Effect of Different Initialization Methods

1. Introduction

Training neural network requires specifying an initial value of the weights. A well chosen initialization method will help learning.

In this notebook, we will see how different initializations (random normal distribution, and He initialization) lead to different results.

A well chosen initialization can:

- · Speed up the convergence of gradient descent
- · Increase the odds of gradient descent converging to a lower training (and generalization) error

2. Import Packages and Set Default Parameters

- util_func provides some necessary functions for the calculations, e.g., Sigmoid, RELU.
- deep_neural_network.py provides the functions to construct deep neural network

```
In [1]: %load_ext autoreload
%autoreload 2
%matplotlib inline

import warnings
warnings.filterwarnings("ignore")
import numpy as np
from deep_neural_network import *
import matplotlib.pyplot as plt
import sklearn
import sklearn.datasets

plt.rcParams['figure.figsize'] = (7.0, 4.0)
plt.rcParams['image.interpolation'] = 'nearest'
plt.rcParams['image.cmap'] = 'gray' # set colormap
```

```
In [2]: np.random.seed(20)
```

3. Data Set

```
In [3]: # dataset: make a large circle containing a smaller circle in 2d.

def load_makecircles_dataset():

    train_x, train_y = sklearn.datasets.make_circles(n_samples = 300, noise = 0.05) # training set
    test_x, test_y = sklearn.datasets.make_circles(n_samples = 100, noise = 0.05) # test set

plt.scatter(train_x[:, 0], train_x[:, 1], c = train_y, s = 40, cmap = plt.cm.Spectral)

plt.title("training set")

# convert the data to proper shape
    train_x = train_x.T
    train_y = train_y.reshape((1, train_y.shape[0]))
    test_x = test_x.T
    test_y = test_y.reshape((1, test_y.shape[0]))

return train_x, train_y, test_x, test_y
```

```
In [4]: train_x, train_y, test_x, test_y = load_makecircles_dataset()
        print("Total number of training examples: " + str(train x.shape[1]))
        print("Total number of test examples: " + str(test_x.shape[1]))
        print("train_x shape: " + str(train_x.shape))
        print("train_y shape: " + str(train_y.shape))
        print("test_x shape: " + str(test_x.shape))
        print("test y shape: " + str(test y.shape))
        print("Example of y values: " + str(train y[0, 0:10]))
        Total number of training examples: 300
        Total number of test examples: 100
        train x shape: (2, 300)
        train y shape: (1, 300)
        test x shape: (2, 100)
        test y shape: (1, 100)
        Example of y values: [1 1 0 1 0 1 0 1 0 0]
                               training set
          1.0
          0.5
```

4. Some Useful Functions

-1.0

-0.5

0.0

0.5

1.0

-0.5

-1.0

```
In [5]: # use the trained params to predict the output

def predict(params, X):
    """
    Arguments:
    X: input features
    params: trained weight matrices and bias vectors of the neural network

    Returns:
    predicted labels for X, with the shape of (1, number of examples)
    """
    Aout, _ = L_layer_forward(X, params)
    m = X.shape[1] # number of examples
    Aout.reshape(1, m)
    Aout = (Aout > 0.5)
    return Aout
```

```
In [6]: # print the accuracy

def accuracy(params, X, Y):
    """
    Arguments:
    X: input features
    params: trained weight matrices and bias vectors of the neural network
    Y: true labels
    """
    Aout = predict(params, X)
    m = X.shape[1]
    print("Accuracy: " + str(np.sum(Aout == Y) / m))
```

```
In [7]: # plot the decision boundary contour
        def plot_dicision_boundary(model, X, Y):
            Arguments:
            model: the function that inputs X and outputs the predicted labels.
            X: the input features
            Y: the true labels
            # set min and max values and give some padding
            x1_{min}, x1_{max} = min(X[0, :]) - 1, max(X[0, :]) + 1
            x2_{min}, x2_{max} = min(X[1, :]) - 1, max(X[1, :]) + 1
            h = 0.01 # interval of the grid
            x1, x2 = np.meshgrid(np.arange(x1_min, x1_max, h), np.arange(x2_min, x2_max, h))
            # flatten x1 and x2 to 1-D arrays, concatenate along second axis, and transpose
            Z = model(np.c_[x1.ravel(), x2.ravel()].T) # Z.shape = (1, total number of grid points)
            Z = Z.reshape(x1.shape)
            # plot the contour
            plt.contourf(x1, x2, Z, cmap = plt.cm.Spectral)
            plt.scatter(X[0, :], X[1, :], c = Y[0, :], s = 40, cmap = plt.cm.Spectral)
            plt.show()
```

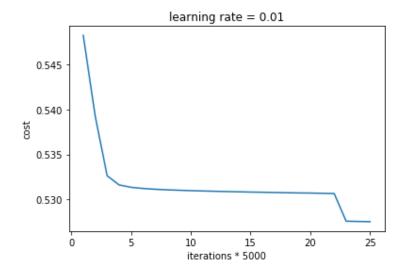
5. Random Initialization

Set weight matrices to random values following standard normal distribution.

5.1 Model Training

 $Set\ params['W'\ +\ str(I)] = np.random.randn(layer_dims[I],\ layer_dims[I-1])\ *\ 10\ in\ deep_neural_network.py$

current iteration: 1, cost: inf current iteration: 5000, cost: 0.5313345951096806 current iteration: 10000, cost: 0.5309626422074155 current iteration: 15000, cost: 0.5308041850528132 current iteration: 20000, cost: 0.5306922261648953 current iteration: 25000, cost: 0.5275040956001885



```
In [9]: print("For the training set:")
    accuracy(params_rand_init, train_x, train_y)
    print("For the test set:")
    accuracy(params_rand_init, test_x, test_y)
```

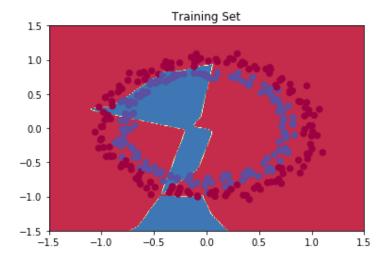
For the training set:

Accuracy: 0.63
For the test set:
Accuracy: 0.59

5.2. Result Analysis

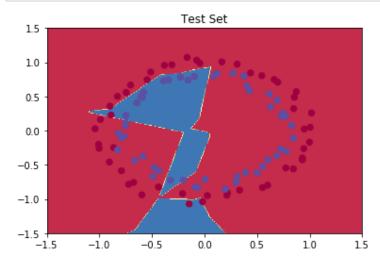
```
In [10]: # plot the decision boundary for the training set

axes = plt.gca()
axes.set_xlim([-1.5,1.5])
axes.set_ylim([-1.5,1.5])
plt.title("Training Set")
plot_dicision_boundary(lambda x : predict(params_rand_init, x), train_x, train_y)
```



```
In [11]: # plot the decision boundary for the test set

axes = plt.gca()
axes.set_xlim([-1.5,1.5])
axes.set_ylim([-1.5,1.5])
plt.title("Test Set")
plot_dicision_boundary(lambda x : predict(params_rand_init, x), test_x, test_y)
```



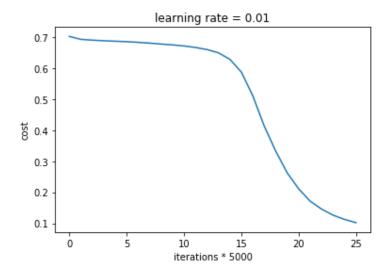
Note that If we use "random_normal" initialization, we can tune the factor, and we find that params['W' + str(I)] = np.random.randn(layer_dims[I], layer_dims[I-1]) * 1.0 gives the best result.

6. "HE" Initialization

6.1 Model Training

Set params['W' + str(I)] = np.random.randn(layer_dims[I], layer_dims[I-1]) * np.sqrt(2 / layer_dims[I-1]) in deep_neural_network.py

current iteration: 1, cost: 0.7023331991177456 current iteration: 5000, cost: 0.6852563590193823 current iteration: 10000, cost: 0.6715161681407267 current iteration: 15000, cost: 0.5875845905330312 current iteration: 20000, cost: 0.21132503139022032 current iteration: 25000, cost: 0.10201128914427847



```
In [13]: print("For the training set:")
    accuracy(params_he_init, train_x, train_y)
    print("For the test set:")
    accuracy(params_he_init, test_x, test_y)
```

For the training set:

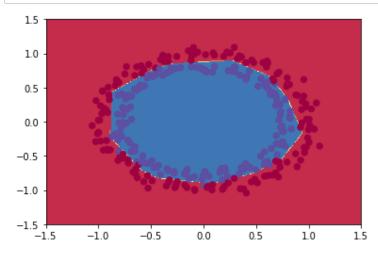
Accuracy: 0.986666666666667

For the test set: Accuracy: 0.98

6.2. Result Analysis

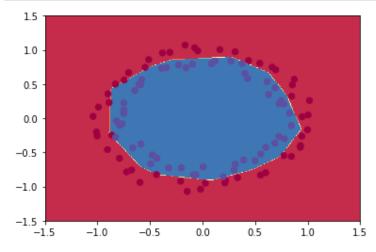
```
In [14]: # plot the decision boundary for the training set

axes = plt.gca()
axes.set_xlim([-1.5,1.5])
axes.set_ylim([-1.5,1.5])
plot_dicision_boundary(lambda x : predict(params_he_init, x), train_x, train_y)
```



```
In [15]: # plot the decision boundary for the test set

axes = plt.gca()
axes.set_xlim([-1.5,1.5])
axes.set_ylim([-1.5,1.5])
plot_dicision_boundary(lambda x : predict(params_he_init, x), test_x, test_y)
```



As we can see, with the same number of iterations and learning rate, He initialization yields much better results.