

# Finding the Higgs boson

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**Abstract**—The Higgs boson is an elementary particle in the Standard Model of physics which explains why other particles have mass. The Higgs boson decays rapidly into other particles and we observe the resulting signature to determine if it was a Higgs boson or some other particle / process. By using binary classification techniques, we can estimate the likelihood that a given signature was the result of a Higgs boson. This will be helpful to narrow the region where the Higgs is expected to be from the overwhelming number of noise in the dataset.

## I. INTRODUCTION

TODO: Find intro, will probably need to have finished the other sections before.

## II. MODELS AND METHODS

### A. Background, Basic methods

1) *general\_gradient\_descent*: is a function who perform the gradient descent algorithm for *any* gradient and loss formula. it's just a loop of a certain length, at each iteration we compute the gradient and move a little bit our weight vector along this gradient. the argument are:

a) *y,tx,w*: namely, the output vector, the input feature matrix and the initial weight vector from where we'll start the gradient descent loop

b) *max\_iter*: the number of gradient descent iteration we'll do

c) *gamma*: the coefficient by which will add the gradient to our current vector  $w^{t+1} = w^t + \text{gamma} * \nabla L(w^t)$

d) *grad\_function*: the function who will return the gradient at point *w* given (*y,tx,w*)

e) *cost\_function*: the function who will return the cost at *w* given (*y,tx,w*)

2) *general\_stochastic\_gradient\_descent*: is the same as the function *general\_gradient\_descent* but it perform a stochastic gradient descent loop, this mean only compute correction on part of the sample. all argument are the same as above, with same meaning except *grad\_function* is gone and we have two new ones.

a) *batch\_size*: the size of the subset on which we'll compute the "partial gradient" (the *g* correction term in the course notes)

b) *stock\_grad\_function*: the function who will compute the stochastic gradient given a subset of *y*, a corresponding subset of *tx* and the current weight vector *w*. selection of the subset is done inside the function with the help of the *batch\_iter* function from the labs exercises

3) *least\_squares\_GD*: perform a gradient descent with cost function MSE and gradient function the gradient of the MSE function. this is the definition of least square GD. argument have same meaning as their homonyms described above

4) *least\_squares\_SGD*: perform a stochastic gradient descent with cost function MSE and gradient function the gradient of the MSE function. this is the definition of least square GD. argument have same meaning as their homonyms described above

5) *least\_squares*: solve directly the least square problem using linear algebra, we use numpy linear system solve function to solve the system presented in the course this way as a better accuracy and less rounding error than if we directly use the solution formula who need to compute an inverse

6) *ridge regression*: same as the one before expect that we introduce a regularization term in the equation. again we solve the linear system derived in course.

7) *logistic regression*: perform logistic regression which is a gradient descent with special cost function and associate gradient function. that what we do we call *general\_gradient\_descent* with the new cost and gradient function

8) *reg logistic regression*: same as the previous one but with a regularization term in the cost function. (and so also in gradient function). As *general\_gradient\_descent* only call the cost and grad function with the argument *y,tx* and *w* we need to fix the *lambda\_* parameter before passing the cost and gradient function to *general\_gradient\_descent*. this is done using the partial function of python

### B. Parameters

There are multiple parameters in the project which can be tuned.

a) *Gamma*: is the parameter describing the step of the gradient descent and of the stochastic gradient descent. At each iteration, the new weight will be calculated by adding to the current weights the derivative of the weights times gamma. It mustn't be too high or we will go in the direction of the good weights but miss them and go further. And at each iteration, we will further ourselves from the weights. If it is too small, we will eventually reach the best weights, but it will take a lot of time. Thus, we searched iteratively for the smallest gamma for which we got a reasonable computing time. Least square GD and Least square SGD :  $\text{gamma}=0.000003$ . Logistic regression and regularized logistic regression :  $0.00003$  Max iterations: this parameter describes how many steps will we take before stopping the algorithm. The higher, the more finely tuned the weights will get, and the smaller, the shorter the

algorithm will be. Least square GD and Least square SGD  
: max iterations=1000. Logistic regression and regularized  
logistic regression : max iterations=800

b) *Lambda*: is the parameter used to avoid overfitting. For this, we did a cross-validation (with 4 folds) and check the test error. We tried it for 30 different lambdas and we saw that there was no lambdas for which the test error was extremely different from the train error. We can then conclude that we don't have overfitting. This is also because we use linear models. So  $\lambda=0$  Which function we will use to estimate the weights. We decided to use the least square method. We chose this before the least square GD and SGD since it yielded a better mse (0.3423 against 0.3616). We chose Least square over logistic regression, since when we applied the weights to the input data, classified it, and compared it to the original y, we saw that Least square was more effective than logistic regression (64695 misclassifications against 68921).

c) *IX itself*: we've tried different preprocessing on the feature before performing machine learning on them. We tried to normalize outliers (feature outside of its 95% interval for each feature. TO COMPLETE

we also tried to rescale the data. this mean that all feature will be in mapped  $[0, 1]$  but with conservation of the relative distance between them. for each feature set we took the max and min and define new value  $x'$  from old value  $x$  by  $x' = \frac{x - \min}{\max - \min}$ . this is know to improve convergence and efficacy of gradient descend algorithm. But even with this the direct least square computation was still the better way.

### C. Final implementation

TODO: When finalized describe here the final implementation with the reasoning behind it.

## III. RESULTS

TODO: Present our results.

## IV. DISCUSSION

TODO: Explain our results strength and weaknesses.

## V. SUMMARY

TODO: Conclusion on how this will be helpful in this research area.