

CS111 Assignment

Nonlinear Regression via Gradient Descent

Prof. Drummond

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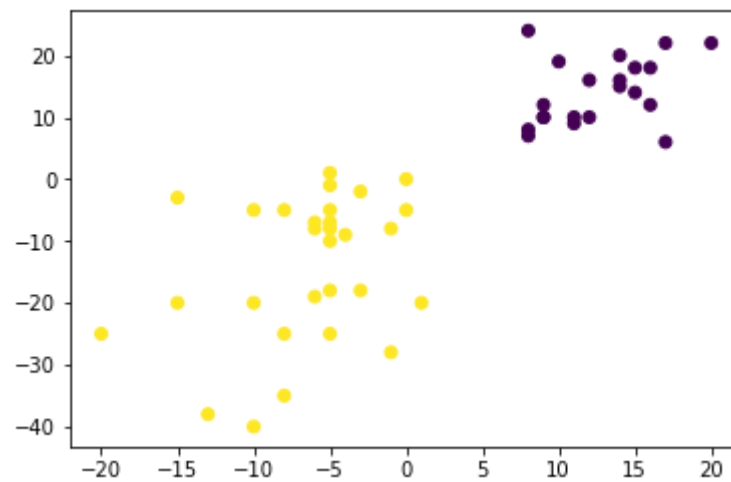
1. Data exploration

```
In [3]: import csv, math
import numpy as np
import matplotlib.pyplot as plt
```

```
In [5]: with open('cs111-svm-dataset---sheet1.csv') as csvfile:
        data = list(csv.reader(csvfile, delimiter=','))

        data = np.array(data[1:], dtype = np.float)
```

```
In [12]: x1 = data[:,0]
x2 = data[:,1]
y = data[:,2]
plt.scatter(x1,x2, c= y)
plt.show()
```



Our original dataset comprises of 50 observations. We have 2 independent variables (X1 and X2) and an independent variable, which classifies our the data into 2 classes (0 and 1). The graph illustrates that the data are groups into 2 clusters, purple (representing data points with value 0) and yellow (representing data points with value 1).

2. Gradient descent

The code below runs the gradient descent for a function of 3 variables (m1, m2, b) to generate an SVM which is the equation of the separating line. We are trying to minimize the value of the error function, which is the average loss across all observations. Instead of using a for loop to examine every data point, I use operations on matrices using the numpy library to implement it efficiently.

```
In [69]: def error_function(m1, m2, b, f):
        """
        Return the value of the error function
        """
        error = np.mean(np.log(1 + np.exp(-y*f)))
        return error

    def gradient(m1, m2, b, f):
        """
        Return a list of gradient corresponding to current
        value of m1, m2, b
        """

        gradient_m1 = np.mean((1/(1 + np.exp(-y*f))) * np.exp(-y*f) * (-y)*x1)
```

```

gradient_m2 = np.mean((1/(1 + np.exp(-y*f))) * np.exp(-y*f) * (-y)*x2)
gradient_b = np.mean((1/(1 + np.exp(-y*f))) * np.exp(-y*f) * (-y))

lis = [gradient_m1, gradient_m2, gradient_b]
return lis

def plotting(m1,m2,b):
    """
    Plotting the data and the current SVM at each step
    """
    global x1, x2, y
    plt.figure()
    plt.scatter(x1,x2, c= y)
    abline_values = [-m1/m2 * x1 - b/m2 for x1 in range(-20,20)]
    plt.plot(range(-20,20), abline_values)
    plt.show()

def gradient_descent(guess, step_size, max_steps):
    global x1, x2, y
    #main program
    m1 = guess[0]
    m2 = guess[1]
    b = guess[2]

    errors = []
    for step in range(max_steps):
        #a vector contains predicted y values for all x1 and x2 based on c
        #urrent m1,m2,b
        f= x1*m1 + x2*m2 +b

        gradient_list = gradient(m1,m2,b,f)
        temp_m1 = m1 - step_size*gradient_list[0]
        temp_m2 = m2 - step_size*gradient_list[1]
        temp_b = b - step_size*gradient_list[2]

        m1, m2, b = temp_m1, temp_m2, temp_b #simultaneous update
        errors.append(error_function(m1, m2, b,f))

        if step% 300 ==0 or step ==999: #plot the graph of the SVM each 30
        0 steps
            print("Current step: {}".format(step))
            print("Current error: {}".format(error_function(m1, m2, b,f)))
            print("m1 = {:f}, m2 = {:f}, b = {:f}".format(m1,m2,b))
            plotting(m1,m2,b)
    return errors

```

3. Implementation

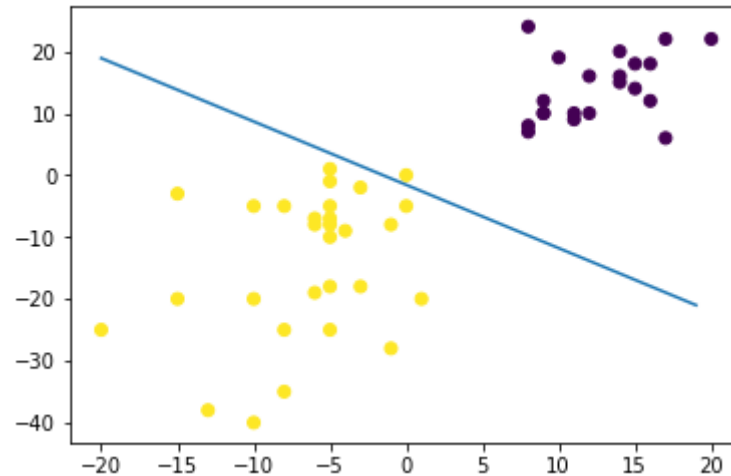
I implement the gradient descent for 1000 steps and a step size of 0.01 to find the SVM for our data set. I plot the line f learned at each 300 steps to visually demonstrate the algorithm. I also plot the error function over time to observe how the algorithm minimize the error values.

```
In [70]: initial_guess = [2,2,3]
errors = gradient_descent(initial_guess, 0.01, 1000)
```

Current step: 0

Current error: 22.572266803508132

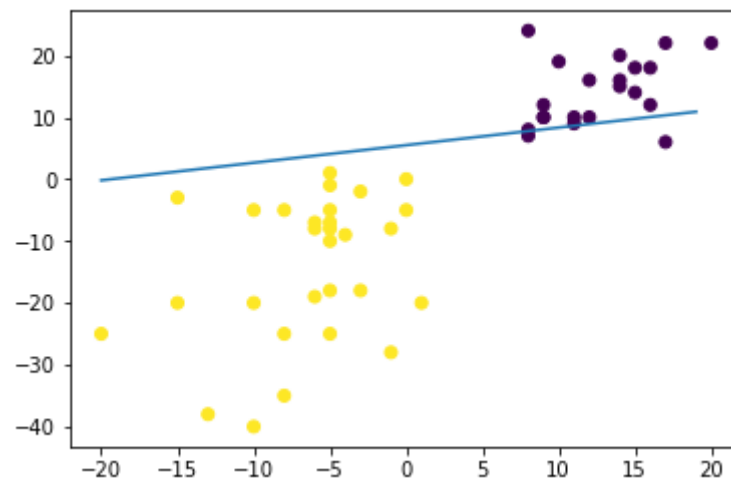
$m_1 = 1.962807$, $m_2 = 1.917400$, $b = 3.005608$



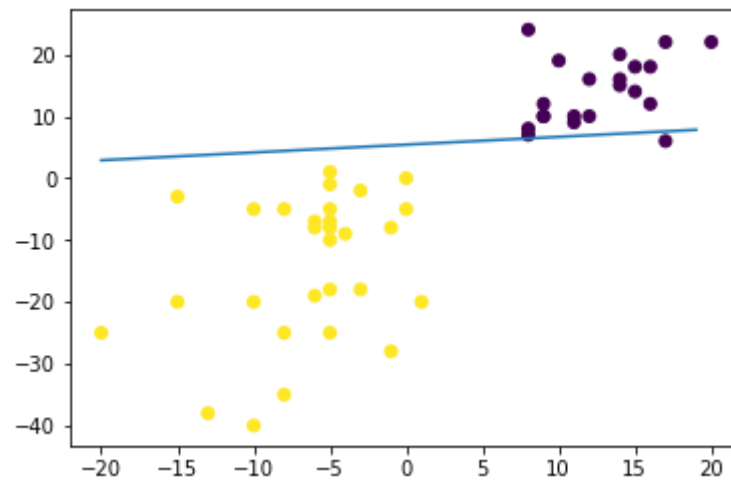
Current step: 300

Current error: 0.2985686528003694

$m_1 = 0.167230$, $m_2 = -0.588498$, $b = 3.232990$



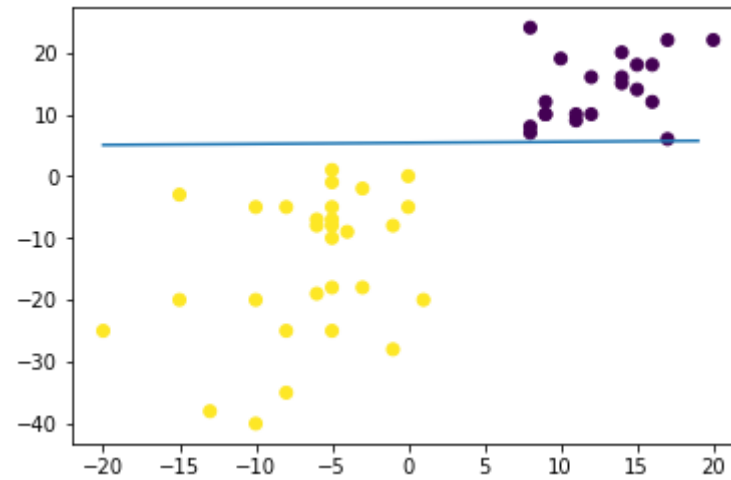
Current step: 600
 Current error: 0.295471264760855
 $m_1 = 0.075748$, $m_2 = -0.600102$, $b = 3.248865$



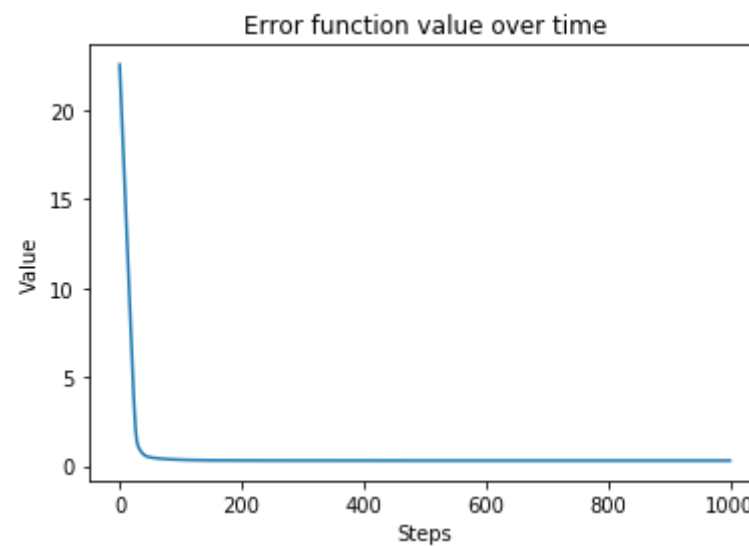
Current step: 900
 Current error: 0.29451616343733794
 $m_1 = 0.024023$, $m_2 = -0.605108$, $b = 3.260000$



Current step: 999
 Current error: 0.29431896117018413
 $m1 = 0.010484$, $m2 = -0.606273$, $b = 3.263185$



```
In [52]: plt.plot([step for step in range(1000)], errors)
plt.title("Error function value over time")
plt.xlabel("Steps")
plt.ylabel("Value")
plt.show()
```



As we can see, the error value decreases sharply and reaches an asymptotic value of approximately 0.30 after about 20 steps. In general, the time it takes for the error value to converge depends on several factors:

- initial guess: choosing an unsuitable initial guess can make the function take longer time to converge. We also run the risks of getting stuck in local optimum. Therefore, we can examine the data by plotting it initially to choose a sensible initial value for m_1 , m_2 and b .
- step size: A small step size takes longer time to converge. However, an exceedingly big step size also causes the algorithm to overshoot and even diverge. We can experiment with different step sizes to find the most optimal one.

4. Interpretation of the Loss Function

- The loss function is an indication of the accuracy of the SVM (our classification algorithm). Suppose that an observation x_o have an y value of 0, the SVM is accurate if it classifies x_o as having the y value of 0 and incorrect otherwise.
- To be more specific, the loss function will return a small value if the classification is correct since $e^{-y \cdot f(x)}$ would be small in this case. On the other hand, it returns a large value if the classification is incorrect since $e^{-y \cdot f(x)}$ would be large in this case.

Due to this property, we can implement gradient descent to minimize error function, which is the average of the loss function across all observations, to find the SVM that can classify the data best.

5. SVM strengths and weaknesses

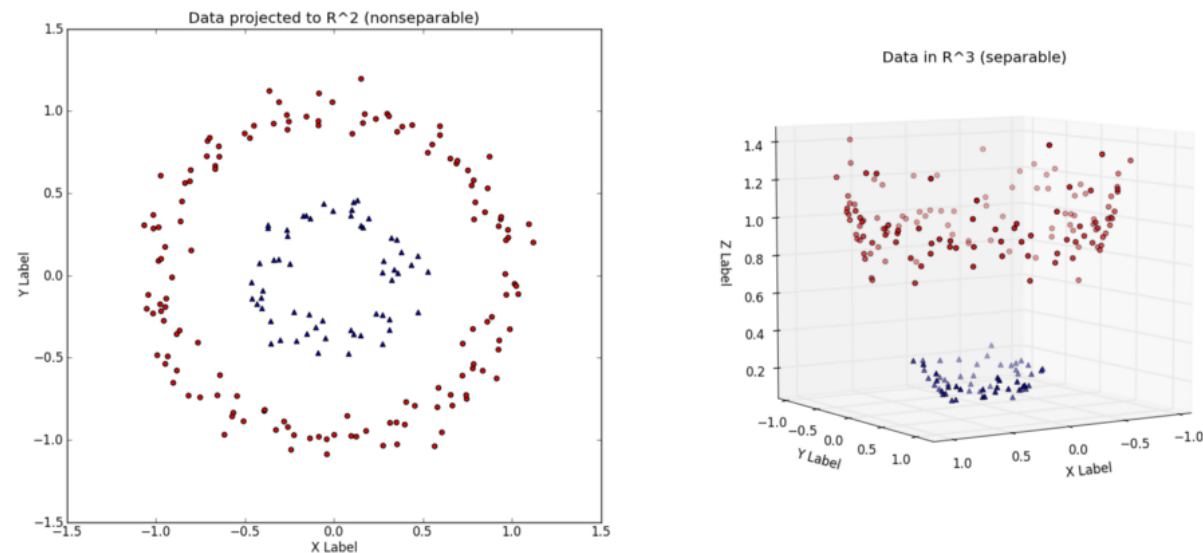
The strength of our algorithm and SVM is that it is quite easy to implement and we do not need to calculate an analytical solution, which can be hard to do when we have millions of data points. Using gradient descent, we can also specify the step size at our disposal.

However, a major disadvantage of SVM is that it can only classify **linearly separable data**. For example in this case, a line is sufficient to classify the data. On the other hand, for data that cannot be separated linearly, SVM classification can be quite inaccurate.

In this case, we can implement **kernel trick** to transform our existing data into a higher dimensional data, which can help us classify the data better. For example in the figure below, the data is not linearly separable (left) but we can map them into a 3-dimensional space (right) where a separating hyperplane can be found.

```
In [71]: from IPython.display import Image  
Image("Kerneltrick.png")
```

Out[71]:



Note: this figure is referenced from <https://towardsdatascience.com/understanding-the-kernel-trick-e0bc6112ef78>

6. SVM for Spams Classification

We can specify some variables that are typical for an email to classify whether it is a spam or not. For example:

- x_1 - Frequency of words: Spams are usually used in promotion. Therefore, they are more likely to contain clickbait words such as "Discounts", "Free Trials", "Click here". We can build a list of all these words and calculate the frequency of those words appearing in the email as a variable in the classifier.
- x_2 - Sender: spam email usually comes from senders who are not in the contact list or have mutual connections with a person. It also comes from dubious companies that usually send out promos. However, we should make the weight on this less important to avoid mistakenly classifying important email from new people.
- x_3 - Number of embedded links: A spam email usually contains many links that users can click on. Usually it's dangerous since it may contain viruses or unwanted content.

To implement the classifier, we can collect a sample of thousands of emails including both spam and normal emails (larger data improve accuracy). The classification variable y (0 or 1) indicates whether an email is spam or not. We can then implement the SVM with the independent variables (x_1 , x_2 , x_3) and the dependent variable y to classify our data. To examine the accuracy, we use the SVM from the training set and test it on a new data set. We can then decide to add more variables or data to improve the accuracy of our algorithm.