SOFTWARE USED: Jupyter Notebook/ Google Colab

Instructions for Use:

- $1.\ Download\ the\ dataset\ from\ -\ https://www.kaggle.com/sartajbhuvaji/brain-tumor-classification-mri?select=Training$
- 2. Give the appropriate file paths to the tumor and notumor variables.
- 3. Run each cell one by one, after the successful execution of the previous cell.

Brain Tumor Detection Using Histogram of Oriented Gradient (HOG) And Support Vector Machine (SVM)

About the Brain MRI Images dataset:

The dataset contains 2 folders: yes and no which contains Brain MRI Images. The folder yes contains 827 Brain MRI Images that are tumorous and the folder no contains 395 Brain MRI Images that are non-tumorous.

Import necessary libraries

```
In [1]: import numpy as np
        import pandas as pd
        import openpyxl
        import os
        import cv2
        from skimage.feature import hog
        from sklearn import svm
        from sklearn.model selection import train test split
        from sklearn.metrics import roc curve, roc auc score, auc
        from sklearn.metrics import f1 score
        import matplotlib.pyplot as plt
        from skimage.feature import hog
        from skimage import exposure
        import imutils
        from matplotlib import pyplot as plt
        from sklearn import svm
        from sklearn.model_selection import train_test_split
        from sklearn.metrics import accuracy_score
        from sklearn.preprocessing import StandardScaler
```

Data Preparation & Preprocessing

```
# Load tumor images
for filename in os.listdir(tumor dir):
    if filename.endswith('.jpg'):
        image = cv2.imread(os.path.join(tumor_dir, filename))
        # Resize the image to the target size
        image = cv2.resize(image, target size)
        # Perform any necessary preprocessing
        image = cv2.cvtColor(image, cv2.COLOR BGR2GRAY) # Convert to grayscale
        # Call the crop brain contour function
        images.append(image)
        labels.append(1) # Tumor images are labeled as 1
# Load non-tumor images
for filename in os.listdir(non tumor dir):
    if filename.endswith('.jpg'):
        image = cv2.imread(os.path.join(non tumor dir, filename))
        # Resize the image to the target size
        image = cv2.resize(image, target_size)
        # Perform any necessary preprocessing
        image = cv2.cvtColor(image, cv2.COLOR BGR2GRAY) # Convert to grayscale
        # Call the crop brain contour function
        images.append(image)
        labels.append(0) # Non-tumor images are labeled as 0
# Ensure that all images have the same dimensions
images = [cv2.resize(image, target size) for image in images]
# Convert image and label lists to numpy arrays
images = np.array(images)
labels = np.array(labels)
```

```
In [3]: import matplotlib.pyplot as plt

# Select a few sample images to display
sample_images = images[:5] # Display the first 5 images as samples

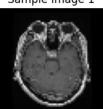
# Define a function to display the images
def display_images(images, titles):
    fig, axes = plt.subplots(1, len(images), figsize=(12, 3))

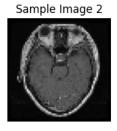
for i, ax in enumerate(axes):
    ax.imshow(images[i], cmap='gray')
    ax.set_title(titles[i])
    ax.axis('off')

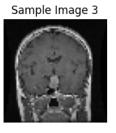
plt.show()

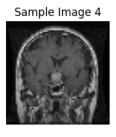
# Titles for the sample images
sample_titles = ["Sample Image 1", "Sample Image 2", "Sample Image 3", "Sample Image
# Display the sample images
display_images(sample_images, sample_titles)
```

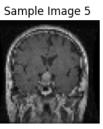






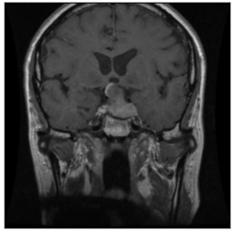




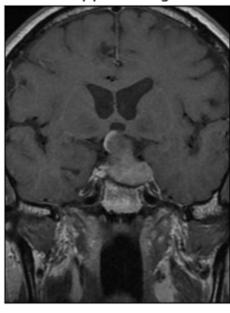


```
In [4]: def crop brain contour(image, plot=False):
        # Convert the image to grayscale, and blur
            if len(image.shape) == 2:
                gray = image
            else:
                gray = cv2.cvtColor(image, cv2.COLOR BGR2GRAY)
                gray = cv2.GaussianBlur(gray, (5, 5), 0)
        # Threshold the image, then perform a series of erosions + dilations to remove any
            thresh = cv2.threshold(gray, 45, 255, cv2.THRESH BINARY)[1]
            thresh = cv2.erode(thresh, None, iterations=2)
            thresh = cv2.dilate(thresh, None, iterations=2)
        # Find contours in thresholded image, then grab the largest one
            cnts = cv2.findContours(thresh.copy(), cv2.RETR_EXTERNAL, cv2.CHAIN_APPROX_SIMP
            cnts = imutils.grab contours(cnts)
        # At Least one contour was found
            if cnts:
                c = max(cnts, key=cv2.contourArea)
                # Find the extreme points
                extLeft = tuple(c[c[:, :, 0].argmin()][0])
                extRight = tuple(c[c[:, :, 0].argmax()][0])
                extTop = tuple(c[c[:, :, 1].argmin()][0])
                extBot = tuple(c[c[:, :, 1].argmax()][0])
        # Crop a new image out of the original image using the four extreme points (left, r
                new_image = image[extTop[1]:extBot[1], extLeft[0]:extRight[0]]
            else:
        # No contours found, return the original image
                new_image = image
            if plot:
                plt.figure()
                plt.subplot(1, 2, 1)
                plt.imshow(image)
                plt.tick_params(axis='both', which='both',
                                top=False, bottom=False, left=False, right=False,
                                 labelbottom=False, labeltop=False, labelleft=False, labelri
                plt.title('Original Image')
                plt.subplot(1, 2, 2)
```









Extracting hog features

```
In [6]: def extract_hog_features(images):
    features = []
    for image in images:
        # Resize and convert to grayscale if needed
        image = cv2.resize(image, (64, 64))
        if len(image.shape) == 3: # Ensure grayscale
            image = cv2.cvtColor(image, cv2.COLOR_BGR2GRAY)

# Extract HOG features parameters
    feature_vector = hog(image, pixels_per_cell=(8, 8), cells_per_block=(2, 2),

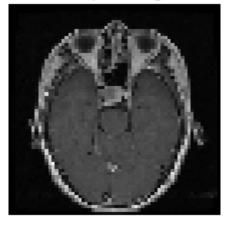
# cell size, block size, block normalization

    features.append(feature_vector)
        # return the feature matrix
    return np.array(features)
```

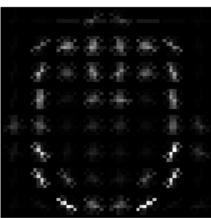
```
# Define X and y based on dataset and extracted features
X = extract_hog_features(images) # X should be the feature matrix
y = labels # y should be the corresponding labels
```

```
In [7]: sample_image = images[100]
        # Check if the sample image is valid and its shape contains the expected dimensions
        if sample_image is not None and len(sample_image.shape) == 2 and sample_image.shape
            # Adjust these parameters based on image size
            pixels_per_cell = (8, 8)
            cells_per_block = (2, 2)
            # Calculate HOG features
            fd, hog_image = hog(sample_image, pixels_per_cell=pixels_per_cell, cells_per_bl
            # Plot the HOG features
            fig, (ax1, ax2) = plt.subplots(1, 2, figsize=(6, 3), sharex=True, sharey=True)
            ax1.imshow(sample_image, cmap=plt.cm.gray)
            ax1.set_title('Sample Image')
            ax1.axis('off')
            ax2.imshow(hog_image, cmap=plt.cm.gray)
            ax2.set title('HOG Features')
            ax2.axis('off')
            plt.show()
        else:
            print("Invalid sample image or dimensions.")
```

Sample Image



HOG Features



```
In [8]: #extracting hog features in .xlsx file
hog_features = extract_hog_features(images)

# Create a DataFrame to store HOG features
data = pd.DataFrame(hog_features)

# Add LabeLs to the DataFrame
data['Labels'] = labels

# Save the DataFrame to an Excel file
excel_file = 'hog_features.xlsx'
data.to_excel(excel_file, index=False)
```

```
print(f'HOG features saved to {excel_file}')
       data.head()
     HOG features saved to hog_features.xlsx
Out[8]:
       1 0.099434 0.029215
                         0.027257  0.040866  0.071981
                                                0.017518  0.014160  0.000000  0.000000
       2 0.027277 0.018087
                         0.013534 0.024894
                                        0.070325
                                                0.015072 0.022019 0.034812 0.013796
       3 0.067036 0.047399
                         0.018749 0.072046 0.129830 0.042744 0.014996 0.058078 0.006348
       4 0.023618 0.029191 0.042352 0.071796 0.050758 0.033817 0.027648 0.089340 0.000000
      5 rows × 1765 columns
```

Splitting into training and testing dataset

```
In [9]: # Ensure that X is a 2D array
         if len(X.shape) == 1:
             X = np.array([x.flatten() for x in X])
         # Split the dataset into training and testing sets
         X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_sta
         print("X_train shape:", X_train.shape)
         print("y_train shape:", y_train.shape)
       X train shape: (977, 1764)
       y_train shape: (977,)
In [10]: print ("number of training examples = " + str(X_train.shape[0]))
         print ("number of test examples = " + str(X_test.shape[0]))
         print ("X train shape: " + str(X train.shape))
         print ("Y_train shape: " + str(y_train.shape))
         print ("X_test shape: " + str(X_test.shape))
         print ("Y_test shape: " + str(y_test.shape))
       number of training examples = 977
       number of test examples = 245
       X_train shape: (977, 1764)
       Y_train shape: (977,)
       X_test shape: (245, 1764)
       Y_test shape: (245,)
```

Implementing SVM

```
In [11]: scaler = StandardScaler()
         #Standardization data and involves scaling features to have a mean and standard dev
         X train = scaler.fit transform(X train)
         X_test = scaler.transform(X_test)
In [12]: from sklearn.model_selection import GridSearchCV
         # Define the parameter grid for hyperparameter tuning
         param_grid = {
             # C represents the regularization parameter in an SVM.
             'C': [0.1, 1, 10],
             #qamma defines how much influence each training example has in determining the
             'gamma': [0.001, 0.01, 0.1],
             #The kernel parameter specifies type of kernel function to be used in the SVM.(
             'kernel': ['rbf'],
         }
         # Create an SVM classifier
         clf = GridSearchCV(svm.SVC(kernel='linear', probability=True), param_grid, cv=3)
```

Hyperparameter Tuning

```
In [15]: # Train the SVM with hyperparameter tuning
         clf.fit(X_train, y_train)
         # Get the best hyperparameters
         best_params = clf.best_params_
         print(f"Best hyperparameters: {best_params}")
         # Make predictions on the test set
         y_pred = clf.predict(X_test)
         y_prob = clf.predict_proba(X_test)[:, 1] # Probability of class 1 (tumor)
         #Evaluate the model using ROC
         fpr, tpr, thresholds = roc_curve(y_test, y_prob)
         roc auc = auc(fpr, tpr)
         # Evaluate the SVM model
         accuracy_SVM = accuracy_score(y_test, y_pred)
         print(f"Accuracy: {accuracy_SVM * 100:.2f}%")
       Best hyperparameters: {'C': 10, 'gamma': 0.001, 'kernel': 'rbf'}
       Accuracy: 98.78%
         1.1.1
In [16]:
         The C parameter is the regularization parameter in an SVM.
         The gamma is kernel coefficient that control shape of decision boundary
         The kernel maps input data in multi-D shape, rbf (radial basis fn)
```

Out[16]: '\nThe C parameter is the regularization parameter in an SVM.\nSmaller values of C (e.g., 0.1) result in a more relaxed model.\nLarger values of C (e.g., 10) make the model stricter, small margin\n'

Implementing KNN

```
In [17]: from sklearn.neighbors import KNeighborsClassifier
         from sklearn.model selection import train test split
         from sklearn.metrics import accuracy score
         from sklearn.preprocessing import StandardScaler
         from sklearn.model selection import GridSearchCV
         # Assuming you have your X_train, X_test, y_train, and y_test defined
         # Standardize the features
         scaler = StandardScaler()
         X_train_KNN = scaler.fit_transform(X_train)
         X test KNN = scaler.transform(X test)
         # Define the parameter grid for hyperparameter tuning
         param grid KNN = {
             'n_neighbors': [3, 5, 7, 9],
             'weights': ['uniform', 'distance'],
             'p': [1, 2], # 1 for Manhattan distance, 2 for Euclidean distance
         }
         # Create a KNN classifier
         knn = KNeighborsClassifier()
         # Create a GridSearchCV object
         clf KNN = GridSearchCV(knn, param grid KNN, cv=3)
         # Fit the model
         clf_KNN.fit(X_train_KNN, y_train)
         # Get the best hyperparameters
         best_params_KNN = clf_KNN.best_params_
         # Make predictions
         y_pred_KNN = clf_KNN.predict(X_test_KNN)
         # Calculate accuracy
         accuracy KNN = accuracy score(y test, y pred KNN)
         print("Best KNN Hyperparameters:", best_params_KNN)
         print(f"Accuracy KNN: {accuracy_KNN * 100:.2f}%")
```

Best KNN Hyperparameters: {'n_neighbors': 3, 'p': 1, 'weights': 'distance'}
Accuracy KNN: 95.51%

Implementing Logistic Regression

```
In [18]: from sklearn.linear_model import LogisticRegression
         from sklearn.model_selection import train_test_split
         from sklearn.metrics import accuracy_score
         from sklearn.preprocessing import StandardScaler
         from sklearn.model_selection import GridSearchCV
         # Assuming you have your X_train, X_test, y_train, and y_test defined
         # Standardize the features
         scaler = StandardScaler()
         X train LR = scaler.fit transform(X train)
         X_test_LR = scaler.transform(X_test)
         # Define the parameter grid for hyperparameter tuning
         param grid LR = {
             'C': [0.1, 1, 10]
         # Create a Logistic Regression classifier
         lr = LogisticRegression()
         # Create a GridSearchCV object
         clf LR = GridSearchCV(lr, param grid LR, cv=3)
         # Fit the model
         clf LR.fit(X train LR, y train)
         # Get the best hyperparameters
         best_params_LR = clf_LR.best_params_
         # Make predictions
         y_pred_LR = clf_LR.predict(X_test_LR)
         # Calculate accuracy
         accuracy_LR = accuracy_score(y_test, y_pred_LR)
         print("Best Hyperparameters:", best params LR)
         print(f"Accuracy LR: {accuracy_LR * 100:.2f}%")
```

Best Hyperparameters: {'C': 1} Accuracy LR: 97.96%

Implementing Decision Tree

```
In [19]: from sklearn.tree import DecisionTreeClassifier
    from sklearn.model_selection import train_test_split
    from sklearn.metrics import accuracy_score
    from sklearn.model_selection import GridSearchCV

# Assuming you have your X_train, X_test, y_train, and y_test defined

# Define the parameter grid for hyperparameter tuning
    param_grid_DT = {
        'max_depth': [None, 10, 20, 30],
```

```
'min_samples_split': [2, 5, 10],
    'min_samples_leaf': [1, 2, 4],
# Create a Decision Tree classifier
dt = DecisionTreeClassifier()
# Create a GridSearchCV object
clf DT = GridSearchCV(dt, param grid DT, cv=3)
# Fit the model
clf DT.fit(X train, y train)
# Get the best hyperparameters
best params DT = clf DT.best params
# Make predictions
y_pred_DT = clf_DT.predict(X_test)
# Calculate accuracy
accuracy_DT = accuracy_score(y_test, y_pred_DT)
print("Best Hyperparameters DT:", best_params)
print(f"Accuracy DT: {accuracy_DT * 100:.2f}%")
```

Best Hyperparameters DT: {'C': 10, 'gamma': 0.001, 'kernel': 'rbf'} Accuracy DT: 88.98%

Implementing Naive Bayes

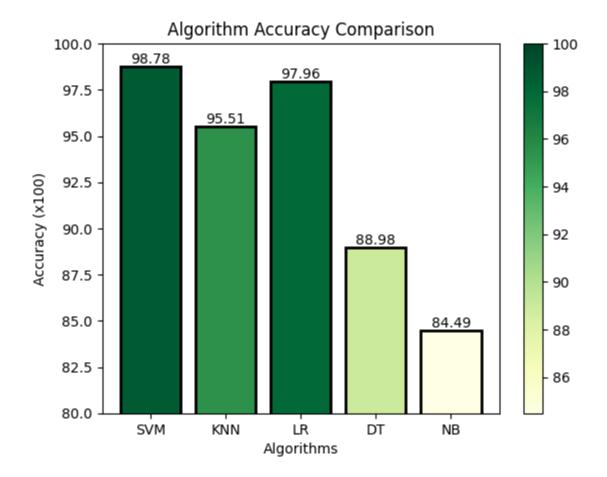
```
In [20]: from sklearn.naive_bayes import GaussianNB
         from sklearn.model_selection import train_test_split
         from sklearn.metrics import accuracy score
         from sklearn.preprocessing import StandardScaler
         # Assuming you have your X_train, X_test, y_train, and y_test defined
         # Standardize the features (optional for Gaussian Naive Bayes)
         scaler = StandardScaler()
         X_train_GNB = scaler.fit_transform(X_train)
         X_test_GNB = scaler.transform(X_test)
         # Create a Gaussian Naive Bayes classifier
         nb = GaussianNB()
         # Fit the model
         nb.fit(X_train_GNB, y_train)
         # Make predictions
         y_pred_GNB = nb.predict(X_test_GNB)
         # Calculate accuracy
         accuracy_GNB = accuracy_score(y_test, y_pred_GNB)
```

```
print(f"Accuracy GNB: {accuracy_GNB * 100:.2f}%")
```

Accuracy GNB: 84.49%

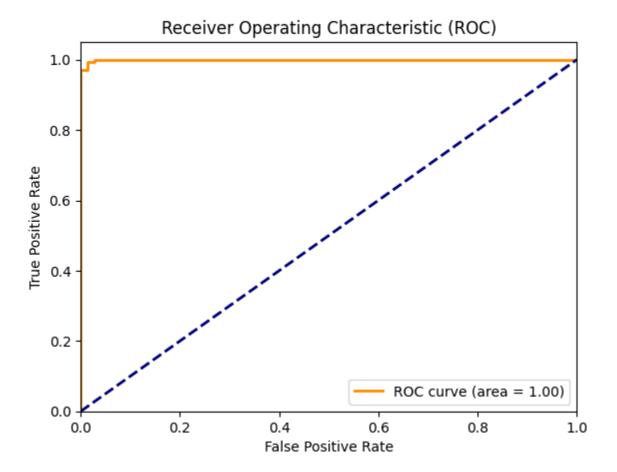
Comparision of Various ML Techniques

```
In [21]: import matplotlib.pyplot as plt
         import numpy as np
         from matplotlib.cm import ScalarMappable
         from matplotlib.colors import Normalize
         # Assuming you have accuracy values for each algorithm
         accuracies = [accuracy_SVM, accuracy_KNN, accuracy_LR, accuracy_DT, accuracy_GNB]
         labels = ['SVM', 'KNN', 'LR', 'DT', 'NB']
         # Increase the magnitude by multiplying with a constant (e.g., 100)
         magnified_accuracies = [accuracy * 100 for accuracy in accuracies]
         # Create a colormap based on the proximity to 100 (closer to 100, greener)
         norm = Normalize(vmin=min(magnified_accuracies), vmax=100)
         cmap = plt.get_cmap('YlGn') # Choose a colormap, e.g., Yellow-Green
         # Map the colors to accuracies
         colors = [cmap(norm(accuracy)) for accuracy in magnified_accuracies]
         # Create a figure and axis
         fig, ax = plt.subplots()
         # Create a bar chart to compare accuracy with color gradient
         bars = ax.bar(labels, magnified_accuracies, color=colors, edgecolor='black', linewi
         ax.set_xlabel('Algorithms')
         ax.set_ylabel('Accuracy (x100)')
         ax.set_title('Algorithm Accuracy Comparison')
         # Add labels to the bars with accuracy values
         for bar, v in zip(bars, magnified_accuracies):
             ax.text(bar.get_x() + bar.get_width() / 2, v, f'{v:.2f}', ha='center', va='bott
         ax.set ylim(80, 100) # Set the y-axis range from 80 to 100
         # Create a colorbar
         sm = ScalarMappable(cmap=cmap, norm=norm)
         sm.set_array([])
         cbar = plt.colorbar(sm, ax=ax,)
         plt.show()
```



Plot the ROC curve

```
In [22]: plt.figure()
   plt.plot(fpr, tpr, color='darkorange', lw=2, label='ROC curve (area = %0.2f)' % roc
   plt.plot([0, 1], [0, 1], color='navy', lw=2, linestyle='--')
   plt.xlim([0.0, 1.0])
   plt.ylim([0.0, 1.05])
   plt.xlabel('False Positive Rate')
   plt.ylabel('True Positive Rate')
   plt.title('Receiver Operating Characteristic (ROC)')
   plt.legend(loc='lower right')
   plt.show()
```



Accuracy and FPR Calculation

```
In [23]: from sklearn.metrics import accuracy_score, confusion_matrix

# Compute accuracy and FPR
accuracy = accuracy_score(y_test, y_pred)

# Compute confusion matrix
conf_matrix = confusion_matrix(y_test, y_pred)
true_positive = conf_matrix[1, 1]
false_positive = conf_matrix[0, 1]

# Calculate FPR (False Positive Rate)
fpr = false_positive / (false_positive + true_positive)

# Print accuracy and FPR
print("Accuracy: {:.2f}".format(accuracy))
print("False Positive Rate (FPR): {:.2f}".format(fpr))

# Measure of error of proportion of negative instances that were incorrectly classing.
```

Fowlkes-Mallows 1 Score (F1 Score)

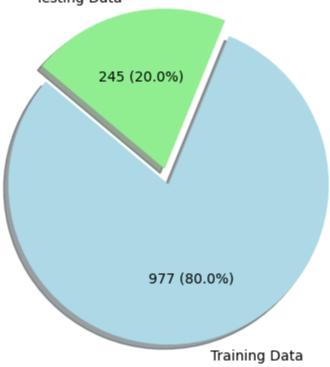
Accuracy: 0.99

False Positive Rate (FPR): 0.02

```
In [24]: def compute_f1_score(y_true, prob):
             # convert the vector of probabilities to a target vector
             y pred = np.where(prob > 0.5, 1, 0)
             score = f1_score(y_true, y_pred)
             return score
         f1score = compute_f1_score(y_test, y_prob)
         print(f"F1 score: {f1score}")
         #true positives between O(low precision) and 1(high precision)
       F1 score: 0.9912023460410556
In [25]: def data percentage(y):
             m=len(y)
             n_positive = np.sum(y)
             n_negative = m - n_positive
             pos prec = (n positive* 100.0)/ m
             neg_prec = (n_negative* 100.0)/ m
             print(f"Number of examples: {m}")
             print(f"Percentage of positive examples: {pos_prec}%, number of pos examples: {
             print(f"Percentage of negative examples: {neg prec}%, number of neg examples: {
In [26]: print("Training Data:")
         data percentage(y train)
         print("Testing Data:")
         data_percentage(y_test)
       Training Data:
       Number of examples: 977
       Percentage of positive examples: 67.24667349027635%, number of pos examples: 657
       Percentage of negative examples: 32.75332650972364%, number of neg examples: 320
       Testing Data:
       Number of examples: 245
       Percentage of positive examples: 69.38775510204081%, number of pos examples: 170
       Percentage of negative examples: 30.612244897959183%, number of neg examples: 75
In [27]: import matplotlib.pyplot as plt
         # Data
         labels = 'Training Data', 'Testing Data'
         sizes = [len(y_train), len(y_test)]
         colors = ['lightblue', 'lightgreen']
         explode = (0.1, 0) # Explode the first slice (Training Data)
         # Create pie chart for training and testing data
         plt.pie(sizes, explode=explode, labels=labels, colors=colors, autopct=lambda p: '{:
         plt.axis('equal') # Equal aspect ratio ensures that pie is drawn as a circle.
```

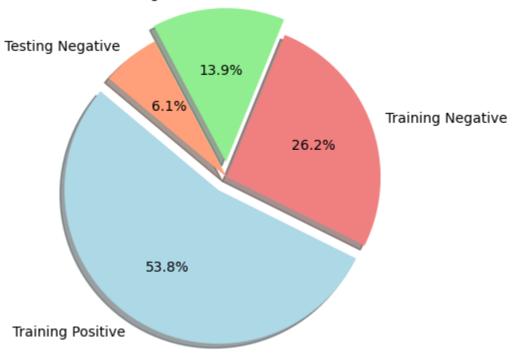
```
plt.title("Distribution of Training and Testing Data")
plt.show()
```

Distribution of Training and Testing Data Testing Data



```
In [28]: import matplotlib.pyplot as plt
         def data_percentage(y):
             m = len(y)
             n_positive = np.sum(y)
             n_negative = m - n_positive
             return n_positive, n_negative
         # Get the number of positive and negative examples for training and testing data
         train_pos, train_neg = data_percentage(y_train)
         test_pos, test_neg = data_percentage(y_test)
         # Data
         labels = ['Training Positive', 'Training Negative', 'Testing Positive', 'Testing Ne
         sizes = [train_pos, train_neg, test_pos, test_neg]
         colors = ['lightblue', 'lightcoral', 'lightgreen', 'lightsalmon']
         explode = (0.1, 0, 0.1, 0) # Explode the Positive slices
         # Create pie chart for training and testing data
         plt.pie(sizes, explode=explode, labels=labels, colors=colors, autopct='%1.1f%%', sh
         plt.axis('equal') # Equal aspect ratio ensures that pie is drawn as a circle.
         plt.title("Distribution of Positive and Negative Examples")
         plt.show()
```

Distribution of Positive and Negative Examples Testing Positive

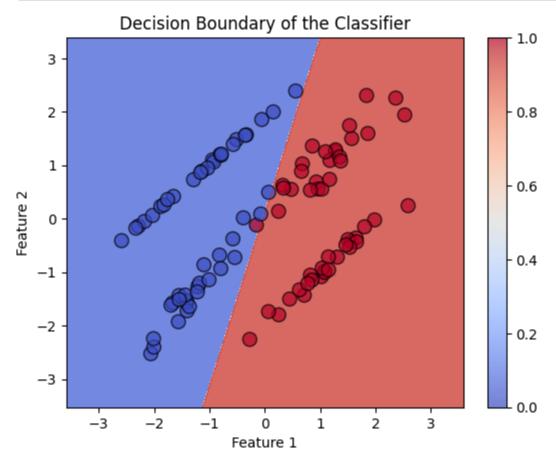


Decision Boundary Graph

```
In [29]: import numpy as np
         import matplotlib.pyplot as plt
         from sklearn import datasets
         from sklearn import svm
         # Generate a simple dataset
         X, y = datasets.make_classification(n_samples=100, n_features=2, n_informative=2, n
         # Train a binary classifier (SVM)
         clf = svm.SVC(kernel='linear')
         clf.fit(X, y)
         # Create a mesh grid over the feature space
         x_{min}, x_{max} = X[:, 0].min() - 1, X[:, 0].max() + 1
         y_{min}, y_{max} = X[:, 1].min() - 1, X[:, 1].max() + 1
         xx, yy = np.meshgrid(np.arange(x_min, x_max, 0.01), np.arange(y_min, y_max, 0.01))
         # Make predictions on the mesh grid
         Z = clf.predict(np.c_[xx.ravel(), yy.ravel()])
         Z = Z.reshape(xx.shape)
         # Plot the decision boundary
         plt.contourf(xx, yy, Z, cmap=plt.cm.coolwarm, alpha=0.8)
         # Plot the data points with different colors for each class
         scatter = plt.scatter(X[:, 0], X[:, 1], c=y, cmap=plt.cm.coolwarm, edgecolors='k',
```

```
# Add a color bar to the right
plt.colorbar(scatter)

plt.xlabel('Feature 1')
plt.ylabel('Feature 2')
plt.title('Decision Boundary of the Classifier')
plt.show()
```



| F1 score | 0.99 |

| Accuracy | 99% |