Problem Set 1: Linear Regression

Grading: This assignment is self-graded. It is extremely important that you complete all assignments (both the programming and written questions), as they will prepare you for quizzes and the final project. Please submit your self-assigned grade in "grade.txt"; we will randomly choose a subset of submissions to verify that the points computation in "grade.txt" is correct.

Total: 185 points.

To run and solve this assignment, one must have a working IPython Notebook installation. The easiest way to set it up for both Windows and Linux is to install Anaconda (https://www.continuum.io/downloads). Then save this file to your computer (use "Raw" link on gist\github), run Anaconda and choose this file in Anaconda's file explorer. Use Python 3 version. Below statements assume that you have already followed these instructions. If you are new to Python or its scientific library, Numpy, there are some nice tutorials here (https://www.learnpython.org/) and here (https://www.scipy-lectures.org/).

To run code in a cell or to render <u>Markdown (https://en.wikipedia.org/wiki/Markdown)+LaTeX (https://en.wikipedia.org/wiki/LaTeX)</u> press Ctr+Enter or [>|] (like "play") button above. To edit any code or text cell [double]click on its content. To change cell type, choose "Markdown" or "Code" in the drop-down menu above.

1. Numpy Tutorial

1.1 [5pt] Modify the cell below to return a 5x5 matrix of ones. Put some code there and press Ctrl+Enter to execute contents of the cell. You should see something like the output above. [1] (https://docs.scipy.org/doc/numpy-1.13.0/user/basics.creation.html#arrays-creation) [2] (https://docs.scipy.org/doc/numpy-1.13.0/reference/routines.array-creation.html#routines-array-creation)

1.2 [5pt] Let A be a 10x10 matrix and x be a 10-element column vector. Your friend writes the following code. How would you vectorize this code to run without any for loops? Compare execution speed for different values of n with https://ipython.readthedocs.io/en/stable/interactive/magics.html#magic-timeit).

```
In [2]:
        n = 10
        def compute_something(A, x):
            v = np.zeros((n, 1))
             for i in range(n):
                 for j in range(n):
                     v[i] += A[i, j] * x[j]
             return v
        A = np.random.rand(n, n)
        x = np.random.rand(n, 1)
        print(compute something(A, x))
        [[2.52666745]
         [2.34701873]
         [1.51475595]
         [2.40778152]
         [2.14875302]
         [1.61496849]
         [1.4716191]
         [2.63789067]
         [2.66146688]
         [1.64305876]]
In [3]: def vectorized(A, x):
             return A @ x
             # raise NotImplementedError('Put your vectorized code here!')
        print(vectorized(A, x))
        assert np.max(abs(vectorized(A, x) - compute_something(A, x))) < 1e-3
        [[2.52666745]
         [2.34701873]
         [1.51475595]
         [2.40778152]
         [2.14875302]
         [1.61496849]
         [1.4716191]
         [2.63789067]
         [2.66146688]
         [1.64305876]]
```

```
In [4]: for n in [5, 10, 100, 500]:
             A = np.random.rand(n, n)
             x = np.random.rand(n, 1)
             %timeit -n 5 compute something(A, x)
             %timeit -n 5 vectorized(A, x)
             # raise NotImplementedError('Put your timeit code here!')
             print('---')
        51.3 \mus \pm 5.16 \mus per loop (mean \pm std. dev. of 7 runs, 5 loops each)
        1.38 \mus \pm 872 ns per loop (mean \pm std. dev. of 7 runs, 5 loops each)
        209 \mu s \pm 26.4 \mu s per loop (mean \pm std. dev. of 7 runs, 5 loops each)
        The slowest run took 5.12 times longer than the fastest. This could m
        ean that an intermediate result is being cached.
         1.67 \mus \pm 1.5 \mus per loop (mean \pm std. dev. of 7 runs, 5 loops each)
         19 ms \pm 361 \mus per loop (mean \pm std. dev. of 7 runs, 5 loops each)
        2.3 \mus \pm 1 \mus per loop (mean \pm std. dev. of 7 runs, 5 loops each)
        467 ms \pm 3.28 ms per loop (mean \pm std. dev. of 7 runs, 5 loops each)
        The slowest run took 732.30 times longer than the fastest. This could
        mean that an intermediate result is being cached.
        830 \mus \pm 2.01 ms per loop (mean \pm std. dev. of 7 runs, 5 loops each)
         - - -
```

2. Linear regression with one variable

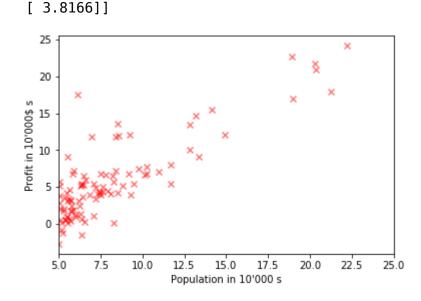
In this part of this exercise, you will implement linear regression with one variable to predict profits for a food truck. Suppose you are the CEO of a restaurant franchise and are considering different cities for opening a new outlet. The chain already has trucks in various cities and you have data for profits and populations from the cities. You would like to use this data to help you select which city to expand to next. The file ex1data.txt contains the dataset for our linear regression problem. The first column is the population of a city and the second column is the profit of a food truck in that city. A negative value for profit indicates a loss.

2.1 [10pt] Get a plot similar to above : [1]

(https://matplotlib.org/devdocs/api/_as_gen/matplotlib.pyplot.scatter.html) [2] (https://matplotlib.org/api/pyplot_api.html?highlight=xlim#matplotlib.pyplot.xlim) [3] (https://matplotlib.org/api/pyplot_api.html?highlight=matplotlib%20pyplot%20xlabel#matplotlib.pyplot.xlabel)

Before starting on any task, it is often useful to understand the data by visualizing it. For this dataset, you can use a scatter plot to visualize the data, since it has only two properties to plot (profit and population). Many other problems that you will encounter in real life are multi-dimensional and can't be plotted on a 2-d plot.

```
data = np.loadtxt('ex1data1.txt', delimiter=',')
X, y = data[:, 0, np.newaxis], data[:, 1, np.newaxis]
n = data.shape[0]
print(X.shape, y.shape, n)
print(X[:10], '\n', y[:10])
plt.scatter(X, y, marker='x', color='r', alpha=0.5)
plt.xlim(5, 25)
plt.xlabel('Population in 10\'000 s')
plt.ylabel('Profit in 10\'000$ s')
# NotImplementedError('Put the visualziation code here.')
plt.show()
(97, 1) (97, 1) 97
[[6.1101]
 [5.5277]
 [8.5186]
 [7.0032]
 [5.8598]
 [8.3829]
 [7.4764]
 [8.5781]
 [6.4862]
 [5.0546]]
 [[17.592]
 [ 9.1302]
 [13.662]
```



[11.854] [6.8233] [11.886] [4.3483] [12.] [6.5987]

2.2 Gradient Descent

In this part, you will fit the linear regression parameter θ to our dataset using gradient descent.

The objective of linear regression is to minimize the cost function

$$J(heta) = rac{1}{2m} \sum_{i=1}^m \left(h(x^{(i)}; heta) - y^{(i)}
ight)^2.$$

where the hypothesis $h(x;\theta)$ is given by the linear model (x' has an additional fake feature always equal to '1')

$$h(x; heta) = heta^T x' = heta_0 + heta_1 x$$

Recall that the parameters of your model are the θ_j values. These are the values you will adjust to minimize cost $J(\theta)$. One way to do this is to use the gradient descent algorithm for single element. In gradient descent algorithm, each iteration performs the update for a single sample.

$$heta_{j}^{(k+1)} = heta_{j}^{(k)} - \eta rac{1}{m} ig(h(x^{(i)}; heta) - y^{(i)} ig) x_{j}^{(i)}$$

With each step of gradient descent, your parameter θ_j come closer to the optimal values that will achieve the lowest cost J(θ).

2.2.1 [5pt] Where does this update rule comes from?

we need to calculate $rac{\partial J(heta)}{\partial heta_j}$ to get the gradient for each $heta_j$

$$rac{\partial J(heta)}{\partial heta_j} = rac{1}{2m} imes rac{\partial J(heta)}{\partial h(x; heta)} imes rac{\partial h(x; heta)}{\partial heta_j} = rac{1}{2m} imes 2ig(h(x^{(i)}; heta) - y^{(i)}ig) imes x_j^{(i)} = rac{1}{m}ig(h(x^{(i)}; heta) - y^{(i)}ig) x$$

We also need to choose a learning rate η to determine the update speed for each step. Then we can update θ_j iteratively to reduce the loss function $J(\theta)$ by

$$heta_{j}^{(k+1)} = heta_{j}^{(k)} - \eta rac{1}{m} ig(h(x^{(i)}; heta) - y^{(i)} ig) x_{j}^{(i)}$$

2.2.2 [30pt] Cost Implementation

As you perform gradient descent to learn to minimize the cost function, it is helpful to monitor the convergence by computing the cost. In this section, you will implement a function to calculate $J(\theta)$ so you can check the convergence of your gradient descent implementation.

In the following lines, we add another dimension to our data to accommodate the intercept term and compute the prediction and the loss. As you are doing this, remember that the variables X and y are not scalar values, but matrices whose rows represent the examples from the training set. In order to get x' add a column (https://docs.scipy.org/doc/numpy/reference/generated/numpy.insert.html) of ones to the data matrix X .

You should expect to see a cost of approximately 33.04.

```
# assertions below are true only for this
In [6]:
         # specific case and are given to ease debugging!
         def add column(X):
             assert len(X.shape) == 2 and X.shape[1] == 1
             # raise NotImplementedError("Insert a column of ones to the left
          side of the matrix")
             return np.insert(X, 0, 1, axis=1)
         def predict(X, theta):
             """ Computes h(x; theta) """
             assert len(X.shape) == 2 and X.shape[1] == 1
             assert theta.shape == (2, 1)
             X \text{ prime} = \text{add column}(X)
             pred = X prime @ theta
             # raise NotImplementedError("Compute the regression predictions")
             return pred
         def loss(X, y, theta):
             assert X.shape == (n, 1)
             assert y.shape == (n, 1)
             assert theta.shape == (2, 1)
             X \text{ prime} = \text{add column}(X)
             assert X prime.shape == (n, 2)
             # raise NotImplementedError("Compute the model loss; use the pred
         ict() function")
             loss = ((predict(X, theta) - y)**2).mean()/2
             return loss
         theta init = np.zeros((2, 1))
         print(loss(X, y, theta_init))
```

32.072733877455676

2.2.3 [40pt] GD Implementation

Next, you will implement gradient descent. The loop structure has been written for you, and you only need to supply the updates to θ within each iteration.

As you program, make sure you understand what you are trying to optimize and what is being updated. Keep in mind that the cost is parameterized by the vector θ not X and y. That is, we minimize the value of by changing the values of the vector θ , not by changing X or y.

A good way to verify that gradient descent is working correctly is to look at the value of and check that it is decreasing with each step. Your value of $J(\theta)$ should never increase, and should converge to a steady value by the end of the algorithm. Other way of making sure your gradient estimate is correct is to check it againts a <u>finite</u> <u>difference (https://en.wikipedia.org/wiki/Finite_difference)</u> approximation.

We also initialize the initial parameters to 0 and the learning rate alpha to 0.01.

```
In [7]:
         import scipy.optimize
         from functools import partial
         def loss gradient(X, y, theta):
             X \text{ prime} = \text{add column}(X)
             loss_grad = ((predict(X, theta) - y)*X_prime).mean(axis=0)[:, np.
               raise NotImplementedError("Compute the model loss gradient; "
         #
                                            "use the predict() function; "
         #
                                            "this also must be vectorized!")
             return loss grad
         assert loss gradient(X, y, theta init).shape == (2, 1)
         def finite diff grad check(f, grad, points, eps=1e-10):
             errs = []
             for point in points:
                 point errs = []
                 grad func val = grad(point)
                  for dim i in range(point.shape[0]):
                      diff v = np.zeros like(point)
                      diff v[dim i] = eps
                      \dim \operatorname{grad} = (f(\operatorname{point}+\operatorname{diff} v) - f(\operatorname{point}-\operatorname{diff} v))/(2*\operatorname{eps})
                      point_errs.append(abs(dim_grad - grad_func_val[dim_i]))
                 errs.append(point errs)
             return errs
         test points = [np.random.rand(2, 1) for in range(10)]
         finite diff errs = finite diff grad check(
             partial(loss, X, y), partial(loss gradient, X, y), test points
         )
         print('max grad comp error', np.max(finite diff errs))
         assert np.max(finite diff errs) < 1e-3, "grad computation error is to
         o large"
         def run gd(loss, loss gradient, X, y, theta init, lr=0.01, n iter=150
         0):
             theta current = theta init.copy()
             loss values = []
             theta_values = []
             for i in range(n iter):
                  loss value = loss(X, y, theta current)
                  theta current = theta current - lr*loss gradient(X, y, theta
         current)
                 loss values.append(loss value)
                 theta values.append(theta current)
             return theta current, loss values, theta values
         result = run gd(loss, loss gradient, X, y, theta init)
         theta_est, loss_values, theta_values = result
         print('estimated theta value', theta est.ravel())
         print('resulting loss', loss(X, y, theta est))
```

```
plt.ylabel('loss')
plt.xlabel('iter_i')
plt.plot(loss_values)
plt.show()
plt.ylabel('log(loss)')
plt.xlabel('iter_i')
plt.semilogy(loss_values)
plt.show()
max grad comp error 1.5638362621750446e-05
estimated theta value [-3.63029144 1.16636235]
resulting loss 4.483388256587726
   30
   25
   20
 055
   15
   10
    5
                         600
                                     1000
        Ò
             200
                   400
                               800
                                          1200
                                                1400
                             iter i
   3 \times 10^{1}
   2 \times 10^{1}
 log(loss)
      10^{1}
   6 \times 10^{\circ}
```

2.2.4 [10pt] After you are finished, use your final parameters to plot the linear fit. The result should look something like on the figure below.

600

800

iter_i

400

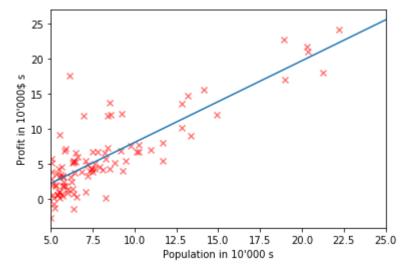
1000

1200

1400

200

```
In [8]: plt.scatter(X, y, marker='x', color='r', alpha=0.5)
    x_start, x_end = 5, 25
    plt.xlim(x_start, x_end)
    X_test = np.array([[5], [25]])
    y_test = predict(X_test, theta_est)
    plt.plot(X_test, y_test)
    plt.xlabel('Population in 10\'000 s')
    plt.ylabel('Profit in 10\'000$ s')
    plt.show()
```

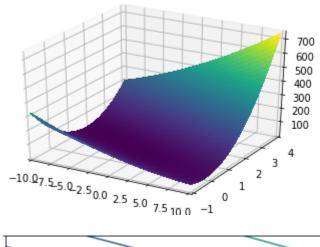


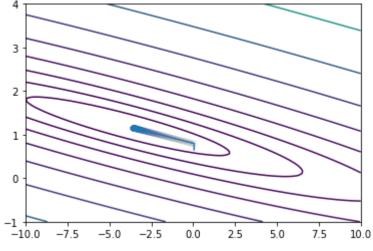
Now use your final values for θ and the predict() function to make predictions on profits in areas of 35,000 and 70,000 people.

```
In [9]: print(predict(np.array([[35000], [70000]]), theta_est))
        [[40819.05197031]
        [81641.73423205]]
```

To understand the cost functionbetter, you will now plot the cost over a 2-dimensional grid of and values. You will not need to code anything new for this part, but you should understand how the code you have written already is creating these images.

```
from mpl toolkits.mplot3d import Axes3D
import matplotlib.cm as cm
# scale all = 1
# scale_x, scale_y = 20*scale all, 5*scale all
# limits = [(theta\ est[0]-scale\ x,\ theta\ est[0]+scale\ x),
            (theta_est[1]-scale_y, theta_est[1]+scale_y)]
limits = [(-10, 10), (-1, 4)]
space = [np.linspace(*limit, 100) for limit in limits]
theta 1 grid, theta 2 grid = np.meshgrid(*space)
theta_meshgrid = np.vstack([theta_1_grid.ravel(), theta_2_grid.ravel
()])
loss_test_vals_flat = (((add\ column(X)\ @\ theta\ meshgrid\ -\ y)**2).mean
(axis=0)/2
loss test vals grid = loss test vals flat.reshape(theta 1 grid.shape)
print(theta 1 grid.shape, theta 2 grid.shape, loss test vals grid.sha
pe)
plt.gca(projection='3d').plot_surface(theta_1_grid, theta_2_grid,
                                       loss_test_vals_grid, cmap=cm.vi
ridis,
                                       linewidth=0, antialiased=False)
xs, ys = np.hstack(theta values).tolist()
zs = np.array(loss values)
plt.gca(projection='3d').plot(xs, ys, zs, c='r')
plt.xlim(*limits[0])
plt.ylim(*limits[1])
plt.show()
plt.contour(theta 1 grid, theta 2 grid, loss test vals grid, levels=n
p.logspace(-2, 3, 20))
plt.plot(xs, ys)
plt.scatter(xs, ys, alpha=0.005)
plt.xlim(*limits[0])
plt.ylim(*limits[1])
plt.show()
```





In [11]: theta_values

```
Out[11]: [array([[0.05839135],
                  [0.6532885 ]]), array([[0.06289175],
                  [0.77000978]]), array([[0.05782293],
                  [0.79134812]]), array([[0.05106363],
                  [0.79572981]]), array([[0.04401438],
                  [0.79709618]]), array([[0.03692413],
                  [0.79792547]]), array([[0.02983712],
                  [0.79865824]]), array([[0.02276118],
                  [0.79937279]]), array([[0.0156977],
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```

```
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```

```
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[ 1.12224687]]), array([[-3.19243049],
[ 1.12237447]]), array([[-3.19369835],
 1.12250184]]), array([[-3.19496392],
[ 1.12262898]]), array([[-3.19622721],
 1.12275589]]), array([[-3.19748822],
[ 1.12288257]]), array([[-3.19874696],
[ 1.12300903]]), array([[-3.20000343],
[ 1.12313525]]), array([[-3.20125764],
[ 1.12326125]]), array([[-3.20250958],
 1.12338702]]), array([[-3.20375927],
```

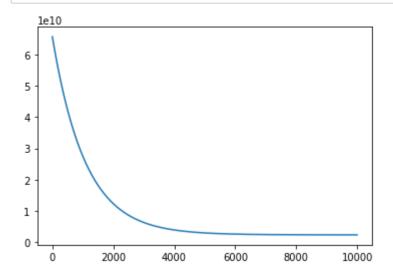
```
[ 1.12351257]]), array([[-3.2050067],
[ 1.12363789]]), array([[-3.20625189],
[ 1.12376298]]), array([[-3.20749483],
[ 1.12388785]]), array([[-3.20873553],
[ 1.12401249]]), array([[-3.209974],
[ 1.1241369]]), array([[-3.21121023],
[ 1.1242611 ]]), array([[-3.21244424],
[ 1.12438507]]), array([[-3.21367602],
[ 1.12450881]]), array([[-3.21490557],
[ 1.12463233]]), array([[-3.21613292],
[ 1.12475563]]), array([[-3.21735805],
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[ 1.12500157]]), array([[-3.21980169],
[ 1.1251242 ]]), array([[-3.22102021],
[ 1.12524662]]), array([[-3.22223653],
[ 1.12536881]]), array([[-3.22345065],
[ 1.12549078]]), array([[-3.22466259],
[ 1.12561253]]), array([[-3.22587235],
[ 1.12573407]]), array([[-3.22707992],
[ 1.12585538]]), array([[-3.22828532],
[ 1.12597648]]), array([[-3.22948855],
[ 1.12609735]]), array([[-3.2306896],
[ 1.12621801]]), array([[-3.23188849],
[ 1.12633845]]), array([[-3.23308522],
[ 1.12645868]]), array([[-3.23427979],
[ 1.12657869]]), array([[-3.23547221],
[ 1.12669848]]), array([[-3.23666248],
[ 1.12681805]]), array([[-3.23785061],
[ 1.12693741]]), array([[-3.23903659],
[ 1.12705656]]), array([[-3.24022043],
[ 1.12717549]]), array([[-3.24140214],
[ 1.1272942 ]]), ...]
```

3. Linear regression with multiple input features

3.1 [20pt] Copy-paste your add_column, predict, loss and loss grad implementations from above and modify your code (if necessarily) of linear regression with one variable to support any number of features (vectorize your code.)

```
data = np.loadtxt('ex1data2.txt', delimiter=',')
X, y = data[:, :-1], data[:, -1, np.newaxis]
n = data.shape[0]
print(X.shape, y.shape, n)
print(X[:10], '\n', y[:10])
(47, 2) (47, 1) 47
[[2.104e+03 3.000e+00]
 [1.600e+03 3.000e+00]
 [2.400e+03 3.000e+00]
 [1.416e+03 2.000e+00]
 [3.000e+03 4.000e+00]
 [1.985e+03 4.000e+00]
 [1.534e+03 3.000e+00]
 [1.427e+03 3.000e+00]
 [1.380e+03 3.000e+00]
 [1.494e+03 3.000e+00]]
 [[399900.]
 [329900.]
 [369000.]
 [232000.]
 [539900.]
 [299900.]
 [314900.]
 [198999.]
 [212000.]
 [242500.]]
```

```
In [14]:
         def add column(X):
              """ Adds a column of ones to a matrix"""
              n = X.shape[0]
              return np.concatenate([X, np.ones((n , 1))], axis=1)
          def predict(X, theta):
              """ Computes h(x; theta) """
              X \text{ prime} = \text{add column}(X)
              return X prime @ theta
          def loss(X, y, theta):
              X \text{ prime} = \text{add column}(X)
              loss = ((predict(X, theta) - y)**2).mean()/2
              return loss
          def loss_gradient(X, y, theta):
              # this also must be vectorized !
              X \text{ prime} = \text{add column}(X)
              loss grad = ((predict(X, theta) - y)*X_prime).mean(axis=0)[:, np.
          newaxis]
              return loss grad
          theta init = np.zeros((3, 1))
          result = run_gd(loss, loss_gradient, X, y, theta_init, n_iter=10000,
          lr=1e-10)
          theta est, loss values, theta values = result
          plt.plot(loss values)
          plt.show()
          # raise NotImplementedError("Put your multivariate regression code he
          re")
```



3.2 [20pt] Draw a histogam of values for the first and second feature. Why is feature normalization important? Normalize features and re-run the gradient decent.

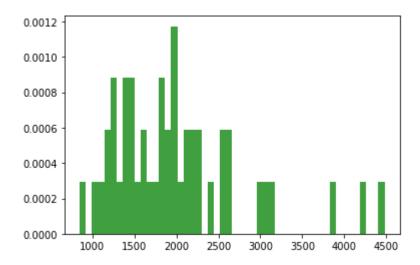
In [15]: plt.hist(X[:, 0], 50, normed=1, facecolor='green', alpha=0.75)
 plt.show()
 plt.hist(X[:, 1], 50, normed=1, facecolor='green', alpha=0.75)
 plt.show()

raise NotImplementedError("Draw histogram and run re-run gd with fe
 ature normalization")

/home/runqi/anaconda3/lib/python3.7/site-packages/ipykernel_launcher.
py:1: MatplotlibDeprecationWarning:

The 'normed' kwarg was deprecated in Matplotlib 2.1 and will be removed in 3.1. Use 'density' instead.

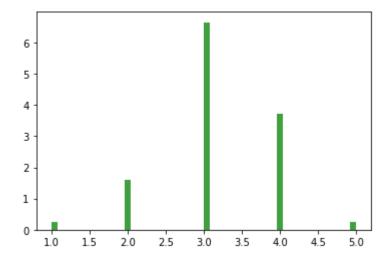
"""Entry point for launching an IPython kernel.



/home/runqi/anaconda3/lib/python3.7/site-packages/ipykernel_launcher.
py:3: MatplotlibDeprecationWarning:

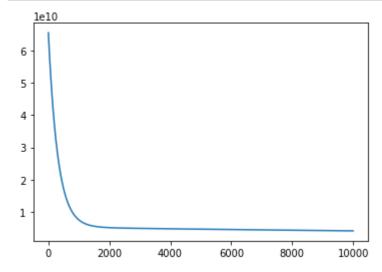
The 'normed' kwarg was deprecated in Matplotlib 2.1 and will be removed in 3.1. Use 'density' instead.

This is separate from the ipykernel package so we can avoid doing i mports until



```
In [16]: theta_init = np.zeros((3, 1))
    X_normed = np.zeros_like(X)
    X_normed[:, 0] = X[:, 0] / X[:, 0].max()
    X_normed[:, 1] = X[:, 1] / X[:, 1].max()
    result = run_gd(loss, loss_gradient, X_normed, y, theta_init, n_iter=
    10000, lr=le-3)
    theta_est, loss_values, theta_values = result

plt.plot(loss_values)
    plt.show()
```



3.3 [10pt] How can we choose an appropriate learning rate? See what will happen if the learning rate is too small or too large for normalized and not normalized cases?

By choosing the best performaning hyperparameter on validation dataset. If learning rate is too small, the model will take a long time to converge to minimum loss. On the other hand, if the learning rate is too large, the loss might be oscillation and might not converge.

4. Written part

These problems are extremely important preparation for the quizzes. Submit solutions to each problem, along with a self-assigned grade. Give yourself full points if you think you have found the correct solution (half points for a partial solution, zero for no attempt.) You may look up definitions online but not the actual answer. Solutions will be released shortly after the due date.

4.1 [10 pt] Maximum Likelihood Estimate for Coin Toss

The probability distribution of a single binary variable that takes value with probability is given by the Bernoulli distribution

$$Bern(x|\mu) = \mu^x (1-\mu)^{1-x}$$

For example, we can use it to model the probability of seeing 'heads' (x=1) or 'tails' (x=0) after tossing a coin, with μ being the probability of seeing 'heads'. Suppose we have a dataset of independent coin flipu $D=\{x^{(1)},\ldots,x^{(m)}\}$ and we would like to estimate μ using Maximum Likelihood. Recall that we can write down the likelihood function as

$$\mathcal{L}(x^{(i)}|\mu) = \mu^{x^{(i)}} (1-\mu)^{1-x^{(i)}} \ P(D|\mu) = \prod_i \mathcal{L}(x^{(i)}|\mu)$$

The log of the likelihood function is

$$\ln p(D|\mu) = \sum_i x^{(i)} \ln \mu + (1-x^{(i)}) \ln (1-\mu)$$

Show that the ML solution for μ is given by $\mu_{ML}=\frac{h}{m}$ where h is the total number of 'heads' in the dataset. Show all of your steps.

4.2 [10 pt] Localized linear regression

Suppose we want to estimate localized linear regression by weighting the contribution of the data points by their distance to the query point x_q , i.e. using the cost

$$E(x_q) = rac{1}{2} \sum_{i}^{m} rac{(y^{(i)} - h(x^{(i)}| heta))^2}{\left|\left|x^{(i)} - x_q
ight|
ight|^2}$$

where $rac{1}{||x^{(i)}-x_q||}=w^{(i)}$ is the inverse Euclidean distance between the training point $x^{(i)}$ and query (test) point x_q .

Derive the modified normal equations (closed form?) for the above cost function $E(x_q)$. (Hint: first, re-write the cost function in matrix/vector notation, using a diagonal matrix to represent the weights $w^{(i)}$).

4.3 [10 pt] Betting on Trick Coins

A game is played with three coins in a jar: one is a normal coin, one has "heads" on both sides, one has "tails" on both sides. All coins are "fair", i.e. have equal probability of landing on either side. Suppose one coin is picked randomly from the jar and tossed, and lands with "heads" on top. What is the probability that the bottom side is also "heads"? Show all your steps.

In []:		

CS542 A1 Machine Learning Problem Set 1 Solutions to Written Problems

4.1 Maximum Likelihood Estimate for Coin Toss

The probability distribution of a single binary variable $x \in \{0,1\}$ that takes value 1 with probability μ is given by the *Bernoulli* distribution

$$Bern(x|\mu) = \mu^{x}(1-\mu)^{1-x}$$

For example, we can use it to model the probability of seeing 'heads' (x = 1) or 'tails' (x = 0) after tossing a coin, with μ being the probability of seeing 'heads'. Suppose we have a dataset of independent coin flips $D = \{x^{(1)}, x^{(2)}, ..., x^{(m)}\}$ and we would like to estimate μ using Maximum Likelihood. Recall that we can write down the likelihood function as

$$p(D|\mu) = \prod_{i=1}^{m} p(x^{(i)}|\mu) = \prod_{i=1}^{m} \mu^{x^{(i)}} (1-\mu)^{1-x^{(i)}}$$

The log of the likelihood function is

$$\ln p(D|\mu) = \sum_{i=1}^{m} x^{(i)} \ln \mu + (1 - x^{(i)}) \ln(1 - \mu)$$

Show that the ML solution for μ is given by $\mu_{ML} = \frac{h}{m}$, where h is the total number of 'heads' in the dataset. Show all of your steps.

Answer: The Maximum Likelihood solution is obtained by finding the value μ_{ML} that maximizes the likelihood function, which is the solution to

$$\frac{d}{d\mu} \left(\sum_{i=1}^{m} x^{(i)} \ln \mu + \left(1 - x^{(i)} \right) \ln(1 - \mu) \right) = 0$$

Taking the derivative, we get

$$\sum_{i=1}^{m} x^{(i)} \frac{1}{\mu} + (-1) (1 - x^{(i)}) \frac{1}{1 - \mu} = 0,$$

$$\sum_{i=1}^{m} x^{(i)} \frac{1}{\mu} = \sum_{i=1}^{m} (1 - x^{(i)}) \frac{1}{1 - \mu},$$

$$\frac{h}{\mu} = \frac{m-h}{1-\mu} , \qquad \mu = \frac{h}{m}$$

4.2 Localized linear regression

Suppose we want to estimate *localized* linear regression by weighting the contribution of the data points by their distance to the query point $x^{(q)}$, i.e. using the cost

$$E(x^{(q)}) = \frac{1}{2} \sum_{i=1}^{m} (y^{(i)} - h_{\theta}(x^{(i)}))^{2} (x^{(i)} - x^{(q)})^{-2}$$

where $(x^{(i)} - x^{(q)})^{-2} = (w^{(i)})^2$ is the inverse Euclidean distance between the training point $x^{(i)}$ and query (test) point $x^{(q)}$.

Derive the modified normal equations for the above cost function $E(x^{(q)})$. (Hint: first, rewrite the cost function in matrix/vector notation, using a diagonal matrix to represent the weights $w^{(i)}$).

Answer: First, re-write the cost as

$$E(\mathbf{x}^{(q)}) = \frac{1}{2} \sum_{i=1}^{m} (y^{(i)} - \boldsymbol{\theta}^T \mathbf{x}^{(i)})^2 (w^{(i)})^2 = \frac{1}{2} \sum_{i=1}^{m} (w^{(i)} y^{(i)} - w^{(i)} \boldsymbol{\theta}^T \mathbf{x}^{(i)})^2$$

Note that, unlike the SSD cost, this cost depends on the query example $x^{(q)}$. Let \mathbf{W} be a diagonal m x m matrix with diagonal entries $w^{(i)}$, i=1,...,m. Let us re-write the above cost equation in terms of matrices and vectors. Notice that multiplying the output vector $\mathbf{Y} = \begin{bmatrix} y^{(1)} & ... & y^{(m)} \end{bmatrix}^T$ by \mathbf{W} results in the vector of weighted outputs,

$$WY = [w^{(1)}y^{(1)} \dots w^{(m)}y^{(m)}]^T$$

and, similarly, multiplying the design matrix by W produces WX, the weighted design matrix. Replacing the original Y and X in the SSD cost with their weighted versions, we can write the weighted cost function (showing vectors/matrices in boldface for clarity) as:

$$E(x^{(q)}) = \frac{1}{2} \sum_{i=1}^{m} (w^{(i)}y^{(i)} - w^{(i)}\theta^{T}x^{(i)})^{2} = (WY - WX\theta)^{2}$$

Substituting Y = WY and X = WX into the normal equations, the solution is

$$\boldsymbol{\theta}_{ML} = \left((WX)^T (WX) \right)^{-1} (WX)^T (WY)$$

4.3 Betting on Trick Coins

A game is played with three coins in a jar: one is a normal coin, one has "heads" on both sides, one has "tails" on both sides. All coins are "fair", i.e. have equal probability of landing on either side. Suppose one coin is picked randomly from the jar and tossed, and lands with "heads" on top. What is the probability that the bottom side is also "heads"? Show all your steps.

Answer: At first, it may seem that the answer is one-half: We know that the tails/tails coin has not been picked, and only one of the remaining two—the heads/heads coin—can have the down-side be heads. However, the correct answer is two thirds, so you should not bet on the intuitive estimate! Let the following variables designate possible outcomes:

BH: bottom side of picked coin is heads

TH: top side of picked coin is heads
H: heads/heads coin is picked, 1/3
T: tails/tails coin is picked, 1/3
N: heads/tails coin is picked, 1/3

Then, using rules of probability

$$\frac{p(BH|TH) =}{\frac{p(BH,TH)}{p(TH)}} = \frac{\frac{p(BH,TH)}{p(TH|N)p(N)}}{\frac{p(BH,TH)}{p(TH|H)p(H) + p(TH|T)p(T) + p(TH|N)p(N)}} = \frac{\frac{1}{3}}{1 * \frac{1}{3} + 0 * \frac{1}{3} + \frac{1}{2} * \frac{1}{3}} = \frac{\frac{1}{3}}{\frac{1}{2}} = \frac{2}{3}$$