## Peer Review 2 for Jared Fisher

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October 13, 2016

### 1 Introduction

Hi Jared,

Thanks for letting me review your implementation of stochastic gradient descent. Thanks for meeting with me. In this review I will provide my overall impressions of comments on both your written report and the R code you provided on you github. This will essenitially be a written version of what we discussed today after class.

# 2 Review of Report

### 2.1 Linear Regression

First off, this section should really be called "Linear Weighted Regression" or something similar. Since we are dealing with matrices and vectors throughout this problem, you don't need to bolden all you variables for the matrices and vectors. Doing so takes extra time and also can lead to confusion. For example, you bolden both y but not  $\beta$  even though they are both vectors. On the bottom of page 1 you say "Taking the partial derivative with respect to  $\beta$  ...," however, we are taking the  $\operatorname{gradient}$  of this expression since  $\beta$  is a vector. Additionally the notation  $\frac{\partial}{\partial \beta}$  is invalid and so you should instead use  $\nabla_{\beta}$  throughout. However, besides these nitpicks your derivation of  $\hat{\beta}$  is very clear.

I suggest trying to code up the SVD and Cholesky decompositions in addition to the QR decomposition. Also, why is Cholesky more unstable than QR? Expand on that. Your pseudocode should look a little more like actual

R code. Here's a generic example:

```
Data: this text

Result: how to write algorithm with LATEX2e initialization;

while not at end of this document do

read current;

if understand then

go to next section;

current section becomes this one;

else

go back to the beginning of current section;

end

end
```

There are several packages you can use to do this. You can learn more about incorporating pseudocode on the LATEX Wikibooks page for Algorithms.

You do a good job of including your R code inline with the rest of your report. However, I would still recommend using syntax highlighting which will make your code more readable at a first glance. Look into the listings package within LATEX.

In your section talking about sparse matrices, perhaps expand a bit more on how exactly R handles spare matrices. I believe that R does this by treating a matrix as an element list of non-zero entries. Also, I did not see any benchmarking done across your various solve functions. This is key to understanding the comparative computational advantages of your functions.

#### 2.2 GLM

I noticed that you have only shown the derivation of the gradient of the likelihood, but I will still review what you have. You can use the \ln command within your LATEX code to have it appear as "ln" instead of "ln". It is generally understood that logarithms are usually natural logarithms unless otherwise noted, so it is more common to see log used instead of ln. Also,  $e^{-x_i^T\beta}$  looks pretty congested. Use \exp(-x\_ i^T\beta) instead to have it appear as  $\exp(-x_i^T\beta)$ . If you use the align environment (read about it on the LATEX Wikibooks page for Advanced Mathematics) then you can align all your equations so that the equal signs all line up. The rest of this section looks good though. I like that you apply the chain rule by first finding the

gradient of  $w_i$  and then using that in finding the full gradient of the log-likelihood.

### 3 Review of R Code

Overall, your R code looks very clean and organized. The spacing is done well so that it is easy to read. A big help, though, would be to include plenty of comments so that a reader can get through your code more efficiently.

You sometimes put a lot of matrix commands in one line. For example, in one line you write betahat = solve(t(X) %\*% W %\*% X) %\*% t(X) %\*% W %\*% y. I'm not sure, but it may be more computationally efficient to break this off into two lines. Do some benchmarking with matrix operations and see what works most efficiently. This page has some general tips on the best ways to perform matrix multiplication in R. For example, crossprod(X,y) is faster than t(X) %\*% y, and A %\*% (B %\*% y) is faster than A %\*% B %\*% y.

Within your function for solving sparse matrices, be sure that the first thing you do is redefine X as  $X \leftarrow \text{matrix}(X, \text{sparse} = T)$ . Otherwise R won't treat X as sparse!

I like the cleanness in how you define your functions. For example, you define a sigmoid function to make calculations of  $w_i$  easier later on. It looks very nice because it makes your subsequent functions less cluttered. It's a good application of function-oriented programming.

There is a very easy way to turn Y into a series of zeros and ones instead of writing an inefficient for loop. Just do Y <- Y == "B". Technically this will be a vector of Boolean values (TRUE and FALSE) but they are still handled mathematically as zeros and ones. You may find it useful to read up on how R handles logicals on the documentation page for Logic in R.

When you are going through the iterations of a for loop and updating a list for each iteration, it is more efficient to instead define a list before your loop of NA values with the same length as the number of iterations, and then changing the *i*th entry in your list during the *i*th iteration of your loop. You can do this with mylist <- rep(NA, n.iterations). This is because it is

inefficient in R to expand a data structure in a loop. Preallocation avoids this pitfall.

# 4 Conclusion

I hope this review has been helpful. As always, you can reach out to me via email or in my office if you would like to further discuss this assignment. If you update report or R code, possibly to include more about the second problem, I am happy to review that as well.