Speed Maintained SVRG

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Abstract—Stochastic gradient descent (SGD) is widely used for large-scale machine learning optimization, but has slow convergence rate due to the highly inherent variance. In recent years, the popular Stochastic Variance Reduced Gradient (SVRG) method mitigates this shortcoming, through computing the full-gradient of the entire dataset occasionally. However, conventional SVRG and its variants usually need a hyper-parameter to identify when to compute such the full gradient, which is essential to the convergence performance. Few previous studies discuss the method to identify such the hyper-parameter, which makes it hard to gain a good convergence performance in practical machine learning tasks. In our paper, we propose a new stochastic gradient descent with variance reduction technique named SMSVRG which computes the full gradient adaptively. Moreover, we propose an improved method denoted by SMSVRG+, which is comparable to and even better than SVRG with best-tuned epoch sizes for smooth and strongly convex functions.

I. INTRODUCTION

Many machine learning tasks such as logistic regression and linear regression can be formulated to be an optimization problem which is described as

$$\min F(\omega), \qquad F(\omega) = \frac{1}{n} \sum_{i=1}^{n} f_i(\omega) + R(\omega). \tag{1}$$

where n is the size of the training data. $F(\omega)$ means the loss function or training loss of a machine model, and ω is its parameter. $R(\omega)$ is the regularizer, which is widely used to avoid overfitting. It is noting that the total number of instances, i.e. n becomes very large with the proliferation of data.

Gradient Descent (GD) is a basic method to solve such the optimization problem. The gradient of $F(\omega)$ is obtained by passing over the entire training data, which is extremely time-consuming when the size of training data, i.e. n becomes large. Besides, GD is an iterative-convergent algorithm, that is, the parameter, i.e. ω , usually needs thousands of iterations to be converged. Since GD needs to compute the gradient of $F(\omega)$ every iteration, when the volume of data is large, the computation cost increases sharply and impairs the convergent performance significantly.

Stochastic Gradient Descent (SGD) mitigates this shortcoming by replacing the calculation of $\nabla F(\omega)$ with a stochastic gradient $\nabla f_i(\omega)$ with $i \in \{1,2,...,n\}$. In SGD, i is selected randomly from the entire training data. Thus, SGD outperforms GD on the time efficiency significantly. Take

the expectation of i, we obtain $\mathbb{E}[\nabla f_i(\omega)] = \nabla F(\omega)$. The difference between $\nabla f_i(\omega)$ and $\nabla F(\omega)$ represents variance which makes it difficult to achieve the optimum. In order to make the loss function, i.e. $F(\omega)$ converge, a decaying learning rate is usually used to reduce the variance. However, value of the learning rate is decayed to be very small after hundreds of iterations, which impedes the loss function to converge. In a nutshell, SGD with a decaying learning rate incurs a sub-linear convergence rate.

In recent years, variance reduced variants of SGD such as SVRG [3] is proposed to reduce the variance and gain the linear convergence performance with a constant learning rate. In SVRG, a full gradient is computed occasionally during the inexpensive SGD steps to reduce the variance, dividing the optimization procedure into many epochs. On the basis of SVRG, many variants have been proposed to improve its performance. SVRG-BB [9] uses the Barzilai and Borwein (BB) method proposed by Barzilai and Borwein in [2] to compute the step size before every epoch, which generally achieves the comparable convergence performance to SVRG with the best-tuned step size. CHEAPSVRG [8] and SAMPLEVR aim at reducing the expensive cost of full gradient computation through using a surrogate with a subset of the training dataset. mS2GD [5] uses mini-batch method to obtain a full gradient to reduce the variance, which shows a clear advantage for parallel computation. EMGD [11], SVR-GHT [7], Prox-SVRG [10] and SVRG with second-order information [4] modify the update rule of stochastic steps, and show advantages to SVRG in some cases. However, there are few studies discussing about how frequently should a full gradient be computed, i.e. how to set the epoch size m.

Most previous researches present that the epoch size, i.e. m should be constant [3, 9, 8] or increased monotonically [5], regardless of the learning rate. It is recommended that m=2n for convex problems and m=5n for non-convex problems in SVRG, without theoretical analysis and further experimental verification.

The epoch size, i.e. m has a great impact on the convergence performance of SVRG. More specifically, when m is too small, it wastes too much time to compute the full gradient frequently. When m is rather large, the variance between the stochastic gradient and the full gradient increases sharply, making the convergence of training loss extremely difficult. According to

the analysis of variance in [YaWei], both the epoch size, i.e. m and learning rate, i.e. η have a significant impact on the convergence performance. However, those previous studies do not provide a practical method to set the value of those hyperparameters. Extensive empirical studies illustrates that the choice of good value for those hyper-parameters costs much time in real machine learning tasks. In this paper, we develop a novel algorithm denoted by SMSVRG which can adjust the epoch size adaptively. SMSVRG applies a new stop condition regarding the change of parameters and checks the condition at the same interval, if the condition is satisfied, we stop the current epoch and step into the next one. Besides, we give guidance of how to set the parameters. Since SMSVRG may stop iterations earlier than expectation when n is quite small. we propose an improved algorithm denoted by SMSVRG+. In a nutshell, our contributions are highlighted as follows:

- SMSVRG, this novel algorithm can adjust the epoch size to a suitable value dynamically.
- SMSVRG+, it is an improvement of SMSVRG which is not sensitive to parameters and more practical than SMSVRG.
- Extensive empirical studies shows the effectiveness of our proposed algorithms which outperform their countparts on the convergence performance significantly.

This paper is organized as follows: Section II reviews the related work. Section III describes the SVRG algorithm. Section IV presents the new variant of SVRG, i.e. SMSVRG and its improved method, i.e. SMSVRG+. Section V demonstrates the numerical results of our algorithm. Section VI concludes this paper.

II. RELATED WORK

Several variants of SVRG discussing the strategy of adjusting epoch size has been proposed, including SVRG++ [1], S2GD [6], SVRG_Auto_Epoch [1] and so on.

SVRG++ adopts a simple strategy that epoch size m doubles between every consecutive two epochs. This method is absolutely heuristic and sometimes not justified. Our experiments show that when η is large or moderate, the exponential growth of m will incur great variance and impairs convergence.

S2GD designs a probability model of m and shows that a large epoch size is used with a high probability. However it needs to know the lower bound on the strong convexity constant of F, which is hard to estimate in practice. Meanwhile, the maximum of stochastic steps per epoch is also a sensitive parameter.

SVRG_Auto_Epoch is introduced as an additional improvement of SVRG++. It determines the termination of epoch through the quality of the snapshot full gradient. It records $diff_t = \|\nabla f_i(\omega_t^s) - \nabla f_i(\tilde{\omega}^{s-1})\|$ every iteration t and uses it as a tight upper bound on the variance of the gradient estimator. Although this method is reasonable, it has too much parameters to tune. Moreover, it takes much additional computation for every iterations, which impairs performances significantly.

Algorithm 1 SVRG

```
Require: learning rate \eta, epoch size m, initial point \tilde{\omega}
 1: for s=0,1,...,m do
2: \tilde{\mu}=\frac{1}{n}\sum_{i=1}^{n}\nabla f_{i}(\tilde{\omega}_{s})
            \omega_0 = \tilde{\omega}_s^i
 3:
 4:
            for t = 0, 1, 2, ... do
                   Randomly pick i_t \in \{1, 2, ..., n\}
  5:
                  \omega_t = \omega_{t-1} - \eta(\nabla f_{i_t}(\omega_{t-1}) - \nabla f_{i_t}(\tilde{\omega}_s) + \tilde{\mu})
  6:
  7:
            Option I: \tilde{\omega}_{s+1} = \omega_m
  8:
            Option II: \tilde{\omega}_{s+1} = \omega_t for randomly chosen t \in
       \{0,...,m-1\}
10: end for
```

Comparing with the above methods, smSVRG is apparently reasonable in intuition and does not need to tune extra parameters. Besides, it takes little additional computation cost and outperform the aforementioned three methods.

III. OVERVIEW

In this section we review the SVRG algorithm proposed by Johnson and Zhang[3]. As is shown in Algorithm 1, there are two loops in SVRG. Each outer loop is called an epoch while each inner loop is called an iteration. In each outer loop, a full gradient $\tilde{\mu}$ is computed at first, and its calculation requires to scan the entire dataset. In each inner loop, every iteration needs to pick $i_t \in \{1, 2, ..., n\}$ randomly. The update rule of the parameters is illustrated in Equation 2:

$$\omega_t = \omega_{t-1} - \eta(\nabla f_{i_t}(\omega_{t-1}) - \nabla f_{i_t}(\tilde{\omega}_s) + \tilde{\mu}). \tag{2}$$

Note that the expectation of $\nabla f_{i_t}(\tilde{\omega}_s)$ over i_t is $\nabla F(\omega_s)$, and the expectation of $\nabla f_{i_t}(\omega_{t-1})$ over i_t is $\nabla F(\omega_{t-1})$. We thus obtain

$$\mathbb{E}[\omega_t | \omega_{t-1}] = \omega_{t-1} - \eta \nabla F(\omega_{t-1}) \tag{3}$$

We can be seen that the variance of the update rule, i.e. Equation 2 is reduced. When both $\tilde{\omega}$ and ω_t converge to the optimum ω^* , then $\tilde{\mu} \to 0$ and $\nabla f_{i_t}(\omega_{t-1}) \to \nabla f_{i_t}(\tilde{\omega}_s)$, therefore

$$\nabla f_{i_t}(\omega_{t-1}) - \nabla f_{i_t}(\tilde{\omega}_s) + \tilde{\mu} \to 0$$

Hence, the learning rate for SVRG is allowed to be set as a relatively large constant against SGD, which results in a high convergence rate. At the end of each epoch, $\tilde{\omega}_{s+1}$ is updated by the output of inner loop. Note that there are two options for the update. Although only the convergence of SVRG with Option I is analyzed in [3], SVRG with Option II has been confirmed numerically to perform better. We adopt Option II in this paper. It is obvious that when m is too large, SGD with variance reducer will degenerate to the basic SGD, which results in huge variance. Thus our work focus on how to set an appropriate epoch size.

IV. SPEED MAINTAINED SVRG

In this section we describe two novel algorithms: SMSVRG and SMSVRG+, which can set the epoch size adaptively and have significant convergence performance superior to the previous studies. We assume the loss function F and the component functions f_i are convex and smooth throughout the paper.

A. smSVRG

1) Idea: It has been proved in [YaWei] that the optimal m is strongly related to η . Our experimental results also report that when η is large, the training loss begins to fluctuate after merely a small number of iterations. In the other hands, SVRG can endure far more than n iterations with a small η . In specific, if we stop the iterations in one epoch just before the training loss begins to fluctuate, the new algorithm will certainly be very efficient and outperform SVRG with constant epoch size. A direct approach is to evaluate the training loss occasionally. However, the training loss computation requires to pass over the entire training dataset, which is rather time consuming.

When we apply gradient descent to convex problem, as the ω_t gradually approaches to the optimal value ω^* , the gradient $\nabla F(\omega_t)$ thus keeps decreasing. We know that

$$\omega_t - \omega_{t-1} = -\eta \nabla F(\omega_{t-1})$$

Then we have $\|\omega_{t+1} - \omega_t\| < \|\omega_t - \omega_{t-1}\|$. Consider the update rule of SVRG, and take the expectation of i_t , we can obtain

$$\mathbb{E}[\omega_{t} - \omega_{t-1}] = \mathbb{E}[-\eta(\nabla f_{i_{t}}(\omega_{t-1}) - \nabla f_{i_{t}}(\tilde{\omega}_{s}) + \tilde{\mu})]$$

$$= -\eta \mathbb{E}[\nabla f_{i_{t}}(\omega_{t-1}) - \nabla f_{i_{t}}(\tilde{\omega}_{s}) + \tilde{\mu}]$$

$$= -\eta \nabla F(\omega_{t-1})$$
(4)

Intuitively, if ω_t keeps approaching to the optimal value ω^* , the $\nabla F(\omega_t)$ decays generally, with slight fluctuation caused by variance. Therefore, if we consider a window, we can believe that $\|\omega_{t+m_0} - \omega_t\| < \|\omega_t - \omega_{t-m_0}\|$ holds during the train of parameters. On the other hand, when ω_t oscillates near the optimal value due to the variance, $\|\omega_{t+m_0} - \omega_t\|$ will keep fluctuating.

Inspired by this observation, SMSVRG sets

$$\|\omega_{t+m_0} - \omega_t\| > \|\omega_t - \omega_{t-m_0}\| \tag{5}$$

as a stop condition in each epoch. When the Inequality 6 holds, we deem that the training loss fails to converge and begins to fluctuate, we have to suspend the iterations in such epoch and compute a full gradient in order to reduce the variance.

2) Details: As illustrated in Algorithm 2, smSVRG just requires two parameters: learning rate η and window size m_0 . It differs from SVRG only in the inner loop. At line 5, the first condition, i.e. $t\%m_0=0$ means that we check Inequality 6 every m_0 iterations. The second condition, i.e. $t>=2m_0$ is trivial as we need at least two windows to check the stop conditions. If the condition holds, we break the inner loop, and step out of the current epoch, and into the next epoch. Note that the epoch size is definitely a multiple of m_0 . Hence we

Algorithm 2 SMSVRG

```
Require: the learning rate \eta, the window size m_0, and an
      initial point \tilde{\omega}
  1: for s = 0, 1, ... do
           \tilde{\mu} = \frac{1}{n} \sum_{i=1}^{n} \nabla f_i(\tilde{\omega}_s)\omega_0 = \tilde{\omega}_s
  3:
 4:
            for t = 0, 1, 2, ... do
                  if t\%m_0 = 0 and t >= 2m_0 and \|\omega_t - \omega_{t-m_0}\| >
  5:
      \|\omega_{t-m_0}-\omega_{t-2m_0}\| then
                        break
  6:
  7:
                  Randomly pick i_t \in \{1, 2, ..., n\}
  8:
                  \omega_t = \omega_{t-1} - \eta(\nabla f_{i_t}(\omega_{t-1}) - \nabla f_{i_t}(\tilde{\omega}_s) + \tilde{\mu})
 9:
10:
            end for
            \tilde{\omega}_{s+1} = \omega_t
11:
12: end for
13: return \tilde{\omega}_{s+1}
```

should set m_0 to be smaller than 0.5n for flexibility. At the same time, m_0 cannot be too small because the stop condition may usually hold, which is caused by the random pick of the instances. We recommend to set m_0 between 0.1n and 0.2n according to extensive empirical sutdies.

3) Advantages:

- SMSVRG is superior to SVRG with prefixed epoch size as it can adapt the epoch size to an appropriate value dynamically. For a large η , SMSVRG will adjust the epoch size to be relatively small to constrain variance. On the other hand, when η is small, each epoch can thus contain more iterations. SMSVRG will make the epoch size larger than 2n to avoid to compute the time-consuming full gradient frequently.
- Comparing with the variants of SVRG such as SVRG_Auto_Epoch and S2GD, smSVRG requires far less additional computation cost. The reason is that it just needs to compute a simple inequality every m₀ iterations, which is very efficient.

B. Optimal Choice of the window size

In SMSVRG, we use $\|\omega_t - \omega_{t-m_0}\|$ to detect when the training loss begins to fluctuate and fails to converge. However, how to choose a suitable value of the window size i.e. m_0 becomes an important issue. In this section, we analyze the convergence performance by varying settings of m_0 , and provide guidance on setting an optimal value.

First, since the variance incurred by SGD iterations cannot be ignored, when we set m_0 to be small, the variance of $\|\omega_t - \omega_{t-m_0}\|$ is too large, and the probability of the stop condition of SMSVRG which is denoted by $C(m_0)$, is relatively high. As a result, the Inequality 6 may often holds due to variance, even though the loss function is still decreased during iterations. Therefore, the full gradient will be computed frequently, which thus wastes much time. It is obvious that $C(m_0)$ is a monotonically increasing function of m_0 .

If we set m_0 to be relatively large, it will be too late to detect the oscillation of training loss. Therefore, it wastes much time, and makes no process to optimize the objective function. We use $D_s(m_0)$ to denote the <u>Absolute Delay</u> of detecting fluctuation. Furthermore, the <u>Absolute Delay</u> makes different effects on the convergence performance, which depends on the epoch size. Thus, it is better to consider the Relative Delay:

$$D_r(m_0) = \frac{D_s(m_0)}{es+n}. (6)$$

Note that es denote the number of iterations in a specific epoch and n denotes the size of the training dataset. Function $D_r(m_0)$ is also monotonically increasing with respect to m_0 .

It is necessary to choice a suitable m_0 which should ensure that $C(m_0)$ is large and $D_r(m_0)$ is small. In order to obtain a trade-off between $C(m_0)$ and $D_r(m_0)$, we convert the parameter-choosing problem to the following maximization problem:

$$m_0 = \max_{m_0} \{ C(m_0) - D_r(m_0) \}$$

$$= \max_{m_0} \{ C(m_0) - \frac{D_s(m_0)}{es + n} \}$$
(7)

s.t.

$$0 < m_0 < n \tag{8}$$

From equation (7) we can obtain the following conclusions:

- 1) When epoch size es is small, the D_s tends to be more important than C. Hence we should set m_0 to be relatively small to maximize the objective function. On the contrary, it is recommended to set m_0 to be large. In spirit of this, we can set m_0 to be proportional to the iteration number of previous epoch.
- 2) According to our experiment on SVRG, when the learning rate η is large, the loss function begin to fluctuate after merely n/10 iterations, so we should set the initial m_0 to the same order of magnitude as $10^{-1} \times n$

C. smSVRG+

- 1) Idea: Our experiments on SMSVRG show that it performs well for large and moderate η . However, when η is rather small, the best epoch size for SVRG will be larger than 10n. And SMSVRG may finish the epoch prematurely while the training loss keeps declining. It is because the inequality (6) may holds due to variance, rather than the fluctuation of training loss. According to the analysis in IV-B, variance of $\|\omega_{t+m_0} \omega_t\|$ has more impact on the performance of our algorithm when the best epoch size is large. Intuitively, we should adjust the window size m_0 dynamically according to current epoch size, instead of a constant value.
- 2) Details: As illustrated in Algorithm 3, we initialize the m_0 to be 0.1n and recompute m_0 after the inner loop in each epoch. m_0 is set to be $(\lfloor es/n \rfloor + 1) \times (0.1n)$, where es denotes the iteration number of current epoch size. It means that m_0 increases when es exceeds estimates n, estimates 2n, estimates 3n and so on.

Algorithm 3 SMSVRG+

```
Require: learning rate \eta, window size m_0, initial point \tilde{\omega}
 1: for s=0,1,... do
2: \tilde{\mu}=\frac{1}{n}\sum_{i=1}^{n}\nabla f_{i}(\tilde{\omega}_{s})
 3:
            for t = 0, 1, 2, ... do
 4:
  5:
                  if t > m_0 and t\%m_0 = 0 and \|\omega_t - \omega_{t-m_0}\| >
      \|\omega_{t-m_0}-\omega_{t-2m_0}\| then
                        break
  6:
                  end if
 7:
                  Randomly pick i_t \in \{1, 2, ..., n\}
 8:
                 \omega_t = \omega_{t-1} - \eta(\nabla f_{i_t}(\omega_{t-1}) - \nabla f_{i_t}(\tilde{\omega}_s) + \tilde{\mu})
  9:
10:
11:
            es = t
12:
            m_0 = (int(es/n) + 1) * (0.1n)
13:
            \tilde{\omega}_{s+1} = \omega_t
14: end for
15: return \tilde{\omega}_{s+1}
```

TABLE I
DETAIL INFORMATION OF DATASETS AND MODELS

Dataset	size	dimension	model	λ
ijenn1	49990	22	logistic	10^{-4}
a9a	32561	123	logistic	10^{-4}
YearPredictionMSD	463715	90	linear	10^{-4}
cadata	20640	8	linear	10^{-4}

3) Advantages:

- SMSVRG+ outperforms SMSVRG as it will enlarge the window size to reduce variance when necessary, avoiding to stop an epoch early. It shows good performance regardless the learning rate and datasets in our experiments.
- SMSVRG+ is not sensitive to the initialized m_0 , as it tunes m_0 to an appropriate size adaptively. By contrast, the performance of SMSVRG fairly depends on the choice of m_0 .

V. NUMERICAL EXPERIMENTS

A. Experimental settings

In this section, we conduct extensive experiments to demonstrate the advantages of our proposed algorithms. We evaluate our algorithms on eight training datasets, which are public on the LIBSVM website¹. More specifically, *l*2-regularized logistic regression on datasets: ijcnn1, a9a, w8a and mushrooms, and *l*2-regularized ridge regression on datasets: earPredictMSD, cadata, mg, and abalone are conducted to evaluate our proposed algorithms and the previous work. The details of those datasets are illustrated in Table I.

The l2-regularized logistic regression task is conducted on datasets: ijcnn1, a9a, w8a and mushrooms where the label of

http://www.csie.ntu.edu.tw/~cjlin/libsvmtools/datasets/

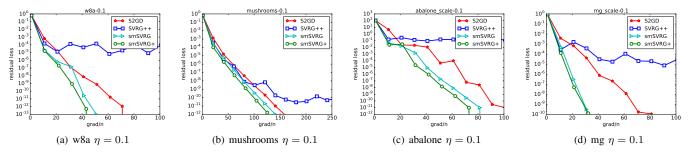


Fig. 1. Comparison of SMSVRG, SMSVRG+, SVRG++, S2GD on four datasets.

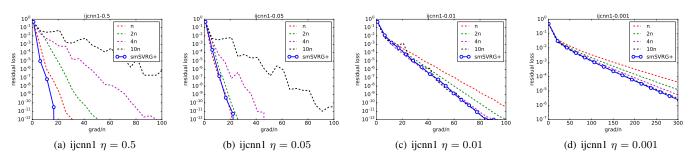


Fig. 2. Generally, SMSVRG can automatically set an appropriate m with different learning rates for the l2-regularized logistic regression

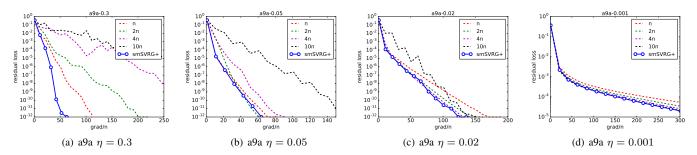


Fig. 3. Generally, SMSVRG can automatically set an appropriate m with different learning rates for the l2-regularized logistic regression

instances is set to be 1 or -1. Thus, the loss function of l2-regularized logistic regression task is formulated:

$$\min_{\omega} \frac{1}{n} \sum_{i=1}^{n} \log(1 + e^{-y_i \omega^{\mathsf{T}} x_i}) + \lambda \parallel \omega \parallel^2. \tag{9}$$

Here, x_i is the instances in the training dataset, and y_i is the label of x_i . λ is the coefficient of regularizer. Additionally, the l2-regularized ridge regression task is conducted on datasets: YearPredictMSD, cadata, mg and abalone. The loss function of l2-regularized ridge regression task is formulated:

$$\min_{\omega} \frac{1}{n} \sum_{i=1}^{n} \left(\omega^{\mathrm{T}} x_i - y_i \right)^2 + \lambda \parallel \omega \parallel^2.$$
 (10)

We scale the value of all features to [-1,1] and set λ to be 10^{-4} for all evaluations. In all figures, the x-axis denotes the computational cost, which is measured by the number of gradient computation divided by the size of training data, i.e. n. The y-axis denotes training loss residual, i.e. $F(\tilde{\omega}_s) - F(\omega^*)$. Note that the optimum $F(\omega^*)$ is estimated

by running the gradient descent for a long time. The x-axis in all figures means that the number of gradient calculations divided by the size of training data, i.e. $\operatorname{grad/n}$. This is a metric for measuring gradient complexity in previous work, which is a similar to measure the time cost during train of parameters [???]. Our numerical experiments include three parts: comparison of convergence performance with previous methods, comparison of convergence performance with SVRG by varying learning rate and sensitivity test varying m_0 . The experimental results report the superior performance of our methods.

B. Comparison of convergence performance with previous methods

In this section, we compare our SMSVRG and SMSVRG+ with two aforementioned existing methods: SVRG++ and S2GD. We do not compare with SVRG_Auto_Epoch in that we find that its termination condition of epoch is never satisfied and thus SVRG_Auto_Epoch keeps doing SGD iteration, resulting in non-convegent. For SVRG++, we initialize

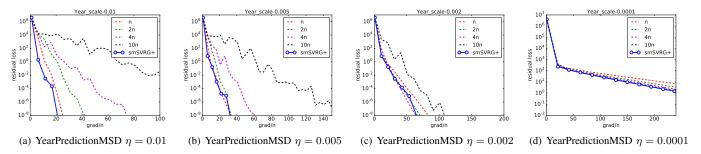


Fig. 4. Generally, SMSVRG can automatically set an appropriate m with different learning rates for the l2-regularized linear regression

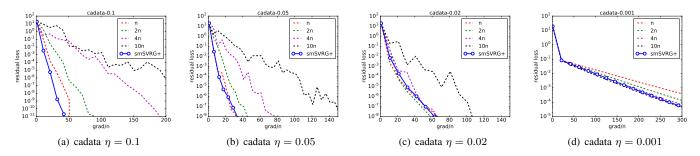


Fig. 5. Generally, SMSVRG can automatically set an appropriate m with different learning rates for the l2-regularized linear regression

m=n. For S2GD, we set the maximum of m to be 4n. For both sMSVRG and sMSVRG+, we set the window size, i.e. m_0 to be 0.1n. We evaluate these methods by running logistic regression on the dataset ijcnn1. As illustrated in Figure 1, we can see that SVRG++ always fluctuates violently and fails to converge due the large variance caused by too large epoch size. It is shown that sMSVRG and sMSVRG+ always outperform SVRG++ and S2GD and converge rapidly. Besides, the performance of sMSVRG+ is superior to that of sMSVRG. The main reason is that sMSVRG+ can adjust the windows size to a suitable value adaptively.

C. Comparison of convergence performance with SVRG by varying learning rate

Since SMSVRG+ performs better than SMSVRG, SMSVRG+ is used to conduct the comparison with SVRG. For SVRG, we increase the epoch size, i.e. m as four different values: n, 2n, 4n, 10n. Those values are used to conduct performance evaluation in SVRG [3]. Besides, the performance with the epoch size larger than 10n is similar to the performance with an extremely large epoch size. It is because that the full gradient is rare to be computed due to such extremely epoch size. The dashed lines represents SVRG with a fixed epoch size; while the green solid lines stands for SMSVRG+.

As illustrated in Figures 2, 3, 4, 5, SMSVRG+ can always have the similar performance as SVRG with most best-tuned epoch size. We observe that when η is large, and m is set to be a small value, e.g. n, can achieve best performance. The main reason is that when η is large, the variance becomes significant simultaneously, so m must be set to be small in order to bound the variance. As η decays, the optimal value of m increases,

which means that the algorithm can tolerate more variance induced by extra iterations. As illustrated in Figures 2(a), 3(a), 4(a), 5(a), our method is significantly better than SVRG with best-tuned epoch sizes when learning rate is large or medium. However, As illustrated in Figures 2(d), 3(d), 4(d), 5(d), if η is set to be too small, SMSVRG+ performs slightly superior to SVRG with large epoch sizes, but outperforms SVRG with recommended epoch sizes, i.e. n and 2n. It is noting that setting η to be too small is not a practical approach when using SVRG or its variants, because the convergence rate will be extremely low. Therefore, the sub-optimal performance of SMSVRG+ with very small η is acceptable.

D. Sensitivity test varying m_0

In this section, we conduct the sensitivity test on both SMSVRG and SMSVRG+ regarding window size m_0 . As analyzed in IV-B, we vary the initial m_0 ranges from 0.1 to 0.25 in order to test the sensitivity of SMSVRG and SMSVRG+. We conduct the experiments by using logistic regression and linear regression on two datasets: a9a and abalone, respectively. As illustrated in Figures 6(a) and 6(c), the performance of SMSVRG varies with the m_0 significantly. And we can see in Figures 6(b) and 6(d) that the performance of SMSVRG+ is not sensitive to the choice of m_0 . The main reason is that SMSVRG+ will adjust the m_0 adaptively regardless of the initialization. Hence SMSVRG+ is more practical than SMSVRG in reality.

VI. CONCLUSION

In this paper we propose a novel stop condition for each epoch in SVRG, which leads to a new variant of SVRG: SMSVRG. It can adjust the epoch size adaptively. We analyze

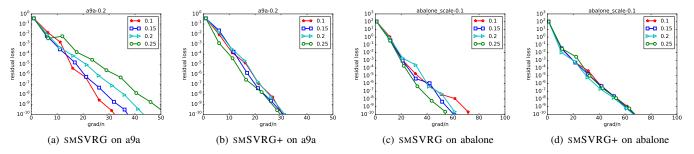


Fig. 6. Sensitivity test on m_0 for SMSVRG and SMSVRG+. The four numbers on the legends means four different choices of m_0 .

how to choose the optimal value of parameters, and thus develops an improved method called SMSVRG+. We conduct numerical experiments on real datasets to demonstrate the performance of SMSVRG and SMSVRG+. The experiments show that both SMSVRG and SMSVRG+ are superior to existing methods. Moreover, the SMSVRG+ is comparable to and sometimes even better than the original SVRG with best-tuned epoch sizes.

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