CSA4705-DEEP LEARNING FOR DATA HANDLING

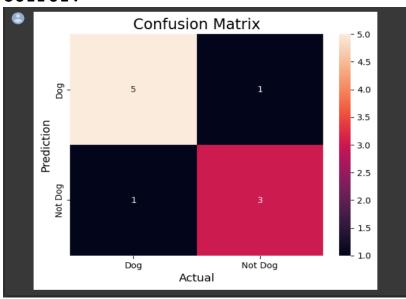
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EXPERIMENT:1(A)

<u>AIM:</u> To demonstrate confusion matrix using python

```
PROGRAM: #Import the necessary libraries
import numpy as np
from sklearn.metrics import confusion matrix
import seaborn as sns
import matplotlib.pyplot as plt
#Create the NumPy array for actual and predicted labels.
actual
         = np.array(
  ['Dog','Dog','Dog','Not Dog','Dog','Dog','Dog','Not Dog','N
ot Dog'])
predicted = np.array(
  ['Dog','Not Dog','Dog','Dog','Dog','Dog','Dog','Not Dog','N
ot Dog'])
#compute the confusion matrix.
cm = confusion matrix(actual, predicted)
#Plot the confusion matrix.
sns.heatmap(cm,
            annot=True,
            fmt='g',
           xticklabels=['Dog','Not Dog'],
            yticklabels=['Dog','Not Dog'])
plt.ylabel('Prediction', fontsize=13)
plt.xlabel('Actual', fontsize=13)
plt.title('Confusion Matrix', fontsize=17)
plt.show()
```

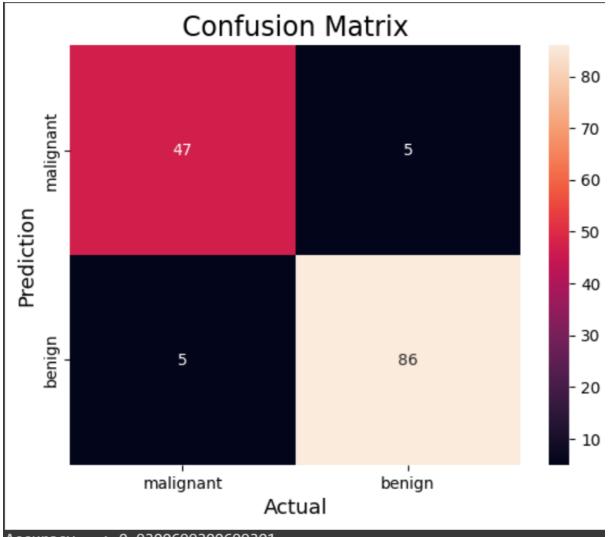


EXPERIMENT:1(B)

AIM: To demonstrate 2 class confusion matrix using python

```
#Import the necessary libraries
from sklearn.datasets import load breast cancer
from sklearn.model selection import train test split
from sklearn.tree import DecisionTreeClassifier
from sklearn.metrics import confusion matrix
import seaborn as sns
import matplotlib.pyplot as plt
from sklearn.metrics import accuracy score, precision score, recall sco
re, fl score
# Load the breast cancer dataset
X, y= load breast cancer(return X y=True)
X train, X test, y train, y test = train test split(X, y,test size=0.25
# Train the model
tree = DecisionTreeClassifier(random state=23)
tree.fit(X train, y train)
# preduction
y pred = tree.predict(X test)
# compute the confusion matrix
cm = confusion matrix(y test, y pred)
#Plot the confusion matrix.
sns.heatmap(cm,
            annot=True,
            fmt='q',
            xticklabels=['malignant', 'benign'],
            yticklabels=['malignant', 'benign'])
plt.ylabel('Prediction', fontsize=13)
plt.xlabel('Actual', fontsize=13)
plt.title('Confusion Matrix', fontsize=17)
plt.show()
# Finding precision and recall
accuracy = accuracy score(y test, y pred)
print("Accuracy :", accuracy)
```

```
precision = precision_score(y_test, y_pred)
print("Precision :", precision)
recall = recall_score(y_test, y_pred)
print("Recall :", recall)
F1_score = f1_score(y_test, y_pred)
print("F1-score :", F1 score)
```

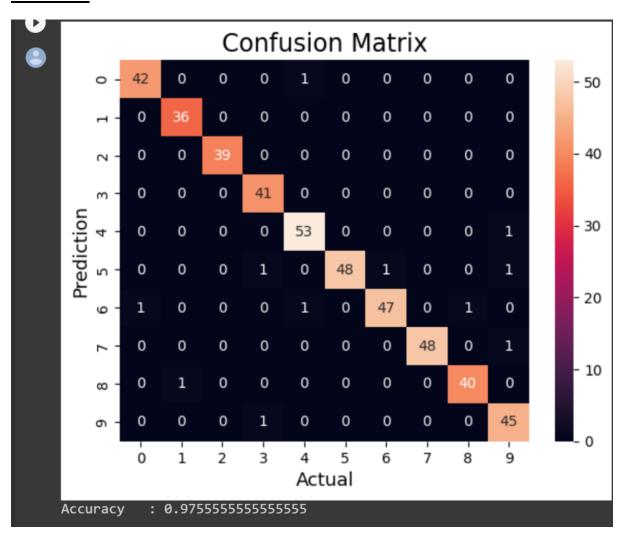


Accuracy : 0.9300699300699301 Precision : 0.945054945054945 Recall : 0.945054945054945 F1-score : 0.945054945054945

EXPERIMENT:2

<u>AIM:</u> Verifying the performance of a multi class confusion matrix by using choosen database with phython code

```
#Import the necessary libraries
from sklearn.datasets import load digits
from sklearn.model selection import train test split
from sklearn.ensemble import RandomForestClassifier
from sklearn.metrics import confusion matrix
import seaborn as sns
import matplotlib.pyplot as plt
from sklearn.metrics import accuracy score, precision score, recall sco
re, f1 score
# Load the breast cancer dataset
X, y= load digits(return X y=True)
X train, X test, y train, y test = train test split(X, y,test size=0.25
# Train the model
clf = RandomForestClassifier(random state=23)
clf.fit(X_train, y_train)
# preduction
y pred = clf.predict(X test)
# compute the confusion matrix
cm = confusion matrix(y test, y pred)
#Plot the confusion matrix.
sns.heatmap(cm,
            annot=True,
            fmt='g')
plt.ylabel('Prediction', fontsize=13)
plt.xlabel('Actual', fontsize=13)
plt.title('Confusion Matrix', fontsize=17)
plt.show()
# Finding precision and recall
accuracy = accuracy score(y test, y pred)
print("Accuracy :", accuracy)
```

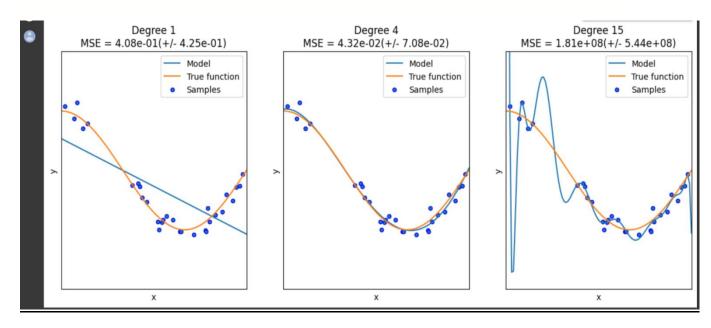


EXPERIMENT:3

<u>AIM:</u>: Verifying the performance of a over fitting by using choosen database with python code

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.pipeline import Pipeline
from sklearn.preprocessing import PolynomialFeatures
from sklearn.linear_model import LinearRegression
from sklearn.model_selection import cross_val_score
def true fun(X):
  return np.cos(1.5 * np.pi * X)
np.random.seed(0)
n_samples = 30
degrees = [1, 4, 15]
X = np.sort(np.random.rand(n_samples))
y = true_fun(X) + np.random.randn(n_samples) * 0.1
plt.figure(figsize=(14, 5))
for i in range(len(degrees)):
  ax = plt.subplot(1, len(degrees), i + 1)
  plt.setp(ax, xticks=(), yticks=())
  polynomial_features = PolynomialFeatures(degree=degrees[i], include_bias=
False)
  linear_regression = LinearRegression()
  pipeline = Pipeline(
    ("polynomial_features", polynomial_features),
       ("linear_regression", linear_regression),
  pipeline.fit(X[:, np.newaxis], y)
```

```
# Evaluate the models using crossvalidation
  scores = cross_val_score(
     pipeline, X[:, np.newaxis], y, scoring="neg_mean_squared_error", cv=10
  X_{\text{test}} = \text{np.linspace}(0, 1, 100)
  plt.plot(X_test, pipeline.predict(X_test[:, np.newaxis]), label="Model")
  plt.plot(X_test, true_fun(X_test), label="True function")
  plt.scatter(X, y, edgecolor="b", s=20, label="Samples")
  plt.xlabel("x")
  plt.ylabel("y")
  plt.xlim((0, 1))
  plt.ylim((-2, 2))
  plt.legend(loc="best")
  plt.title(
     "Degree { \text{NMSE} = \{:.2e\}(+/-\{:.2e\})$".format(
       degrees[i], -scores.mean(), scores.std()
     )
plt.show()
```



EXPERIMENT:4

<u>AIM:</u> To demonstrate the performance of a linear regression by using choosen database with python code

PROGRAM: LINEAR REGRESSION

```
import numpy as np
import matplotlib.pyplot as plt
def estimate_coef(x, y):
  # number of observations/points
  n = np.size(x)
  # mean of x and y vector
  m_x = np.mean(x)
  m_y = np.mean(y)
  # calculating cross-deviation and deviation about x
  SS_xy = np.sum(y*x) - n*m_y*m_x
  SS_x = np.sum(x*x) - n*m_x*m_x
  # calculating regression coefficients
  b_1 = SS_xy / SS_xx
  b_0 = m_y - b_1 * m_x
  return (b_0, b_1)
def plot_regression_line(x, y, b):
  # plotting the actual points as scatter plot
  plt.scatter(x, y, color = "r",
         marker = "o", s = 30)
  # predicted response vector
  y_pred = b[0] + b[1]*x
  # plotting the regression line
  plt.plot(x, y_pred, color = "b")
  # putting labels
  plt.xlabel('x')
  plt.ylabel('y')
```

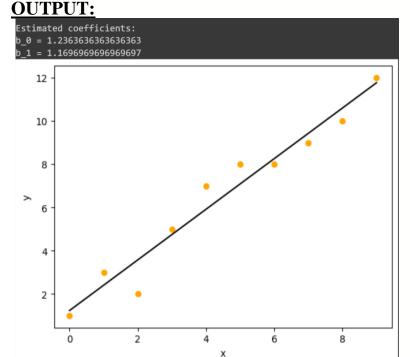
```
# function to show plot
plt.show()

def main():
    # observations / data
    x = np.array([0, 1, 2, 3, 4, 5, 6, 7, 8, 9])
    y = np.array([1, 3, 2, 5, 7, 8, 8, 9, 10, 12])

# estimating coefficients
b = estimate_coef(x, y)
print("Estimated coefficients:\nb_0 = {}
    \nb_1 = {}".format(b[0], b[1]))

# plotting regression line
plot_regression_line(x, y, b)

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



EXPERIMENT:5

<u>AIM:</u>: To demonstrate the performance of a logistic regression by using choosen database with python code.

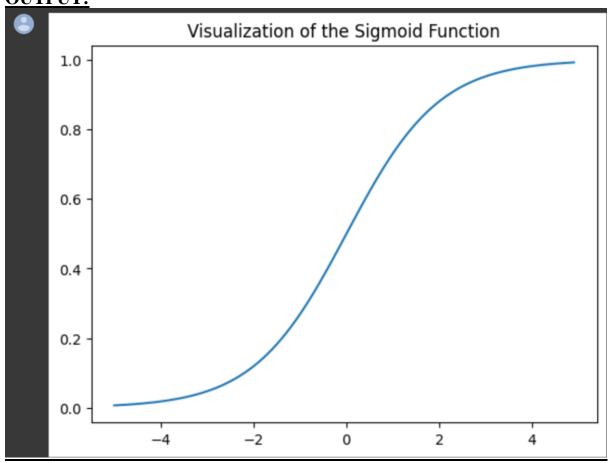
PROGRAM:

```
import numpy as np
import matplotlib.pyplot as plt

def sigmoid(z):
    return 1 / (1 + np.exp( - z))

plt.plot(np.arange(-5, 5, 0.1), sigmoid(np.arange(-5, 5, 0.1)))
plt.title('Visualization of the Sigmoid Function')

plt.show()
```



EXPERIMENT:6(a)KNN

AIM: Finding accuracy value of iris data set using KNN algorithm

```
import numpy as np
import pandas as pd
dataset = pd.read_csv("/content/IRIS.csv")
The breast cancer dataset has the following features: Sample code number, Clump
Thickness, Uniformity of Cell Size,
Uniformity of Cell Shape, Marginal Adhesion, Single Epithelial Cell Size, Bare Nuclei,
Bland Chromatin,
Normal Nucleoli, Mitosis, Class.
X = dataset.iloc[:, :-1].values
y = dataset.iloc[:, -1].values
dataset.shape
#splitting the dataset into the Training set and Test set
from sklearn.model selection import train test split
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size = 0.20, random_state
= 42)
#Feature Scaling
Feature scaling is the process of converting the data into a given range.
In this case, the standard scalar technique is used.
from sklearn.preprocessing import StandardScaler
sc = StandardScaler()
X train = sc.fit transform(X train)
X test = sc.transform(X test)
#Training the K-Nearest Neighbors (K-NN) Classification model on the Training set
Once the dataset is scaled, next, the K-Nearest Neighbors (K-NN) classifier algorithm
is used to create a model.
The hyperparameters such as n neighbors, metric, and p are set to 5, Minkowski,
and 2 respectively.
The remaining hyperparameters are set to default values.
from sklearn.neighbors import KNeighborsClassifier
classifier = KNeighborsClassifier(n neighbors = 5, metric = 'minkowski', p = 2)
classifier.fit(X_train, y_train)
Display the results (confusion matrix and accuracy)
```

Here evaluation metrics such as confusion matrix and accuracy are used to evaluate the performance of the model built using a decision tree classifier.

 $from \ sklearn.metrics \ import \ confusion_matrix, \ accuracy_score$ y_pred = classifier.predict(X_test) cm = confusion_matrix(y_test, y_pred) print(cm) accuracy_score(y_test, y_pred)

OUTPUT:



[[85 0] [2 50]] 0.9854014598540146

EXPERIMENT:6(B)NAVIE

<u>AIM: : finding accuracy value of iris data set using NAVIE BAYES algorithm</u>

PROGRAM:

```
import numpy as np
import pandas as pd
#Importing the dataset
```

targets into y.

Next, we import or read the dataset. Click here to download the breast cancer dataset used in this implementation.

After reading the dataset, divide the dataset into concepts and targets. Store the concepts into X and

```
"""
dataset = pd.read_csv("/content/IRIS.csv ")
X = dataset.iloc[:, :-1].values
y = dataset.iloc[:, -1].values
```

Splitting the dataset into the Training set and Test set

Once the dataset is read into the memory, next, divide the dataset into two parts, training and

testing using the train_test_split function from sklearn.

The test_size and random_state attributes are set to 0.25 and 0 respectively.

You can change these attributes as per your requirements.

```
from sklearn.model_selection import train_test_split

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size = 0.25, random_state = 0)
```

```
#Feature Scaling
```

Feature scaling is the process of converting the data into a min-max range. In this case, the standard scalar method is used.

from sklearn.preprocessing import StandardScaler
sc = StandardScaler()
X_train = sc.fit_transform(X_train)
X_test = sc.transform(X_test)

....

Training the Naive Bayes Classification model on the Training set

Once the dataset is scaled, next, the Naive Bayes classifier algorithm is used to create a model.

The GaussianNB function is imported from sklearn.naive_bayes library. The hyperparameters such as kernel,

and random_state to linear, and 0 respectively. The remaining hyperparameters of the support vector machine

algorithm are set to default values.

```
from sklearn.naive_bayes import GaussianNB classifier = GaussianNB()
```

```
classifier.fit(X_train, y_train)
#Naive Bayes classifier model
GaussianNB(priors=None, var_smoothing=1e-09)
#Display the results (confusion matrix and accuracy)
Here evaluation metrics such as confusion matrix and accuracy are used to evaluate the
performance of
the model built using a decision tree classifier.
from sklearn.metrics import confusion_matrix, accuracy_score
y_pred = classifier.predict(X_test)
cm = confusion_matrix(y_test, y_pred)
print(cm)
```



[[114 2] [2 53]] 0.9766081871345029

accuracy_score(y_test, y_pred)

EXPERIMENT:6(C)LOGISTIC

<u>AIM:</u>: finding accuracy value of iris data set using LOGISTIC REGRESSION algorithm

PROGRAM:

```
import numpy as np
import pandas as pd
#"Importing the dataset
After importing the necessary libraries, next, we import or read the dataset.
Click here to download the breast cancer dataset used in this implementation.
The breast cancer dataset has the following features:
Sample code number, Clump Thickness, Uniformity of Cell Size, Uniformity of Cell
Shape, Marginal Adhesion,
Single Epithelial Cell Size, Bare Nuclei, Bland Chromatin, Normal Nucleoli, Mitosis,
Class.
111111
# divide the dataset into concepts and targets. Store the concepts into X and targets
dataset = pd.read csv("/content/IRIS.csv ")
X = dataset.iloc[:, :-1].values
y = dataset.iloc[:, -1].values
#Splitting the dataset into the Training set and Test
from sklearn.model selection import train test split
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size = 0.30, random_state
= 2
#Feature Scaling
Feature scaling is the process of converting the data into a given range. In this case,
the standard scalar technique is used.
from sklearn.preprocessing import StandardScaler
from sklearn.preprocessing import StandardScaler
sc = StandardScaler()
X train = sc.fit transform(X train)
X test = sc.transform(X test)
```

Training the Logistic Regression (LR) Classification model on the Training set Once the dataset is scaled, next, the Logistic Regression (LR) classifier algorithm is used to create a model.

```
The hyperparameters such as random state to 0 respectively.
The remaining hyperparameters Logistic Regression (LR) are set to default values.
from sklearn.linear model import LogisticRegression
classifier = LogisticRegression(random state = 0)
classifier.fit(X train, y train)
#Logistic Regression (LR) classifier model
LogisticRegression(C=1.0, class weight=None, dual=False, fit intercept=True,
          intercept_scaling=1, l1_ratio=None, max_iter=100,
          multi class='warn', n jobs=None, penalty='l2',
          random state=0, solver='warn', tol=0.0001, verbose=0,
          warm start=False)
#Display the results (confusion matrix and accuracy)
Here evaluation metrics such as confusion matrix and accuracy are used to evaluate
the performance of the model
built using a decision tree classifier.
from sklearn.metrics import confusion_matrix, accuracy_score
y pred = classifier.predict(X test)
cm = confusion matrix(y test, y pred)
print(cm)
accuracy_score(y_test, y_pred)
```



[[117 8] [6 74]] 0.9317073170731708

EXPERIMENT:6(D)DECISION

AIM: : finding accuracy value of iris data set using DECISION TREE algorithm

PROGRAM:

```
import numpy as np
```

import pandas as pd

```
# Importing the dataset
```

dataset = pd.read_csv("/content/IRIS.csv ")

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

y = dataset.iloc[:, -1].values

Splitting the dataset into the Training set and Test set

from sklearn.model_selection import train_test_split

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size = 0.25,
random_state = 8)

Feature Scaling

from sklearn.preprocessing import StandardScaler

sc = StandardScaler()

X_train = sc.fit_transform(X_train)

X_test = sc.transform(X_test)

Training the Decision Tree Classification model on the Training set

from sklearn.tree import DecisionTreeClassifier

```
classifier = DecisionTreeClassifier(criterion = 'entropy', random state = 5)
classifier.fit(X_train, y_train)
# Display the Decision Tree
from sklearn.tree import plot_tree
import matplotlib.pyplot as plt
plt.figure(figsize=(20,10))
plot tree(classifier, filled=True, rounded=True,
feature_names=dataset.columns[:-1])
plt.show()
# Predicting the Test set results
y_pred = classifier.predict(X_test)
# Display the results (confusion matrix and accuracy)
from sklearn.metrics import confusion_matrix, accuracy_score
cm = confusion_matrix(y_test, y_pred)
print(cm)
accuracy_score(y_test, y_pred)
output:
```

```
petal_width <= -0.533
                             \overline{\text{entropy}} = 1.585
                          samples = 112
value = [37, 38, 37]
                                                      petal_width <= 0.71
entropy = 0.0
samples = 37
value = [37, 0, 0]
                                                          \bar{\text{entropy}} = 1.0
                                                          samples = 75
                                                       value = [0, 38, 37]
                        petal length <= 0.732
                                                                                    entropy = 0.0
samples = 36
value = [0, 0, 36]
                             entropy = 0.172
samples = 39
                            value = [0, 38, 1]
                                                    sepal_width <= -0.658
  entropy = 0.0
samples = 37
                                                          entropy = 1.0
                                                           samples = 2
value = [0, 37, 0]
                                                         value = [0, 1, 1]
                              entropy = 0.0
samples = 1
                                                                                      entropy = 0.0
samples = 1
                             value = [0, 1, 0]
                                                                                     value = [0, 0, 1]
```

```
0.8947368421052632
```

[[13 0 0] [0 11 1]

[0 3 10]]

EXPERIMENT:6(E)SVM

AIM: : finding accuracy value of iris data set using SVM algorithm

PROGRAM:

```
import numpy as np
```

import pandas as pd

Importing the dataset

dataset = pd.read_csv("/content/IRIS.csv ")

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

y = dataset.iloc[:, -1].values

Splitting the dataset into the Training set and Test set

from sklearn.model_selection import train_test_split

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.25, random_state=32)

Feature Scaling

from sklearn.preprocessing import StandardScaler

sc = StandardScaler()

X_train = sc.fit_transform(X_train)

X test = sc.transform(X test)

Training the SVM model on the Training set

```
from sklearn.svm import SVC
classifier = SVC(kernel='linear', random_state=0)
classifier.fit(X_train, y_train)
# Predicting the Test set results
y_pred = classifier.predict(X_test)
# Evaluating the performance of the model using confusion matrix and
accuracy
from sklearn.metrics import confusion_matrix, accuracy_score
cm = confusion_matrix(y_test, y_pred)
print(cm)
print('Accuracy: {:.2f}%'.format(accuracy_score(y_test, y_pred) * 100))
OUTPUT:
```



[[108 1] 5 57]] Accuracy: 96.49%

EXPERIMENT:6(F)RANDOM

AIM: : finding accuracy value of iris data set using RANDOM FOREST algorithm

PROGRAM:

```
import numpy as np
```

import pandas as pd

```
# Importing the dataset
```

dataset = pd.read_csv("/content/IRIS.csv ")

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

y = dataset.iloc[:, -1].values

Splitting the dataset into the Training set and Test set

from sklearn.model_selection import train_test_split

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.25, random_state=39)

Feature Scaling

from sklearn.preprocessing import StandardScaler

sc = StandardScaler()

X_train = sc.fit_transform(X_train)

X_test = sc.transform(X_test)

Training the Random Forest Classification model on the Training set

from sklearn.ensemble import RandomForestClassifier

```
classifier = RandomForestClassifier(n_estimators=100, random_state=42)
classifier.fit(X_train, y_train)

# Predicting the Test set results

y_pred = classifier.predict(X_test)

# Making the Confusion Matrix

from sklearn.metrics import confusion_matrix, accuracy_score

cm = confusion_matrix(y_test, y_pred)

print(cm)

print('Accuracy:', accuracy_score(y_test, y_pred))
CUTDLE:
```



[[111 1] [2 57]] Accuracy: 0.9824561403508771

EXPERIMENT:7(A)

<u>AIM:</u> To demonstrate gradient descent using python(actual data)

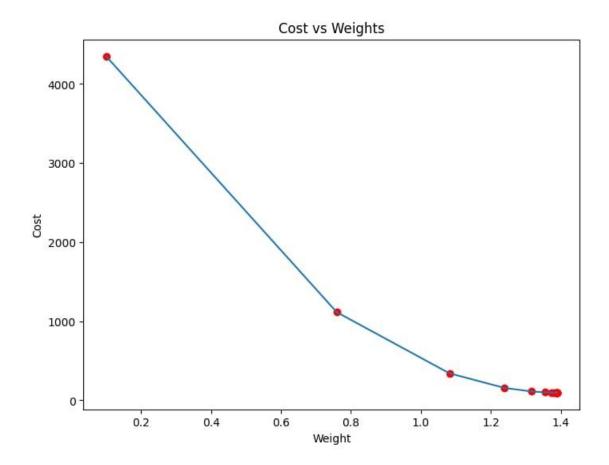
```
# Importing Libraries
import numpy as np
import matplotlib.pyplot as plt
def mean squared error (y true, y predicted):
    # Calculating the loss or cost
    cost = np.sum((y true-y predicted)**2) / len(y true)
   return cost
# Gradient Descent Function
# Here iterations, learning rate, stopping threshold
# are hyperparameters that can be tuned
def gradient descent(x, y, iterations = 1000, learning rate = 0.0001,
                     stopping threshold = 1e-6):
    # Initializing weight, bias, learning rate and iterations
    current weight = 0.1
    current bias = 0.01
    iterations = iterations
    learning rate = learning rate
    n = float(len(x))
    costs = []
    weights = []
    previous cost = None
    # Estimation of optimal parameters
    for i in range(iterations):
        # Making predictions
        y_predicted = (current_weight * x) + current_bias
        # Calculating the current cost
        current_cost = mean_squared_error(y, y_predicted)
        # If the change in cost is less than or equal to
        # stopping threshold we stop the gradient descent
        if previous cost and abs(previous cost-
current cost) <= stopping threshold:</pre>
           break
        previous cost = current cost
        costs.append(current cost)
        weights.append(current weight)
        # Calculating the gradients
        weight derivative = -(2/n) * sum(x * (y-y predicted))
        bias derivative = -(2/n) * sum(y-y predicted)
```

```
# Updating weights and bias
        current weight = current weight - (learning rate *
weight derivative)
       current bias = current bias - (learning rate * bias derivative)
        # Printing the parameters for each 1000th iteration
        print(f"Iteration {i+1}: Cost {current cost}, Weight \
        {current weight}, Bias {current bias}")
    # Visualizing the weights and cost at for all iterations
    plt.figure(figsize = (8,6))
    plt.plot(weights, costs)
    plt.scatter(weights, costs, marker='o', color='red')
   plt.title("Cost vs Weights")
   plt.ylabel("Cost")
   plt.xlabel("Weight")
   plt.show()
    return current weight, current bias
def main():
    # Data
   X = np.array([32.50234527, 53.42680403, 61.53035803, 47.47563963,
59.81320787,
           55.14218841, 52.21179669, 39.29956669, 48.10504169,
52.55001444,
           45.41973014, 54.35163488, 44.1640495 , 58.16847072,
56.72720806,
           48.95588857, 44.68719623, 60.29732685, 45.61864377,
38.81681754])
   Y = np.array([31.70700585, 68.77759598, 62.5623823, 71.54663223,
87.23092513,
           78.21151827, 79.64197305, 59.17148932, 75.3312423 ,
71.30087989,
           55.16567715, 82.47884676, 62.00892325, 75.39287043,
81.43619216,
           60.72360244, 82.89250373, 97.37989686, 48.84715332,
56.87721319])
    # Estimating weight and bias using gradient descent
    estimated weight, estimated bias = gradient descent(X, Y,
iterations=2000)
    print(f"Estimated Weight: {estimated weight} \nEstimated Bias:
{estimated bias}")
    # Making predictions using estimated parameters
    Y pred = estimated weight*X + estimated bias
    # Plotting the regression line
    plt.figure(figsize = (8,6))
   plt.scatter(X, Y, marker='o', color='red')
   plt.plot([min(X), max(X)], [min(Y pred), max(Y pred)],
color='blue', markerfacecolor='red',
            markersize=10,linestyle='dashed')
```

```
plt.xlabel("X")
plt.ylabel("Y")
plt.show()

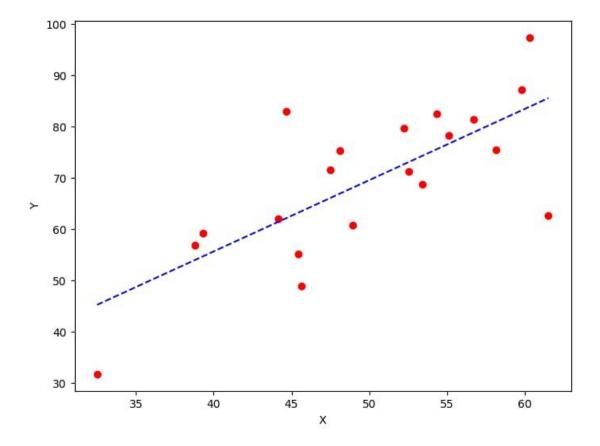
if __name__=="__main__":
    main()
```

output:



Estimated Weight: 1.393097090459544 Estimated Bias: 0.035349609417819915

100 -



Experiment:7(b)

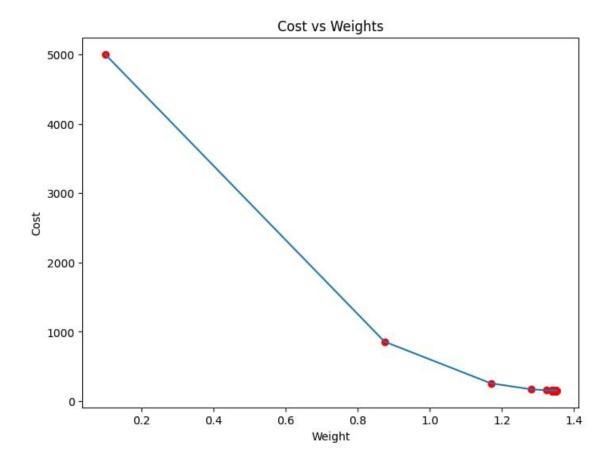
<u>AIM:</u> To demonstrate gradient descent using python(modified data)

```
# Importing Libraries
import numpy as np
import matplotlib.pyplot as plt
def mean squared error (y true, y predicted):
    # Calculating the loss or cost
    cost = np.sum((y true-y predicted)**2) / len(y true)
   return cost
# Gradient Descent Function
# Here iterations, learning rate, stopping threshold
# are hyperparameters that can be tuned
def gradient_descent(x, y, iterations = 1000, learning rate = 0.0001,
                     stopping threshold = 1e-6):
    # Initializing weight, bias, learning rate and iterations
    current weight = 0.1
    current bias = 0.01
    iterations = iterations
    learning rate = learning rate
    n = float(len(x))
    costs = []
    weights = []
    previous cost = None
    # Estimation of optimal parameters
    for i in range(iterations):
        # Making predictions
        y_predicted = (current_weight * x) + current_bias
        # Calculating the current cost
        current_cost = mean_squared_error(y, y_predicted)
        # If the change in cost is less than or equal to
        # stopping threshold we stop the gradient descent
        if previous cost and abs(previous cost-
current cost) <= stopping threshold:</pre>
           break
        previous cost = current cost
        costs.append(current cost)
        weights.append(current weight)
        # Calculating the gradients
        weight derivative = -(2/n) * sum(x * (y-y predicted))
        bias derivative = -(2/n) * sum(y-y predicted)
```

```
# Updating weights and bias
        current weight = current weight - (learning rate *
weight derivative)
       current bias = current bias - (learning rate * bias derivative)
        # Printing the parameters for each 1000th iteration
        print(f"Iteration {i+1}: Cost {current cost}, Weight \
        {current weight}, Bias {current bias}")
    # Visualizing the weights and cost at for all iterations
    plt.figure(figsize = (8,6))
    plt.plot(weights, costs)
    plt.scatter(weights, costs, marker='o', color='red')
    plt.title("Cost vs Weights")
   plt.ylabel("Cost")
   plt.xlabel("Weight")
   plt.show()
    return current weight, current bias
def main():
    # Data
   X = np.array([52.50234527, 63.42680403, 81.53035803, 47.47563963,
89.81320787,
           55.14218841, 52.21179669, 39.29956669, 48.10504169,
52.55001444,
           45.41973014, 54.35163488, 44.1640495 , 58.16847072,
56.72720806,
           48.95588857, 44.68719623, 60.29732685, 45.61864377,
38.81681754])
   Y = np.array([41.70700585, 78.77759598, 82.5623823, 91.54663223,
77.23092513,
           78.21151827, 79.64197305, 59.17148932, 75.3312423 ,
71.30087989,
           55.16567715, 82.47884676, 62.00892325, 75.39287043,
81.43619216,
           60.72360244, 82.89250373, 97.37989686, 48.84715332,
56.87721319])
    # Estimating weight and bias using gradient descent
    estimated weight, estimated bias = gradient descent(X, Y,
iterations=2000)
    print(f"Estimated Weight: {estimated weight} \nEstimated Bias:
{estimated bias}")
    # Making predictions using estimated parameters
    Y pred = estimated weight*X + estimated bias
    # Plotting the regression line
    plt.figure(figsize = (8,6))
   plt.scatter(X, Y, marker='orange', color='pink')
   plt.plot([min(X), max(X)], [min(Y pred), max(Y pred)],
color='blue', markerfacecolor='red',
            markersize=10,linestyle='dashed')
```

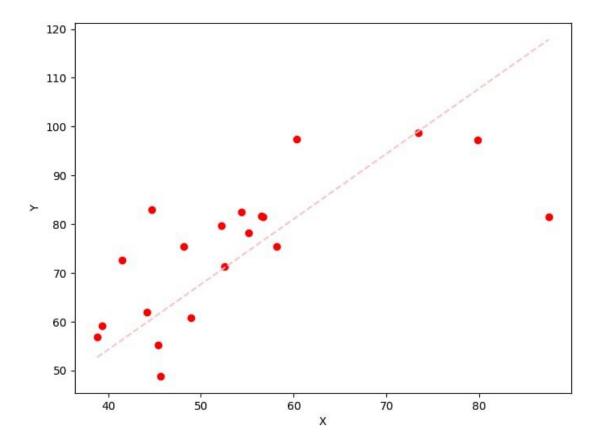
```
plt.xlabel("X")
plt.ylabel("Y")
plt.show()

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



Estimated Weight: 1.393097090459544 Estimated Bias: 0.035349609417819915

100 -



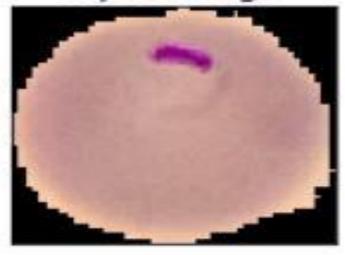
EXPERIMENT:8(A)SEGMENTATION

<u>AIM:</u>: Verifying the performance of a image processing by using choosen database with phython code

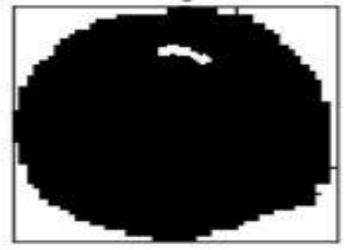
PROGRAM:

```
# SEGMENTATION
import numpy as np
import cv2
from matplotlib import pyplot as plt
img = cv2.imread(r'C33P1thinF IMG 20150619 114756a cell 181.png')
b,g,r = cv2.split(img)
rgb img = cv2.merge([r,g,b])
gray = cv2.cvtColor(img,cv2.COLOR BGR2GRAY)
ret, thresh =
cv2.threshold(gray,0,255,cv2.THRESH BINARY INV+cv2.THRESH OTSU)
# noise removal
kernel = np.ones((2,2),np.uint8)
#opening = cv2.morphologyEx(thresh,cv2.MORPH OPEN,kernel, iterations = 2)
closing = cv2.morphologyEx(thresh,cv2.MORPH CLOSE,kernel, iterations = 2)
# sure background area
sure_bg = cv2.dilate(closing, kernel, iterations=3)
# Finding sure foreground area
dist transform = cv2.distanceTransform(sure bg,cv2.DIST L2,3)
# Threshold
ret, sure fg = cv2.threshold(dist transform, 0.1*dist transform.max(), 255,0)
# Finding unknown region
sure fg = np.uint8(sure fg)
unknown = cv2.subtract(sure bg, sure fg)
# Marker labelling
ret, markers = cv2.connectedComponents(sure fg)
# Add one to all labels so that sure background is not 0, but 1
markers = markers+1
# Now, mark the region of unknown with zero
markers[unknown==255] = 0
markers = cv2.watershed(img,markers)
img[markers == -1] = [255, 0, 0]
plt.subplot(211),plt.imshow(rgb img)
plt.title('Input Image'), plt.xticks([]), plt.yticks([])
plt.subplot(212),plt.imshow(thresh, 'gray')
plt.imsave(r'thresh.png',thresh)
plt.title("Otsu's binary threshold"), plt.xticks([]), plt.yticks([])
plt.tight_layout()
plt.show()
```

Input Image



Otsu's binary threshold

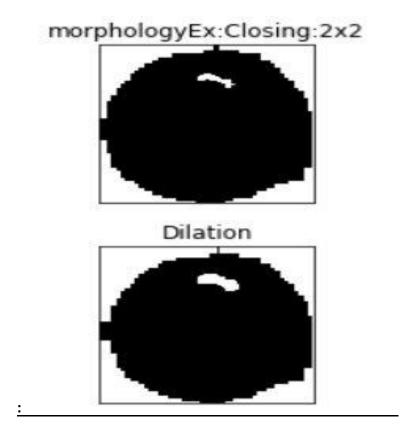


EXPERIMENT:8(B)

<u>AIM:</u>: Verifying the performance of a image processing by using water shed database with python code

PROGRAM:

```
# SEGMENTATION
import numpy as np
import cv2
from matplotlib import pyplot as plt
imq =
cv2.imread(r'C33P1thinF IMG 20150619 114756a cell 181.png')
b,g,r = cv2.split(img)
rgb img = cv2.merge([r,g,b])
gray = cv2.cvtColor(img,cv2.COLOR BGR2GRAY)
ret, thresh =
cv2.threshold(gray,0,255,cv2.THRESH BINARY INV+cv2.THRESH OTSU
plt.subplot(211),plt.imshow(closing, 'gray')
plt.title("morphologyEx:Closing:2x2"), plt.xticks([]),
plt.yticks([])
plt.subplot(212),plt.imshow(sure bg, 'gray')
plt.imsave(r'dilation.png', sure bg)
plt.title("Dilation"), plt.xticks([]), plt.yticks([])
plt.tight layout()
plt.show()
```



EXPERIMENT:9 (a) TANH

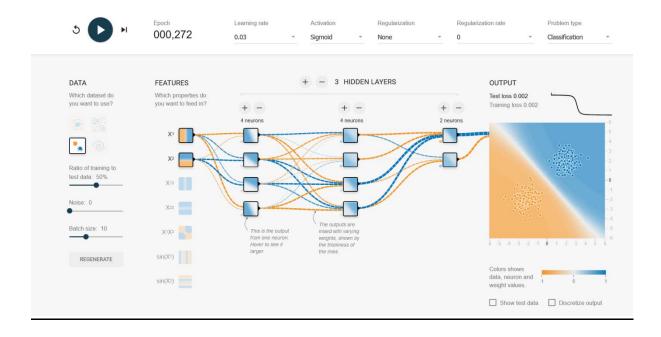
AIM: Neural network analysis using TANH activation

OUTPUT:



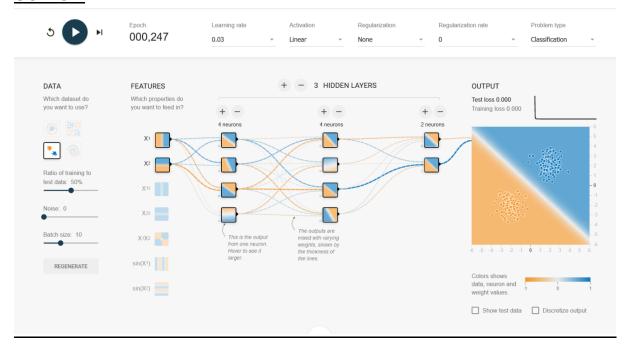
EXPERIMENT:9(B) SIGMIOD

AIM: Neural network analysis using SIGMOID activation



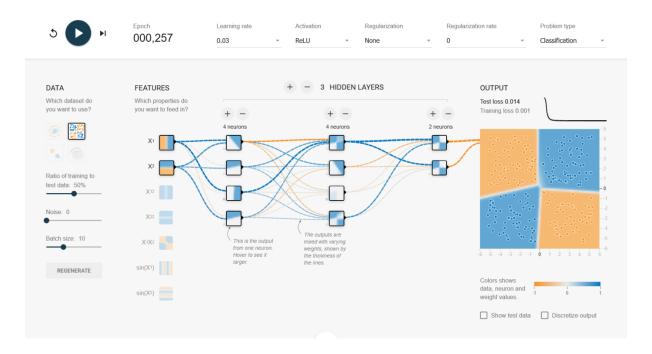
EXPERIMENT:9(C) LINEAR

AIM: Neural network analysis using LINEAR activation



EXPERIMENT:9(D)RELU

<u>AIM:</u> Neural network analysis using ReLU activation



EXPERIMENT:10

AIM: To demonstrate linear separability using python code

```
import numpy as np
import matplotlib.pyplot as plt
def estimate_coef(x, y):
  # number of observations/points
  n = np.size(x)
  # mean of x and y vector
  m_x = np.mean(x)
  m_y = np.mean(y)
  # calculating cross-deviation and deviation about x
  SS_xy = np.sum(y*x) - n*m_y*m_x
  SS_x = np.sum(x*x) - n*m_x*m_x
  # calculating regression coefficients
  b_1 = SS_xy / SS_xx
  b_0 = m_y - b_1 * m_x
  return (b_0, b_1)
def plot_regression_line(x, y, b):
  # plotting the actual points as scatter plot
  plt.scatter(x, y, color = "r",
         marker = "o", s = 30)
  # predicted response vector
  y_pred = b[0] + b[1]*x
  # plotting the regression line
  plt.plot(x, y_pred, color = "b")
  # putting labels
  plt.xlabel('x')
  plt.ylabel('y')
  # function to show plot
```

```
plt.show()
def main():
  # observations / data
  x = np.array([0, 1, 2, 3, 4, 5, 6, 7, 8, 9])
  y = np.array([1, 3, 2, 5, 7, 8, 8, 9, 10, 12])
  # estimating coefficients
  b = estimate\_coef(x, y)
  print("Estimated coefficients:\nb_0 = {} \
      nb_1 = {}".format(b[0], b[1]))
  # plotting regression line
  plot_regression_line(x, y, b)
if __name__ == "__main__":
  main()
```

