

On Distributed Adaptive Optimization with Gradient Compression

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Abstract

We study COMP-AMS, a distributed optimization framework based on gradient averaging and adaptive AMSGrad algorithm. Gradient compression is applied to reduce the communication in the gradient transmission process, whose bias is corrected by the tool of error feedback. Our convergence analysis of COMP-AMS shows two novel results in adaptive optimization: (i) like SGD, error feedback can also fix the convergence issue of compressed gradients for adaptive method, achieving same convergence rate as the full-gradient counterpart; (ii) distributed COMP-AMS has a linear speedup w.r.t. the number of local workers. Numerical experiments are conducted to justify the theoretical findings, and demonstrate significant communication reduction of the proposed method. Since COMP-AMS is simple and efficient in both communication and memory, it can serve as a convenient distributed training protocol for adaptive methods.

1 Introduction

Deep neural network has achieved the state-of-the-art learning performance on numerous AI applications, e.g., computer vision [Goodfellow et al. \(2014\)](#); [He et al. \(2016\)](#); [Voulodimos et al. \(2018\)](#), Natural Language Processing [Graves et al. \(2013\)](#); [Young et al. \(2018\)](#); [Zhang et al. \(2018\)](#), Reinforcement Learning [Mnih et al. \(2013\)](#); [Silver et al. \(2017\)](#) and recommendation systems [Covington et al. \(2016\)](#); [Wei et al. \(2017\)](#). With the sheer size of data observations and the increasing complexity of deep neural networks, standard single-machine training procedures encounter at least two major challenges:

- Due to the limited computing power of a single-machine, processing the massive number of data samples takes a long time — training is too slow. Many real-world applications can not even afford spending that much time on training.
- In many practical scenarios, data are typically stored in multiple servers, possibly at different locations, due to the storage constraints (massive user behavior data, Internet images, etc.) or privacy reasons [Chang et al. \(2018\)](#). Hence, transmitting data among servers might be costly.

Distributed learning framework [Dean et al. \(2012\)](#) has been a common training strategy to tackle the above two issues. For example, in centralized distributed stochastic gradient descent (SGD) protocol, data are located at n local nodes, at which the gradients of the model are computed in parallel. In each iteration, a central server aggregates the local gradients, updates the global model, and transmits back the updated model to the local nodes for subsequent gradient computation. As we can see, this setting naturally solves aforementioned issues: 1) We use n computing nodes to train the model, so the time per training epoch can be largely reduced; 2) There is no need to transmit the local data to central server. Besides, distributed training also provides stronger error tolerance since the training process could continue even one local machine breaks down. As a result of these advantages, there has been a surge of study and applications on distributed systems [Boyd et al. \(2011\)](#); [Nedic and Ozdaglar \(2009\)](#); [Duchi et al. \(2011\)](#); [Goyal et al. \(2017\)](#); [Hong et al. \(2017\)](#); [Lu et al. \(2019\)](#); [Koloskova et al. \(2019\)](#).

Among many optimization strategies, SGD is still the most popular prototype in distributed training for its simplicity and effectiveness [Chilimbi et al. \(2014\)](#); [Agarwal et al. \(2018\)](#); [Mikami et al. \(2018\)](#). Yet, when the deep learning model is very large, the communication between local nodes and central server could be expensive. Burdensome gradient transmission would slow down the whole training system, or even be impossible because of the limited bandwidth in some applications. Thus, reducing the communication cost in distributed SGD has become an active topic, and an important ingredient of large-scale distributed systems (e.g. [Seide et al. \(2014\)](#)). Solutions based on quantization, sparsification and other compression techniques of the local gradients are proposed, e.g., [Alistarh et al. \(2017\)](#); [Wen et al. \(2017\)](#); [Wangni et al. \(2018\)](#); [Stich et al. \(2018\)](#); [Aji and Heafield \(2017\)](#); [Bernstein et al. \(2018\)](#); [De Sa et al. \(2017\)](#); [Yang et al. \(2019\)](#); [Ivkin et al. \(2019\)](#). As one would expect, in most approaches, there exists a trade-off between compression and learning performance. In general, larger bias and variance of the compressed gradients usually bring more significant performance downgrade in terms of convergence [Stich et al. \(2018\)](#); [Ajallooeian and Stich \(2020\)](#). Interestingly, studies (e.g., [Karimireddy et al. \(2019\)](#)) show that the technique of *error feedback* can to a large extent remedy the issue of such biased compressors, achieving same convergence rate as full-gradient SGD.

On the other hand, in recent years, adaptive optimization algorithms (e.g. AdaGrad [Duchi et al. \(2010\)](#), Adam [Kingma and Ba \(2014\)](#) and AMSGrad [Reddi et al. \(2018\)](#)) have become popular because of their superior empirical performance. These methods use different implicit learning rates

for different coordinates that keep changing adaptively throughout the training process, based on the learning trajectory. In many learning problems, adaptive methods have been shown to converge faster than SGD, sometimes with better generalization as well. Despite of the great popularity of adaptive methods, the body of literature that extends them to distributed training is still very limited. In particular, even the simple gradient averaging approach, though appearing standard, has not been considered for adaptive optimization algorithms. Furthermore, adopting gradient compression in adaptive methods has also been rarely studied in literature. We try to fill this gap in this paper, by studying COMP-AMS, a distributed adaptive optimization framework using the gradient averaging protocol. Gradient compression is incorporated to reduce the communication cost. Theoretical and empirical results are presented to demonstrate the effectiveness of the proposed method.

1.1 Our Contributions

Specifically, in this paper, we focus on a *simple optimization framework* leveraging the *adaptivity* of AMSGrad and the computational virtue of *local gradient compression*. Our contributions summarize as follows:

- We consider COMP-AMS, a synchronous distributed adaptive optimization framework based on global averaging with gradient compression, which is efficient in both communication and memory as no local moment estimation is needed. Our scheme is coupled with an error-feedback technique to reduce the bias implied by the compression step.
- We provide the general convergence rate of distributed COMP-AMS (with n workers) in smooth non-convex optimization, where data heterogeneity is allowed. In the special case of $n = 1$, we derive the first result in literature that similar to SGD, gradient compression with error feedback in adaptive method achieves same convergence rate $\mathcal{O}(\frac{1}{\sqrt{T}} + \frac{1}{T})$ (omitting all constants) as the standard full-gradient counterpart. Moreover, we show that with a properly chosen learning rate, COMP-AMS achieves $\mathcal{O}(\frac{1}{\sqrt{nT}})$ convergence, implying a linear speedup in terms of the number of local workers to attain $\mathcal{O}(\delta)$ -stationary point. This is also the first result in literature regarding the linear speedup property of distributed adaptive learning under gradient compression.
- We present numerical experiments on three tasks that validate our theoretical findings on the merit of error feedback and the linear speedup effect, and show that COMP-AMS with **Top- k** or **Block-Sign** compressors can reduce ~ 30 -100x communication overhead, without losing much accuracy compared with using full gradient. Finally, we provide more discussions on the comparison to other related methods and potential future directions.

In Section 2 we provide a comprehensive summary of the contributions to date, related to gradient compression, adaptive methods and the error feedback technique. We develop in Section 3 our communication-efficient COMP-AMS method, using AMSGrad as a prototype optimization algorithm. In Section 4, we establish the convergence analysis and demonstrate several implications of our result. Numerical results are illustrated in Section 5 to justify the theory and show the effectiveness of the proposed approach.

2 Related Work

2.1 Distributed SGD with Compressed Gradients

Quantization. As we mentioned before, SGD is the most commonly adopted optimization method in distributed training of deep neural nets. To reduce the expensive communication in large-scale distributed systems, extensive works have considered various compression techniques applied to the gradient transaction procedure. The first strategy is quantization. Dettmers (2016) condenses 32-bit floating numbers into 8-bits when representing the gradients. Seide et al. (2014); Bernstein et al. (2018); Karimireddy et al. (2019); Bernstein et al. (2019) use the extreme 1-bit information (sign) of the gradients, combined with tricks like momentum, majority vote and memory. Other quantization-based methods include QSGD Alistarh et al. (2017); Wu et al. (2018); Zhang et al. (2017) and LPC-SVRG Yu et al. (2019b), leveraging unbiased stochastic quantization. The saving in communication of quantization methods is moderate: for example, 8-bit quantization reduces the cost to 25% (compared with 32-bit full-precision). Even in the extreme 1-bit case, the largest compression ratio is around $1/32 \approx 3.1\%$.

Sparsification. Gradient sparsification is another popular solution which may provide higher compression rate. Instead of commuting the full gradient, each local worker only passes a few coordinates to the central server and zeros out the others. Thus, we can more freely choose higher compression ratio (e.g., 1%, 0.1%), still achieving impressive performance in many applications Lin et al. (2018). Stochastic sparsification methods, including uniform and magnitude based sampling Wangni et al. (2018), select coordinates based on some sampling probability, yielding unbiased gradient compressors with proper scaling. Deterministic methods are simpler, e.g., Random- k , Top- k Stich et al. (2018); Shi et al. (2019) (selecting k elements with largest magnitude), Deep Gradient Compression Lin et al. (2018), but usually lead to biased gradient estimation. In Ivkin et al. (2019), the central server identifies heavy-hitters from the count-sketch Charikar et al. (2002) of the local gradients, which can be regarded as a noisy variant of Top- k strategy. More applications and analysis of compressed distributed SGD can be found in Jiang and Agrawal (2018); Shen et al. (2018); Alistarh et al. (2018); Basu et al. (2019); Jiang et al. (2018), among others.

Error Feedback (EF). Biased gradient estimation, which is a consequence of many aforementioned methods (e.g., signSGD, Top- k), undermines the model training, both theoretically and empirically, with slower convergence and worse generalization Ajalloeian and Stich (2020); Beznosikov et al. (2020). The technique of *error feedback* is able to “correct for the bias” and fix the convergence issues. In this procedure, the difference between the true stochastic gradient and the compressed one is accumulated locally, which is then added back to the local gradients in later iterations. Stich et al. (2018); Karimireddy et al. (2019) prove the $\mathcal{O}(\frac{1}{T})$ and $\mathcal{O}(\frac{1}{\sqrt{T}})$ convergence rate of EF-SGD in strongly convex and non-convex setting respectively, matching the rates of vanilla SGD Nemirovski et al. (2009); Ghadimi and Lan (2013). More works on the convergence rate of SGD with error feedback include Zheng et al. (2019); Stich and Karimireddy (2019), among other related papers.

2.2 Adaptive Optimization

Algorithm 1 AMSGRAD optimization method

- 1: **Input:** parameters β_1 , β_2 , and η_t .
 - 2: Initialize: $\theta_1 \in \Theta$ and $v_0 = \epsilon \mathbf{1} \in \mathbb{R}^d$.
 - 3: **for** $t = 1, \dots, T$ **do**
 - 4: Compute stochastic gradient g_t at θ_t .
 - 5: $m_t = \beta_1 m_{t-1} + (1 - \beta_1) g_t$.
 - 6: $v_t = \beta_2 v_{t-1} + (1 - \beta_2) g_t^2$.
 - 7: $\hat{v}_t = \max(\hat{v}_{t-1}, v_t)$.
 - 8: $\theta_{t+1} = \theta_t - \eta_t \frac{\theta_t}{\sqrt{\hat{v}_t}}$.
 - 9: **end for**
-

In each SGD update, all the gradient coordinates share the same learning rate. This latter is either constant or decreasing through the iterations. Adaptive optimization methods cast different learning rate on each dimension. For instance, AdaGrad, developed in [Duchi et al. \(2010\)](#), divides the gradient element-wisely by $\sqrt{\sum_{t=1}^T g_t^2} \in \mathbb{R}^d$, where $g_t \in \mathbb{R}^d$ is the gradient vector at time t and d is the model dimensionality. Thus, it intrinsically assigns different learning rates to different coordinates throughout the training – elements with smaller previous gradient magnitude tend to move a larger step via larger learning rate. AdaGrad has been shown to perform well especially under some sparsity structure in the model and data. Other adaptive methods include AdaDelta [Zeiler \(2012\)](#) and Adam [Kingma and Ba \(2014\)](#), which introduce momentum and moving average of second moment estimation into AdaGrad hence leading to better performances. AMSGrad [Reddi et al. \(2018\)](#) fixes the potential convergence issue of Adam, which will serve as the prototype in this paper. We present the pseudocode in Algorithm 1.

In general, adaptive optimization methods in many cases exhibit faster convergence than SGD. Thus, they have been widely used in training deep learning models in language and computer vision applications, e.g., [Choi et al. \(2019\)](#); [You et al. \(2020\)](#); [Zhang et al. \(2020\)](#). In distributed setting, the work [Nazari et al. \(2019\)](#) proposes a decentralized system in online optimization. However, communication efficiency is not considered. The recent work [Chen et al. \(2020\)](#) is the most relevant to our paper. Yet, their method is based on Adam, and requires every local node to store a local estimation of the moments of the gradient. Thus, one has to keep extra two more tensors of the model size on each local worker, which may be less feasible in terms of memory particularly with large models. We will present more detailed comparison in Section 3.

3 Communication-Efficient Adaptive Optimization

Consider the distributed optimization task where n workers jointly solve a large finite-sum optimization problem written as:

$$\min_{\theta \in \Theta} \frac{1}{n} \sum_{i=1}^n f_i(\theta) := \frac{1}{n} \sum_{i=1}^n \mathbb{E}_{x \sim \mathcal{X}_i} [F_i(\theta; x)], \quad (1)$$

where the non-convex function f_i represents the average loss (over the local data samples) for worker $i \in [n]$ and θ the global model parameter taking value in Θ , a subset of \mathbb{R}^d . \mathcal{X}_i is the data distribution on each local node.

3.1 Gradient Compressors

In this paper, we mainly consider deterministic q -deviate compressors defined as below.

Assumption 1. *The gradient compressor $\mathcal{C} : \mathbb{R}^d \mapsto \mathbb{R}^d$ is q -deviate: for $\forall x \in \mathbb{R}^d$, $\exists 0 \leq q < 1$ such that $\|\mathcal{C}(x) - x\| \leq q \|x\|$.*

Note that, larger q indicates an important compression while smaller q implies better approximation of the true gradient. Hence, $q = 0$ implies no compression, i.e., $\mathcal{C}(x) = x$. We give two popular and highly efficient q -deviate compressors that will be adopted in this paper.

Definition 1 (Top- k). *For $x \in \mathbb{R}^d$, denote \mathcal{S} as the size- k set of $i \in [d]$ with largest k magnitude $|x_i|$. The **Top- k** compressor is defined as $\mathcal{C}(x)_i = x_i$, if $i \in \mathcal{S}$; $\mathcal{C}(x)_i = 0$ otherwise.*

Definition 2 (Block-Sign). *For $x \in \mathbb{R}^d$, define M blocks indexed by \mathcal{B}_i , $i = 1, \dots, M$, with $d_i := |\mathcal{B}_i|$. The **Block-Sign** compressor is defined as $\mathcal{C}(x) = [\text{sign}(x_{\mathcal{B}_1}) \frac{\|x_{\mathcal{B}_1}\|_1}{d_1}, \dots, \text{sign}(x_{\mathcal{B}_M}) \frac{\|x_{\mathcal{B}_M}\|_1}{d_M}]$.*

Remark 1. *It is well-known [Stich et al. \(2018\)](#); [Zheng et al. \(2019\)](#) that for **Top- k** , $q^2 = 1 - \frac{k}{d}$; for **Block-Sign**, by Cauchy-Schwartz inequality we have $q^2 = 1 - \min_{i \in [M]} \frac{1}{d_i}$ where M and d_i are defined in Definition 2.*

The intuition behind **Top- k** is that, it has been observed during many deep neural networks training procedure, most gradients are typically very small and can be regarded as redundant—gradients with large magnitude contain most information. The **Block-Sign** compressor is a simple extension of the 1-bit **SIGN** compressor [Seide et al. \(2014\)](#); [Bernstein et al. \(2018\)](#), adapted to different gradient magnitude in different blocks, which, for neural nets, are usually set as the distinct network layers. The scaling factor in Definition 2 is to preserve the (possibly very different) gradient magnitude in each layer. In principle, **Top- k** would perform the best when the gradient is sparse, or only has a few very large absolute values, while **Block-Sign** compressor is beneficial when most gradients have similar magnitude within each layer.

3.2 COMP-AMS for Distributed Optimization

We present in Algorithm 2 our proposed communication-efficient distributed adaptive method in this paper, COMP-AMS. This framework can be regarded as an analogue to the standard synchronous distributed SGD protocol: in each iteration, each local worker transmits to the central server the compressed stochastic gradient computed using local data. Then the central server takes the average of local gradients, and performs an AMSGrad update. Despite that this method seems a straightforward extension of distributed SGD, no formal analysis of COMP-AMS has been conducted in literature, and training AMSGrad by compressed gradients has not been considered either.

In Algorithm 2, line 7-8 depict the error feedback operation at local nodes. $e_{t,i}$ is the accumulated error from gradient compression on the i -th worker up to time $t - 1$. This residual is added back to $g_{t,i}$ to get the “correct” gradient. In Section 4 and Section 5, we will show that error feedback, similar to the case of SGD, also brings good convergence behavior under gradient compression in adaptive optimization methods.

Comparison with Quantized Adam [Chen et al. \(2020\)](#). The first crucial difference of COMP-AMS compared with ([Chen et al., 2020](#)) which develops a quantized variant of Adam [Kingma and Ba \(2014\)](#) is that, in COMP-AMS, only compressed gradients are transmitted from the workers to the central server. In ([Chen et al., 2020](#)), each worker keeps a local copy of the moment estimates commonly noted m and v , and compresses and transmits the ratio $\frac{m}{v}$ as a whole to the server. Thus, that method is very much like the compressed distributed SGD, with the exception that the ratio $\frac{m}{v}$

Algorithm 2 Distributed COMP-AMS with error-feedback

```
1: Input: parameters  $\beta_1, \beta_2$ , learning rate  $\eta_t$ .
2: Initialize: central server parameter  $\theta_1 \in \Theta \subseteq \mathbb{R}^d$ ;  $e_{1,i} = 0$  the error accumulator for each worker;
   sparsity parameter  $k$ ;  $n$  local workers;  $m_0 = 0, v_0 = 0, \hat{v}_0 = 0$ 
3: for  $t = 1, \dots, T$  do
4:   parallel for worker  $i \in [n]$  do:
5:     Receive model parameter  $\theta_t$  from central server
6:     Compute stochastic gradient  $g_{t,i}$  at  $\theta_t$ 
7:     Compute  $\tilde{g}_{t,i} = \mathcal{C}(g_{t,i} + e_{t,i}, k)$ 
8:     Update the error  $e_{t+1,i} = e_{t,i} + g_{t,i} - \tilde{g}_{t,i}$ 
9:     Send  $\tilde{g}_{t,i}$  back to central server
10:  end parallel
11:  Central server do:
12:     $\bar{g}_t = \frac{1}{n} \sum_{i=1}^n \tilde{g}_{t,i}$ 
13:     $m_t = \beta_1 m_{t-1} + (1 - \beta_1) \bar{g}_t$ 
14:     $v_t = \beta_2 v_{t-1} + (1 - \beta_2) \bar{g}_t^2$ 
15:     $\hat{v}_t = \max(v_t, \hat{v}_{t-1})$ 
16:    Update the global model  $\theta_{t+1} = \theta_t - \eta_t \frac{m_t}{\sqrt{\hat{v}_t + \epsilon}}$ 
17: end for
```

plays the role of the gradient vector g communication-wise. Thus, two local moment estimators are additionally required, which have same size as the deep learning model. In our optimization method in Algorithm 2, the moment estimates m and v are kept and updated only at the central server, thus not introducing any extra variables (tensors) on local nodes during training (except for the error accumulator). In other words, COMP-AMS is not only effective in communication reduction, but also efficient in terms of memory (space), which is favorable for the distributed adaptive training of large-scale learners like BERT and CTR prediction models, e.g. Devlin et al. (2019); Zhao et al. (2020), to lower the hardware consumption happening in practice.

The second key difference is that, Adam is used as the underlying algorithm in Chen et al. (2020). It does not use the variable \hat{v} (line 15 of Algorithm 2) ensuring a non-decreasing second moment estimation, which has been shown to be an important factor for the convergence guarantee, see Reddi et al. (2018); Chen et al. (2019). Indeed, the convergence rate given by Chen et al. (2020) does not match the rate of vanilla AMSGrad and even does not converge to 0, when a decreasing learning rate, e.g., $\mathcal{O}(\frac{1}{\sqrt{T}})$, is employed, see (Chen et al., 2020, Th. 1). This would be intuitively problematic and given the fact that decreasing learning rate is standard in theoretical analysis and practical implementation. In next section, we prove the fast convergence of COMP-AMS using compressed gradients that matches the rate of full-gradient AMSGrad.

4 Convergence Analysis

For the convergence analysis of COMP-AMS we will make following additional assumptions.

Assumption 2. (*Smoothness*) For $\forall i \in [n]$, f_i is L -smooth: $\|\nabla f_i(\theta) - \nabla f_i(\vartheta)\| \leq L \|\theta - \vartheta\|$.

Assumption 3. (*Unbiased and bounded stochastic gradient*) For $\forall t > 0, \forall i \in [n]$, the stochastic gradient is unbiased and uniformly bounded: $\mathbb{E}[g_{t,i}] = \nabla f_i(\theta_t)$ and $\|g_{t,i}\| \leq G$.

Assumption 4. (*Bounded variance*) For $\forall t > 0, \forall i \in [n]$: (i) the **local variance** of the stochastic gradient is bounded: $\mathbb{E}[\|g_{t,i} - \nabla f_i(\theta_t)\|^2] < \sigma^2$; (ii) the **global variance** is bounded by $\frac{1}{n} \sum_{i=1}^n \|\nabla f_i(\theta_t) - \nabla f(\theta_t)\|^2 \leq \sigma_g^2$.

In Assumption 3, the uniform bound on the stochastic gradient is common in the convergence analysis of adaptive methods, e.g., Reddi et al. (2018); Zhou et al. (2018); Chen et al. (2019). The global variance bound σ_g^2 characterizes the difference among local objective functions, which, is mainly caused by different local data distribution \mathcal{X}_i in (1). In classical distributed setting where all the workers can access the same dataset and local data are assigned randomly, $\sigma_g^2 \equiv 0$. The scenario where \mathcal{X}_i 's are different gives rise to the recently proposed Federated Learning (FL) McMahan et al. (2017) framework where local data can be non-i.i.d. While typical FL method with periodical model averaging is beyond the scope of this present paper, we consider the global variance in our analysis to shed some light on the impact of non-i.i.d. data distribution in the federated setting for broader interest and future investigation.

Under the mild assumptions stated above, we derive the following general convergence rate of COMP-AMS in the distributed setting.

Theorem 1. Denote $C_0 = \sqrt{\frac{4(1+q^2)^3}{(1-q^2)^2} G^2} + \epsilon$, $C_1 = \frac{\beta_1}{1-\beta_1} + \frac{2q}{1-q^2}$, $\theta^* = \arg \min f(\theta)$. Under Assumption 1 to Assumption 4, with $\eta_t = \eta \leq \frac{\epsilon}{3C_0 \sqrt{2L \max\{2L, C_2\}}}$, COMP-AMS satisfies

$$\begin{aligned} \frac{1}{T} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] &\leq 2C_0 \left(\frac{\mathbb{E}[f(\theta_1) - f(\theta^*)]}{T\eta} + \frac{\eta L \sigma^2}{n\epsilon} + \frac{3\eta^2 L C_0 C_1 \sigma^2}{\epsilon^2} \right. \\ &\quad \left. + \frac{12\eta^2 q^2 L C_0 \sigma_g^2}{(1-q^2)^2 \epsilon^2} + \frac{(1+C_1)G^2 d}{T\sqrt{\epsilon}} + \frac{\eta(1+2C_1)C_1 L G^2 d}{T\epsilon} \right). \end{aligned}$$

The LHS of Theorem 1 is the expected squared norm of the gradient from a uniformly chosen iterate $t \in [T]$, which is a common convergence measure. From Theorem 1, we see that the more compression we apply to the gradient vectors (i.e., larger q), the larger the the gradient magnitude is, i.e., the slower the algorithm converges. This is intuitive as heavier compression loses more gradient information which would slower down the learner to find a good solution.

In the following paragraphs, we provide two interesting extension of our main result Theorem 1. We begin with the single-machine case, when $n = 1$ and then provide a linear speedup of our methods in the general distributed optimization case.

Single machine rate ($n = 1$). Note that, COMP-AMS with $n = 1$ naturally reduces to the single machine (sequential) AMSGrad (Algorithm 1) with compressed gradients instead of full-precision ones. The paper Karimireddy et al. (2019) shows that for SGD, error feedback fixes the convergence issue of biased compressors in the sense that SGD-EF (with compressed gradients) has the same convergence rate as vanilla SGD using full gradients. For AMSGrad, we have a similar result.

Corollary 1. Assume $n = 1$. Under Assumption 1 to Assumption 4, setting the stepsize as $\eta = \min\{\frac{\epsilon}{3C_0 \sqrt{2L \max\{2L, C_2\}}}, \frac{1}{\sqrt{T}}\}$, the sequence of iterates $\{\theta_t\}_{t>0}$ output from Algorithm 2 satisfies:

$$\frac{1}{T} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] \leq \mathcal{O}\left(\frac{1}{\sqrt{T}} + \frac{\sigma^2}{\sqrt{T}} + \frac{d}{T}\right).$$

Corollary 1 states that with error feedback, single machine AMSGrad with biased compressed gradients can also match the convergence rate $\mathcal{O}(\frac{1}{\sqrt{T}} + \frac{d}{T})$ of standard AMSGrad Zhou et al. (2018)

in non-convex optimization. It also achieves the same rate $\mathcal{O}(\frac{1}{\sqrt{T}})$ of vanilla SGD [Karimireddy et al. \(2019\)](#) when T is sufficiently large. In other words, EF also fixes the convergence issue of using compressed gradients in AMSGrad. To the best of our knowledge, this is the first result in literature showing that compressed adaptive methods with EF converges as fast as the standard counterpart. We will validate this benefit of EF in our numerical experiments.

Linear Speedup ($n > 1$). In Theorem 1, the convergence rate is derived assuming constant learning rate. By carefully choosing a decreasing learning rate dependent on the number of workers, we have the following simplified statement.

Corollary 2. *Under the same setting as Theorem 1, set $\eta = \min\{\frac{\epsilon}{3C_0\sqrt{2L\max\{2L, C_2\}}}, \frac{\sqrt{n}}{\sqrt{T}}\}$. Ignoring irrelevant quantities, we have*

$$\frac{1}{T} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] \leq \mathcal{O}\left(\frac{1}{\sqrt{nT}} + \frac{\sigma^2}{\sqrt{nT}} + \frac{n(\sigma^2 + \sigma_g^2)}{T}\right). \quad (2)$$

In Corollary 2, we see that the global variance σ_g^2 appears in the $\mathcal{O}(\frac{1}{T})$ term, which says that it asymptotically has no impact on the convergence. This matches the result of momentum SGD [Yu et al. \(2019a\)](#). When $T \geq \mathcal{O}(n^3)$ is sufficiently large, the third term in (2) vanishes, and the convergence rate becomes $\mathcal{O}(\frac{1}{\sqrt{nT}})$. Therefore, to reach a $\mathcal{O}(\delta)$ stationary point, one worker ($n = 1$) needs $T = \mathcal{O}(\frac{1}{\delta^2})$ iterations, while distributed training with n workers requires only $T = \mathcal{O}(\frac{1}{N\delta^2})$ iterations, which is n times faster than single machine training. That is, COMP-AMS has a linear speedup in terms of the number of the local workers. Such acceleration effect has also been reported for compressed SGD with error feedback [Zheng et al. \(2019\)](#); [Jiang and Agrawal \(2018\)](#) and momentum SGD [Yu et al. \(2019a\)](#). To our knowledge, this is also the first result showing the linear speedup for distributed adaptive methods under compression.

5 Experiments

In this section, we provide numerical results on several real-world datasets. Our main objective is to validate the theoretical results, and demonstrate that the proposed COMP-AMS can give satisfactory learning performance with significantly reduced communication costs.

5.1 Error Feedback Fixes the Convergence of Compressed AMSGrad

In Corollary 1, we established that for standard single machine AMSGrad under compression, EF helps achieve same convergence rate as the full-gradient AMSGrad. We implement COMP-AMS with a single worker and different gradient compressors, with or without EF, to justify this claim. We train MNIST [LeCun et al. \(1998\)](#) classification using a simple Convolutional Neural Network (CNN). The model has two convolutional layers followed by two fully connected layers with ReLU activation. Dropout is applied after the max-pooled convolutional layer with rate 0.5. The training batch size is 200. The parameters in COMP-AMS are set as default $\beta_1 = 0.9$ and $\beta_2 = 0.999$, which is true for all the experiments in this paper. The learning rate is searched over a fine grid and the best result is reported averaged over three runs. In Figure 1, the methods are (all performed on single worker):

- TopK-EF-0.01: COMP-AMS (Algorithm 2) and **Top- k** compressor, with sparsity 0.01 (i.e., keeping 1% gradient coordinates with largest magnitude).

- TopK-EF-0.001: COMP-AMS with **Top- k** compressor, with sparsity 0.001.
- BkSign-EF: COMP-AMS with **Block-Sign** compression.
- Methods without “-EF”: AMSGrad using corresponding compressors without error feedback (directly trained with compressed gradients).
- MVSS-0.01: AMSGrad with Minimal Variance Stochastic Sparsification (MVSS) [Wangni et al. \(2018\)](#) at 0.01 sparsity. This method probabilistically chooses gradient coordinates according to their magnitude, and divides each selected coordinate by its sampling probability to generate unbiased output (of the full gradient). Therefore, error feedback is not used for MVSS¹.

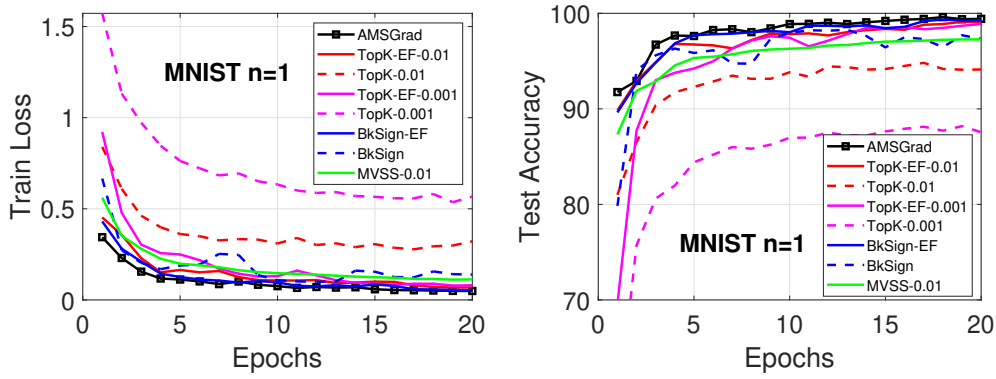


Figure 1 Training loss and test accuracy of COMP-AMS using a single machine.

From the train loss and test accuracy in Figure 1, we observe:

- AMSGrad without EF (dash curves) performs very poorly in terms of both convergence speed (more fluctuations in later epochs) and generalization (much worse test accuracy). With error feedback, the training loss and test accuracy both approach those of full-gradient AMSGrad with faster convergence. The issue of biased compression is fixed.
- **Top- k -EF-0.01** performs better than **Top- k -EF-0.001**, which justifies the influence of q in Theorem 1 that higher compression ratio would undermine the learning performance.
- MVSS-0.01 is outperformed by the proposed EF-corrected **Block-Sign** and **Top- k** even with 0.001 sparsity. This suggests that using biased compressors with EF in COMP-AMS is more effective than using unbiased stochastic compressors.

¹We also tested MVSS with error feedback, but the performance is uncompetitive with other methods.

5.2 Linear Speedup of COMP-AMS

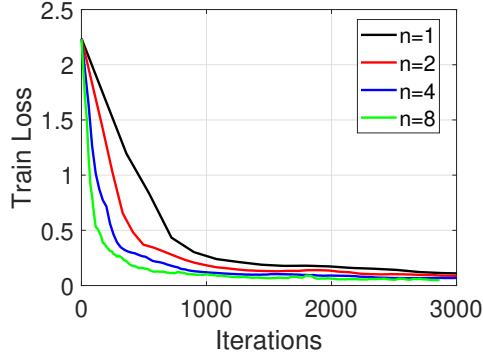


Figure 2 Training loss of COMP-AMS with **Top-k**-0.01 on MNIST.

Corollary 2 reveals the linear speedup of COMP-AMS in distributed training. We validate this claim in Figure 2, where the training loss on MNIST against the number of iterations is provided. Here we use COMP-AMS with **Top-k** and 0.01 sparsity. We test $n = 1, 2, 4, 8$, where the local mini-batch size is 100. As suggested by the theory, we use $10^{-4}\sqrt{n}$ as the learning rate. From Figure 2, we observe that the number of iterations for multiple workers to achieve a certain loss decreases as n increases. For example, to achieve 0.5 loss, we need around 700, 400, 240, 120 iterations for $n = 1, 2, 4, 8$ respectively, which decreases approximately linearly. This numerically justifies the linear speedup effect ($\mathcal{O}(\frac{1}{\sqrt{nT}})$ convergence rate) of COMP-AMS.

5.3 General Evaluation and Communication Efficiency

In this section, we present general evaluation of distributed training on more datasets and compare the communication efficiency. For CIFAR-10 dataset Krizhevsky et al. (2009), we train a larger CNN model with 3 convolutional layers. For IMDB movie review Maas et al. (2011) sentiment analysis, we train a LSTM network with 64 cells, equipped with an embedding layer which embeds top 1000 words in the movie reviews into 32-dimensional vectors. The local batch size is 100. The detailed architecture can be found in the supplement. We compare COMP-AMS with full gradient, **Top-k** and **Block-Sign** compression, with Quantized Adam (QADAM) Chen et al. (2020) (see supplement for the details). To match the compression ratio of COMP-AMS, we present 1-bit QAdam for highest communication reduction.

In Figure 3, we provide the test accuracy on three tasks against the number of epochs and the number of bits transmitted per local worker, respectively. On all three datasets, COMP-AMS with **Top-k**-0.01 and **Block-Sign** compression achieves same accuracy as using full gradients, with substantial 100x and 30x communication saving, respectively. While performing well on MNIST, more extreme compression (i.e., **Top-k**-0.001) leads to unnegligible loss in convergence speed and accuracy on the other two datasets. These results demonstrate the trade-off between compression and accuracy, and that under COMP-AMS framework, we can achieve significant (around 30-100x) communication reduction with similar model accuracy.

5.4 Discussion on COMP-AMS and QADAM

Noticeably, compared with QADAM, we see that on CIFAR-10 and IMDB, COMP-AMS converges to a worse stationary point/local optimum (worse generalization). Firstly, we should note that according to our theory and numerical results (e.g., Figure 2), distributed COMP-AMS in principle would at least have same learning performance as standard AMSGrad (trained on single machine) by properly choosing the learning rate. While the generalization of deep networks is still mysterious from rigorous theoretical perspective, we suspect that the performance gap of COMP-AMS might be a consequence of larger variance of the effective update, which is the price of abandoning local moment estimation. For the ease of illustration, we consider Adam (Algorithm 1 without line 7). In distributed adaptive training, we are essentially using the effective update ratio $\frac{m_t}{\sqrt{v_t}}$ by aggregating local information. For COMP-AMS, we first take the average (for the 1st and 2nd moments separately) and then take the ratio, while in QADAM we first take the ratio and then compute the average. Hence, in QADAM, there is only one additive error coming from linear averaging, which may better approximate the true ratio itself. On the other hand, since COMP-AMS is simpler and does not keep local moment estimates, we have two “sources” of error coming from both 1st and 2nd moment estimation, and the later introduces multiplicative error (from the denominator) which may lead to higher variance and more unstable estimation, from the statistical point of view. As a result, we can use larger learning rate for QADAM (indeed, the found optimal learning rate of QADAM is usually larger than that of COMP-AMS), which may speedup the algorithm and help escape local optima.

Nevertheless, as mentioned before, COMP-AMS is a very simple strategy that does not require extra local moment estimators (tensors), which is memory and hardware-friendly when training very large models in practice. Together with its high communication efficiency and linear speedup property, COMP-AMS would be a useful protocol in many applications. While in this paper we mainly consider the generic algorithm, we would like to place more investigation on the generalization ability (of both COMP-AMS and QADAM) and possible heuristics to further improve the performance of COMP-AMS (e.g., by warm-up or variance reduction techniques Liu et al. (2020)) as future work.

6 Conclusion

In this paper, we study the simple, convenient, yet unexplored gradient averaging strategy for distributed adaptive optimization called COMP-AMS. **Top- k** and **Block-Sign** compressor are incorporated for communication efficiency, whose biased is compensated by the error feedback strategy. We develop the convergence rate of COMP-AMS, and for the first time in literature show that for AMSGrad, gradient averaging with error feedback matches the convergence of full-gradient AMSGrad, and linear speedup can be obtained in distributed training. Numerical experiments are conducted to justify our theory, and demonstrate that similar performance can be achieved with significantly reduced communication. Though COMP-AMS generalizes slightly worse than QADAM since no local moment estimation is required, this in turn brings high memory/hardware efficiency that makes it a useful distributed learning scheme in practice. We hope that our work could originate more related research on the theory and application of distributed adaptive optimization.

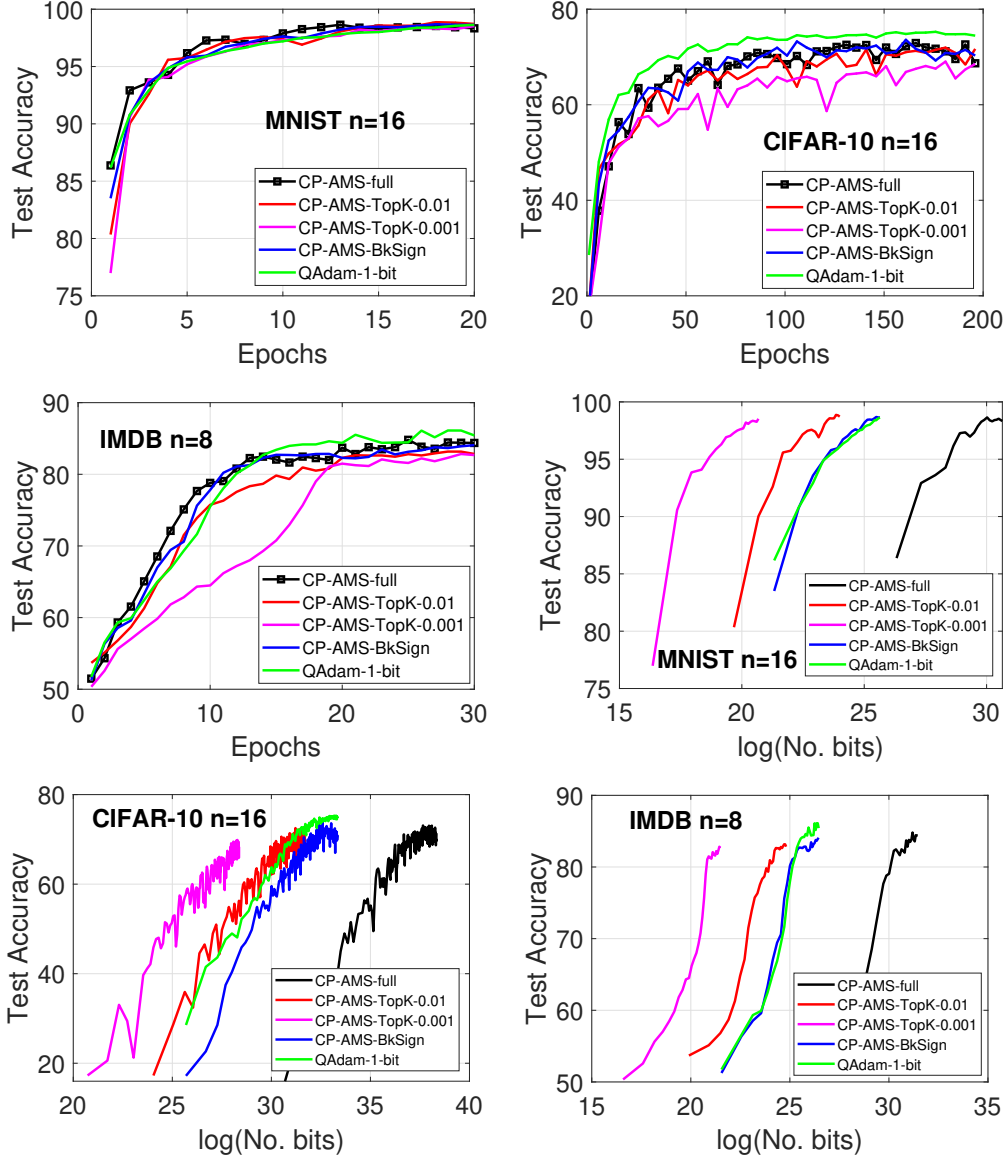


Figure 3 Test accuracy of distributed COMP-AMS. Top row: accuracy vs. epochs. Bottom row: accuracy vs. number of bits transmitted per worker.

Appendix

The supplementary material of this paper is organized in three main parts. Section A contains additional content and discussion such as the algorithmic formulation of the single-machine COMP-AMS and QADAM. Section B includes the proof of the main theoretical result. Section C contains more details on the experiments.

A Additional content

A.1 Extension to the Single-Machine Setting

In Corollary 1 we obtain the convergence rate of COMP-AMS in the single machine setting. Such setting has been fully considered in detail for SGD Karimireddy et al. (2019). For clarity, we provide in this subsection the formulation of our method in the single-worker setting, see Algorithm 3. Here, the computations, of the stochastic gradient and the various moment estimates, are all performed on a single-machine and the data is stored in this same worker.

Algorithm 3 COMP-AMS for a single-machine

- 1: **Input:** parameter β_1, β_2 , learning rate η_t .
 - 2: Initialize: central server parameter $\theta_1 \in \Theta \subseteq \mathbb{R}^d$; $e_1 = 0$ the error accumulator; sparsity parameter k ; $m_0 = 0, v_0 = 0, \hat{v}_0 = 0$
 - 3: **for** $t = 1, \dots, T$ **do**
 - 4: Compute stochastic gradient $g_t := g_{t,i_t}$ at θ_t for randomly sampled index i_t among the available observations indices
 - 5: Compute $\tilde{g}_t = \mathcal{C}(g_t + e_t)$
 - 6: Update the error $e_{t+1} = e_t + g_t - \tilde{g}_t$
 - 7: $m_t = \beta_1 m_{t-1} + (1 - \beta_1) \tilde{g}_t$
 - 8: $v_t = \beta_2 v_{t-1} + (1 - \beta_2) \tilde{g}_t^2$
 - 9: $\hat{v}_t = \max(v_t, \hat{v}_{t-1})$
 - 10: Update the model $\theta_{t+1} = \theta_t - \eta_t \frac{m_t}{\sqrt{\hat{v}_t + \epsilon}}$
 - 11: **end for**
-

A.2 QADAM Method

The closely related work to ours, QADAM discussed in Chen et al. (2020), is presented in Algorithm 4. Note that, the original method also compresses the model parameters in the server-to-worker communication, so we adapt it to one-way compression (only for the gradients) as our COMP-AMS. Here, $Q(\cdot)$ is a uniform quantization function that represents the effective update ratio m/\sqrt{v} using low bits. It is formally defined as

$$Q_b(g) = \|g\|_\infty \tilde{Q}_b(g/\|g\|_\infty),$$

where $\tilde{Q}_b(x) = \arg \min_{y \in M_b} \|y - x\|_2$, with $M_b := \{-1, -\frac{2^{b-1}-2}{2^{b-1}-1}, \dots, 0, \dots, \frac{2^{b-1}-2}{2^{b-1}-1}, 1\}$. As we can see, QADAM does not contain the \hat{v}_t term, and needs local moment estimations $m_{t,i}$ and $v_{t,i}$, for $i = 1, \dots, n$ on each worker. As discussed in the main paper, this costs substantially more memory and space when training large deep learning models.

Algorithm 4 QADAM [Chen et al. \(2020\)](#)

```

1: Input: parameters  $\beta_1, \beta_2$ , learning rate  $\eta_t$ .
2: Initialize: central server parameter  $\theta_1 \in \Theta \subseteq \mathbb{R}^d$ ;  $e_{1,i} = 0$  the error accumulator for each worker;
   sparsity parameter  $k$ ;  $n$  local workers; local moment estimate  $m_{0,i} = 0, v_{0,i} = 0$ 
3: for  $t = 1, \dots, T$  do
4:   parallel for worker  $i \in [n]$  do:
5:     Receive model parameter  $\theta_t$  from central server
6:     Compute stochastic gradient  $g_{t,i}$  at  $\theta_t$ 
7:      $m_{t,i} = \beta_1 m_{t-1,i} + (1 - \beta_1) g_{t,i}$ 
8:      $v_{t,i} = \beta_2 v_{t-1,i} + (1 - \beta_2) g_{t,i}^2$ 
9:      $a_{t,i} = \frac{m_{t,i}}{\sqrt{v_{t,i} + \epsilon}}$ 
10:    Compute  $\tilde{a}_{t,i} = Q(a_{t,i} + e_{t,i})$ 
11:    Update the error  $e_{t+1,i} = e_{t,i} + a_{t,i} - \tilde{a}_{t,i}$ 
12:    Send  $\tilde{a}_{t,i}$  back to central server
13:   end parallel
14:   Central server do:
15:    $\bar{a}_t = \frac{1}{n} \sum_{i=1}^n \tilde{a}_{t,i}$ 
16:   Update the global model  $\theta_{t+1} = \theta_t - \eta_t \bar{a}_t$ 
17: end for

```

B Proof of the Convergence Result

B.1 Proof of Theorem 1

Theorem. Denote $C_0 = \sqrt{\frac{4(1+q^2)^3}{(1-q^2)^2} G^2 + \epsilon}$, $C_1 = \frac{\beta_1}{1-\beta_1} + \frac{2q}{1-q^2}$. Under Assumption 1 to Assumption 4, with $\eta_t = \eta \leq \frac{\epsilon}{3C_0 \sqrt{2L \max\{2L, C_2\}}}$, for any $T > 0$, COMP-AMS satisfies

$$\begin{aligned}
\frac{1}{T} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] &\leq 2C_0 \left(\frac{\mathbb{E}[f(\theta_1) - f(\theta^*)]}{T\eta} + \frac{\eta L \sigma^2}{n\epsilon} + \frac{3\eta^2 L C_0 C_1 \sigma^2}{\epsilon^2} \right. \\
&\quad \left. + \frac{12\eta^2 q^2 L C_0 \sigma_g^2}{(1-q^2)^2 \epsilon^2} + \frac{(1+C_1)G^2 d}{T\sqrt{\epsilon}} + \frac{\eta(1+2C_1)C_1 L G^2 d}{T\epsilon} \right).
\end{aligned}$$

Proof. We first clarify some notations. At time t , let the full-precision gradient of the j -th worker be $g_{t,j}$, the error accumulator be $e_{t,j}$, and the compressed gradient be $\tilde{g}_{t,j} = \mathcal{C}(g_{t,j} + e_{t,j})$. Denote $\bar{g}_t = \frac{1}{n} \sum_{j=1}^n g_{t,j}$, $\bar{\tilde{g}}_t = \frac{1}{n} \sum_{j=1}^n \tilde{g}_{t,j}$ and $\bar{e}_t = \frac{1}{n} \sum_{j=1}^n e_{t,j}$. The second moment computed by the compressed gradients is denoted as $v_t = \beta_2 v_{t-1} + (1 - \beta_2) \bar{\tilde{g}}_t^2$, and $\hat{v}_t = \max\{\hat{v}_{t-1}, v_t\}$. Also, the first order moving average sequence

$$m_t = \beta_1 m_{t-1} + (1 - \beta_1) \bar{\tilde{g}}_t \quad \text{and} \quad m'_t = \beta_1 m'_{t-1} + (1 - \beta_1) \bar{g}_t.$$

By construction we have $m'_t = (1 - \beta_1) \sum_{i=1}^t \beta_1^{t-i} \bar{g}_i$.

Denote the following auxiliary sequences,

$$\mathcal{E}_{t+1} := (1 - \beta_1) \sum_{\tau=1}^{t+1} \beta_1^{t+1-\tau} \bar{e}_\tau$$

$$\theta'_{t+1} := \theta_{t+1} - \eta \frac{\mathcal{E}_{t+1}}{\sqrt{\hat{v}_t + \epsilon}}.$$

Then,

$$\begin{aligned} \theta'_{t+1} &= \theta_{t+1} - \eta \frac{\mathcal{E}_{t+1}}{\sqrt{\hat{v}_t + \epsilon}} \\ &= \theta_t - \eta \frac{(1 - \beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \bar{g}_\tau + (1 - \beta_1) \sum_{\tau=1}^{t+1} \beta_1^{t+1-\tau} \bar{e}_\tau}{\sqrt{\hat{v}_t + \epsilon}} \\ &= \theta_t - \eta \frac{(1 - \beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} (\bar{g}_\tau + \bar{e}_{\tau+1}) + (1 - \beta) \beta_1^t \bar{e}_1}{\sqrt{\hat{v}_t + \epsilon}} \\ &= \theta_t - \eta \frac{(1 - \beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \bar{e}_\tau}{\sqrt{\hat{v}_t + \epsilon}} - \eta \frac{m'_t}{\sqrt{\hat{v}_t + \epsilon}} \\ &= \theta_t - \eta \frac{\mathcal{E}_t}{\sqrt{\hat{v}_{t-1} + \epsilon}} - \eta \frac{m'_t}{\sqrt{\hat{v}_t + \epsilon}} + \eta \left(\frac{1}{\sqrt{\hat{v}_{t-1} + \epsilon}} - \frac{1}{\sqrt{\hat{v}_t + \epsilon}} \right) \mathcal{E}_t \\ &\stackrel{(a)}{=} \theta'_t - \eta \frac{m'_t}{\sqrt{\hat{v}_t + \epsilon}} + \eta \left(\frac{1}{\sqrt{\hat{v}_{t-1} + \epsilon}} - \frac{1}{\sqrt{\hat{v}_t + \epsilon}} \right) \mathcal{E}_t \\ &:= \theta'_t - \eta a'_t + \eta D_t \mathcal{E}_t, \end{aligned}$$

where (a) uses the fact that for every $j \in [n]$, $\tilde{g}_{t,j} + e_{t+1,j} = g_{t,j} + e_{t,j}$, and $e_{t,1} = 0$ at initialization. Further define the virtual iterates:

$$x_{t+1} := \theta'_{t+1} - \eta \frac{\beta_1}{1 - \beta_1} a'_t = \theta'_{t+1} - \eta \frac{\beta_1}{1 - \beta_1} \frac{m'_t}{\sqrt{\hat{v}_t + \epsilon}},$$

which follows the recurrence:

$$\begin{aligned} x_{t+1} &= \theta'_{t+1} - \eta \frac{\beta_1}{1 - \beta_1} \frac{m'_t}{\sqrt{\hat{v}_t + \epsilon}} \\ &= \theta'_t - \eta \frac{m'_t}{\sqrt{\hat{v}_t + \epsilon}} - \eta \frac{\beta_1}{1 - \beta_1} \frac{m'_t}{\sqrt{\hat{v}_t + \epsilon}} + \eta D_t \mathcal{E}_t \\ &= \theta'_t - \eta \frac{\beta_1 m'_{t-1} + (1 - \beta_1) \bar{g}_t + \frac{\beta_1^2}{1 - \beta_1} m'_{t-1} + \beta_1 \bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} + \eta D_t \mathcal{E}_t \\ &= \theta'_t - \eta \frac{\beta_1}{1 - \beta_1} \frac{m'_{t-1}}{\sqrt{\hat{v}_t + \epsilon}} - \eta \frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} + \eta D_t \mathcal{E}_t \\ &= x_t - \eta \frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} + \eta \frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} + \eta D_t \mathcal{E}_t. \end{aligned}$$

When summing over $t = 1, \dots, T$, the difference sequence D_t satisfies the bounds of Lemma 5. By Assumption 2 we have

$$f(x_{t+1}) \leq f(x_t) - \eta \langle \nabla f(x_t), x_{t+1} - x_t \rangle + \frac{L}{2} \|x_{t+1} - x_t\|^2.$$

Taking expectation w.r.t. the randomness at time t , we obtain

$$\begin{aligned} &\mathbb{E}[f(x_{t+1})] - f(x_t) \\ &\leq -\eta \mathbb{E}[\langle \nabla f(x_t), \frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} \rangle] + \eta \mathbb{E}[\langle \nabla f(x_t), \frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} + D_t \mathcal{E}_t \rangle] \end{aligned}$$

$$\begin{aligned}
& + \frac{\eta^2 L}{2} \mathbb{E}[\|\frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} - \frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} - D_t \mathcal{E}_t\|^2] \\
= & \underbrace{-\eta \mathbb{E}[\langle \nabla f(\theta_t), \frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} \rangle]}_I + \underbrace{\eta \mathbb{E}[\langle \nabla f(x_t), \frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} + D_t \mathcal{E}_t \rangle]}_{II} \\
& + \underbrace{\frac{\eta^2 L}{2} \mathbb{E}[\|\frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} - \frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} - D_t \mathcal{E}_t\|^2]}_{III} + \underbrace{\eta \mathbb{E}[\langle \nabla f(\theta_t) - \nabla f(x_t), \frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}} \rangle]}_{IV}, \quad (3)
\end{aligned}$$

Bounding term I. We have

$$\begin{aligned}
I &= -\eta \mathbb{E}[\langle \nabla f(\theta_t), \frac{\bar{g}_t}{\sqrt{\hat{v}_{t-1} + \epsilon}} \rangle] - \eta \mathbb{E}[\langle \nabla f(\theta_t), (\frac{1}{\sqrt{\hat{v}_t + \epsilon}} - \frac{1}{\sqrt{\hat{v}_{t-1} + \epsilon}}) \bar{g}_t \rangle] \\
&\leq -\eta \mathbb{E}[\langle \nabla f(\theta_t), \frac{\nabla f(\theta_t)}{\sqrt{\hat{v}_{t-1} + \epsilon}} \rangle] + \eta G^2 \mathbb{E}[\|D_t\|] \\
&\leq -\frac{\eta}{\sqrt{\frac{4(1+q^2)^3}{(1-q^2)^2} G^2 + \epsilon}} \mathbb{E}[\|\nabla f(\theta_t)\|^2] + \eta G^2 \mathbb{E}[\|D_t\|_1], \quad (4)
\end{aligned}$$

where we use Assumption 3, Lemma 4 and the fact that l_2 norm is no larger than l_1 norm.

Bounding term II. It holds that

$$\begin{aligned}
II &\leq \eta \mathbb{E}[\langle \nabla f(\theta_t), \frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} + D_t \mathcal{E}_t \rangle] + \mathbb{E}[\langle \nabla f(x_t) - \nabla f(\theta_t), \frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} + D_t \mathcal{E}_t \rangle] \\
&\leq \eta \mathbb{E}[\|\nabla f(\theta_t)\| \|\frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} + D_t \mathcal{E}_t\|] + \eta^2 L \mathbb{E}[\|\frac{\frac{\beta_1}{1 - \beta_1} m'_{t-1} + \mathcal{E}_t}{\sqrt{\hat{v}_{t-1} + \epsilon}}\| \|\frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} + D_t \mathcal{E}_t\|] \\
&\leq \eta C_1 G^2 \mathbb{E}[\|D_t\|_1] + \frac{\eta^2 C_1^2 L G^2}{\sqrt{\epsilon}} \mathbb{E}[\|D_t\|_1], \quad (5)
\end{aligned}$$

where $C_1 := \frac{\beta_1}{1 - \beta_1} + \frac{2q}{1 - q^2}$. The second inequality is because of smoothness of $f(\theta)$, and the last inequality is due to Lemma 2, Assumption 3 and the property of norms.

Bounding term III. This term can be bounded as follows:

$$\begin{aligned}
III &\leq \eta^2 L \mathbb{E}[\|\frac{\bar{g}_t}{\sqrt{\hat{v}_t + \epsilon}}\|^2] + \eta^2 L \mathbb{E}[\|\frac{\beta_1}{1 - \beta_1} D_t m'_{t-1} - D_t \mathcal{E}_t\|^2] \\
&\leq \frac{\eta^2 L}{\epsilon} \mathbb{E}[\|\frac{1}{n} \sum_{j=1}^i g_{t,j} - \nabla f(\theta_t) + \nabla f(\theta_t)\|^2] + \eta^2 L \mathbb{E}[\|D_t (\frac{\beta_1}{1 - \beta_1} m'_{t-1} - \mathcal{E}_t)\|^2] \\
&\stackrel{(a)}{\leq} \frac{\eta^2 L}{\epsilon} \mathbb{E}[\|\nabla f(\theta_t)\|^2] + \frac{\eta^2 L \sigma^2}{n \epsilon} + \eta^2 C_1^2 L G^2 \mathbb{E}[\|D_t\|^2], \quad (6)
\end{aligned}$$

where (a) follows from $\nabla f(\theta_t) = \frac{1}{n} \sum_{j=1}^n \nabla f_j(\theta_t)$ and Assumption 4 that $g_{t,j}$ is unbiased of $\nabla f_j(\theta_t)$ and has bounded variance σ^2 .

Bounding term IV. We have

$$\begin{aligned}
IV &= \eta \mathbb{E}[\langle \nabla f(\theta_t) - \nabla f(x_t), \frac{\bar{g}_t}{\sqrt{\hat{v}_{t-1} + \epsilon}} \rangle] + \eta \mathbb{E}[\langle \nabla f(\theta_t) - \nabla f(x_t), (\frac{1}{\sqrt{\hat{v}_t + \epsilon}} - \frac{1}{\sqrt{\hat{v}_{t-1} + \epsilon}}) \bar{g}_t \rangle] \\
&\leq \eta \mathbb{E}[\langle \nabla f(\theta_t) - \nabla f(x_t), \frac{\nabla f(\theta_t)}{\sqrt{\hat{v}_{t-1} + \epsilon}} \rangle] + \eta^2 L \mathbb{E}[\|\frac{\frac{\beta_1}{1 - \beta_1} m'_{t-1} + \mathcal{E}_t}{\sqrt{\hat{v}_{t-1} + \epsilon}}\| \|D_t g_t\|]
\end{aligned}$$

$$\begin{aligned}
&\stackrel{(a)}{\leq} \frac{\eta\rho}{2\epsilon} \mathbb{E}[\|\nabla f(\theta_t)\|^2] + \frac{\eta}{2\rho} \mathbb{E}[\|\nabla f(\theta_t) - \nabla f(x_t)\|^2] + \frac{\eta^2 C_1 L G^2}{\sqrt{\epsilon}} \mathbb{E}[\|D_t\|] \\
&\stackrel{(b)}{\leq} \frac{\eta\rho}{2\epsilon} \mathbb{E}[\|\nabla f(\theta_t)\|^2] + \frac{\eta^3 L}{2\rho} \mathbb{E}[\|\frac{\beta_1}{1-\beta_1} m'_{t-1} + \mathcal{E}_t\|^2] + \frac{\eta^2 C_1 L G^2}{\sqrt{\epsilon}} \mathbb{E}[\|D_t\|_1],
\end{aligned} \tag{7}$$

where (a) is due to Young's inequality and (b) is based on Assumption 2.

Regarding the second term in (7), by Lemma 3 and Lemma 1, summing over $t = 1, \dots, T$ we have

$$\begin{aligned}
&\sum_{t=1}^T \frac{\eta^3 L}{2\rho} \mathbb{E}[\|\frac{\beta_1}{1-\beta_1} m'_{t-1} + \mathcal{E}_t\|^2] \\
&\leq \sum_{t=1}^T \frac{\eta^3 L}{2\rho\epsilon} \mathbb{E}[\|\frac{\beta_1}{1-\beta_1} m'_{t-1} + \mathcal{E}_t\|^2] \\
&\leq \sum_{t=1}^T \frac{\eta^3 L}{\rho\epsilon} \left[\frac{\beta_1^2}{(1-\beta_1)^2} \mathbb{E}[\|m'_t\|^2] + \mathbb{E}[\|\mathcal{E}_t\|^2] \right] \\
&\leq \frac{T\eta^3 \beta_1^2 L \sigma^2}{\rho(1-\beta_1)^2 \epsilon} + \frac{\eta^3 \beta_1^2 L}{\rho(1-\beta_1)^2 \epsilon} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] \\
&\quad + \frac{4T\eta^3 q^2 L}{\rho(1-q^2)^2 \epsilon} (\sigma^2 + \sigma_g^2) + \frac{4\eta^3 q^2 L}{\rho(1-q^2)^2 \epsilon} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] \\
&= \frac{T\eta^3 L C_2 \sigma^2}{\rho\epsilon} + \frac{4T\eta^3 q^2 L \sigma_g^2}{\rho(1-q^2)^2 \epsilon} + \frac{\eta^3 L C_2}{\rho\epsilon} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2],
\end{aligned} \tag{8}$$

with $C_2 := \frac{\beta_1^2}{(1-\beta_1)^2} + \frac{4q^2}{(1-q^2)^2}$. Now integrating (4), (5), (6), (7) and (8) into (3), taking the telescoping summation over $t = 1, \dots, T$, we obtain

$$\begin{aligned}
&\mathbb{E}[f(x_{T+1}) - f(x_1)] \\
&\leq \left(-\frac{\eta}{C_0} + \frac{\eta^2 L}{\epsilon} + \frac{\eta\rho}{2\epsilon} + \frac{\eta^3 L C_2}{\rho\epsilon}\right) \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] + \frac{T\eta^2 L \sigma^2}{n\epsilon} + \frac{T\eta^3 L C_2 \sigma^2}{\rho\epsilon} + \frac{4T\eta^3 q^2 L \sigma_g^2}{\rho(1-q^2)^2 \epsilon} \\
&\quad + (\eta(1+C_1)G^2 + \frac{\eta^2(1+C_1)C_1 L G^2}{\sqrt{\epsilon}}) \sum_{t=1}^T \mathbb{E}[\|D_t\|_1] + \eta^2 C_1^2 L G^2 \sum_{t=1}^T \mathbb{E}[\|D_t\|^2].
\end{aligned}$$

with $C_0 := \sqrt{\frac{4(1+q^2)^3}{(1-q^2)^2} G^2} + \epsilon$. Setting $\eta \leq \frac{\epsilon}{3C_0 \sqrt{2L \max\{2L, C_2\}}}$ and choosing $\rho = \frac{\epsilon}{3C_0}$, we obtain

$$\begin{aligned}
&\mathbb{E}[f(x_{T+1}) - f(x_1)] \\
&\leq -\frac{\eta}{2C_0} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] + \frac{T\eta^2 L \sigma^2}{n\epsilon} + \frac{3T\eta^3 L C_0 C_2 \sigma^2}{\epsilon^2} + \frac{12T\eta^3 q^2 L C_0 \sigma_g^2}{(1-q^2)^2 \epsilon^2} \\
&\quad + \frac{\eta(1+C_1)G^2 d}{\sqrt{\epsilon}} + \frac{\eta^2(1+2C_1)C_1 L G^2 d}{\epsilon}.
\end{aligned}$$

where the last inequality follows from Lemma 5. Re-arranging terms, we get that

$$\frac{1}{T} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2] \leq 2C_0 \left(\frac{\mathbb{E}[f(x_1) - f(x_{T+1})]}{T\eta} + \frac{\eta L \sigma^2}{n\epsilon} + \frac{3\eta^2 L C_0 C_2 \sigma^2}{\epsilon^2} \right)$$

$$\begin{aligned}
& + \frac{12\eta^2 q^2 L C_0 \sigma_g^2}{(1-q^2)^2 \epsilon^2} + \frac{(1+C_1)G^2 d}{T\sqrt{\epsilon}} + \frac{\eta(1+2C_1)C_1 L G^2 d}{T\epsilon} \\
\leq & 2C_0 \left(\frac{\mathbb{E}[f(\theta_1) - f(\theta^*)]}{T\eta} + \frac{\eta L \sigma^2}{n\epsilon} + \frac{3\eta^2 L C_0 C_1 \sigma^2}{\epsilon^2} \right. \\
& \left. + \frac{12\eta^2 q^2 L C_0 \sigma_g^2}{(1-q^2)^2 \epsilon^2} + \frac{(1+C_1)G^2 d}{T\sqrt{\epsilon}} + \frac{\eta(1+2C_1)C_1 L G^2 d}{T\epsilon} \right),
\end{aligned}$$

where $C_0 = \sqrt{\frac{4(1+q^2)^3}{(1-q^2)^2} G^2} + \epsilon$, $C_1 = \frac{\beta_1}{1-\beta_1} + \frac{2q}{1-q^2}$. The last inequality is because $\theta'_1 = \theta_1$, $\theta^* := \arg \min_{\theta} f(\theta)$ and the fact that $C_2 \leq C_1$. This completes the proof. \square

Proofs of Corollary 2 and Corollary 1 follow naturally from the above.

B.2 Intermediary Lemmas

Lemma 1. *Under Assumption 1 to Assumption 4 we have:*

$$\sum_{t=1}^T \mathbb{E} \|\bar{m}'_t\|^2 \leq T\sigma^2 + \sum_{\tau=1}^t \mathbb{E} [\|\nabla f(\theta_t)\|^2].$$

Proof. Firstly, the expected squared norm of average stochastic gradient can be bounded by

$$\begin{aligned}
\mathbb{E} [\|\bar{g}_t^2\|] &= \mathbb{E} \left[\left\| \frac{1}{n} \sum_{i=1}^n g_{t,i} - \nabla f(\theta_t) + \nabla f(\theta_t) \right\|^2 \right] \\
&= \mathbb{E} \left[\left\| \frac{1}{n} \sum_{i=1}^n (g_{t,i} - \nabla f_i(\theta_t)) \right\|^2 \right] + \mathbb{E} [\|\nabla f(\theta_t)\|^2] \\
&\leq \sigma^2 + \mathbb{E} [\|\nabla f(\theta_t)\|^2],
\end{aligned}$$

where we use Assumption 4 that $g_{t,i}$ is unbiased and has bounded variance. Let $\bar{g}_{t,i}$ denote the i -th coordinate of \bar{g}_t . By the updating rule of COMP-AMS

$$\begin{aligned}
\mathbb{E} [\|\bar{m}'_t\|^2] &= \mathbb{E} \left[\left\| (1-\beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \bar{g}_\tau \right\|^2 \right] \\
&\leq (1-\beta_1)^2 \sum_{i=1}^d \mathbb{E} \left[\left(\sum_{\tau=1}^t \beta_1^{t-\tau} \bar{g}_{\tau,i} \right)^2 \right] \\
&\stackrel{(a)}{\leq} (1-\beta_1)^2 \sum_{i=1}^d \mathbb{E} \left[\left(\sum_{\tau=1}^t \beta_1^{t-\tau} \right) \left(\sum_{\tau=1}^t \beta_1^{t-\tau} \bar{g}_{\tau,i}^2 \right) \right] \\
&\leq (1-\beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \mathbb{E} [\|\bar{g}_\tau\|^2] \\
&\leq \sigma^2 + (1-\beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \mathbb{E} [\|\nabla f(\theta_t)\|^2],
\end{aligned}$$

where (a) is due to Cauchy-Schwartz inequality. Summing over $t = 1, \dots, T$, we obtain

$$\sum_{t=1}^T \mathbb{E} \|\bar{m}'_t\|^2 \leq T\sigma^2 + \sum_{t=1}^T \mathbb{E} [\|\nabla f(\theta_t)\|^2].$$

This completes the proof. □

Lemma 2. *Under Assumption 4, we have for $\forall t$ and each local worker $\forall i \in [n]$,*

$$\begin{aligned}\|e_{t,i}\|^2 &\leq \frac{4q^2}{(1-q^2)^2} G^2, \\ \mathbb{E}[\|e_{t+1,i}\|^2] &\leq \frac{4q^2}{(1-q^2)^2} \sigma^2 + \frac{2q^2}{1-q^2} \sum_{\tau=1}^t \left(\frac{1+q^2}{2}\right)^{t-\tau} \mathbb{E}[\|\nabla f_i(\theta_\tau)\|^2].\end{aligned}$$

Proof. We start by using Assumption 1 and Young's inequality to get

$$\begin{aligned}\|e_{t+1,i}\|^2 &= \|g_{t,i} + e_{t,i} - \mathcal{C}(g_{t,i} + e_{t,i})\|^2 \\ &\leq q^2 \|g_{t,i} + e_{t,i}\|^2 \\ &\leq q^2(1+\rho) \|e_{t,i}\|^2 + q^2\left(1 + \frac{1}{\rho}\right) \|g_{t,i}\|^2 \\ &\leq \frac{1+q^2}{2} \|e_{t,i}\|^2 + \frac{2q^2}{1-q^2} \|g_{t,i}\|^2,\end{aligned}\tag{9}$$

by choosing $\rho = \frac{1-q^2}{2q^2}$. Now by recursion and the initialization $e_{1,i} = 0$, we have

$$\begin{aligned}\mathbb{E}[\|e_{t+1,i}\|^2] &\leq \frac{2q^2}{1-q^2} \sum_{\tau=1}^t \left(\frac{1+q^2}{2}\right)^{t-\tau} \mathbb{E}[\|g_{\tau,i}\|^2] \\ &\leq \frac{4q^2}{(1-q^2)^2} \sigma^2 + \frac{2q^2}{1-q^2} \sum_{\tau=1}^t \left(\frac{1+q^2}{2}\right)^{t-\tau} \mathbb{E}[\|\nabla f_i(\theta_\tau)\|^2],\end{aligned}$$

which proves the second argument. Meanwhile, the absolute bound $\|e_{t,i}\|^2 \leq \frac{4q^2}{(1-q^2)^2} G^2$ follows directly from (9). □

Lemma 3. *For the moving average error sequence \mathcal{E}_t , it holds that*

$$\sum_{t=1}^T \mathbb{E}[\|\mathcal{E}_t\|^2] \leq \frac{4Tq^2}{(1-q^2)^2} (\sigma^2 + \sigma_g^2) + \frac{4q^2}{(1-q^2)^2} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2].$$

Proof. Let $\bar{e}_{t,i}$ be the j -th coordinate of \bar{e}_t . Denote $K_{t,i} := \sum_{\tau=1}^t \left(\frac{1+q^2}{2}\right)^{t-\tau} \mathbb{E}[\|\nabla f_i(\theta_\tau)\|^2]$ and $K_{t,i} = 0, \forall i \in [n]$. Using the same technique as in the proof of Lemma 1, we have

$$\begin{aligned}\mathbb{E}[\|\mathcal{E}_t\|^2] &= \mathbb{E}[\|(1-\beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \bar{e}_\tau\|^2] \\ &\leq (1-\beta_1)^2 \sum_{j=1}^d \mathbb{E}[\left(\sum_{\tau=1}^t \beta_1^{t-\tau} \bar{e}_{\tau,j}\right)^2] \\ &\stackrel{(a)}{\leq} (1-\beta_1)^2 \sum_{j=1}^d \mathbb{E}[\left(\sum_{\tau=1}^t \beta_1^{t-\tau}\right) \left(\sum_{\tau=1}^t \beta_1^{t-\tau} \bar{e}_{\tau,j}^2\right)]\end{aligned}$$

$$\begin{aligned}
&\leq (1 - \beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \mathbb{E}[\|\bar{e}_\tau\|^2] \\
&\leq (1 - \beta_1) \sum_{\tau=1}^t \beta_1^{t-\tau} \mathbb{E}\left[\frac{1}{n} \sum_{i=1}^n \|e_{\tau,i}\|^2\right] \\
&\stackrel{(b)}{\leq} \frac{4q^2}{(1 - q^2)^2} \sigma^2 + \frac{2q^2(1 - \beta_1)}{(1 - q^2)} \sum_{\tau=1}^t \beta_1^{t-\tau} \left(\frac{1}{n} \sum_{i=1}^n K_{\tau,i}\right),
\end{aligned}$$

where (a) is due to Cauchy-Schwartz and (b) is a result of Lemma 2. Summing over $t = 1, \dots, T$ and using the technique of geometric series summation leads to

$$\begin{aligned}
\sum_{t=1}^T \mathbb{E}[\|\mathcal{E}_t\|^2] &= \frac{4Tq^2}{(1 - q^2)^2} \sigma^2 + \frac{2q^2(1 - \beta_1)}{(1 - q^2)} \sum_{t=1}^T \sum_{\tau=1}^t \beta_1^{t-\tau} \left(\frac{1}{n} \sum_{i=1}^n K_{\tau,i}\right) \\
&\leq \frac{4Tq^2}{(1 - q^2)^2} \sigma^2 + \frac{2q^2}{(1 - q^2)} \sum_{t=1}^T \sum_{\tau=1}^t \left(\frac{1 + q^2}{2}\right)^{t-\tau} \mathbb{E}\left[\frac{1}{n} \sum_{i=1}^n \|\nabla f_i(\theta_\tau)\|^2\right] \\
&\leq \frac{4Tq^2}{(1 - q^2)^2} \sigma^2 + \frac{4q^2}{(1 - q^2)^2} \sum_{t=1}^T \mathbb{E}\left[\frac{1}{n} \sum_{i=1}^n \|\nabla f_i(\theta_t)\|^2\right] \\
&\stackrel{(a)}{\leq} \frac{4Tq^2}{(1 - q^2)^2} \sigma^2 + \frac{4q^2}{(1 - q^2)^2} \sum_{t=1}^T \mathbb{E}\left[\left\|\frac{1}{n} \sum_{i=1}^n \nabla f_i(\theta_t)\right\|^2 + \frac{1}{n} \sum_{i=1}^n \|\nabla f_i(\theta_t) - \nabla f(\theta_t)\|^2\right] \\
&\leq \frac{4Tq^2}{(1 - q^2)^2} (\sigma^2 + \sigma_g^2) + \frac{4q^2}{(1 - q^2)^2} \sum_{t=1}^T \mathbb{E}[\|\nabla f(\theta_t)\|^2],
\end{aligned}$$

where (a) is derived by the variance decomposition and the last inequality holds due to Assumption 4. The desired result is obtained. \square

Lemma 4. *It holds that $\forall t \in [T], \forall i \in [d], \hat{v}_{t,i} \leq \frac{4(1+q^2)^3}{(1-q^2)^2} G^2$.*

Proof. For any t , by Lemma 2 and Assumption 3 we have

$$\begin{aligned}
\|\tilde{g}_t\|^2 &= \|\mathcal{C}(g_t + e_t)\|^2 \\
&\leq \|\mathcal{C}(g_t + e_t) - (g_t + e_t) + (g_t + e_t)\|^2 \\
&\leq 2(q^2 + 1)\|g_t + e_t\|^2 \\
&\leq 4(q^2 + 1)(G^2 + \frac{4q^2}{(1 - q^2)^2} G^2) \\
&= \frac{4(1 + q^2)^3}{(1 - q^2)^2} G^2.
\end{aligned}$$

It's then easy to show by the updating rule of \hat{v}_t ,

$$\hat{v}_{t,i} = (1 - \beta_2) \sum_{\tau=1}^t \beta_2^{t-\tau} \tilde{g}_{\tau,i}^2 \leq \frac{4(1 + q^2)^3}{(1 - q^2)^2} G^2.$$

\square

Lemma 5. Let $D_t := \frac{1}{\sqrt{\hat{v}_{t-1} + \epsilon}} - \frac{1}{\sqrt{\hat{v}_t + \epsilon}}$ be defined as above. Then,

$$\sum_{t=1}^T \|D_t\|_1 \leq \frac{d}{\sqrt{\epsilon}}, \quad \sum_{t=1}^T \|D_t\|^2 \leq \frac{d}{\epsilon}.$$

Proof. By the updating rule of COMP-AMS, $\hat{v}_{t-1} \leq \hat{v}_t$ for $\forall t$. Therefore, by the initialization $\hat{v}_0 = 0$, we have

$$\begin{aligned} \sum_{t=1}^T \|D_t\|_1 &= \sum_{t=1}^T \sum_{i=1}^d \left(\frac{1}{\sqrt{\hat{v}_{t-1,i} + \epsilon}} - \frac{1}{\sqrt{\hat{v}_{t,i} + \epsilon}} \right) \\ &= \sum_{i=1}^d \left(\frac{1}{\sqrt{\hat{v}_{0,i} + \epsilon}} - \frac{1}{\sqrt{\hat{v}_{T,i} + \epsilon}} \right) \\ &\leq \frac{d}{\sqrt{\epsilon}}. \end{aligned}$$

For the sum of squared l_2 norm, note the fact that for $a \geq b > 0$, it holds that

$$(a - b)^2 \leq (a - b)(a + b) = a^2 - b^2.$$

Thus,

$$\begin{aligned} \sum_{t=1}^T \|D_t\|^2 &= \sum_{t=1}^T \sum_{i=1}^d \left(\frac{1}{\sqrt{\hat{v}_{t-1,i} + \epsilon}} - \frac{1}{\sqrt{\hat{v}_{t,i} + \epsilon}} \right)^2 \\ &\leq \sum_{t=1}^T \sum_{i=1}^d \left(\frac{1}{\hat{v}_{t-1,i} + \epsilon} - \frac{1}{\hat{v}_{t,i} + \epsilon} \right) \\ &\leq \frac{d}{\epsilon}, \end{aligned}$$

which gives the desired result. □

C Model Architecture of the Experiments

In Figure 4, we provide the detailed description of the data and model architectures used in our numerical study. MNIST [LeCun et al. \(1998\)](#) is a popular hand-written letter recognition dataset, where each training sample is a 28×28 black and white image belonging to a class (digits 0-9). CIFAR-10 [Krizhevsky et al. \(2009\)](#) is a benchmark image classification dataset consisting of natural images from 10 classes. The image size is a $3 \times 32 \times 32$. In IMDB dataset [Maas et al. \(2011\)](#), each sample is a movie review, and the task is to classify the reviews as positive or negative. The reviews are tokenized by words and transformed into integer vectors. We threshold at 300 for the length of each review. Zero-padding is applied to reviews that have less than 300 words. All our experiments are trained on a Linux server equipped with four Nvidia Tesla V100 cards. We use two Convolutional Neural Networks (CNN) for MNIST and CIFAR-10. For IMDB dataset, we use a LSTM network. For all three models, ReLu activation is adopted. For LSTM, each input movie review is a 300-dimensional vector, and the embedding layer embeds top 1000 most frequent words into 32-dimensional vectors. 64 LSTM cells are used, where the last hidden state is connected to two fully connected layers before the output.

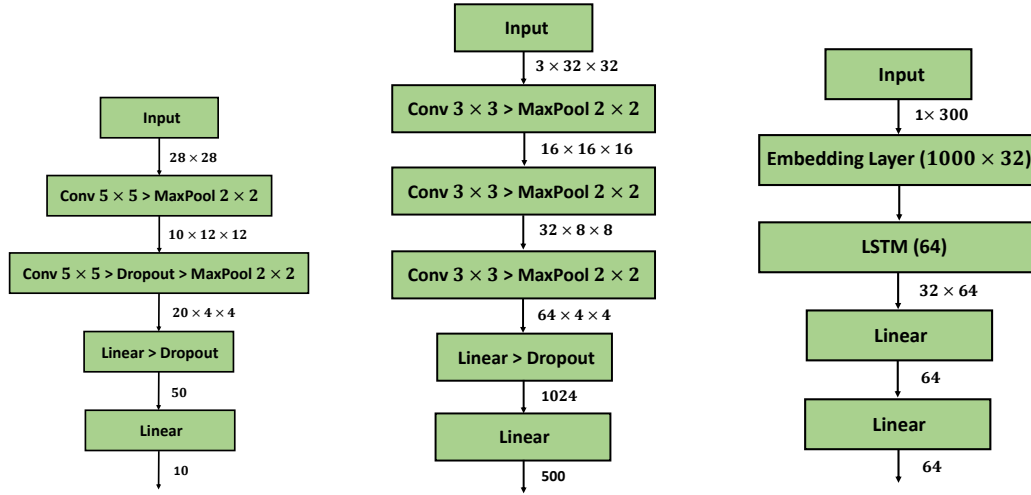


Figure 4 Model architectures used in the experiments. Left: MNIST + CNN. Middle: CIFAR-10 + CNN. Right: IMDB + LSTM. In the last figure, the penultimate linear layer takes the last hidden state (64-dim vector) as the input.

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