9th Homework

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Exercise 82

```
P_y(\omega_1)=1/3 and P_y(\omega_2)=2/3 P_N(\omega_1)=1/2 and P_N(\omega_2)=2/3 I_y=0.918 I_N=1
```

0.01997309402197489

Exercise 83

The problem below can be solved using decision trees

```
import matplotlib.pyplot as plt

# Given points
points = [(-5,0), (-3,0), (0,0), (-1,0), (1,0), (3,0)]

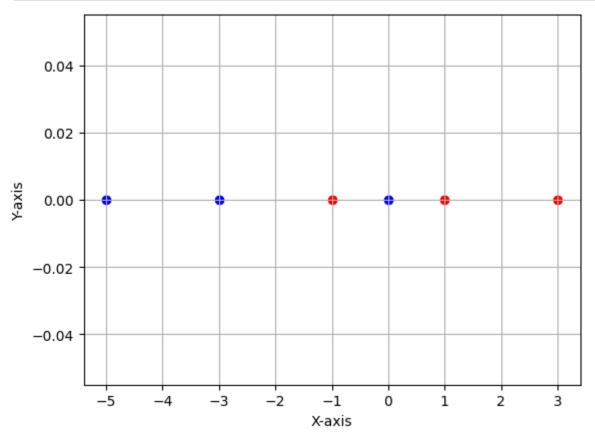
# Separate points into two groups for color coding
blue_points = points[:3]
red_points = points[3:]

# Extract x and y coordinates for plotting
x_blue, y_blue = zip(*blue_points)
x_red, y_red = zip(*red_points)

# Plot the points
plt.scatter(x_blue, y_blue, color='blue', label='Blue Points')
plt.scatter(x_red, y_red, color='red', label='Red Points')
```

```
# Set labels and title
plt.xlabel('X-axis')
plt.ylabel('Y-axis')

# Show the plot
plt.grid()
plt.show()
```



Exercise 84

$$egin{aligned} rac{\partial L(heta, heta_0,\lambda)}{\partial heta} &= rac{\partial \left(rac{1}{2} heta^T heta - \sum_{i=1}^N \lambda_i [y_i(heta^Tx_i + heta_0) - 1]
ight)}{\partial heta} &= heta - \sum_{i=1}^N \lambda_i y_i x_i = 0 \Rightarrow \ \Rightarrow heta &= \sum_{i=1}^N \lambda_i y_i x_i heta_i \end{aligned}$$

$$egin{aligned} rac{\partial L(heta, heta_0,\lambda)}{\partial heta_0} &= rac{\partial \left(rac{1}{2} heta^T heta - \sum_{i=1}^N \lambda_i [y_i(heta^Tx_i + heta_0) - 1]
ight)}{\partial heta_0} = -\sum_{i=1}^N \lambda_i y_i = 0 \Rightarrow \\ &\Rightarrow \sum_{i=1}^N \lambda_i y_i = 0 \ ext{ (2)} \end{aligned}$$

Using (1), (2) and the Lagrangian

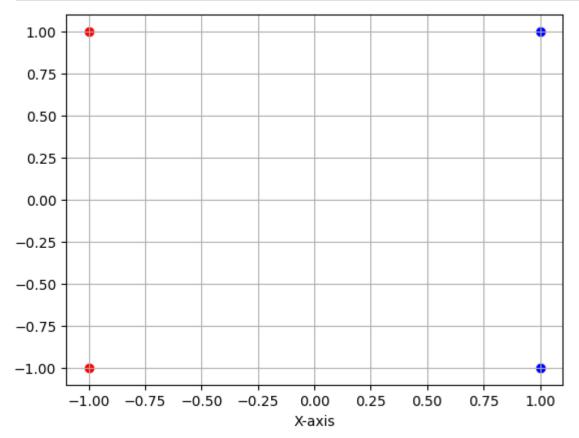
$$L(heta, heta_0,\lambda) = rac{1}{2} heta^T heta - \sum_{i=1}^N \lambda_i[y_i(heta^Tx_i+ heta_0)-1] = rac{1}{2}\sum_{i=1}^N (\lambda_iy_ix_i^T) \sum_{j=1}^N (\lambda_jy_jx_j) - \ = \sum_{i=1}^N \lambda_i - rac{1}{2}\sum_{i,j}\lambda_i\lambda_jy_iy_jx_i^Tx_i$$

Thus $min_{\theta,\lambda}L(\theta,\lambda)\Leftrightarrow max_{\lambda\geq 0}L(\theta,\lambda)$

Exercise 85

Looking at the plot below, it is expected that x=0 would be the line separating the 2 classes.

```
In [18]: import matplotlib.pyplot as plt
         import numpy as np
         # Given points
         x1 = np.array([-1, 1])
         x2 = np.array([-1, -1])
         x3 = np.array([1, -1])
         x4 = np.array([1, 1])
         # Plot the points
         plt.scatter(x1[0], x1[1], color='red', label='x1')
         plt.scatter(x2[0], x2[1], color='red', label='x2')
         plt.scatter(x3[0], x3[1], color='blue', label='x3')
         plt.scatter(x4[0], x4[1], color='blue', label='x4')
         # Set labels and title
         plt.xlabel('X-axis')
         plt.ylabel
         plt.grid()
```



$$J(\lambda) = \sum_{i=1}^N \lambda_i - rac{1}{2} \lambda_i \lambda_j y_j y_i x_i^T x_j$$
, thus

```
= \text{a lot of calculations later}
= \lambda_1 + \lambda_2 + \lambda_3 + \lambda_4 - \lambda_1^2 - \lambda_2^2 - \lambda_3^2 - \lambda_4^2 - 2\lambda_1\lambda_3 - 2\lambda_2\lambda_4
\frac{\partial J(\lambda)}{\partial \lambda_1} = 1 - 2\lambda_1 - 2\lambda_3 = 0 \Rightarrow \lambda_1 + \lambda_3 = \frac{1}{2}
\frac{\partial J(\lambda)}{\partial \lambda_2} = \lambda_2 + \lambda_4 = \frac{1}{2}
Further it stands that \sum_{i=1}^N \lambda_i y_i = 0 \Rightarrow \lambda_1 + \lambda_2 = \lambda_3 + \lambda_4 and using \theta = \sum_{i=1}^N \lambda_i y_i x_i and using the 4th Karush-Kuhn-Tucker condition \lambda_1 [y_1(\theta^T x_i + \theta_0) - 1] = 0 \Rightarrow \theta_0 = 0
Thus g(x) = \theta^T x + \theta_0 = 0 \Rightarrow x_1 = 0
```

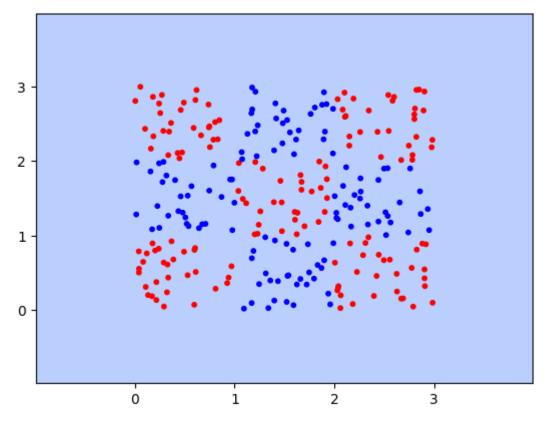
Exercise 86

```
In [28]: import scipy.io as sio
         import numpy as np
         import matplotlib.pyplot as plt
         from sklearn.metrics import accuracy_score, fbeta_score, make_scorer, roc_au
         Dataset a = sio.loadmat('HW9a.mat')
         train_x_a = Dataset_a['train_X']
         train y a = Dataset a['train y']
         test_x_a = Dataset_a['test_X']
         test_y_a = Dataset_a['test_y']
         from sklearn import svm
         def make_meshgrid(x, y, h=.02):
             """Create a mesh of points to plot in
             Parameters
             x: data to base x-axis meshgrid on
             y: data to base y-axis meshgrid on
             h: stepsize for meshgrid, optional
             Returns
             xx, yy : ndarray
             x_{min}, x_{max} = x.min() - 1, x.max() + 1
             y_{min}, y_{max} = y_{min}() - 1, y_{max}() + 1
             xx, yy = np.meshgrid(np.arange(x_min, x_max, h),
                                   np.arange(y_min, y_max, h))
             return xx, yy
```

```
def plot_contours(ax, clf, xx, yy, **params):
    """Plot the decision boundaries for a classifier.

Parameters
    _____
    ax: matplotlib axes object
    clf: a classifier
    xx: meshgrid ndarray
    yy: meshgrid ndarray
    params: dictionary of params to pass to contourf, optional
    """

Z = clf.predict(np.c_[xx.ravel(), yy.ravel()])
    Z = Z.reshape(xx.shape)
    out = ax.contourf(xx, yy, Z, **params)
    return out
```



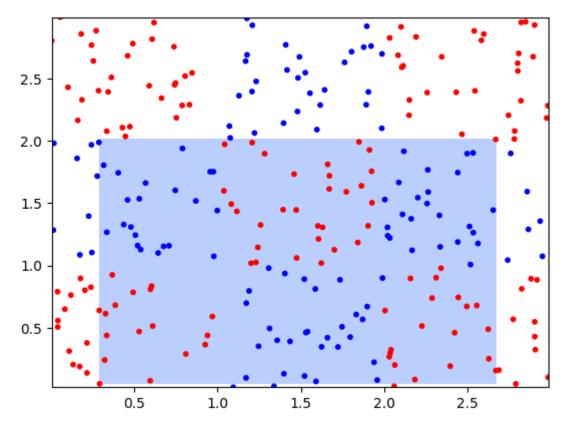
Accuracy: 50.00%

```
In [30]: # Polynomial Kernel
# Select kernel function: {'rbf', 'poly'}
clf_poly = svm.SVC(kernel='poly', C=10, degree=2, gamma=1)
clf_poly.fit(train_x_a, train_y_a.reshape(270))

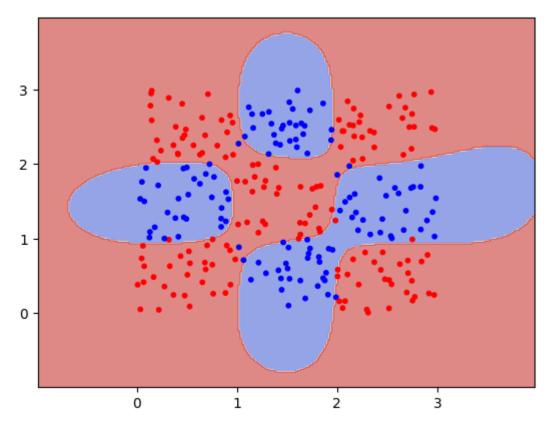
X00, X11 = test_x_a[:, 0], test_x_a[:, 1]
xx, yy = np.meshgrid(X00, X11)

fig, ax = plt.subplots(1, 1)
color = ['red' if l == 1 else 'blue' for l in test_y_a.reshape(270)]
plot_contours(ax, clf_poly, xx, yy, cmap=plt.cm.coolwarm, alpha=0.6)
ax.scatter(X00, X11, c=color, s=10, edgecolors='face')
plt.show()

accuracy = roc_auc_score(test_y_a, clf_poly.predict(test_x_a).flatten()) * 1
print(f"Accuracy: {accuracy:.2f}%")
```



Accuracy: 50.00%



```
1,
              1, 1,
                    1,
                      1, 1, 1, 1, 1,
                                    1,
                                      1,
                                           1, -1, -1, -1, -1,
           -1, -1, -1, -1, -1, -1, -1, -1,
                                    1, -1,
                                         1, -1,
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           1, -1, -1, -1, -1, -1, -1, -1,
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                            1, -1, -1, -1, -1, -1, -1, -1, -1, 1,
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                   1, -1, -1, -1, -1, -1,
                                      1, 1, 1, 1, -1, 1,
           -1, -1,
                    1, -1, 1, 1, -1, 1, 1, 1, 1, 1, 1, 1, 1,
           1, 1,
                        1,
                     1,
           -1, -1, -1,
                   -1, -1, -1, -1, -1, -1, -1, -1, -1, -1, -1, -1, -1,
                   1,
                1,
                      1, 1, 1, 1, 1, 1, -1, 1, 1, 1, 1, 1,
           -1, -1,
           1, 1, 1,
                    1,
                      1,
                         1, 1, 1, 1,
                                   1, 1, 1, 1, 1, -1],
          dtype=int16)
```

Trying out different parameters

```
In [32]: # Linear Kernel with Grid Search
C_values = [0.1, 1, 10]
degree_values = [2, 3, 4]
gamma_values = [0.1, 1, 10]

best_accuracy = 0
best_params = {}

for C in C_values:
    for degree in degree_values:
        for gamma in gamma_values:
```

```
clf_poly = svm.SVC(kernel='linear', C=C, degree=degree, gamma=ga
                     clf_poly.fit(train_x_a, train_y_a.reshape(270))
                     accuracy = roc_auc_score(test_y_a, clf_poly.predict(test_x_a).fl
                     print(f"Parameters: C={C}, Degree={degree}, Gamma={gamma}, Accur
                     if accuracy > best_accuracy:
                         best accuracy = accuracy
                         best_params = {'C': C, 'degree': degree, 'gamma': gamma}
         print(f"Best Parameters: {best params}")
         print(f"Best Accuracy: {best accuracy:.2%}")
        Parameters: C=0.1, Degree=2, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=2, Gamma=1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=2, Gamma=10, Accuracy: 50.00%
        Parameters: C=0.1, Degree=3, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=3, Gamma=1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=3, Gamma=10, Accuracy: 50.00%
        Parameters: C=0.1, Degree=4, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=4, Gamma=1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=4, Gamma=10, Accuracy: 50.00%
        Parameters: C=1, Degree=2, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=1, Degree=2, Gamma=1, Accuracy: 50.00%
        Parameters: C=1, Degree=2, Gamma=10, Accuracy: 50.00%
        Parameters: C=1, Degree=3, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=1, Degree=3, Gamma=1, Accuracy: 50.00%
        Parameters: C=1, Degree=3, Gamma=10, Accuracy: 50.00%
        Parameters: C=1, Degree=4, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=1, Degree=4, Gamma=1, Accuracy: 50.00%
        Parameters: C=1, Degree=4, Gamma=10, Accuracy: 50.00%
        Parameters: C=10, Degree=2, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=10, Degree=2, Gamma=1, Accuracy: 50.00%
        Parameters: C=10, Degree=2, Gamma=10, Accuracy: 50.00%
        Parameters: C=10, Degree=3, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=10, Degree=3, Gamma=1, Accuracy: 50.00%
        Parameters: C=10, Degree=3, Gamma=10, Accuracy: 50.00%
        Parameters: C=10, Degree=4, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=10, Degree=4, Gamma=1, Accuracy: 50.00%
        Parameters: C=10, Degree=4, Gamma=10, Accuracy: 50.00%
        Best Parameters: {'C': 0.1, 'degree': 2, 'gamma': 0.1}
        Best Accuracy: 50.00%
In [33]: # Polynomial Kernel with Grid Search
         C_{values} = [0.1, 1, 10]
         degree_values = [2, 3, 4]
         gamma_values = [0.1, 1, 10]
         best accuracy = 0
         best_params = {}
         for C in C values:
             for degree in degree_values:
                 for gamma in gamma_values:
                     clf_poly = svm.SVC(kernel='poly', C=C, degree=degree, gamma=gamm
```

```
clf_poly.fit(train_x_a, train_y_a.reshape(270))
                     accuracy = roc auc score(test y a, clf poly.predict(test x a).fl
                     print(f"Parameters: C={C}, Degree={degree}, Gamma={gamma}, Accur
                     if accuracy > best_accuracy:
                         best_accuracy = accuracy
                         best params = {'C': C, 'degree': degree, 'gamma': gamma}
         print(f"Best Parameters: {best_params}")
         print(f"Best Accuracy: {best accuracy:.2%}")
        Parameters: C=0.1, Degree=2, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=2, Gamma=1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=2, Gamma=10, Accuracy: 50.00%
        Parameters: C=0.1, Degree=3, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=3, Gamma=1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=3, Gamma=10, Accuracy: 50.00%
        Parameters: C=0.1, Degree=4, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=0.1, Degree=4, Gamma=1, Accuracy: 67.58%
        Parameters: C=0.1, Degree=4, Gamma=10, Accuracy: 66.33%
        Parameters: C=1, Degree=2, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=1, Degree=2, Gamma=1, Accuracy: 50.00%
        Parameters: C=1, Degree=2, Gamma=10, Accuracy: 50.00%
        Parameters: C=1, Degree=3, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=1, Degree=3, Gamma=1, Accuracy: 50.00%
        Parameters: C=1, Degree=3, Gamma=10, Accuracy: 50.00%
        Parameters: C=1, Degree=4, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=1, Degree=4, Gamma=1, Accuracy: 68.00%
        Parameters: C=1, Degree=4, Gamma=10, Accuracy: 70.25%
        Parameters: C=10, Degree=2, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=10, Degree=2, Gamma=1, Accuracy: 50.00%
        Parameters: C=10, Degree=2, Gamma=10, Accuracy: 50.00%
        Parameters: C=10, Degree=3, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=10, Degree=3, Gamma=1, Accuracy: 50.00%
        Parameters: C=10, Degree=3, Gamma=10, Accuracy: 44.00%
        Parameters: C=10, Degree=4, Gamma=0.1, Accuracy: 50.00%
        Parameters: C=10, Degree=4, Gamma=1, Accuracy: 66.75%
        Parameters: C=10, Degree=4, Gamma=10, Accuracy: 71.92%
        Best Parameters: {'C': 10, 'degree': 4, 'gamma': 10}
        Best Accuracy: 71.92%
In [36]: # RBF Kernel Grid Search
         # Best parameters and accuracy for RBF Kernel
         best accuracy rbf = 0
         best_params_rbf = {}
         for C in C values:
             for gamma in gamma_values:
                 clf_rbf = svm.SVC(kernel='rbf', C=C, gamma=gamma)
                 clf_rbf.fit(train_x_a, train_y_a.reshape(270))
                 accuracy = roc_auc_score(test_y_a, clf_rbf.predict(test_x_a).flatter
```

```
print(f"RBF Kernel - Parameters: C={C}, Gamma={gamma}, Accuracy: {ac
         if accuracy > best accuracy rbf:
             best_accuracy_rbf = accuracy
             best_params_rbf = {'C': C, 'gamma': gamma}
 print(f"Best Parameters for RBF Kernel: {best params rbf}")
 print(f"Best Accuracy for RBF Kernel: {best_accuracy_rbf:.2%}")
RBF Kernel - Parameters: C=0.1, Gamma=0.1, Accuracy: 50.00%
RBF Kernel - Parameters: C=0.1, Gamma=1, Accuracy: 55.83%
RBF Kernel - Parameters: C=0.1, Gamma=10, Accuracy: 72.50%
RBF Kernel - Parameters: C=1, Gamma=0.1, Accuracy: 50.00%
RBF Kernel - Parameters: C=1, Gamma=1, Accuracy: 89.50%
RBF Kernel - Parameters: C=1, Gamma=10, Accuracy: 92.42%
RBF Kernel - Parameters: C=10, Gamma=0.1, Accuracy: 67.92%
RBF Kernel - Parameters: C=10, Gamma=1, Accuracy: 94.17%
RBF Kernel - Parameters: C=10, Gamma=10, Accuracy: 89.92%
Best Parameters for RBF Kernel: {'C': 10, 'gamma': 1}
Best Accuracy for RBF Kernel: 94.17%
```

We observe that the best accuracy is achieved with the RBF kernel followed by the polynomial and lastly the linear which is to be expected. Also the linear kernel does not seem to improve with different parameters in contrast to the polynomial and the rbf.

```
In []:
```