Problem Set 2: Classification

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. Then save this file to your computer, run Anaconda and choose this file in Anaconda's file explorer. Use Python 3 version. The statements below 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 and here.

To run code in a cell or to render Markdown+LaTeX 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. Here are some useful resources for Markdown guide and LaTeX tutorial if you are not familiar with the basic syntax.

If a certain output is given for some cells, that means that you are expected to get similar results in order to receive full points (small deviations are fine). For some parts we have already written the code for you. You should read it closely and understand what it does.

Only **PDF** files are accepted for ps1 submission. To print this notebook to a pdf file, you can go to "File" -> "Download as" -> "PDF via LaTex(.pdf)" or simply use "print" in browser.

Total: 85 points (35 (Q1) + 20 (Q2) + 30 (Q3)).

1. Logistic Regression

In this part of the exercise, you will build a logistic regression model to predict whether a student gets admitted into a university.

Suppose that you are the administrator of a university department and you want to determine each applicant's chance of admission based on their results on two exams. You have historical data from previous applicants in *ex2data1.txt* that you can use as a training set for logistic regression. For each training example, you have the applicant's scores on two exams and the admissions decision.

Your task is to build a classification model that estimates an applicant's probability of admission based on the scores from those two exams. This outline and code framework will guide you through the exercise.

1.1 Implementation

```
import sys
import numpy as np
import matplotlib
import matplotlib.pyplot as plt
print('Tested with:')
print('Python', sys.version)
print({x.__name__: x.__version__ for x in [np, matplotlib]})

Tested with:
Python 3.8.8 (default, Apr 13 2021, 15:08:03) [MSC v.1916 64 bit (AMD64)]
{'numpy': '1.20.1', 'matplotlib': '3.3.4'}
```

1.1.1 Visualizing the data

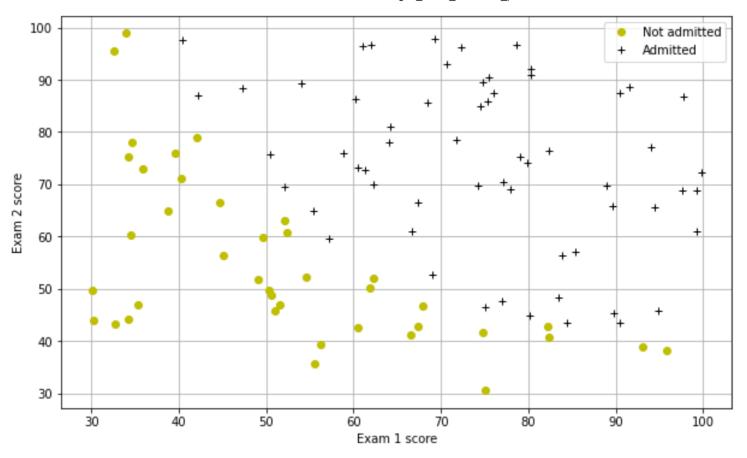
Before starting to implement any learning algorithm, it is always good to visualize the data if possible. This first part of the code will load the data and display it on a 2-dimensional plot by calling the function plotData. The axes are the two exam scores, and the positive and negative examples are shown with different markers.

```
# Try to avoid reassinging/mutating variables because when you encounter an # unexplainable bugs (you will) it is easier to have the whole history # of values to reason about.
# Using %debug magic to run pdb might be useful for debugging.
# It is just like gdb but for Python.
```

```
In [159...
          # it is good to isolate logical parts to avoid variables leaking into the
          # global scope and messing up your logic later in weird ways
          def read classification csv data(fn, add ones=False):
              # read comma separated data
              data = np.loadtxt(fn, delimiter=',')
              X, y = data[:, :-1], data[:, -1, None] # a fast way to keep last dim
              # printing statistics of data before working with it might have saved
              # hundreds hours of of my time, do not repeat my errors :)
              print(X .shape, X .min(), X .max(), X .dtype)
              print(y .shape, y .min(), y .max(), y .dtype)
              # aha, y is float! this is not what we expected
              # what might go wrong with further y == 0 checks?
              # A: floating point equality comparison, that's what!
              # insert the column of 1's into the "X" matrix (for bias)
              X = np.insert(X , X .shape[1], 1, axis=1) if add ones else X
              y = y \cdot astype(np \cdot int32)
              return X, y
          X data, y data = read classification csv data('ex2data1.txt', add ones=True)
          print(X data.shape, X data.min(), X data.max(), X data.dtype)
          print(y data.shape, y data.min(), y data.max(), y data.dtype)
          (100, 2) 30.05882244669796 99.82785779692128 float64
```

```
(100, 2) 30.05882244669796 99.82785779692128 float64 (100, 1) 0.0 1.0 float64 (100, 3) 1.0 99.82785779692128 float64 (100, 1) 0 1 int32
```

```
In [160... # how does the *X[v.ravel()==1, :2].T trick work?
          # https://docs.pvthon.org/3/tutorial/controlflow.html#unpacking-argument-lists
          def plot data(X, v, labels, markers, xlabel, vlabel, figsize=(10, 6), ax=None):
              if figsize is not None:
                  plt.figure(figsize=figsize)
              ax = ax or plt.gca()
              for label id, (label, marker) in enumerate(zip(labels, markers)):
                  ax.plot(*X[v.ravel()==label id, :2].T, marker, label=label)
              ax.set xlabel(xlabel)
              ax.set ylabel(ylabel)
              plt.legend()
              ax.grid(True)
          student plotting spec = {
              'X': X data,
              'y': y data,
              'xlabel': 'Exam 1 score',
              'vlabel': 'Exam 2 score',
              'labels': ['Not admitted', 'Admitted'],
              'markers': ['yo', 'k+'],
              'figsize': (10, 6)
          plot data(**student plotting spec)
          plt.show()
```



1.1.2 [5pts] Sigmoid function

Before you start with the actual cost function, recall that the logistic regression hypothesis is defined as:

$$h_{ heta}(x) = g(heta^T x)$$

where function g is the sigmoid function. The sigmoid function is defined as:

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

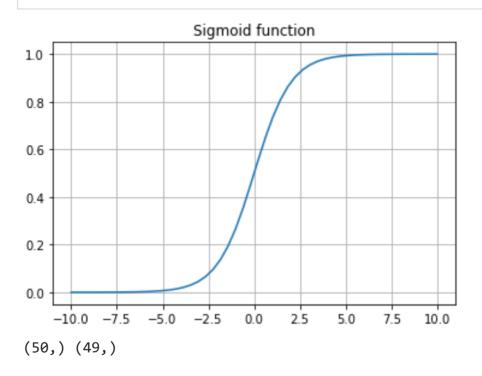
Your first step is to implement/find a sigmoid function so it can be called by the rest of your program. Your code should also work with vectors and matrices. For a matrix, your function should perform the sigmoid function on every element.

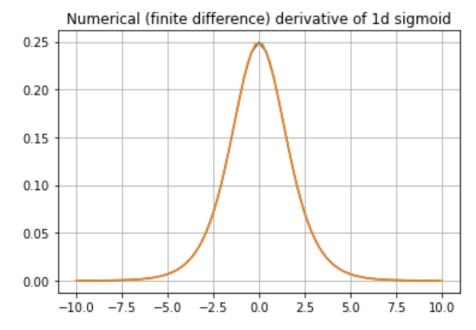
When you are finished, (a) plot the sigmoid function, and (b) test the function with a scalar, a vector, and a matrix. For scalar large positive values of x, the sigmoid should be close to 1, while for scalar large negative values, the sigmoid should be close to 0. Evaluating sigmoid(0) should give you exactly 0.5.

```
In [161...
          # check out scipy.special for great variaty of vectorized functions
          # remember that sigmoid is the inverse of logit function
          # maybe worth checking out scipy.special.logit first
          from scipy.special import expit
           sigmoid = expit
          def check that sigmoid f(f):
               # don't use np.arange with float step because it works as
               # val {i+1} = val i + step while val i < end</pre>
               # what might do wrong with float precision?
               x \text{ test} = \text{np.linspace}(-10, 10, 50)
               sigm test = f(x test)
               plt.plot(x test, sigm test)
               plt.title("Sigmoid function")
               plt.grid(True)
               plt.show()
               # why should analytical diff almost == finite diff for sigmoid?
               analytical diff = sigm test*(1-sigm test)
               finite step = x \text{ test}[1]-x \text{ test}[0]
               finite diff = np.diff(sigm test) / finite step
               print(x test.shape, finite diff.shape)
               plt.plot(x test[:-1]+finite step/2, finite diff)
               plt.plot(x test, analytical diff)
               plt.title("Numerical (finite difference) derivative of 1d sigmoid")
               plt.grid(True)
```

plt.show()

check that sigmoid f(sigmoid)





1.1.3 [15pts] Cost function and gradient

Now you will implement the cost function and gradient for logistic regression. Complete the code in the functions *hyposesis_function* and *binary_logistic_loss* below to return the value of the hypothesis function and the cost, respectively. Recall that the cost function in logistic regression is

$$J(heta) \ = \ rac{1}{m} \ \sum_{i=1}^m \ [\ -y^{(i)}log(h_{ heta}(x^{(i)})) \ - \ (1-y^{(i)})log(1-h_{ heta}(x^{(i)})) \]$$

and the gradient of the cost is a vector of the same length as θ where the j^{th} element (for $j=0,1,\ldots,n$) is defined as follows:

$$rac{\partial J(heta)}{\partial heta_i} \; = \; rac{1}{m} \; \sum_{i=1}^m \; (h_ heta(x^{(i)}) - y^{(i)}) x_j^{(i)}$$

where m is the number of points and n is the number of features. Note that while this gradient looks identical to the linear regression gradient, the formula is actually different because linear and logistic regression have different definitions of $h_{\theta}(x)$.

What should be the value of the loss for $\theta=\bar{0}$ regardless of input? Why? Make sure your code also outputs this value

```
In [162...
          # we are trying to fit a function that would return a
          # "probability of "
          # hyposesis function describes parametric family of functions that we are
          # going to pick our "best fitting function" from. It is parameterized by
          # real-valued vector theta, i.e. we are going to pick
               h best = argmin \{h \in H\} logistic loss h(x, y, h)
          # but because there exist a bijection between theta's and h's it is
          # eavivalent to choosina
               theta best = argmin {theta \in H} logistic loss theta(x, y, theta)
          def hyposesis function(x, theta):
              #raise NotImplementedError('Implement it yourself.')
              thetaTX = np.dot(x, theta)
              #print(thetaTX.shape)
              return sigmoid(thetaTX)
          # negative log likelihood of observing sequence of integer
          # y's given probabilities y pred's of each Bernoulli trial
          # recommentation: convert both variables to float's
          # or weird sign stuff might happen like -1*v != -v for uint8
          # use np.mean and broadcastina
          def binary logistic loss(y, y pred):
              assert v pred.shape == v.shape
              # or weird sign stuff happens! like -1*y != -y
              y, y pred = y.astype(np.float64), y pred.astype(np.float64)
              # When y pred = 0, log(0) = -inf,
              # we could add a small constant to avoid this case
              CONSTANT = 0.000001
              y pred = np.clip(y pred, 0+CONSTANT, 1-CONSTANT)
              #TODO: Calculate the log likelihoods
              #raise NotImplementedError('Implement it yourself.')
              loglikelihood = np.mean((-y)*np.log(y pred) - (1-y) * np.log(1-y pred))
```

```
return loglikelihood
          def logistic loss theta grad(x, y, h, theta):
              Arguments (np arrays of shape):
                   x : [m, n] ground truth data
                   v : [m, 1] ground truth prediction
                   h : [m, n] -> [m, 1] our guess for a prediction function
               .....
              # reshape theta: n bv 1
              theta = theta.reshape((-1,1))
              y \text{ pred} = h(x, \text{ theta})
              point wise grads = (v \text{ pred } - v)*x
              grad = np.mean(point wise grads, axis=0)[:, None]
              assert grad.shape == theta.shape
              return grad.ravel()
          def logistic loss theta(x, y, h, theta):
              # reshape theta: n by 1
              theta = theta.reshape((-1,1))
               return binary logistic loss(y, h(x, theta))
In [163...
          # Check that with theta as zeros, cost is about 0.693:
          theta init = np.zeros((X data.shape[1], 1))
          print(logistic loss theta(X data, y data, hyposesis function, theta init))
          print(logistic loss theta grad(X data, y data, hyposesis function, theta init))
         0.6931471805599453
         [-12.00921659 -11.26284221 -0.1
```

1.1.4 Learning parameters using scipy.optimize

In the previous assignment, you found the optimal parameters of a linear regression model by implementing gradient descent. You wrote a cost function and calculated its gradient, then took a gradient descent step accordingly. This time, instead of taking gradient descent steps, you will use a scipy.optimize built-in function called *scipy.optimize.minimize*. In this case, we will use the *conjugate gradient algorithm*.

The final θ value will then be used to plot the decision boundary on the training data, as seen in the figure below.

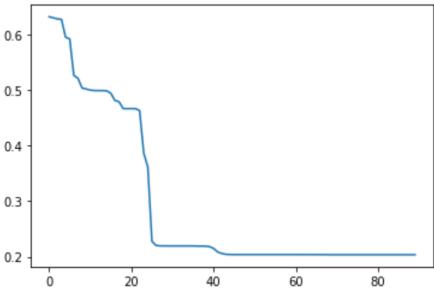
```
In [164...
          import scipy.optimize
          from functools import partial
In [165...
          def optimize(theta init, loss, loss grad, max iter=10000, print every=1000, optimizer fn=None, sho
              theta = theta init.copy()
              opt args = {'x0': theta init, 'fun': loss, 'jac': loss grad, 'options': {'maxiter': max iter}}
              loss curve = []
              def scipy callback(theta):
                  f value = loss(theta)
                  loss curve.append(f value)
              if optimizer fn is None:
                  optimizer fn = partial(scipy.optimize.minimize, method='CG', callback=scipy callback)
              opt result = optimizer fn(**opt args)
              if show:
                  plt.plot(loss curve)
                  plt.show()
              return opt result['x'].reshape((-1, 1)), opt result['fun']
```

loss = partial(logistic loss theta, X data, y data, hyposesis function)

theta init = np.zeros((3, 1))

In [166...

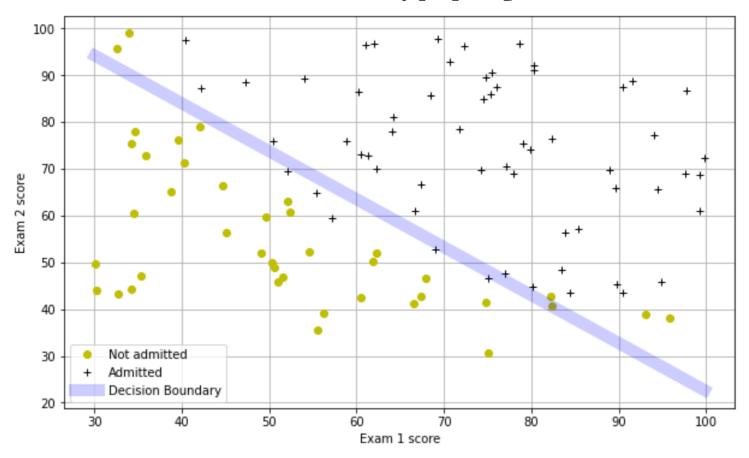
```
loss_grad = partial(logistic_loss_theta_grad, X_data, y_data, hyposesis_function)
theta, best_cost = optimize(theta_init, loss, loss_grad, show=True)
print(best_cost)
```



0.203497703218614

```
# Plotting the decision boundary: two points, draw a line between
# Decision boundary occurs when h = 0, or when
# theta_0*x1 + theta_1*x2 + theta_2 = 0
# y=mx+b is replaced by x2 = (-1/theta1)(theta2 + theta0*x1)

line_xs = np.array([np.min(X_data[:,0]), np.max(X_data[:,0])])
line_ys = (-1./theta[1])*(theta[2] + theta[0]*line_xs)
plot_data(**student_plotting_spec)
plt.plot(line_xs, line_ys, 'b-', lw=10, alpha=0.2, label='Decision Boundary')
plt.legend()
plt.show()
```



1.1.5 [15pts] Evaluating logistic regression

After learning the parameters, you can use the model to predict whether a particular student will be admitted.

(a) [5 pts] Show that for a student with an Exam 1 score of 45 and an Exam 2 score of 85, you should expect to see an admission probability of 0.776.

Another way to evaluate the quality of the parameters we have found is to see how well the learned model predicts on our training set.

(b) [10 pts] In this part, your task is to complete the code in *makePrediction*. The predict function will produce "1" or "0" predictions given a dataset and a learned parameter vector θ . After you have completed the code, the

script below will proceed to report the training accuracy of your classifier by computing the percentage of examples it got correct. You should also see a Training Accuracy of 89.0.

```
In [168...
          # For a student with an Exam 1 score of 45 and an Exam 2 score of 85.
          # you should expect to see an admission probability of 0.776.
          check data = np.array([[45., 85., 1]])
          print(check data.shape)
          print(hyposesis function(check data, theta))
         (1, 3)
         [[0.77631812]]
In [169...
          # use hyposesis function and broadcast compare operator
          def predict(x, theta):
              #raise NotImplementedError('Implement it yourself.')
              prediction = hyposesis function(x, theta)
              prediction = np.apply along axis(lambda x: 0 if x < 0.5 else 1, 1, prediction)
              return prediction
          def accuracy(x, y, theta):
              #raise NotImplementedError('Implement it yourself.')
              count = 0;
              y predict = predict(x, theta)
              for i in range(0, y predict.shape[0]):
                  if y predict[i] == y[i]:
                       count +=1
              accuracy = count / y predict.shape[0]
              return accuracy
          print(accuracy(X data, y data, theta))
```

0.89

2. Regularized logistic regression

In this part of the exercise, you will implement regularized logistic regression to predict whether microchips from a fabrication plant pass quality assurance (QA). During QA, each microchip goes through various tests to ensure it is functioning correctly. Suppose you are the product manager of the factory and you have the test results for some microchips on two different tests. From these two tests, you would like to determine whether the microchips should be accepted or rejected. To help you make the decision, you have a dataset of test results on past microchips in *ex2data2.txt*, from which you can build a logistic regression model.

2.1 Visualizing the data

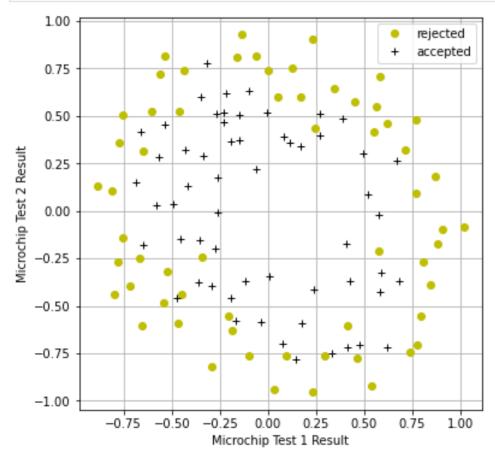
Similar to the previous parts of this exercise, plotData is used to generate the figure below, where the axes are the two test scores, and the positive (y = 1, accepted) and negative (y = 0, rejected) examples are shown with different markers

The figure below shows that our dataset cannot be separated into positive and negative examples by a straight line. Therefore, a straightforward application of logistic regression will not perform well on this dataset since logistic regression will only be able to find a linear decision boundary.

```
In [225...
          X data , y data = read classification csv data('ex2data2.txt')
          X data = X data - X data .mean(axis=0)[None, :]
          print(X data.shape, X data.min(), X data.max(), X data.dtype)
          print(y data.shape, y_data.min(), y_data.max(), y_data.dtype)
          (118, 2) -0.83007 1.1089 float64
          (118, 1) 0.0 1.0 float64
          (118, 2) -0.9528415593220338 1.0161210915254237 float64
          (118, 1) 0 1 int32
In [171...
          chip plotting spec = {
              'X': X data,
               'y': y data,
              'xlabel': 'Microchip Test 1 Result',
               'ylabel': 'Microchip Test 2 Result',
               'labels': ['rejected', 'accepted'],
```

```
'markers': ['yo', 'k+'],
  'figsize': (6, 6)
}

plot_data(**chip_plotting_spec)
plt.show()
```



2.2 Nonlinear feature mapping

One way to fit the data better is to create more features from each data point. In *mapFeature* below, we will map the features into all polynomial terms of x_1 and x_2 up to the sixth power as follows:

As a result of this mapping, our vector of two features (the scores on two QA tests) has been transformed into a 28-dimensional vector. A logistic regression classifier trained on this higher-dimension feature vector will have a more complex decision boundary and will appear nonlinear when drawn in our 2-dimensional plot. While the feature mapping allows us to build a more expressive classifier, it is also more susceptible to overfitting. In the next parts of the exercise, you will implement regularized logistic regression to fit the data and also see for yourself how regularization can help combat the overfitting problem.

Either finite dimentional (or even infinite-dimentional, as you would see in the SVM leacture and the corresponding home assingment) feature mappings are usually denoted by Φ and therefore our hyposesis is now that the Bernoulli probability of chip matfunctioning might be described as

$$p_i = \sigma(\Phi(x_i)^T \theta)$$

```
from itertools import combinations_with_replacement

def polynomial_feature_map(X_data, degree=20, show_me_ur_powers=False):
    assert len(X_data.shape) == 2
    group_size = X_data.shape[1]
```

```
assert group size == 2
    # hm.. how to get all ordered pairs (c, d) of non-negative ints
    # such that their sum is c + d <= deraee?</pre>
    # it is eavivalent to getting all aroups of integers (a, b) such that
    \# 0 \le a \le b \le degree  and defininta c = a. d = b - a
    # their sum is below dearee. both are >= 0
    # then feature i = (x \ 0 \ ^c) * (x \ 1 \ ^d)
    comb iterator = combinations with replacement(range(degree+1), group size)
    not quite powers = np.arrav(list(comb iterator))
    powers bad order = not guite powers.copy()
    powers bad order[:, 1] -= not quite powers[:, 0]
    # Let's reoder them so that lower power monomials come first
    rising power idx = np.argsort(powers bad order.sum(axis=1))
    powers = powers bad order[rising power idx]
    if show me ur powers is True:
        print(powers.T)
        print('total power per monomial', powers.sum(axis=1))
    X with powers = np.power(X data[:, :, None], powers.T[None])
    # tu tu power rangers (with replacement)
    X poly = np.prod(X with powers, axis=1)
    return X poly
X pf = polynomial feature map(X data, show me ur powers=True)
print(X pf.shape)
[[0 \ 0 \ 1 \ 0 \ 2 \ 1 \ 0 \ 2 \ 3 \ 1 \ 4 \ 2 \ 1 \ 3 \ 0 \ 5 \ 3 \ 0 \ 2]
    4 3 1 4 0 1 2 6 7 3 5 2 3 4 6
                                              8
    9 8 0 2 1 5 9 3 5 7 8 4 10 6 1 0
                                    6 2 8 11 10
        8 6 11 3 7 0 1 9 12 4
                                                 3 5
       4 12 2 10 0 11 3 0 12 11 14 10 9 7
                                                 8 13
       9 0 15 14 7 1 10 8 3 2 11 6 12 13
                                               8
                                                 3 11
 10 16 4 0 6 5 1 15 7 12 5 2 13 11 6 15 3 10 7
  4 9 17 5 12 2 4 8 14 13 16 11 9 6 15 0 1 10 3 18 7 17 15 13
 14 17 19 18 16 2 12 3 1 4
                             5 0
                                       8 6 9 10 11 11 18 3 1 17 10
                                   7
  4 16 13 5 0 19 15 12 6 9 7 14 2 8 20]
[0 1 0 2 0 1 3 1 0 2 0 2 3
                                       1 4 0 2 5 3
                         1 0 4 2 6
                                      5 4 2 0 3 1
          9 7 8 4 1 7 5 3 2
                                    6 0 4 9 10 8 11
```

8 5 12 11 3 0 8 6 10 4 1 2

```
4 5 9 1 11 3 13 2 10 14 2 3 0 4 5 7 8 6
10 11 6 15 0 1 8 14
          5 7 12 13 4 9 3 2 8 13
                     5 14
 0 12 16 10 11 15 1 9 5 12 15 4 6 11 2 14 7 10 17 16
13 8 0 13 6 16 14 10 4 5 2 7 9 12 3 18 17 8 15 0 11 1 4 6
   0 1 3 17 7 16 18 15 14 19 12 11 13 10 9 8 9
                      2 17 19
 4 7 15 20 1 5 8 14 11 13 6 18 12 0]]
total power per monomial [ 0 1 1 2 2 2 3 3 3 3 4
  6 6 7 7 7 7 7 7 7 7 8 8 8 8 8 8 8 8
  (118, 231)
```

2.3 Cost function and gradient

Now you will implement code to compute the cost function and gradient for regularized logistic regression. Recall that the regularized cost function in logistic regression is:

$$j(heta) \ = \ [\ rac{1}{m} \ \sum_{i=1}^m \ [\ -y^{(i)}log(h_{ heta}(x^{(i)})) \ - \ (1-y^{(i)})log(1-h_{ heta}(x^{(i)})) \]\] \ + rac{\lambda}{2m} \sum_{j=2}^n heta_j^2$$

Note that you should not regularize the parameter θ_0 (Why not? Think about why that would be a bad idea).

The gradient of the cost function is a vector where the j element is defined as follows (you should understand how to obtain this expression):

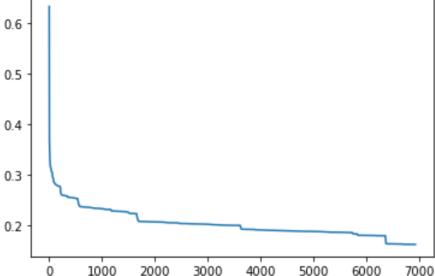
2.3.1 [10pts] Implementing regularized logistic regression

Re-implement computeCost with regularization.

```
In Γ194...
          # Cost function. default Lambda (regularization) 0
          def logistic loss theta w reg(x, y, h, theta, lambda = 0.0):
               #raise NotImplementedError('Implement it vourself.')
              #Loglikelihood = np.mean((-v)*np.log(h(x, theta)) - (1-v) * np.log(1-h(x, theta))) + (lambda / (lambda))
              binary lost = logistic loss theta(x, y, h, theta)
               reg = (lambda /(2*x.shape[0]))*np.sum(theta[1:]**2)
              loglikelihood = binarv lost + reg
              return loglikelihood
          def logistic loss theta w reg grad(x, y, h, theta, lambda = 0.0):
              #raise NotImplementedError('Implement it yourself.')
               binary grad = logistic loss theta grad(x, y, h, theta)
              theta0 = (lambda *theta)/x.shape[0]
              theta0[0] = 0
              \#np.mean((h(x, theta) - y)*x, axis = 0) + theta0/x.shape[0]
              return binary grad + theta0
```

Once you are done, you will call your cost function using the initial value of θ (initialized to all zeros). You should see that the cost is about 0.693

```
3.41810853e-05 7.95432475e-03]
[ 8.47457627e-03 -1.47393694e-03 1.83238652e-02 ... 5.15053246e-04 3.41810853e-05 7.95432475e-03]
...
[ 8.47457627e-03 -1.47393694e-03 1.83238652e-02 ... 5.15053246e-04 3.41810853e-05 7.95432475e-03]
[ 8.47457627e-03 -1.47393694e-03 1.83238652e-02 ... 5.15053246e-04 3.41810853e-05 7.95432475e-03]
[ 8.47457627e-03 -1.47393694e-03 1.83238652e-02 ... 5.15053246e-04 3.41810853e-05 7.95432475e-03]
[ 8.47457627e-03 -1.47393694e-03 1.83238652e-02 ... 5.15053246e-04 3.41810853e-05 7.95432475e-03]]
```



best loss 0.1627064882074352 best acc 0.9322033898305084

2.4 Plotting the decision boundary

To help you visualize the model learned by this classifier, we have provided the function *plotBoundary* which plots the (non-linear) decision boundary that separates the positive and negative examples.

```
def plot_boundary(theta, ax=None):
    Function to plot the decision boundary for arbitrary theta, X, y, lambda value
    Inside of this function is feature mapping, and the minimization routine.
```

```
It works by making a grid of x1 ("xvals") and x2 ("yvals") points,
And for each, computing whether the hypothesis classifies that point as
True or False. Then, a contour is drawn with a built-in pyplot function.
"""

ax = ax or plt.gca()
x_range = np.linspace(-1,1.5,50)
y_range = np.linspace(-1,1.5,50)
xx, yy = np.meshgrid(x_range, y_range)
X_fake = np.stack([xx, yy]).reshape(2, -1).T
X_fake_fm = polynomial_feature_map(X_fake)
y_pred_fake = hyposesis_function(X_fake_fm, theta)
return ax.contour( x_range, y_range, y_pred_fake.reshape(50, 50).T, [0.5])
```

2.4.1 [10pts] Plot Decision Boundaries

- (a) [2 pts] Use *plotBoundary* to obtain four subplots of the decision boundary for the following values of the regularization parameter: $\lambda=0,1,5,10$
- (b) [2 pts] Comment on which plots are overfitting and which plots are underfitting.

Plot 1 is Over fitting

Plot 3 and 4 is underfitting

(c) [2 pts] Which is the model with the highest bias? The highest variance?

Highest bias is model 4

Highest variance is also model 4

(d) [2 pts] What is another way to detect overfitting?

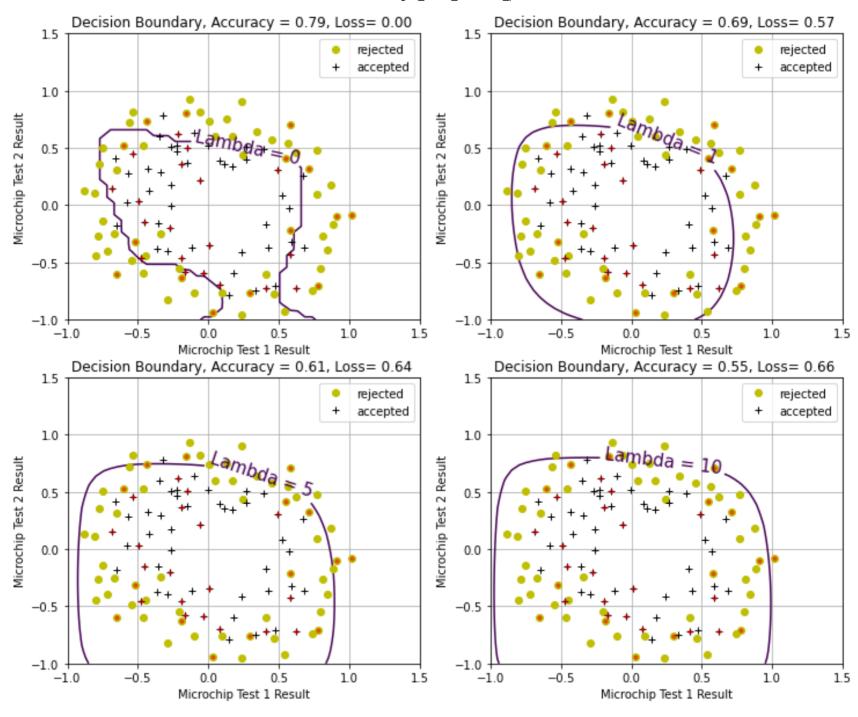
If Loss is 0

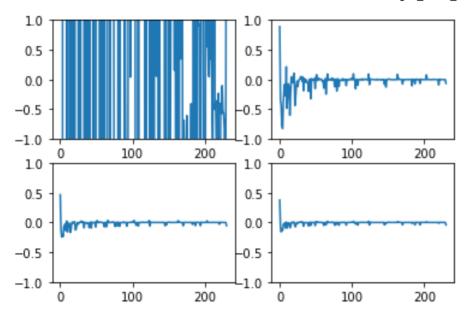
(e) [2 pts] Considering that later components of theta correspond to higher powers of monomials, plot values of theta and commend on effects of regularization

It makes theta magnitude smaller

```
In [292...
          # (a) Build a figure showing contours for various values of regularization parameter. Lambda
          np.random.seed(2)
          train idx mask = np.random.rand(X pf.shape[0]) < 0.3</pre>
          X pf train, y train = X pf[train idx mask], y data[train idx mask]
          X pf test, y test = X pf[~train idx mask], y data[~train idx mask]
          print([x.shape for x in (X pf train, y train, X pf test, y test)])
          def silent optimize w lambda(lambda ):
              theta init = np.zeros((X pf.shape[1], 1))
              data = (X pf train, y train, hyposesis function)
              loss = partial(logistic loss theta w reg, *data, lambda =lambda )
              loss grad = partial(logistic loss theta w reg grad, *data, lambda =lambda )
              theta, final loss = optimize(
                  theta init, loss, loss grad, optimizer fn=None,
                  max iter=1000, print every=0, show=False
              return theta, final loss
          thetas = []
          plt.figure(figsize=(12,10))
          # wow, I mutates an object used in the scope of another function (plot data)
          # don't do that! it is really hard to debug later
          chip plotting spec['figsize'] = None
          # you might find following lines useful:
          #cnt fmt = {0.5: 'Lambda = %d' % Lambda }
          #ax.clabel(cnt, inline=1, fontsize=15, fmt=cnt fmt)
          # red dots indicate training samples
```

```
for id . lambda in enumerate([0, 1, 5, 10]):
    theta, final loss = silent optimize w lambda(lambda)
    #plt.plot(plot boundary(theta))
    thetas.append(theta)
    ax = plt.subplot(2, 2, id +1)
    cnt = plot boundary(theta, ax)
    plot data(**chip plotting spec)
    cnt fmt = {0.5: 'Lambda = %d' % lambda }
    accuracy value = accuracy(X pf test, y test, theta)
    ax.set title('Decision Boundary, Accuracy = %.2f, Loss= %.2f ' %(accuracy value, final loss))
    ax.clabel(cnt, inline=1, fontsize=15, fmt=cnt fmt)
    plt.plot(*X data[train idx mask].T, 'r.', alpha = 0.5)
    #raise NotImplementedError('Implement it yourself.')
plt.show()
# (e) [2 pts] Considering that Later components of theta correspond to higher powers
# of monomials, plot values of theta and commend on effects of regularization
ax = None
for th id, theta in enumerate(thetas):
    ax = plt.subplot(2, 2, th id+1, sharey=ax)
    plt.ylim([-1,1])
    plt.plot(thetas[th id])
    #raise NotImplementedError('Implement it vourself.')
plt.show()
```





3. Written part

These problems are extremely important preparation for the exam. Submit solutions to each problem by filling the markdown cells below (add a cell if necessary).

- 3.1 [10pts] Maximum likelihood for Logistic Regression
- **3.1.1 [2pts]** Suppose we have a dataset D of m points, each consisting of input x and output y, and the probability of observing the output for a given input is $p(y|x;\theta)$. θ refers to the parameters of the distribution p. In your own terms, explain how we can use the principle of Maximum Likelihood to derive a solution for this problem.

$$L(D) = \prod_{i=1}^m p(y|x; heta) = \prod_{i=1}^m p(y^{(i)}|x^{(i)}; heta^{(i)})$$

3.1.2 [8pts] In Logistic Regression, the outputs y are discrete binary variables. The Logistic Regression cost function is given by

$$J(heta) \ = \ -rac{1}{m} \ \sum_{i=1}^m \ [\ y^{(i)}log(h_ heta(x^{(i)})) \ + \ (1-y^{(i)})log(1-h_ heta(x^{(i)})) \]$$

Assume that the probability of y given x is described by:

$$p(|y=1||x|;| heta|)=h_{ heta}(x)$$

$$p(y = 0 \mid x ; \theta) = 1 - h_{\theta}(x)$$

Showing all steps, and in more detail than we went over in class, derive the Logistic Regression cost function using maximum likelihood.

$$L(D) = \prod_{i=1}^m h_{ heta}(x)^{y^{(i)}} (1 - h_{ heta}(x))^{(1-y^{(i)})}$$

Take log both side:

$$egin{align} log L(D) &= log (\prod_{i=1}^m h_{ heta}(x)^{y^{(i)}} (1-h_{ heta}(x))^{(1-y^{(i)})}) \ & log(a.\,b) = log a + log b \ & log a^b = b.\, log a \ \end{gathered}$$

Thus, we have:

$$log L(d) = \sum_{i=1}^{m} y^{(i)} log(h_{ heta}(x)) + (1-y^{(i)}) log(1-h_{ heta}(x))$$

3.2 [10pts] Logistic Regression Classification with Label Noise



Suppose you are building a classifier for images of dogs to classify them into one of two categories $y \in \{0,1\}$, where 0 is "terrier" and 1 is "beagle." Each dog image is represented by a feature vector x consisting of image pixels. You decide to use the logistic regression model $p(y=1 \mid x) = h_{\theta}(x) = \sigma(\theta^T x)$. You collected an image dataset and labeled each image with 0 or 1, however, you are not a dog breed expert and you made some mistakes in assigning labels, so they may be different from true labels. Denote your labels as $t^{(i)}$, and the true labels as $y^{(i)}$, your final dataset is $\mathbf{D} = \{x^{(i)}, t^{(i)}\}$.

You then consulted with a friend who is a dog expert and she estimated that you were correct in about τ fraction of all cases.

3.2.1 [2pts] Write down the equation for the posterior probability $p(t = 1 \mid x)$ of your label being 1 for some point x, in terms of the probability of the true class, $p(y = 1 \mid x)$.

$$egin{aligned} p(t=1 \mid x) &= p(t=1, y=1 \mid x) + p(t=1, y=0 \mid x) \ &= p(t=1 \mid y=1).\, p(y=1 \mid x) + p(t=1 \mid y=0).\, p(y=0 \mid x) \ &= au.\, p(y=1 \mid x) + (1- au).\, (1-p(y=1 \mid x)) \end{aligned}$$

3.2.2 [8pts] Derive the modified cost function in terms of $\theta, x^{(i)}, t^{(i)}$ and au.

$$-rac{1}{m}\sum_{i=1}^{m}[t^{(i)}.\,log(p(t=1\mid x))+(1-t^{(i)})log(1-p(t=1\mid x))$$

$$=-rac{1}{m}\sum_{i=1}^{m}[t^{(i)}.\,log(au.\,p(y=1\mid x)+(1- au).\,(1-p(y=1\mid x)))+(1-t^{(i)})log(1-p(au.\,p(y=1\mid x)+(1- au)))]$$

$$=-rac{1}{m}\sum_{i=1}^{m}[t^{(i)}.log(au.\sigma(heta^{T}x)+(1- au).(1-\sigma(heta^{T}x)))+(1-t^{(i)})log(1-p(au.\sigma(heta^{T}x)+(1- au).(1-\sigma(heta^{T}x))))$$

3.3 [10pts] Cross-entropy loss for multiclass classification

•

This problem asks you to derive the cross-entropy loss for a multiclass classification problem using maximum likelihood. Consider the multiclass classification problem in which each input is assigned to one of K mutually exclusive classes. The binary target variables $y_k \in \{0, 1\}$ have a "one-hot" coding scheme, where the value is 1 for the indicated class and 0 for all others. For example, for K=4 the second-class label is encoded as $[y_1,y_2,y_3,y_4]=[0,1,0,0]$.

Assume that we can interpret the network outputs as $h_k(x,\theta) = p(y_k = 1|x)$, or the probability of the kth class.

Show that the maximum likelihood estimate of the parameters θ can be obtained by minimizing the multiclass cross-entropy loss function

$$L(heta) = -rac{1}{N} \sum_{i=1}^{N} \sum_{k=1}^{K} y_{ik} \log(h_k(x_i, heta))$$

where N is the number of examples $\{x_i, y_i\}$.

Assuming N = 1, where there is only one set of x^i, y^i . Then the maximum likelihood estimate will be:

$$L(D) = \prod_{k=1}^K p(y_k|x) = \prod_{k=1}^K p(y_k=1|x)^{y_k}.\, p(y_k=0|x)^{y_k}.$$

Since y_k is onehot, thus at kth_term, y_k is always 1.

Therefore,

$$L(D) = \prod_{k=1}^K p(y_k = 1|x)^{y_k}.\, p(y_k = 0|x)^{y_k} = \prod_{k=1}^K p(y_k = 1|x)^{y_k}$$

Take log both side, we get:

$$log L(D) = log (\prod_{k=1}^{K} p(y_k = 1|x)^{y_k}) = \sum_{k=1}^{K} y_k log (p(y_k = 1|x)) = \sum_{k=1}^{K} y_k log (h_k(x, heta))$$

With N sets of x^i, y^i we get:

$$logL(D) = \sum_{i=1}^{N} \sum_{k=1}^{K} y_{ik} log(h_k(x_i, heta))$$

In []:			