    #recalculate cluster mean and variance
    for i in range(len(c)):
        recalculate_cluster_mean_variance(c[i])
    #display recalculated mean and variance of cluster
    for i in range(k):
        print('----cluster',i,'-----')
        c[i].display_head()
    #clear all the values of the cluster list for the next iteration
    for i in range(k):
        c[i].l=[]
    #calculate the error
    error_ratio=0
```

```
    for i in range(len(c)):
        if abs(c[i].variance - prevc[i].variance)>0.1:
            error_ratio=error_ratio+1
#display the cluster elements
#clear all the values of the cluster list for the next iteration
for i in range(k):
    c[i].l=[]
#calculate the nearest prob for each element and include it in the resultant cluster
for i in range(len(db)):
    find_nearest(c,db[i])

#display cluster elements
for i in range(len(c)):
    print(c[i].l)
```

Output

```
----cluster 0 -----
m= 2.7213340517024402
v= 2.7213340517024402
----cluster 1 -----
m= 12.504957731742246
v= 12.504957731742246
pxb for cluster i 0 = 0.14030103544739295
pxb for cluster i 1 = 0.0005673529507970876
index 0 nearest_prob= 0.14030103544739295
pxb for cluster i 0 = 0.23839358908896524
pxb for cluster i 1 = 0.003044550243209151
index 0 nearest_prob= 0.23839358908896524
pxb for cluster i 0 = 0.09314788538736465
pxb for cluster i 1 = 0.01186516649339238
index 0 nearest_prob= 0.09314788538736465
pxb for cluster i 0 = 0.0014456877959066348
pxb for cluster i 1 = 0.05011024116755676
index 1 nearest_prob= 0.05011024116755676
pxb for cluster i 0 = 0.2197718847381247
pxb for cluster i 1 = 0.0013678961391086488
index 0 nearest_prob= 0.2197718847381247
pxb for cluster i 0 = 8.216307351891605e-07
pxb for cluster i 1 = 0.10304124381579008
index 1 nearest_prob= 0.10304124381579008
```

```
pxb for cluster i 0 = 1.709934543851912e-11
pxb for cluster i 1 = 0.10316387232255203
index 1 nearest_prob= 0.10316387232255203
pxb for cluster i 0 = 5.707857566318809e-20
pxb for cluster i 1 = 0.033728682873350235
index 1 nearest_prob= 0.033728682873350235
[1.0, 3.0, 5.0, 2.0]
[8.0, 11.0, 14.0, 18.0]
>>>
```

K-Means in Python:

```
import csv
import math
import copy
k=3

class cluster:
    def __init__(self,cluster_head_data):
        self.head=cluster_head_data
        self.l=[]
    def display_head(self):
        print(self.head)
    def add_ele_cluster(self,data):
        self.l.append(data)
    def display_ele(self):
        print('list contains',self.l)

def compare_the_values(first,second):
    x=float(first[0])
    y=float(first[1])
    x1=float(second[0])
    x2=float(second[1])
    val=math.sqrt( math.pow(math.fabs(x-x1),2)+math.pow(math.fabs(x-x1),2) )
    return val

def compare_the_nearest_cluster(cluster,data):
    #intialize the nearest cluster to 0
    dist_measure=None
    nearest=0
    for i in range(len(cluster)):
        dist=compare_the_values(cluster[i].head,data)
        if dist_measure is None:
```

```
        dist_measure=dist
        nearest=i
    if dist<dist_measure:
        dist_measure=dist
        nearest=i
    return nearest

def recal_head(cluster):
    for i in range(len(cluster)):
        l1=cluster[i].l
        xval=0.0
        yval=0.0
        for j in l1:
            xval=xval+float(j[0])
            yval=yval+float(j[1])
        xavg=xval/len(l1)
        yavg=yval/len(l1)
        avgl=[]
        avgl.append(xavg)
        avgl.append(yavg)
        cluster[i].head=avgl

#read the contents of CSV file
with open('cluster.csv') as csvfile:
    spamreader = csv.reader(csvfile, delimiter=',')
    #inserting elements in to the list
    db=[]
    for row in spamreader:
        db.append(row)
    #creating individual cluster heads
#displaying elements of the list
print('Db entries')
print(db)

#initialize the cluster
c=[]

#initialize the cluster heads
for i in range(k):
    new_clust=cluster(db[i])
```

```
c.append(new_clust)
print('initial cluster head values')
#print display heads
for i in range(k):
    print('----cluster',i,'-----')
    c[i].display_head()

error_ratio=1
# Iteration and including elements in the cluster
while error_ratio>0:
    prevc=copy.deepcopy(c)
    for ele in db:
        r=compare_the_nearest_cluster(c,ele)
        c[r].add_ele_cluster(ele)

    # display all the elements
    for clust in c:
        clust.display_ele()
    #recalculate the avg value
    recal_head(c)

    for i in range(k):
        print('----cluster',i,'-----')
        c[i].display_head()
    #remove the ele of cluster head for the next iter
    for i in range(k):
        c[i].l=[]
    #calculate the error
    error_ratio=0
    for i in range(k):
        if c[i].head != prevc[i].head:
            error_ratio=error_ratio+1
#final cluster ele
```

Output

```
Db entries
[['1', '1'], ['3', '3'], ['5', '5'], ['8', '8'], ['2', '2'], ['11', '11'], ['14', '14'], ['18', '18']]
initial cluster head values
----cluster 0 -----
['1', '1']
----cluster 1 -----
```


[3, 3]
----cluster 2 -----
[5, 5]
list contains [['1', '1'], [2, '2']]
list contains [['3', '3']]
list contains [['5', '5'], [8, '8'], [11, '11'], [14, '14'], [18, '18']]
----cluster 0 -----
[1.5, 1.5]
----cluster 1 -----
[3.0, 3.0]
----cluster 2 -----
[11.2, 11.2]
list contains [['1', '1'], [2, '2']]
list contains [['3', '3'], [5, '5']]
list contains [['8', '8'], [11, '11'], [14, '14'], [18, '18']]
----cluster 0 -----
[1.5, 1.5]
----cluster 1 -----
[4.0, 4.0]
----cluster 2 -----
[12.75, 12.75]
list contains [['1', '1'], [2, '2']]
list contains [['3', '3'], [5, '5'], [8, '8']]
list contains [['11', '11'], [14, '14'], [18, '18']]
----cluster 0 -----
[1.5, 1.5]
----cluster 1 -----
[5.333333333333333, 5.333333333333333]
----cluster 2 -----
[14.333333333333334, 14.333333333333334]
list contains [['1', '1'], [3, '3'], [2, '2']]
list contains [['5', '5'], [8, '8']]
list contains [['11', '11'], [14, '14'], [18, '18']]
----cluster 0 -----
[2.0, 2.0]
----cluster 1 -----
[6.5, 6.5]
----cluster 2 -----
[14.333333333333334, 14.333333333333334]
list contains [['1', '1'], [3, '3'], [2, '2']]
list contains [['5', '5'], [8, '8']]

```
list contains [['11', '11'], ['14', '14'], ['18', '18']]
----cluster 0 -----
[2.0, 2.0]
----cluster 1 -----
[6.5, 6.5]
----cluster 2 -----
[14.333333333333334, 14.333333333333334]
>>>
```

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Program 9

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.

TASK: The task of this program is to classify the IRIS data set examples by using the k-Nearest Neighbour algorithm. The new instance has to be classified based on its k nearest neighbors.

KNN falls in the supervised learning family of algorithms. Informally, this means that we are given a labeled dataset consisting of training observations (x,y) and would like to capture the relationship between x and y . More formally, our goal is to learn a function $h:X \rightarrow Y$ so that given an unseen observation x , $h(x)$ can confidently predict the corresponding output y .

The KNN classifier is also a non parametric and instance-based learning algorithm.

- Non-parametric means it makes no explicit assumptions about the functional form of h , avoiding the dangers of mismodeling the underlying distribution of the data. For example, suppose our data is highly non-Gaussian but the learning model we choose assumes a Gaussian form. In that case, our algorithm would make extremely poor predictions.
- Instance-based learning means that our algorithm doesn't explicitly learn a model. Instead, it chooses to memorize the training instances which are subsequently used as "knowledge" for the prediction phase. Concretely, this means that only when a query to our database is made (i.e. when we ask it to predict a label given an input), will the algorithm use the training instances to spit out an answer.

In the classification setting, the K-nearest neighbor algorithm essentially boils down to forming a majority vote between the K most similar instances to a given "unseen" observation. Similarity is defined according to a distance metric between two data points. A popular choice is the Euclidean distance given by

$$d(x, x') = \sqrt{(x_1 - x'_1)^2 + (x_2 - x'_2)^2 + \dots + (x_n - x'_n)^2}$$

but other measures can be more suitable for a given setting and include the Manhattan, Chebyshev and Hamming distance.

More formally, given a positive integer K , an unseen observation x and a similarity metric d , KNN classifier performs the following two steps:

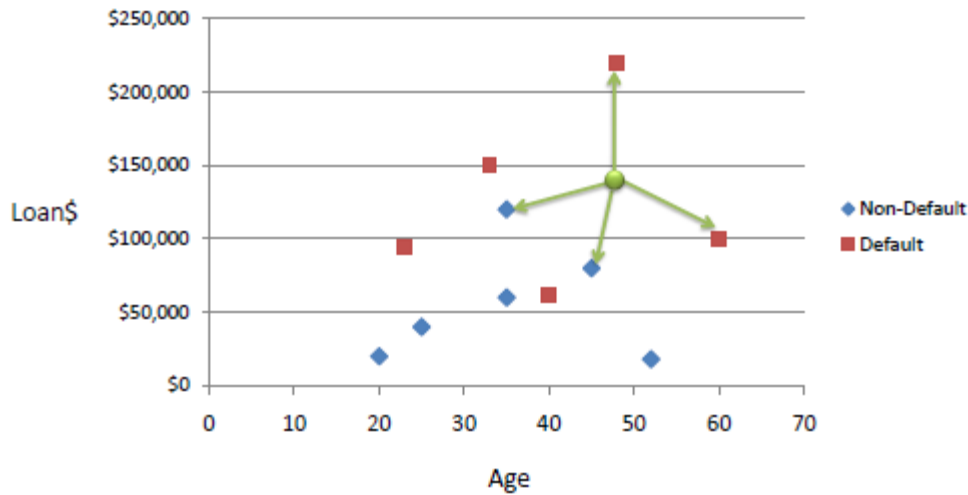
- It runs through the whole dataset computing d between x and each training observation. The K points in the training data that are closest to x are called the set A . Note that K is usually odd to prevent tie situations.
- It then estimates the conditional probability for each class, that is, the fraction of points in A with that given class label. (Note $I(x)$ is the indicator function which evaluates to 1 when the argument x is true and 0 otherwise)

$$P(y = j | X = x) = \frac{1}{K} \sum_{i \in A} I(y^{(i)} = j)$$

Finally, the input x gets assigned to the class with the largest probability.

EXAMPLE:

Consider the following data concerning credit default. Age and Loan are two numerical variables (predictors) and Default is the target.



We can now use the training set to classify an unknown case (Age=48 and Loan=\$142,000) using Euclidean distance. If K=1 then the nearest neighbor is the last case in the training set with Default=Y.

$$D = \text{Sqrt}[(48-33)^2 + (142000-150000)^2] = 8000.01 \gg \text{Default}=Y$$

Age	Loan	Default	Distance
25	\$40,000	N	102000
35	\$60,000	N	82000
45	\$80,000	N	62000
20	\$20,000	N	122000
35	\$120,000	N	22000
52	\$18,000	N	124000
23	\$95,000	Y	47000
40	\$62,000	Y	80000
60	\$100,000	Y	42000
48	\$220,000	Y	78000
33	\$150,000	Y	8000
48	\$142,000	?	

Euclidean Distance

$$D = \sqrt{(x_1 - y_1)^2 + (x_2 - y_2)^2}$$

With $K=3$, there are two Default=Y and one Default=N out of three closest neighbors. The prediction for the unknown case is again Default=Y.

ALGORITHM:

Let m be the number of training data samples. Let p be an unknown point.

1. Store the training samples in an array of data points $arr[]$. This means each element of this array represents a tuple (x, y) .
2. for $i=0$ to m :
 Calculate Euclidean distance $d(arr[i], p)$.
3. Make set S of K smallest distances obtained. Each of these distances correspond to an already classified data point.
4. Return the majority label among S .

KNN in Python:

```
import csv
import random
import math
import operator

def loadDataset(filename, split, trainingSet=[], testSet=[]):
    with open(filename) as csvfile:
        lines = csv.reader(csvfile)
        dataset = list(lines)
        for x in range(len(dataset)-1):
            for y in range(4):
                dataset[x][y] = float(dataset[x][y])
                if random.random() < split:
                    trainingSet.append(dataset[x])
                else:
                    testSet.append(dataset[x])

def euclideanDistance(instance1, instance2, length):
    distance = 0
    for x in range(length):
        distance += pow((instance1[x] - instance2[x]), 2)
    return math.sqrt(distance)

def getNeighbors(trainingSet, testInstance, k):
    distances = []
    length = len(testInstance)-1
    for x in range(len(trainingSet)):
        dist = euclideanDistance(testInstance, trainingSet[x], length)
        distances.append((trainingSet[x], dist))
    distances.sort(key=operator.itemgetter(1))
```

```
neighbors = []
for x in range(k):
    neighbors.append(distances[x][0])
return neighbors

def getResponse(neighbors):
    classVotes = { }
    for x in range(len(neighbors)):
        response = neighbors[x][-1]
        if response in classVotes:
            classVotes[response] += 1
        else:
            classVotes[response] = 1
    sortedVotes = sorted(classVotes.items(), key=operator.itemgetter(1), reverse=True)
    return sortedVotes[0][0]

def getAccuracy(testSet, predictions):
    correct = 0
    for x in range(len(testSet)):
        if testSet[x][-1] == predictions[x]:
            correct += 1
    return (correct/float(len(testSet))) * 100.0

def main():
    # prepare data
    trainingSet=[]
    testSet=[]
    split = 0.67
    loadDataset('iris_data.csv', split, trainingSet, testSet)
    print ('\n Number of Training data: ' + (repr(len(trainingSet))))
    print (' Number of Test Data: ' + (repr(len(testSet))))
    # generate predictions
    predictions=[]
    k = 3
    print("\n The predictions are: ")
    for x in range(len(testSet)):
        neighbors = getNeighbors(trainingSet, testSet[x], k)
        result = getResponse(neighbors)
        predictions.append(result)
        print(' predicted=' + repr(result) + ', actual=' + repr(testSet[x][-1]))
    accuracy = getAccuracy(testSet, predictions)
    print("\n The Accuracy is: ' + repr(accuracy) + '%')

main()
```


predicted='Iris-virginica', actual='Iris-virginica'
predicted='Iris-virginica', actual='Iris-virginica'
predicted='Iris-virginica', actual='Iris-virginica'
predicted='Iris-virginica', actual='Iris-virginica'
The Accuracy is: 100.0%

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Program 10

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

Nonparametric regression: is a category of regression analysis in which the predictor does not take a predetermined form but is constructed according to information derived from the data. Nonparametric regression requires larger sample sizes than regression based on parametric models because the data must supply the model structure as well as the model estimates.

Nonparametric regression is used for prediction and is reliable even if hypotheses of linear regression are not verified.

Locally weighted Learning also known as memory-based learning, instance-based learning, lazy-learning, and closely related to kernel density estimation, similarity searching and case-based reasoning.

LOWESS (Locally Weighted Scatterplot Smoothing), sometimes called LOESS (locally weighted smoothing), is a popular tool used in regression analysis that creates a smooth line through a timeplot or scatter plot to help you to see relationship between variables and foresee trends.

Locally weighted regression is a very powerful non-parametric model used in statistical learning.

Introduction :

Scatter-diagram smoothing (e.g. using the **lowess()** or **loess()** functions) involves drawing a smooth curve on a scatter diagram to summarize a relationship, in a fashion that makes few assumptions initially about the form or strength of the relationship. It is related to (and is a special case of) *nonparametric regression*, in which the objective is to represent the relationship between a response variable and one or more predictor variables, again in way that makes few assumptions about the form of the relationship. In other words, in contrast to “standard” linear regression analysis, no assumption is made that the relationship is represented by a straight line (although one could certainly think of a straight line as a special case of nonparametric regression).

If the basic decomposition-of-the-data model is:

$$data = \text{predictable component} + \text{noise},$$

then for the standard bivariate or multiple (linear) regression, the model is

$$data = \text{straight-line, polynomial or linearizable function} + \text{noise},$$

while for nonparametric regression, the model is

$$data = \text{smooth function determined by data} + \text{noise}.$$

Another way of looking at scatter diagram smoothing is as a way of depicting the “local” relationship between a response variable and a predictor variable over parts of their ranges, which may differ from a “global” relationship determined using the whole data set. Nonparametric regression can be thought of as generalizing the scatter plot smoothing idea to the multiple-regression context.

Locally Weighted Learning is a class of function approximation techniques, where a prediction is done by using an approximated local model around the current point of interest.

The **goal** of function approximation and regression is to find the underlying relationship between input and output. In a supervised learning problem training data, where each input is associated to one output, is used to create a model that predicts values which come close to the true function. All of these models use complete training data to derive global function.

Locally weighted regression

Local means using nearby points (i.e. a nearest neighbors approach)

Weighted means we value points based upon how far away they are.

Regression means approximating a function

This is **an instance-based learning method**

The idea: whenever you want to classify a sample:

- Build a local model of the function (using a linear function, quadratic, neural network, etc.)
- Use the model to predict the output value
- Throw the model away.

Locally Weighted Regression

Our final method combines advantages of parametric methods with non-parametric. The idea is to fit a regression model locally, weighting examples by the kernel K .

Locally Weighted Regression Algorithm

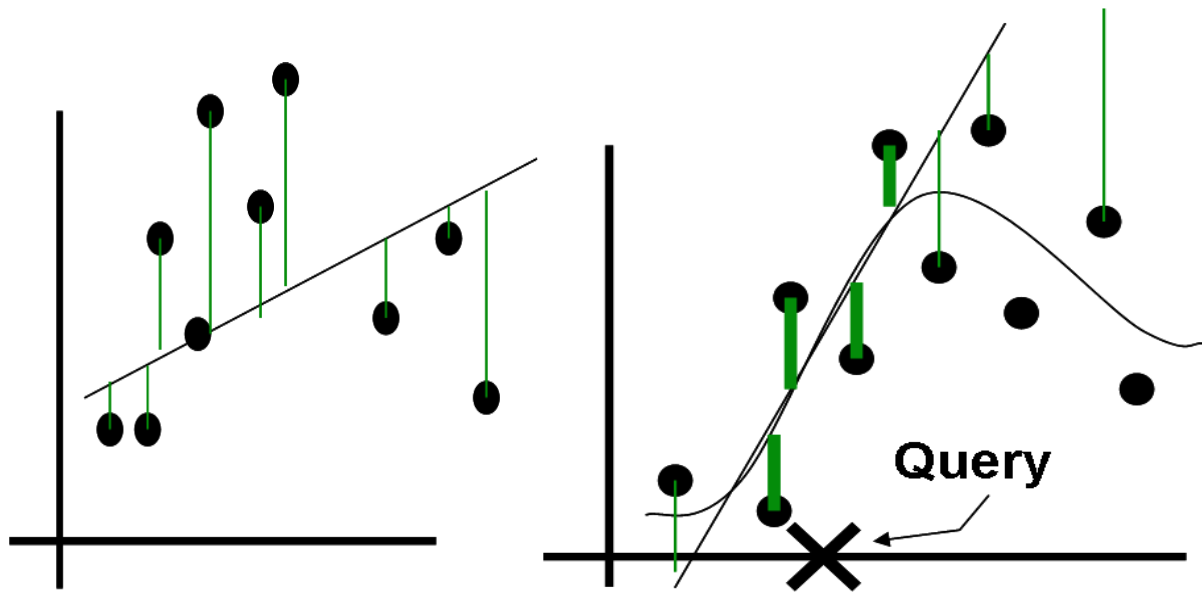
1. Given training data $D = \{\mathbf{x}_i, y_i\}$, Kernel function $K(\cdot, \cdot)$ and input \mathbf{x}
2. Fit weighted regression $\mathbf{w}_*(\mathbf{x}) = \operatorname{argmin}_{\mathbf{w}} \sum_{i=1}^n K(\mathbf{x}, \mathbf{x}_i) (\mathbf{w}^\top \mathbf{x}_i - y_i)^2$
3. Return regression prediction $\mathbf{w}_*(\mathbf{x})^\top \mathbf{x}$.

Note that we can do the same for classification, fitting a locally weighted logistic regression:

Locally Weighted Logistic Regression Algorithm

1. Given training data $D = \{\mathbf{x}_i, y_i\}$, Kernel function $K(\cdot, \cdot)$ and input \mathbf{x}
2. Fit weighted logistic regression $\mathbf{w}_*(\mathbf{x}) = \operatorname{argmin}_{\mathbf{w}} \sum_{i=1}^n K(\mathbf{x}, \mathbf{x}_i) \log(1 + \exp\{-y_i \mathbf{w}^\top \mathbf{x}_i\})$
3. Return logistic regression prediction $\operatorname{sign}(\mathbf{w}_*(\mathbf{x})^\top \mathbf{x})$.

The difference between regular linear regression and locally weighted linear regression can be visualized as follows:



x ---- is an instance,

D ---- is the set of possible instances

$a_i(x)$ ---- is the value of the i th attribute

w_i ----The weights form our hypothesis

f ----is the target function

f^{\wedge} ----is our approximation to the target function

In this case, we use a linear model to do the local approximation f^{\wedge} :

$$f^{\wedge}(x) = w_0 + w_1 a_1(x) + \dots + w_n a_n(x)$$

LOWESS is typically used for:

- Fitting a line to a scatter plot or time plot where noisy data values, sparse data points or weak interrelationships interfere with your ability to see a line of best fit.
- Linear regression where least squares fitting doesn't create a line of good fit or is too labor-intensive to use.
- Data exploration and analysis in the social sciences, particularly in elections and voting behavior.

Parametric and Non-Parametric Fitting

LOWESS, and least squares fitting in general, are **non-parametric** strategies for fitting a smooth curve to data points. "Parametric" means that the researcher or analyst assumes in advance that the data fits

some type of distribution (i.e. the normal distribution). Because some type of distribution is assumed in advance, parametric fitting can lead to fitting a smooth curve that misrepresents the data. In those cases, non-parametric smoothers may be a better choice. Non-parametric smoothers like LOESS try to find a curve of best fit without assuming the data must fit some distribution shape. In general, both types of smoothers are used for the same set of data to offset the advantages and disadvantages of each type of smoother.

Benefits of Non-Parametric Smoothing

- Provides a flexible approach to representing data.
- Ease of use.
- Computations are relatively easy.

Disadvantages of Non-Parametric Smoothing

- Can't be used to obtain a simple equation for a set of data.
- Less well understood than parametric smoothers.
- Requires the analyst to use a little guesswork to obtain a result.

Locally Weighted Regression in Python:

```
from math import ceil
import numpy as np
from scipy import linalg

def lowess(x, y, f=2./3., iter=3):
    n = len(x)
    r = int(ceil(f*n))
    h = [np.sort(np.abs(x - x[i]))[r] for i in range(n)]
    w = np.clip(np.abs((x[:,None] - x[None,:]) / h), 0.0, 1.0)
    w = (1 - w**3)**3
    yest = np.zeros(n)
    delta = np.ones(n)
    for iteration in range(iter):
        for i in range(n):
            weights = delta * w[:,i]
            b = np.array([np.sum(weights*y), np.sum(weights*y*x)])
            A = np.array([[np.sum(weights), np.sum(weights*x)],
                          [np.sum(weights*x), np.sum(weights*x*x)]])
            beta = linalg.solve(A, b)
            yest[i] = beta[0] + beta[1]*x[i]

        residuals = y - yest
        s = np.median(np.abs(residuals))
        delta = np.clip(residuals / (6.0 * s), -1, 1)
        delta = (1 - delta**2)**2

    return yest
```

```
if __name__ == '__main__':
    import math
    n = 100
    x = np.linspace(0, 2 * math.pi, n)
    print("=====values of x=====")
    print(x)
    y = np.sin(x) + 0.3*np.random.randn(n)
    print("=====Values of y=====")
    print(y)
    f = 0.25
    yest = lowess(x, y, f=f, iter=3)

    import pylab as pl
    pl.clf()
    pl.plot(x, y, label='y noisy')
    pl.plot(x, yest, label='y pred')
    pl.legend()
    pl.show()
```

Output

```
=====values of x=====
[0.    0.06346652 0.12693304 0.19039955 0.25386607 0.31733259
 0.38079911 0.44426563 0.50773215 0.57119866 0.63466518 0.6981317
 0.76159822 0.82506474 0.88853126 0.95199777 1.01546429 1.07893081
 1.14239733 1.20586385 1.26933037 1.33279688 1.3962634  1.45972992
 1.52319644 1.58666296 1.65012947 1.71359599 1.77706251 1.84052903
 1.90399555 1.96746207 2.03092858 2.0943951  2.15786162 2.22132814
 2.28479466 2.34826118 2.41172769 2.47519421 2.53866073 2.60212725
 2.66559377 2.72906028 2.7925268  2.85599332 2.91945984 2.98292636
 3.04639288 3.10985939 3.17332591 3.23679243 3.30025895 3.36372547
 3.42719199 3.4906585  3.55412502 3.61759154 3.68105806 3.74452458
 3.8079911  3.87145761 3.93492413 3.99839065 4.06185717 4.12532369
 4.1887902  4.25225672 4.31572324 4.37918976 4.44265628 4.5061228
 4.56958931 4.63305583 4.69652235 4.75998887 4.82345539 4.88692191
 4.95038842 5.01385494 5.07732146 5.14078798 5.2042545  5.26772102
 5.33118753 5.39465405 5.45812057 5.52158709 5.58505361 5.64852012
 5.71198664 5.77545316 5.83891968 5.9023862  5.96585272 6.02931923
 6.09278575 6.15625227 6.21971879 6.28318531]
=====Values of y=====
[-0.12131282 0.00704817 0.16084684 0.48415764 0.5790074  0.53758041
 0.41471799 0.31435684 0.42541349 0.56873662 0.64857746 0.56416606
 0.66906299 0.62626102 0.78385793 0.52007101 0.57661707 0.50246257
 0.59223656 1.19576243 0.9781296  0.76416587 0.85825567 0.68360755]
```

```
0.88965032 0.63976862 0.90159422 0.87464446 1.60256473 1.26886601
0.4200674 1.43656644 0.48378272 1.20891979 1.02955476 0.85766616
0.96519144 0.58303717 0.74996692 0.36357818 0.38481967 0.87027853
0.11686969 0.15356422 0.30403899 0.28014845 -0.15378752 0.56020242
0.20123909 -0.68626527 -0.19912738 -0.20357156 -0.1440503 -0.26493741
-0.32157543 -0.22139059 -0.42434559 -0.11447119 -0.67075934 -0.62973277
-1.18204429 -0.43386851 -1.12424791 -0.80806532 -0.85818104 -1.14149911
-0.23009694 -0.95399329 -0.62350893 -0.29176666 -0.55523556 -0.87957338
-1.78651723 -0.55784149 -1.1570492 -0.91445927 -0.99808369 -1.05836202
-0.55710891 -0.57815623 -0.60534368 -1.23588355 -1.34863514 -1.22421771
-1.01493553 -1.22922209 -0.48436674 -0.25055899 -0.61074483 -0.69197299
-0.52765904 -0.31451055 -0.55313683 -0.03399613 -0.53184918 0.33832978
-0.28385608 0.03227412 -0.38242201 0.1601836 ]
>>>
```

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