```
from google.colab import drive drive.mount('/content/drive')

→ Mounted at /content/drive
```

Some Helper Function:

→ Softmax Function:

```
import numpy as np
def softmax(z):
   Compute the softmax probabilities for a given input matrix.
   Parameters:
   z (numpy.ndarray): Logits (raw scores) of shape (m, n), where
                       - m is the number of samples.
                       - n is the number of classes.
   numpy.ndarray: Softmax probability matrix of shape (m, n), where
                   each row sums to 1 and represents the probability
                   distribution over classes.
   - The input to softmax is typically computed as: z = XW + b.
    - Uses numerical stabilization by subtracting the max value per row.
   # Numerical stability trick: subtract max value in each row
   z_{max} = np.max(z, axis=1, keepdims=True)
   z_stable = z - z_max
   exp_z = np.exp(z_stable) # Compute exponentials
   softmax_probs = exp_z / np.sum(exp_z, axis=1, keepdims=True) # Normalize
   return softmax_probs
```

Softmax Test Case:

This test case checks that each row in the resulting softmax probabilities sums to 1, which is the fundamental property of softmax.

```
WS02_MNIST_Digit_Classification_with_Logistic_Regression_Worksheet_Starter_Code.ipynb - Colab
    softmax_probs = exp_z / np.sum(exp_z, axis=1, keepdims=True) # Normalize
    return softmax_probs
# Example test case
z_{\text{test}} = \text{np.array}([[2.0, 1.0, 0.1], [1.0, 1.0, 1.0]])
softmax_output = softmax(z_test)
# Verify if the sum of probabilities for each row is 1
row_sums = np.sum(softmax_output, axis=1)
# Assert that the sum of each row is 1
assert np.allclose(row_sums, 1), f"Test failed: Row sums are {row_sums}"
print("Softmax function passed the test case! <a href="">✓"</a>)
→ Softmax function passed the test case!
Prediction Function:
import numpy as np
def predict_softmax(X, W, b):
    Predict the class labels for a set of samples using the trained softmax model.
    Parameters:
    X (numpy.ndarray): Feature matrix of shape (n, d), where n is the number of samples and d is the number of features.
    W (numpy.ndarray): Weight matrix of shape (d, c), where c is the number of classes.
    b (numpy.ndarray): Bias vector of shape (c,).
    Returns:
    numpy.ndarray: Predicted class labels of shape (n,), where each value is the index of the predicted class.
```

Test Function for Prediction Function:

return predicted_classes

logits = np.dot(X, W) + b # Compute raw scores (logits) softmax_probs = softmax(logits) # Apply softmax function

The test function ensures that the predicted class labels have the same number of elements as the input samples, verifying that the model produces a valid output shape.

predicted_classes = np.argmax(softmax_probs, axis=1) # Get index of max probability for each sample

```
import numpy as np
def softmax(z):
   Compute the softmax probabilities for a given input matrix.
   Parameters:
   z (numpy.ndarray): Logits (raw scores) of shape (m, n), where
                       - m is the number of samples.
                       - n is the number of classes.
   Returns:
   numpy.ndarray: Softmax probability matrix of shape (m, n), where
                  each row sums to 1.
   z_max = np.max(z, axis=1, keepdims=True) # Numerical stability
   z_stable = z - z_max
   exp_z = np.exp(z_stable)
   return exp_z / np.sum(exp_z, axis=1, keepdims=True)
def predict_softmax(X, W, b):
   Predict the class labels for a set of samples using the trained softmax model.
   Parameters:
```

```
X (numpy.ndarray): Feature matrix of shape (n, d).
    W (numpy.ndarray): Weight matrix of shape (d, c).
    b (numpy.ndarray): Bias vector of shape (c,).
    Returns:
    numpy.ndarray: Predicted class labels of shape (n,).
    logits = np.dot(X, W) + b # Compute raw scores (logits)
    softmax_probs = softmax(logits) # Apply softmax
    return np.argmax(softmax_probs, axis=1) # Get the class with the highest probability
# **Test Case for Prediction Function**
X_{\text{test}} = \text{np.array}([[0.2, 0.8], [0.5, 0.5], [0.9, 0.1]]) # 3 samples, 2 features
W_{\text{test}} = \text{np.array}([[0.4, 0.2, 0.1], [0.3, 0.7, 0.5]]) # 2 features, 3 classes
b_test = np.array([0.1, 0.2, 0.3]) # Bias for 3 classes
# Expected output: Array of predicted class labels (each between 0 and 2)
y_pred_test = predict_softmax(X_test, W_test, b_test)
# Validate output shape
assert y_pred_test.shape == (3,), f"Test failed: Expected shape (3,), got {y_pred_test.shape}"
# Print the predicted labels
print("Predicted class labels:", y_pred_test)
→ Predicted class labels: [1 1 0]
Loss Function:
import numpy as np
def loss_softmax(y_pred, y):
    Compute the cross-entropy loss for a single sample.
    y_pred (numpy.ndarray): Predicted probabilities of shape (c,) for a single sample,
                             where c is the number of classes.
    y (numpy.ndarray): True labels (one-hot encoded) of shape (c,), where c is the number of classes.
    float: Cross-entropy loss for the given sample.
    # Add a small value (epsilon) to prevent log(0) errors
    epsilon = 1e-12
    y_pred = np.clip(y_pred, epsilon, 1.0 - epsilon)
    # Compute the cross-entropy loss
    loss = -np.sum(y * np.log(y_pred))
    return loss
```

Test case for Loss Function:

This test case Compares loss for correct vs. incorrect predictions.

- · Expects low loss for correct predictions.
- · Expects high loss for incorrect predictions.

```
[0.85, 0.1, 0.05]])
# Compute loss for both cases
loss_correct = loss_softmax(y_pred_correct, y_true_correct)
loss_incorrect = loss_softmax(y_pred_incorrect, y_true_correct)
# Validate that incorrect predictions lead to a higher loss
assert loss_correct < loss_incorrect, f"Test failed: Expected loss_correct < loss_incorrect, but got {loss_correct:.4f} >= {loss_incorrect:.
# Print results
print(f"Cross-Entropy Loss (Correct Predictions): {loss_correct:.4f}")
print(f"Cross-Entropy Loss (Incorrect Predictions): {loss_incorrect:.4f}")
Tross-Entropy Loss (Correct Predictions): 0.4304
     Cross-Entropy Loss (Incorrect Predictions): 8.9872
Cost Function:
import numpy as np
def cost_softmax(X, y, W, b):
    Compute the average softmax regression cost (cross-entropy loss) over all samples.
    Parameters:
    X (numpy.ndarray): Feature matrix of shape (n, d), where n is the number of samples and d is the number of features.
    y (numpy.ndarray): True labels (one-hot encoded) of shape (n, c), where n is the number of samples and c is the number of classes.
    W (numpy.ndarray): Weight matrix of shape (d, c).
    b (numpy.ndarray): Bias vector of shape (c,).
    Returns:
    float: Average softmax cost (cross-entropy loss) over all samples.
    # Compute logits: Z = XW + b
    logits = np.dot(X, W) + b
    # Compute softmax probabilities
    softmax_probs = softmax(logits)
    # Numerical stability: clip probabilities to avoid log(0) errors
    epsilon = 1e-12
    softmax_probs = np.clip(softmax_probs, epsilon, 1.0 - epsilon)
    # Compute cross-entropy loss for each sample
    total_loss = -np.sum(y * np.log(softmax_probs))
    # Compute the average loss
    n = X.shape[0] # Number of samples
```

```
# Example input data
X_test = np.array([[0.2, 0.8], [0.5, 0.5], [0.9, 0.1]]) # (3 samples, 2 features)
W_test = np.array([[0.4, 0.2, 0.1], [0.3, 0.7, 0.5]]) # (2 features, 3 classes)
b_test = np.array([0.1, 0.2, 0.3]) # (3 classes)

# True labels (one-hot encoded)
y_test = np.array([[0, 1, 0], [1, 0, 0], [0, 0, 1]])

# Compute cost
cost_value = cost_softmax(X_test, y_test, W_test, b_test)
print("Softmax cost:", cost_value)
```

→ Softmax cost: 1.09862371811174

return total_loss / n

Test Case for Cost Function:

The test case assures that the cost for the incorrect prediction should be higher than for the correct prediction, confirming that the cost function behaves as expected.

```
import numpy as np
```

```
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# Example 1: Correct Prediction (Closer predictions)
X_{correct} = np.array([[1.0, 0.0], [0.0, 1.0]]) # Feature matrix for correct predictions
y_{correct} = np.array([[1, 0], [0, 1]]) # True labels (one-hot encoded, matching predictions)
W_{correct} = np.array([[5.0, -2.0], [-3.0, 5.0]]) # Weights for correct prediction
b_correct = np.array([0.1, 0.1]) # Bias for correct prediction
# Example 2: Incorrect Prediction (Far off predictions)
X_{incorrect} = np.array([[0.1, 0.9], [0.8, 0.2]]) # Feature matrix for incorrect predictions
y_{incorrect} = np.array([[1, 0], [0, 1]]) # True labels (one-hot encoded, incorrect predictions)
W_incorrect = np.array([[0.1, 2.0], [1.5, 0.3]]) # Weights for incorrect prediction
b_incorrect = np.array([0.5, 0.6]) # Bias for incorrect prediction
# Compute cost for correct predictions
cost_correct = cost_softmax(X_correct, y_correct, W_correct, b_correct)
# Compute cost for incorrect predictions
cost_incorrect = cost_softmax(X_incorrect, y_incorrect, W_incorrect, b_incorrect)
# Check if the cost for incorrect predictions is greater than for correct predictions
assert cost_incorrect > cost_correct, f"Test failed: Incorrect cost {cost_incorrect} is not greater than correct cost {cost_correct}"
# Print the costs for verification
print("Cost for correct prediction:", cost_correct)
print("Cost for incorrect prediction:", cost_incorrect)
print("Test passed!")
→ Cost for correct prediction: 0.0006234364133349324
     Cost for incorrect prediction: 0.29930861359446115
     Test passed!
Computing Gradients:
import numpy as np
def compute_gradient_softmax(X, y, W, b):
    Compute the gradients of the cost function with respect to weights and biases.
```

Commute anadienta

```
Parameters:
    X (numpy.ndarray): Feature matrix of shape (n, d).
    y (numpy.ndarray): True labels (one-hot encoded) of shape (n, c).
    W (numpy.ndarray): Weight matrix of shape (d, c).
    b (numpy.ndarray): Bias vector of shape (c,).
    Returns:
    tuple: Gradients with respect to weights (d, c) and biases (c,).
    # Compute logits
    logits = np.dot(X, W) + b
    # Compute softmax probabilities
    softmax_probs = softmax(logits)
    # Compute the error (difference between predicted probabilities and true labels)
    error = softmax_probs - y # Shape: (n, c)
    # Compute gradients
    n = X.shape[0] # Number of samples
    grad_W = np.dot(X.T, error) / n # Gradient w.r.t. weights, shape: (d, c)
    grad_b = np.sum(error, axis=0) / n # Gradient w.r.t. biases, shape: (c,)
    return grad_W, grad_b
# Example input data
X_{\text{test}} = \text{np.array}([[0.2, 0.8], [0.5, 0.5], [0.9, 0.1]]) # (3 samples, 2 features)
W_{\text{test}} = \text{np.array}([[0.4, 0.2, 0.1], [0.3, 0.7, 0.5]]) # (2 features, 3 classes)
b_test = np.array([0.1, 0.2, 0.3]) # (3 classes)
# True labels (one-hot encoded)
y_{test} = np.array([[0, 1, 0], [1, 0, 0], [0, 0, 1]])
```

```
# compute gradients
grad_W, grad_b = compute_gradient_softmax(X_test, y_test, W_test, b_test)

# Print gradients
print("Gradient w.r.t Weights:\n", grad_W)
print("Gradient w.r.t Biases:\n", grad_b)

Gradient w.r.t Weights:

[[ 0.0031051   0.11805685 -0.12116196]
   [-0.03600547 -0.09320977   0.12921524]]
Gradient w.r.t Biases:
   [-0.03290036   0.02484708   0.00805328]
```

Test case for compute_gradient function:

The test checks if the gradients from the function are close enough to the manually computed gradients using np.allclose, which accounts for potential floating-point discrepancies.

```
import numpy as np
# Define a simple feature matrix and true labels
X_{\text{test}} = \text{np.array}([[0.2, 0.8], [0.5, 0.5], [0.9, 0.1]]) # Feature matrix (3 samples, 2 features)
y_{test} = np.array([[1, 0, 0], [0, 1, 0], [0, 0, 1]]) # True labels (one-hot encoded, 3 classes)
# Define weight matrix and bias vector
W_{test} = np.array([[0.4, 0.2, 0.1], [0.3, 0.7, 0.5]]) # Weights (2 features, 3 classes)
b_{\text{test}} = \text{np.array}([0.1, 0.2, 0.3]) \# \text{Bias} (3 \text{ classes})
# Compute the gradients using the function
grad_W, grad_b = compute_gradient_softmax(X_test, y_test, W_test, b_test)
# Manually compute the predicted probabilities (using softmax function)
z test = np.dot(X test, W test) + b test
y_pred_test = softmax(z_test)
# Compute the manually computed gradients
grad_W_manual = np.dot(X_test.T, (y_pred_test - y_test)) / X_test.shape[0]
grad_b_manual = np.sum(y_pred_test - y_test, axis=0) / X_test.shape[0]
# Assert that the gradients computed by the function match the manually computed gradients
assert np.allclose(grad_W, grad_W_manual), f"Test failed: Gradients w.r.t. W are not equal.\nExpected: {grad_W_manual}\nGot: {grad_W}"
assert np.allclose(grad_b, grad_b_manual), f"Test failed: Gradients w.r.t. b are not equal.\nExpected: {grad_b_manual}\nGot: {grad_b}"
# Print the gradients for verification
print("Gradient w.r.t. W:", grad_W)
print("Gradient w.r.t. b:", grad_b)
print("Test passed!")
 → Gradient w.r.t. W: [[ 0.1031051 0.01805685 -0.12116196]
      [-0.13600547 0.00679023 0.12921524]]
     Gradient w.r.t. b: [-0.03290036 0.02484708 0.00805328]
     Test passed!
   Implementing Gradient Descent:
import numpy as no
def gradient_descent_softmax(X, y, W, b, alpha, n_iter, show_cost=False):
    Perform gradient descent to optimize the weights and biases.
    Parameters:
    X (numpy.ndarray): Feature matrix of shape (n, d).
    y (numpy.ndarray): True labels (one-hot encoded) of shape (n, c).
```

W (numpy.ndarray): Weight matrix of shape (d, c). b (numpy.ndarray): Bias vector of shape (c,).

show_cost (bool): Whether to display the cost at intervals.

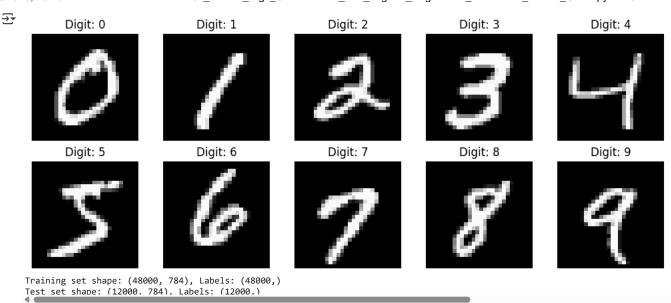
alpha (float): Learning rate.
n_iter (int): Number of iterations.

Returns:

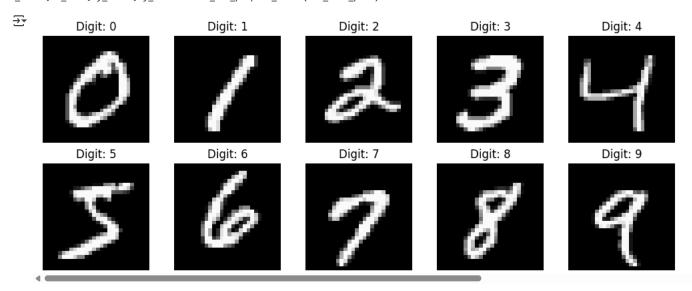
```
tuple: Optimized weights, biases, and cost history.
    cost_history = []
    for i in range(n_iter):
        # Compute gradients
        grad_W, grad_b = compute_gradient_softmax(X, y, W, b)
        # Update weights and biases using gradient descent
        W -= alpha * grad_W
       b -= alpha * grad_b
        # Compute cost
        cost = cost_softmax(X, y, W, b)
        cost_history.append(cost)
        # Display cost at intervals
        if show_cost and (i % 100 == 0 or i == n_iter - 1):
            print(f"Iteration {i}: Cost = {cost}")
    return W, b, cost_history
# Example input data
X_{\text{test}} = \text{np.array}([[0.2, 0.8], [0.5, 0.5], [0.9, 0.1]]) # (3 samples, 2 features)
W_test = np.random.rand(2, 3) # Random initialization (2 features, 3 classes)
b_test = np.random.rand(3) # Random initialization (3 classes)
y_{test} = np.array([[0, 1, 0], [1, 0, 0], [0, 0, 1]]) # One-hot encoded labels
# Hyperparameters
alpha = 0.1 # Learning rate
n_iter = 1000 # Number of iterations
# Run gradient descent
W_opt, b_opt, cost_hist = gradient_descent_softmax(X_test, y_test, W_test, b_test, alpha, n_iter, show_cost=True)
# Print final optimized weights and biases
print("Optimized Weights:\n", W_opt)
print("Optimized Biases:\n", b_opt)
→ Iteration 0: Cost = 1.1798757910598126
     Iteration 100: Cost = 0.7870631729938221
     Iteration 200: Cost = 0.6268705497085009
     Iteration 300: Cost = 0.5367725728401082
     Iteration 400: Cost = 0.4755970867696704
     Iteration 500: Cost = 0.42951381754933693
     Iteration 600: Cost = 0.3926222985407403
     Iteration 700: Cost = 0.36194683735713107
     Iteration 800: Cost = 0.3357930781515659
     Iteration 900: Cost = 0.31310300343208225
     Iteration 999: Cost = 0.29335414772356894
     Optimized Weights:
      [[ 0.76754981 -3.46676061 4.09294856]
      [ 0.93835218  4.42541883 -3.6698721 ]]
     Optimized Biases:
      [ 0.68905651  0.03164296 -0.09163724]
```

Preparing Dataset:

```
random_state (int) : Random seed for reproducibility (default: 42).
    Returns:
    X_train, X_test, y_train, y_test : Split dataset.
    # Load dataset
    df = pd.read_csv(csv_file)
    # Separate labels and features
    y = df.iloc[:, 0].values # First column is the label
    X = df.iloc[:, 1:].values # Remaining columns are pixel values
    # Normalize pixel values (optional but recommended)
    X = X / 255.0 # Scale values between 0 and 1
    # Split data into train and test sets
    X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=test_size, random_state=random_state)
    # Plot one sample image per class
    plot_sample_images(X, y)
    return X_train, X_test, y_train, y_test
def plot_sample_images(X, y):
    Plots one sample image for each digit class (0-9).
    Arguments:
    X (np.ndarray): Feature matrix containing pixel values.
    y (np.ndarray): Labels corresponding to images.
    plt.figure(figsize=(10, 4))
    unique_classes = np.unique(y) # Get unique class labels
    for i, digit in enumerate(unique_classes):
        index = np.where(y == digit)[0][0] # Find first occurrence of the class
        image = X[index].reshape(28, 28) # Reshape 1D array to 28x28
        plt.subplot(2, 5, i + 1)
        plt.imshow(image, cmap='gray')
        plt.title(f"Digit: {digit}")
       plt.axis('off')
    plt.tight_layout()
    plt.show()
# Load and process the MNIST dataset
X_train, X_test, y_train, y_test = load_and_prepare_mnist(csv_file)
# Print dataset shapes
print(f"Training set shape: {X_train.shape}, Labels: {y_train.shape}")
print(f"Test set shape: {X_test.shape}, Labels: {y_test.shape}")
```



csv_file_path = "_content/drive/MyDrive/AI/week2/worksheet-2/mnist_dataset.csv" # Path to saved dataset
X_train, X_test, y_train, y_test = load_and_prepare_mnist(csv_file_path)



→ A Quick debugging Step:

```
# Assert that X and y have matching lengths assert len(X_{train}) == len(y_{train}), f"Error: X and y have different lengths! X=\{len(X_{train})\}, y=\{len(y_{train})\}" print("Move forward: Dimension of Feture Matrix X and label vector y matched.")
```

→ Move forward: Dimension of Feture Matrix X and label vector y matched.

Train the Model:

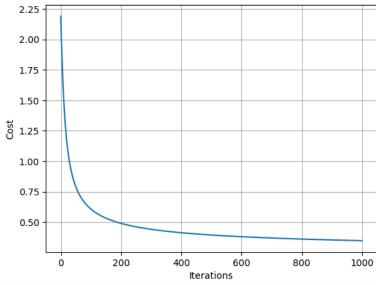
```
print(f"Training data shape: {X_train.shape}")
print(f"Test data shape: {X_test.shape}")

Training data shape: (48000, 784)
    Test data shape: (12000, 784)

from sklearn.preprocessing import OneHotEncoder

# Check if y_train is one-hot encoded
if len(y_train.shape) == 1:
    encoder = OneHotEncoder(sparse_output=False) # Use sparse_output=False for newer versions of sklearn
    y_train = encoder.fit_transform(y_train.reshape(-1, 1)) # One-hot encode labels
    y_test = encoder.transform(y_test.reshape(-1, 1)) # One-hot encode test labels
```

```
# Now y_train is one-hot encoded, and we can proceed to use it
d = X_{train.shape[1]} # Number of features (columns in X_{train})
c = y_train.shape[1] # Number of classes (columns in y_train after one-hot encoding)
# Initialize weights with small random values and biases with zeros
W = np.random.randn(d, c) * 0.01 # Small random weights initialized
b = np.zeros(c) # Bias initialized to 0
# Set hyperparameters for gradient descent
alpha = 0.1 # Learning rate
n_iter = 1000 # Number of iterations to run gradient descent
# Train the model using gradient descent
W_opt, b_opt, cost_history = gradient_descent_softmax(X_train, y_train, W, b, alpha, n_iter, show_cost=True)
# Plot the cost history to visualize the convergence
plt.plot(cost_history)
plt.title('Cost Function vs. Iterations')
plt.xlabel('Iterations')
plt.ylabel('Cost')
plt.grid(True)
plt.show()
→ Iteration 0: Cost = 2.1907523908880377
     Iteration 100: Cost = 0.6071497628881436
     Iteration 200: Cost = 0.4896775892376777
     Iteration 300: Cost = 0.44108426078870444
     Iteration 400: Cost = 0.4129891218247169
     Iteration 500: Cost = 0.39410246103548874
     Iteration 600: Cost = 0.3802628935679442
     Iteration 700: Cost = 0.36954004085975195
     Iteration 800: Cost = 0.3609019197961151
     Iteration 900: Cost = 0.35374051873910445
     Iteration 999: Cost = 0.34772746084789385
                                Cost Function vs. Iterations
         2.25
```



Evaluating the Model:

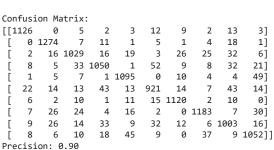
```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.metrics import confusion_matrix, precision_score, recall_score, f1_score

# Evaluation Function
def evaluate_classification(y_true, y_pred):
    """
    Evaluate classification performance using confusion matrix, precision, recall, and F1-score.

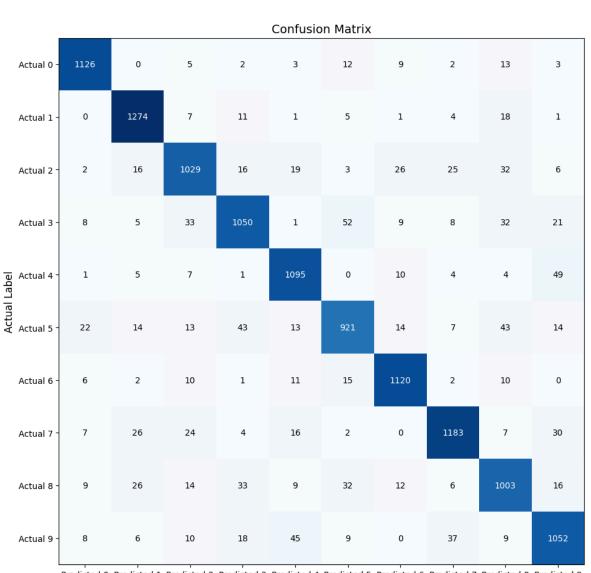
Parameters:
    y_true (numpy.ndarray): True labels
    y_pred (numpy.ndarray): Predicted labels
```

```
Returns:
    tuple: Confusion matrix, precision, recall, F1 score
    # Compute confusion matrix
    cm = confusion_matrix(y_true, y_pred)
    # Compute precision, recall, and F1-score
    precision = precision_score(y_true, y_pred, average='weighted')
    recall = recall_score(y_true, y_pred, average='weighted')
    f1 = f1_score(y_true, y_pred, average='weighted')
    return cm, precision, recall, f1
# Predict on the test set
y_pred_test = predict_softmax(X_test, W_opt, b_opt)
# Evaluate accuracy
y_test_labels = np.argmax(y_test, axis=1) # True labels in numeric form
# Evaluate the model
cm, precision, recall, f1 = evaluate_classification(y_test_labels, y_pred_test)
# Print the evaluation metrics
print("\nConfusion Matrix:")
print(cm)
print(f"Precision: {precision:.2f}")
print(f"Recall: {recall:.2f}")
print(f"F1-Score: {f1:.2f}")
# Visualizing the Confusion Matrix
fig, ax = plt.subplots(figsize=(12, 12))
cax = ax.imshow(cm, cmap='Blues') # Use a color map for better visualization
# Dynamic number of classes
num_classes = cm.shape[0]
ax.set_xticks(range(num_classes))
ax.set_yticks(range(num_classes))
ax.set_xticklabels([f'Predicted {i}' for i in range(num_classes)])
ax.set_yticklabels([f'Actual {i}' for i in range(num_classes)])
# Add labels to each cell in the confusion matrix
for i in range(cm.shape[0]):
    for j in range(cm.shape[1]):
        ax.text(j, i, cm[i, j], ha='center', va='center', color='white' if cm[i, j] > np.max(cm) / 2 else 'black')
# Add grid lines and axis labels
ax.grid(False)
plt.title('Confusion Matrix', fontsize=14)
plt.xlabel('Predicted Label', fontsize=12)
plt.ylabel('Actual Label', fontsize=12)
# Adjust layout
plt.tight_layout()
plt.colorbar(cax)
plt.show()
```

∓*



Recall: 0.90 F1-Score: 0.90



Predicted 0 Predicted 1 Predicted 2 Predicted 3 Predicted 5 Predicted 5 Predicted 6 Predicted 7 Predicted 8 Predicted 9 Predicted Label

1200

- 1000

800

600

400

200

0

Linear Seperability and Logistic Regression:

```
import numpy as np
import matplotlib.pvplot as plt
from sklearn.datasets import make_classification, make_circles
from sklearn.model selection import train test split
from sklearn.linear_model import LogisticRegression
# Set random seed for reproducibility
np.random.seed(42)
# Generate a synthetic dataset that is linearly separable
X_linear_separable, y_linear_separable = make_classification(
    n_samples=200, n_features=2, n_informative=2, n_redundant=0, n_clusters_per_class=1, random_state=42
# Split the dataset into training and testing sets (80% train, 20% test)
X_train_linear, X_test_linear, y_train_linear, y_test_linear = train_test_split(
    X_linear_separable, y_linear_separable, test_size=0.2, random_state=42
# Train a Logistic Regression model on the linearly separable dataset
logistic_model_linear_separable = LogisticRegression()
logistic_model_linear_separable.fit(X_train_linear, y_train_linear)
# Generate a synthetic dataset that is non-linearly separable (circles pattern)
X_non_linear_separable, y_non_linear_separable = make_circles(
    n_samples=200, noise=0.1, factor=0.5, random_state=42
# Split the dataset into training and testing sets (80% train, 20% test)
X_train_non_linear, X_test_non_linear, y_train_non_linear, y_test_non_linear = train_test_split(
    X_non_linear_separable, y_non_linear_separable, test_size=0.2, random_state=42
# Train a Logistic Regression model on the non-linearly separable dataset
logistic_model_non_linear_separable = LogisticRegression()
logistic_model_non_linear_separable.fit(X_train_non_linear, y_train_non_linear)
# Function to plot the decision boundary of a trained model
def plot_decision_boundary(ax, model, X, y, title):
    h = 0.02 # Step size for mesh grid
    x_{min}, x_{max} = X[:, 0].min() - 1, X[:, 0].max() + 1
    y_{min}, y_{max} = X[:, 1].min() - 1, X[:, 1].max() + 1
    # Create a mesh grid over the feature space
    xx, yy = np.meshgrid(np.arange(x_min, x_max, h), np.arange(y_min, y_max, h))
    # Predict class labels for each point in the grid
    Z = model.predict(np.c_[xx.ravel(), yy.ravel()])
    Z = Z.reshape(xx.shape)
    # Plot the decision boundary using contour plot
    ax.contourf(xx, yy, Z, alpha=0.8, cmap=plt.cm.Paired)
    # Scatter plot of the actual data points
    ax.scatter(X[:, 0], X[:, 1], c=y, edgecolors='k', cmap=plt.cm.Paired)
    # Formatting the plot
    ax.set_title(title)
    ax.set xlabel('Feature 1')
    ax.set_ylabel('Feature 2')
# Create subplots to visualize decision boundaries
fig, axes = plt.subplots(2, 2, figsize=(12, 10))
# Plot decision boundary for linearly separable data (Training set)
plot_decision_boundary(axes[0, 0], logistic_model_linear_separable, X_train_linear, y_train_linear, 'Linearly Separable Data (Training)')
# Plot decision boundary for linearly separable data (Testing set)
plot_decision_boundary(axes[0, 1], logistic_model_linear_separable, X_test_linear, y_test_linear, 'Linearly Separable Data (Testing)')
# Plot decision boundary for non-linearly separable data (Training set)
plot_decision_boundary(axes[1, 0], logistic_model_non_linear_separable, X_train_non_linear, y_train_non_linear, 'Non-Linearly Separable Data
```

```
# Plot decision boundary for non-linearly separable data (Testing set)
plot_decision_boundary(axes[1, 1], logistic_model_non_linear_separable, X_test_non_linear, y_test_non_linear, 'Non-Linearly Separable Data (Tour Adjust layout for better spacing
plt.tight_layout()

# Save the plot as a PNG file
plt.savefig('decision_boundaries.png')

# Display the plots
plt.show()
```

