Stochastic Gradient Descent with Momentum and Line Searches

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Abstract

In recent years, tailored line search approaches have proposed to define the step-size, or learning rate, in SGD-type algorithms for finite-sum problems. In particular, a stochastic extension of standard Armijo line search has been proposed in Vaswani, Mishkin, Laradji et al. [1]. The development of this kind of techniques is relevant, because it shall allow to enforce a stronger converging behaviour (due to the Armijo condition), similar to that of standard GD, within SGD methods that are commonly employed with large scale training problems.

However, the stochastic line search is not immediately employable when the momentum term is part of the update equation, as the search direction might not be a descent direction (which is a necessary condition for the Armijo condition). This problem is addressed in Fan, Vaswani, Thrampoulidis *et al.* [2], where a strategy is proposed to guarantee the descent property with momentum.

Contents

1	Inti	roduction	2
	1.1	Classification task	2
	1.2	Optimization problem	3
2	Sto	chastic gradient descent variants	4
	2.1	chastic gradient descent variants Basic stochastic gradient descent	5
		Adding momentum term	
3	Exp	periments and results discussion	9
	3.1	Solving the optimization problem	
		Performance of the objective function	
	Ref	erences	17

1 Introduction

Different SGD-type algorithms proposed by the literature were implemented and tested on different benchmark datasets for training the ℓ_2 -regularized Logistic Regression model.

For the purpose of this work, those algorithms were grouped into one, see algorithm 5 on page 8, follows a list of the variants

- SGD with fixed or decreasing step-size, and line search, see section 2.1 on page 5;
- SGD with momentum term and line search, see section 2.2 on page 6.

This section describes the Machine Learning (ML) problem and the related optimization problem, then section 2 on page 4 summarizes the approaches proposed from the retrieved papers. Section 3 on page 9 describes the experiments performed for showing the performance of the algorithms on different datasets.

1.1 Classification task

Given a dataset as follows

$$\mathcal{D} = \{ (x^{(i)}, y^{(i)}) \mid x^{(i)} \in \mathcal{X}, y^{(i)} \in \mathcal{Y}, i = 1, 2, \dots, N \}$$

the general machine learning optimization problem in the context of supervised learning is

$$\min_{w} f(w) = L(w) + \lambda \Omega(w) \longrightarrow \begin{cases} L(w) = \frac{1}{N} \sum_{i=1}^{N} \ell_i(w) \\ \Omega_{\ell_2} = \frac{1}{2} ||w||_2^2 \end{cases}$$

where L(w) is the loss function which for scaling issues is dived by the total number of samples in the dataset and $\Omega(w)$ is the regularization term with its coefficient λ . There are three regularization possible choices, the ℓ_2 regularization was chosen for the problem that we want to address. The vector w contains the model weights associated to the dataset features.

The task performed is the binary classification (so the allowed values for the response variable are $\mathcal{Y} = \{-1, 1\}$), using the Logistic Regression model. The selected loss function is the log-loss, for one dataset sample is

$$\ell_i(w) = \log(1 + \exp(-y^{(i)}w^T x^{(i)})) \tag{1}$$

figure 1a on the next page shows a plot of the loss function $\ell(uv) = \log(1 + \exp(-uv))$ where $u = y^{(i)}$ and $v = w^T x^{(i)}$.

Prediction

The sigmoid function, see figure 1b on the following page, is used for predicting the labels (positive or negative class) of unseen samples as follows

$$y^{(i)} = \begin{cases} 1 & \text{if } \sigma(w^T x^{(i)}) > 0.5\\ -1 & \text{if } \sigma(w^T x^{(i)}) \le 0.5 \end{cases}$$

the threshold is set according to the Bayes classifier.

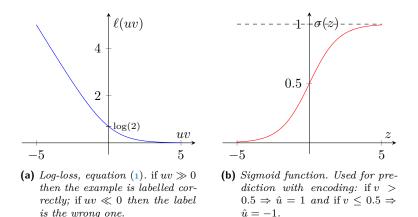


Figure 1: Loss function and sigmoid

1.2 Optimization problem

Putting together the loss function and the regularization term, we can obtain the optimization problem that we want to solve using the Stochastic Gradient Descent (SGD) algorithm variants

$$\min_{w \in \mathbb{R}^{(p+1)}} f(w) = \frac{1}{N} \sum_{i=1}^{N} \log(1 + \exp(-y^{(i)} w^T x^{(i)})) + \lambda \frac{1}{2} ||w||^2$$
 (2)

where $i=1,\ldots,N$ are the dataset samples, $\mathcal{X}\subseteq\mathbb{R}^{(p+1)}$ where p+1 means that there are p features from the dataset and the intercept. We define the matrix associated to the dataset and the model weights as follows

$$X^{T} = \begin{pmatrix} 1 & x_{1}^{(1)} & x_{2}^{(1)} & \dots & x_{p}^{(1)} \\ 1 & x_{1}^{(2)} & x_{2}^{(2)} & \dots & x_{p}^{(2)} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 1 & x_{1}^{(N)} & x_{2}^{(N)} & \dots & x_{p}^{(N)} \end{pmatrix} \in \mathbb{R}^{N \times (p+1)} \qquad x^{(i)} = \begin{pmatrix} 1 \\ x_{1}^{(i)} \\ x_{2}^{(i)} \\ \vdots \\ x_{p}^{(i)} \end{pmatrix} \quad w = \begin{pmatrix} b \\ w_{1} \\ w_{2} \\ \vdots \\ w_{p} \end{pmatrix}$$

the constant column is meant for the intercept, known as bias the b weight in vector w. The dataset matrix can be written as $X = (x^{(1)}, x^{(2)}, \dots, x^{(N)})$.

The objective function $f: \mathbb{R}^{(p+1)} \to \mathbb{R}$ is at least of class $f \in C^2(\mathbb{R}^{(p+1)})$, we can compute the first and second order derivatives

$$f(w) = \frac{1}{N} \sum_{i=1}^{N} \log(1 + \exp(-y^{(i)} w^{T} x^{(i)})) + \lambda \frac{1}{2} ||w||^{2}$$
(3a)

$$\nabla f(w) = \frac{1}{N} X r + \lambda w \tag{3b}$$

$$\nabla^2 f(w) = \frac{1}{N} X D X^T + \lambda I_{(p+1)}$$
(3c)

where $r \in \mathbb{R}^N$ is a vector of the same length as the total number of samples, whose elements are $r_i = -y^{(i)}\sigma(-y^{(i)}w^Tx^{(i)})$, note that $\sigma(z)$ is the sigmoid function; $D \in \mathbb{R}^{N \times N}$ is a diagonal matrix

whose elements are $d_{ii} = \sigma(y^{(i)}w^Tx^{(i)})\sigma(-y^{(i)}w^Tx^{(i)})$ which implies $d_{ii} \in (0,1)$; and $I_{(p+1)}$ is the identity matrix with size p+1. Dividing by N means dividing by the total number of samples involved.

The next proposition allows to solve the optimization problem.

Proposition 1. Problem (2) admits a unique optimal solution.

Proof. We need to prove the existence and the uniqueness of the global minimum.

(i) Existence of a optimal solution. The problem is quadratic and the objective function (3a) is coercive, that is $\forall \{w^k\}$ s.t. $\lim_{k\to\infty} ||w^k|| = \infty$ holds

$$\lim_{k \to \infty} f(w^k) \ge \lim_{k \to \infty} \lambda \frac{1}{2} \|w^k\|^2 = \infty \Rightarrow \lim_{k \to \infty} f(w^k) = \infty$$

hence by a corollary of the Weirstrass theorem, the problem admits global minimum in $\mathbb{R}^{(p+1)}$. (ii) Unicity of the optimal solution. We now prove that the hessian matrix (3c) is positive definite

$$w^T \nabla^2 f(w) w = w^T X D X^T w + \lambda w^T I w = \underbrace{y^T D y}_{>0} + \lambda \|w\|^2 \ge \lambda \|w\|^2 > 0 \quad \forall w$$

the 1/N is omitted. The hessian matrix positive definite implies that the objective function is *strictly convex* and that implies that the global minimum, if exists, is unique. Being in the convex case, the global minimum is a $w^* \in \mathbb{R}^{(p+1)}$ s.t. $\nabla f(w^*) = 0$ for first-order optimality conditions.

Remark 1. Since the log-loss is convex, the regularization term makes the objective function also *strongly convex*, this should speed up the optimization process.

2 Stochastic gradient descent variants

In this section we tackle the algorithmic part, specifically the SGD-type is the Mini-batch Gradient Descent where the mini-batch size M is greater than 1 and much less than the dataset size, i.e. $1 < |B| = M \ll N$, however, we will call it SGD anyway.

In order to use the algorithm, it is necessary to make further assumptions on the objective function and the gradients (like how far the mini-batch gradient are from the *true gradients*)

- the objective function from problem (2) is a convex loss function plus a quadratic regularization term, since f admits global minimum in $\mathbb{R}^{(p+1)}$ the function is bounded below by some value f^* ;
- for some constant G > 0 the magnitude of all gradients samples is bounded $\forall w \in \mathbb{R}^{(p+1)}$, by $\|\nabla f_i(w)\| \leq G$;
- other than twice continuously differentiable, we assume that f has Lipschitz-continuous gradients with constant L > 0, one can also say that f is L-smooth.

The algorithm is globally convergent, so the starting solution will be an arbitrary $w^0 \in \mathbb{R}^{(p+1)}$, that choice will is discussed in the experiments section.

Stopping criterion and failures

Regarding the implementation of the algorithm, it is essential to define a stopping criterion. Given a small $\varepsilon > 0$ the chosen criterion is

$$\|\nabla f(w^k)\| \le \varepsilon \tag{4}$$

note that the criterion uses the full gradient. In machine learning, one does not seek a very low precision. Other than the stopping criterion, one can add conditions of premature termination like

- exceeding a threshold for the epochs number k^* ;
- internal failures when computing w^{k+1} , for example exceeding q^* iterations during the line search (as you will se later, for the step-size α in the Armijo method as well as the momentum term β in the momentum correction).

Mini-batch gradient

Now we spend few words about the notation and the computation of the gradient evaluated on certain mini-batch. Being on *epoch* k with weights w^k , for every t iteration a (internal) model update has the following form

$$z^{t+1} = z^t + \alpha_t d_t, \qquad z^0 = w^k \tag{5}$$

until all the mini-batch are processed. The update uses information from the mini-batch B_t when calculating the direction d_t , the step-size α_t follows a certain rule that may use mini-batch information.*

The direction makes use of the gradient, so we want to compute the gradient w.r.t. z^t using information from the mini-batch B_t whose indices are randomly chosen from the full dataset $i_t \subset \{1, ..., N\}$ using a uniform distribution

$$\nabla f_{i_t}(z^t) = \frac{1}{M} \sum_{i \in B_t} \nabla \ell_i(z^t) + \lambda \nabla \Omega(z^t)$$

$$= \frac{1}{M} \underbrace{Xr}_{i \in B_t} + \lambda z^t$$
(6)

the expression is the same as the full gradient (3b) except that the dataset matrix contains just the mini-batch samples, so the r vector.

2.1 Basic stochastic gradient descent

The basic SGD version has the following iteration update rule

$$z^{t+1} = z^t - \alpha_t \nabla f_{i_t}(z^t) \tag{7}$$

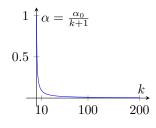
so the direction is defined as $d_t = -\nabla f_{i_t}(z^t)$ that is the negative gradient evaluated on the considered mini-batch, we know that on average is a descent direction so the objective function doesn't decrease necessarily at each step.

^{*} Iterations is defined as the total number of mini-batches extracted from the dataset, while one epoch is when the entire dataset is passed forward. The counter for the mini-batch currently processed is t while k is for the epoch.

Given an initial $\alpha_0 \in \mathbb{R}^+$, the first two basic versions are

- SGD-Fixed: step-size s.t. $\alpha_t = \alpha_0$;
- SGD-Decreasing: step-size s.t. $\alpha_t = \frac{\alpha_0}{k+1}$, see on the right.

The first choice sees the same step-size between epochs and so iterations. The second choice changes the step-size on every epoch, while being constant between iterations, that particular form ensures the convergence. For this two algorithms the momentum term in algorithm 5 is set to $\beta_0 = 0$.



2.1.1 Stochastic line search

Now we move to the approach by Vaswani, Mishkin, Laradji et al. [1]. For using the algorithm proposed by the paper, one more assumption is needed, that is, the model is able to *interpolate* the data, this property requires that the gradient evaluated on each sample converges to zero at the optimal solution

if
$$w^* \mid \nabla f(w^*) = 0 \Rightarrow \nabla f_i(w^*) = 0 \ \forall i = 1, \dots, N$$

The proposed approach applies the Armijo line search to the SGD algorithm at every iteration, specializing the sufficient reduction condition in the context of finite-sum problems. Hence the *Armijo condition* has the following form

$$f_{i_t}(z^t - \alpha_t \nabla f_{i_t}(z^t)) \le f_{i_t}(z^t) - \gamma \alpha_t \|\nabla f_{i_t}(z^t)\|^2$$
 (8)

the coefficient γ is the hyper-parameter controlling the aggressiveness of the condition, the paper suggests to set 1/2 as its maximum value.

As the standard Armijo method, the proposed line search uses a backtracking technique that iteratively decreases the initial step-size $\alpha_0 \in \mathbb{R}^+$ by a constant factor δ usually set to 1/2 until the condition is satisfied.

The authors also gave heuristics in order to avoid unnecessary function evaluations by restarting at each iteration the step-size to the previous one multiplied by the factor $a^{M/N}/\delta$, see algorithm 1 on page 8. Like the basic version, the momentum term is set to $\beta_0 = 0$.

2.2 Adding momentum term

The iteration performed over the mini-batches is still (5) what is the direction that is

$$d_t = -((1 - \beta_0)\nabla f_{i_t}(z^t) + \beta_0 d_{t-1})$$

in a finite-sum problem the momentum term lies in a specific range $\beta_0 \in (0,1)$ and is a constant value, the algorithm that uses this direction is the SGDM, the resulting iteration will be

$$z^{t+1} = z^t - \alpha_t ((1 - \beta) \nabla f_{i_t}(z^t) + \beta d_{t-1})$$
(9)

which is applied as the general update rule in algorithm 5, in this case the momentum term is set to a constant value $\beta = \beta_0$. To be clear we have the following cases

$$z^{t+1} = z^t - \alpha_t \left((1 - \beta_0) \nabla f_{i_t}(z^t) + \beta_0 d_{t-1} \right)$$

$$\beta_0 = 0$$

$$(7) \text{ SGD-}$$

$$\beta_0 \in (0, 1)$$

$$(9) \text{ SGDM-}$$

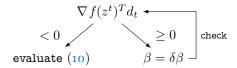
2.2.1 Stochastic line search

As the paper by Fan, Vaswani, Thrampoulidis et al. [2] says, when using the momentum term together with a line search, β_0 complicates the selection of a suitable step-size. The Armijo line search applied to the SGDM algorithm has the following condition

$$f_{i_t}(z^{t+1}) \le f_{i_t}(z^t) - \gamma \alpha_t \nabla f_{i_t}(z^t)^T ((1 - \beta_0) \nabla f_{i_t}(z^t) + \beta_0 d_{t-1})$$
(10)

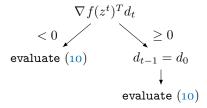
but the line search is not robust to the choice of the momentum term as the paper stated.

The problem is that $\nabla f_{i_t}(z^t)^T d_t < 0$ isn't always guaranteed, i.e. the direction is not descent, therefore the line search doesn't converge. Starting from an initial $\beta_0 \in (0,1)$, there are two situations that can be resolved as follows



in algorithmic terms, until the direction is descent, damp the momentum term by a factor δ , which is usually set to 0.5 like in the line search. When using this procedure, a descent direction d_t is guaranteed and it is possible to apply the algorithm 2, the procedure is called *momentum* correction, see algorithm 3 on the following page. The resulting algorithm is MSL-SGDM-C.

The momentum correction procedure can be expensive, so the paper suggests another approach called momentum restart. When the descent direction condition for d_t isn't satisfied, the procedure restarts that direction by setting $d_{t-1} = d_0$, the paper suggests $d_0 = 0$, in general



so if $d_0 = 0$ the direction will be $d_t = -(1 - \beta_0)\nabla f_{i_t}(z^t)$ that is a descent direction on the considered mini-batch, see algorithm 4 on the next page. The algorithm that uses this procedure is MSL-SGDM-R.

The authors suggest to set the momentum term to $\beta_0 = 0.9$.

D. Nardi

```
Algorithm 1: reset
                                                                         Algorithm 2: armijo-method
  Input: \alpha, \alpha_0, M, N, t, a \in \mathbb{R}^+,
                                                                           Data: \gamma \in (0,1), \delta \in (0,1), q^*
               \mathtt{opt} \in \{0,1,2\}
                                                                           Input: z^t, d_t, \alpha
_{1} if t=0 or opt =1 then
                                                                        \alpha \leftarrow \alpha/\delta;
_{2} | return \alpha_{0}
                                                                        q \leftarrow 0;
3 else if opt = 0 then
                                                                        3 repeat
                                                                                \alpha \leftarrow \delta \alpha;
\alpha \leftarrow \alpha
                                                                                z^{t+1} \leftarrow z^t + \alpha d_t;
_{5} else if opt = 2 then
  \alpha \leftarrow \alpha a^{M/N}
                                                                               q \leftarrow q + 1;
_7 end
                                                                        7 until
                                                                             f_{i_t}(z^{t+1}) \le f_{i_t}(z^t) + \gamma \alpha \nabla f_{i_t}(z^t)^T d_t or
   Output: \alpha
                                                                             q \geq q^*;
                                                                           Output: \alpha
```

```
Algorithm 3: momentum-correction
                                                                       Algorithm 4: momentum-restart
  Data: \delta \in (0,1), q^*
                                                                         Data: d_0
                                                                         Input: \beta_0, \nabla f_{i_t}(z^t), d_{t-1}
  Input: \beta_0, \nabla f_{i_t}(z^t), d_{t-1}
_{1} \beta \leftarrow \beta_{0};
                                                                      q \leftarrow 0;
q \leftarrow 0;
                                                                      d_t \leftarrow -((1-\beta_0)\nabla f_{i_t}(z^t) + \beta_0 d_{t-1});
3 repeat
                                                                        if not \nabla f_{i_t}(z^t)^T d_t < 0 then
      \beta \leftarrow \delta \beta;
                                                                              d_{t-1} \leftarrow d_0;
       d_t \leftarrow -((1-\beta)\nabla f_{i_t}(z^t) + \beta d_{t-1});
                                                                              compute d_t;
     q \leftarrow q + 1;
                                                                      6 end
7 until \nabla f_{i_t}(z^t)^T d_t < 0 or q \geq q^*;
                                                                         Output: d_t
   Output: d_t
```

Algorithm 5: SGD variants

```
Input: w^0 \in \mathbb{R}^{(p+1)}, M > 1, k^*, \varepsilon > 0, \alpha_0 \in \mathbb{R}^+, \beta_0 \in (0,1)
 1 k \leftarrow 0;
 while \|\nabla f(w^k)\| > \varepsilon and k < k^* do
          create mini-batches B_0, \ldots, B_{N/M-1};
          z^0 \leftarrow w^k:
          d_{-1} \leftarrow 0;
          \alpha_{-1} \leftarrow \begin{cases} \frac{\alpha_0}{k+1} & \text{if SGD-Decreasing} \\ \alpha_0 & \text{otherwise} \end{cases};
 6
          for t = 0 to N/M - 1 do
 7
                get indices i_t from B_t then get the samples;
 8
                \nabla f_{i_t}(z^t) \leftarrow \sum_{j \in B_t} \nabla f_j(z^t);
 9
                           \int -((1-\beta_0)\nabla f_{i_t}(z^t) + \beta_0 d_{t-1})
                                                                                                      if SGD-, SGDM
                          \left\{ \text{ momentum-correction}(\beta_0, \nabla f_{i_t}(z^t), d_{t-1}) \right\} \text{ if MSL-SGDM-C}; 
10
                          momentum-restart (\beta_0, \nabla f_{i_t}(z^t), d_{t-1})
                                                                                                     if MSL-SGDM-R
                if SGD-Armijo, MSL-SGDM-C/R then
11
                      \alpha \leftarrow \mathtt{reset}(\alpha_{t-1}, \alpha_0, M, N, t, a, \mathtt{opt});
                     \alpha_t \leftarrow \operatorname{armijo-method}(z^t, d_t, \alpha);
13
                else
14
                 \alpha_t \leftarrow \alpha_{-1};
15
                end
16
               z^{t+1} \leftarrow z^t + \alpha_t d_t;
17
18
          w^{k+1} \leftarrow z^{N/M}: k \leftarrow k+1:
19
20 end
     Output: w^*
```

3 Experiments and results discussion

To test the efficiency the algorithms, a benchmark of six datasets retrieved from LIBSVM is used, see table 1 on page 11 for details. Every dataset comes already pre-processed, with every sample scaled in range [-1,1] and the response variable in $\{-1,1\}$; many features are categorical with values 0,1,2..., this implies that the dataset matrix can be stored in sparse format, so the SciPy CSR matrix format was used.

Compared to the available benchmark dataset, those chosen are not that large, the choice is due to the hardware available (Intel® CoreTM i7, memory 16GB). As can be seen few dataset are unbalanced, this will affect the accuracy.

3.1 Solving the optimization problem

In order to solve the optimization problem, an initial guess w^0 for the model weights is given: we set a null *bias* and the other features are drawn from an uniform distribution in range [-0.5, 0.5].

Then a hyper-parameters tuning is performed. We set fixed values for the λ regularization coefficient from (2), the ε tolerance from the stopping criterion (4), the initial momentum term β_0 and the aggressiveness of the Armijo condition γ to a small value. Regarding failures for exceeding epochs and iterations, k^* and q^* for both Armijo method and momentum correction were set. Follow the values

$$\lambda = 0.5$$
 $\varepsilon = 10^{-3}$ $\beta_0 = 0.9$
 $\gamma = 10^{-3}$ $k^* = 600$ $q^* = 100$

Moving to the other hyper-parameters, a *grid search* is applied to find the best combination for each algorithm. The procedure confronts different combinations of the mini-batch size, the learning rate in the basic SGD version and the ones with line search, and in the latter are confronted also different values for the damping both in the Armijo method and momentum correction.

The mini-batch size grid depends on the dataset being considered. As said, the retrieved datasets vary in size, but the general rule is to stay around 100 iterations using values that are powers of 2, the grid is composed by at least two values. For the first five datasets, the size starts at 32. Follows the grids used

```
1, 0.5, 0.1, 0.01, 0.001 and 0.0005
     SGD-Fixed, SGDM
\alpha_0
     SGD-Decreasing
                                        1, 0.8, 0.5, 0.1, 0.05, 0.01 and 0.005
\alpha_0
\alpha_0
     SGD-Armijo, MSL-SGDM-C/R
                                       1, 0.5, 0.1, 0.05, 0.01 and 0.005
     SGD-, SGDM-
                                        depends on dataset
M
\delta_a
     SGD-Armijo, MSL-SGDM-C/R
                                       0.3, 0.5, 0.7 \text{ and } 0.9
    MSL-SGDM-C
                                        0.3, 0.5 and 0.7
```

where δ_a is the damping for the Armijo line search and δ_m for the momentum correction, the combinations for the MSL-SGDM-C are three times those of SGD-Armijo and MSL-SGDM-R.

The grid search chooses the best combination for a certain solver based on the greatest *test accuracy* and lowest *objective function* value. The results can be seen in tables from 2 on page 12 to 7 ordered descending by accuracy on the test dataset and ascending by objective function on reached solution. The other displayed values are the number of epochs and the run-time, the solution norm and the gradient norm.

For benchmarking purposes the optimization problem is solved also using the Full-batch Gradient Descent and three solvers from SciPy which are L-BFGS, Conjugate Gradient and Newton-CG

Regarding the grid search implementation, the Joblib module was used to parallelize on multiple cores the for loop needed to evaluate the algorithm with various combinations, see algorithm 6 on the following page.

First thing to say, the regularization term has a great influence on the final model, as you can see in the solution norm column the values are not that high, however lowering the λ coefficient would have led to a smaller solution norm, so the null model, possibly when the stochastic line search is used.

Speaking of the algorithms that use the Armijo line search, first of all one notices that their execution time is obviously longer than the others. More important, in each dataset you can see that the ε tolerance value is never reached though the limit of 600 epochs, especially the SGD-Armijo algorithm. However, in machine learning a very low tolerance is not necessarily required. Nonetheless the SLS solvers prove to achieve a higher accuracy.

The full-batch method for these dataset sizes is still a valid choice and could reach also a lower tolerance value, same for the SciPy solvers that reach the smallest gradient norm.

As expected the accuracy is greatly influenced by the class distribution of each datasets, for which different metrics that deal with unbalanced datasets could be used. Between all solvers the test score is very similar, as the $f(w^*)$ except in some cases for the SGD-Armijo solver that ended on a different solution with even higher accuracy.

3.2 Performance of the objective function

Now, as done by the authors of both articles, we want to show how the value of the objective function decreases with each epoch and also during time. For this purpose, another *grid search* is performed, again the fixed hyper-parameters are

$$\lambda = 0.5$$
 $\varepsilon = 10^{-3}$ $\beta_0 = 0.9$
 $\gamma = 10^{-3}$ $k^* = 200$ $q^* = 100$

What differs now is that the grid search finds the best hyper-parameters for each learning rate in 1, 0.1 and 0.01, hence the grids are

α_0	SGD-, SGDM-	1 or 0.1 or 0.01
M	SGD-, SGDM-	depends on dataset
δ_a	SGD-Armijo, MSL-SGDM-C/R	0.3, 0.5, 0.7 and 0.9
δ_m	MSL-SGDM-C	0.3, 0.5 and 0.7

where the mini-batch size M as in the previous grid search is composed by at least two values.

The algorithm finds the best values for each α_0 , then the objective function for every epoch and for the time took by each epoch is plotted in figures from 2 on page 14 to 4.

Although the first two datasets are unbalanced, the performance of the objective function tends to the minimum without significant fluctuations. In figure 4b the SGD-Armijo shows important oscillations, that may also be due to small dataset size.

In figure 3 we note that the SGD-Armijo reaches a low objective function value quickly but struggles to stay around and terminate, unlike the momentum versions that may take more epochs but oscillate less. Perhaps a different line search like a Wolfe-type may reduce this behaviour.

As pointed out by the authors, an important conclusion is that the methods with line search have a low sensitivity to the initial learning rate, thus having a similar performance. In contrast to the basic versions where significant differences in the trend can be seen, for example in figure 4a the SGD-Fixed and SGDM have very different behaviours for the each value in the α_0 grid, note that the momentum term tries to correct this abnormal behaviour.

Among the basic versions, the SGD-Decreasing seems to be the least affected by different initial learning rates. However if α_0 is too small, the algorithm takes smaller steps and requires more epochs to reach a solution.

Algorithm 6: Grid search for hyper-parameters tuning

Input: $\alpha_0, M, \delta_a, \delta_m$

- prepare the combinations using the given grids;
- $_{\mathtt{2}}$ for param in params \mathbf{do}
- 3 | model training using param;
- confront metrics with current best-model and update best-param if necessary;
- $_{5}$ end

Output: best-model, best-param

 Table 1: Benchmark datasets

Name	Train	Test	Features	Distribution
w1a	2477	47272	300	-1:0.97 1:0.03
w3a	4912	44837	300	-1:0.97 1:0.03
Phishing	8844	2211	68	-1:0.45 1:0.55
a2a	2265	30296	119	-1:0.75 1:0.25
Mushrooms	6499	1625	112	-1:0.48 1:0.52
German	800	200	24	-1:0.70 1:0.30

Table 2: w1a dataset

Solver	Epochs	Run-time	$ w^* $	$f(w^*)$	$\ \nabla f(w^*)\ $	Test score
SGD-Armijo	600	3.2954	0.455603	0.536138	3.63×10^{-1}	0.971 400
Newton-CG	6	NaN	0.667394	0.464614	4.60×10^{-5}	0.970236
CG	7	NaN	0.667395	0.464614	9.00×10^{-6}	0.970236
L-BFGS-B	7	NaN	0.667406	0.464614	2.30×10^{-5}	0.970236
BatchGD-Fixed	12	0.0060	0.667389	0.464614	5.64×10^{-4}	0.970236
SGD-Decreasing	27	0.0895	0.667400	0.464614	7.92×10^{-4}	0.970236
SGD-Fixed	27	0.1787	0.667311	0.464615	8.52×10^{-4}	0.970236
SGDM	386	2.9202	0.667383	0.464615	9.78×10^{-4}	0.970236
MSL-SGDM-R	600	3.1091	0.665659	0.464693	9.14×10^{-3}	0.970236
MSL- $SGDM$ - C	600	3.0976	0.665656	0.464693	9.15×10^{-3}	0.970236

Table 3: w3a dataset

Solver	Epochs	Run-time	$ w^* $	$f(w^*)$	$\ \nabla f(w^*)\ $	Test score
SGD-Armijo	600	8.0974	0.478986	0.500431	2.68×10^{-1}	0.971 006
Newton-CG	6	NaN	0.666640	0.462742	1.10×10^{-5}	0.970203
CG	7	NaN	0.666648	0.462742	2.20×10^{-5}	0.970203
L-BFGS-B	7	NaN	0.666658	0.462742	3.30×10^{-5}	0.970203
BatchGD-Fixed	12	0.0241	0.666635	0.462742	5.64×10^{-4}	0.970203
SGD-Decreasing	19	0.1657	0.666731	0.462743	8.76×10^{-4}	0.970203
SGD-Fixed	23	0.1549	0.666893	0.462743	9.49×10^{-4}	0.970203
SGDM	45	0.4911	0.666594	0.462743	8.95×10^{-4}	0.970203
MSL-SGDM-C	600	3.5907	0.667576	0.462787	6.92×10^{-3}	0.970203
MSL- $SGDM$ - R	600	3.5150	0.667576	0.462787	6.92×10^{-3}	0.970203

Table 4: Phishing dataset

Solver	Epochs	Run-time	$ w^* $	$f(w^*)$	$\ \nabla f(w^*)\ $	Test score
SGD-Armijo	600	7.2636	0.144155	0.687736	6.65×10^{-2}	0.865219
MSL-SGDM-C	600	9.4775	0.150617	0.686392	4.88×10^{-2}	0.709181
MSL-SGDM-R	600	17.7458	0.152190	0.685660	3.26×10^{-2}	0.568521
Newton-CG	5	NaN	0.164188	0.685065	0.00	0.567616
L-BFGS-B	5	NaN	0.164196	0.685065	8.00×10^{-6}	0.567616
CG	6	NaN	0.164214	0.685065	2.30×10^{-5}	0.567616
SGD-Decreasing	6	0.0407	0.164270	0.685065	5.08×10^{-4}	0.567616
SGDM	22	0.2663	0.163872	0.685065	5.75×10^{-4}	0.567616
BatchGD-Fixed	11	0.0448	0.164001	0.685065	5.34×10^{-4}	0.567616
SGD-Fixed	13	0.1701	0.163727	0.685065	9.27×10^{-4}	0.567616

Table 5: a2a dataset

Solver	Epochs	Run-time	$ w^* $	$f(w^*)$	$\ \nabla f(w^*)\ $	Test score
SGD-Armijo	600	3.9613	0.344862	0.603503	3.97×10^{-1}	0.825 191
BatchGD-Fixed	600	0.2068	0.355752	0.594416	3.64×10^{-1}	0.822386
MSL-SGDM-C	600	7.0367	0.406144	0.585491	2.90×10^{-1}	0.810569
SGD-Fixed	600	3.5597	0.425177	0.602741	3.56×10^{-1}	0.807136
MSL-SGDM-R	600	7.6314	0.413366	0.577575	2.28×10^{-1}	0.790236
SGDM	600	1.8107	0.438444	0.564030	2.63×10^{-3}	0.760298
Newton-CG	5	NaN	0.438972	0.564027	4.00×10^{-6}	0.760265
CG	12	NaN	0.438961	0.564027	1.50×10^{-5}	0.760265
L-BFGS-B	8	NaN	0.438969	0.564027	1.20×10^{-5}	0.760265
SGD-Decreasing	59	0.1687	0.438522	0.564028	7.26×10^{-4}	0.760265

Table 6: Mushrooms dataset

Solver	Epochs	Run-time	$ w^* $	$f(w^*)$	$\ \nabla f(w^*)\ $	Test score
SGD-Armijo	600	11.4039	0.644938	0.535765	2.34×10^{-1}	0.953 231
MSL-SGDM-C	600	6.7131	0.652698	0.529380	2.50×10^{-1}	0.942769
MSL-SGDM-R	600	6.4450	0.642551	0.527069	2.32×10^{-1}	0.940308
SGD-Fixed	600	3.3183	0.646431	0.525499	2.00×10^{-1}	0.926154
SGDM	600	3.6847	0.658660	0.557069	4.79×10^{-1}	0.924308
SGD-Decreasing	26	0.2326	0.635898	0.517727	7.79×10^{-4}	0.893538
Newton-CG	7	NaN	0.635933	0.517726	3.00×10^{-6}	0.892923
$\overline{\mathrm{CG}}$	11	NaN	0.635939	0.517726	2.40×10^{-5}	0.892923
L-BFGS-B	10	NaN	0.635930	0.517726	1.70×10^{-5}	0.892923
BatchGD-Fixed	26	0.0433	0.635906	0.517727	7.57×10^{-4}	0.892923

 $\textbf{Table 7:} \ \operatorname{German} \ \operatorname{dataset}$

Solver	Epochs	Run-time	$ w^* $	$f(w^*)$	$\ \nabla f(w^*)\ $	Test score
MSL-SGDM-C	600	4.9993	0.365021	0.629212	3.65×10^{-1}	0.755000
SGD-Armijo	600	5.0955	0.423356	0.632348	3.63×10^{-1}	0.750000
SGDM	600	2.3890	0.313661	0.616375	3.14×10^{-1}	0.745000
MSL-SGDM-R	600	5.6525	0.373225	0.599779	7.32×10^{-2}	0.730000
SGD-Decreasing	600	2.2918	0.348660	0.607993	1.14×10^{-1}	0.720000
Newton-CG	5	NaN	0.358504	0.597303	1.00×10^{-5}	0.710000
CG	12	NaN	0.358506	0.597303	4.00×10^{-6}	0.710000
L-BFGS-B	7	NaN	0.358506	0.597303	1.40×10^{-5}	0.710000
SGD-Fixed	58	0.2345	0.358265	0.597303	6.75×10^{-4}	0.710000
BatchGD-Fixed	20	0.0050	0.358324	0.597303	8.82×10^{-4}	0.710000

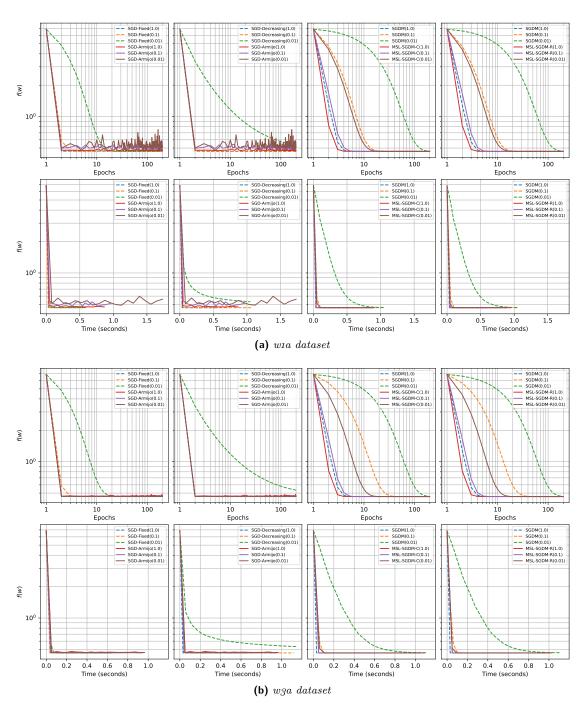


Figure 2: w1a and w3a datasets

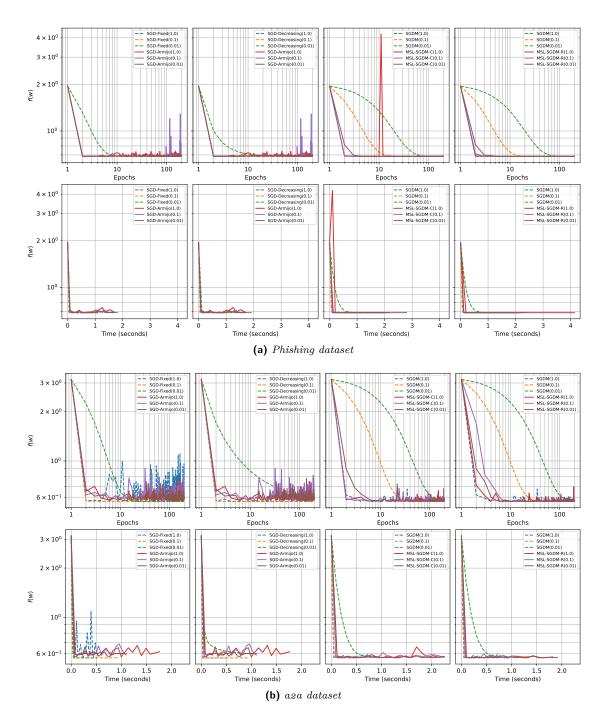


Figure 3: Phishing and a2a datasets

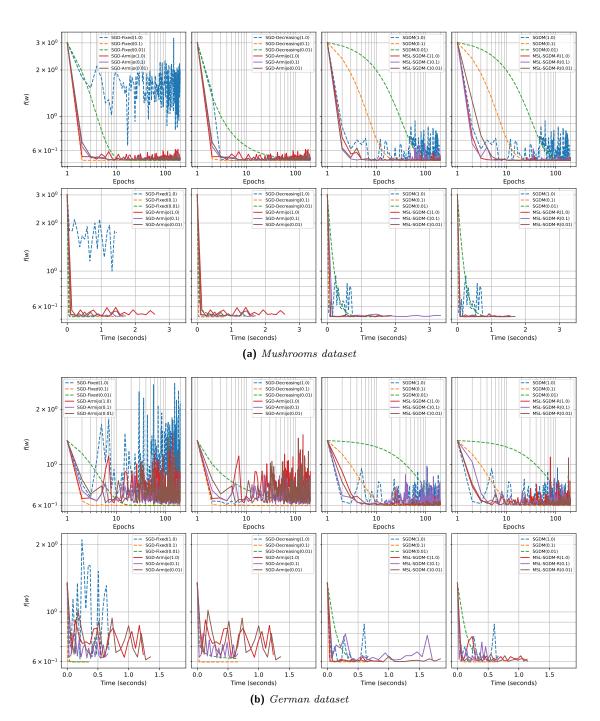


Figure 4: Mushrooms and German datasets

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