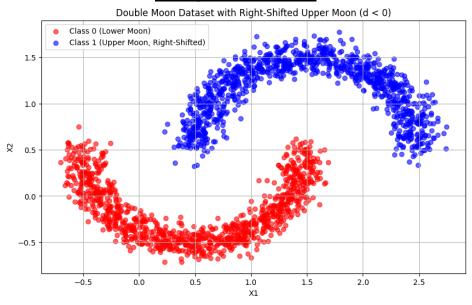
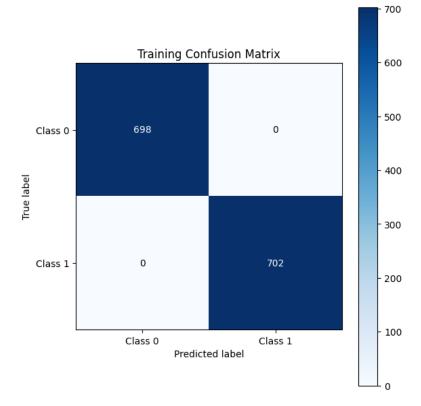
NAME: SHAILESH KUMAR

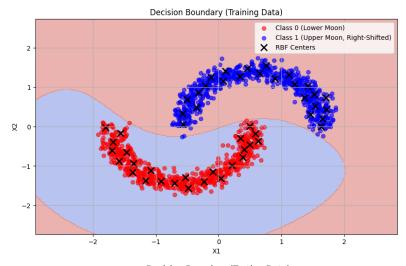
ROLL NO: 224EC6013

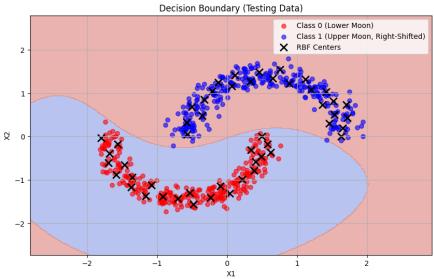
SIGNAL AND IMAGE PROCESSING

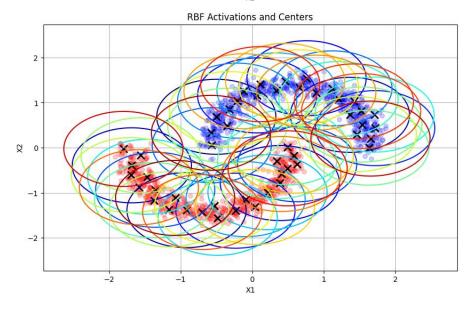
Assignment No: 3



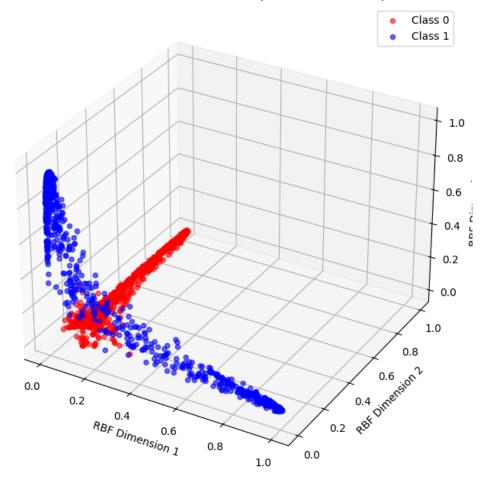


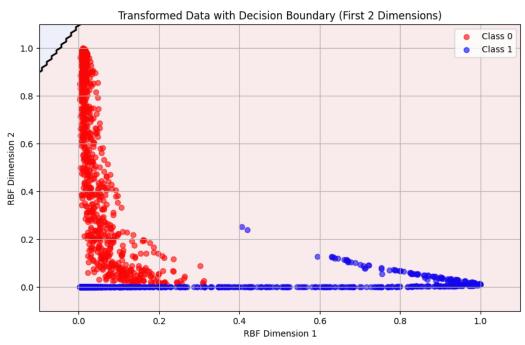






Data After RBF Transformation (First 3 Dimensions)





```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.model selection import train test split
from sklearn.metrics import accuracy score, confusion matrix
from sklearn.preprocessing import StandardScaler
from sklearn.cluster import KMeans
from sklearn.base import BaseEstimator, ClassifierMixin
from scipy.spatial.distance import cdist
from mpl toolkits.mplot3d import Axes3D
class RBFNetwork(BaseEstimator, ClassifierMixin):
  def init (self, n rbf=20, sigma=1.0, lr=0.01, epochs=1000):
    self.n rbf = n rbf
                          # Number of RBF neurons
    self.sigma = sigma
                           # Width of RBFs
    self.lr = lr
                      # Learning rate
    self.epochs = epochs # Training epochs
    self.centers = None
                           # RBF centers
    self.weights = None
                          # Output layer weights
    self.bias = None
                          # Output layer bias
  def rbf(self, X, center):
    return np.exp(-self.sigma * np.sum((X - center)**2, axis=1))
  def fit(self, X, y):
    # Step 1: Find RBF centers using K-means clustering
    kmeans = KMeans(n clusters=self.n rbf, random state=42)
    kmeans.fit(X)
    self.centers = kmeans.cluster centers
    # Step 2: Calculate RBF activations
    rbf activations = np.zeros((len(X), self.n rbf))
     for i, center in enumerate(self.centers):
       rbf activations[:, i] = self. rbf(X, center)
    # Step 3: Train output layer weights (using gradient descent)
    n_samples, n_features = rbf activations.shape
    self.weights = np.random.randn(n features)
    self.bias = np.random.randn(1)
    # Convert y to {-1, 1} for perceptron learning
    y_{-} = np.where(y == 0, -1, 1)
     for in range(self.epochs):
       # Forward pass
       outputs = np.dot(rbf activations, self.weights) + self.bias
       predictions = np.where(outputs \geq 0, 1, -1)
       # Compute error
       errors = y_ - predictions
       # Update weights
```

```
self.weights += self.lr * np.dot(rbf activations.T, errors)
       self.bias += self.lr * np.sum(errors)
    return self
  def predict(self, X):
    rbf activations = np.zeros((len(X), self.n rbf))
    for i, center in enumerate(self.centers):
       rbf activations[:, i] = self. rbf(X, center)
    outputs = np.dot(rbf activations, self.weights) + self.bias
    return np.where(outputs \geq 0, 1, 0)
  def predict from activations(self, rbf activations):
    outputs = np.dot(rbf activations, self.weights) + self.bias
    return np.where(outputs \geq 0, 1, 0)
# 1. Generate Double Moon Dataset with Right-Shifted Upper Moon
def generate shifted double moon(n samples=1000, radius=1.0, width=0.5, d=-0.5,
                   shift=1.0, noise=0.1):
  # Generate upper moon (shifted right)
  theta = np.linspace(0, np.pi, n samples//2)
  upper x = radius * np.cos(theta) + radius/2 + shift
  upper y = radius * np.sin(theta) - d
  # Generate lower moon (original position)
  lower x = radius * np.cos(theta) + radius/2
  lower y = -radius * np.sin(theta) + width
  # Add noise
  upper x \neq = np.random.normal(0, noise, n samples//2)
  upper y \neq = np.random.normal(0, noise, n samples//2)
  lower x += np.random.normal(0, noise, n samples//2)
  lower y += np.random.normal(0, noise, n samples//2)
  # Combine data
  X = np.vstack([np.column stack((lower x, lower y)),
           np.column stack((upper_x, upper_y))])
  y = np.hstack([np.zeros(n samples//2), np.ones(n samples//2)])
  return X, y
# 2. Generate and visualize the data
X, y = generate shifted double moon(n samples=2000, d=-0.5, shift=1.0, noise=0.1)
plt.figure(figsize=(10, 6))
plt.scatter(X[y==0, 0], X[y==0, 1], color='red', label='Class 0 (Lower Moon)', alpha=0.6)
plt.scatter(X[y==1, 0], X[y==1, 1], color='blue', label='Class 1 (Upper Moon, Right-Shifted)',
alpha=0.6)
plt.title("Double Moon Dataset with Right-Shifted Upper Moon (d < 0)")
plt.xlabel("X1")
```

```
plt.ylabel("X2")
plt.legend()
plt.grid(True)
plt.show()
# 3. Split data into training and testing sets
X train, X test, y train, y test = train test split(X, y, test size=0.3, random state=42)
# 4. Standardize the data
scaler = StandardScaler()
X train = scaler.fit transform(X train)
X \text{ test} = \text{scaler.transform}(X \text{ test})
# 5. Create and train RBF network
rbf net = RBFNetwork(n rbf=50, sigma=1.0, lr=0.01, epochs=1000)
rbf net.fit(X train, y train)
# 6. Evaluate the model
train pred = rbf net.predict(X train)
test pred = rbf net.predict(X test)
train acc = accuracy score(y train, train pred)
test acc = accuracy score(y test, test pred)
print(f"\nTraining Accuracy: {train acc:.4f}")
print(f"Testing Accuracy: {test acc:.4f}")
# 7. Plot confusion matrix
def plot confusion matrix(y true, y pred, title):
  cm = confusion_matrix(y_true, y_pred)
  plt.figure(figsize=(6, 6))
  plt.imshow(cm, interpolation='nearest', cmap=plt.cm.Blues)
  plt.title(title)
  plt.colorbar()
  tick marks = np.arange(2)
  plt.xticks(tick marks, ['Class 0', 'Class 1'])
  plt.yticks(tick marks, ['Class 0', 'Class 1'])
  thresh = cm.max() / 2.
  for i in range(cm.shape[0]):
     for j in range(cm.shape[1]):
       plt.text(j, i, format(cm[i, j], 'd'),
             ha="center", va="center",
             color="white" if cm[i, j] > thresh else "black")
  plt.ylabel('True label')
  plt.xlabel('Predicted label')
  plt.tight layout()
plot_confusion_matrix(y_train, train_pred, "Training Confusion Matrix")
plot confusion matrix(y test, test pred, "Testing Confusion Matrix")
```

```
plt.show()
#8. Plot decision boundary
def plot decision boundary(X, y, model, title):
  # Create a mesh grid
  h = 0.02 # step size
  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, h),
                np.arange(y min, y max, h))
  # Predict for each point in the grid
  Z = model.predict(np.c [xx.ravel(), yy.ravel()])
  Z = Z.reshape(xx.shape)
  # Plot
  plt.figure(figsize=(10, 6))
  plt.contourf(xx, yy, Z, alpha=0.4, cmap='coolwarm')
  plt.scatter(X[y==0, 0], X[y==0, 1], color='red', label='Class 0 (Lower Moon)', alpha=0.6)
  plt.scatter(X[y==1, 0], X[y==1, 1], color='blue', label='Class 1 (Upper Moon, Right-Shifted)',
alpha=0.6)
  # Plot RBF centers
  if hasattr(model, 'centers'):
     plt.scatter(model.centers[:, 0], model.centers[:, 1],
            color='black', marker='x', s=100, linewidths=2,
            label='RBF Centers')
  plt.title(title)
  plt.xlabel("X1")
  plt.ylabel("X2")
  plt.legend()
  plt.grid(True)
plot decision boundary(X train, y train, rbf net, "Decision Boundary (Training Data)")
plot decision boundary(X test, y test, rbf net, "Decision Boundary (Testing Data)")
plt.show()
# 9. Plot RBF activations (for visualization)
def plot rbf activations(model, X, y, title):
  if not hasattr(model, 'centers'):
     return
  plt.figure(figsize=(10, 6))
  # Create a grid for visualization
  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.linspace(x min, x max, 100),
                np.linspace(y_min, y_max, 100))
```

```
# Plot each RBF activation
  for i, center in enumerate(model.centers):
     grid points = np.c [xx.ravel(), yy.ravel()]
     activations = model. rbf(grid points, center)
     activations = activations.reshape(xx.shape)
     # Plot contour of this RBF
     cs = plt.contour(xx, yy, activations, levels=[0.5],
                colors=[plt.cm.jet(i/len(model.centers))])
  plt.scatter(X[y==0, 0], X[y==0, 1], color='red', alpha=0.2)
  plt.scatter(X[y==1, 0], X[y==1, 1], color='blue', alpha=0.2)
  plt.scatter(model.centers[:, 0], model.centers[:, 1],
         color='black', marker='x', s=100, linewidths=2)
  plt.title(title)
  plt.xlabel("X1")
  plt.ylabel("X2")
  plt.grid(True)
plot rbf activations(rbf net, X train, y train, "RBF Activations and Centers")
plt.show()
# 10. Plot the data after RBF transformation (first 3 dimensions for visualization)
def plot transformed data(X, y, model, title):
  if not hasattr(model, 'centers'):
     return
  # Transform the data using RBF activations
  rbf activations = np.zeros((len(X), model.n rbf))
  for i, center in enumerate(model.centers):
     rbf_activations[:, i] = model. rbf(X, center)
  # For visualization, we'll plot the first 3 dimensions
  fig = plt.figure(figsize=(12, 8))
  ax = fig.add subplot(111, projection='3d')
  # Plot class 0
  ax.scatter(rbf activations[y==0, 0],
         rbf activations[y==0, 1],
         rbf activations[y==0, 2],
         color='red', label='Class 0', alpha=0.6)
  # Plot class 1
  ax.scatter(rbf activations[y==1, 0],
         rbf activations[y==1, 1],
         rbf activations[y==1, 2],
         color='blue', label='Class 1', alpha=0.6)
  ax.set title(title)
  ax.set xlabel("RBF Dimension 1")
  ax.set ylabel("RBF Dimension 2")
```

```
ax.set zlabel("RBF Dimension 3")
  ax.legend()
  plt.show()
# Plot transformed training data
plot_transformed_data(X_train, y_train, rbf_net, "Data After RBF Transformation (First 3
Dimensions)")
# 11. Plot the transformed data with the separating hyperplane (first 2 dimensions)
def plot transformed decision boundary(X, y, model, title):
  if not hasattr(model, 'centers'):
    return
  # Transform the data using RBF activations
  rbf activations = np.zeros((len(X), model.n rbf))
  for i, center in enumerate(model.centers):
    rbf activations[:, i] = model. rbf(X, center)
  # We'll visualize just the first 2 dimensions
  dim1, dim2 = 0, 1 # Can change these indices to view different dimensions
  plt.figure(figsize=(10, 6))
  # Plot the transformed data points
  plt.scatter(rbf activations[y==0, dim1], rbf activations[y==0, dim2],
          color='red', label='Class 0', alpha=0.6)
  plt.scatter(rbf activations[y==1, dim1], rbf activations[y==1, dim2],
         color='blue', label='Class 1', alpha=0.6)
  # Plot the decision boundary (works well for 2D visualization)
  if model.weights is not None:
    # Create a grid for decision boundary
    x min, x max = rbf activations[:, dim1].min() - 0.1, rbf_activations[:, dim1].max() + 0.1
    y min, y max = rbf activations[:, dim2].min() - 0.1, rbf activations[:, dim2].max() + 0.1
    xx, yy = np.meshgrid(np.linspace(x min, x max, 100),
                  np.linspace(y min, y max, 100))
    # For the decision boundary, we need to consider all dimensions but we'll
    # set other dimensions to their mean values for this 2D visualization
    grid points = np.zeros((xx.ravel().shape[0], model.n rbf))
    grid points[:, dim1] = xx.ravel()
    grid points[:, dim2] = yy.ravel()
    # Set other dimensions to their mean values
    for dim in range(model.n rbf):
       if dim not in [dim1, dim2]:
          grid points[:, dim] = np.mean(rbf activations[:, dim])
    # Make predictions
    Z = model.predict from activations(grid points)
    Z = Z.reshape(xx.shape)
```