

Accommodating LLM Training over Decentralized Computational Resources

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05.06.24



THE HONG KONG
UNIVERSITY OF SCIENCE
AND TECHNOLOGY

Amazing Progress of ML/AI

DATA
LAB
6TH EDITION



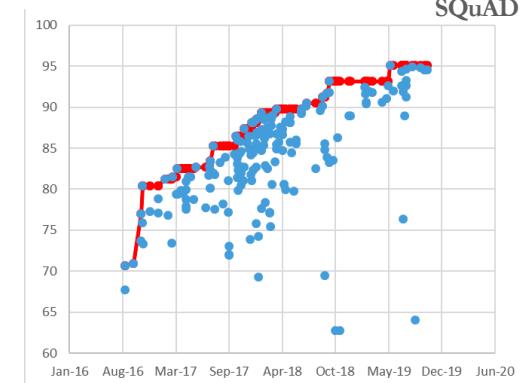
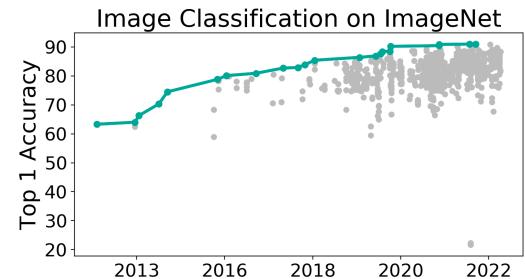
"space robot studying a book in front of Stanford"



Write a haiku from the perspective of a copywriter who is feeling sad that AI might diminish the value of the written word



Words on a screen,
Once valued, now just a blur
Machine takes the pen.



The challenge of Today:

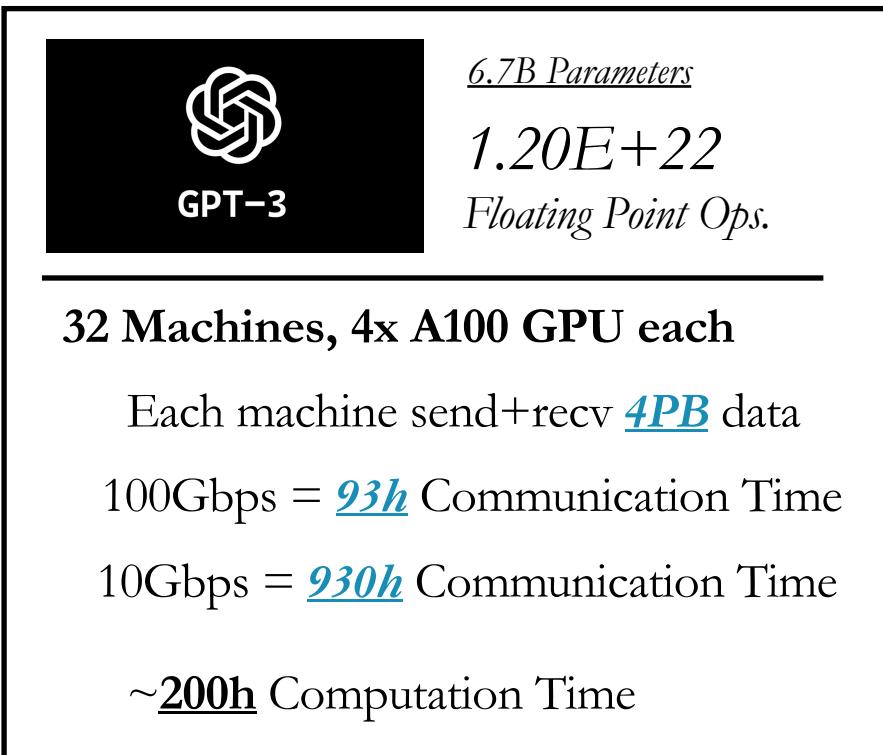
(Million \$)

Building ML Applications at SOTA scale is expensive!

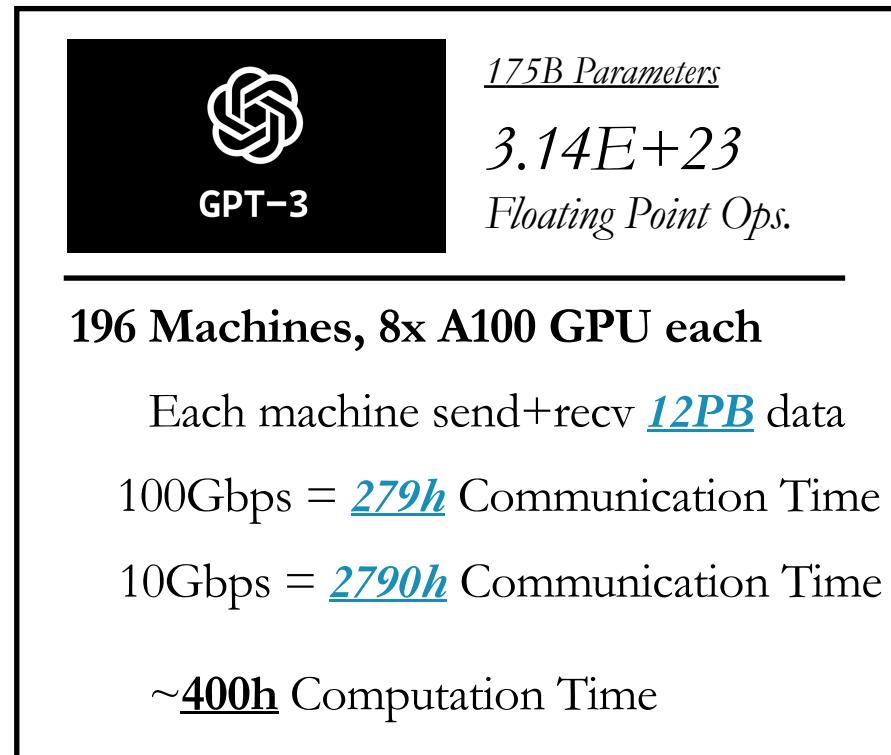
Further Scaling is facing non-linear bottlenecks.

Bottleneck: Communications & Data Movement

Distributed training at scale is communication-intensive.



(Today) Model training today is largely restricted to centralized data centers with fast network connections. Hard to use cheaper alternatives (Tier 2-4 clouds, Spot Instances, Volunteer Computes, etc.).



(Future) 10x further scaling requires fast connections between 10x machines. Becoming challenging even for data center.



NVIDIA DGX SuperPOD:
Up to 256 GPUs

Optimizing Communications for Distributed and Decentralized Learning.



Communication Bottlenecks across Infrastructure

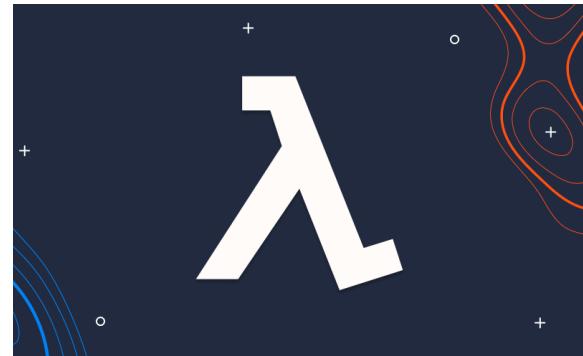
communication becomes slower, open up more choices (and some can be cheaper)



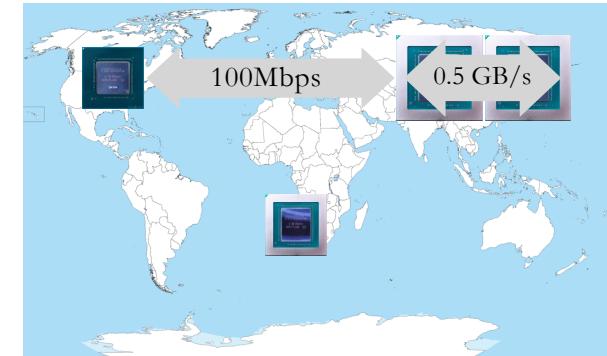
Data Center



(Multi-cloud) Spot Instances



Serverless Environment



Decentralized Network

The more we can optimize communications, the more choices we have when building our infrastructure.

$$\min_x \mathbb{E}_\xi f(\xi, x)$$

$$\min_x \mathbb{E}_\xi f(\xi, x)$$

Data

- (ImageNet) 1.3M Images (est. 160+ GB)
- (GPT-3) 300 Billion Tokens (est. 2+ TB)

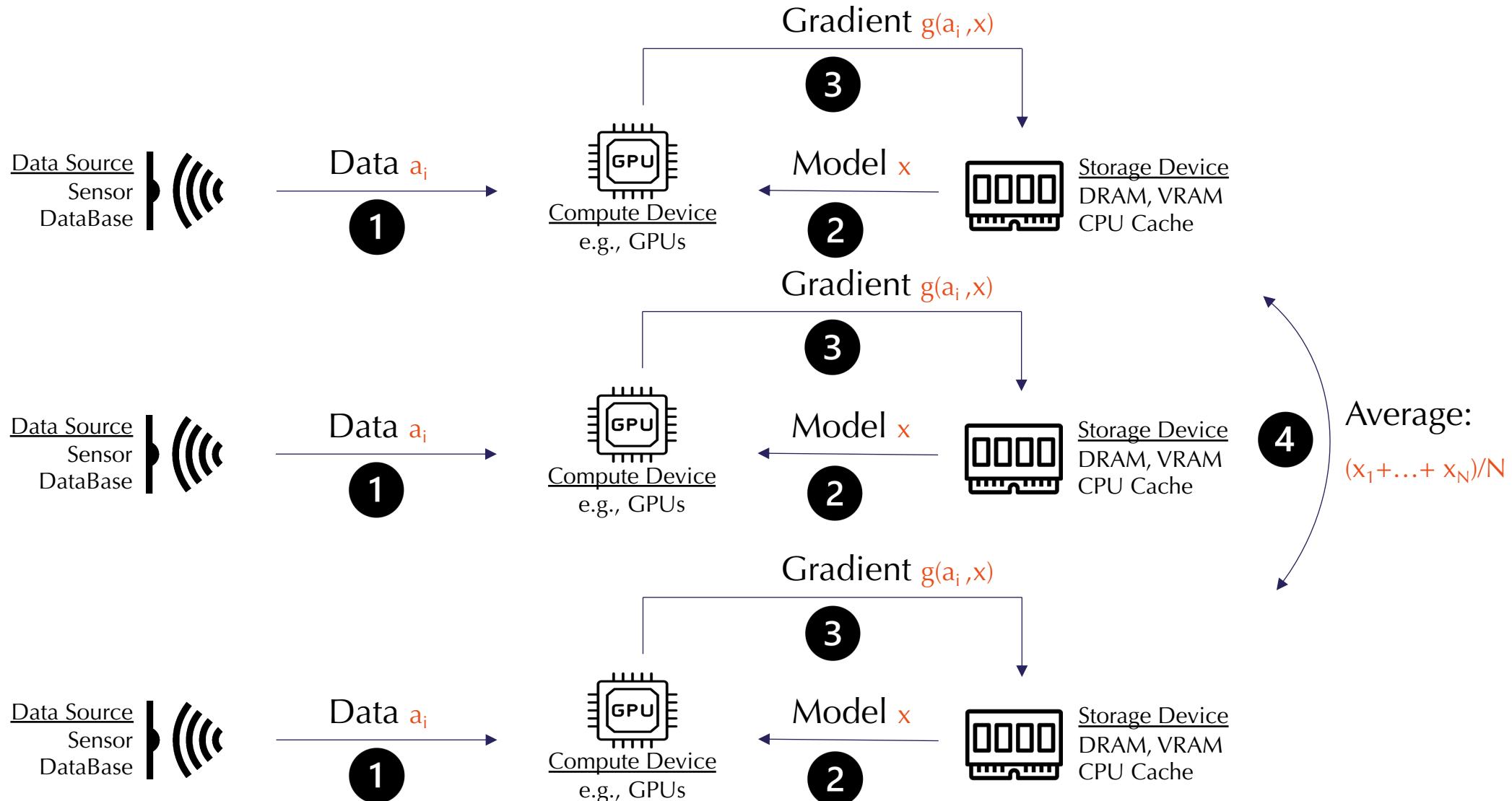
Model

- (GPT-2) 1.3 Billion Parameters (2.6 GB fp16)
- (GPT-3) 175 Billion Parameters (350GB fp16)

Compute

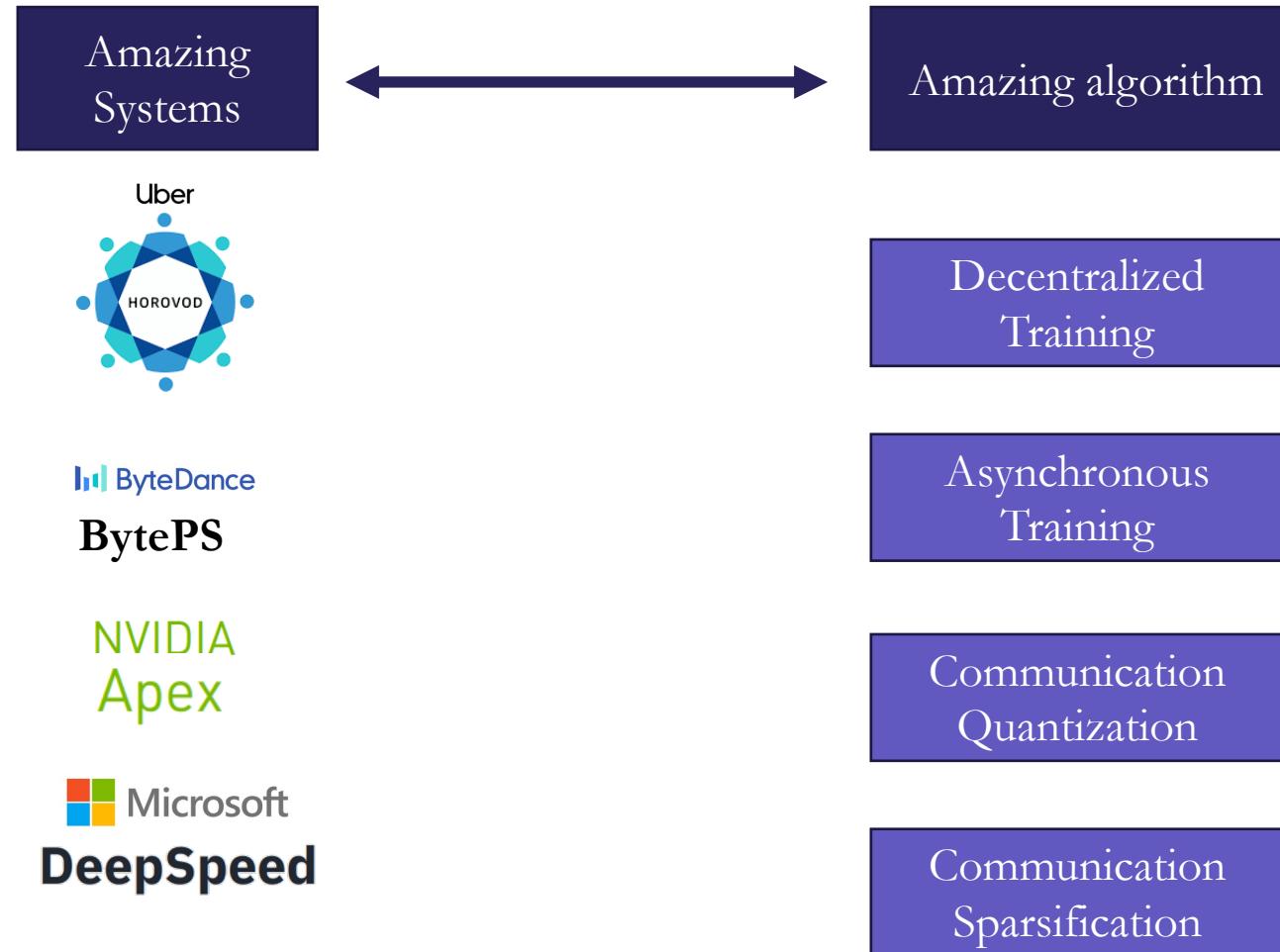
- (GPT-2) est. 2.5 GFLOPS/token
- (GPT-3) est. 0.4 TFLOPS/token

Data Parallel SGD

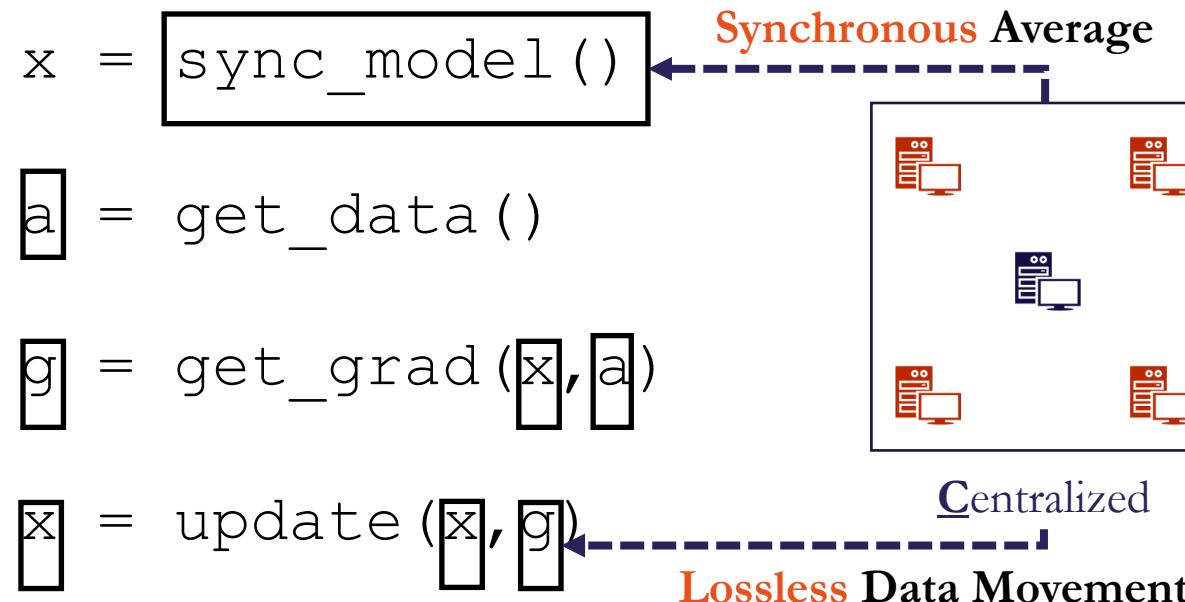


System Optimizations and Relaxed Algorithms

LocalLab@
ETH Zurich



Baseline: Centralized, Synchronous, Lossless, SGD



Idea

- Distribute batch gradient calculation to multiple workers;
- Synchronize workers with a central server (or AllReduce).

Mathematical Formulation

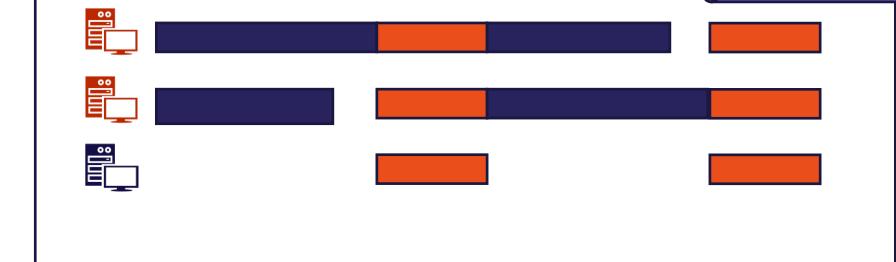
$$x_{t+1} = x_t - \gamma \sum_{i=1..n} g_i(x_t; a_i)$$

Convergence

$$O(1/\sqrt{nT})$$

Goal 1: Keep This Similar

System Profile

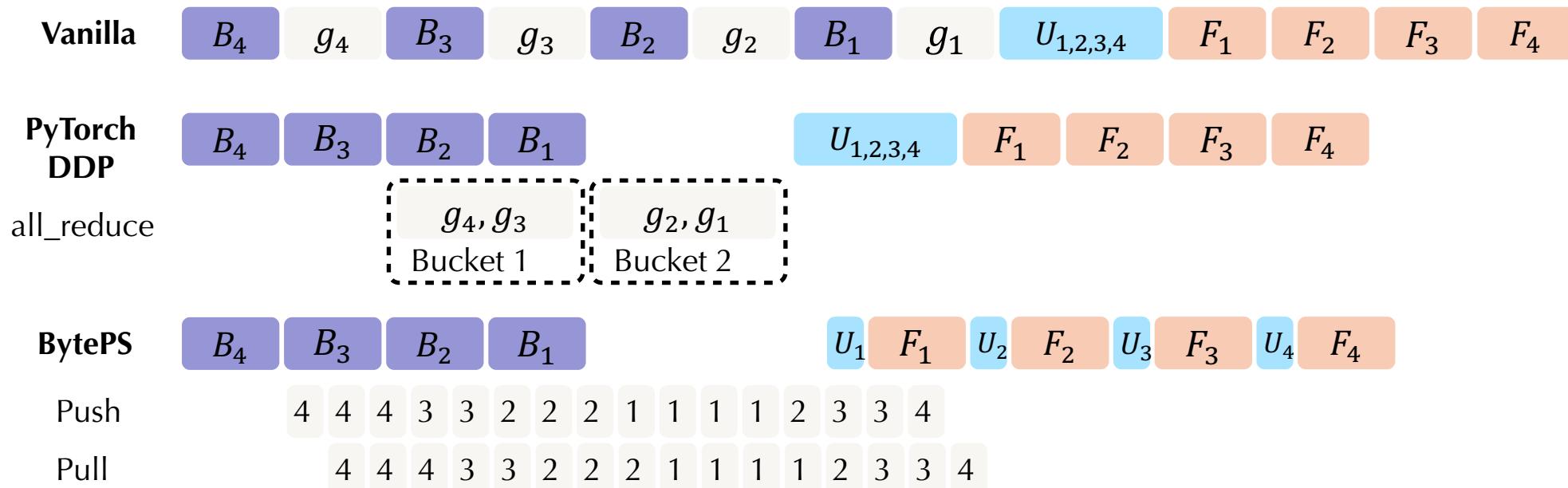


■ Computation
■ Communication

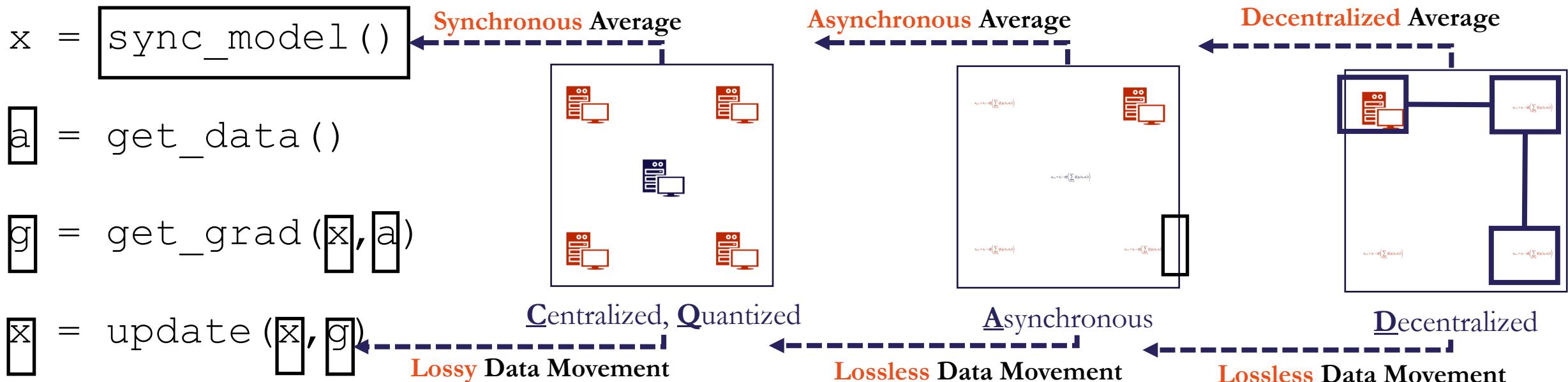
System Optimizations



- **Existing Systems:**
Optimize the standard DP-SGD computation:



Relaxed Algorithms



Mathematical Formulation

$$x_{t+1} = x_t - \gamma Q\left(\sum_{i=1..n} Q(g_i(x_t, a_i))\right)$$

Convergence

$$O(1/\sqrt{nT} + \boxed{\epsilon/\sqrt{T}})$$

Quantization error: ϵ

$$x_{t+1} = x_t - \gamma g(x_{t-\tau_t}; a_i)$$

↑
staleness caused by async

$$x_{t+1,i} = \frac{x_{t,i-1} + x_{t,i} + x_{t,i+1}}{3} - \gamma g(x_{t,i}; a_i)$$

$$O(1/\sqrt{nT} + \boxed{\rho/T^{1.5}})$$

ρ : network topology constant

Attempt 1

Automatic System Optimization for Relaxed Algorithms

Amazing Systems



ByteDance
BytePS

NVIDIA
Apex

Microsoft
DeepSpeed

GAP: Current Amazing Systems Don't Support Recently Developed Amazing Techniques



OUR GOAL:
Distributed Learning with SOTA
Communication Optimization Techniques.

Amazing algorithm

Decentralized Training

Asynchronous Training

Communication Quantization

Communication Sparsification

BAGUA: Scaling up Distributed Learning with System Relaxations

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ABSTRACT

Recent years have witnessed a growing list of systems for distributed data-parallel training. Existing systems largely fit into two paradigms, i.e., parameter server and MPI-style collective operations. On the algorithmic side, researchers have proposed a wide range of techniques to lower the communication via "system relaxations": quantization, decentralization, and communication delay. However, most, if not all, existing systems only rely on standard synchronous and asynchronous stochastic gradient (SG) based optimization, therefore, cannot take advantage of all possible optimizations that the machine learning domain has been developing recently. Given this emerging gap between the current landscapes of systems and theory, we build BAGUA, a MPI-style communication library, providing a collection of primitives, that is both flexible and modular to support state-of-the-art system relaxation techniques of distributed training. Powered by this design, BAGUA has a great ability to implement and extend various state-of-the-art distributed learning algorithms. In a production cluster with up to 16 machines (128 GPUs), BAGUA can outperform PyTorch-DDP, Horovod and BytePS in the end-to-end training time by a significant margin (up to 2X) across a diverse range of tasks. Moreover, we conduct a rigorous tradeoff exploration showing that different algorithms and system relaxations achieve the best performance over different network conditions.

PVLDB Reference Format:

Shaoduo Gan, Xiangru Lian, Rui Wang, Jianbin Chang, Chengjun Liu, Hongmei Shi, Shengzhou Zhang, Xianghong Li, Tengxu Sun, Sen Yang, Ji Liu, Ce Zhang. BAGUA: Scaling up Distributed Learning with System Relaxations. PVLDB, 15(4): 804 - 813, 2022.
doi:10.14778/3503585.3503590

PVLDB Artifact Availability:

The source code, data, and/or other artifacts have been made available at <https://github.com/BAGUASys/bagua>.

1 INTRODUCTION

The increasing performance of distributed machine learning systems has been one of the main driving forces behind the rapid

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*Equal contribution.
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[VLDB 2022]

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ENDOWMENT
WITH SUPPORT FROM THE SWISS NATIONAL SCIENCE FOUNDATION

It is not easy to translate *algorithmic flexibility* into *system performance gain*.

Bagua: System Design & Implementation

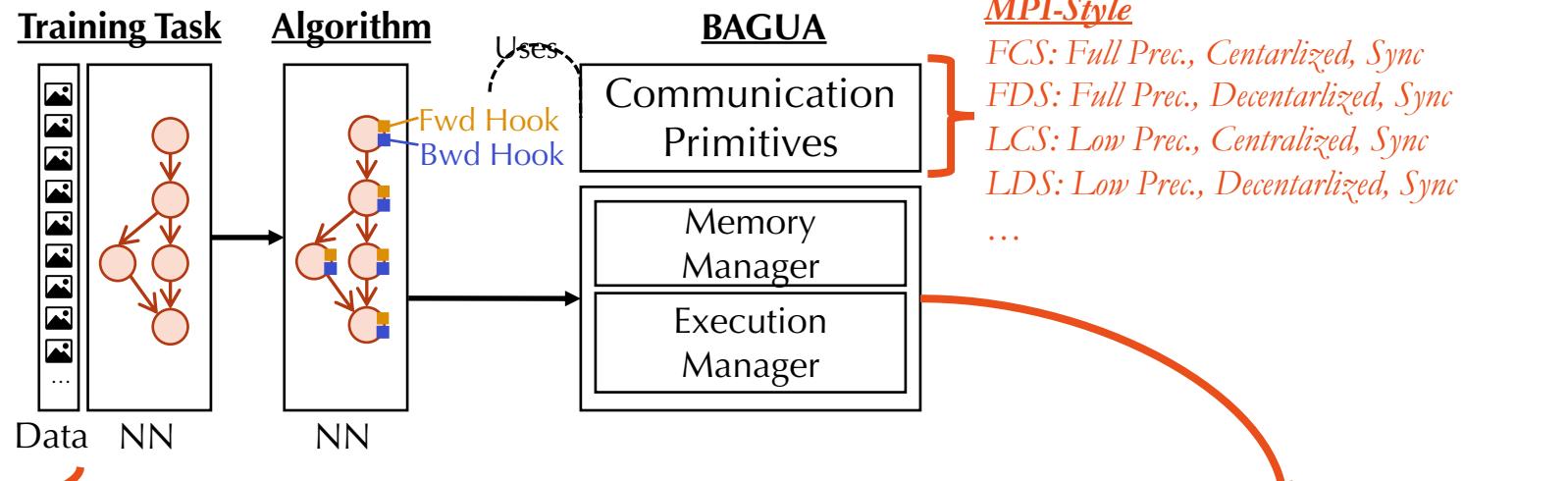
- A modular design to accommodate the diversity of different algorithms and communication patterns.
- An optimization framework that applies automatically to an algorithm implemented in BAGUA.

End user: simply wrap up your training script with BAGUA. Specify the algorithm you want to use

```

1 import torch
2 from bagua import bagua_init, DefaultAlgo
3
4 def main():
5     args = parse_args()
6
7     # define model and optimizer
8     model = MyNet().to(args.device)
9     optimizer = torch.optim.SGD(model.parameters(), lr=args.lr)
10    # transform to BAGUA wrapper
11    model, optimizer = bagua_init(model, optimizer, DefaultAlgo,
12        is_intra)
13
14    # train the model over the dataset
15    for epoch in range(args.epochs):
16        for b_idx, (inputs, targets) in enumerate(train_loader):
17            outputs = model(inputs)
18            loss = torch.nn.CrossEntropyLoss(outputs, targets)
19            optimizer.zero_grad()
20            loss.backward()
21            optimizer.step()

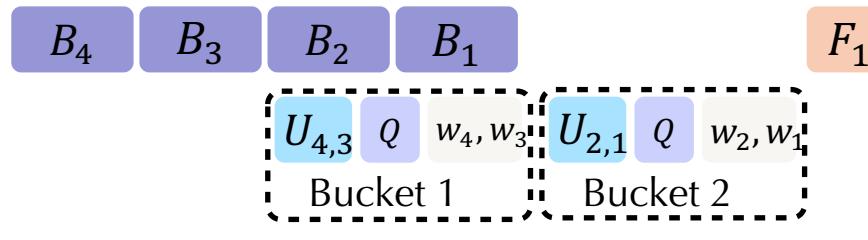
```



E.g., Decentralized, Low Precision Alg.



Automatic

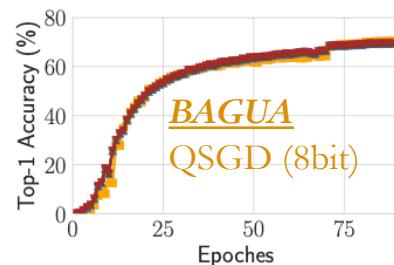


	Forward Computation
	Backward Computation
	Gradient Communication
	Model Update

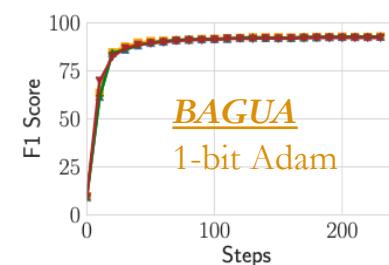
Bagua Results



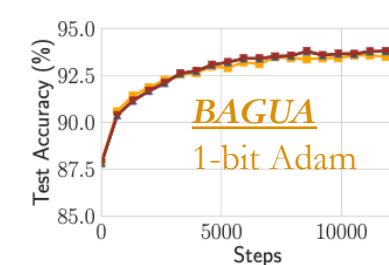
Setup: 16 machines, each 8 V100 GPUs. Connected via $\{10\text{Gbps}, 25\text{Gbps}, 100\text{Gbps}\}$ networks.



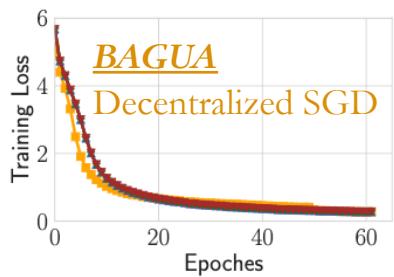
(a) VGG16



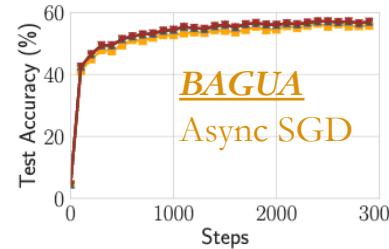
(b) BERT-LARGE Finetune



(c) BERT-BASE Finetune



(d) Transformer



(e) LSTM+AlexNet

- Bagua
- ▲— PyTorch-DDP
- ▲— Horovod
- ▼— BytePS

Network Conditions	VGG16	BERT-LARGE	BERT-BASE	Transformer	LSTM+AlexNet
100 Gbps	1.1×	1.05×	1.27×	1.2×	1.34×
25 Gbps	1.1×	1.05×	1.27×	1.2×	1.34×
10 Gbps	1.94×	1.95×	1.27×	1.2×	1.34×

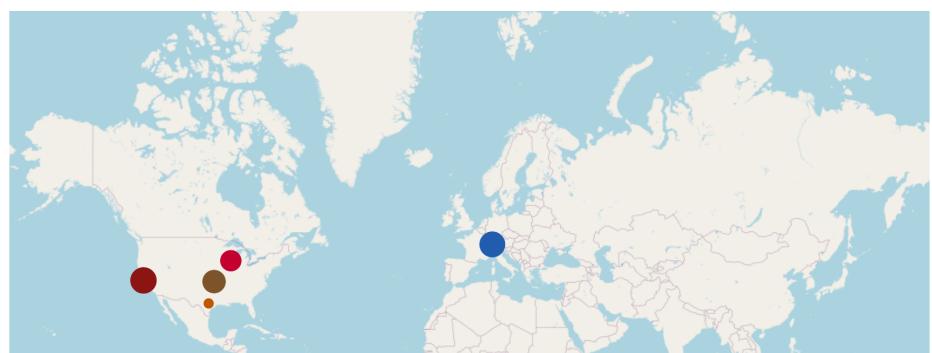
Significant speed-up over {Torch-DDP, Horovod 32bits, Horovod 16bits, BytePS}

Supporting a diverse set of algorithms can provide significant improvements over existing systems.

Same Convergence with Relaxed Algorithms

From Cloud to Decentralized Compute Resource

Instance Size	vCPUs	Instance Memory (GiB)	GPU – A100	GPU memory	Network Bandwidth (Gbps)	GPU Direct RDMA	GPU Direct RDMA	Bandwidth (GB)	Bandwidth (Gbps)	Demand	Price/hr
p4d.24xlarge	96	1152	8	320 GB HBM2	400 ENA and EFA	Yes	Yes	60 NVLink	60 NVLink	Low	\$4.09/hr
p4de.24xlarge (preview)	96	1152	8	640 GB HBM2e	400 ENA and EFA	Yes	Yes	60 NVLink	60 NVLink	Low	\$4.09/hr



 Status Global View

Interruptible • On-Demand #GPUs: ANY 0X 1X 2X 4X 8X 8X+

ID	Type	Location	GPU Model	Cores	Memory	Processor	Storage	Ports	Bandwidth	Price	
m:7424	vast.ai Type #5693570	datacenter:40660 Netherlands, NL	1x A100 SXM4	19.5 TFLOPS Max CUDA: 11.7	80 GB 1401.7 GB/s	Motherboard PCIE 4.0,16x	AMD EPYC 7542 ... 24.0/24 cpu	Storage nvme 583 MB/s	0 ports 270.0 GB	↑628 Mbps ↓602 Mbps verified	\$0.500/hr
m:7207	vast.ai Type #5552288	host:33081	1x A100 SXM4	19.5 TFLOPS Max CUDA: 11.8	39 GB 1140.6 GB/s 300.0 GB/s	Not Specified PCIE 4.0,16x	AMD EPYC 7763 ... 64.0/256 cpu	Storage nvme 1008 MB/s	250 ports 813.1 GB	↑11 Mbps ↓321 Mbps	ability
m:5308	vast.ai Type #5493102	host:33081 Texas, US	1x A100 SXM4	19.5 TFLOPS Max CUDA: 11.7	40 GB 1130.8 GB/s 300.0 GB/s	08XP3P PCIE 4.0,16x	AMD EPYC 7513 ... 32.0/128 cpu	Storage DELL PERC 1218 MB/s	4 ports 238.4 GB	↑11 Mbps ↓317 Mbps	44.4 DLPerf Reliability 48.9 DLP/\$/hr 99.69%

This is what you can get from a decentralized GPU pool!

MAKE BID

Attempt 2

These algorithmic
building blocks need
to *be put together!*

CocktailSGD: Fine-tuning Foundation Models over 500Mbps Networks

Jue Wang ^{*1} Yucheng Lu ^{*2} Binhang Yuan ¹ Beidi Chen ³ Percy Liang ⁴ Christopher De Sa ² Christopher Re ⁴
Ce Zhang ¹

Abstract

Distributed training of foundation models, especially large language models (LLMs), is communication-intensive and so has heavily relied on centralized data centers with fast interconnects. *Can we train on slow networks and unlock the potential of decentralized infrastructure for foundation models?* In this paper, we propose COCKTAILSGD, a novel communication-efficient training framework that combines three distinct compression techniques—random sparsification, top-K sparsification, and quantization—to achieve much greater compression than each individual technique alone. We justify the benefit of such a hybrid approach through a theoretical analysis of convergence. Empirically, we show that COCKTAILSGD achieves up to 117 \times compression in fine-tuning LLMs up to 20 billion parameters without hurting convergence. On a 500Mbps network, COCKTAILSGD only incurs ~ 1.2 \times slowdown compared with data center networks.

1. Introduction

In recent years, foundation models (Bommasani et al., 2021), including large language models (Brown et al., 2020; Chowdhery et al., 2022; Bommasani et al., 2021; Zhang et al., 2022; Liang et al., 2022; Scao et al., 2022), have enabled rapid advancement for various machine learning tasks, especially in natural language processing (Brants et al., 2007; Austin et al., 2021). Such a significant improvement on quality has been fueled by an ever-increasing amount of data and computes that are required in training these models (Kaplan et al., 2020). Today, training even modest scale models requires a significant amount of compute: For example, fine-tuning GPT-J-6B (6 billion parameters) over

^{*}Equal contribution. ¹ETH Zürich, Switzerland ²Cornell University, USA ³Carnegie Mellon University, USA ⁴Stanford University, USA. Correspondence to: Jue Wang <juewang@inf.ethz.ch>.

Proceedings of the 40th International Conference on Machine Learning, Honolulu, Hawaii, USA. PMLR 202, 2023. Copyright 2023 by the author(s).

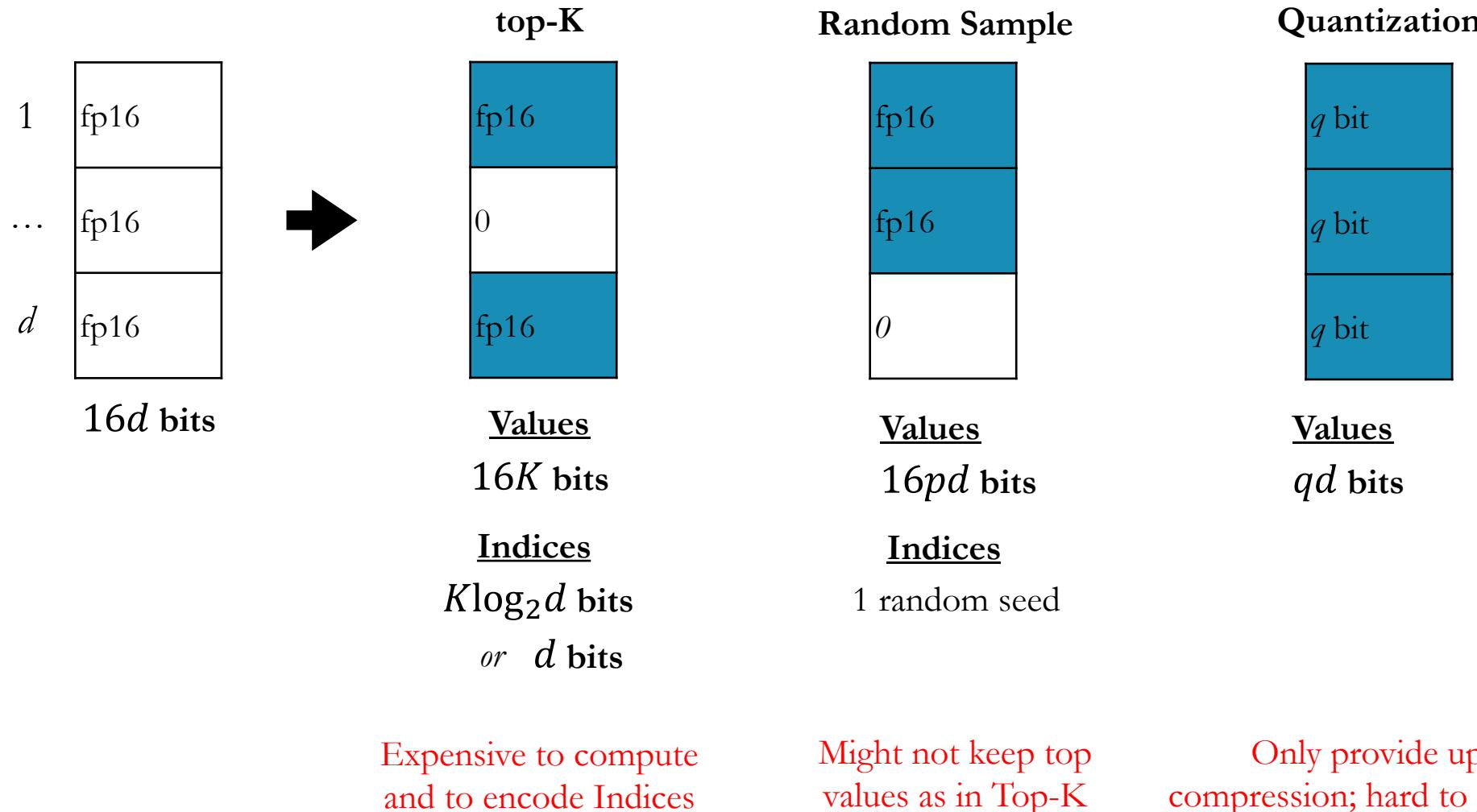
merely 10 billion tokens would require 6 petaflops-days: 8 A100 GPUs running at 50% capacity for 5 days!

When training foundation models in a distributed way, *communication* is the key bottleneck in scaling. As an example, fine-tuning GPT-J-6B over 10 billion tokens with a batch size of 262K tokens over 4 machines (each with 2 A100 GPUs) would require 915.5 TB data being communicated during the whole training process! The computation time for such a workload is 114 hours, which means that we need to have at least 20 Gbps connections between these machines to bring the communication overhead to the same scale as the computation time. Not surprisingly, today's infrastructure for training and fine-tuning foundation models are largely *centralized*, with GPUs connected via fast 100Gbps–400Gbps connections (Microsoft, 2020).

Such a heavy reliance on centralized networks increases the cost of infrastructure, and makes it incredibly hard to take advantage of cheaper alternatives, including tier 2 to tier 4 clouds, spot instances and volunteer compute. For example, while volunteering compute projects such as Folding@Home can harvest significant amount of computes for embarrassingly parallelizable workloads (e.g. 2.43exaflops in April 2020 (Larson et al., 2009)), it is challenging to harvest these cycles for foundation model training due to the communication bottleneck. Recently, there has been an exciting collection of work focusing on the decentralized training of neural networks, including those that are algorithmic (Lian et al., 2017; Ryabinin & Gusev, 2020; Diskin et al., 2021; Ryabinin et al., 2021; Yuan et al., 2022; Jue et al.) as well as system efforts such as Training Transformer Together (Borzunov et al., 2022b), and PETALS (Borzunov et al., 2022a). However, despite of these recent efforts, communication is still a significant bottleneck, and one can only compress the communication by at most 10-30 \times in these recent efforts without hurting convergence. To fully close the gap between centralized infrastructure (100Gbps) and decentralized infrastructure (100Mbps-1Gbps), we need to decrease the communication overhead by at least 100 \times !

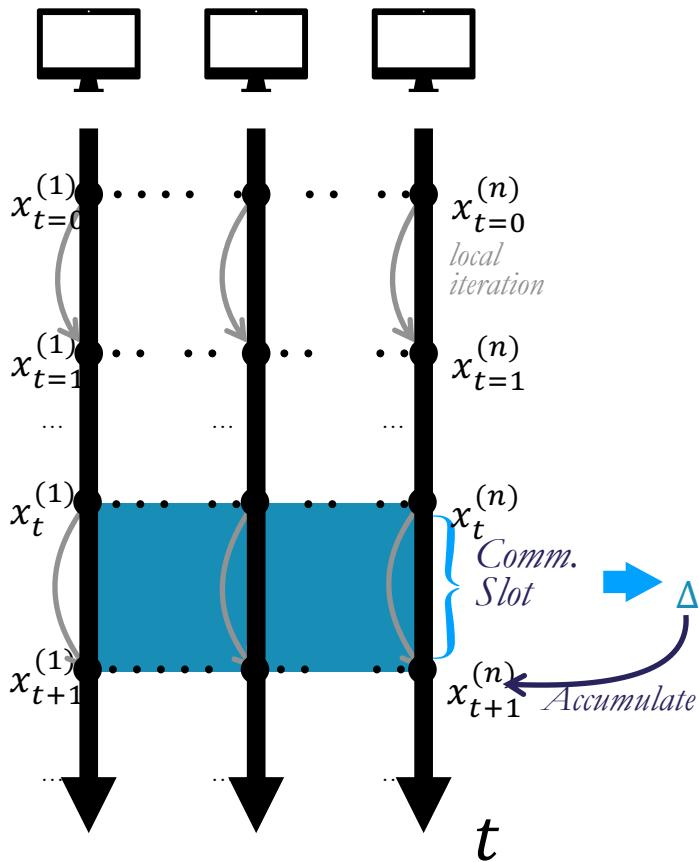
Luckily, there have also been rapid development of communication-efficient optimization algorithms and these efforts provide the foundational building blocks of this paper. Researchers have proposed a wide range of

Three Methods of Compression



It is very hard to reach *100X compression* ratio with a single method.

CocktailSGD: Mixture of Compression Methods



Idea: A Mixture of communication compression techniques.

Looking at Δ_t :

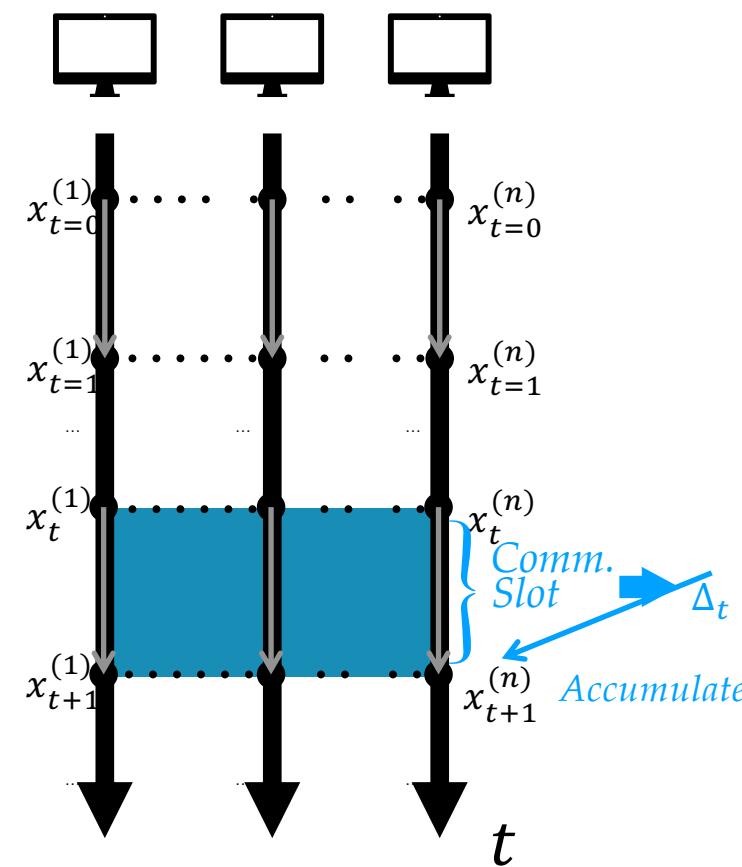
- It has 1-step staleness // asynchrony
- At t , randomly pick $p\%$ parameters to communicate // local training: compress $\sim \frac{1}{p\%} \times$
- For selected parameters, let $\delta_t^{(i)}$ be local model updates since last communication:
 - $\tilde{\delta}_t^{(i)} = \text{top-}K\%(\delta_t^{(i)})$ // topK: compress $\sim \frac{1}{K\%} \times$
 - $\hat{\delta}_t^{(i)} = \text{Quantize}(\tilde{\delta}_t^{(i)}, q\text{bits})$ // Quantization: compress $\sim \frac{16}{b} \times$
- Communicate: $\Delta_t = \sum_i \hat{\delta}_t^{(i)}$

As long as **Communication** fully fills the **Comm. Slot**, no slow down caused by communication.

We need to compose different communication compression techniques together!

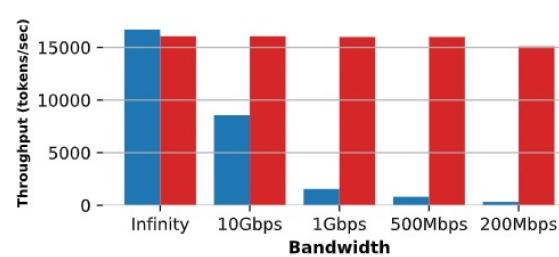
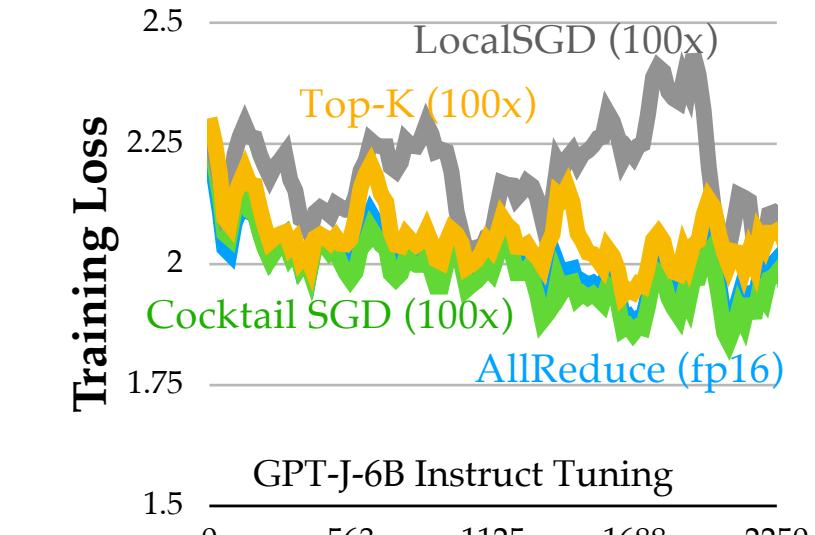
“Cocktail SGD”: Data Parallel over 1Gbps

DS-GLoLo
ETH Zurich

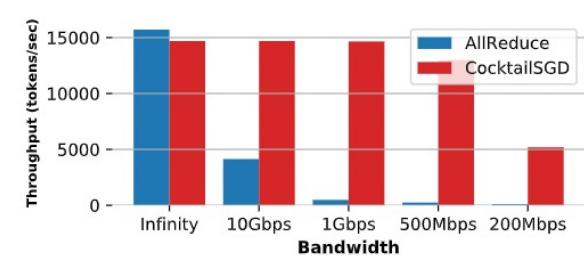
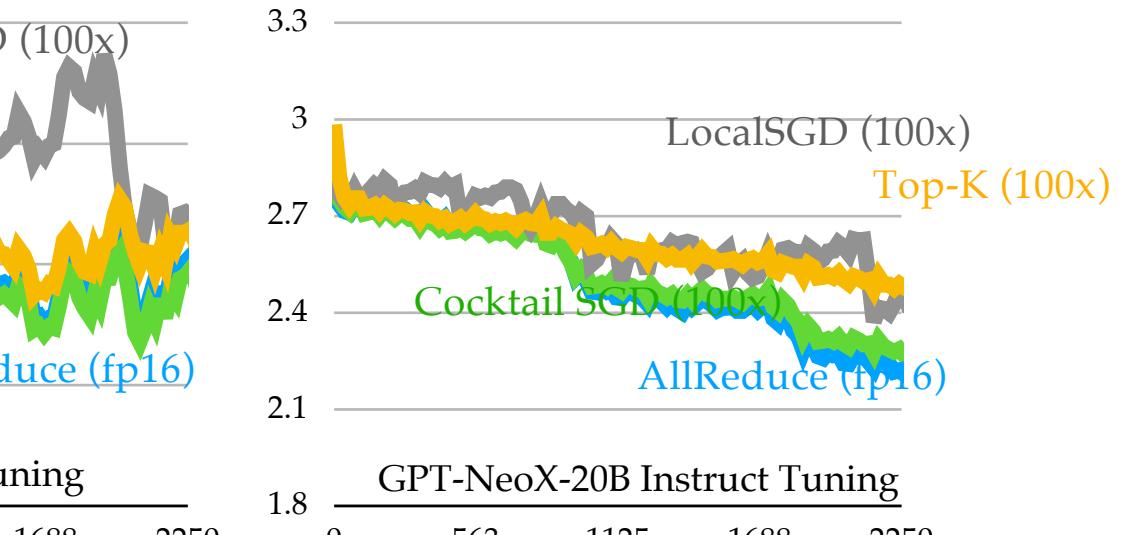


As long as **Communication** fully fills the **Comm. Slot**, no slow down caused by communication.

Different communication compression techniques complement each other and compose well!



(b) GPT-J-6B

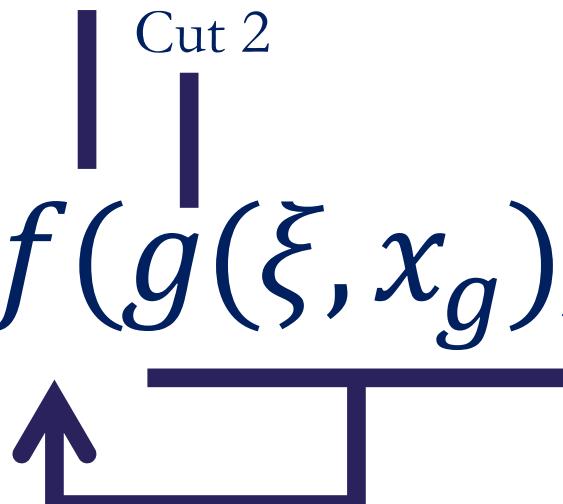


(c) GPT-NeoX-20B

Data parallel over ~1Gbps network!

Large language model training goes
beyond data parallelism.

$$\min_x \mathbb{E}_\xi f(\xi, x) \rightarrow$$

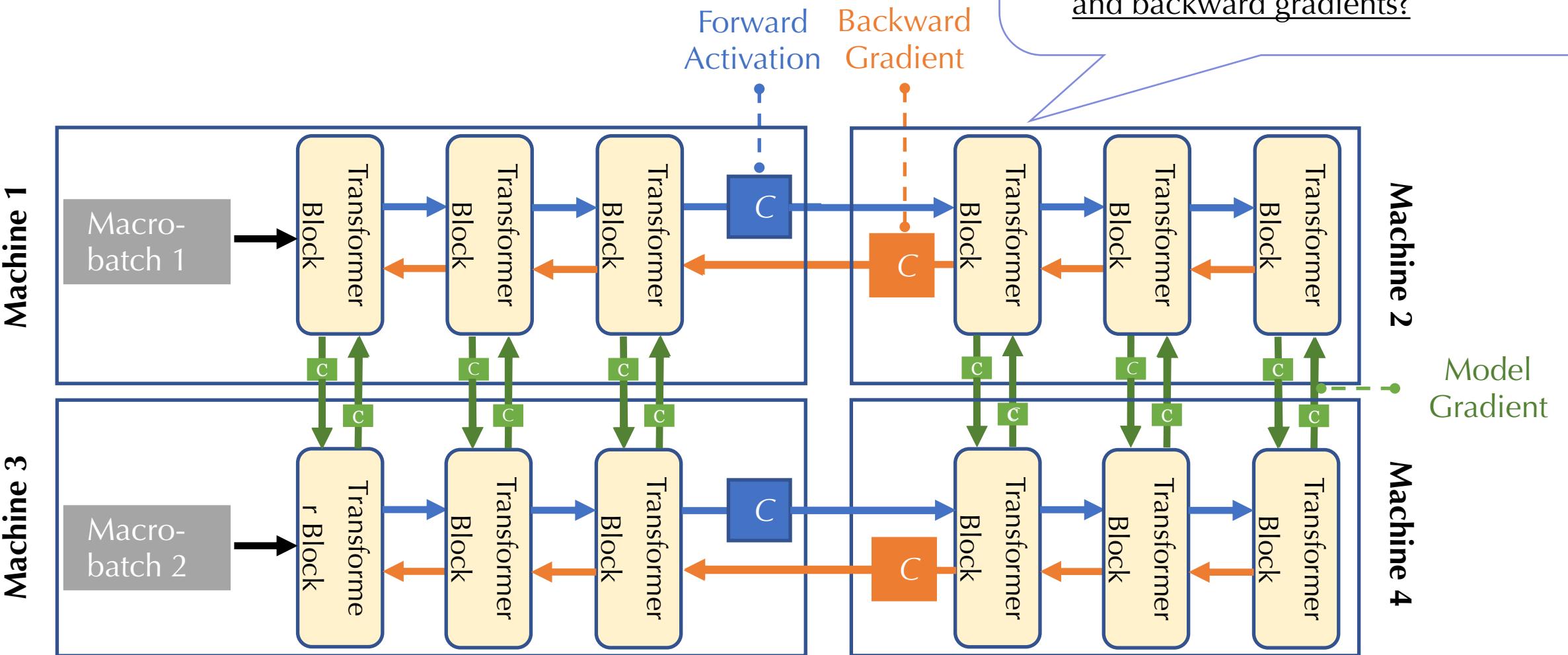
$$\min_{x_f, x_g} \mathbb{E}_\xi f(g(\xi, x_g), x_f)$$


Cut 1
Cut 2

Forward Activation

- (GPT-3) 24MB / 1000tokens

Pipeline Parallelism



Decentralized Training of Foundation Models

- Decentralized training of FM: the network is 100× slower, but the pre-training throughput is only 1.7~3.5× slower!
- Decentralized fine-tuning of FM: AQ-SGD communication-efficient pipeline training with activation compression.

Decentralized Training of Foundation Models in Heterogeneous Environments

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Abstract

Training foundation models, such as GPT-3 and PaLM, can be extremely expensive, often involving tens of thousands of GPU hours, especially for large-scale language models that are typically trained in centralized clusters featuring fast, homogeneous interconnects and using carefully designed software systems that support both data parallelism and model/pipeline parallelism. Such dedicated clusters can be costly and difficult to obtain. Can we instead leverage the much greater amount of decentralized, heterogeneous, and lower-bandwidth interconnected compute? Previous works examining the heterogeneous, decentralized setting focus on relatively small models that can be trained in a purely data parallel manner. State-of-the-art schemes for model parallel foundation model training, such as Megatron, only consider the homogeneous data center setting. In this paper, we present the first study of training large foundation models with model parallelism in a decentralized regime over a heterogeneous network. Our key technical contribution is a scheduling algorithm that allocates different computational “tasks” in the training of foundation models to a group of decentralized GPU devices connected by a slow heterogeneous network. We provide a formal cost model and further propose an efficient evolutionary algorithm to find the optimal allocation strategy. We conduct extensive experiments that represent different scenarios for learning over geo-distributed devices simulated using real-world network measurements. In the most extreme case, across 8 different cities spanning 3 continents, our approach is 4.8× faster than prior state-of-the-art training systems (Megatron).

Code Availability: <https://github.com/DS3Lab/DT-FM>

1 Introduction

Recent years have witnessed the rapid development of deep learning models, particularly foundation models (FMs) [1] such as GPT-3 [2] and PaLM [3]. Along with these rapid advancements, however, comes computational challenges in training these models: the training of these FMs can be very expensive — a single GPT-3-175B training run takes 3.6K Petaflops-days [2] — this amounts to \$4M on today’s AWS on demand instances, even assuming 50% device utilization (V100 GPUs peak at 125 TeraFLOPS!) Even the smaller scale language models, e.g., GPT-3-XL (1.3 billion parameters), on which this paper evaluates, require 64 Tesla V100 GPUs to run for one week, costing \$32K on AWS. As a result, speeding up training and decreasing the cost of FMs have been active research areas. Due to their vast number of model parameters, state-of-the-art systems (e.g., Megatron[4], Deepspeed[5], Fairscale[6]) leverage multiple forms of parallelism [4, 7, 8, 9, 10, 11]. However, their design is only tailored to *fast, homogeneous* data center networks.

* Equal contribution.

[NeurIPS 2022-(a)]

Fine-tuning Language Models over Slow Networks using Activation Compression with Guarantees

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Abstract

Communication compression is a crucial technique for modern distributed learning systems to alleviate their communication bottlenecks over slower networks. Despite recent intensive studies of gradient compression for data parallel style training, compressing the activations for models trained with pipeline parallelism is still an open problem. In this paper, we propose AC-SGD, a novel activation compression algorithm for communication-efficient pipeline parallelism training over slow networks. Different from previous efforts in activation compression, instead of compressing activation values directly, AC-SGD compresses the *changes of the activations*. This allows us to show, to the best of our knowledge for the first time, that one can still achieve $O(1/\sqrt{T})$ convergence rate for non-convex objectives under activation compression, without making assumptions on gradient unbiasedness that do not hold for deep learning models with non-linear activation functions. We then show that AC-SGD can be optimized and implemented efficiently, without additional end-to-end network overhead. Our AC-SGD implementation provides up to 1.5 billion parameters, compressing activations to 2.4 bits, and provides up to 4.3× end-to-end speed-up over slow networks, without sacrificing model quality. Moreover, we also show that AC-SGD can be combined with state-of-the-art gradient compression algorithms to enable “end-to-end communication compression”: All communications between machines, including *model gradients, forward activations, and backward gradients* are compressed into *lower precision*. This provides up to 4.9× end-to-end speed-up, without sacrificing model quality.

Code Availability: <https://github.com/DS3Lab/AC-SGD>

1 Introduction

Recent efforts in improving communication efficiency for distributed learning have significantly decreased the dependency of training deep learning models on fast data center networks — the *gradient* can be compressed to lower precision or sparsified [1, 2, 3, 4], which speeds up training over low bandwidth networks, whereas the *communication topology* can be decentralized [5, 6, 7, 8, 9, 10], which speeds up training over high latency networks. Indeed, today’s state-of-the-art training systems, such as Pytorch [11, 12], Horovod [13], Bagua [14], and BytePS [15], already support many of these communication-efficient training paradigms.

However, with the rise of large foundation models [16] (e.g., BERT [17], GPT-3 [18], and CLIP [19]), improving communication efficiency via compression becomes more challenging. Current training systems for foundation models such as Megatron [20], Deepspeed [21], and FairScale [22], allocate different layers of the model onto multiple devices and need to communicate — *in addition to* the gradients on the models — the

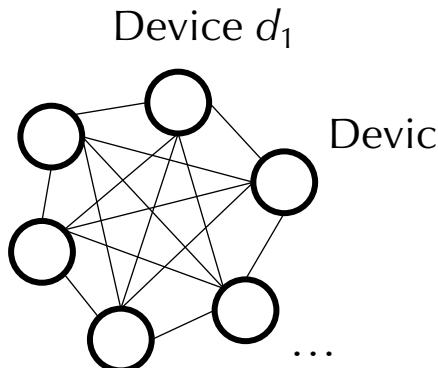
* Equal contribution.

[NeurIPS 2022-(b)]

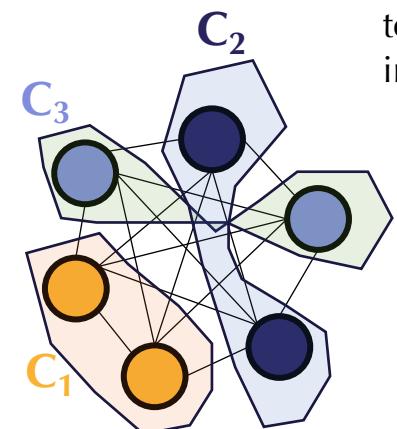
Accommodate Communication in a Decentralized network

A bi-level scheduling algorithm based on an extended balanced graph partition to estimate the communication cost:

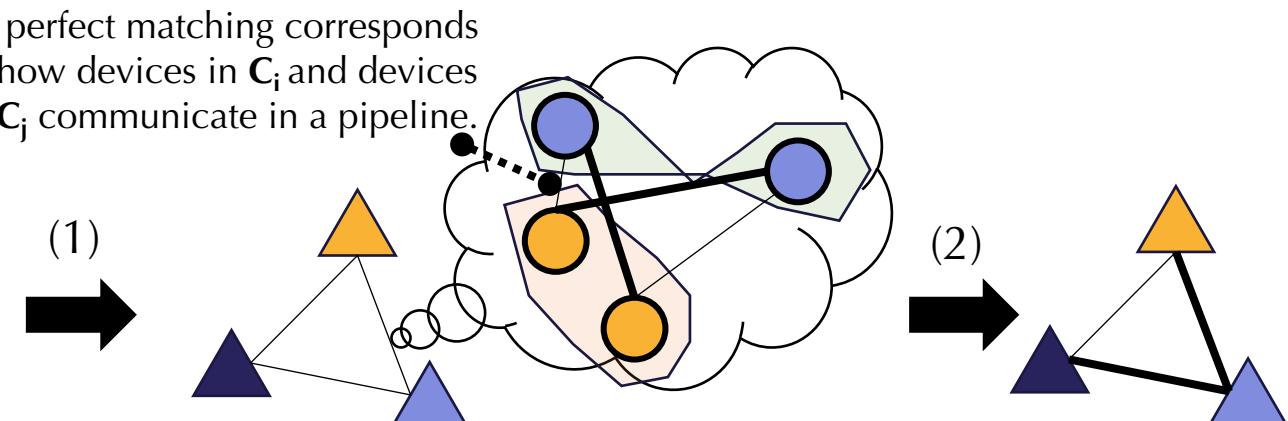
- Data parallel communication cost: nodes handling the same stage need to exchange gradients;
- Pipeline parallel communication cost: nodes handling nearby stages for the same micro-batch need to communicate activation in the forward propagation and gradients of the activation in the backward propagation.



(a) Communication Topology Graph \mathbf{G} over N devices



(b) Each partition \mathbf{C}_i deals with one stage, running data parallel within each partition



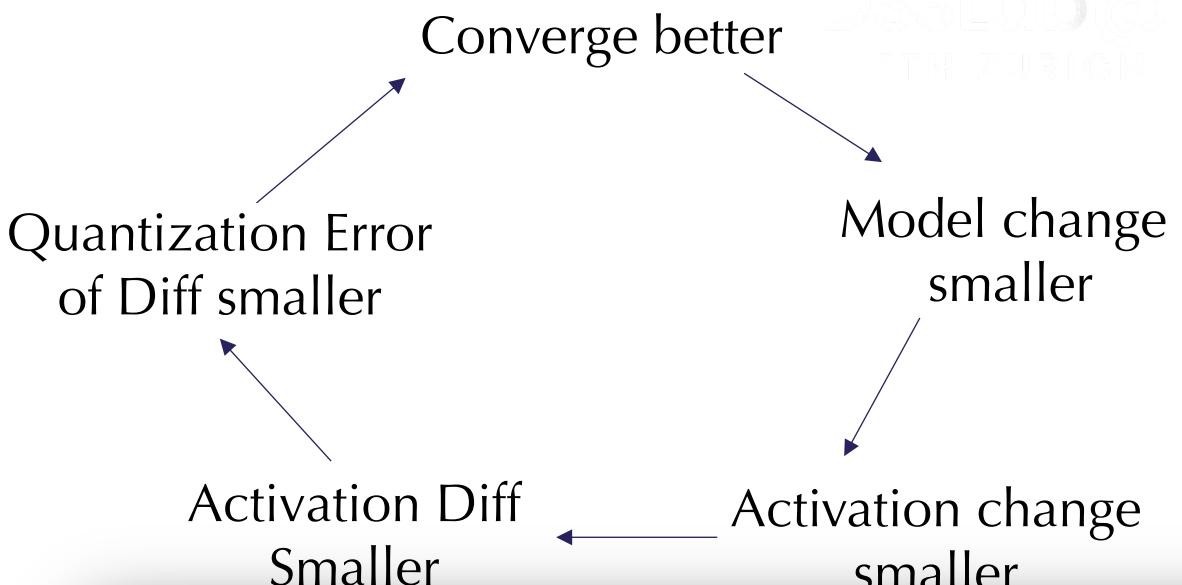
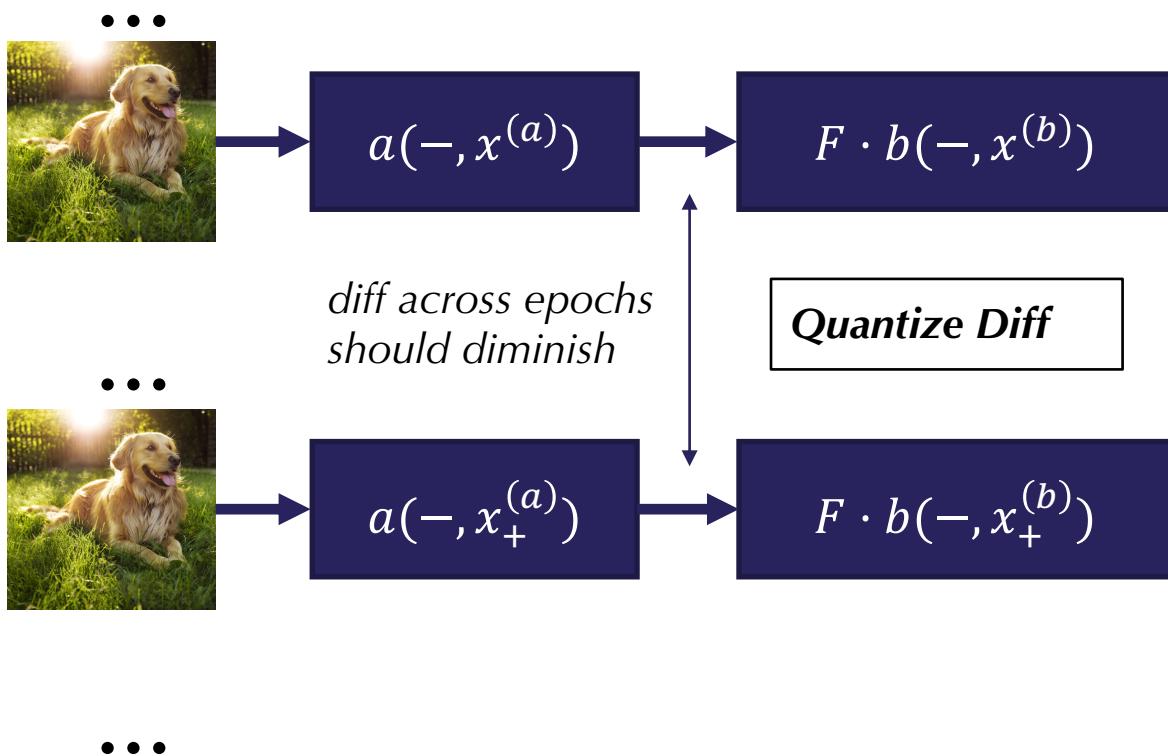
(c) Coarsened graph $\widehat{\mathbf{G}}$ decoding the cost of pipeline parallel

(e) Open-loop-traveling-salesman provides a pipeline structure

AQ-SGD

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

Direct quantization only works to some degree.



- **(A1: Lipschitz assumptions)** We assume that ∇f , $\nabla(f \circ b)$ and a are L_f , $L_{f \circ b}$ -, and ℓ_a -Lipschitz, respectively, recalling that a function g is L_g -Lipschitz if

$$\|g(x) - g(y)\| \leq L_g \|x - y\|, \quad \forall x, \forall y.$$

Furthermore, we assume that a and $f \circ b$ have gradients bounded by C_a and $C_{f \circ b}$, respectively, i.e. $\|\nabla a(x)\| \leq C_a$, and $\|\nabla(f \circ b)(x)\| \leq C_{f \circ b}$.

- **(A2: SGD assumptions)** We assume that the stochastic gradient g_ξ is unbiased, i.e. $\mathbb{E}_\xi[g_\xi(x)] = \nabla f(x)$, for all x , and with bounded variance, i.e. $\mathbb{E}_\xi\|\nabla f(x) - \nabla f(x)\|^2 \leq \sigma^2$, for all x .

Theorem 3.1. Suppose that Assumptions A1, A2 hold, and consider an unbiased quantization function $Q(x)$ which satisfies that there exists $c_Q < \sqrt{1/2}$ such that $\mathbb{E}\|x - Q(x)\| \leq c_Q \|x\|$, for all x .¹ Let $\gamma = \frac{1}{3(C+3L_f)\sqrt{T}}$ be the learning rate, where

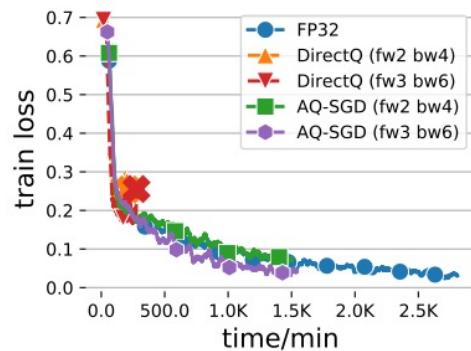
$$C = \frac{4c_Q\ell_a(1+C_a)L_{f \circ b}N}{\sqrt{1-2c_Q^2}}.$$

Then after performing T updates one has

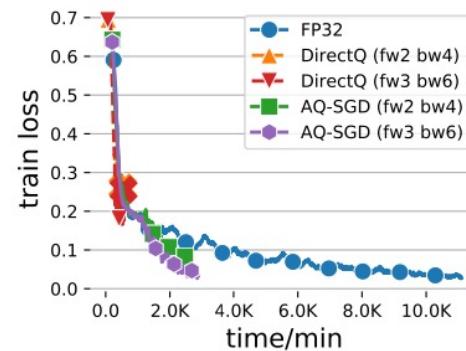
$$\frac{1}{T} \sum_{t \in [T]} \mathbb{E}\|\nabla f(x_t)\|^2 \lesssim \frac{(C+L_f)(f(x_1) - f^*)}{\sqrt{T}} + \frac{\sigma^2 + (c_Q C_a C_{f \circ b})^2}{\sqrt{T}}. \quad (3.1)$$

AQ-SGD Results

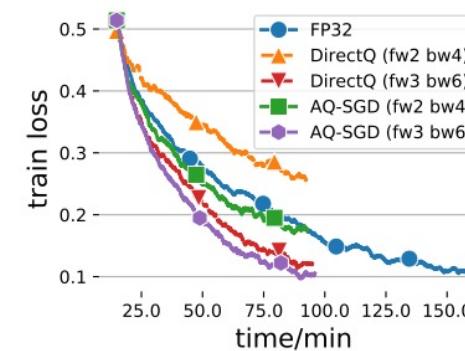
- End-to-end training performance over different networks. x represents divergence.



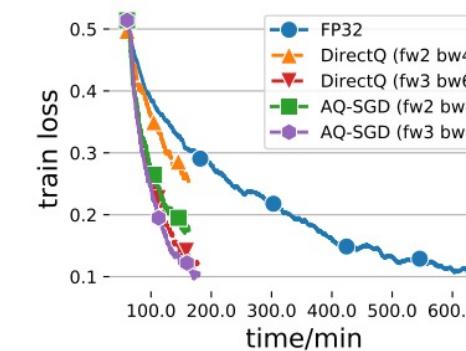
(a) QNLI, 500Mbps



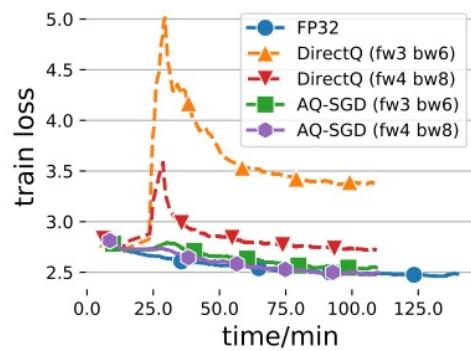
(b) QNLI, 100Mbps



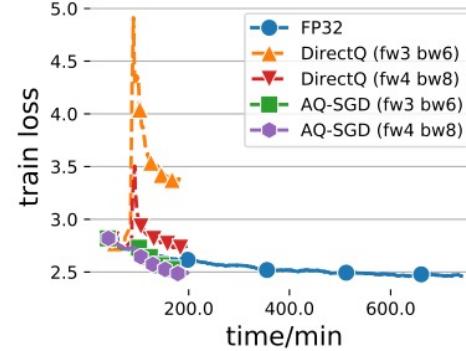
(c) CoLA, 500Mbps



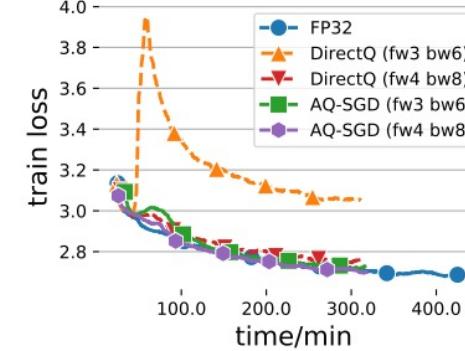
(d) CoLA, 100Mbps



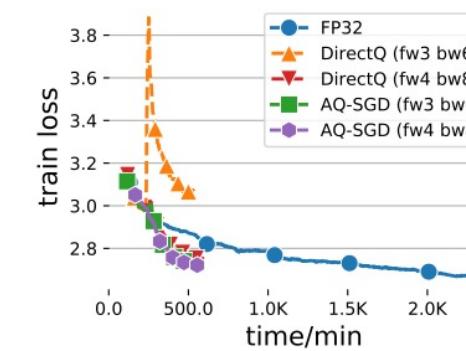
(e) WikiText2, 500Mbps



(f) WikiText2, 100Mbps



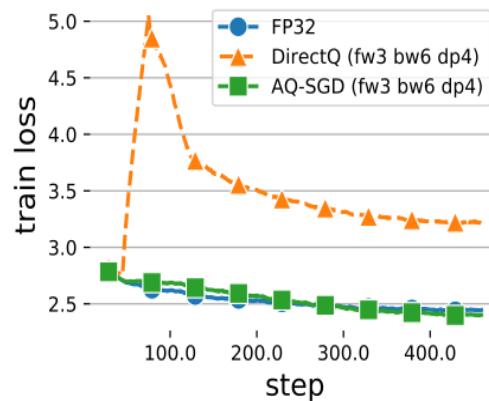
(g) arXiv, 500Mbps



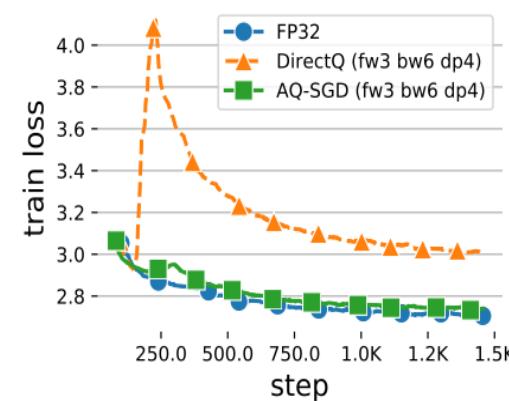
(h) arXiv, 100Mbps

AQ-SGD Results

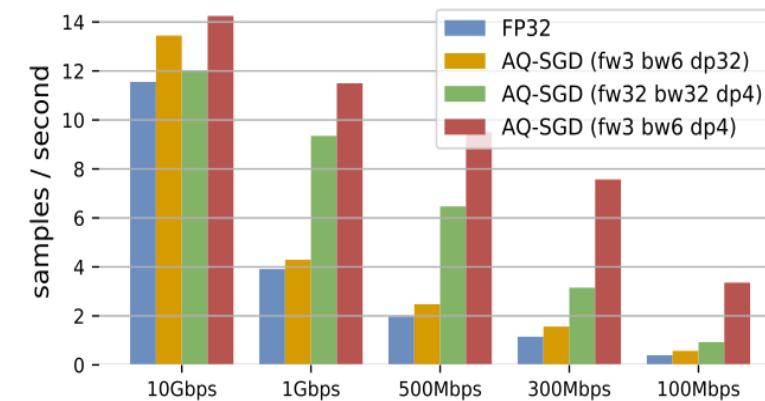
- Convergence and Throughput of AQ-SGD combined with gradient compression.



(a) WikiText2, GPT2-1.5B



(b) arXiv, GPT2-1.5B



(c) Training Throughput

Computes are there but *scattered*, we have made
some small steps towards decentralized ML.

GPT-JT: Instruct Tuned GPT-J (6B) over 1Gbps Network

Data Sources

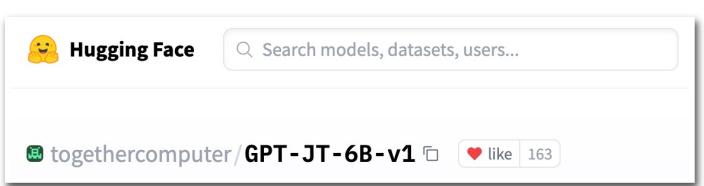
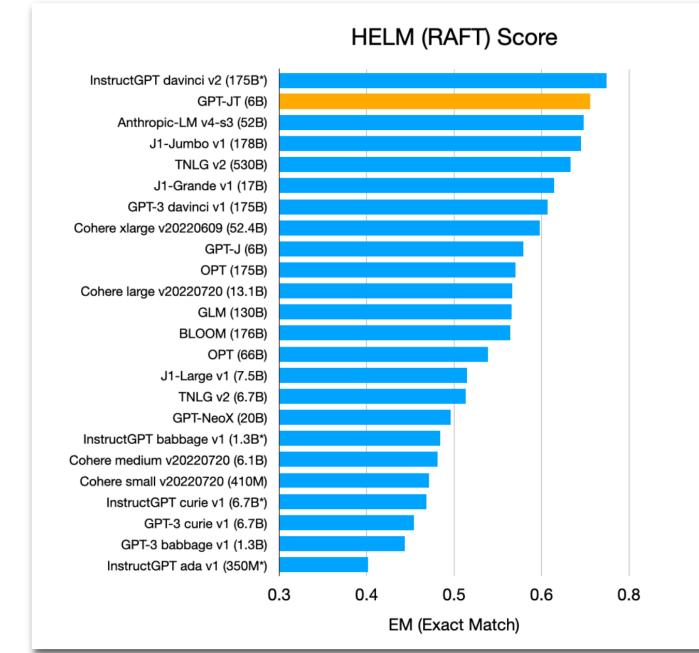
- UL2, Chain of thought
- Natural Instruction
- Public Pool of Prompts (P3)

Model & Training

- GPT-J 6B
- Cocktail SGD

1Gbps network; 4-way data parallel; 2x A100 each

30% end-to-end overhead, compared with
100Gbps data-center network

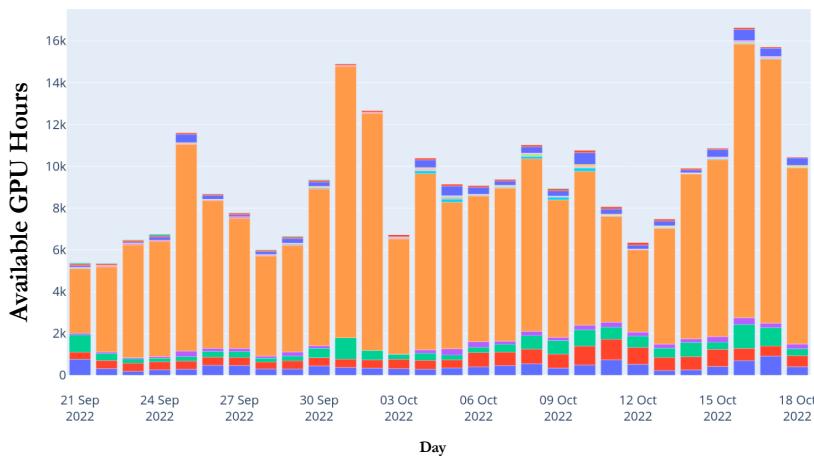


We are able to do useful things over slow networks!

This is just the beginning, give us feedback, and let's train open model together!

Open Research on the Together Decentralized Cloud

Connecting idle computes across academic institutions.



11 billion tokens
60K GPU Hours
10 Open Models

BLOOM	176B	July 2022
T0pp	11B	October 2021
GPT-J	6B	July 2021
GPT-NeoX	20B	February 2022
GLM	130B	August 2022
UL2	20B	October 2022
T5	11B	February 2020
OPT	175B	June 2022
OPT	66B	June 2022
YaLM	100B	June 2022

Summary

- Communication is a key bottleneck of distributed learning, both for centralized data center network and decentralized environments.
- We can develop Algorithms to alleviate communication bottlenecks:
 - Data Parallel: {asynchronous, local training, compression, quantization, decentralized topology} & their combinations.
 - Model Parallel: Careful error compensation.
- Innovation of Systems is need to unleash the full potential Algorithms:
 - Bagua: Automatic optimization framework.
 - System Scheduling of communication in decentralized environments.
- We can build awesome things with it.



Personal page:
<https://binhangyuan.github.io/site/>

Thank you!