Before you turn this problem in, make sure everything runs as expected. First, **restart the kernel** (in the menubar, select Kernel \rightarrow Restart) and then **run all cells** (in the menubar, select Cell \rightarrow Run All).

Make sure you fill in any place that says YOUR CODE HERE or "YOUR ANSWER HERE", as well as your name and collaborators below:

```
In [1]:

NAME = "Kanawut Kaewnoparat"
ID = "st122109"
```

Lab 05: Optimization Using Newton's Method

In this lab, we'll explore an alternative to gradient descent for nonlinear optimization problems: Newton's method.

Newton's method in one dimension

Consider the problem of finding the *roots* x of a nonlinear function $f: \mathbb{R}^N \to \mathbb{R}$. A root of f is a point x that satisfies f(x) = 0.

In one dimension, Newton's method for finding zeroes works as follows:

```
1. Pick an initial guess x_0
```

```
2. Let x_{i+1} = x_i + \frac{f(x_i)}{f(x_i)}
```

3. If not converged, go to #2.

```
Convergence occurs when |\mathit{I}(x_j)| \le \epsilon_1 or when |\mathit{I}(x_{j+1}) - \mathit{I}(x_j)| \le \epsilon_2
```

Let's see how this works in practice.

Example 1: Root finding for a cubic polynomial

Let's begin by using Newton's method to find roots of a simple cubic polynomial

```
f(x) = x^3 + x^2.
```

```
In [2]:
import matplotlib.pyplot as plt
import numpy as np
from mpl toolkits.mplot3d import Axes3D
import pandas as pd
Here's a function to evaluate a polynomial created with Numpy's polyld function at a particular point x
In [3]:
def fx(x, p):
    f_x = np.polyval(p, x)
    return f x
In [4]:
np.polyval([3,5,2], 2)
Out[4]:
24
np.polyval([2,1,5, 3], 2)
Out[5]:
33
In [6]:
np.polyval([3,5,2], 1)
Out[6]:
And here's some code to create the polynomial x^3 + x^2
, get its derivative, and evalute the derivative at 200 points along the \,x\,
axis::
```

```
In [7]:
print(np.poly1d([2,1,0,1]))
2 x + 1 x + 1
In [8]:
np.polyval([3,5,2], [1,2,4])
Out[8]:
array([10, 24, 70])
In [9]:
# Create the polynomial f(x) = x^3 + x^2
p = np.polyld([1, 1, 0, 0]) # [1 * x^3, 1 * x^2, 0 * x^1, 0 * 1]
# Get f'(x) (the derivative of f(x) in polynomial form) # We know it's 2x^2 + 2x, which is [3, 2, 0] in poly1d form
p_d = np.polyder(p)
print('f(x):')
print('----')
print(p)
print('----')
print("f'(x):")
print('----')
print(p_d)
print('----')
\# Get 200 points along the x axis between -3 and 3
n = 200
x = np.linspace(-3, 3, n)
\# Get values for f(x) and f'(x) in order to graph them later
y = fx(x, p)
y d = fx(x, p d)
f(x):
  3 2
1 x + 1 x
f'(x):
   2
3 x + 2 x
Next, let's try three possible guesses for x_0
: -3, 1, and 3, and in each case, run Newton's root finding method from that initial guess.
In [10]:
# Initial guesses
x0_arr = [-3.0, 1.0, 3.0]
# Parameters for Newton: number of iterations,
# threshold for identifying a point as a zero max iters = 30
\frac{-}{\text{threshold}} = 0.0001
# Set up plot
fig1 = plt.figure(figsize=(8,8))
ax = plt.axes()
plt.plot(x, y, 'g-', label='f(x)')
plt.plot(x, y_d, 'b--', label="f'(x)")
roots = []
for x0 in x0_arr:
     i = 0
     xi = x0
     fxi = fx(xi, p)
     # Plot initial data point
     plt.plot(xi, fxi, '*', label=("iter 0: x=%.4f" % x0))
     while i < max_iters:</pre>
         # x_i + 1 = x_i - f(x_i)/f'(x_i)

xi = xi - fx(xi, p) / fx(xi, p_d) #NEWTON'S METHOD
         fxi = fx(xi, p) # Plot (xi, fxi) and add a legend entry every 5 iterations if (i+1) % 5 == 0:
              plt.plot(xi, fxi, '.', label=("iter %d: x=%.4f" % (i+1, xi)))
         else:
              plt.plot(xi, fxi, '.')
         # Check if |f(x)| < threshold if np.abs(fxi) < threshold:
```

roots.append(xi)

plt.legend(bbox_to_anchor=(1.5, 1.0), loc='upper right')

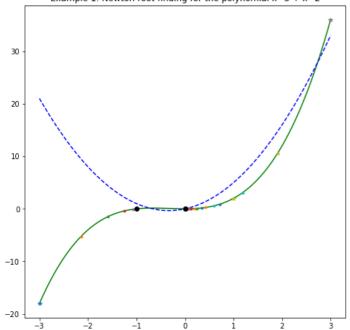
plt.plot(xi, fx(xi, p), 'ko', label=("converged at iter %d" % (i+1)))

plt.title('Example 1: Newton root finding for the polynomial $x^3 + x^2$ ')

 $\begin{array}{c} \textbf{break} \\ \textbf{i} = \textbf{i} + 1 \end{array}$

plt.show()

Example 1: Newton root finding for the polynomial $x^3 + x^2$



```
f(x)

f(x)

iter 0: x=-3.0000

iter 5: x=-1.0085

converged at iter 7

iter 0: x=1.0000

iter 5: x=0.0543

converged at iter 8

iter 0: x=3.0000

iter 5: x=0.2393

iter 10: x=0.0091

converged at iter 10
```

Example 2: Root finding for the sine function

```
Next, consider the function f(x) = \sin(x)
```

```
In [11]:

def fx_sin(x):
    return np.sin(x)

def fx_dsin(x):
    return np.cos(x)
```

Let's get 200 points in the range $[-\pi, \pi]$ for plotting:

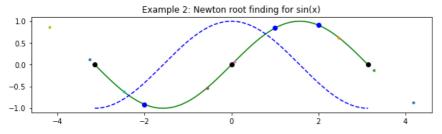
```
In [12]:
```

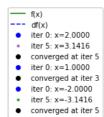
```
# Get f(x)=sin(x) and f'(x) at 200 points for plotting
n = 200
x = np.linspace(-np.pi, np.pi, n)
y = fx_sin(x)
y d = fx dsin(x)
```

In [13]:

```
# Initial guesses
x0_arr = [2.0, 1.0, -2.0]
# Parameters for Newton: number of iterations,
# threshold for identifying a point as a zero
max_iters = 30
threshold = 0.0001
# Set up plot
fig1 = plt.figure(figsize=(10,10))
ax = plt.axes()
ax.set_aspect(aspect='equal', adjustable='box')
plt.plot(x, y, 'g-', label='f(x)')
plt.plot(x, y_d, 'b--', label='df(x)')
roots = []
for x0 in x0_arr:
    i = 0;
    xi = x0
    fxi = fx_sin(xi)
     # Plot initial data point
    plt.plot(xi, fxi, 'bo', label=("iter 0: x=%.4f" % x0))
     while i < max_iters:</pre>
         # x_i + 1 = x_i - f(x_i)/f'(x_i)
xi = xi - fx_sin(xi) / fx_dsin(xi)
         fxi = fx_sin(xi)
         # Plot (xi, fxi) and add a legend entry every 5 iterations if (i+1) % 5 == 0:
              plt.plot(xi, fxi, '.', label=("iter %d: x=%.4f" % (i+1, xi)))
              plt.plot(xi, fxi, '.')
          \# Check if |f(x)| < threshold
         if np.abs(fxi) < threshold:</pre>
              roots.append(xi)
              break
```







Roots: 3.141593, -0.000096, -3.141593

Notice that we get some extreme values of x for some cases. For example, when $x_0 = -2$, where the slope is pretty close to 0, the next iteration gives a value less than -4.

Newton's method for optimization

Now, consider the problem of minimizing a scalar function $\mathit{J} : R^\mathit{n} \mapsto R$

. We would like to find

$$\theta^* = \operatorname{argmin}_{\theta} J(\theta)$$

We already know gradient descent:

$$\theta^{(i+1)} \leftarrow \theta^{(i)} - \alpha \nabla (\theta^{(i)}).$$

But Newton's method gives us a potentially faster way to find θ^* as a zero of the system of equations

$$\nabla_f(\theta^*) = \mathbf{0}.$$

In one dimension, to find the zero of f'(x)

- , obviously, we would apply Newton's method to f(x)
- , obtaining the iteration

$$x_{i+1} = x_i - f(x_i) / f''(x_i).$$

The multivariate extension of Newton's optimization method is

$$\mathbf{x}_{i+1} = \mathbf{x}_i - \mathbf{H}_f^{-1}(\mathbf{x}_i) \nabla_f(\mathbf{x}_i),$$

where $H_f(\mathbf{x})$ is the *Hessian* of f evaluated at \mathbf{x} :

$$\mathbf{H}_{f}(\mathbf{x}) = \begin{bmatrix} \frac{\partial^{2} f}{\partial x_{1}^{2}} & \frac{\partial^{2} f}{\partial x_{1} x_{2}} & \dots & \frac{\partial^{2} f}{\partial x_{1} x_{n}} \\ \frac{\partial^{2} f}{\partial x_{2} x_{1}} & \frac{\partial^{2} f}{\partial x_{2}^{2}} & \dots & \frac{\partial^{2} f}{\partial x_{2} x_{n}} \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\partial^{2} f}{\partial x_{n} x_{1}} & \frac{\partial^{2} f}{\partial x_{n} x_{2}} & \dots & \frac{\partial^{2} f}{\partial x_{n}^{2}} \end{bmatrix}$$

This means, for the minimization of $\mathcal{J}(\theta)$, we would obtain the update rule

$$\theta^{(i+1)} \leftarrow \theta^{(i)} - \mathbf{H}_I^{-1}(\theta^{(i)}) \nabla_f \theta^{(i)}$$
.

Application to logistic regression

Let's create some difficult sample data as follows:

Class 1: Two features x_1 and x_2

jointly distributed as a two-dimensional spherical Gaussian with parameters

$$\mu = \begin{bmatrix} x_{1c} \\ x_{2c} \end{bmatrix}, \Sigma = \begin{bmatrix} \sigma_1 & \sigma_1 \\ 0 & \sigma_1^2 \end{bmatrix}.$$

```
Class 2: Two features x_1
```

and x

in which the data are generated by first sampling an angle θ according to a uniform distribution, sampling a distance d according to a one-dimensional Gaussian with a mean of $(3\,\sigma_{\rm l})^2$

and a variance of $({}^2\sigma_1)^2$, then outputting the point

$$\mathbf{x} = \begin{bmatrix} x_{1c} + d\cos\theta \\ x_{2c} + d\sin\theta \end{bmatrix}.$$

Generate 100 samples for each of the classes, guided by the following exercises.

Exercise 1.1 (5 points)

Generate data for class 1 with 100 samples:

$$\mu = \begin{bmatrix} x_{1c} \\ x_{2c} \end{bmatrix}, \Sigma = \begin{bmatrix} \sigma_1^2 & 0 \\ 0 & \sigma_1^2 \end{bmatrix}.$$

► Hint:

```
In [14]:
mu_1 = np.array([1.0, 2.0])
sigma_1 =
num\_sample = 100
sigma vector = np.array([sigma 1, sigma 1])
cov mat = np.diag(sigma vector)
print(f"covariance matrix:\n {cov mat}")
print()
X1 = np.random.multivariate normal(mean = mu 1, cov = cov mat, size = num sample)
print(X1[:5])
X1.mean(axis=0)
# YOUR CODE HERE
# raise NotImplementedError()
covariance matrix:
 [[1 0]
 [0 1]]
[[ 0.93961715 2.09174582]
  0.41114958 2.48695983]
 [ 0.5398311
                3.226419071
 [-0.24316161 0.87464345]
 [ 0.67288163  0.98057817]]
Out[14]:
array([0.89755737, 1.84060126])
In [15]:
print(X1[:5])
# Test function: Do not remove
assert X1.shape == (100, 2), 'Size of X1 is incorrect'
assert cov_mat.shape == (2, 2), 'Size of x_test is incorrect'
count = 0
for i in range(2):
    for j in range(2):
         if i==j and cov_mat[i,j] != 0:
             if cov_mat[i,j] == sigma_1:
                 count +=
            if cov_mat[i,j] == 0:
assert count == 4, 'cov mat data is incorrect'
print("success!")
# End Test function
[[ 0.93961715 2.09174582]
 [ 0.41114958 2.48695983]
 [ 0.5398311
                3.22641907]
 [-0.24316161
               0.87464345]
 [ 0.67288163  0.98057817]]
success!
```

Expected result (or something similar):\ [[-0.48508229 2.65415886]\ [1.17230227 1.61743589]\ [-0.61932146 3.53986541]\ [0.70583088 1.45944356]\ [-0.93561505 0.2042285]]

Exercise 1.2 (5 points)

Generate data for class 2 with 100 samples:

[8.5494541 . 8.16187663].

```
\mathbf{x} = \begin{bmatrix} x_{1c} + d\cos\theta \\ x_{2c} + d\sin\theta \end{bmatrix}
```

```
where \theta
is sampled uniformly from [0, 2\pi]
is sampled from a one-dimensional Gaussian with a mean of (3\sigma_1)^2
and a variance of (2 \sigma_1)^2
► Hint:
In [16]:
np.pi
Out[16]:
3.141592653589793
In [17]:
 # 1. Create sample angle from 0 to 2pi with 100 samples
angle = np.random.uniform(0, 2*np.pi, size= num sample)
print(angle.shape)
# 2. Create sample with normal distribution of d with mean and variance
d = np.random.normal(loc = (3*sigma_1)**2, scale = (sigma_1 /2)**2, size =num_sample)
print(d.shape)
 # 3 Create X2
X2 = np.vstack((X1[:, 0] + (d * np.cos(angle)), X1[:, 1] + (d * np.sin(angle)))).T
 # YOUR CODE HERE
 # raise NotImplementedError()
(100,)
(100,)
In [18]:
X1[:, 0].shape
Out[18]:
(100,)
In [19]:
\texttt{np.vstack}((\texttt{X1[:, 0] - (d * np.cos(angle)), X1[:, 1] - (d * np.sin(angle))))}.\texttt{T}
Out[19]:
array([[ 2.96008865, -7.06612894],
             [-5.48175108, 8.9010257], [9.14287679, 0.73877222], [7.89868875, 4.87379543], [-7.60228533, 5.03025906], [6.33487392, 8.91623875],
                0.68080784, -6.70786588],
            [ 0.68080784, -6.70786588],
[-7.59470607, -1.5011998],
[-7.25443377, 3.78844097],
[ 5.38799381, -4.83738222],
[ 8.73332188, 8.55823444],
[-7.99497509, 2.44369681],
[ 8.35112927, 9.56425151],
[ 5.64041528, -5.38478839],
[-6.00294295, -3.13119899],
             [ 5.71505013, 10.75831802],
             [ 1.98128536, 11.89507898],
             [-7.82714798, -2.2790989],
             [ 3.070826 , -8.32430685],
[10.53823534 , 2.48982195],
[ 3.93647954 , 10.49708356],
            [-6.29442054, 6.63731663],
[-8.16166793, 1.56427131],
[-4.77078483, 8.73999081],
[-2.28672369, 11.1264415],
            [ 6.36392317, 9.00619679],
[ 9.80311426, -3.00724707],
[-1.91448507, 9.97443549],
            [-6.10000181, 5.72628506], [-4.56408541, 6.80233849], [-3.50828964, -6.83761767], [1.66319046, -7.05662394],
```

```
4.79984827, -6.92310068],
8.47556971, -4.41761017],
                1.47457415, 11.35696446],
             [-8.8064449 , -0.05024625],
             [-3.19222599, -6.26923287], [6.93778699, 8.89082898], [-4.0516323, -6.4752065],
             [-4.0516323 , -6.4752065 ],
[8.2223447 , 8.61843156],
[7.69870515 , 7.70331835],
[-4.57527108 , 7.63321615],
[-8.38435525 , 1.67318255],
[-1.12893806 , 8.47670486],
[7.91848835 , 6.2941658 ],
[1.39061144 , -7.87804416],
[8.95658259 , -2.99675471]
             [ 8.95658259, -2.99675471],
[10.71916783, 2.922222 ],
             [ 9.96650694, -0.99715112],
                3.36678814, -7.35659057],
             [-1.29727051, -6.02381103],
[-1.89966253, -6.1056425],
              [-5.74155946, 9.11803404],
[-8.84745901, 2.6033943],
                4.47796487, 8.73686196],
6.60975967, 8.3842056],
7.46008352, 3.57342453],
9.27301918, -2.85873306],
             [-1.50957143, -6.71848346],
                7.78053374, -3.30588691],
               -4.67476273, -5.89532808],
             [-2.21056537, -5.40678576],
                 6.98008082, -1.31691101],
                 1.85905295, -6.10904963],
                9.82158256, 3.45661419],
6.76404173, -3.93265668],
             [ 8.4182352 , -1.29316894],
[-2.30513179, 8.44716296],
                 7.76308979, -2.99808243],
6.56307976, -3.76381842],
             [ 8.6173021 , -3.91876901],
[ 2.59266962, -8.09747415],
            [ 2.59266962, -8.09747415],

[-1.05208527, -6.09779469],

[ 7.88250761, 8.17430896],

[ -9.39813043, 1.06528924],

[ 9.43857626, 2.3363296],

[ -8.07203751, -1.9304691],

[ 5.82128112, 9.38601624],

[ -3.41405792, 9.7525935],

[ 10.11511579, 4.5104464],

[ 10.91843046, -1.53802303],

[ 8.49961022, 4.05223443],

[ 5.36197447, -5.6762687],

[ 7.2853866, 9.87501112],

[ 0.28686481, -7.3834579],

[ 3.73685992, -6.76286351],
                3.73685992, -6.76286351],
7.69704719, -4.01015001],
             [-2.22176399, 11.01859379],
             [ 5.84625407, 7.81957616],
[-8.75984463, 3.0484368],
[ 5.85121394, -3.95730919],
             [-5.84688946, -1.96234631],
[-0.02772957, -7.83334652],
[5.31036742, -4.43395419],
             [-7.03039116, 8.11645926],
[-8.86955959, 1.81589595],
             [-7.166953 , -2.9941708 ]])
In [20]:
print('angle:',angle[:5])
print('d:', d[:5])
print('X2:', X2[:5])
 # Test function: Do not remove
assert angle.shape == (100,) or angle.shape == (100,1) or angle.shape == 100, 'Size of angle is incorrect'
assert d.shape == (100,) or d.shape == (100,1) or d.shape == 100, 'Size of d is incorrect'
assert X2.shape == (100,2), 'Size of X2 is incorrect'
assert angle.min() >= 0 and angle.max() <= 2*np.pi, 'angle generate incorrect'
assert d.min() >= 8 and d.max() <= 10, 'd generate incorrect'</pre>
assert X2[:,0].min() >= -13 and X2[:,0].max() <= 13, 'X2 generate incorrect'
assert X2[:,1].min() >= -10 and X2[:,1].max() <= 13.5, 'X2 generate incorrect'</pre>
# End Test function
angle: [1.7879443 5.45546526 2.8601113 3.59816315 5.82807172]
d: [9.37811149 8.71013887 8.95548893 9.07099465 9.21294217]
X2: [[-1.08085435 11.24962057]
  [ 6.30405024 -3.92710604]
[-8.06321458 5.71406592]
  [-8.38501197 -3.12450854]
  [ 8.94804858 -3.06910271]]
success!
```

Expected result (or something similar):\ angle: [4.77258271 3.19733552 0.71226709 2.11244845 6.06280915]\ d: [9.13908279 8.84218552 9.24427852 8.74831667 8.85727588]\ X2: [[0.064701 -6.46837219]\ [-7.65614929 1.12480234]\ [6.37750805 9.58147629]\ [-3.80438416 8.95550952]\ [7.70745021 -1.73194274]]

Exercise 1.3 (5 points)

In [21]:

Combine X1 and X2 into single dataset

```
np.concatenate((X1, X2)).shape
Out[21]:
(200, 2)
In [22]:
np.concatenate((([0] * 100), ([1]*100)))
Out[221:
0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1,
                                  1, 1, 1, 1,
                                           1, 1,
                                                1, 1,
     1, 1, 1, 1,
     1, 1])
In [23]:
# 1. concatenate X1, X2 together
X = np.concatenate((X1, X2))
# 2. Create y with class 1 as 0 and class 2 as 1 y = np.concatenate((([0] * 100), ([1]*100)))
# YOUR CODE HERE
# raise NotImplementedError()
In [24]:
#alternative method
\# X = np.concatenate([X1, X2], axis = 0)
\# y = np.append(np.zeros(num_sample), np.ones(num_sample))
In [25]:
print("shape of X:", X.shape)
print("shape of y:", y.shape)
# Test function: Do not remove
assert X.shape == (200, 2), 'Size of X is incorrect'
assert y.shape == (200,) or y.shape == (200,1) or y.shape == 200, 'Size of y is incorrect'
assert y.min() == 0 and y.max() == 1, 'class type setup is incorrect'
print("success!")
# End Test function
shape of X: (200, 2)
shape of y: (200,)
```

Expect result (or looked alike):\ shape of X: (200, 2)\ shape of y: (200, 1)

Exercise 1.4 (5 points)

Plot the graph between class1 and class2 with difference color and point style.

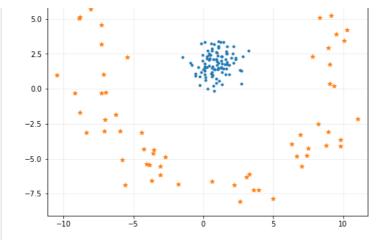
```
In [26]:

fig1 = plt.figure(figsize=(8,8))
ax = plt.axes()
plt.title('Sample data for classification problem')
plt.grid(axis='both', alpha=.25)

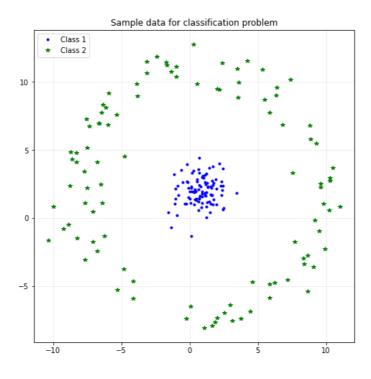
plt.scatter(X1[:, 0], X1[:, 1], marker = '.', label = 'class1')
plt.scatter(X2[:, 0], X2[:, 1], marker = '*', label = 'class2')

plt.legend()
plt.axis('equal')
plt.show()
```





Expect result (or looked alike):



Exercise 1.5 (5 points)

train_size = 0.8
m = len(X)

start_stop_idx = round(train_size * m)

idx_train = all_idx[: start_stop_idx]
idx_test = all_idx[start_stop_idx:]

Split data into training and test datasets with 80% of training set and 20% of test set

```
In [27]:
all_idx = np.arange(len(X))
np.random.seed(16)
np.random.shuffle(all idx)
In [28]:
all idx
Out[28]:
array([154,
               0, 174,
                         59, 112, 170,
                                         73,
                                                    29,
                                                          32,
                                                               90,
                         72,
       134,
              56,
                   81,
                             187,
                                   51, 100,
                                               79, 171,
                                                         113,
                                                              118,
                                                                     47,
                                                                          52,
                       105,
       180,
              39,
                  135,
                             124,
                                    19,
                                         71,
                                               99, 153,
                                                         189,
                                                                9,
                                                                    142,
                                                                         155,
       156,
              45,
                  195,
                        178,
                             104, 190,
                                        183,
                                               82,
                                                   166,
                                                          20,
                                                                     83,
                                                                         131,
                                    11,
         61,
              77, 181,
                         48, 102,
                                         53,
                                                2,
                                                    42,
                                                               60,
                                                                   193,
                                                                          87,
       117,
            157, 122,
                        167,
                              66, 163,
                                       149,
                                               26,
                                                   147,
                                                          78,
                                                               35,
                                                                   126,
                                                                          13,
        63,
               5,
                   75,
                         70,
                             136,
                                    55,
                                         50,
                                              141,
                                                    94,
                                                          95, 103,
                                                                   132,
       194,
            182, 197,
                         49,
                             172,
                                    76,
                                        140,
                                               34,
                                                    89,
                                                         138,
                                                               25,
                                                                     10,
                                                                   109,
       133,
              30,
                   84, 111,
                              17, 164, 173,
                                              161,
                                                     3,
                                                          68,
                                                               98,
                                                                         179,
         86, 176,
                  151,
                         57,
                              33,
                                    38, 198,
                                               97,
                                                  137,
                                                              139,
         91, 150, 188,
                       130, 192,
                                    36,
                                         80,
                                               27, 143,
                                                          22, 184,
                                                                    23,
                                                                          92,
             24, 129,
                         12,
                             199,
                                     6,
                                         46, 191, 119, 107,
                                                               88,
                                                                          44,
       148,
             18,
                   21,
                        16, 168, 186,
                                         15, 85,
                                                    37,
                                                          14, 108, 127,
                                                                          40,
       145, 165,
                    8, 115, 114, 175,
                                         28, 101, 128,
                                                          93, 162,
       196, 160, 116, 146, 110,
                                   58, 120, 158, 144,
                                                                4, 159,
       123,
                   69, 121, 169])
In [29]:
```

```
X train = X[idx train]
X_{\text{test}} = X[idx_{\text{test}}]
y_train = y[idx_train]
y_{test} = y[idx_{test}]
# YOUR CODE HERE
# raise NotImplementedError()
In [30]:
print('idx_train:', idx_train[:10])
print("train size, X:", X_train.shape, ", y:", y_train.shape)
print("test size, X:", X_test.shape, ", y:", y_test.shape)
# Test function: Do not remove
assert X_train.shape == (160, 2), 'Size of X_train is incorrect'
assert y_train.shape == (160,) or y_train.shape == (160,1) or y.shape == 160, 'Size of y_train is incorrect' assert X_test.shape == (40, 2), 'Size of X_test is incorrect'
assert y_test.shape == (40,) or y_test.shape == (40,1) or y.shape == 40, 'Size of y_test is incorrect'
print("success!")
# End Test function
idx_train: [154  0 174  59 112 170  73  7  29  32]
train size, X: (160, 2) , y: (160,)
test size, X: (40, 2) , y: (40,)
success!
Expected reult (or something similar):\idx_train: [ 78 61 28 166 80 143 6 76 98 133]\ train size, X: (160, 2), y: (160, 1) \ test size, X: (40, 2), y: (40, 1)
Exercise 1.6 (5 points)
Write a function to normalize your X
data
Practice yourself (No grade, but has extra score 3 points)
Try to use Jupyter notebook's LaTeX equation capabilities to write the normalization equations for your dataset.
Normalization Function
                                                                           \bar{X} = \frac{X - \mu}{\sigma}
```

where $\bar{\boldsymbol{X}}$ is the output normalizedX is the original input

X is the original input X,

is the mean,

and σ

is the standard deviation

```
In [31]:
X.mean(axis= 0)
Out[31]:
array([0.5046732 , 2.14205036])
In [32]:
X.std(axis= 0)
Out[32]:
array([4.49453967, 4.65625837])
In [33]:
a = (X - X.mean(axis= 0)) / X.std(axis= 0)
In [34]:
a.mean(axis=0)
Out[34]:
array([-4.32986980e-17, -9.76996262e-17])
In [35]:
a.std(axis=0)
Out[35]:
array([1., 1.])
In [36]:
def normalization(X):
```

```
X norm = (X - X mean) / X std
     return X_norm
      # YOUR CODE HERE
        raise NotImplementedError()
In [37]:
X norm = normalization(X)
In [38]:
X \text{ norm.mean}(axis = 0)
Out[38]:
array([-4.32986980e-17, -9.76996262e-17])
In [39]:
X norm[:, 0].min()
Out[39]:
-2.442884229663548
In [40]:
X norm[:, 0].max()
Out[40]:
2.349357093561492
In [41]:
XX = normalization(X)
X train norm = XX[idx train]
X_test_norm = XX[idx_test]
# Add 1 at the first column of training dataset (for bias) and use it when training
X_design_train = np.insert(X_train_norm,0,1,axis=1)
X_design_test = np.insert(X_test_norm,0,1,axis=1)
m,n = X design train.shape
print(X_train_norm.shape)
print(X_design_train.shape)
print(X_test_norm.shape)
print(X_design_test.shape)
# Test function: Do not remove assert XX[:,0].min() >= -2.5 and XX[:,0].max() <= 2.5, 'Does the XX is normalized?' assert XX[:,1].min() >= -2.5 and XX[:,1].max() <= 2.5, 'Does the XX is normalized?'
print("success!")
# End Test function
(160, 2)
 (160, 3)
 (40, 2)
 (40, 3)
success!
```

Take in numpy array of X values and return normalize X values,

the mean and standard deviation of each feature

X_mean = np.mean(X, axis= 0)
X_std = np.std(X, axis = 0)

Exercise 1.7 (10 points)

define class for logistic regression: batch gradient descent

The class includes:

• Sigmoid function

$$sigmoid(z) = \frac{1}{1 + e^{-z}}$$

Softmax function

$$softmax(z) = \frac{e^{Z_i}}{\sum_{n} e^{z_n}}$$

• Hyperthesis (h) function

$$\hat{y} = h(X; \theta) = softmax(\theta. X)$$

• Gradient (Negative likelihood) function

$$gradient = -X. \frac{y - \hat{y}}{n}$$

· Cost function

$$\sum ((-y\log^{\hat{y}}) - ((1-y)\log(1-\hat{y})))$$

$$cost = n$$

- · Gradient ascent function
- Prediction function
- · Get accuracy funciton

```
In [42]:
```

```
class Logistic BGD:
    def __init__(self):
        pass
    def sigmoid(self,z):
        s = 1 / (1 + np.exp(-1 * z))
        return s
    def softmax(self, z):
        sm = np.exp(z) / np.exp(z).sum()
        return sm
    def h(self, X, theta):
        hf = np.dot(X, theta)
        yhat = self.sigmoid(hf)
        return yhat
    def gradient(self, X, y, y pred):
        error = (y - y_pred) / len(y)
grad = -1 * (np.dot(X.T, error))
        return grad
    def costFunc(self, theta, X, y):
        yhat = self.h(X, theta)
error = -1*(y * np.log(yhat)) - ((1-y)*np.log(1-yhat))
cost = np.sum(error) / len(yhat)
        grad = self.gradient(X, y , yhat)
        return cost, grad
    def gradientAscent(self, X, y, theta, alpha, num_iters):
        m = len(y)
        J_history = []
         theta_history = []
         for i in range(num_iters):
             # 1. calculate cost, grad function
             cost, grad = self.costFunc(theta, X,y)
             # 2. update new theta
theta = theta - (grad * alpha)
             J_history.append(cost)
             theta_history.append(theta)
        J_min_index = np.argmin(J_history)
print("Minimum at iteration:",J_min_index)
        return theta_history[J_min_index] , J_history
    def predict(self, X, theta):
         labels=[]
         # 1. take y_predict from hyperthesis function
         # 2. classify y predict that what it should be class1 or class2
# 3. append the output from prediction
         for i in range(X.shape[0]):
             y1= self.h(X[i], theta)
             if y1 >= 0.5:
                  labels.append(1)
             else:
                  labels.append(0)
        labels=np.asarray(labels)
         return labels
    def getAccuracy(self, X, y, theta):
        y = y.ravel()
         yhat = self.predict(X, theta)
         correct_match = np.sum(yhat == y)
        percent_correct = (correct_match / len(y)) * 100
        return percent_correct
```

In [43]:

```
# Test function: Do not remove
lbgd = Logistic_BGD()
test_x = np.array([[1,2,3,4,5]]).T
out_x1 = lbgd.sigmoid(test_x)
out_x2 = lbgd.sigmoid(test_x.T)
print('out_x1', out_x1.T)
assert np.array_equal(np.round(out_x1.T, 5), np.round([[0.73105858, 0.88079708, 0.95257413, 0.98201379, 0.99330715]],
5)), "sigmoid function is incorrect"
```

```
assert np.array_equal(np.round(out_x2, 5), np.round([[0.73105858, 0.88079708, 0.95257413, 0.98201379, 0.99330715]], 5)
), "sigmoid function is incorrect"
out x1 = lbgd.softmax(out x1)
out x2 = lbgd.softmax(out x2)
print('out_x1', out_x1.T)
assert np.array_equal(np.round(out_x1.T, 5), np.round([[0.16681682, 0.19376282, 0.20818183, 0.21440174, 0.21683678]],
5)), "softmax function is incorrect
assert np.array_equal(np.round(out_x2, 5), np.round([[0.16681682, 0.19376282, 0.20818183, 0.21440174, 0.21683678]], 5)
), "softmax function is incorrect'
test_t = np.array([[0.3, 0.2]]).T
test x = np.array([[1,2,3,4,5,6],[2,9,4,3,1,0]]).T
test_y = np.array([[0,1,0,1,0,1]]).T
test_y_p = lbgd.h(test_x, test_t)
print('test_y_p', test_y_p.T)
assert np.array equal(np.round(test y p.T, 5), np.round([[0.66818777, 0.9168273, 0.84553473, 0.85814894, 0.84553473, 0
.85814894]], 5)), "hyperthesis function is incorrect"
test_g = lbgd.gradient(test_x, test_y, test_y_p)
print('test_g', test_g.T)
assert np.array_equal(np.round(test_g.T, 5), np.round([[0.9746016, 0.73165696]], 5)), "gradient function is incorrect"
test_c, test_g = lbgd.costFunc(test_t, test_x, test_y)
print('test c', test c.T)
assert np.round(test_c, 5) == np.round(0.87192491, 5), "costFunc function is incorrect"
test_t_out , test_j = lbgd.gradientAscent(test_x, test_y, test_t, 0.001, 3)
print('test_t_out', test_t_out.T)
print('test_j', test_j)
assert np.array_equal(np.round(test_t_out.T, 5), np.round([[0.29708373, 0.19781153]], 5)), "gradientAscent function is
incorrect"
assert np.round(test_j[2], 5) == np.round(0.86896665, 5), "gradientAscent function is incorrect"
test 1 = lbgd.predict(test_x, test_t)
print('test_l', test_l)
assert np.array_equal(np.round(test_1, 1), np.round([1,1,1,1,1,1], 1)), "gradientAscent function is incorrect"
test_a = lbgd.getAccuracy(test_x,test_y,test_t)
print('test_a', test_a)
assert np.round(test_a, 1) == 50.0, "getAccuracy function is incorrect"
print("success!")
# End Test function
out_x1 [[0.73105858 0.88079708 0.95257413 0.98201379 0.99330715]]
out_x1 [[0.16681682 0.19376282 0.20818183 0.21440174 0.21683678]]
test_y_p [[0.66818777 0.9168273 0.84553473 0.85814894 0.84553473 0.85814894]]
test_g [[0.9746016 0.73165696]]
test c 0.8719249134773479
Minimum at iteration: 2
test_t_out [[0.29708373 0.19781153]]
       [0.8719249134773479, 0.870441756946089, 0.8689666485816598]
test_l [1 1 1 1 1 1]
```

 $\begin{tabular}{ll} \textbf{Expected result:} \out_x1 & [[0.73105858 0.88079708 0.95257413 0.98201379 0.99330715]] \out_x1 & [[0.16681682 0.19376282 0.20818183 0.21440174 0.21683678]] \out_x1 & [[0.66818777 0.9168273 0.84553473 0.85814894 0.84553473 0.85814894]] \out_x1 & [[0.9746016 0.73165696]] \out_x1 & [[0.9746016 0.7316599]] \o$

Exercise 1.8 (5 points)

Training the data using Logistic_BGD class.

- Input: X_design_train
- Output: y_train

alpha = 0.001

test_a 50.0 success!

• Use 50,000 iterations

Find the initial_theta yourself

```
In [44]:
m, n = X design train.shape
In [45]:
n
Out[45]:
3
In [46]:
y train
Out[46]:
array([1, 0, 1, 0, 1, 1, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 1, 0, 1, 0, 1,
       1, 1, 0, 0, 1, 0, 1, 1, 1, 0, 0, 0, 1, 1, 0, 1, 1, 1, 0, 1, 1, 1,
       1, 1, 0, 1, 0, 1, 0, 1, 0, 0, 1, 0, 1, 0, 0, 0, 0, 0, 0, 1, 0, 1,
       1, 1, 1, 0, 1, 1, 0, 1, 0, 0, 1, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 0,
       1, 1, 0, 1, 1, 1, 0, 1, 0, 1, 0, 0, 1, 0, 0, 1, 0, 0, 1, 0, 1,
       1, 1, 0, 0, 0, 1, 1, 0, 1,
                                  1, 0, 0, 0,
                                               1,
                                                  0, 1, 0,
       1, 1, 1, 0, 0, 0, 1, 0, 1, 0, 0, 0, 0,
       1, 0, 1, 0, 0, 0])
In [47]:
```

```
iterations = 50000

BGD_model = Logistic_BGD()
initial_theta = np.zeros(n)
bgd_theta, bgd_cost = BGD_model.gradientAscent(X_design_train, y_train, initial_theta, alpha, iterations)

# YOUR CODE HERE
# raise NotImplementedError()
Minimum at iteration: 49999
```

```
In [48]:
```

```
bgd cost[np.argmin(bgd cost)]
Out[48]:
0.6679280712052752
```

```
In [49]:

np.argmin(bgd cost)
```

Out[49]:

49999

In [50]:

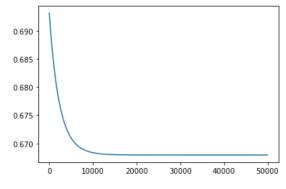
```
print(bgd_theta)
print(len(bgd_cost))

print(bgd_cost[0])
plt.plot(bgd_cost)
plt.show()

# Test function: Do not remove
assert bgd_theta.shape == (X_train.shape[1] + 1,1) or bgd_theta.shape == (X_train.shape[1] + 1,) or bgd_theta.shape == X_train.shape[1] + 1, "theta shape is incorrect"
assert len(bgd_cost) == iterations, "cost data size is incorrect"

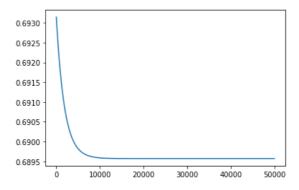
print("success!")
# End Test function
```

```
[-0.17665893 -0.24858832 0.38691536]
50000
0.6931471805599453
```



success!

Expected result (or look alike): \ [[-0.07328673]\ [-0.13632896]\ [0.05430939]]\ 50000



In lab exercises

- 1. Verify that the gradient descent solution is correct. Plot the optimal decision boundary you obtain.
- 2. Write a new class that uses Newton's method for the optmization rather than simple gradient descent.
- 3. Verify that you obtain a similar solution with Newton's method. Plot the optimal decision boundary you obtain.
- 4. Compare the number of iterations required for gradient descent vs. Newton's method. Do you observe other issues with Newton's method such as a singular or nearly singular Hessian matrix?

```
Plot the optimal decision boundary of gradient ascent
In [51]:
test2 = Logistic BGD()
In [52]:
y test
Out[52]:
In [53]:
X test[y test ==0]
Out[53]:
array([[ 1.05221555, 3.06298781], [ 0.28616858, -0.01488312],
         [ 0.7240154 , 2.09116226], [ 1.57829571, 1.05445336],
         [ 0.60630287, 1.4932225 ], [ 0.75026607, -0.13935944],
           1.85090771,
                          1.03255797],
          0.09793746,
                          3.35395429],
                          2.21044956],
           1.99830485,
         0.24511715,
                          1.55425676],
                          2.28283045],
         [-1.50744168,
         [ 2.32762906,
                           2.12881803],
          0.67288163,
                          0.98057817],
         [ 2.13696742,
                          3.016733511
         [ 1.32898485,
                          0.30462115]])
In [54]:
# Retrieve the model parameters.
b = bgd_theta[0]
w1 = bgd_theta[1]
w2 = bgd\_theta[2]
# Calculate the intercept and gradient of the decision boundary.
c = b/w2
m = w1/w2
# Plot the data and the classification with the decision boundary.
xmin, xmax = -15, 15
ymin, ymax = -15, 15
xd = np.array([xmin, xmax])
yd = m*xd + c
plt.plot(xd, yd, 'k', lw=1, ls='--')
plt.fill_between(xd, yd, ymin, color='tab:blue', alpha=0.2)
plt.fill_between(xd, yd, ymax, color='tab:orange', alpha=0.2)
plt.scatter(X_test[y_test ==0][:, 0], X_test[y_test ==0][:, 1], marker = '.', label = 'class1')
plt.scatter(X_test[y_test ==1][:, 0], X_test[y_test ==1][:, 1], marker = '*', label = 'class2')
plt.xlim(xmin, xmax)
plt.ylim(ymin, ymax)
plt.ylabel(r'$x 2$')
plt.xlabel(r'$x_1$')
plt.legend()
plt.show()
    15
                                                 dass1
    10
     5
     0
    -5
   -10
   -15 +
-15
              -10
                      -5
                                       5
                                              10
                               0
                                                      15
In [55]:
from sklearn.metrics import classification report
In [56]:
print(classification report(y true =y test , y pred = BGD model.predict(X design test,bgd theta) ))
```

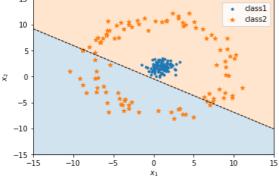
```
int(classification report(y true = y test , y pred = BGD model.predict(X design test,bgd theta) ))

precision recall f1-score support

0 0.44 1.00 0.61 15
1 1.00 0.24 0.39 25
```

```
0.72
                                     0.62
                                                  0.50
                                                                 40
   macro avg
                                     0.53
                                                                 40
                                                  0.47
weighted avg
In [57]:
xd * m
Out[57]:
array([ 9.63731389, -9.63731389])
In [58]:
w1/w2
Out[58]:
-0.6424875926047316
In [59]:
bgd_theta[0] / bgd_theta[2]
Out[59]:
-0.4565828902414966
In [60]:
# Retrieve the model parameters.
b = bgd theta[0]
w1 = bgd_theta[1]
w2 = bgd_theta[2]
# Calculate the intercept and gradient of the decision boundary.
c = b/w2
m = w1/w2
\# w[1] * y = w[0] * x + b
# # to solve for y
\# y = (w[0] * x)/w[1] + b / w[1]
# Plot the data and the classification with the decision boundary.
xmin, xmax = -15, 15

ymin, ymax = -15, 15
xd = np.array([xmin, xmax])
yd = (m*xd) + c
plt.plot(xd, yd, 'k', lw=1, ls='--')
plt.fill_between(xd, yd, ymin, color='tab:blue', alpha=0.2)
plt.fill_between(xd, yd, ymax, color='tab:orange', alpha=0.2)
plt.scatter(X1[:, 0], X1[:, 1], marker = '.', label = 'class1')
plt.scatter(X2[:, 0], X2[:, 1], marker = '*', label = 'class2')
plt.xlim(xmin, xmax)
plt.ylim(ymin, ymax)
plt.ylabel(r'$x_2$')
plt.xlabel(r'$x_1$')
plt.legend()
plt.show()
     15
                                                    dass1
                                                   dass2
     10
```



In [61]:

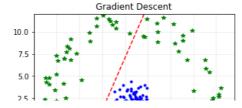
асситасу

BGD model.getAccuracy(X design test,y test,bgd theta)

Out[61]:

52.5

Expected result (or look alike): \



```
-2.5

-5.0

-7.5

-10.0 -7.5 -5.0 -2.5 0.0 2.5 5.0 7.5 10.0
```

```
In [62]:
```

```
print("Accuracy =",BGD model.getAccuracy(X design test,y test,bgd theta))
Accuracy = 52.5
```

```
Exercise 2.1 (10 points)
Write Newton's method class
In [63]:
У
Out[63]:
0, 0, 0,
     0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1,
     1, 1])
In [64]:
class Logistic NM: #logistic regression for newton's method
   def
       _init__(self):
      pass
   def sigmoid(self,z):
     s = 1 / (1 + np.exp(-z))
      return s
   def h(self, X, theta):
      hf = np.dot(X, theta)
      yhat = self.sigmoid(hf)
      return yhat
   def gradient(self, X, y, y_pred):
      m = len(y)
      grad = (1/m) * np.dot(X.T, (y_pred - y))
      return grad
   def hessian(self, X, y, theta):
    yhat = self.h(X, theta) #sigmoid output
     s = yhat * (1-yhat)
s = s.reshape(-1, 1)
hess_mat = X.T @ ( X * s)
      return hess_mat
   def costFunc(self, theta, X, y):
      m = len(y)
      yhat = self.h(X, theta)
      error = (y * np.log(yhat)) + ((1-y) * np.log(1-yhat))
cost = (-1/m) * np.sum(error)
      grad = self.gradient(X, y, yhat)
      return cost, grad
```

def newtonsMethod(self, X, y, theta, num_iters):

yhat = self.h(X, theta)
cost, grad = self.costFunc(theta, X , y)
hess mat = self.hessian(X, y , theta)

theta = theta - (np.dot(np.linalg.inv(hess_mat) , grad))

for i in range(num_iters):

print(i)
print(cost)

J_history.append(cost)
theta history.append(theta)

m = len(y)
J_history = []
theta history = []

```
print (grad)
                print (hess mat)
                print(theta)
          J_min_index = np.argmin(J_history)
         print("Minimum at iteration:", J_min_index)
          return theta history[J min index] , J history
     def predict(self, X, theta):
         y_pred = self.h(X, theta)
          labels=[]
         for i in range(X.shape[0]):
              if y_pred[i] >= 0.5:
                   Tabels.append(1)
                   labels.append(0)
          labels=np.asarray(labels)
         return labels
     def getAccuracy(self, X, y, theta):
         y = y.ravel()
         predicted list = self.predict(X, theta)
          total correct = np.sum(predicted_list == y)
         m = len(predicted_list)
print(predicted_list)
            print(total_correct)
            print(m)
         percent\_correct = (total\_correct / m) *100
         return percent correct
In [65]:
test_x.T @ (test_x * (test_y_p * (1- test_y_p)))
Out[65]:
array([[11.29725396, 5.49706551],
        [ 5.49706551, 10.37936818]])
In [66]:
test y p.shape
Out[66]:
(6, 1)
In [67]:
 # Test function: Do not remove
lbgd = Logistic NM()
test x = np.array([[1,2,3,4,5]]).T
out x1 = lbgd.sigmoid(test x)
out_x2 = lbgd.sigmoid(test_x.T)
print('out x1', out x1.T)
assert np.array equal(np.round(out x1.T, 5), np.round([[0.73105858, 0.88079708, 0.95257413, 0.98201379, 0.99330715]],
5)), "sigmoid function is incorrect
assert np.array_equal(np.round(out_x2, 5), np.round([[0.73105858, 0.88079708, 0.95257413, 0.98201379, 0.99330715]], 5)
), "sigmoid function is incorrect'
test t = np.array([[0.3, 0.2]]).T
test x = \text{np.array}([[1,2,3,4,5,6], [2, 9, 4, 3, 1, 0]]).T
test_x = np.array([[0,1,0,1,0,1]]).T

test_y = np.array([[0,1,0,1,0,1]]).T

test_y_p = lbgd.h(test_x, test_t)

print('test_y_p', test_y_p.T)
assert np.array_equal(np.round(test_y_p.T, 5), np.round([[0.66818777, 0.9168273, 0.84553473, 0.85814894, 0.84553473, 0.85814894]], 5)), "hyperthesis function is incorrect"
test_g = lbgd.gradient(test_x, test_y, test_y_p)
print('test_g', test_g.T)
assert np.array equal(np.round(test g.T, 5), np.round([[0.9746016, 0.73165696]], 5)), "gradient function is incorrect"
test_h = lbgd.hessian(test_x, test_y, test_t)
print('test_h', test_h)
assert test_h.shape == (2, 2), "hessian matrix function is incorrect"
assert np.array equal(np.round(test h.T, 5), np.round([[12.17334371, 6.55487738],[6.55487738, 14.84880387]], 5)), "he ssian matrix function is incorrect"
test_c, test_g = lbgd.costFunc(test_t, test_x, test_y)
print('test_c', test_c.T)
assert np.round(test_c, 5) == np.round(0.87192491, 5), "costFunc function is incorrect"
test_t_out, test_j = lbgd.newtonsMethod(test_x, test_y, test_t, 3)
print('test_t_out', test_t_out.T)
print('test_j', test_j)
assert np.array_equal(np.round(test_t_out.T, 5), np.round([[0.14765747, 0.15607017]], 5)), "newtonsMethod function is
incorrect"
assert np.round(test_j[2], 5) == np.round(0.7534506190845247, 5), "newtonsMethod function is incorrect"
test_l = lbgd.predict(test_x, test_t)
print('test_l', test_l)
\textbf{assert np.array\_equal(np.round(test\_1, 1), np.round([1,1,1,1,1,1], 1)), "gradientAscent function is incorrect"}
test_a = lbgd.getAccuracy(test_x,test_y,test_t)
print('test_a', test_a)
assert np.round(test_a, 1) == 50.0, "getAccuracy function is incorrect"
```

print("success!")
End Test function

Expect result: out_x1 [[0.73105858 0.88079708 0.95257413 0.98201379 0.99330715]]\ test_y_p [[0.66818777 0.9168273 0.84553473 0.85814894 0.84553473 0.85814894]]\ test_g [[0.9746016 0.73165696]]\ test_h [[12.17334371 6.55487738]\ [6.55487738 14.84880387]]\ test_c 0.8719249134773479\ Minimum at iteration: 2\ test_t_out [[0.14765747 0.15607017]]\ test_j [0.8719249134773479, 0.7967484437157274, 0.7534506190845247]\ test_l [1 1 1 1 1 1]\ test_a 50.0

```
In [68]:

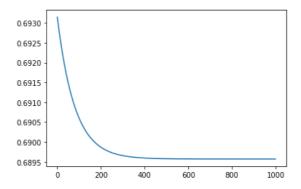
NM_model = Logistic_NM()
iterations = 1000

nm_theta, nm_cost = NM_model.newtonsMethod(X_design_train, y_train, initial_theta, iterations)
print("theta:",nm_theta)

print(nm_cost[0])
plt.plot(nm_cost)
plt.show()

Minimum at iteration: 999
theta: [-0.17631047 -0.24806574  0.38613374]
0.6931471805599454
0690
0685
```

Expected result (or look alike): \ Minimum at iteration: 999\ theta: [[-0.07313861]\ [-0.13605172]\ [0.05419746]]\ 0.6931471805599453



Exercise 2.2 (5 points)

0.675

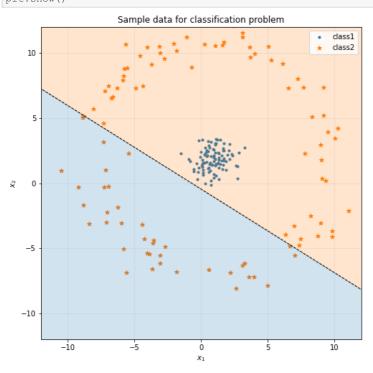
0.670

Plot the optimal decision boundary of Newton method

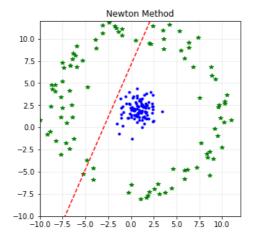
```
In [69]:
nm theta
Out[69]:
array([-0.17631047, -0.24806574,  0.38613374])
In [70]:
X[y==0][: , 0].shape
Out[70]:
(100,)
```

```
±n [/±]:
# plot
fig1 = plt.figure(figsize=(8,8))
ax = plt.axes()
plt.title('Sample data for classification problem')
plt.grid(axis='both', alpha=.25)
plt.scatter(X[y==0][: , 0], X[y==0][: , 1], marker = '.', label = 'class1')
plt.scatter(X[y==1][: , 0], X[y==1][: , 1], marker = '*', label = 'class2')
b = nm_theta[0]
w1 = nm_theta[1]
w2 = nm_theta[2]
# Calculate the intercept and gradient of the decision boundary.
c = b/w2

m = w1/w2
\# w[1] * y = w[0] * x + b
# # to solve for y
# y = (w[0] * x)/w[1] + b / w[1]
# Plot the data and the classification with the decision boundary.
xmin, xmax = -15, 15
ymin, ymax = -15, 15
# xd = np.array([xmin, xmax])
xd = np.linspace(-15, 15)
yd = (m*xd) + c
plt.plct(xd, yd, 'k', lw=1, ls='--')
plt.fill_between(xd, yd, ymin, color='tab:blue', alpha=0.2)
plt.fill_between(xd, yd, ymax, color='tab:orange', alpha=0.2)
plt.xlim(-12, 12)
plt.ylim(-12, 12)
plt.ylabel(r'$x_2$')
plt.xlabel(r'$x_1$')
plt.legend()
plt.show()
```



Expected result (or look alike):



```
In [72]:
print("Accuracy =",NM model.getAccuracy(X design test,y test,bgd theta))
Accuracy = 52.5

In [73]:
print(np.argmin(bgd_cost))
print(bgd_cost[np.argmin(bgd_cost)])
print()
print(np.argmin(nm_cost))
print(bgd_cost[np.argmin(nm_cost)])

49999
0.6679280712052752
999
0.6842329412851584
```

Exercise 2.3 (5 points)

Compare the number of iterations required for gradient descent vs. Newton's method. Do you observe other issues with Newton's method such as a singular or nearly singular Hessian matrix?

- For gradient descent method, it takes as many as almost 50,000 iterations to achieve the lowest cost at around 0.69
- As for the Newton's method, the number of iterations dwarves to only around 1,000 times to achieve almost the same cost at 0.69
- . This means that considering the computational expense, the Newton's method is way more effective than gradient descent

Take-home exercises

- 1. Perform a polar transformation on the data above to obtain a linearly separable dataset. (5 points)
- 2. Verify that you obtain good classification accuracy for logistic regression with GD or Netwon's method after the polar transformation (10 points)
- 3. Apply Newton's method to the dataset you used for the take home exercises in Lab 03. (20 points)

The report

Tn [79].

Write a brief report covering your experiments (both in lab and take home) and submit the Jupyter notebook via JupyterHub at https://puffer.cs.ait.ac.th before the next lab.

In your solution, be sure to follow instructions!

Perform the polar transformation on normalized data!!

```
In [74]:
X norm.mean(axis= 0)
Out[74]:
array([-4.32986980e-17, -9.76996262e-17])
In [75]:
angles = np.arctan((X_norm[:, 1] / X_norm[:, 0])) \#X1 over X0 radius = np.sqrt(X_norm[:, 0]**2 + X_norm[:, 1]**2)
print(angles[:5])
print(radius[:5])
[-0.11118019 -1.29694349 1.53721989 1.02212267 -1.42187275]
[0.09737283 0.07694153 0.23301547 0.31902115 0.25223514]
In [76]:
angles.shape
Out[76]:
(200,)
In [77]:
radius.shape
Out[77]:
(200,)
In [78]:
newX = np.vstack((angles, radius)).T
newX.shape
Out[78]:
(200, 2)
```

```
. زنان بند
newXtrain = newX[idx_train]
newXtest = newX[idx Test]
In [80]:
newXtrain = np.insert(newXtrain, 0 ,1 ,axis =1)
newXtest = np.insert(newXtest, 0 ,1 ,axis =1)
In [81]:
newXtrain.shape
Out[81]:
(160, 3)
In [82]:
plt.scatter(newX[y == 0][:, 0], newX[y == 0][:, 1], label = 'class 1')
plt.scatter(newX[y == 1][:, 0], newX[y == 1][:, 1], label = 'class 2')
plt.xlabel('angles')
plt.ylabel('radius')
plt.legend()
Out[82]:
<matplotlib.legend.Legend at 0x7f9fc79ebe20>
  2.5
   2.0
  1.5
                                            class 1
                                            class 2
   1.0
  0.5
   0.0
       -1.5
              -1.0
                    -0.5
                                  0.5
                                        1.0
In [83]:
alpha = 0.001
iterations = 50000
m,n = newXtrain.shape
polar_BGD_model = Logistic_BGD()
initial_theta = np.zeros(n)
polar bgd theta, polar bgd cost = polar BGD model.gradientAscent(newXtrain, y train, initial theta, alpha, iterations)
Minimum at iteration: 49999
In [84]:
polar_NM_model = Logistic_NM()
iterations = 1000
polar nm theta, polar nm cost = polar NM model.newtonsMethod(newXtrain, y train, initial theta, iterations)
Minimum at iteration: 999
In [85]:
plt.plot(np.arange(1000), polar_nm_cost , label = 'Newton Method')
plt.plot(np.arange(1000), polar bgd cost[:1000] , label = 'Batch Gradient Descent' )
plt.legend()
Out[85]:
<matplotlib.legend.Legend at 0x7f9fc7943cd0>
 0.6
 0.5
 0.4
                               Newton Method
                               Batch Gradient Descent
 0.3
 0.2
 0.1
 0.0
                                     800
                                             1000
In [86]:
y train
```

Out[86]:

```
1, 1,
       1, 1, 0, 1, 0, 1, 0, 1, 0, 0, 1, 0, 1,
                                                0, 0, 0, 0,
                                                             0, 0, 1, 0, 1,
       1, 1, 1, 0, 1, 1, 0, 1, 0, 0, 1, 0, 0,
                                                0, 0, 0, 1,
                                                             0, 0, 1, 0, 0,
       1, 1, 0, 1, 1, 1, 0, 1, 0, 1, 0, 0, 1, 0, 0, 0, 1, 0, 0, 1, 0, 1,
       1, 1, 0, 0, 0, 1, 1, 0, 1,
1, 1, 1, 0, 0, 0, 1, 0, 1,
                                   1,
                                                             1,
                                       0, 0, 0,
                                                1, 0, 1, 0,
                                   0,
                                       0, 0, 0,
          0, 1, 0, 0, 01)
Validate the improved accuracy classification on Training Data
from sklearn.metrics import confusion_matrix
import seaborn as sns
In [88]:
bgd_pred_train = polar_BGD_model.predict(newXtrain, polar_bgd_theta)
polar_BGD_model.getAccuracy(newXtrain, y train, polar_bgd_theta)
Out[88]:
100.0
In [89]:
nm_pred_train = polar_NM_model.predict(newXtrain, polar_nm_theta)
polar NM model.getAccuracy(newXtrain, y train, polar nm theta )
Out[89]:
100.0
Validate the improved accuracy classification on TestData
In [901:
bgd pred test = polar BGD model.predict(newXtest, polar bgd theta)
polar BGD model.getAccuracy(newXtest, y test, polar bgd theta)
Out[90]:
100.0
In [91]:
nm_pred_test = polar_NM_model.predict(newXtest, polar_nm_theta)
polar NM_model.getAccuracy(newXtest, y test, polar nm_theta)
Out[91]:
100.0
In [921:
nm pred train.shape
Out[92]:
(160.)
In [93]:
y train.shape
Out[93]:
(160,)
confusion_matrix(y_true = y_train, y_pred = nm_pred_train)
Out[94]:
array([[85, 0], [ 0, 75]])
In [95]:
bgd pred train
Out[95]:
array([1, 0, 1, 0, 1, 1, 0, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 1, 0, 1,
       1, 1, 0, 0, 1, 0, 1, 1, 1, 0, 0, 0, 1, 1,
                                                    0, 1, 1,
       1, 1, 0, 1, 0, 1, 0, 1, 0, 0, 1, 0, 1, 0,
                                                    0, 0, 0, 0, 0, 1, 0, 1,
       1, 1, 1, 0, 1, 1, 0, 1, 0, 0, 1, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 0,
       1, 1, 0, 1, 1, 1, 0, 1, 0, 1, 0, 0, 1, 0, 0, 1, 0, 0, 1, 0, 1,
       1, 1, 0, 0, 0, 1, 1, 0, 1, 1, 0, 0, 0, 1, 0, 1, 0,
                                                             1, 1, 1, 0, 1,
       1, 1, 1, 0, 0, 0, 1, 0, 1,
                                   0, 0, 0, 0,
       1, 0, 1, 0, 0, 0])
```

sns.heatmap(confusion_matrix(y_true = y_train, y_pred = bgd_pred_train), cmap = 'coolwarm', annot =True, ax = axes[0,0]

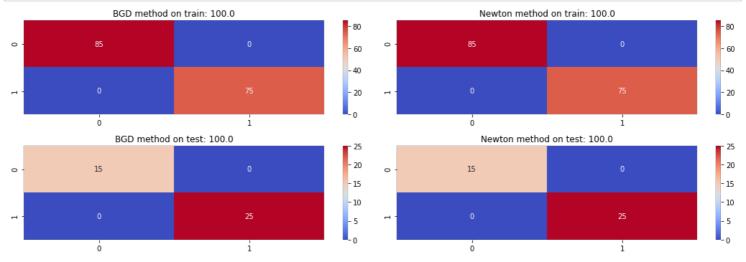
In [96]:

fig ,axes = plt.subplots(2,2,figsize = (15,5))

```
sns.heatmap(confusion_matrix(y_true = y_train, y_pred = nm_pred_train), cmap = 'coolwarm', annot =True, ax = axes[0,1]

sns.heatmap(confusion_matrix(y_true = y_test, y_pred = bgd_pred_test), cmap = 'coolwarm', annot =True, ax = axes[1,0])
sns.heatmap(confusion_matrix(y_true = y_test, y_pred = nm_pred_test), cmap = 'coolwarm', annot =True, ax = axes[1,1])

axes[0,0].set_title(f'BGD method on train: {polar_BGD_model.getAccuracy(newXtrain, y_train, polar_bgd_theta)}')
axes[0,1].set_title(f'Newton method on train: {polar_NM_model.getAccuracy(newXtrain, y_train, polar_nm_theta)}')
axes[1,0].set_title(f'BGD method on test: {polar_BGD_model.getAccuracy(newXtest, y_test, polar_bgd_theta)}')
axes[1,1].set_title(f'Newton method on test: {polar_NM_model.getAccuracy(newXtest, y_test, polar_nm_theta)}')
plt.tight_layout()
plt.show()
```



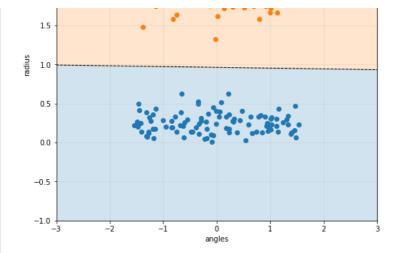
```
In [97]:
```

```
# plot
fig1 = plt.figure(figsize=(8,8))
ax = plt.axes()
plt.title('Sample data for classification problem')
plt.grid(axis='both', alpha=.25)
plt.ylabel('radius')
plt.legend()
b = polar_nm_theta[0]
w1 = polar_nm_theta[1]
w2 = polar_nm_theta[2]
# Calculate the intercept and gradient of the decision boundary.
c = b/w2
m = w1/w2
xmin, xmax = -3, 3
ymin, ymax = -1, 3
xd = np.array([xmin, xmax])
\# xd = np.linspace(-3, 3, 100)
yd = (m*xd) + c
yd += 2
plt.plot(xd, yd, 'k', lw=1, ls='--')
plt.fill_between(xd, yd, ymin, color='tab:blue', alpha=0.2)
plt.fill_between(xd, yd, ymax, color='tab:orange', alpha=0.2)
plt.xlim(-3, 3)
plt.ylim(-1, 3)
plt.xlabel('angles')
plt.ylabel('radius')
plt.legend()
```

Out[97]:

<matplotlib.legend.Legend at 0x7f9fc73560d0>





Summary from polar transformation

- After using the polar transformation, the accuracy rate on both training and test sets skyrockets to 100%, meaning the model can find the linearly seperable line that classify data into 2 classes
- In terms of iterations, the Newton's method performs significantly better as it needs to iterate fewer than 1000 times to achieve perfect classification, compared to almost 10,000 iterations for Batch Gradient Descent method with learning rate at 0.001

Perform Newton's Method on Lab03dataset

```
In [981:
import pandas as pd
In [99]:
data_train = pd.read_csv('clean_train.csv')
data test = pd.read csv('clean test.csv')
In [100]:
data train = data train.iloc[:, 1:]
In [101]:
data_test = data_test.iloc[:, 1:]
In [102]:
X = data_train.iloc[:, :-1]
y = data_train.iloc[:, -1]
In [103]:
from sklearn.model_selection import train_test_split
X train, X test, Y train, y test = train test split(X, y , test size = 0.3, random state = 16)
In [104]:
data train.shape
Out[104]:
(614, 12)
In [105]:
X_train.shape
Out[105]:
(429, 11)
In [106]:
X test.shape
Out[106]:
(185, 11)
```

Normalize the data

```
In [107]:
numerical cols = ['Dependents', 'ApplicantIncome', 'CoapplicantIncome', 'LoanAmount', 'Loan Amount Term']
In [108]:
from sklearn preprocessing import StandardScaler
```

```
std scaler = StandardScaler()
```

In [109]:

normalized_X_train = X_train
normalized_X_test = X_test
normalized_data test = data test

In [110]:

normalized X train[numerical cols] = std_scaler.fit_transform(normalized X_train[numerical_cols])

/tmp/ipykernel_307/3168317029.py:1: SettingWithCopyWarning:

A value is trying to be set on a copy of a slice from a DataFrame.

Try using .loc[row_indexer,col_indexer] = value instead

See the caveats in the documentation: $https://pandas.pydata.org/pandas-docs/stable/user_guide/indexing.html#returning-a-view-versus-a-copy$

normalized X_train[numerical_cols] = std_scaler.fit_transform(normalized_X_train[numerical_cols])

opt/conda/lib/python3.8/site-packages/pandas/core/indexing.py:966: SettingWithCopyWarning:

A value is trying to be set on a copy of a slice from a DataFrame.

Try using .loc[row_indexer,col_indexer] = value instead

See the caveats in the documentation: https://pandas.pydata.org/pandas-docs/stable/user_guide/indexing.html#returning-a-view-versus-a-copy self.obj[item] = s

In [111]:

normalized X train

Out[111]:

	Gender	Married	Dependents	Education	Self_Employed	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_History	Property_Area
119	0	0	-0.705357	0	0	0.802648	-0.497606	1.289912	0.247680	1.0	2
199	1	0	-0.705357	0	1	0.900462	-0.497606	-0.770259	0.247680	1.0	2
20	1	1	-0.705357	1	0	0.348609	-0.497606	-0.524444	0.247680	0.0	2
253	1	1	0.119161	1	0	-0.477352	1.652278	1.524022	-2.553407	1.0	1
548	0	0	-0.705357	0	0	-0.090890	-0.497606	-0.536149	0.247680	0.0	1
452	1	1	-0.705357	0	0	-0.264707	0.027073	0.002305	0.247680	0.0	0
321	1	1	2.592716	0	0	-0.199608	-0.440385	-0.290333	0.247680	1.0	1
581	1	0	-0.705357	0	0	-0.613663	9.746813	-0.688321	0.247680	1.0	2
121	0	0	-0.705357	0	0	-0.228688	-0.497606	-1.226775	0.247680	1.0	1
238	0	0	0.119161	0	0	-0.287178	-0.497606	-0.430799	0.247680	1.0	0

429 rows × 11 columns

In [112]:

normalized X test[numerical cols] = std scaler.transform(normalized X test[numerical cols])

/tmp/ipykernel_307/1923492441.py:1: SettingWithCopyWarning:

A value is trying to be set on a copy of a slice from a DataFrame.

Try using .loc[row_indexer,col_indexer] = value instead

See the caveats in the documentation: https://pandas.pydata.org/pandas-docs/stable/user_guide/indexing.html#returning-a -view-versus-a-copy

normalized_X_test[numerical_cols] = std_scaler.transform(normalized_X_test[numerical_cols])

/opt/conda/lib/python3.8/site-packages/pandas/core/indexing.py:966: SettingWithCopyWarning: A value is trying to be set on a copy of a slice from a DataFrame.

Try using .loc[row_indexer,col_indexer] = value instead

See the caveats in the documentation: https://pandas.pydata.org/pandas-docs/stable/user_guide/indexing.html#returning-a-view-versus-a-copy self.obj[item] = s

In [113]:

normalized X test

Out[113]

	Gender	Married	Dependents	Education	Self_Employed	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_History	Property_Area
274	1	1	0.943680	0	0	-0.272638	-0.497606	-0.688321	0.247680	1.0	1
315	1	1	0.119161	1	0	-0.355416	-0.001083	-0.442505	-2.553407	1.0	2
175	1	1	-0.705357	0	0	-0.339224	0.097010	-0.383977	0.247680	1.0	0
598	1	1	-0.705357	0	1	0.729123	-0.497606	0.365176	0.247680	1.0	0
390	1	0	2.592716	0	0	0.597604	-0.497606	0.423703	0.247680	1.0	0
436	1	0	-0.705357	0	0	-0.598793	0.062799	-1.156542	0.247680	1.0	1
523	1	1	0.943680	0	1	0.396194	1.671957	3.876831	0.247680	1.0	0
213	1	1	2.592716	1	1	0.025263	-0.497606	-0.220100	0.247680	1.0	0
172	1	1	-N 705357	n	n	บ บวยบชด	1 205407	N 447114	N 24768N	1 0	1

Gender Married Dependents Education Self_Employed ApplicantIncome CoapplicantIncome LoanAmount Loan_Amount_Term Credit_History Property_Area

185 rows × 11 columns

In [1141:

normalized data test[numerical cols] = std scaler.transform(normalized data test[numerical cols])

In [115]:

normalized data test

Out[115]:

	Gender	Married	Dependents	Education	Self_Employed	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_History	Property_Area
0	1	1	-1.286937	0	0	-0.917012	-0.497757	-1.747134	-5.350639	1.0	2
1	1	1	-0.607107	0	0	-0.917084	-0.497620	-1.744942	-5.350639	1.0	2
2	1	1	0.072724	0	0	-0.917031	-0.497592	-1.733706	-5.350639	1.0	2
3	1	1	0.072724	0	0	-0.917104	-0.497524	-1.748504	-5.350639	1.0	2
4	1	0	-1.286937	1	0	-0.917079	-0.497757	-1.751519	-5.350639	1.0	2
362	1	1	1.432385	1	1	-0.917058	-0.497594	-1.746723	-5.350639	1.0	2
363	1	1	-1.286937	0	0	-0.917054	-0.497692	-1.746449	-5.350639	1.0	2
364	1	0	-1.286937	0	0	-0.917079	-0.497574	-1.744942	-5.350639	0.0	1
365	1	1	-1.286937	0	0	-0.917031	-0.497538	-1.740557	-5.350639	1.0	0
366	1	0	-1.286937	0	1	-0.916917	-0.497757	-1.748779	-5.394229	1.0	0

367 rows x 11 columns

Turn pandas dataframe into numpy for easier manipulation and Train the model

```
In [116]:
np_Xtrain = np.asarray(normalized_X_train)
np_Xtest = np.asarray(normalized_X_test)
np_ytrain = np.asarray(y_train)
np_ytest = np.asarray(y_test)
```

```
In [117]:
```

```
np_Xtrain = np.insert(np_Xtrain, 0,1, axis=1)
np_Xtest = np.insert(np_Xtest, 0,1, axis=1)
```

In [118]:

np Xtrain.shape

Out[118]:

(429, 12)

In [119]:

```
alpha = 0.001
iterations = 50000
m,n = np_Xtrain.shape
BGD_model = Logistic_BGD()
initial_theta = np.zeros(n)
bgd theta, bgd cost = BGD model.gradientAscent(np Xtrain, np ytrain, initial theta, alpha, iterations)
```

Minimum at iteration: 49999

In [120]:

bgd cost[-1]

Out[120]:

0.4794791327357925

In [121]:

bgd yhat = BGD model.predict(np Xtest, bgd theta)

```
In [122]:
```

```
m,n = np_Xtrain.shape
initial_theta = np.zeros(n)
iterations = 1000
NM_model = Logistic_NM()
nm_theta, nm_cost = NM_model.newtonsMethod(np_Xtrain, np_ytrain, initial_theta, iterations)
# print("theta:",nm theta)
```

Minimum at iteration: 999

```
In [123]:
nm_cost[-1]
Out[123]:
0.4635955902680171
In [1241:
nm_yhat = NM_model.predict(np_Xtest, nm_theta)
In [125]:
iteration = 1000
plt.plot(np.arange(iteration), bgd_cost[:iteration], label = 'BGD cost')
plt.plot(np.arange(iteration), nm_cost[:iteration], label = 'NM cost')
plt.legend()
plt.show()
 0.70
                                         BGD cost
                                      NM cost
 0.65
 0.60
 0.55
 0.50
                                            1000
             200
                     400
                             600
                                     800
In [126]:
from sklearn.metrics import accuracy score
In [127]:
bgd yhat.shape
Out[127]:
(185,)
In [128]:
np.sum(bgd yhat)
Out[128]:
156
In [129]:
confusion matrix(y true =np ytest, y pred = bgd yhat) / len(bgd yhat)
Out[129]:
array([[0.14054054, 0.19459459], [0.01621622, 0.64864865]])
In [130]:
fig, axes = plt.subplots(1,2, figsize = (15,5))
sns.heatmap(confusion_matrix(y_true =np_ytest, y_pred = bgd_yhat) / len(bgd_yhat) , cmap = 'coolwarm', annot =True,
fmt = ".2%", ax = axes[0])
axes[0].set title(f'Confusion matrix of test set with Batch Gradient Method\nwith accuracy score of {np.round(accuracy
_score(y_true = np_ytest, y_pred = bgd_yhat ), 2)}')
axes[0].set_ylabel('True value')
axes[0].set_xlabel('Predicted value')
t = ".2%", ax = axes[1])
axes[1].set_title(f"Confusion matrix of test set with Newton's Method \nwith accuracy score of {np.round(accuracy_scor
e(y_true = np_ytest, y_pred = nm_yhat ), 2)}")
axes[1].set_ylabel('True value')
axes[1].set_xlabel('Predicted value')
plt.show()
  Confusion matrix of test set with Batch Gradient Method
                                                               Confusion matrix of test set with Newton's Method
              with accuracy score of 0.79
                                                                         with accuracy score of 0.79
                                                 - 0.6
                                                                                                            - 0.6
                                                                                                            - 0.5
                                                 - 0.5
  0
                                19.46%
                                                                                           19.46%
                                                 - 0.4
                                                                                                            - 0.4
 True
                                                            Free
```

- 0.3

- 0.3

Summary from running BGD and NW methods on Lab03 data (Loan Prediction)

- The accuracy rate from both approaches result in almost the same prediction accuracy rate, at 79%
- However, effectivenes-wise, it takes only 1,000 iterations for Newton's method to achieve such rate, compared to the computationally expensive 50,000 times of iteration of normal Batch Gradient Descent method
- In conclusion, the Newton's Method is the preferred choice as it manages to yield the same classification accuracy with far fewer computations