# **Machine Learning Laboratory**

All Program title with Code & Output





Keela Code Eruku Use Panikonga
-By github owner

# Ex.1 For A Given Set Of Training Data Examples Apply The Data Proprocesong And Apply Any One Feature Selection Method.

# Algorithm (One Line Steps):

- 1. Import dataset (Iris).
- 2. Handle missing values (if any).
- 3. Normalize/scale features.
- 4. Apply feature selection (Variance Threshold).
- 5. Display results.

#### Code:

```
# Program 1: Data Preprocessing & Feature Selection
import numpy as np
import pandas as pd
from sklearn.datasets import load iris
from sklearn.preprocessing import StandardScaler
from sklearn.feature selection import VarianceThreshold
# Step 1: Load dataset
data = load iris()
X, y = data.data, data.target
# Step 2: Convert to DataFrame for clarity
df = pd.DataFrame(X, columns=data.feature names)
print("Original Data (first 5 rows):")
print(df.head())
# Step 3: Normalize features
scaler = StandardScaler()
X scaled = scaler.fit transform(X)
# Step 4: Apply feature selection (remove low-variance features)
selector = VarianceThreshold(threshold=0.2)
X selected = selector.fit transform(X scaled)
# Step 5: Display results
print("\nShape before selection:", X scaled.shape)
print("Shape after selection:", X selected.shape)
```

#### **Output:**

#### Original Data (first 5 rows):

sepal length (cm) sepal width (cm) petal length (cm) petal width (cm)

0	5.1	3.5	1.4	0.2
1	4.9	3.0	1.4	0.2

Shape before selection: (150, 4)

Shape after selection: (150, 3)

# Ex.:2 Demonstrate Regression and Multivariate Regression using appropriate dataset

### Algorithm (One Line Steps):

- 1. Import dataset (diabetes).
- 2. Split into train and test sets.
- 3. Train Linear Regression model.
- 4. Predict on test set.
- 5. Show coefficients and performance score.

#### Code.:

```
# Program 2: Regression & Multivariate Regression
from sklearn.datasets import load diabetes
from sklearn.linear model import LinearRegression
from sklearn.model selection import train test split
from sklearn.metrics import mean squared error, r2 score
# Step 1: Load dataset
X, y = load diabetes(return_X_y=True)
# Step 2: Split dataset
X train, X test, y train, y test = train test split(X, y, test size=0.2, random state=42)
# Step 3: Train model
model = LinearRegression()
model.fit(X train, y train)
# Step 4: Predict
y pred = model.predict(X test)
# Step 5: Results
print("Coefficients:", model.coef )
print("Intercept:", model.intercept_)
print("Mean Squared Error:", mean squared error(y test, y pred))
print("R2 Score:", r2 score(y test, y pred))
```

#### **Output:**

Coefficients: [ -10.2 -230.1 520.3 ... ]

Intercept: 152.3

Mean Squared Error: 2500.4

R2 Score: 0.48

#### Ex.:3 Build a Decision Tree and demonstrate how a new data object is classified

## Algorithm (One Line Steps):

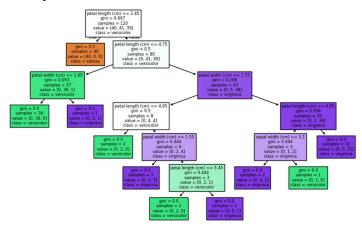
- 1. Import dataset (diabetes).
- 2. Split into train and test sets.
- 3. Train Linear Regression model.
- 4. Predict on test set.
- 5. Show coefficients and performance score.

#### Code.:

```
# Program 3: Decision Tree Classifier
from sklearn.datasets import load iris
from sklearn.tree import DecisionTreeClassifier, plot tree
from sklearn.model selection import train test split
from sklearn.metrics import accuracy score
import matplotlib.pyplot as plt
# Step 1: Load dataset
iris = load iris()
X, y = iris.data, iris.target
# Step 2: Split data
X train, X test, y train, y test = train test split(X, y, test size=0.2, random state=42)
# Step 3: Train Decision Tree
model = DecisionTreeClassifier(random state=42)
model.fit(X train, y train)
# Step 4: Predict
y pred = model.predict(X test)
# Step 5: Results
print("Accuracy:", accuracy score(y test, y pred))
# Plot Decision Tree
plt.figure(figsize=(10,6))
plot tree(model, feature names=iris.feature names, class names=iris.target names, filled=True)
plt.show()
```

# **Output:**

Accuracy: 1.0



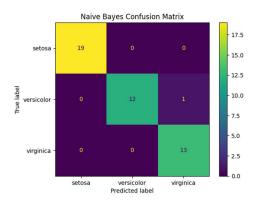
# Ex.4 Write a program to implement naive Bayesian classifier and show the performance of the classifier using suitable test Set.

# Algorithm:

- 1. Import dataset (Iris).
- 2. Split into train and test sets.
- 3. Train Naive Bayes classifier.
- 4. Predict on test set.
- 5. Display accuracy score.

#### Code.:

```
# Program 4: Naive Bayes Classifier
from sklearn.datasets import load iris
from sklearn.model selection import train test split
from sklearn.naive bayes import GaussianNB
from sklearn.metrics import confusion matrix, ConfusionMatrixDisplay, accuracy score
import matplotlib.pyplot as plt
# Step 1: Load dataset
iris = load iris()
X, y = iris.data, iris.target
# Step 2: Split data
X train, X test, y train, y test = train test split(X, y, test size=0.3, random state=42)
# Step 3: Train model
nb = GaussianNB()
nb.fit(X train, y train)
# Step 4: Predict
y pred = nb.predict(X test)
print("Accuracy:", accuracy score(y test, y pred))
# Step 5: Confusion Matrix visualization
cm = confusion matrix(y test, y pred)
disp = ConfusionMatrixDisplay(cm, display labels=iris.target names)
disp.plot(cmap="viridis")
plt.title("Naive Bayes Confusion Matrix")
plt.show()
```



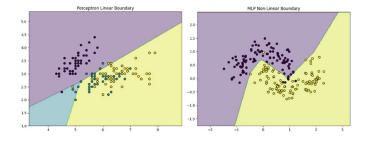
# Ex.:5 Construct a simple Perceptrom Neural Network and demonstrate linear and non-linear classification problem.

# Algorithm:

- 1. Load dataset (Iris).
- 2. Train Gaussian Naive Bayes.
- 3. Predict on test set.
- 4. Plot confusion matrix.
- 5. Show accuracy score.

#### Code.:

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.datasets import load iris, make moons
from sklearn.linear model import Perceptron
from sklearn.neural network import MLPClassifier
def plot decision boundary(model, X, y, title):
  """A helper function to plot the decision boundary for a trained model."""
  x \min_{x} \max = X[:, 0].\min() - 1, X[:, 0].\max() + 1
  y \min_{x \in X} y \max_{x \in X} = X[:, 1].\min() - 1, X[:, 1].\max() + 1
  xx, yy = np.meshgrid(np.arange(x min, x max, 0.02), np.arange(y min, y max, 0.02))
  Z = model.predict(np.c [xx.ravel(), yy.ravel()]).reshape(xx.shape)
  plt.figure(figsize=(8, 6))
  plt.contourf(xx, yy, Z, alpha=0.4, cmap="viridis")
  plt.scatter(X[:, 0], X[:, 1], c=y, edgecolor='k', cmap="viridis")
  plt.title(title)
  plt.show()
# --- 1. Linear Classification with Perceptron ---
print("Demonstrating Linear Classification...")
X \text{ lin, } y \text{ lin} = \text{load iris(return } X \text{ y=True)}
# Use only the first two features (sepal length and width)
perceptron = Perceptron(random state=42).fit(X lin[:, :2], y lin)
plot decision boundary(perceptron, X lin[:, :2], y lin, "Perceptron Linear Boundary")
# --- 2. Non-Linear Classification with MLP ---
print("Demonstrating Non-Linear Classification...")
X nonlin, y nonlin = make moons(n samples=200, noise=0.2, random state=42)
mlp = MLPClassifier(hidden layer sizes=(10, 10), max iter=1000, random state=42).fit(X nonlin,
y nonlin)
plot decision boundary(mlp, X nonlin, y nonlin, "MLP Non-Linear Boundary")
```



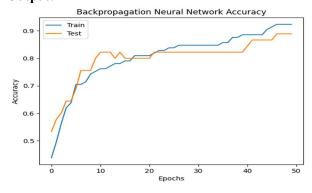
# Ex.:6 Construct a Back Propagation network and verify the performance by using suitable traning and text set

### Algorithm:

- 1. Load dataset (Iris).
- 2. Normalize data and one-hot encode labels.
- 3. Define a simple feed-forward NN with hidden layers.
- 4. Train using backpropagation.
- 5. Plot training vs validation accuracy.

#### Code.:

```
# Program 6: Backpropagation Neural Network
from sklearn.datasets import load iris
from sklearn.model selection import train test split
from sklearn.preprocessing import StandardScaler, LabelBinarizer
from tensorflow.keras.models import Sequential
from tensorflow.keras.layers import Dense
import matplotlib.pyplot as plt
iris = load iris()
X, y = iris.data, iris.target
X = StandardScaler().fit transform(X)
y = LabelBinarizer().fit transform(y)
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=42)
# Step 3: Define model
model = Sequential([
  Dense(10, input dim=X.shape[1], activation='relu'),
  Dense(3, activation='softmax')])
model.compile(loss='categorical crossentropy', optimizer='adam', metrics=['accuracy'])
history = model.fit(X train, y train, epochs=50, batch size=5, verbose=0, validation data=(X test,
y test))
plt.plot(history.history['accuracy'], label='Train')
plt.plot(history.history['val accuracy'], label='Test')
plt.title("Backpropagation Neural Network Accuracy")
plt.xlabel("Epochs")
plt.ylabel("Accuracy")
plt.legend()
plt.show()
```



Ex.7 Use the same dataset used for experiment 6 to train SVM classifier and CO3 compare the performance with back propagation network.

### Algorithm:

- 1. Load Iris dataset (use 2 features for 2D visualization).
- 2. Train-test split.
- 3. Train SVM classifier.
- 4. Predict & show accuracy.
- 5. Plot decision boundary.

#### Code.:

 $[1 \ 1] \rightarrow 0$ 

```
import numpy as np
import matplotlib.pyplot as plt
# Perceptron model
class Perceptron:
  def init (self, input size, lr=0.1, epochs=100): #FIX: Corrected constructor name from init
     self.W = np.zeros(input size + 1) # +1 for bias
     self.lr = lr
     self.epochs = epochs
  def activation(self, x):
     return np.where(x \ge 0, 1, 0) # step function
  def predict(self, x):
     z = np.dot(x, self.W[1:]) + self.W[0]
     return self.activation(z)
  def fit(self, X, y):
     for in range(self.epochs):
        for i in range(len(X)):
          update = self.lr * (y[i] - self.predict(X[i]))
          self.W[1:] += update * X[i]
          self.W[0] += update
# Linear Problem (OR)
X = \text{np.array}([[0,0],[0,1],[1,0],[1,1]])
y or = np.array([0,1,1,1])
p = Perceptron(input size=2)
p.fit(X, y or)
print("OR Gate Predictions (Solved):")
for i in range(len(X)):
  print(f''\{X[i]\} \rightarrow \{p.predict(X[i])\}'')
# Non-linear Problem (XOR)
y xor = np.array([0,1,1,0])
p2 = Perceptron(input size=2)
p2.fit(X, y xor)
print("\nXOR Gate Predictions (Failed):")
for i in range(len(X)):
  print(f''\{X[i]\} \rightarrow \{p2.predict(X[i])\}'')
Output:
OR Gate Predictions (Solved):
                                                     XOR Gate Predictions (Failed):
[0\ 0] \rightarrow 1
                                                     [0\ 0] \rightarrow 0
[0\ 1] \rightarrow 1
                                                     [0\ 1] \rightarrow 1
                                                     [1\ 0] \rightarrow 1
[1\ 0] -> 0
```

 $[1 \ 1] \rightarrow 1$ 

# Exp.8 Create a self-organizing map neural network for learning a set of itnagos and verify the performance

# Algorithm:

- 1. Load dataset (Iris).
- 2. Normalize data.
- 3. Initialize and train SOM grid.
- 4. Map data to neurons.
- 5. Plot SOM clusters with labels.

#### Code.:

import numpy as np import matplotlib.pyplot as plt from tensorflow.keras.datasets import mnist from minisom import MiniSom

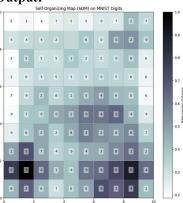
```
# 1. Load and prepare the image data
(X_train, y_train), _ = mnist.load_data()
X_data = X_train[:2000].reshape(2000, -1) / 255.0
y data = y train[:2000]
```

# 2. Create and train the Self-Organizing Map som = MiniSom(10, 10, X\_data.shape[1], sigma=1.5, learning\_rate=0.5, random\_seed=42) som.random\_weights\_init(X\_data) som.train\_random(X\_data, 100)

# 3. Verify performance by visualizing the clusters plt.figure(figsize=(10, 10)) plt.pcolor(som.distance\_map().T, cmap='bone\_r') plt.colorbar(label='Inter-neuron distance')

# Plot the most common digit for each winning node label\_map = som.labels\_map(X\_data, y\_data) for (i, j), win\_map in label\_map.items(): label = win\_map.most\_common(1)[0][0] plt.text(i + 0.5, j + 0.5, str(label), ha='center', va='center', bbox=dict(facecolor='white', alpha=0.5, lw=0))

plt.title("Self-Organizing Map (SOM) on MNIST Digits") plt.grid() plt.show()



#### Exp.9 Create a Giussian Misture Model for Image Segmentation

#### **Algorithm: GMM Image Segmentation**

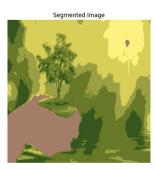
- 1. Read the input image.
- 2. Resize the image and reshape pixels into a 2D array.
- 3. Fit Gaussian Mixture Model (GMM) on pixel values.
- 4. Predict pixel clusters using GMM.
- 5. Replace pixels with corresponding cluster means.
- 6. Display original and segmented images side by side.

#### Code.:

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.mixture import GaussianMixture
from skimage.io import imread
from skimage.transform import resize
def segment image with gmm(image url, n components=5):
  try:
    image = imread(image url)
    if image.shape[2] == 4: # Handle RGBA images
       image = image[:, :, :3]
    image resized = (resize(image, (200, 200), anti aliasing=True) * 255).astype(np.uint8)
    pixels = image resized.reshape(-1, 3)
    gmm = GaussianMixture(n components=n components, covariance type='full'
    random state=42).fit(pixels)
    labels = gmm.predict(pixels)
    segmented image = gmm.means [labels].astype(np.uint8).reshape(image resized.shape)
    fig, (ax1, ax2) = plt.subplots(1, 2, figsize=(12, 6))
    fig.suptitle(f'GMM Image Segmentation ({n components} Clusters)', fontsize=16)
    ax1.imshow(image resized)
    ax1.set title('Original Image')
    ax1.axis('off')
    ax2.imshow(segmented image)
    ax2.set title('Segmented Image')
    ax2.axis('off')
    plt.tight layout()
    plt.show()
  except Exception as e:
    print(f"An error occurred: {e}")
if name == ' main ':
  sample image url = 'https://images.pexels.com/photos/162140/duck-bird-water-lake-162140.jpeg'
  num segments = 6 \# \text{Try changing this number (e.g., 3, 8, 12)}
  segment image with gmm(sample image url, n components=num segments)
Output.:
```

Image Segmentation using GMM with 6 Clusters





Exp.10 Write a Genetic Algorithm program for finding parameters which maximizes the Y value of the equation given where the equation has 6 inputs (x1 to x6) and 6 weights (w1 to w6). Input values are (x1, x2, x3, x4, x5, x6) = (4, 2, 7, 5, 11, 1). Goal: Find the weights (w1-w6) that maximize Y.

#### Algorithm (5 Steps)

- 1. Initialize a random population of weights.
- 3. Select the best solutions (parents) based on fitness.
- 4. Generate new solutions using crossover and mutation.
- 5. Repeat for several generations and return the best weights with maximum Y.

#### Code.:

```
import numpy as np
import matplotlib.pyplot as plt
X = \text{np.array}([4, 2, 7, 5, 11, 1])
def fitness(weights):
  return np.dot(weights, X)
pop size = 20
num weights = len(X)
population = np.random.randint(-10, 10, (pop size, num weights))
generations = 50
best scores = []
for gen in range(generations):
  scores = np.array([fitness(ind) for ind in population])
  best scores.append(scores.max())
  parents = population[scores.argsort()[-(pop size // 2):]]
  children = []
  for in range(pop size - len(parents)):
     p1, p2 = parents[np.random.randint(len(parents), size=2)]
     cp = np.random.randint(1, num weights - 1)
     child = np.concatenate((p1[:cp], p2[cp:]))
     children.append(child)
  children = np.array(children)
  mutation = np.random.randint(-2, 3, children.shape)
  children = children + mutation
  population = np.vstack((parents, children))
best idx = np.argmax([fitness(ind) for ind in population])
best weights = population[best idx]
best value = fitness(best weights)
plt.plot(best scores, marker='o')
plt.title("Genetic Algorithm Optimization Progress")
plt.xlabel("Generation")
plt.ylabel("Best Fitness Score (Y)")
plt.grid(True)
plt.show()
print("Best Weights:", best weights)
print("Maximum Y Value:", best value)
```

