```
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```

```
In [189]: import numpy as np
   import sklearn.datasets as ds
   import matplotlib.pyplot as plt
   from matplotlib import style
   style.use('ggplot')
   %matplotlib inline
   import re
   import pandas as pd
```

Numpy is library for scientific computing in Python. It has efficient implementation of n-dimensional array (tensor) manupulations, which is useful for machine learning applications.

```
In [190]: import numpy as np
```

We can convert a list into numpy array (tensor)

```
In [191]: b = [[1, 2, 4], [2, 6, 9]]
a = np.array(b)
a
```

```
Out[191]: array([[1, 2, 4], [2, 6, 9]])
```

We can check the dimensions of the array

```
In [192]: a.shape
Out[192]: (2, 3)
```

We can apply simple arithmetic operation on all element of a tensor

You can transpose a tensor

You can apply aggregate functions on the whole tensor

```
In [195]: np.sum(a)
Out[195]: 24
```

or on one dimension of it

```
In [196]: np.sum(a, axis=0)
Out[196]: array([ 3,  8, 13])
In [197]: np.sum(a, axis=1)
Out[197]: array([ 7, 17])
```

We can do element-wise arithmetic operation on two tensors (of the same size)

If you want to multiply all columns of a tensor by vector (for example if you want to multiply all data features by their lables) you need a trick. This multiplication shows up in calculating the gradients.

```
In [199]: a = np.array([[1, 2, 4], [2, 6, 9]])
b = np.array([1,-1])
print(a)
print(b)

[[1 2 4]
      [2 6 9]]
      [1 -1]
```

Here we want to multiply the first row of a by 1 and the second row of a by -1. Simply multiplying a by b does not work because a and b do not have the same dimension

To do this multiplication we first have to assume b has one column and then repeat the column of b with the number of columns in a. We use tile function to do that

Now we can multiply each column of a by b:

You can create inital random vector using numpy (using N(0,1)):

```
In [203]: mu = 0 #mean
    sigma = 1 #standard deviation
    r = np.random.normal(mu,sigma, 1000) #draws 1000 samples from a normal distribution
```

We can apply functions on tensors

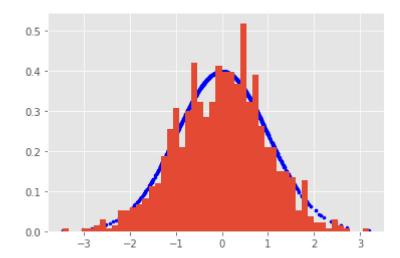
```
In [204]: #implementation of Normal distribution
def normal(x, mu, sigma):
    return np.exp( -0.5 * ((x-mu)/sigma)**2)/np.sqrt(2.0*np.pi*sigma**2)

#probability of samples on the Normal distribution
probabilities = normal(r, mu, sigma)
```

Numpy has useful APIs for analysis. Here we plot the histogram of samples and also plot the probabilies to see if the samples follow the normal distribution.

```
In [205]: counts, bins = np.histogram(r,50,density=True)
    plt.hist(bins[:-1], bins, weights=counts)
    plt.scatter(r, probabilities, c='b', marker='.')
```

Out[205]: <matplotlib.collections.PathCollection at 0x132a0a830d0>



```
In [206]: def read_data(filename):
    f = open(filename, 'r')
    p = re.compile(',')
    xdata = []
    ydata = []
    header = f.readline().strip()
    varnames = p.split(header)
    namehash = {}
    for l in f:
        li = p.split(l.strip())
        xdata.append([float(x) for x in li[:-1]])
        ydata.append(float(li[-1]))

    return np.array(xdata), np.array(ydata)
```

Assuming our data is x is available in numpy we use numpy to implement logistic regression

```
In [207]: (xtrain_whole, ytrain_whole) = read_data('spambase-train.csv')
    (xtest, ytest) = read_data('spambase-test.csv')

In [208]: print("The shape of xtrain:", xtrain_whole.shape)
    print("The shape of ytrain:", ytrain_whole.shape)
    print("The shape of xtest:", xtest.shape)
    print("The shape of ytest:", ytest.shape)

The shape of xtrain: (3601, 54)
    The shape of xtest: (1000, 54)
    The shape of ytest: (1000,)
```

before training make we normalize the input data (features)

```
In [209]: xmean = np.mean(xtrain_whole, axis=0)
    xstd = np.std(xtrain_whole, axis=0)
    xtrain_normal_whole = (xtrain_whole-xmean) / xstd
    xtest_normal = (xtest-xmean) / xstd
```

We need to create a validation set. We create an array of indecies and permute it.

```
In [210]: premute_indicies = np.random.permutation(np.arange(xtrain_whole.shape[0]))
```

We keep the first 2600 data points as the training data and rest as the validation data

```
In [211]: xtrain_normal = xtrain_normal_whole[premute_indicies[:2600]]
    ytrain = ytrain_whole[premute_indicies[:2600]]
    xval_normal = xtrain_normal_whole[premute_indicies[2600:]]
    yval = ytrain_whole[premute_indicies[2600:]]
```

Initiallizing the weights and bias with random values from N(0,1)

```
In [212]: weights = np.random.normal(0, 1, xtrain_normal.shape[1]);
bias = np.random.normal(0,1,1)
```

```
In [213]: #the sigmoid function
def sigmoid(z):
    s = 1 / (1 + np.exp(-z))
    return s
```

We can use dot-product from numpy to calculate the margin and pass it to the sigmoid function

```
In [214]: #w: weight vector (numpy array of size n)
#b: numpy array of size 1
#returns p(y=1/x, w, b)
def prob(x, w, b):
    z = sigmoid(np.dot(x,w) + b)
    #print(z)
    return z
```

You can also calculate l_2 penalty using linalg library of numpy

```
In [215]: np.linalg.norm(weights)
```

Out[215]: 7.412178281536248

```
Cross Entropy Loss = -\sum_{(y^i, \mathbf{x}^i) \in \mathcal{D}} y^i \log p(y = 1 | \mathbf{x}^i; \mathbf{w}, b) + (1 - y^i) \log(1 - p(y = 1 | \mathbf{x}^i; \mathbf{w}, b) - y^i) \log(1 - p(y = 1 | \mathbf{x}^i; \mathbf{w}, b))
```

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```
In [216]: #w: weight vector (numpy array of size n)
    #y_prob: p(y|x, w, b)
    #y_true: class variable data
    #lambda_: L2 penalty coefficient
    #returns the cross entropy loss
    def loss(w, y_prob, y_true, lambda_):
        Cross_Entropy_loss = -(1) * np.sum((y_true * np.log10(y_prob)) + ((1 - y_true)))
        return Cross_Entropy_loss
```

```
In [217]: #x: input variables (data of size m x n with m data point and n features)
    #w: weight vector (numpy array of size n)
    #y_prob: p(y|x, w, b)
    #y_true: class variable data
    #lambda_: l2 penalty coefficient
    #returns tuple of gradient w.r.t w and w.r.t to bias

def grad_w_b(x, w, y_prob, y_true, lambda_):
    grad_w = np.zeros(w.shape)
    #m = len(x)
    grad_b = 0.0
    grad_w = np.dot(x.T, (y_prob - y_true)) + (2 * lambda_ * w)
    grad_b = np.sum(y_prob - y_true)
    return(grad_w, grad_b)
```

```
In [224]:
          #lambda is the coeffienct of L2 norm penalty
          #learning rate is learning rate of gradient descent algorithm
          #max iter determines the maximum number of iterations if the gradients descent de
          #continue the training while gradient > 0.1 or the number steps is less max iter
          #returns model as tuple of (weights, bias)
          def fit(x, y_true, learning_rate, lambda_, max_iter, verbose=0):
              weights = np.random.normal(0, 1, x.shape[1]);
              bias = np.random.normal(0,1,1)
              #raise NotImplementedError
              costs = []
              i = 0
              while(i < max iter):</pre>
                  #To calculate y_prob
                  y_prob = prob(x, weights, bias)
                  Cross_Entropy_loss = loss(weights, y_prob, y_true, lambda_)
                  #costs = np.append(cost)
                  grad w, grad b = grad w b(x, weights, y prob, y true, lambda)
                  #print(grad w)
                  #print(grad b)
                  #Condition to break if grad w and grad b is less than 0.1
                  if((np.linalg.norm(grad w) < 0.1) and (np.linalg.norm(grad b) < 0.1)):</pre>
                      break
                  else:
                      weights = weights - learning rate * grad w
                      bias = bias - learning rate * grad b
                  i = i+1
              return (weights, bias)
In [225]: def accuracy(x, y_true, model):
              w, b = model
              return np.sum((prob(x, w, b)>0.5).astype(float) == y true) / y true.shape[0]
In [226]: | learning rate = 0.001
          lambda_ = 1.0
          model = fit(xtrain normal, ytrain, learning rate, lambda , 10000, verbose=1)
In [227]: | print("Train accuracy: ", accuracy(xtrain_normal, ytrain, model))
```

Train accuracy: 0.9338461538461539

```
In [228]: #grid search for finding the best hyperparams and model
          best_model = None
          best val = -1
          for lr in [0.01, 0.001, 0.0001, 0.00001]:
              for la in [5, 2, 1, 0.1, 0.01]:
                  model = fit(xtrain_normal, ytrain, lr, la, 10000)
                  val_acc = accuracy(xval_normal, yval, model)
                  print(lr, la, val_acc)
                  if val_acc > best_val:
                       best val = val acc
                       best_model = model
          <ipython-input-216-e5e81e887299>:7: RuntimeWarning: divide by zero encountered
          in log10
            Cross_Entropy_loss = -(1) * np.sum((y_true * np.log10(y_prob)) + ((1 - y_true))
          e) * np.log10(1 - y_prob))) + ((lambda_/2) * (w^{**}2))
          <ipython-input-216-e5e81e887299>:7: RuntimeWarning: invalid value encountered i
          n multiply
            Cross_Entropy_loss = -(1) * np.sum((y_true * np.log10(y_prob)) + ((1 - y_true))
          e) * np.log10(1 - y_prob))) + ((lambda_/2) * (w^{**}2))
          0.01 5 0.7922077922077922
          0.01 2 0.903096903096903
          0.01 1 0.9150849150849151
          0.01 0.1 0.9050949050949051
          0.01 0.01 0.8561438561438561
          0.001 5 0.9300699300699301
          0.001 2 0.9330669330669331
          0.001 1 0.9340659340659341
          0.001 0.1 0.9340659340659341
          0.001 0.01 0.9340659340659341
          0.0001 5 0.9300699300699301
          0.0001 2 0.9330669330669331
          0.0001 1 0.9340659340659341
          0.0001 0.1 0.9340659340659341
          0.0001 0.01 0.9340659340659341
          1e-05 5 0.9300699300699301
          1e-05 2 0.9330669330669331
          1e-05 1 0.9340659340659341
          1e-05 0.1 0.936063936063936
          1e-05 0.01 0.9340659340659341
In [229]: print("Test accuracy: ", accuracy(xtest normal, ytest, best model))
          Test accuracy: 0.941
 In [ ]:
```