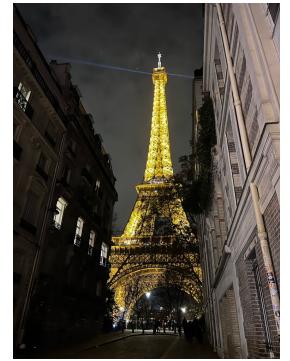


KAUST

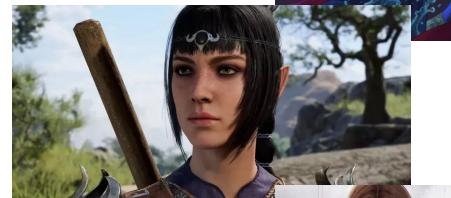
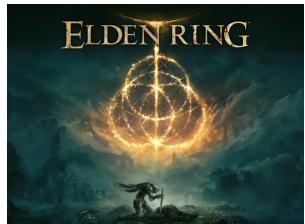


Paris

The First Optimal Parallel SGD (in the Presence of Data, Compute and Communication Heterogeneity)

Peter Richtárik

King Abdullah University of Science and Technology
Kingdom of Saudi Arabia



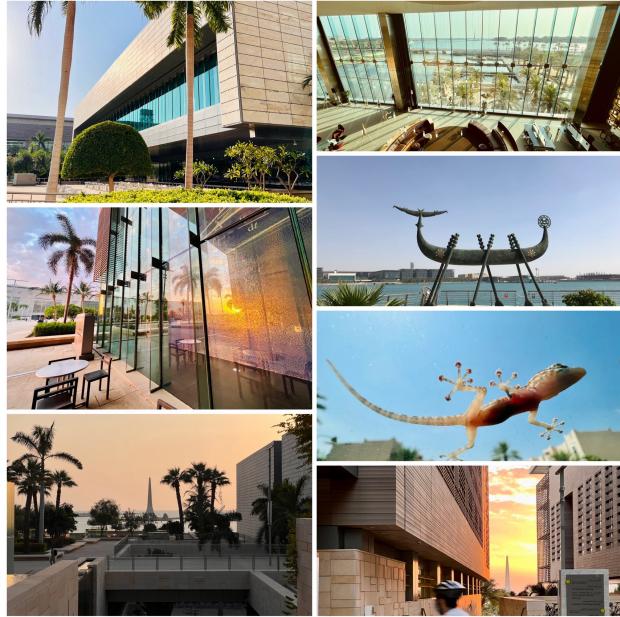
Applied Algorithms for
Machine Learning

A WORKSHOP ON FUTURE OF COMPUTATION

Paris

June 10-12, 2024

Optimization & Machine Learning Lab @ KAUST

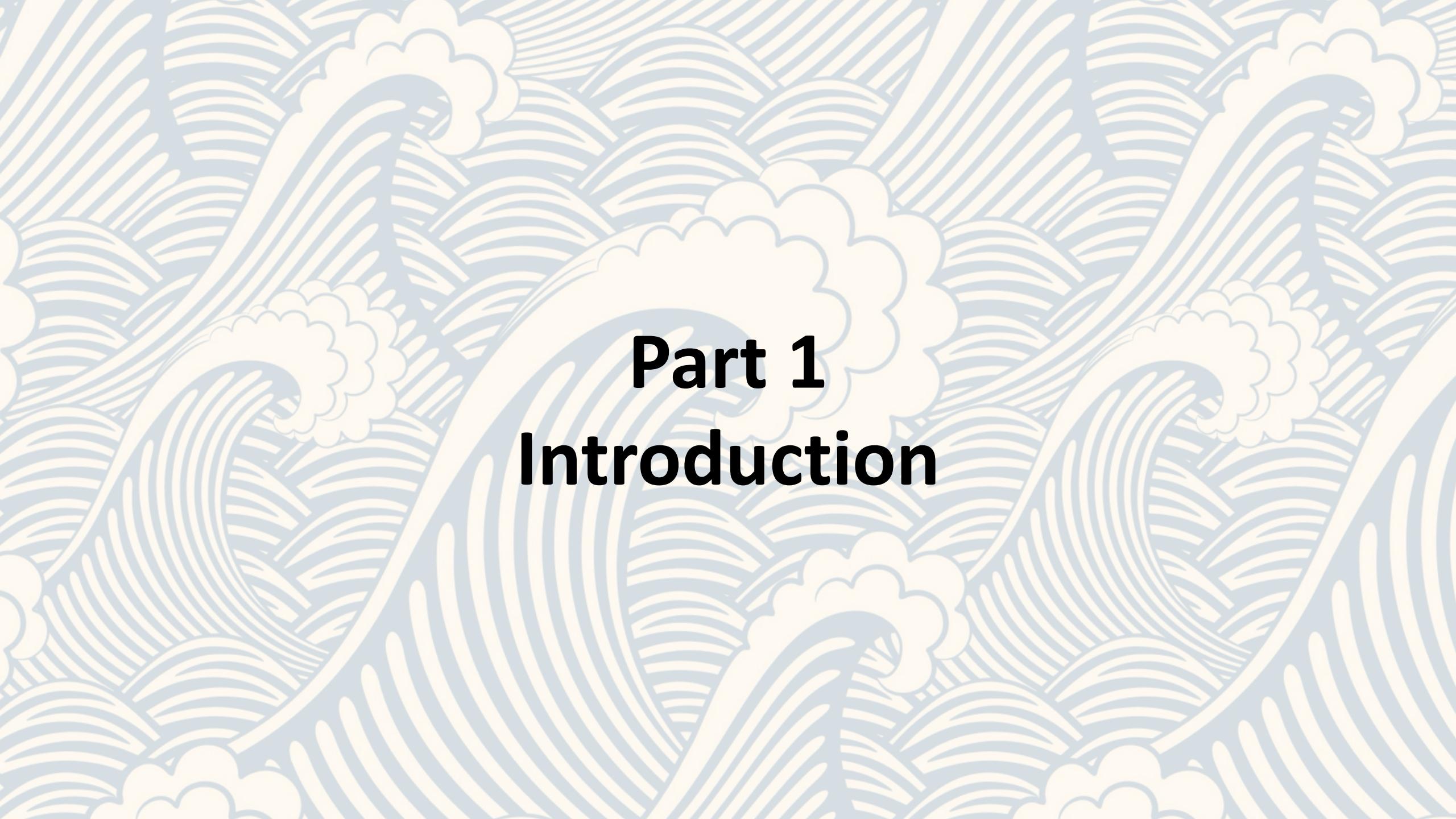


Kaja Gruntkowska

Ivan Ilin

Alexander Tyurin





Part 1

Introduction

Optimization Problem

$$\min_{x \in \mathbb{R}^d} f(x) \stackrel{\text{def}}{=} \frac{1}{n} \sum_{i=1}^n f_i(x)$$

model parameters / features

parallel machines

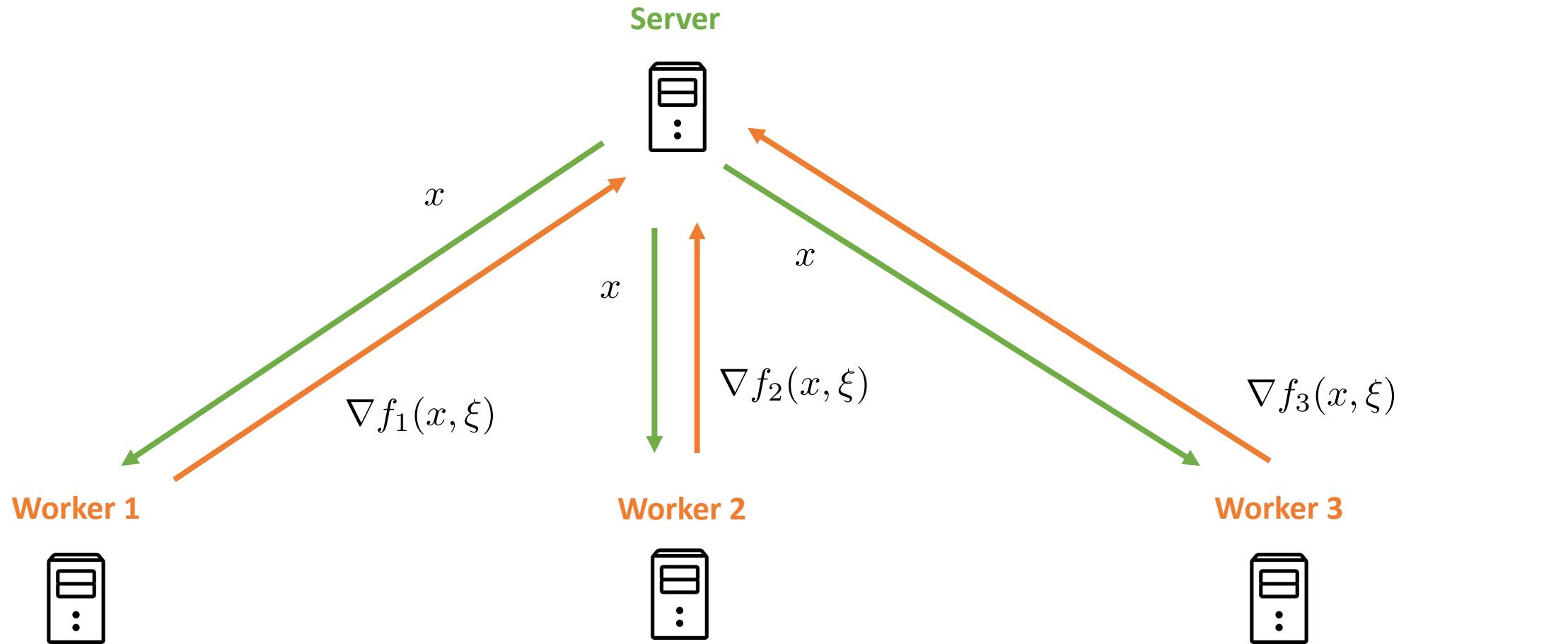
Loss on local data \mathcal{D}_i stored on machine i
 $f_i(x) := \mathbb{E}_{\xi \sim \mathcal{D}_i} [f_i(x, \xi)]$

- ! It takes τ_i seconds for worker i to compute $\nabla f_i(x, \xi)$, where $\xi \sim \mathcal{D}_i$ $0 < \tau_1 \leq \tau_2 \leq \dots \leq \tau_n$
- ! It takes θ_i seconds for worker i to communicate $g \in \mathbb{R}^d$ to the server

Find a (possibly random) vector $\hat{x} \in \mathbb{R}^d$ such that $\mathbb{E} [\|\nabla f(\hat{x})\|^2] \leq \varepsilon$

Parallel Computing Architecture

x gets updated by the server



$$f_1(x) := \mathbb{E}_{\xi \sim \mathcal{D}_1} [f_1(x, \xi)]$$

$$f_2(x) := \mathbb{E}_{\xi \sim \mathcal{D}_2} [f_2(x, \xi)]$$

$$f_3(x) := \mathbb{E}_{\xi \sim \mathcal{D}_3} [f_3(x, \xi)]$$

$\nabla f_1(x, \xi)$ compute time = τ_1 secs $\nabla f_2(x, \xi)$ compute time = τ_2 secs $\nabla f_3(x, \xi)$ compute time = τ_3 secs

Three Types of Heterogeneity

Data	data distributions $\mathcal{D}_1, \dots, \mathcal{D}_n$ can be different
Compute	compute times τ_1, \dots, τ_n are nonzero and can be different
Communication	communication times $\theta_1, \dots, \theta_n$ are nonzero and can be different

Typical Assumptions

1

$$\inf f \in \mathbb{R}$$

2

$$f_i(x) := \mathbb{E}_{\xi \sim \mathcal{D}_i} [f_i(x, \xi)]$$

Gradient of local functions is Lipschitz:

$$\max_{i \in \{1, \dots, n\}} \sup_{x \neq y} \frac{\|\nabla f_i(x) - \nabla f_i(y)\|}{\|x - y\|} \leq L$$



Stochastic gradients have bounded variance:

$$\max_{i \in \{1, \dots, n\}} \sup_{x \in \mathbb{R}^d} \mathbb{E}_{\xi \sim \mathcal{D}_i} [\|\nabla f_i(x, \xi) - \mathbb{E}_{\xi \sim \mathcal{D}_i} [\nabla f_i(x, \xi)]\|^2] \leq \sigma^2$$

Our Papers on Optimal Parallel SGD

Optimal Time Complexities of Parallel Stochastic Optimization Methods Under a Fixed Computation Model

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Abstract

Parallelization is a popular strategy to improve the performance of iterative algorithms. Optimization theory and its empirical design of efficient parallel optimization methods and tight analysis of their theoretical properties are important research endeavors. While the minimax complexities are well known for sequential optimization methods, the time of parallel optimization methods is less explored. In this paper, we propose a new framework for analyzing the time complexities of parallel optimization methods under a fixed computation model. We establish *minimax complexities* for parallel optimization methods that have access to an unbiased stochastic gradient oracle with bounded variance. We consider a fixed computation model characterized by two worker requirements: the broadcast time to calculate a stochastic gradient. We prove lower bounds and develop optimal algorithms that attain them. Our results have surprising consequences for the literature of asynchronous optimization methods.

1 Introduction

We consider the nonconvex optimization problem

$$\min_{x \in \mathbb{R}^d} \{f(x) := \mathbb{E}_{\xi \sim \mathcal{D}} [f(x; \xi)]\}, \quad (1)$$

where $f: \mathbb{R}^d \times S_\xi \rightarrow \mathbb{R}$, $S_\xi \subseteq \mathbb{R}^d$ and ξ is a random variable with some distribution \mathcal{D} on S_ξ . In machine learning, S_ξ could be the space of all possible data, \mathcal{D} is the distribution of the training dataset, and $f(\cdot, \xi)$ is the loss of a data sample ξ . In this paper we address the following natural setup:

- (i) n workers are able to work in parallel,
- (ii) the i^{th} worker requires τ_i seconds¹ to calculate a stochastic gradient of f .

The function f is L -smooth and lower-bounded (see Assumptions 7.1–7.2), and stochastic gradients are unbiased and σ^2 -variance-bounded (see Assumption 7.3).

1.1 Classical theory

In the nonconvex setting, gradient descent (GD) is an optimal method with respect to the number of gradient (∇f) calls (Lan, 2020; Nesterov, 2018; Carmon et al., 2020) for finding an approximately stationary point of f . Obviously, a key issue with GD is that it requires access to the exact gradients

¹Or any other unit of time.

37th Conference on Neural Information Processing Systems (NeurIPS 2023).

Shadowheart SGD: Distributed Asynchronous SGD with Optimal Time Complexity Under Arbitrary Computation and Communication Heterogeneity

Alexander Tyurin¹ Marta Pozzi^{1,2} Ivan Ilin¹ Peter Richtárik¹

Abstract

We consider *nonconvex stochastic optimization problems* in the *asynchronous distributed setting* in which the computation times from workers to a server can be ignored, and the computation and communication times are potentially different for all workers. Using an unbiased compression technique, we develop a new *asynchronous SGD* algorithm that significantly improves the time complexities of all previous centralized methods.

Moreover, we show that the time complexity of *Shadowheart SGD* is optimal in the family of centralized methods with compressed communication. We also consider the decentralized setup, where broadcasting from the server to the workers is non-negligible, and develop a corresponding method.

1. Introduction

We consider the nonconvex smooth optimization problem

$$\min_{x \in \mathbb{R}^d} \{f(x) := \mathbb{E}_{\xi \sim \mathcal{D}_\xi} [f(x; \xi)]\}, \quad (1)$$

where $f(\cdot, \cdot): \mathbb{R}^d \times S_\xi \rightarrow \mathbb{R}$, and \mathcal{D}_ξ is a distribution on $S_\xi \neq \emptyset$. Given $\varepsilon > 0$, we seek to find a probability random point \hat{x} such that $\mathbb{E}[\|\nabla f(\hat{x})\|^2] \leq \varepsilon$. Such a point is called an ε -stationary point. We focus on solving the problem in a distributed setting.

We rely on assumptions which are standard in the literature on stochastic gradient methods: smoothness, lower-boundedness and bounded variance.

Assumption 1.1. f is differentiable and L -smooth, i.e., $\|\nabla f(x) - \nabla f(y)\| \leq L \|x - y\|, \forall x, y \in \mathbb{R}^d$.

Assumption 1.2. There exist $f^*: \mathbb{R} \rightarrow \mathbb{R}$ such that $f(x) \geq f^*$ for all $x \in \mathbb{R}^d$. We define $\hat{f}(x) := f(x^*)$, where $x^* \in \mathbb{R}^d$ is a starting point of all algorithms we consider.

Assumption 1.3. For all $x \in \mathbb{R}^d$, the stochastic gradients $\nabla f(x; \xi)$ are unbiased, and their variance is bounded by $\sigma^2 \geq 0$, i.e., $\mathbb{E}[\nabla f(x; \xi)] = \nabla f(x)$ and $\mathbb{E}[\|\nabla f(x; \xi) - \nabla f(x)\|^2] \leq \sigma^2$.

To simplify the exposition, in what follows (up to Sec. 7) we first focus on the regime in which the broadcast cost can be ignored. We describe a strategy for extending our algorithm to the more general regime in Sec. 8.

arXiv:2402.04785v1 [math.OC] 7 Feb 2024

2/2024

Freya PAGE: First Optimal Time Complexity for Large-Scale Nonconvex Finite-Sum Optimization with Heterogeneous Asynchronous Computations

Alexander Tyurin¹ Kaja Gruntkowska¹ Peter Richtárik¹

Abstract

In practical distributed systems, workers are typically not homogeneous, and due to differences in hardware configurations and network conditions, can have highly varying processing times. We consider smooth nonconvex finite-sum (empirical risk minimization) problems that have been introduced in the nearly-optimal method Freya PAGE, designed to handle arbitrarily heterogeneous and asynchronous computations. By being robust to “stragglers” and adaptively ignoring slow computations, Freya PAGE offers significantly improved *time complexity* guarantees compared to all previously proposed asynchronous optimization methods, Fragile SGD and PAGE, while requiring weaker assumptions. The algorithm relies on novel generic stochastic gradient collection strategies with theoretical guarantees that can be of interest on their own, and may be used in the design of future optimization methods. Furthermore, we prove that Freya PAGE solves the nonconvex finite-sum problems in the asynchronous setup, providing a fundamental time complexity limit. This lower bound is tight and demonstrates the optimality of Freya PAGE in the large-scale regime, i.e., when $\sqrt{m} \geq n$, where n is # of workers, and m is # of data samples.

1 Introduction

In real-world distributed systems used for large-scale machine learning tasks, it is common to encounter device heterogeneity and variations in processing times among different computational units. These can stem from GPU computational delays, disparities in hardware configurations, network conditions, and the number of available devices or distributed across multiple devices (Chen et al., 2016; Tyurin and Richtárik, 2023). As a result, some clients may execute computations faster, while others experience delays or even fail to participate in the training altogether.

Due to the above reasons, we aim to address the challenges posed by device heterogeneity in the context of solving finite-sum nonconvex optimization problems of the form

$$\min_{x \in \mathbb{R}^d} \{f(x) := \frac{1}{m} \sum_{i=1}^m f_i(x)\}, \quad (1)$$

where $f_i: \mathbb{R}^d \rightarrow \mathbb{R}$ can be viewed as the loss of a machine learning model x on the i^{th} example in a training dataset with m samples. Our goal is to find an ε -stationary point, i.e., a (possibly random) point \hat{x} such that $\mathbb{E}[\|\nabla f(\hat{x})\|^2] \leq \varepsilon$. We focus on the homogeneous distributed setup:

- there are n workers/devices able to work in parallel,
- each worker has access to stochastic gradients $\nabla f_i, j \in [m]$,
- worker i calculates $\nabla f_i(\cdot)$ in less or equal to $\tau_i \in [0, \infty)$ seconds for all $i \in [n], j \in [m]$.

5/2024

On the Optimal Time Complexities in Decentralized Stochastic Asynchronous Optimization

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(alexanderturiu, richtarik)@gmail.com

Abstract

We consider the decentralized stochastic asynchronous optimization setup, where many workers asynchronously calculate stochastic gradients and asynchronously send them to a central server. We prove new time complexity lower bounds under the assumption that computation and communication speeds are bounded. We develop a new nearly optimal method, Fragile SGD, and a new optimal method, Amelie SGD, that converge under arbitrary heterogeneous computation and communication speeds. We find that the lower bounds are much better than the homogeneous setting. Our time complexities are new, nearly optimal, and provably improve all previous asynchronous/stochastic methods in the decentralized setup.

1 Introduction

We consider the smooth nonconvex optimization problem

$$\min_{x \in \mathbb{R}^d} \{f(x) := \mathbb{E}_{\zeta \sim D_\zeta} [f(x; \zeta)]\}, \quad (1)$$

where $f: \mathbb{R}^d \times S_\zeta \rightarrow \mathbb{R}$, and D_ζ is a distribution on a non-empty set S_ζ . For a given $\varepsilon > 0$, we want to find a possibly random point \hat{x} , called an ε -stationary point, such that $\mathbb{E}[\|\nabla f(\hat{x})\|^2] \leq \varepsilon$. We analyze the heterogeneous setup and the convex setup with smooth and non-smooth functions in Sections B and C.

1.1 Decentralized setup with times

We investigate the following decentralized asynchronous setup. Assume that we have n workers/nodes with the associated computation times $\{\bar{\tau}_i\}$, and communications times $\{\bar{\rho}_{i,j}\}$. It takes less or equal to $\bar{\tau}_i \in [0, \infty)$ seconds to compute a stochastic gradient by the i^{th} node, and less or equal to $\bar{\rho}_{i,j} \in [0, \infty)$ seconds to send a vector $v \in \mathbb{R}^d$ from the i^{th} node to the j^{th} node (it is possible that $\bar{\tau}_i = \infty$ and $\bar{\rho}_{i,j} = \infty$). All computation and communication times can be drawn independently and in parallel. We would like to emphasize that $\bar{\tau}_i \in [0, \infty]$ and $\bar{\rho}_{i,j} \in [0, \infty]$ are only upper bounds, and the real and effective computation and communication times can be arbitrarily heterogeneous and independently of ε . We assume that the upper bounds are static; however, in Section 5.5, we explain that our result can be trivially extended to the case when the upper bounds are dynamic.

We consider any *weighted directed multigraph* parameterized by a vector $h \in \mathbb{R}^n$ such that $h_i \in [0, \infty]$, and a matrix of distances $\{\rho_{i,j}\}_{i,j} \in \mathbb{R}^{n \times n}$ such that $\rho_{i,j} \in [0, \infty]$ for all $i, j \in [n]$ and $\rho_{i,i} = 0$ for all $i \in [n]$. Every worker i is connected to any other worker j with two edges $i \rightarrow j$ and $j \rightarrow i$. For this setup, it would be convenient to define the *distance of the shortest path*

5/2024

*First optimal
parallel SGD under...*

5/2023

Rennala SGD
Malenia SGD
Acc. Rennala SGD



Our Papers

Alexander Tyurin and P.R.

**Optimal time complexities of parallel stochastic optimization
methods under a fixed computation model**

NeurIPS 2023

*... computation
(and/or data) heterogeneity*

2/2024

Shadowheart SGD



Alexander Tyurin, Marta Pozzi, Ivan Ilin and P.R.

**Shadowheart SGD: Distributed asynchronous SGD with optimal
time complexity under arbitrary computation and
communication heterogeneity**

arXiv:2402.04785, 2024

*... communication
(and computation) heterogeneity*

[Rennala SGD as a special case]

5/2024

Freya PAGE
Freya SGD



Alexander Tyurin, Kaja Gruntkowska, and P.R.

**Freya PAGE: First optimal time complexity for large-scale
nonconvex finite-sum optimization with heterogeneous
asynchronous computations**

arXiv:2405.1554, 2024

*... computation heterogeneity for
finite-sum problems*

in the large-scale regime: $m \geq n^2$

5/2024

Fragile SGD, Amelie SGD
+ accelerated variants



Alexander Tyurin and P.R.

**On the optimal time complexities in decentralized stochastic
asynchronous optimization**

arXiv:2405.16218, 2024

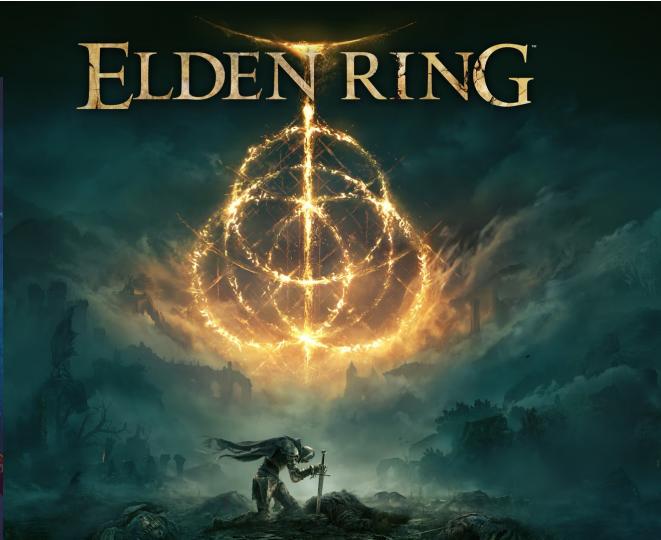
*... computation and
communication heterogeneity in
the decentralized setup*

Peter, What About the Weird Algorithm Names?



Rennala

Rennala, Queen of the Full Moon is a Legend Boss in Elden Ring. Though not a demigod, Rennala is one of the shardbearers who resides in the Academy of Raya Lucaria. Rennala is a powerful sorceress, head of the Carian Royal family, and erstwhile leader of the Academy.



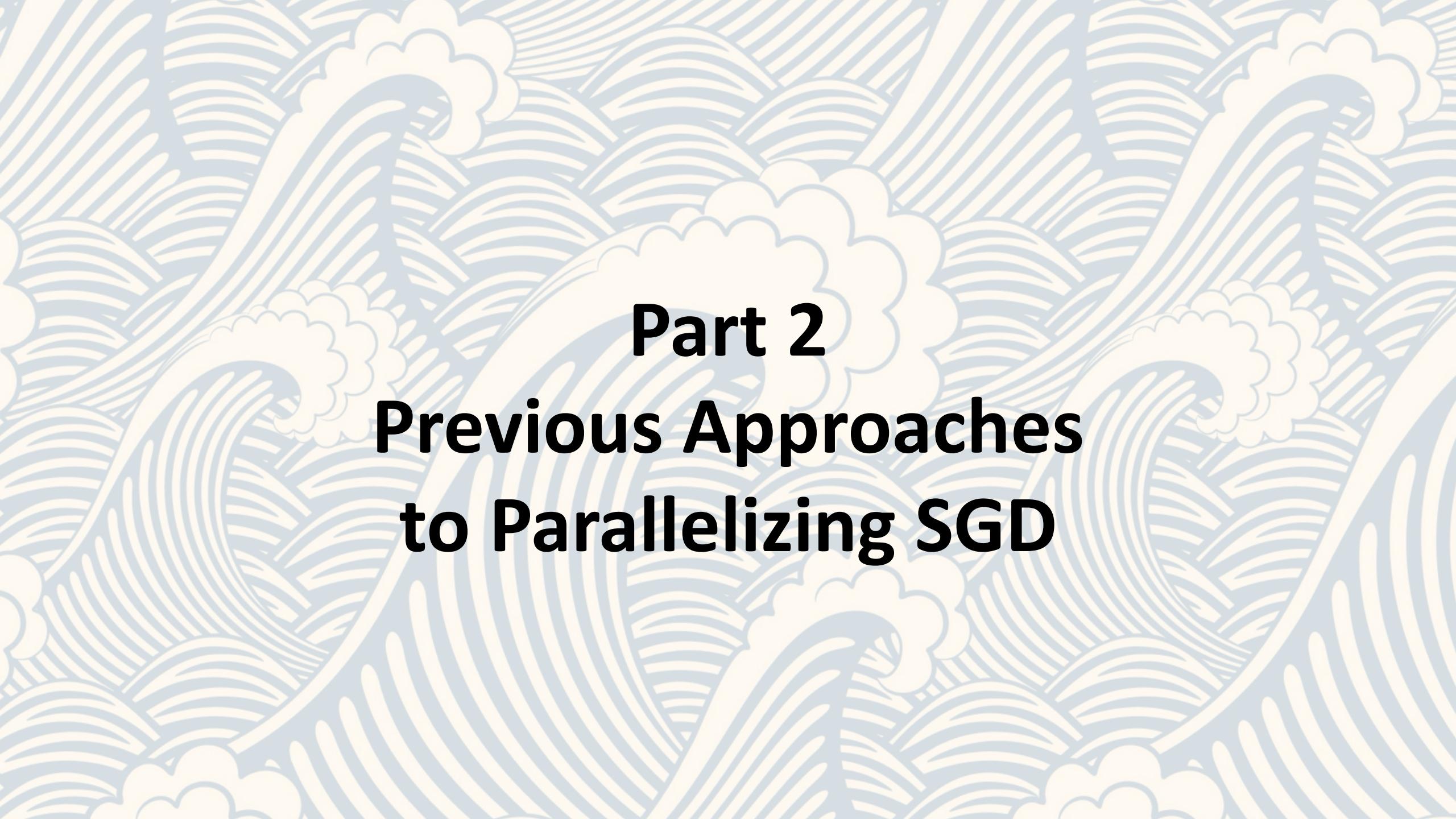
Amelie



Shadowheart

Optimal Parallel Stochastic Gradient Methods

	Data Heterogeneity (\mathcal{D}_i different)	Compute Heterogeneity (τ_i different)	Communication Heterogeneity (θ_i different)	Smooth Nonconvex	Smooth Convex	Infinite / Finite Sum?	Supports Decentralized Setup?	Optimal Time Complexity?
Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0	✓		Inf	✗	✓
Malenia SGD Tyurin & R (NeurIPS '23)	✓	✓	0	✓		Inf	✗	✓
Accelerated Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0		✓	Inf	✗	✓
Shadowheart SGD Tyurin, Pozzi, Ilin & R '24	✗	✓	✓	✓		Inf	✗	✓
Freya PAGE Tyurin, Gruntkowska & R '24	✗	✓	0	✓		Finite	✗	✓ big data regime
Freya SGD Tyurin, Gruntkowska & R '24	✗	✓	0	✓		Finite	✗	✗
Fragile SGD Tyurin & R '24	✗	✓	✓	✓		Inf	✓	nearly
Amelie SGD Tyurin & R '24	✓	✓	✓	✓		Inf	✓	✓

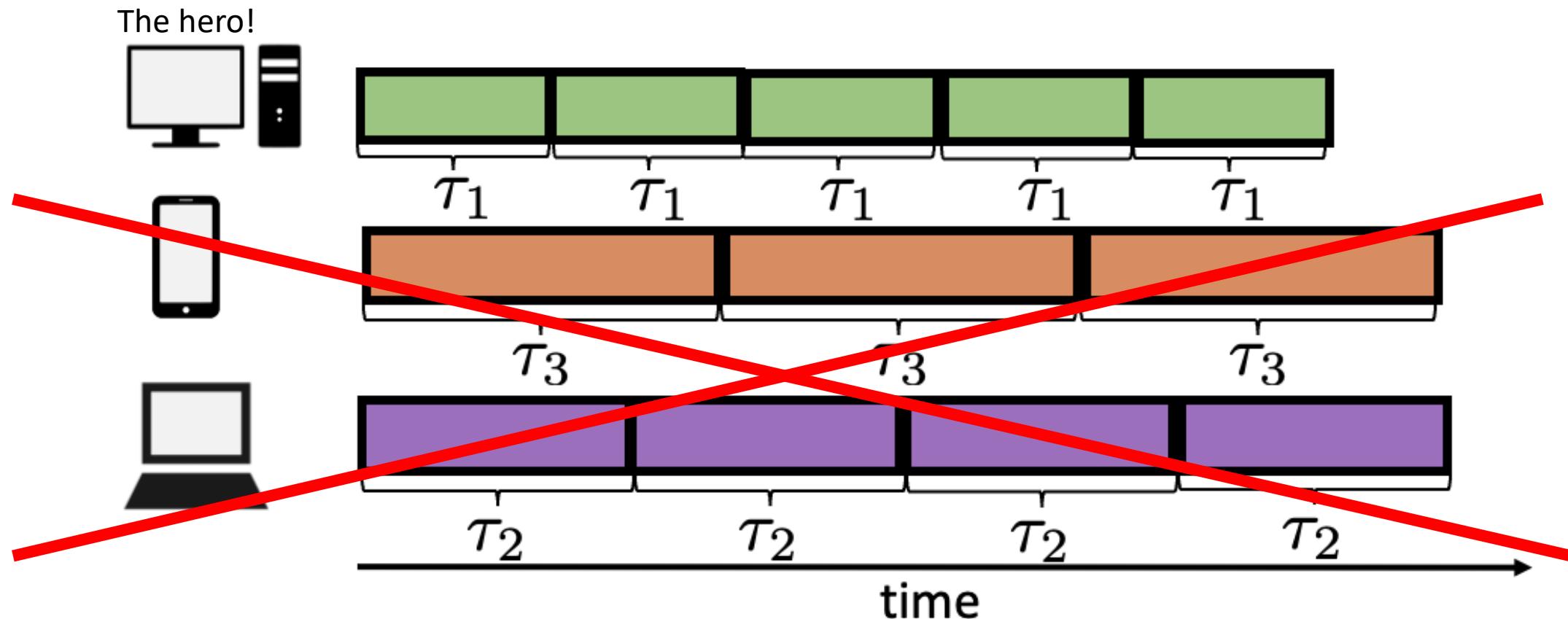


Part 2

Previous Approaches to Parallelizing SGD

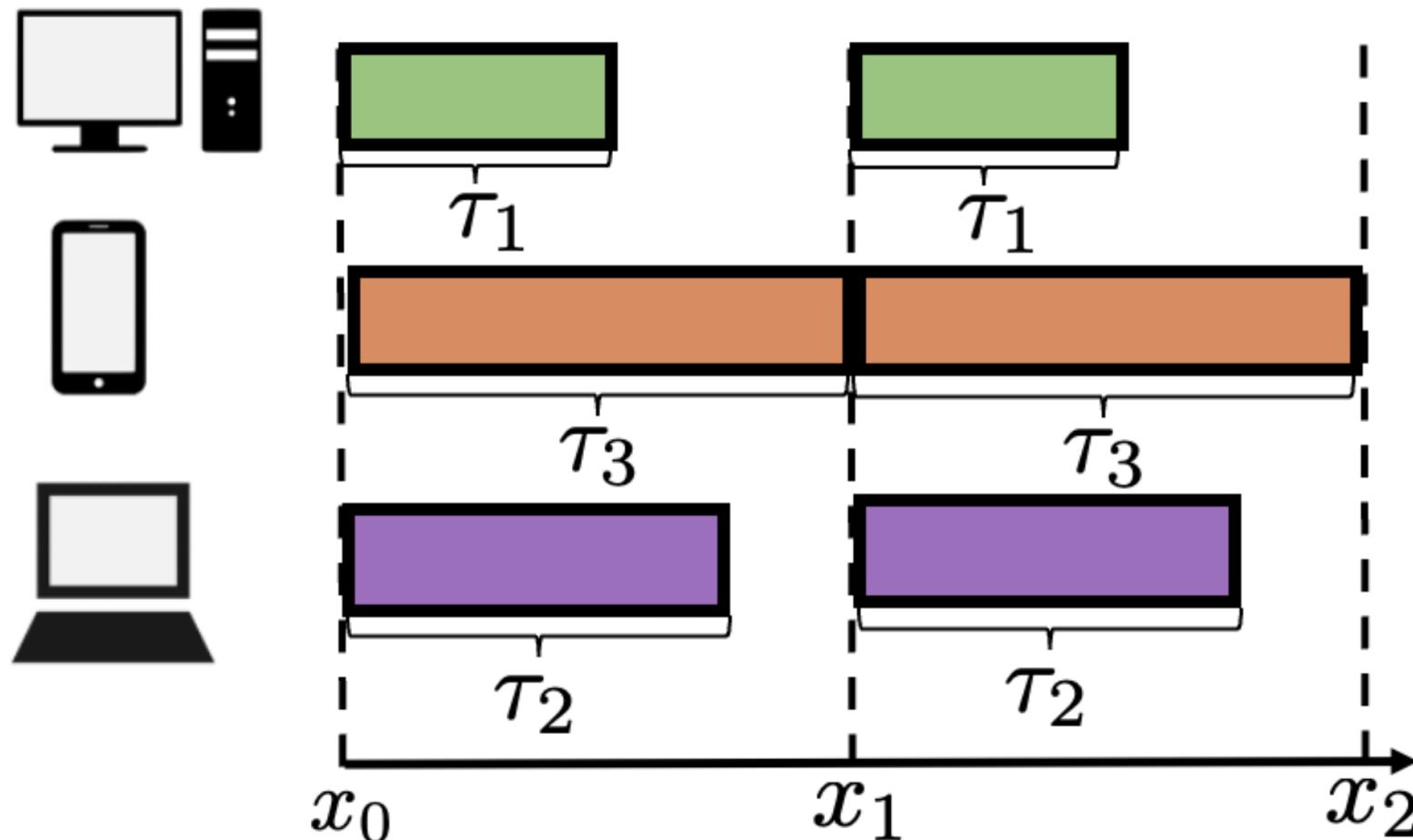
Hero SGD

Algorithmic idea: The fastest worker does it all!



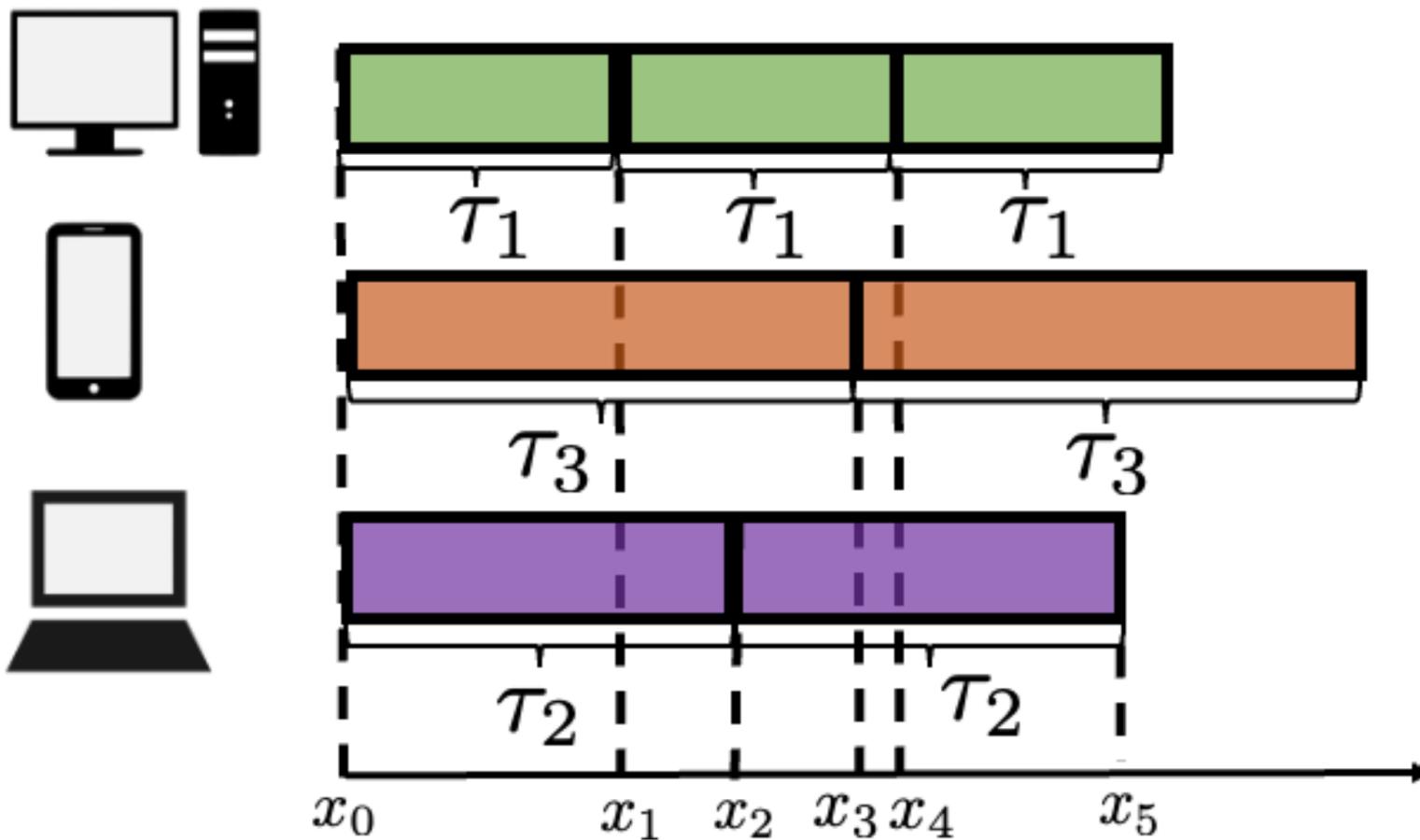
(Fair) Minibatch SGD

Algorithmic idea: Each worker does one job only!



Asynchronous SGD

Algorithmic idea: All workers are slaves and useful



published in NIPS 2011

HOGWILD!: A Lock-Free Approach to Parallelizing Stochastic Gradient Descent

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Abstract

Stochastic Gradient Descent (SGD) is a popular algorithm that can achieve state-of-the-art performance on a variety of machine learning tasks. Several researchers have recently proposed schemes to parallelize SGD, but all require performance-destroying memory locking and synchronization. This work aims to show using novel theoretical analysis, algorithms, and implementation that SGD can be implemented *without any locking*. We present an update scheme called HOGWILD! which allows processors access to shared memory with the possibility of overwriting each other's work. We show that when the associated optimization problem is *sparse*, meaning most gradient updates only modify small parts of the decision variable, then HOGWILD! achieves a nearly optimal rate of convergence. We demonstrate experimentally that HOGWILD! outperforms alternative schemes that use locking by an order of magnitude.

1 Introduction

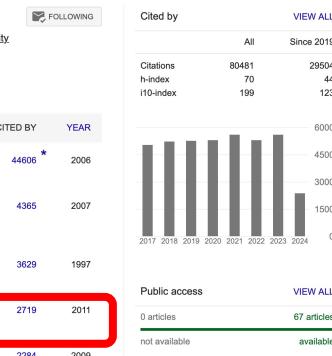
With its small memory footprint, robustness against noise, and rapid learning rates, Stochastic Gradient Descent (SGD) has proved to be well suited to data-intensive machine learning tasks [3, 5, 24]. However, SGD's scalability is limited by its inherently sequential nature; it is difficult to parallelize. Nevertheless, the recent emergence of inexpensive multicore processors and mammoth, web-scale data sets has motivated researchers to develop several clever parallelization schemes for SGD [4, 10, 12, 16, 27]. As many large data sets are currently pre-processed in a MapReduce-like parallel-processing framework, much of the recent work on parallel SGD has focused naturally on MapReduce implementations. MapReduce is a powerful tool developed at Google for extracting information from huge logs (e.g., "find all the urls from a 100TB of Web data") that was designed to ensure fault tolerance and to simplify the maintenance and programming of large clusters of machines [9]. But MapReduce is not ideally suited for online, numerically intensive data analysis. Iterative computation is difficult to express in MapReduce, and the overhead to ensure fault tolerance can result in dismal throughput. Indeed, even Google researchers themselves suggest that other systems, for example Dremel, are more appropriate than MapReduce for data analysis tasks [20].

For some data sets, the sheer size of the data dictates that one use a cluster of machines. However, there are a host of problems in which, after appropriate preprocessing, the data necessary for statistical analysis may consist of a few terabytes or less. For such problems, one can use a single inexpensive work station as opposed to a hundred thousand dollar cluster. Multicore systems have significant performance advantages, including (1) low latency and high throughput shared main memory (a processor in such a system can write and read the shared physical memory at over 12GB/s with latency in the tens of nanoseconds); and (2) high bandwidth off multiple disks (thousand-dollar RAID



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Optimization

TITLE	CITED BY	YEAR
Numerical Optimization (2nd edition) J Nocedal, SJ Wright Springer	44606 *	2006
Gradient projection for sparse reconstruction: Application to compressed sensing and other inverse problems MAT Figueiredo, RD Nowak, SJ Wright IEEE Journal of selected topics in signal processing 1 (4), 586-597	4365	2007
Primal-dual interior-point methods SJ Wright Society for Industrial and Applied Mathematics	3629	1997
Hogwild: A lock-free approach to parallelizing stochastic gradient descent B Recht, C Re, S Wright, F Niu Advances in neural information processing systems, 693-701	2719	2011
Sparse reconstruction by separable approximation SJ Wright, RD Nowak, MAT Figueiredo IEEE Transactions on signal processing 57 (7), 2479-2493	2284	2009



NeurIPS 2020 Test of Time Award

Hogwild: A lock-free approach to parallelizing stochastic gradient descent

Authors Benjamin Recht, Christopher Re, Stephen Wright, Feng Niu

Publication date 2011

Conference Advances in Neural Information Processing Systems

Pages 693-701

Description Stochastic Gradient Descent (SGD) is a popular algorithm that can achieve state-of-the-art performance on a variety of machine learning tasks. Several researchers have recently proposed schemes to parallelize SGD, but all require performance-destroying memory locking and synchronization. This work aims to show using novel theoretical analysis, algorithms, and implementation that SGD can be implemented without any locking. We present an update scheme called Hogwild which allows processors access to shared memory with the possibility of overwriting each other's work. We show that when the associated optimization problem is sparse, meaning most gradient updates only modify small parts of the decision variable, then Hogwild achieves a nearly optimal rate of convergence. We demonstrate experimentally that Hogwild outperforms alternative schemes that use locking by an order of magnitude.

Total citations Cited by 2719

Scholar articles **Hogwild: A lock-free approach to parallelizing stochastic gradient descent**
B Recht, C Re, S Wright, F Niu - Advances in neural information processing systems, 2011
Cited by 2718 Related articles All 35 versions

Hogwild: A lock-free approach to parallelizing stochastic gradient descent *
RB NiuF - ... Systems. Granada, Spain, 2011
Cited by 2 Related articles

Our Inspiration: Two Beautiful Papers

Asynchronous SGD Beats Minibatch SGD Under Arbitrary Delays

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Abstract

The existing analysis of asynchronous stochastic gradient descent (SGD) degrades dramatically when any delay is large, giving the impression that performance depends primarily on the delay. On the contrary, we prove much better guarantees for the same asynchronous SGD algorithm regardless of the delays in the gradients, depending instead just on the number of parallel devices used to implement the algorithm. Our guarantees are strictly better than the existing analyses, and we also argue that asynchronous SGD outperforms synchronous minibatch SGD in the settings we consider. For our analysis, we introduce a novel recursion based on “virtual iterates” and delay-adaptive stepsizes, which allow us to derive state-of-the-art guarantees for both convex and non-convex objectives.

1 Introduction

We consider solving stochastic optimization problems of the form

$$\min_{\mathbf{x} \in \mathbb{R}^d} \{F(\mathbf{x}) := \mathbb{E}_{\xi \sim \mathcal{D}} f(\mathbf{x}; \xi)\}, \quad (1)$$

which includes machine learning (ML) training objectives, where $f(\mathbf{x}; \xi)$ represents the loss of a model parameterized by \mathbf{x} on the datum ξ . Depending on the application, \mathcal{D} could represent a finite dataset of size n or a population distribution. In recent years, such stochastic optimization problems have continued to grow rapidly in size, both in terms of the dimension d of the optimization variable—i.e., the number of model parameters in ML—and in terms of the quantity of data—i.e., the number of samples $\xi_1, \dots, \xi_n \sim \mathcal{D}$ being used. With d and n regularly reaching the tens or hundreds of billions, it is increasingly necessary to use parallel optimization algorithms to handle the large scale and to benefit from data stored on different machines.

There are many ways of employing parallelism to solve (1), but the most popular approaches in practice are first-order methods based on stochastic gradient descent (SGD). At each iteration, SGD employs stochastic estimates of ∇F to update the parameters as $\mathbf{x}_k = \mathbf{x}_{k-1} - \gamma_k \nabla/(x_{k-1}; \xi_{k-1})$ for an i.i.d. sample $\xi_{k-1} \sim \mathcal{D}$. Given M machines capable of computing these stochastic gradient estimates $\nabla/(x; \xi)$ in parallel, one approach to parallelizing SGD is what we call “Minibatch SGD.” This refers to a synchronous, parallel algorithm that dispatches the current parameters \mathbf{x}_{k-1} to each of the M machines, waits while they compute and communicate back their gradient estimates $\mathbf{g}_{k-1}^1, \dots, \mathbf{g}_{k-1}^M$, and then takes a minibatch SGD step $\mathbf{x}_k = \mathbf{x}_{k-1} - \gamma_k \cdot \frac{1}{M} \sum_{m=1}^M \mathbf{g}_{k-1}^m$. This is a natural idea with long history [16, 18, 55] and it is a commonly used in practice [e.g., 22]. However, since Minibatch SGD waits for all M of the machines to finish computing their gradient estimates before updating, it proceeds only at the speed of the *slowest* machine.

There are several possible sources of delays: nodes may have heterogeneous hardware with different computational throughputs [23, 25], network latency can slow the communication of gradients, and

36th Conference on Neural Information Processing Systems (NeurIPS 2022).

Sharper Convergence Guarantees for Asynchronous SGD for Distributed and Federated Learning

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Abstract

We study the asynchronous stochastic gradient descent algorithm for distributed training over n workers which have varying computation and communication frequency over time. In this algorithm, workers compute stochastic gradients in parallel at their own pace and return those to the server without any synchronization. Existing convergence rates for this algorithm for non-convex smooth objectives depend on the maximum gradient delay τ_{\max} and show that an ε -stationary point is reached after $\mathcal{O}(\sigma^2 \varepsilon^{-2} + \tau_{\max} \varepsilon^{-1})$ iterations, where σ denotes the variance of stochastic gradients.

In this work we obtain (i) a tighter convergence rate of $\mathcal{O}(\sigma^2 \varepsilon^{-2} + \sqrt{\tau_{\max} \tau_{avg} \varepsilon^{-1}})$ without any change in the algorithm, where τ_{avg} is the average delay, which can be significantly smaller than τ_{\max} . We also provide (ii) a simple delay-adaptive learning rate scheme, under which asynchronous SGD achieves a convergence rate of $\mathcal{O}(\sigma^2 \varepsilon^{-2} + \tau_{avg} \varepsilon^{-1})$, and does not require any extra hyperparameter tuning nor extra communications. Our result allows to show for the first time that asynchronous SGD is always faster than mini-batch SGD. In addition, (iii) we consider the case of heterogeneous functions motivated by federated learning applications and improve the convergence rate by proving a weaker dependence on the maximum delay compared to prior works. In particular, we show that the heterogeneity term in convergence rate is only affected by the average delay within each worker.

1 Introduction

The stochastic gradient descent (SGD) algorithm [43, 13] and its variants (momentum SGD, Adam, etc.) form the foundation of modern machine learning and frequently achieve state of the art results. With recent growth in the size of models and available training data, parallel and distributed versions of SGD are becoming increasingly important [57, 17, 16]. Without those, modern state-of-the-art language models [44], generative models [40, 41], and many others [50] would not be possible. In the distributed setting, also known as data-parallel training, optimization is distributed over many computer devices working in parallel (e.g. cores, or GPUs on a cluster) in order to speed up training. Every worker computes gradients on a subset of the training data, and the resulting gradients are aggregated (averaged) on a server.

The same type of SGD variants also form the core algorithms for federated learning applications [34, 24] where the training process is naturally distributed over many user devices, or clients, that keep their local data private, and only transfer (e.g. encrypted or differentially private) gradients to the server.

A rich literature exists on the convergence theory of above mentioned parallel SGD methods, see e.g. [17, 13] and references therein. Plain parallel SGD still faces many challenges in practice, motivat-

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Optimal Time Complexities of Parallel Stochastic Optimization Methods Under a Fixed Computation Model

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Abstract

Parallelization is a popular strategy for improving the performance of iterative algorithms. Optimal time complexities are no exception. Most efficient parallel optimization methods and tight analysis of their theoretical properties are in sequential research endeavours. While the minimax complexities are well known for sequential optimization methods, the theory of parallel optimization methods is less explored. In this paper, we propose a fixed computation model for parallel optimization framework approach. Using this protocol, we establish *minimax complexities for parallel optimization methods* that have access to an unbiased stochastic gradient oracle with bounded variance. Our results are general enough to be applied to any method by each worker requiring a fixed but worker-dependent time to calculate stochastic gradients. We prove lower bounds and develop optimal algorithms that attain them. Our results have surprising consequences for the literature of asynchronous optimization methods.

1 Introduction

We consider the nonconvex optimization problem

$$\min_{x \in Q} \left\{ f(x) := \mathbb{E}_{\xi \sim P} [f(x; \xi)] \right\}, \quad (1)$$

where $f: \mathbb{R}^n \times S \rightarrow \mathbb{R}$, $S \subseteq \mathbb{R}^d$, and ξ is a random variable with some distribution P on S . In machine learning, S could be the set of all possible data, P is the distribution of the training dataset, and $f(\cdot; \cdot)$ is the loss of a data sample ξ . In this paper we address the following natural setup:

- (i) n workers are available to work in parallel;
- (ii) the i^{th} worker requires τ_i seconds¹ to calculate a stochastic gradient of f .

The function f is L -smooth and lower bounded (see Assumptions 7.1–7.2), and stochastic gradients are unbiased and σ^2 -variance-bounded (see Assumption 7.3).

1.1 Classical Theory

In the notorious setting, gradient descent (GD) is an optimal method with respect to the number of gradient ∇f calls (Lin, 2020; Nesterov, 2018; Caron et al., 2020) for finding an approximately stationary point of f . Obviously, a key issue with GD is that it requires access to the exact gradients

¹Or any other unit of time.

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Part 3

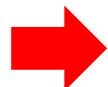
Rennala SGD



Alexander Tyurin and P.R.
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NeurIPS 2023

Setup

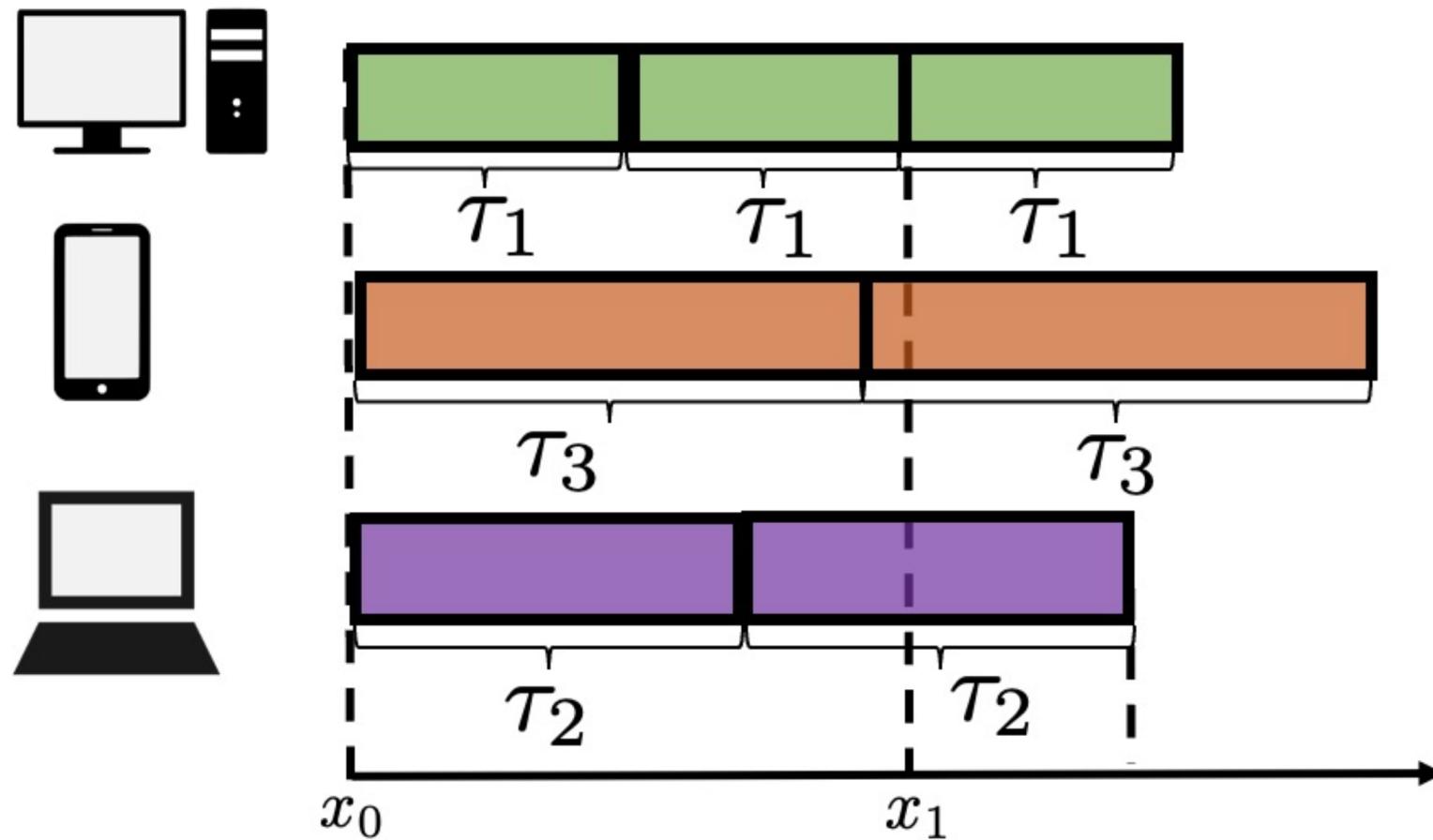
Optimal Parallel Stochastic Gradient Methods



	Data Heterogeneity (\mathcal{D}_i different)	Compute Heterogeneity (τ_i different)	Communication Heterogeneity (θ_i different)	Smooth Nonconvex	Smooth Convex	Infinite / Finite Sum?	Supports Decentralized Setup?	Optimal Time Complexity?
Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0	✓		Inf	✗	✓
Malenia SGD Tyurin & R (NeurIPS '23)	✓	✓	0	✓		Inf	✗	✓
Accelerated Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0		✓	Inf	✗	✓
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Fragile SGD Tyurin & R '24	✗	✓	✓	✓		Inf	✓	nearly
Amelie SGD Tyurin & R '24	✓	✓	✓	✓		Inf	✓	✓

Rennala SGD

Algorithmic idea: Minibatch SGD with asynchronous minibatch collection



Upper Bound

Theorem (informal)

Assume data homogeneity and zero communication times.

Then Rennala SGD solves the problem in

Number of parallel machines

$$96 \times \min_{m \in \{1, \dots, n\}} \left(\frac{1}{m} \sum_{i=1}^m \frac{1}{\tau_i} \right)^{-1} \left(\frac{L\Delta}{\varepsilon} + \frac{L\Delta\sigma^2}{\varepsilon^2 m} \right)$$

seconds.

Compute times

$$0 < \tau_1 \leq \tau_2 \leq \dots \leq \tau_n$$

Algorithm outputs \hat{x} such that $\mathbb{E} [\|\nabla f(\hat{x})\|^2] \leq \varepsilon$

Gradient of f is L -Lipschitz

$$\Delta := f(x^0) - \inf f$$

$$\sup_{x \in \mathbb{R}^d} \mathbb{E}_{\xi \sim \mathcal{D}} [\|\nabla f(x, \xi) - \nabla f(x)\|^2] \leq \sigma^2$$

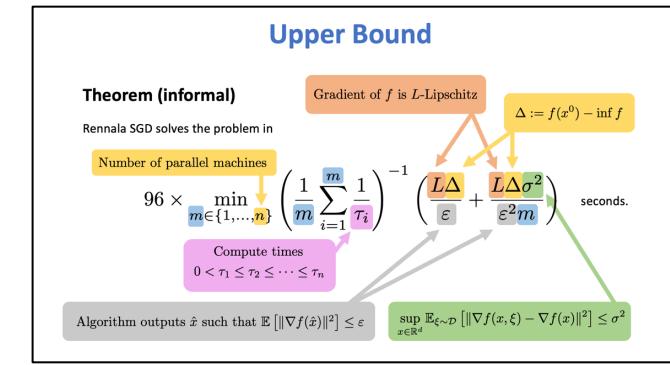
Matching Lower Bound

Theorem (informal)

It is not possible to design a method that will find a solution faster than in

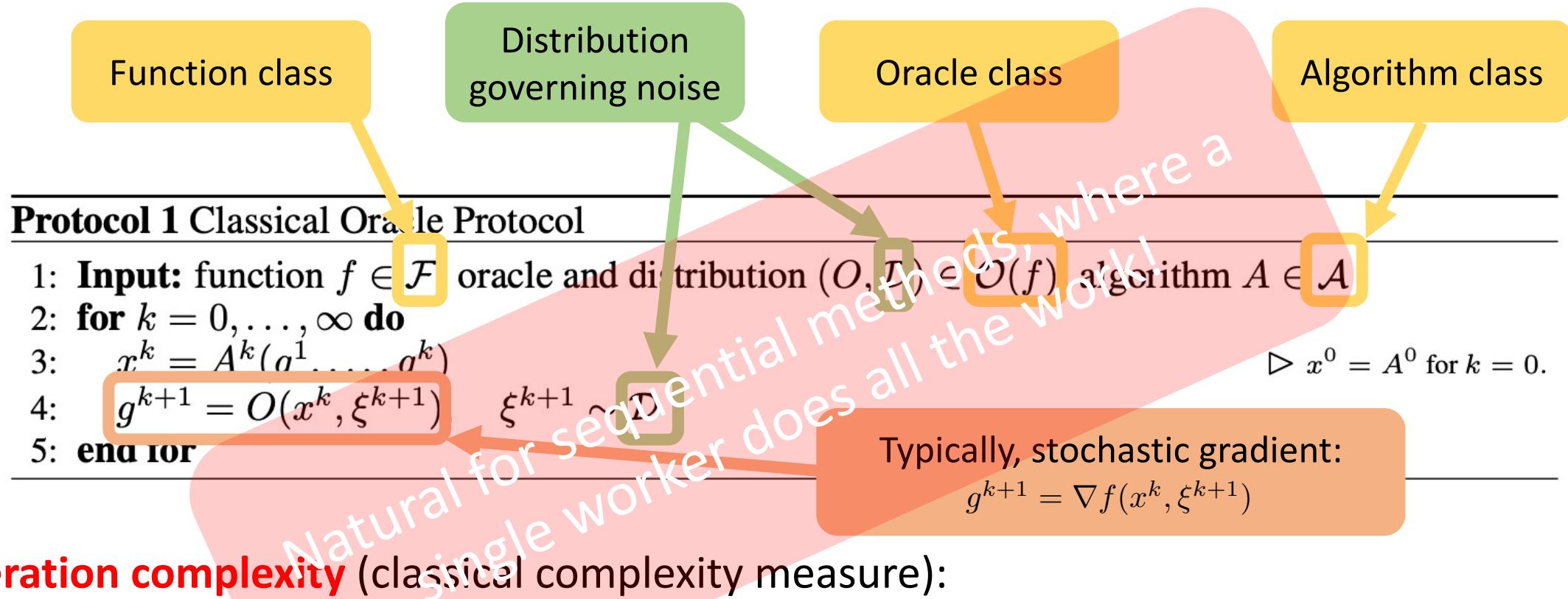
$$\Omega \left(\min_{m \in \{1, \dots, n\}} \left(\frac{1}{m} \sum_{i=1}^m \frac{1}{\tau_i} \right)^{-1} \left(\frac{L\Delta}{\varepsilon} + \frac{L\Delta\sigma^2}{\varepsilon^2 m} \right) \right)$$

seconds.



Rennala SGD = first optimal parallel SGD

Classical Oracle: Keeps Track of # Iterations



Iteration complexity (classical complexity measure):

$$m_{\text{oracle}}(\mathcal{A}, \mathcal{F}) := \inf_{A \in \mathcal{A}} \sup_{f \in \mathcal{F}} \sup_{(O, \mathcal{D}) \in \mathcal{O}(f)} \inf \{k \in \mathbb{N} \mid \mathbb{E} [\|\nabla f(x^k)\|^2] \leq \varepsilon\}$$

[Nemirovsky and Yudin, 1983] [Nesterov, 2018]
[Carmon et al, 2020] [Arjevani et al, 2022]

New Oracle: Keeps Track of Time

Protocol 2 Time Oracle Protocol

```
1: Input: functions  $f \in \mathcal{F}$ , oracle and distribution  $(O, \mathcal{D}) \in \mathcal{O}(f)$ , algorithm  $A \in \mathcal{A}$ 
2:  $s^0 = 0$ 
3: for  $k = 0, \dots, \infty$  do
4:    $(t^{k+1}, x^k) = A^k(g^1, \dots, g^k),$ 
5:    $(s^{k+1}, g^{k+1}) = O(t^{k+1}, x^k, s^k, \xi^{k+1}), \quad \xi^{k+1} \sim \mathcal{D}$ 
6: end for
```

$\triangleright t^{k+1} \geq t^k$

Iteration complexity (classical complexity measure):

$$m_{\text{oracle}}(\mathcal{A}, \mathcal{F}) := \inf_{A \in \mathcal{A}} \sup_{f \in \mathcal{F}} \sup_{(O, \mathcal{D}) \in \mathcal{O}(f)} \inf \left\{ k \in \mathbb{N} \mid \mathbb{E} [\|\nabla f(x^k)\|^2] \leq \varepsilon \right\}$$

Time complexity (new complexity measure):

$$m_{\text{time}}(\mathcal{A}, \mathcal{F}) := \inf_{A \in \mathcal{A}} \sup_{f \in \mathcal{F}} \sup_{(O, \mathcal{D}) \in \mathcal{O}(f)} \inf \left\{ t \geq 0 \mid \mathbb{E} \left[\inf_{k \in S_t} \|\nabla f(x^k)\|^2 \right] \leq \varepsilon \right\}$$

$$S_t := \{k \in \mathbb{N} \cup \{0\} \mid t^k \leq t\}$$

Data Homogeneous Regime

Method	Time Complexity
Minibatch SGD	$\tau_n \left(\frac{L\Delta}{\varepsilon} + \frac{\sigma^2 L\Delta}{n\varepsilon^2} \right)$
Asynchronous SGD (Cohen et al., 2021) (Koloskova et al., 2022) (Mishchenko et al., 2022)	$\left(\frac{1}{n} \sum_{i=1}^n \frac{1}{\tau_i} \right)^{-1} \left(\frac{L\Delta}{\varepsilon} + \frac{\sigma^2 L\Delta}{n\varepsilon^2} \right)$
Rennala SGD (Theorem 7.5)	$\min_{m \in [n]} \left[\left(\frac{1}{m} \sum_{i=1}^m \frac{1}{\tau_i} \right)^{-1} \left(\frac{L\Delta}{\varepsilon} + \frac{\sigma^2 L\Delta}{m\varepsilon^2} \right) \right]$
Lower Bound (Theorem 6.4)	$\min_{m \in [n]} \left[\left(\frac{1}{m} \sum_{i=1}^m \frac{1}{\tau_i} \right)^{-1} \left(\frac{L\Delta}{\varepsilon} + \frac{\sigma^2 L\Delta}{m\varepsilon^2} \right) \right]$

Experimental Results (Sample)

$$\tau_i = \sqrt{i} \text{ seconds}$$

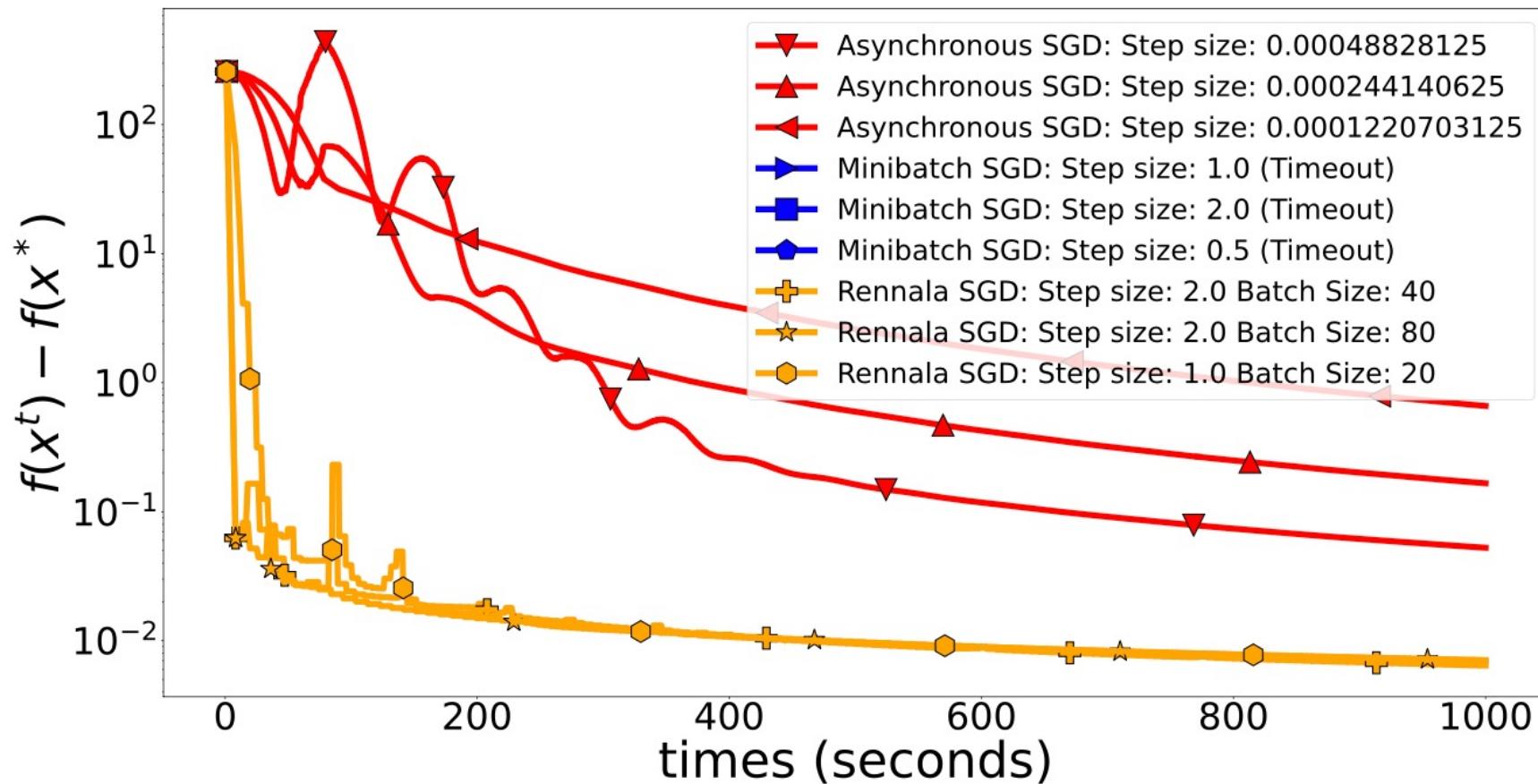


Figure 3: # of workers $n = 10000$.

Optimal Time Complexities of Parallel Stochastic Optimization Methods Under a Fixed Computation Model

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Part 4

Two Extensions

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Extension 1

Handling Data Heterogeneity

(Malenia SGD)

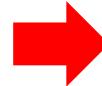
Malenia SGD: Setup

$$\min_{x \in \mathbb{R}^d} \left\{ f(x) = \frac{1}{n} \sum_{i=1}^n f_i(x) \right\}$$

$$f_i(x) := \mathbb{E}_{\xi \sim \mathcal{D}_i} [f_i(x, \xi)]$$

Optimal Parallel Stochastic Gradient Methods

	Data Heterogeneity (\mathcal{D}_i different)	Compute Heterogeneity (τ_i different)	Communication Heterogeneity (θ_i different)	Smooth Nonconvex	Smooth Convex	Infinite / Finite Sum?	Supports Decentralized Setup?	Optimal Time Complexity?
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Malenia SGD Tyurin & R (NeurIPS '23)	✓	✓	0	✓		Inf	✗	✓
Accelerated Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0		✓	Inf	✗	✓
Shadowheart SGD Tyurin, Pozzi, Ilin & R '24	✗	✓	✓	✓		Inf	✗	✓
Freya PAGE Tyurin, Gruntkowska & R '24	✗	✓	0	✓		Finite	✗	✓ big data regime
Freya SGD Tyurin, Gruntkowska & R '24	✗	✓	0	✓		Finite	✗	✗
Fragile SGD Tyurin & R '24	✗	✓	✓	✓		Inf	✓	nearly
Amelie SGD Tyurin & R '24	✓	✓	✓	✓		Inf	✓	✓



The distributions $\mathcal{D}_1, \dots, \mathcal{D}_n$ are allowed to be different

Malenia SGD

Method 6 Malenia SGD

```

1: Input: starting point  $x^0$ , stepsize  $\gamma$ , parameter  $S$ 
2: Run Method 7 in all workers
3: for  $k = 0, 1, \dots, K - 1$  do
4:   Init  $g_i^k = 0$  and  $B_i = 0$ 
5:   while  $\left( \frac{1}{n} \sum_{i=1}^n \frac{1}{B_i} \right)^{-1} < \frac{S}{n}$  do
6:     Wait for the next worker
7:     Receive gradient, iteration index, worker's index  $(g, k', i)$ 
8:     if  $k' = k$  then
9:        $g_i^k = g_i^k + g$ 
10:       $B_i = B_i + 1$ 
11:    end if
12:    Send  $(x^k, k)$  to the worker
13:  end while
14:   $g^k = \frac{1}{n} \sum_{i=1}^n \frac{1}{B_i} g_i^k$ 
15:   $x^{k+1} = x^k - \gamma g^k$ 
16: end for

```

Minibatch size

$$S = \max \left\{ \left\lceil \frac{\sigma^2}{\varepsilon} \right\rceil, n \right\}$$

Method 7 Worker's Infinite Loop

```

1: Init  $g = 0$ ,  $k' = -1$ , and worker's index  $i$ 
2: while True do
3:   Send  $(g, k', i)$  to the server
4:   Receive  $(x^k, k)$  from the server
5:    $k' = k$ 
6:    $g = \hat{\nabla} f_i(x^k; \xi)$ ,  $\xi \sim \mathcal{D}$ 
7: end while

```

(Nonconvex) Data Heterogeneous Regime

Method	Time Complexity
Minibatch SGD	$\tau_n \left(\frac{L\Delta}{\varepsilon} + \frac{\sigma^2 L\Delta}{n\varepsilon^2} \right)$
Malenia SGD (Theorem A.4)	$\tau_n \frac{L\Delta}{\varepsilon} + \left(\frac{1}{n} \sum_{i=1}^n \tau_i \right) \frac{\sigma^2 L\Delta}{n\varepsilon^2}$
Lower Bound (Theorem A.2)	$\tau_n \frac{L\Delta}{\varepsilon} + \left(\frac{1}{n} \sum_{i=1}^n \tau_i \right) \frac{\sigma^2 L\Delta}{n\varepsilon^2}$

Extension 2

Handling the Convex Regime

(Accelerated Rennala SGD)

Accelerated Rennala SGD: Setup

Optimal Parallel Stochastic Gradient Methods

	Data Heterogeneity (\mathcal{D}_i different)	Compute Heterogeneity (τ_i different)	Communication Heterogeneity (θ_i different)	Smooth Nonconvex	Smooth Convex	Infinite / Finite Sum?	Supports Decentralized Setup?	Optimal Time Complexity?
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Accelerated Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0		✓	Inf	✗	✓
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Amelie SGD Tyurin & R '24	✓	✓	✓	✓		Inf	✓	✓



Convex (Data Homogeneous) Regime

Method	Time Complexity
Minibatch SGD	$\tau_n \left(\min \left\{ \frac{\sqrt{L}R}{\sqrt{\varepsilon}}, \frac{M^2 R^2}{\varepsilon^2} \right\} + \frac{\sigma^2 R^2}{n \varepsilon^2} \right)$
Asynchronous SGD (Mishchenko et al., 2022)	$\left(\frac{1}{n} \sum_{i=1}^n \frac{1}{\tau_i} \right)^{-1} \left(\frac{LR^2}{\varepsilon} + \frac{\sigma^2 R^2}{n \varepsilon^2} \right)$
(Accelerated) Rennala SGD (Theorems B.9 and B.11)	$\min_{m \in [n]} \left[\left(\frac{1}{m} \sum_{i=1}^m \frac{1}{\tau_i} \right)^{-1} \left(\min \left\{ \frac{\sqrt{L}R}{\sqrt{\varepsilon}}, \frac{M^2 R^2}{\varepsilon^2} \right\} + \frac{\sigma^2 R^2}{m \varepsilon^2} \right) \right]$
Lower Bound (Theorem B.4)	$\min_{m \in [n]} \left[\left(\frac{1}{m} \sum_{i=1}^m \frac{1}{\tau_i} \right)^{-1} \left(\min \left\{ \frac{\sqrt{L}R}{\sqrt{\varepsilon}}, \frac{M^2 R^2}{\varepsilon^2} \right\} + \frac{\sigma^2 R^2}{m \varepsilon^2} \right) \right]$
Lower Bound (Section M) (Woodworth et al., 2018)	$\tau_1 \min \left\{ \frac{\sqrt{L}R}{\sqrt{\varepsilon}}, \frac{M^2 R^2}{\varepsilon^2} \right\} + \left(\frac{1}{n} \sum_{i=1}^n \frac{1}{\tau_i} \right)^{-1} \frac{\sigma^2 R^2}{n \varepsilon^2}$

∇f is L -Lipschitz, f is M -Lipschitz, and $\|x^0 - x^\star\| \leq R$



The End

Further Extensions



Shadowheart



Shadowheart SGD

Optimal Parallel SGD under Compute Heterogeneity & Communication Heterogeneity

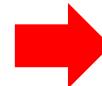
Shadowheart SGD: Setup

$$\min_{x \in \mathbb{R}^d} \left\{ f(x) = \frac{1}{n} \sum_{i=1}^n f_i(x) \right\}$$

$$f_i(x) := \mathbb{E}_{\xi \sim \mathcal{D}_i} [f_i(x, \xi)]$$

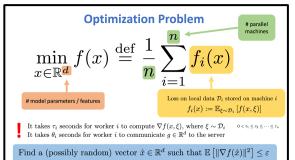
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Fragile SGD Tyurin & R '24	✗	✓	✓	✓		Inf	✓	nearly
Amelie SGD Tyurin & R '24	✓	✓	✓	✓		Inf	✓	✓



$\mathcal{D}_1 = \dots = \mathcal{D}_n$

Communication costs $\theta_1, \dots, \theta_n$ are nonzero (and possibly different)



Shadowheart SGD

Unbiased compressor:

$$\mathbb{E} [\mathcal{C}_{ij}(g)] = g \quad \& \quad \mathbb{E} [\|\mathcal{C}_{ij}(g) - g\|^2] \leq \omega \|g\|^2 \quad \forall g \in \mathbb{R}^d$$

Aggregation weight associated with worker i

$$w_i = \left(\omega b_i + \omega \frac{\sigma^2}{\varepsilon} + m_i \frac{\sigma^2}{\varepsilon} \right)^{-1}$$

$$x^{k+1} = x^k - \gamma \cdot \frac{\sum_{i=1}^n w_i}{\sum_{i=1}^n w_i m_i b_i} \mathcal{C}_{ij} \left(\sum_{l=1}^n \nabla f(x^k, \xi_{il}^k) \right)$$

$$\gamma = \frac{1}{2L}$$

Equilibrium time: $t^* : \left(\omega, \frac{\sigma^2}{\varepsilon}, (\tau_i)_{i=1}^n, (\theta_i)_{i=1}^n \right) \mapsto \mathbb{R}_+$

of compressed batches sent by worker i to the server

$$m_i = \left\lfloor \frac{t^*}{\theta_i} \right\rfloor$$

Batch size to compress by worker i

$$b_i = \left\lfloor \frac{t^*}{\tau_i} \right\rfloor$$

Table 1: Time Complexities of Centralized Distributed Algorithms. Assume that it takes at most h_i seconds to worker i to calculate a stochastic gradient and $\dot{\tau}_i$ seconds to send *one coordinate/float* to server. Abbreviations: L = smoothness constant, ε = error tolerance, $\Delta = f(x^0) - f^*$, n = # of workers, d = dimension of the problem. We take the RandK compressor with $K = 1$ (Def. C.1) (as an example) in QSGD and Shadowheart SGD. Due to Property 5.2, the choice $K = 1$ is optimal for Shadowheart SGD up to a constant factor.

Method	Time Complexity	Time Complexities in Some Regimes		Numerical Comparison ^(b) $\sigma^2/\varepsilon =$		
		$\max\{h_n, \dot{\tau}_n\} \rightarrow \infty$, $\max\{h_i, \dot{\tau}_i\} < \infty \forall i < n$ (the last worker is slow)	$h_i = h, \dot{\tau}_i = \dot{\tau} \forall i \in [n]$ (equal performance)			
Minibatch SGD (see (3))	$\max_{i \in [n]} \max\{h_i, d\dot{\tau}_i\} \left(\frac{L\Delta}{\varepsilon} + \frac{\sigma^2 L\Delta}{n\varepsilon^2} \right)$	∞ (non-robust)	$\max\{h, d\dot{\tau}, \frac{d\dot{\tau}\sigma^2}{n\varepsilon}, \frac{h\sigma^2}{n\varepsilon}\} \frac{L\Delta}{\varepsilon}$ (worse, e.g., when $\dot{\tau}, d$ or n large)	$\times 10^3$	$\times 10^3$	$\times 10^4$
QSGD (see (7)) (Alistarh et al., 2017) (Khaled & Richtárik, 2020)	$\max_{i \in [n]} \max\{h_i, \dot{\tau}_i\} \left(\left(\frac{d}{n} + 1 \right) \frac{L\Delta}{\varepsilon} + \frac{d\sigma^2 L\Delta}{n\varepsilon^2} \right)$	∞ (non-robust)	$\geq \frac{dh\sigma^2}{n\varepsilon} \frac{L\Delta}{\varepsilon}$ (worse, e.g., when ε small)	$\times 3$	$\times 10^2$	$\times 10^4$
Rennala SGD (Tyurin & Richtárik, 2023c), Asynchronous SGD (e.g., (Mishchenko et al., 2022))	$\geq \min_{j \in [n]} \max \left\{ h_{\bar{\pi}_j}, d\dot{\tau}_{\bar{\pi}_j}, \frac{\sigma^2}{\varepsilon} \left(\sum_{i=1}^j \frac{1}{h_{\bar{\pi}_i}} \right)^{-1} \right\} \frac{L\Delta}{\varepsilon}$ ^(a)	$< \infty$ (robust)	$\geq \max \left\{ h, d\dot{\tau}, \frac{h\sigma^2}{n\varepsilon} \right\} \frac{L\Delta}{\varepsilon}$ (worse, e.g., when $\dot{\tau}, d$ or n large)	$\times 10^2$	$\times 10$	$\times 1.5$
Shadowheart SGD (see (9) and Alg. 1) (Corollary 4.4)	$t^*(d - 1, \sigma^2/\varepsilon, [h_i, \dot{\tau}_i]_1^n) \frac{L\Delta}{\varepsilon}$ ^(c)	$< \infty$ (robust)	$\max \left\{ h, \dot{\tau}, \frac{d\dot{\tau}}{n}, \sqrt{\frac{d\dot{\tau}h\sigma^2}{n\varepsilon}}, \frac{h\sigma^2}{n\varepsilon} \right\} \frac{L\Delta}{\varepsilon}$	$\times 1$	$\times 1$	$\times 1$

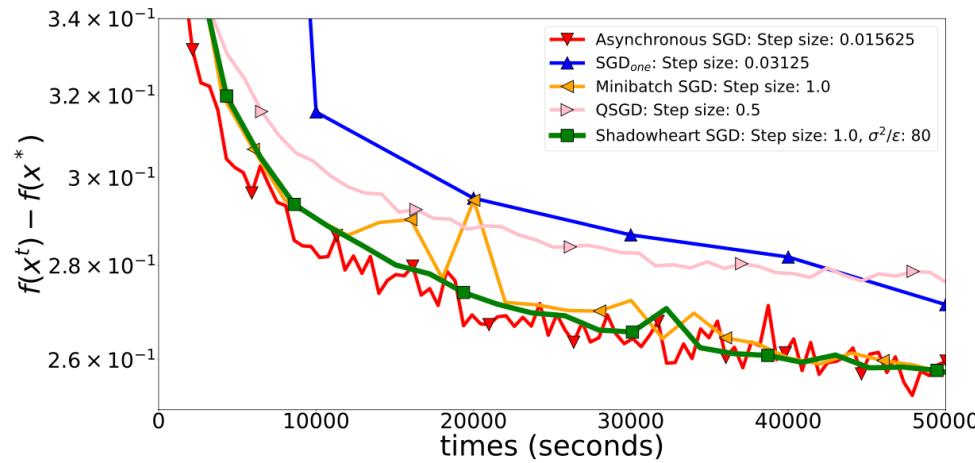
The time complexity of Shadowheart SGD is not worse than the time complexity of the competing centralized methods (see Sec. 6), and is *strictly* better in many regimes.

We show that (12) is the *optimal time complexity* in the family of centralized methods with compression (see Sec. 7).

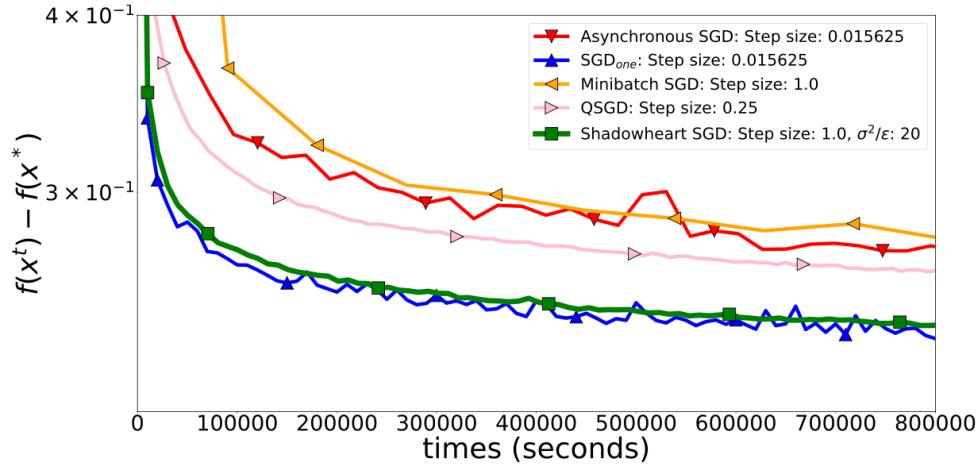
^(a) Upper bound time complexities are not derived for Rennala SGD and Asynchronous SGD. However, we can derive the lower bound using Theorem N.5 with $\omega = 0$. One should take $d\dot{\tau}_i$ instead of τ_i when apply Theorem N.5 because these methods send d coordinates. $\bar{\pi}$ is a permutation that sorts $\max\{h_i, d\dot{\tau}_i\} : \max\{h_{\bar{\pi}_1}, d\dot{\tau}_{\bar{\pi}_1}\} \leq \dots \leq \max\{h_{\bar{\pi}_n}, d\dot{\tau}_{\bar{\pi}_n}\}$

^(b) We numerically compute time complexities for $d = 10^6$, $n = 10^3$, $h_i \sim U(0.1, 1)$, $\dot{\tau}_i \sim U(0.1, 1)$ (uniform i.i.d.), and three noise regimes $\sigma^2/\varepsilon \in \{1, 10^3, 10^6\}$. We report the factors by which the time complexities of the competing methods are worse compared to the time complexity of our method Shadowheart SGD. So, for example, Minibatch SGD, QSGD and Asynchronous SGD can be worse by the factors $\times 10^4$, $\times 10^4$, and $\times 10^2$, respectively.

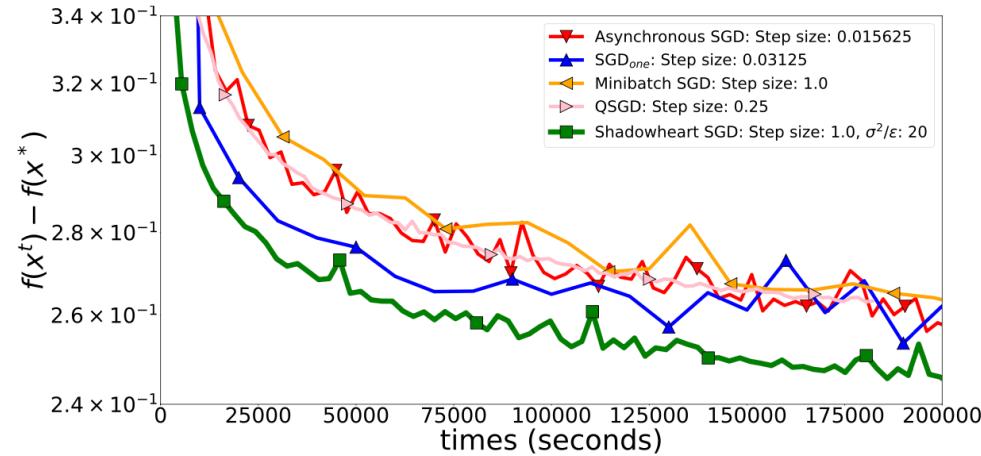
^(c) The mapping t^* is defined in Def. 4.2.



Fast communication: $\dot{\theta}_i = \frac{\sqrt{i}}{d}$



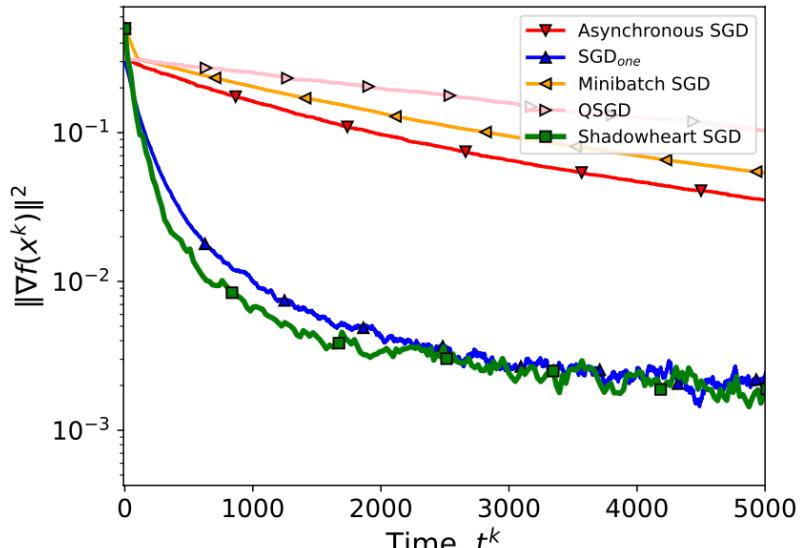
Slow communication: $\dot{\theta}_i = \frac{\sqrt{i}}{d^{1/2}}$



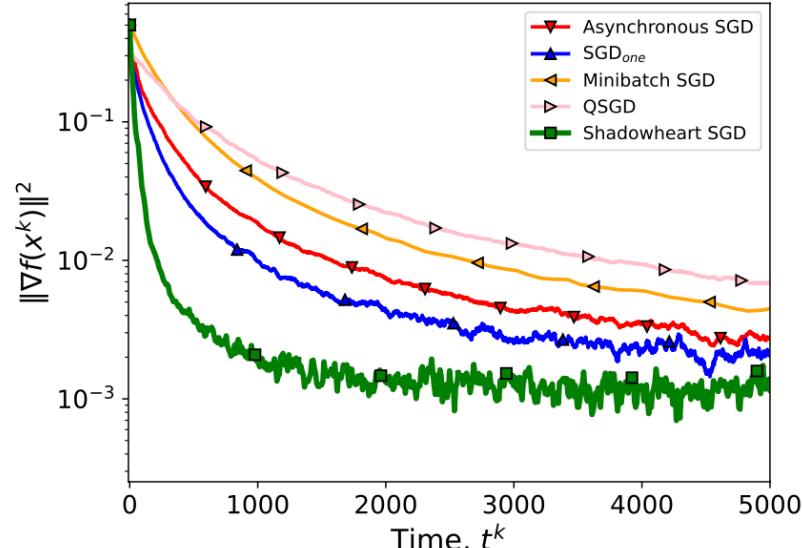
Medium-speed communication: $\dot{\theta}_i = \frac{\sqrt{i}}{d^{3/4}}$

Computation times: $\tau_i = \sqrt{i}$ for all machines $i = 1, \dots, n$

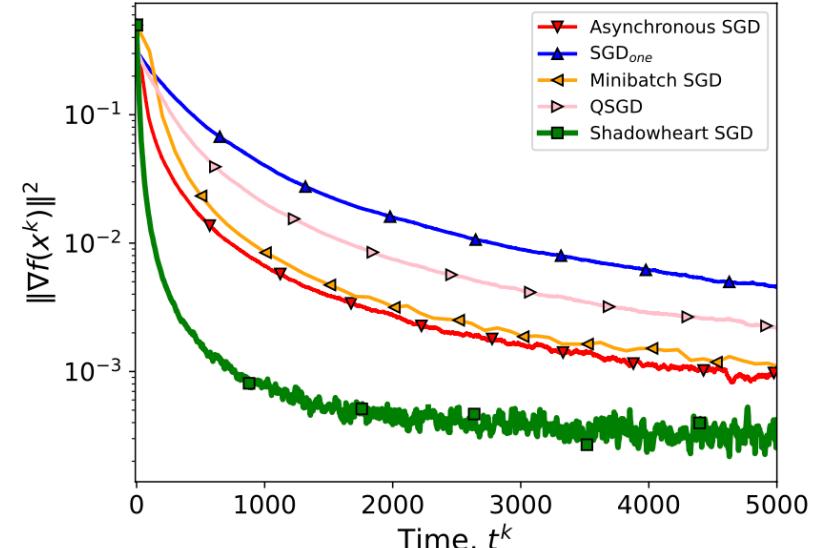
Shadowheart SGD: Adding More Workers...



(a) $n = 10$



(b) $n = 10^2$



(c) $n = 10^3$

$\tau_i^k, \dot{\theta}_i^k \sim \text{Uniform}(0.1, 1)$ for all $i \in \{1, \dots, n\}$ and $k \geq 0$

Freya PAGE: First Optimal Parallel SGD for Large-Scale Nonconvex Finite-Sum Optimization with Heterogeneous Asynchronous Computations

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Abstract

In practical distributed systems, workers are typically not homogeneous, and due to differences in hardware conditions and network conditions, can have highly varying computation times. This motivates the need for parallel SGD algorithms that can handle heterogeneous workers. In this paper, we propose Freya PAGE, designed to handle arbitrary heterogeneous asynchronous computations. Freya PAGE offers significantly improved time complexity guarantees compared to prior work, while maintaining the same asymptotic convergence rate as PAGE, while requiring weaker assumptions. The algorithm relies on novel gossip mechanisms that allow workers to exchange information about their local gradients on their own, and may be used in the design of Stochastic optimization methods for large-scale finite-sum problems. We also show that Freya PAGE generalizes to the asynchronous setup, providing a fundamental time complexity guarantee for the first time. Finally, we show that Freya PAGE is optimal for the data-sparse regime, i.e., when $\gamma \in (0, 1)$, where γ is a weight, and n is # of data samples.

1 Introduction

In real-world distributed systems (and for large-scale machine learning tasks), it is common to encounter device heterogeneity and variations in processing times among different computational nodes, communication links, and network conditions, as well as different hardware conditions, and other factors, resulting in different computational capabilities and speeds across different workers. This heterogeneity is a challenge for distributed optimization, as some workers may complete their local optimization task faster than others, while others experience delayed cost due to latency in the network (shifting the computation burden to other workers). This heterogeneity is the central challenge in solving finite-sum nonconvex optimization problems of the size encountered in the context of machine learning.

$$\max_{\theta} \left\{ f(\theta) + \frac{\lambda}{n} \| \theta \|_2^2 \right\}, \quad (1)$$

where $f_i: \mathbb{R}^d \rightarrow \mathbb{R}$ can be viewed as the loss of a machine learning model, i.e., $f = \sum_i f_i$ example is a linear regression problem, $\theta \in \mathbb{R}^d$ is the parameter vector, $\lambda > 0$ is a regularization parameter, and n is the number of training samples. We focus on the heterogeneous distributed setup,

- each worker has access to stochastic gradients $\nabla_i f_i(\cdot)$;
- worker i calculates $\nabla_i f_i(\cdot)$ in less or equal to $\tau_i \in [0, \infty)$ seconds for all $i \in [n], j \in [m]$.

Project: Under review

Shadowheart



Freya PAGE

Optimal Parallel SGD for Large-Scale Finite-Sum Problems

Freya PAGE: Setup

$$\min_{x \in \mathbb{R}^d} \left\{ f(x) = \frac{1}{n} \sum_{i=1}^n f_i(x) \right\}$$

$$f_i(x) := \mathbb{E}_{\xi \sim \mathcal{D}_i} [f_i(x, \xi)]$$

Optimal Parallel Stochastic Gradient Methods

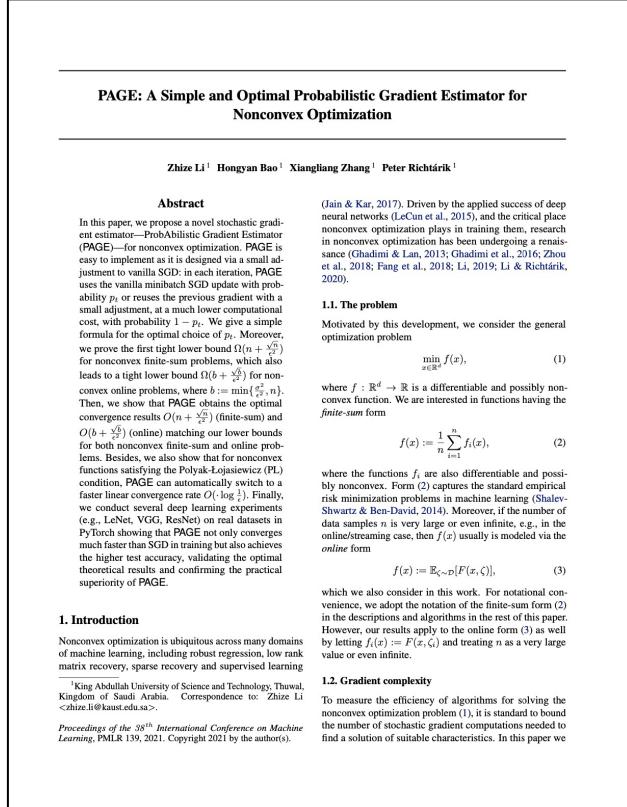
	Data Heterogeneity (\mathcal{D}_i different)	Compute Heterogeneity (τ_i different)	Communication Heterogeneity (θ_i different)	Smooth Nonconvex	Smooth Convex	Infinite / Finite Sum?	Supports Decentralized Setup?	Optimal Time Complexity?
Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0	✓		Inf	✗	✓
Malenia SGD Tyurin & R (NeurIPS '23)	✓	✓	0	✓		Inf	✗	✓
Accelerated Rennala SGD Tyurin & R (NeurIPS '23)	✗	✓	0		✓	Inf	✗	✓
Shadowheart SGD Tyurin, Pozzi, Ilin & R '24	✗	✓	✓	✓		Inf	✗	✓
Freya PAGE Tyurin, Gruntkowska & R '24	✗	✓	0	✓		Finite	✗	✓ big data regime
Freya SGD Tyurin, Gruntkowska & R '24	✗	✓	0	✓		Finite	✗	✗
Fragile SGD Tyurin & R '24	✗	✓	✓	✓		Inf	✓	nearly
Amelie SGD Tyurin & R '24	✓	✓	✓	✓		Inf	✓	✓



$\mathcal{D}_1 = \dots = \mathcal{D}_n$

$\mathcal{D}_i =$ uniform distribution over m outcomes

PAGE: Optimal Serial SGD for Finite-Sum Nonconvex Optimization



$$\min_{x \in \mathbb{R}^d} \left\{ f(x) = \frac{1}{n} \sum_{i=1}^n f_i(x) \right\}$$

$$f_i(x) := \mathbb{E}_{\xi \sim \mathcal{D}_i} [f_i(x, \xi)]$$

$$\mathcal{D}_1 = \dots = \mathcal{D}_n$$

\mathcal{D}_i = uniform distribution over m outcomes

$$\min_{x \in \mathbb{R}^d} \left\{ f(x) = \frac{1}{m} \sum_{i=1}^m f_i(x) \right\}$$

(after butchering/redefining notation)



Table 1: Comparison of the *worst-case time complexity* guarantees of methods that work with asynchronous computations in the setup from Section 1 (up to smoothness constants). We assume that $\tau_i \in [0, \infty]$ is the bound on the times required to calculate one stochastic gradient ∇f_j by worker i , $\tau_1 \leq \dots \leq \tau_n$, and $m \geq n \log n$.
 Abbr: $\delta^0 := f(x^0) - f^*$, $m = \#$ of data samples, $n = \#$ of workers, $\varepsilon = \text{error tolerance}$.

Method	Worst-Case Time Complexity	Comment
Hero GD (Soviet GD)	$\tau_1 m \frac{\delta^0}{\varepsilon} \quad (\tau_n \frac{m}{n} \frac{\delta^0}{\varepsilon})$	Suboptimal
Hero PAGE (Soviet PAGE) [Li et al., 2021]	$\tau_1 m + \tau_1 \frac{\delta^0}{\varepsilon} \sqrt{m} \quad (\tau_n \frac{m}{n} + \tau_n \frac{\delta^0}{\varepsilon} \frac{\sqrt{m}}{n})$	Suboptimal
SYNTHESIS [Liu et al., 2022]	—	Limitations: bounded gradient assumption, calculates the full gradients ^(a) , suboptimal. ^(b)
Asynchronous SGD [Koloskova et al., 2022] [Mishchenko et al., 2022]	$\frac{\delta^0}{\varepsilon} \left(\left(\sum_{i=1}^n \frac{1}{\tau_i} \right)^{-1} \left(\frac{\sigma^2}{\varepsilon} + n \right) \right)$	Limitations: σ^2 -bounded variance assumption, suboptimal when ε is small.
Rennala SGD [Tyurin and Richtárik, 2023]	$\frac{\delta^0}{\varepsilon} \min_{j \in [n]} \left(\left(\sum_{i=1}^j \frac{1}{\tau_i} \right)^{-1} \left(\frac{\sigma^2}{\varepsilon} + j \right) \right)$	Limitations: σ^2 -bounded variance assumption, suboptimal when ε is small.
Freya PAGE (Theorems 7 and 8)	$\begin{aligned} & \min_{j \in [n]} \left(\left(\sum_{i=1}^j \frac{1}{\tau_i} \right)^{-1} (m + j) \right) \\ & + \frac{\delta^0}{\varepsilon} \min_{j \in [n]} \left(\left(\sum_{i=1}^j \frac{1}{\tau_i} \right)^{-1} (\sqrt{m} + j) \right) \end{aligned}$	Optimal in the large-scale regime, i.e., $\sqrt{m} \geq n$ (see Section 5)
Lower bound (Theorem 10)	$\begin{aligned} & \min_{j \in [n]} \left(\left(\sum_{i=1}^j \frac{1}{\tau_i} \right)^{-1} (m + j) \right) \\ & + \frac{\delta^0}{\sqrt{m\varepsilon}} \min_{j \in [n]} \left(\left(\sum_{i=1}^j \frac{1}{\tau_i} \right)^{-1} (m + j) \right) \end{aligned}$	—

Freya PAGE has *universally better guarantees* than all previous methods: the dependence on ε is $\mathcal{O}(1/\varepsilon)$ (unlike Rennala SGD and Asynchronous SGD), the dependence on $\{\tau_i\}$ is harmonic-like and robust to slow workers (robust to $\tau_n \rightarrow \infty$) (unlike Soviet PAGE and SYNTHESIS), the assumptions are weak, and the time complexity of Freya PAGE is optimal when $\sqrt{m} \geq n$.

^(a) In Line 3 of their Algorithm 3, they calculate the full gradient, assuming that it can be done for free and not explaining how.

^(b) Their convergence rates in Theorems 1 and 3 depend on a bound on the delays Δ , which in turn depends on the performance of the slowest worker. Our method does not depend on the slowest worker if it is too slow (see Section 4.3), which is required for optimality.

^(c) We prove better time complexity in Theorem 6, but this result requires the knowledge of $\{\tau_i\}$ in advance, unlike Theorems 7 and 8.

Algorithm 1 Freya PAGE

- 1: **Parameters:** starting point $x^0 \in \mathbb{R}^d$, learning rate $\gamma > 0$, minibatch size $S \in \mathbb{N}$, probability $p \in (0, 1]$, initialization $g^0 = \nabla f(x^0)$ using **ComputeGradient**(x^0) (Alg. 2)
- 2: **for** $k = 0, 1, \dots, K - 1$ **do**
- 3: $x^{k+1} = x^k - \gamma g^k$
- 4: Sample $c^k \sim \text{Bernoulli}(p)$
- 5: **if** $c^k = 1$ **then** (with probability p)
- 6: $\nabla f(x^{k+1}) = \text{ComputeGradient}(x^{k+1})$ (Alg. 2)
- 7: $g^{k+1} = \nabla f(x^{k+1})$
- 8: **else** (with probability $1 - p$)
- 9: $\frac{1}{S} \sum_{i \in \mathcal{S}^k} (\nabla f_i(x^{k+1}) - \nabla f_i(x^k)) = \text{ComputeBatchDifference}(S, x^{k+1}, x^k)$ (Alg. 3)
- 10: $g^{k+1} = g^k + \frac{1}{S} \sum_{i \in \mathcal{S}^k} (\nabla f_i(x^{k+1}) - \nabla f_i(x^k))$
- 11: **end if**
- 12: **end for**

(note): \mathcal{S}^k is a set of i.i.d. indices that are sampled from $[m]$, *uniformly with replacement*, $|\mathcal{S}^k| = S$

Algorithm 2 ComputeGradient(x)

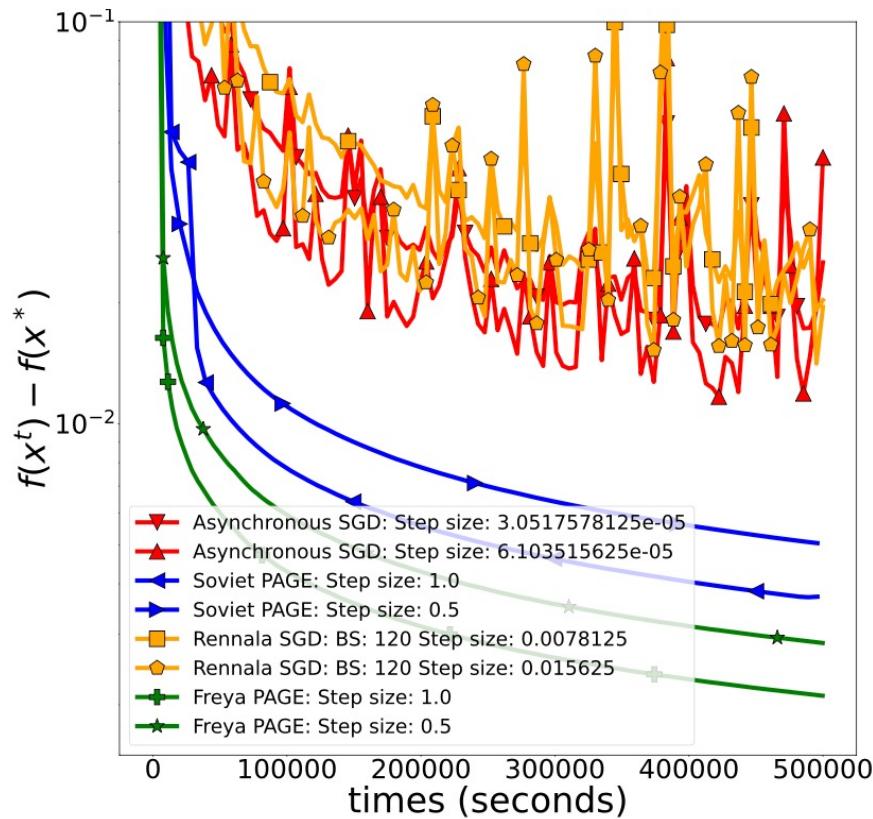
```
1: Input: point  $x \in \mathbb{R}^d$ 
2: Init  $g = 0 \in \mathbb{R}^d$ , set  $\mathcal{M} = \emptyset$ 
3: Broadcast  $x$  to all workers
4: For each worker  $i \in [n]$ , sample  $j$  from  $[m]$  uniformly and ask it to calculate  $\nabla f_j(x)$ 
5: while  $\mathcal{M} \neq [m]$  do
6:   Wait for  $\nabla f_p(x)$  from a worker
7:   if  $p \in [m] \setminus \mathcal{M}$  then
8:      $g \leftarrow g + \frac{1}{m} \nabla f_p(x)$ 
9:     Update  $\mathcal{M} \leftarrow \mathcal{M} \cup \{p\}$ 
10:  end if
11:  Sample  $j$  from  $[m] \setminus \mathcal{M}$  uniformly and ask this worker to calculate  $\nabla f_j(x)$ 
12: end while
13: Return  $g = \frac{1}{m} \sum_{i=1}^m \nabla f_i(x)$ 
```

Algorithm 3 ComputeBatchDifference(S, x, y)

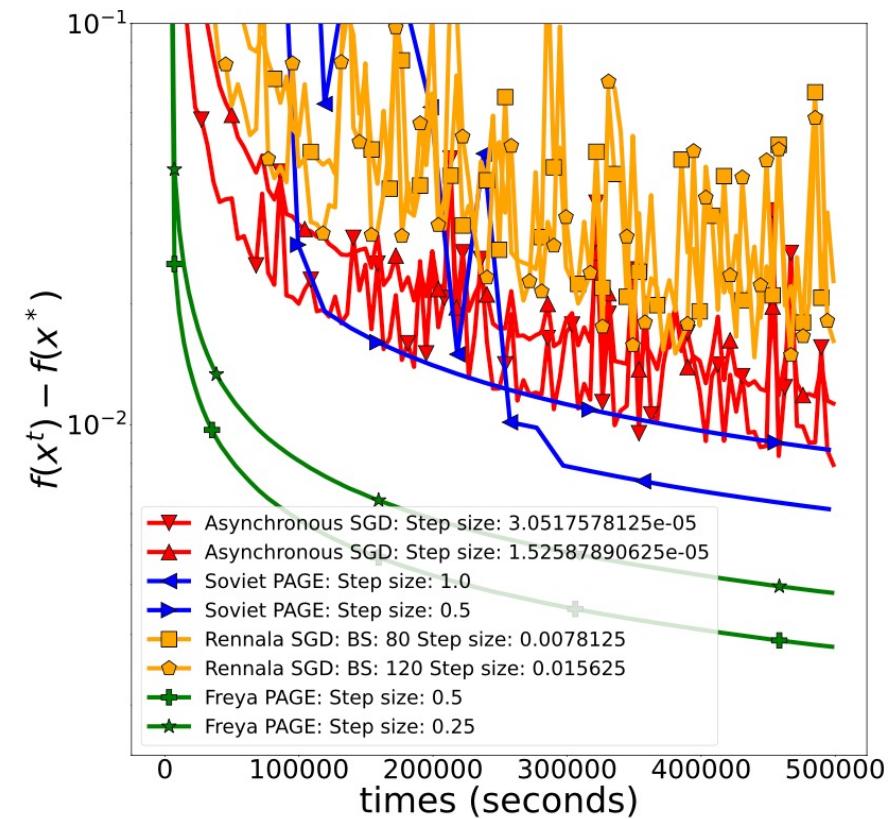
```
1: Input: batch size  $S \in \mathbb{N}$ , points  $x, y \in \mathbb{R}^d$ 
2: Init  $g = 0 \in \mathbb{R}^d$ 
3: Broadcast  $x, y$  to all workers
4: For each worker, sample  $j$  from  $[m]$  uniformly and ask it to calculate  $\nabla f_j(x) - \nabla f_j(y)$ 
5: for  $i = 1, 2, \dots, S$  do
6:   Wait for  $\nabla f_p(x) - \nabla f_p(y)$  from a worker
7:    $g \leftarrow g + \frac{1}{S} (\nabla f_p(x) - \nabla f_p(y))$ 
8:   Sample  $j$  from  $[m]$  uniformly and ask this worker to calculate  $\nabla f_j(x) - \nabla f_j(y)$ 
9: end for
10: Return  $g$ 
```

Notes: i) the workers can aggregate ∇f_p locally, and the algorithm can call AllReduce once to collect all calculated gradients. ii) By splitting $[m]$ into blocks, instead of one ∇f_p , we can ask the workers to calculate the sum of one block in Alg. 2 (and use a similar idea in Alg. 3).

Freya PAGE: Experiment 1



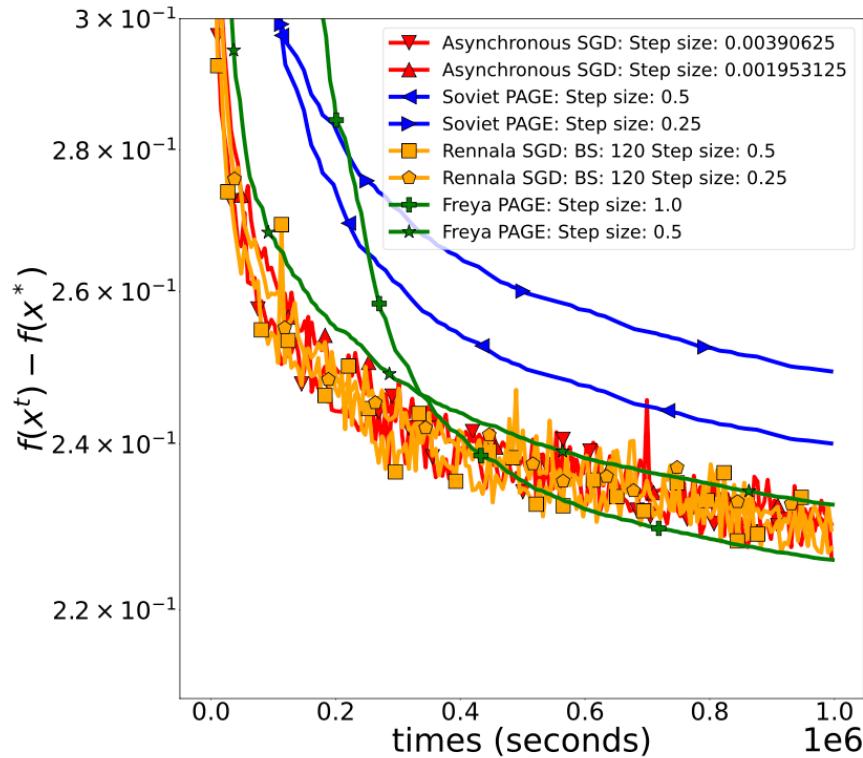
(a) $n = 1000$



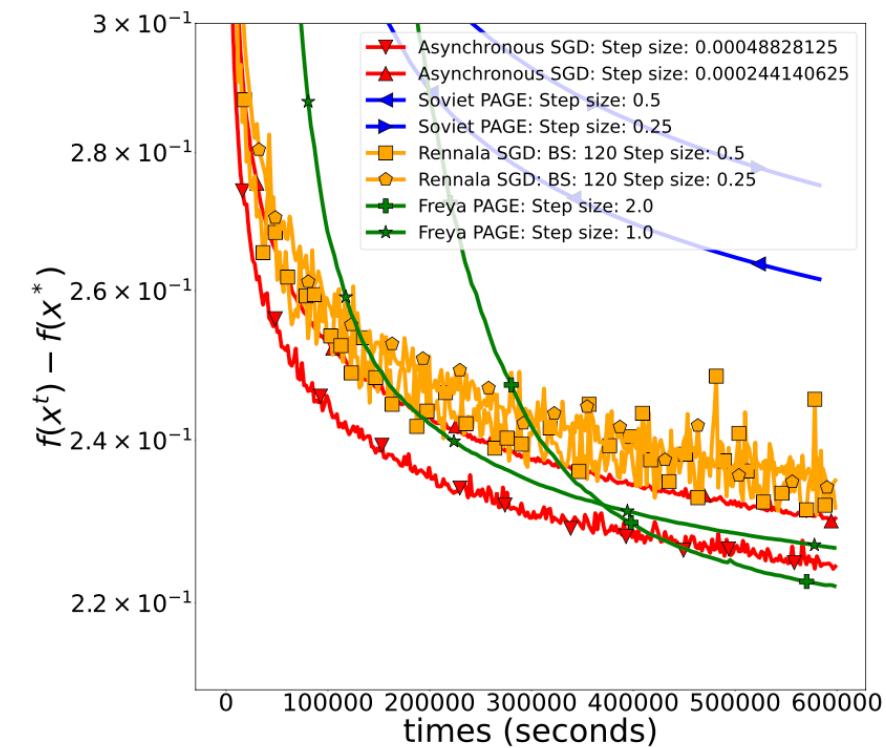
(b) $n = 10000$

Figure 1: Experiments with nonconvex quadratic optimization tasks. We plot function suboptimality against elapsed time.

Freya PAGE: Experiment 2



(a) $n = 100$



(b) $n = 10000$

Figure 2: Experiments with the logistic regression problem on the MNIST dataset.

Freya PAGE: Experiment 2

Table 2: Mean and variance of algorithm accuracies on the MNIST test set during the final 100K seconds of the experiments from Figure 2b.

Method	Accuracy	Variance of Accuracy
Asynchronous SGD [Koloskova et al., 2022] [Mishchenko et al., 2022]	92.60	5.85e-07
Soviet PAGE [Li et al., 2021]	92.31	1.62e-07
Rennala SGD [Tyurin and Richtárik, 2023]	92.37	3.12e-06
Freya PAGE	92.66	1.01e-07

On the Optimal Time Complexities in Decentralized Stochastic Asynchronous Optimization

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Abstract

We consider the decentralized stochastic asynchronous optimization setup, where there are n nodes connected via a directed graph. Each node performs local computation with each other using objects in a multigraph. The local computation at the i -th node is denoted by $f_i(\cdot)$. We propose a new optimal method, Amelie SGD, which is based on a novel decomposition of the total time complexity. Our analysis shows that Amelie SGD is faster than the state-of-the-art SGD and a new nearly optimal method, Fregia SGD, and a new optimal method, Amelie SGD, has a lower variance than Fregia SGD. The proposed method is robust to heterogeneity and matches our lower bounds up to a logarithmic factor in the heterogeneous setting. Our analysis also shows that Amelie SGD is faster than all the state-of-the-art previous asynchronous/stochastic methods in the decentralized setup.

1. Introduction

We consider the smooth convex optimization problem

$$\min_{x \in \mathbb{R}^d} f(x) := \mathbb{E}_{i \sim \mathcal{D}_n} [f_i(x)],$$

where $\mathcal{D}_n = \mathbb{R}^d \times S_n = \mathbb{R}^d$, and D_n is a distribution on a non-empty set S . For a given $\varepsilon \geq 0$, we want to find a ε -possibly stochastic point \hat{x} , called an ε -suboptimal point, such that $\mathbb{E}[f(\hat{x})] \leq \varepsilon$.

There are two main ways to approach this problem: a central server approach with round-trip communication with clients, and a decentralized approach without a central server.

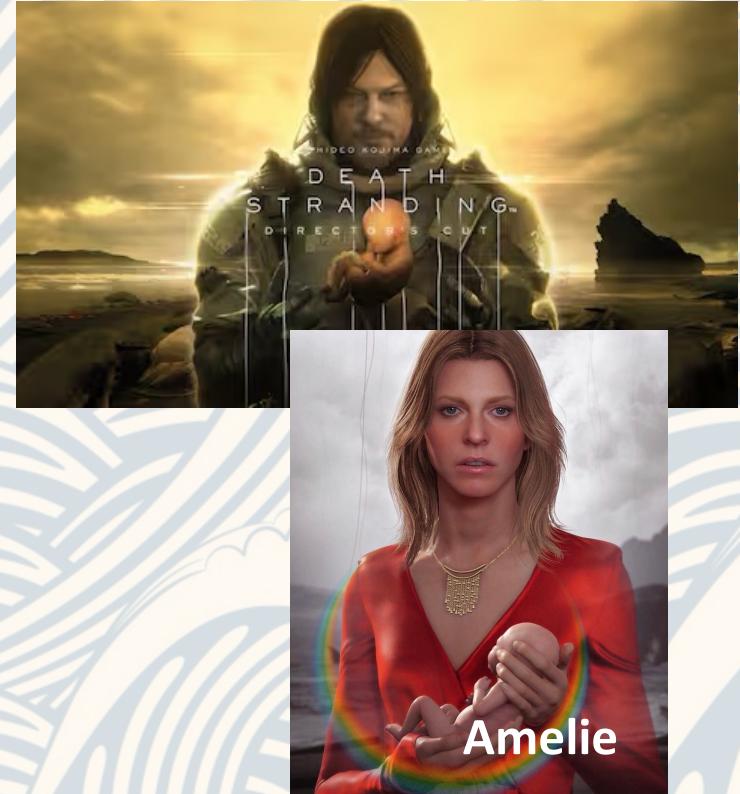
1.1. Decentralized setup with direct communication

We investigate the following decentralized asynchronous setup. Assume that we have n workstations with the associated computation times $\{\lambda_i\}_{i \in [n]}$, and communication times $\{\mu_{i,j}\}_{i,j \in [n]}$. It takes time λ_i to compute a function $f_i(\cdot)$ at the i -th node, and it takes time $\mu_{i,j} \in [0, \infty)$ seconds to send and receive a vector $v \in \mathbb{R}^d$ from the i -th node to the j -th node if it is possible for the i -th node to communicate with the j -th node. We assume that the network is fully connected and is parallel. We would like to emphasize that $\lambda_i \in [0, \infty)$, and $\mu_{i,j} \in [0, \infty)$ are only upper bounds, and the actual computation and communication times can be smaller than the bounds. The nodes are not necessarily homogeneous. For simplicity of presentation, we assume the upper bounds are static; however, in practice, they can change over time due to various factors such as network latency, power consumption, and resource usage. For simplicity of presentation, we assume the upper bounds are static; however, in practice, they can change over time due to various factors such as network latency, power consumption, and resource usage.

We consider any weighted directed multigraph parameterized by a vector $b \in \mathbb{R}^n$ such that $b_i \in [0, n]$, and a matrix of distances $\{\rho_{i,j}\}_{i,j \in [n]} \in \mathbb{R}^{n \times n}$ such that $\rho_{i,j} \in [0, \infty)$ for all $i, j \in [n]$ and $\rho_{i,i} = 0$ for all $i \in [n]$. We denote the distance between the i -th node and the j -th node by $d_{i,j} = \rho_{i,j}$ if $b_j > i$ and $d_{i,j} = \rho_{j,i}$ if $b_i > j$. For this setup, it would be convenient to define the distance of the observed path from

Amelie SGD

Optimal Decentralized SGD under Computation & Communication Heterogeneity



Decentralized Setup: Amelie SGD

Method	The Worst-Case Time Complexity Guarantees	Comment
Minibatch SGD	$\frac{L\Delta}{\varepsilon} \max \left\{ \left(1 + \frac{\sigma^2}{n\varepsilon}\right) \max\left\{ \max_{i,j \in [n]} \tau_{i \rightarrow j}, \max_{i \in [n]} h_i \right\} \right\}$	suboptimal if σ^2/ε is large
RelaySGD, Gradient Tracking <small>(Vogels et al., 2021) (Liu et al., 2024)</small>	$\geq \frac{\max_{i \in [n]} L_i \Delta}{\varepsilon} \frac{\sigma^2}{n\varepsilon} \max_{i \in [n]} h_i$	requires local L_i -smooth. of f_i , suboptimal if σ^2/ε is large (even if $\max_{i \in [n]} L_i = L$)
Asynchronous SGD <small>(Even et al., 2024)</small>	—	requires similarity of the functions $\{f_i\}$, requires local L_i -smooth. of f_i
Amelie SGD and Lower Bound <small>(Thm. 7 and Cor. 2)</small>	$\frac{L\Delta}{\varepsilon} \max \left\{ \max_{i,j \in [n]} \tau_{i \rightarrow j}, \max_{i \in [n]} h_i, \frac{\sigma^2}{n\varepsilon} \left(\frac{1}{n} \sum_{i=1}^n h_i \right) \right\}$	Optimal up to a constant factor



**The End
(for real)**