Training Logistic Regression via Stochastic Gradient Ascent

The goal of this assignment is to implement a logistic regression classifier using stochastic gradient ascent. You will:

- Extract features from Amazon product reviews.
- Convert an SFrame into a NumPy array.
- Write a function to compute the derivative of log likelihood function (with L2 penalty) with respect to a single coefficient.
- Implement stochastic gradient ascent with L2 penalty
- Compare convergence of stochastic gradient ascent with that of batch gradient ascent

If you are doing the assignment with IPython Notebook

An IPython Notebook has been provided below to you for this assignment. This notebook contains the instructions, quiz questions and partially-completed code for you to use as well as some cells to test your code.

Make sure that you are using GraphLab Create 1.8.3. See this post for installing the correct version of GraphLab Create.

What you need to download

If you are using GraphLab Create:

- Download the Amazon product review dataset (subset) in SFrame format. Notice the subset suffix: amazon_baby_subset.gl.zip
- Download the companion IPython notebook: module-10-online-learning-assignment-blank.ipynb
- Download the list of 193 significant words: important words.json
- If you are using Amazon EC2, download the binary files for NumPy arrays: module-10-assignment-numpy-arrays.npz. See the companion notebook for the instructions.
- Save both of these files in the same directory (where you are calling IPython notebook from) and unzip the data file.

If you are not using GraphLab Create:

- If you are using SFrame, download the the Amazon product review dataset (subset) in SFrame format: amazon_baby_subset.gl.zip
- If you are using a different package, download the Amazon product review dataset (subset) in CSV format: amazon_baby_subset.csv.zip
- Download the list of 193 significant words: important_words.json

If you are using GraphLab Create and the companion IPython Notebook

Open the companion IPython notebook and follow the instructions in the notebook.

If you are using other tools

This section is designed for people using tools other than GraphLab Create. You will not need any machine learning packages since we will be implementing decision trees from scratch. We highly suggest you use SFrame since it is open source. In this part of the assignment, we describe general instructions, however we will tailor the instructions for SFrame.

- If you choose to use SFrame, you should be able to follow the instructions in the next section and complete the assessment. All code samples given here will be applicable to SFrame .
- You are free to experiment with any tool of your choice, but some many not produce correct numbers for the quiz questions.

If you are using SFrame

Make sure to download the companion IPython notebook: module-10-online-learning-assignment-blank.ipynb. You will be able to follow along exactly if you replace the first two lines of code with these two lines:

```
import sframe
products = sframe.SFrame('amazon_baby_subset.gl/')
```

After running this, you can follow the rest of the IPython notebook and disregard the rest of this reading.

Note: To install SFrame (without installing GraphLab Create), run

```
pip install sframe
```

If you are NOT using SFrame

Load and process review dataset

1. For this assignment, we will use the same subset of the Amazon product review dataset that we used in Module 3 assignment. The subset was chosen to contain similar numbers of positive and negative reviews, as the original dataset consisted of mostly positive reviews. Load the data file into a data frame products.

Apply text cleaning on the review data

- 2. In this section, we will perform some simple feature cleaning using data frames. The last assignment used all words in building bag-of-words features, but here we limit ourselves to 193 words (for simplicity). We compiled a list of 193 most frequent words into the JSON file named important_words.json . Load the words into a list important_words .
- 3. Let us perform 2 simple data transformations:
- Remove punctuation
- Compute word counts (only for important_words)

We start with the first item as follows:

• If your tool supports it, fill n/a values in the review column with empty strings. The n/a values indicate empty reviews. For instance, Pandas's the fillna() method lets you replace all N/A's in the review columns as follows:

```
products = products.fillna({'review':''}) # fill in N/A's in the review column
```

• Write a function remove_punctuation that takes a line of text and removes all punctuation from that text. The function should be analogous to the following Python code:

```
def remove_punctuation(text):
  import string
  return text.translate(None, string.punctuation)
```

- Apply the remove_punctuation function on every element of the review column and assign the result to the new column review_clean . Note. Many data frame packages support apply operation for this type of task. Consult appropriate manuals.
- 4. Now we proceed with the second item. For each word in important_words , we compute a count for the number of

times the word occurs in the review. We will store this count in a separate column (one for each word). The result of this feature processing is a single column for each word in important_words which keeps a count of the number of times the respective word occurs in the review text.

Note: There are several ways of doing this. One way is to create an anonymous function that counts the occurrence of a particular word and apply it to every element in the review_clean column. Repeat this step for every word in important_words . Your code should be analogous to the following:

```
for word in important_words:
    products[word] = products['review_clean'].apply(lambda s : s.split().count(word))
```

5. After #4 and #5, the data frame products should contain one column for each of the 193 important_words. As an example, the column perfect contains a count of the number of times the word perfect occurs in each of the reviews.

Split data into training and validation sets

6. We will now split the data into a 90-10 split where 90% is in the training set and 10% is in the validation set. We use seed=1 so that everyone gets the same result.

```
train_data, validation_data = products.random_split(.9, seed=1)
```

If you are not using SFrame, download the list of indices for the training and validation sets: module-10-assignment-trainidx.json, module-10-assignment-validation-idx.json. IMPORTANT: If you are using a programming language with 1-based indexing (e.g. R, Matlab), make sure to increment all indices by 1.

Call the training and validation sets train_data and validation_data, respectively.

7. Convert train_data and validation_data into multi-dimensional arrays.

Using the function given in #8 of Module 3 assignment, extract two arrays feature_matrix_train and sentiment_train from train_data. The 2D array feature_matrix_train would contain the content of the columns given by the list important_words. The 1D array sentiment_train would contain the content of the column sentiment. Do the same for validation_data, producing the arrays feature_matrix_valid and sentiment_valid. The code should be analogous to this cell:

```
feature_matrix_train, sentiment_train = get_numpy_data(train_data, important_words, 'sentiment')
feature_matrix_valid, sentiment_valid = get_numpy_data(validation_data, important_words, 'sentiment')
```

Quiz question: In Module 3 assignment, there were 194 features (an intercept + one feature for each of the 193 important words). In this assignment, we will use stochastic gradient ascent to train the classifier using logistic regression. How does the changing the solver to stochastic gradient ascent affect the number of features?

Building on logistic regression

8. Let us now build on Module 3 assignment. Recall from lecture that the link function for logistic regression can be defined as:

$$P(y_i = +1|\mathbf{x}_i, \mathbf{w}) = \frac{1}{1 + \exp(-\mathbf{w}^T h(\mathbf{x}_i))}$$

where the feature vector $h(x_i)$ is given by the word counts of important_words in the review x_i .

We will use the same code as in Module 3 assignment to make probability predictions, since this part is not affected by using stochastic gradient ascent as a solver. Only the way in which the coefficients are learned is affected by using stochastic gradient ascent as a solver. Refer to #10 of Module 3 assignment in order to obtain the function predict_probability .

Derivative of log likelihood with respect to a single coefficient

9. Let us now work on making minor changes to how the derivative computation is performed for logistic regression.

Recall from the lectures and Module 3 assignment that for logistic regression, the derivative of log likelihood with respect to a single coefficient is as follows:

$$\frac{\partial \ell}{\partial w_j} = \sum_{i=1}^N h_j(\mathbf{x}_i) \left(\mathbf{1}[y_i = +1] - P(y_i = +1|\mathbf{x}_i, \mathbf{w}) \right)$$

Recall that, in the Module 3 assignment, we wrote the function feature_derivative to compute the derivative of log likelihood with respect to a single coefficient w_j. Refer to #11 of Module 3 assignment to obtain the function feature derivative .

Note. We are not using regularization in this assignment, but, as discussed in the optional video, stochastic gradient can also be used for regularized logistic regression.

10. To verify the correctness of the gradient computation, we provide a function for computing average log likelihood (which we recall from the last assignment was a topic detailed in an advanced optional video, and used here for its numerical stability).

To track the performance of stochastic gradient ascent, write yourself a function to compute average log likelihood . The average log likelihood is given by the formula

$$\ell\ell_A(\mathbf{w}) = \frac{1}{N} \sum_{i=1}^{N} \left((\mathbf{1}[y_i = +1] - 1) \mathbf{w}^T h(\mathbf{x}_i) - \ln\left(1 + \exp(-\mathbf{w}^T h(\mathbf{x}_i)) - \exp(-\mathbf{w}^T h(\mathbf{x}_i)) \right) \right)$$

Call this function compute_avg_log_likelihood . It should be analogous to the following Python function:

```
def compute_avg_log_likelihood(feature_matrix, sentiment, coefficients):
    indicator = (sentiment==+1)
    scores = np.dot(feature_matrix, coefficients)
    logexp = np.log(1. + np.exp(-scores))

# Simple check to prevent overflow
    mask = np.isinf(logexp)
    logexp[mask] = -scores[mask]

    lp = np.sum((indicator-1)*scores - logexp)/len(feature_matrix)
    return lp
```

Note. We made one tiny modification to the log likelihood function (called compute_log_likelihood) in our earlier assignments. We added a 1/N term which averages the log likelihood accross all data points. The 1/N term makes it easier for us to compare stochastic gradient ascent with batch gradient ascent. We will use this function to generate plots that are similar to those you saw in the lecture.

Quiz question: Recall from the lecture and the earlier assignment, the log likelihood (without the averaging term) is given by

$$\ell\ell(\mathbf{w}) = \sum_{i=1}^{N} \left((\mathbf{1}[y_i = +1] - 1)\mathbf{w}^T h(\mathbf{x}_i) - \ln\left(1 + \exp(-\mathbf{w}^T h(\mathbf{x}_i)\right) \right)$$

How are the functions II(w) and II_A(w) related?

Modifying the derivative for stochastic gradient ascent

11. Recall from the lecture that the gradient for a single data point x_i can be computed using the following formula:

$$\frac{\partial \ell_{\mathbf{i}}(\mathbf{w})}{\partial w_j} = h_j(\mathbf{x}_i) \left(\mathbf{1}[y_i = +1] - P(y_i = +1 | \mathbf{x}_i, \mathbf{w}) \right)$$

Computing the gradient for a single data point

Do we really need to re-write all our code to modify

$$\frac{\partial \ell(\mathbf{w})}{\partial w_i}$$

to

$$\frac{\partial \ell_{\mathbf{i}}(\mathbf{w})}{\partial w_j}$$

? Thankfully No! We access x_i in the training data using feature_matrix_train[i:i+1,:] . Similarly, we access y_i in the training data using sentiment_train [i:i+1]. We can compute

$$\frac{\partial \ell_{\mathbf{i}}(\mathbf{w})}{\partial w_j}$$

by re-using all the code written in feature_derivative and predict_probability .

Carry out the following steps:

- First, compute P(y_i = +1 | x_i, w) using the predict_probability function with feature_matrix_train [i:i+1,:] as the first parameter.
- Next, compute the indicator value 1[y_i = +1] using sentiment_train [i:i+1].
- Finally, call the feature_derivative function with feature_matrix_train [i:i+1,j] as one of the parameters.

The following Python cell carries out the steps for j=1 and i=10. (Try this in your own tool.)

Quiz Question: The code block above computed the derivative

$$\frac{\partial \ell_{\mathbf{i}}(\mathbf{w})}{\partial w_{\mathbf{i}}}$$

for j = 1 and i = 10. Is this a scalar or a 194-dimensional vector?

Modifying the derivative for using a batch of data points

12. Stochastic gradient estimates the ascent direction using 1 data point, while gradient uses N data points to decide how to update the the parameters. In an optional video, we discussed the details of a simple change that allows us to use a mini-batch of B<=N data points to estimate the ascent direction. This simple approach is faster than regular gradient but less noisy than stochastic gradient that uses only 1 data point. Although we encorage you to watch the optional video on the topic to better understand why mini-batches help stochastic gradient, in this assignment, we will simply use this technique, since the approach is very simple and will improve your results.

Given a mini-batch (or a set of data points) $x_i, x_i + 1$, ..., $x_i + 1$, the gradient function for this mini-batch of data points is given by:

$$\sum_{s=i}^{i+B} \frac{\partial \ell_s}{\partial w_j} = \sum_{s=i}^{i+B} h_j(\mathbf{x}_s) \left(\mathbf{1}[y_s = +1] - P(y_s = +1|\mathbf{x}_s, \mathbf{w}) \right)$$

Computing the gradient for a "mini-batch" of data points

We access the points $x_i, x_i + 1, ..., x_$

Write come code to compute

$$\sum_{s=i}^{i+B} \frac{\partial \ell_s}{\partial w_j}$$

for i = 10, j = 1, and B = 10.

Your code should be equivalent to the following Python cell:

Quiz Question: The code block above computed

$$\sum_{s=i}^{i+B} \frac{\partial \ell_s}{\partial w_j}$$

for j = 10, i = 10, and B = 10. Is this a scalar or a 194-dimensional vector?

Quiz Question: For what value of B is the term

$$\sum_{s=1}^{B} \frac{\partial \ell_s(\mathbf{w})}{\partial w_j}$$

the same as the full gradient

$$\frac{\partial \ell(\mathbf{w})}{\partial w_i}$$

?

Averaging the gradient across a batch

13. It is a common practice to normalize the gradient update rule by the batch size B:

$$\frac{\partial \ell_{\mathbf{A}}(\mathbf{w})}{\partial w_j} \approx \frac{1}{B} \sum_{s=i}^{i+B} h_j(\mathbf{x}_s) \left(\mathbf{1}[y_s = +1] - P(y_s = +1 | \mathbf{x}_s, \mathbf{w}) \right)$$

In other words, we update the coefficients using the average gradient over data points (instead of using a summation). By using the average gradient, we ensure that the magnitude of the gradient is approximately the same for all batch sizes. This way, we can more easily compare various batch sizes of stochastic gradient ascent (including a batch size of all the data points), and study the effect of batch size on the algorithm as well as the choice of step size.

Implementing stochastic gradient ascent

14. Now we are ready to implement our own logistic regression with stochastic gradient ascent. The function logistic_regression_SG should accept the following parameters:

• feature_matrix : 2D array of features

• sentiment: 1D array of class labels

• initial_coefficients : 1D array containing initial values of coefficients

• step_size: a parameter controlling the size of the gradient steps

• batch_size : size of mini-batch

• max_iter: number of iterations to run stochastic gradient ascent

The function should return the final set of coefficients, along with the list of log likelihood values over time. (In practice, the final set of coefficients is rarely used; it is better to use the average of the last K sets of coefficients instead, where K should be adjusted depending on how fast the log likelihood oscillates around the optimum.)

The function logistic_regression_SG carries out the steps shown in the following pseudocode:

```
* Create an empty list called log_likelihood_all
* Initialize coefficients to initial_coefficients
* Set random seed = 1
* Shuffle the data before starting the loop below
* Set i = 0, the index of current batch
* Run the following steps max_iter times, performing linear scans over the data:
  * Predict P(y_i = +1|x_i,w) using your predict_probability() function
    Make sure to slice the i-th row of feature_matrix with [i:i+batch_size,:]
  * Compute indicator value for (y_i = +1)
   Make sure to slice the i-th entry with [i:i+batch_size]
  * Compute the errors as (indicator - predictions)
  * For each coefficients[j]:
    - Compute the derivative for coefficients[j] and save it to derivative.
     Make sure to slice the i-th row of feature_matrix with [i:i+batch_size,j]
    - Compute the product of the step size, the derivative, and (1./batch_size).
    - Increment coefficients[j] by the product just computed.
 * Compute the average log likelihood over the current batch.
   Add this value to the list log_likelihood_all.
  \ensuremath{^{*}} Increment i by batch_size, indicating the progress made so far on the data.
  st Check whether we made a complete pass over data by checking
    whether (i+batch_size) exceeds the data size. If so, shuffle the data. If not, do nothing.
* Return the final set of coefficients, along with the list log_likelihood_all.
```

At the end of the day, your logistic_regression_SG function should be analogous to this Python function:

```
def logistic_regression_SG(feature_matrix, sentiment, initial_coefficients, step_size, batch_size, max_iter):
   log_likelihood_all = []
   # make sure it's a numpy array
   coefficients = np.array(initial_coefficients)
   # set seed=1 to produce consistent results
   np.random.seed(seed=1)
   # Shuffle the data before starting
   permutation = np.random.permutation(len(feature_matrix))
   feature_matrix = feature_matrix[permutation,:]
   sentiment = sentiment[permutation]
   i = 0 # index of current batch
   # Do a linear scan over data
   for itr in xrange(max_iter):
       # Predict P(y_i = +1|x_i,w) using your predict_probability() function
       # Make sure to slice the i-th row of feature_matrix with [i:i+batch_size,:]
       ### YOUR CODE HERE
       predictions = ...
       # Compute indicator value for (y_i = +1)
       # Make sure to slice the i-th entry with [i:i+batch_size]
       ### YOUR CODE HERE
        indicator = ...
       # Compute the errors as indicator - predictions
        errors = indicator - predictions
        for j in xrange(len(coefficients)): # loop over each coefficient
           # Recall that feature_matrix[:,j] is the feature column associated with coefficients[j]
           # Compute the derivative for coefficients[j] and save it to derivative.
           # Make sure to slice the i-th row of feature_matrix with [i:i+batch_size,j]
           ### YOUR CODE HERE
                  # Compute the product of the step size, the derivative, and
           # the **normalization constant** (1./batch_size)
           ### YOUR CODE HERE
           coefficients[j] += ...
       # Checking whether log likelihood is increasing
```

```
# Print the log likelihood over the *current batch*
    lp = compute_avg_log_likelihood(feature_matrix[i:i+batch_size,:], sentiment[i:i+batch_size],
                                    coefficients)
    log_likelihood_all.append(lp)
    if itr <= 15 or (itr <= 1000 and itr % 100 == 0) or (itr <= 10000 and itr % 1000 == 0) \
     or itr % 10000 == 0 or itr == max_iter-1:
        data_size = len(feature_matrix)
        print 'Iteration %*d: Average log likelihood (of data points [%0*d:%0*d]) = %.8f' % \
            (int(np.ceil(np.log10(max_iter))), itr, \
             int(np.ceil(np.log10(data_size))), i, \
             int(np.ceil(np.log10(data_size))), i+batch_size, lp)
    # if we made a complete pass over data, shuffle and restart
    i += batch size
    if i+batch_size > len(feature_matrix):
        permutation = np.random.permutation(len(feature_matrix))
        feature_matrix = feature_matrix[permutation,:]
        sentiment = sentiment[permutation]
        i = 0
# We return the list of log likelihoods for plotting purposes.
return coefficients, log_likelihood_all
```

Compare convergence behavior of stochastic gradient ascent

15. For the remainder of the assignment, we will compare stochastic gradient ascent against batch gradient ascent. For this, we need a reference implementation of batch gradient ascent. But do we need to implement this from scratch?

Quiz Question: For what value of batch size B above is the stochastic gradient ascent function logistic_regression_SG act as a standard gradient ascent algorithm?

Running gradient ascent using the stochastic gradient ascent implementation

16. Instead of implementing batch gradient ascent separately, we save time by re-using the stochastic gradient ascent function we just wrote — to perform gradient ascent , it suffices to set batch_size to the number of data points in the training data. Yes, we did answer above the quiz question for you, but that is an important point to remember in the future:)

Small Caveat . The batch gradient ascent implementation here is slightly different than the one in the earlier assignments, as we now normalize the gradient update rule.

We now run stochastic gradient ascent over the feature_matrix_train for 10 iterations using:

- initial_coefficients = np.zeros(194)
- step_size = 5e-1
- batch_size = 1
- max iter = 10

Quiz Question . When you set batch_size = 1, as each iteration passes, how does the average log likelihood in the batch change?

17. Now run batch gradient ascent over the feature_matrix_train for 200 iterations using:

- initial_coefficients = np.zeros(194)
- step_size = 5e-1
- batch_size = number of rows in feature_matrix_train
- max_iter = 200

Quiz Question . When you set batch_size = len(train_data), as each iteration passes, how does the average log likelihood in

the batch change?

Make "passes" over the dataset

18. To make a fair comparison betweeen stochastic gradient ascent and batch gradient ascent, we measure the average log likelihood as a function of the number of passes (defined as follows):

$$[\# \text{ of passes}] = \frac{[\# \text{ of data points touched so far}]}{[\text{size of dataset}]}$$

Quiz Question Suppose that we run stochastic gradient ascent with a batch size of 100. How many gradient updates are performed at the end of two passes over a dataset consisting of 50000 data points?

Log likelihood plots for stochastic gradient ascent

19. With the terminology in mind, let us run stochastic gradient ascent for 10 passes. We will use

- step_size =1e-1
- batch_size =100
- initial_coefficients set to all zeros.

20. Write yourself a function to generate a plot of the average log likelihood as a function of the number of passes. The function should accept the following parameters:

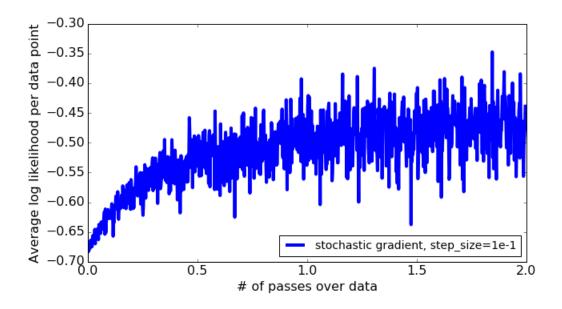
- log_likelihood_all , the list of average log likelihood over time
- len_data, number of data points in the training set
- batch_size , size of each mini-batch
- smoothing_window , a parameter for computing moving averages

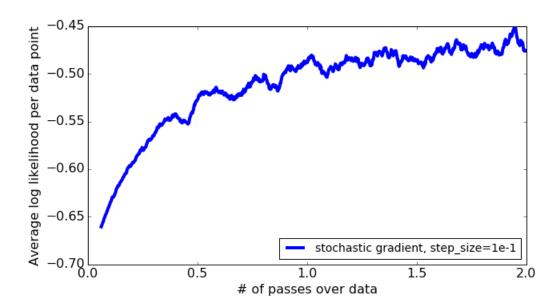
The function should first compute moving averages of log_likelihood_all . To do this efficiently, convolve log_likelihood_all with the vector of length smoothing_window that is filled with the value 1/smoothing_window .

The function then plot the moving averages over the number of passes over the data. Use len_data and batch_size to convert iteration number of (fractional) number of passes.

Using matplotlib, the plot-making function would be as follows:

The resulting plot should look like one of the two figures below:





Stochastic gradient ascent vs batch gradient ascent

21. To compare convergence rates for stochastic gradient ascent with batch gradient ascent, we call make_plot () multiple times.

We are comparing:

- stochastic gradient ascent : step_size = 0.1, batch_size=100
- batch gradient ascent : step_size = 0.5, batch_size=[# rows in feature_matrix_train]

Write code to run stochastic gradient ascent for 200 passes using:

- step_size =1e-1
- batch_size =100
- initial_coefficients set to all zeros.

For batch gradient ascent, use the results obtained from #17.

22. We compare the convergence of stochastic gradient ascent and batch gradient ascent by calling the make_plot function. Apply smoothing with smoothing_window =30.

Quiz Question: In the figure above, how many passes does batch gradient ascent need to achieve a similar log likelihood as stochastic gradient ascent?

Explore the effects of step sizes on stochastic gradient ascent

23. In previous sections, we chose step sizes for you. In practice, it helps to know how to choose good step sizes yourself.

To start, we explore a wide range of step sizes that are equally spaced in the log space. Run stochastic gradient ascent with step_size set to 1e-4, 1e-3, 1e-2, 1e-1, 1e0, 1e1, and 1e2. Use

- initial_coefficients =np.zeros(194)
- batch_size =100
- max_iter initialized so as to run 10 passes over the data.

Plotting the log likelihood as a function of passes for each step size

- 24. Now, we will plot the change in log likelihood using the make_plot for each of the following values of step_size:
- step_size = 1e-4
- step_size = 1e-3
- step_size = 1e-2
- step_size = 1e-1
- step_size = 1e0
- step_size = 1e1
- step_size = 1e2

For consistency, use smoothing_window =30.

Quiz Question: Which of the following is the worst step size? Pick the step size that results in the lowest log likelihood in the end.

Quiz Question: Which of the following is the best step size? Pick the step size that results in the highest log likelihood in the end.