Different Aspects of Communication Organization in Distributed and Federated learning Optimization methods

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MIPT

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Modern trends in machine learning

Exponential growth in model sizes and data volumes.

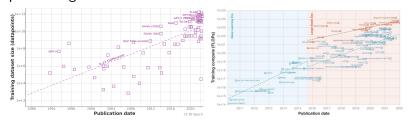


Figure: Trends in machine learning tasks

Varieties of distributed learning

- Cluster learning (big players): training within one large and powerful computing cluster
- Collaborative learning (all players): pooling computing resources over the Internet

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- Cluster learning (big players): training within one large and powerful computing cluster
- Collaborative learning (all players): pooling computing resources over the Internet
- Federated learning (another paradigm): learning from users' local data using their computing powers



Figure: Federated Learning



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Distributed optimization problem

Distributed optimization problem:

$$\min_{x \in \mathbb{R}^d} \left[f(x) := \frac{1}{M} \sum_{m=1}^M f_m(x) = \frac{1}{M} \sum_{m=1}^M \frac{1}{N} \sum_{i=1}^N I(g(x, a_i^m), b_i^m) \right],$$

where x – optimization variables, (a_i^m, b_i^m) – data sample, I – loss function, g - model.

 The problem is divided into M pieces, each node m having access only to the function f_m (and its derivatives).

Output: x_K

Communicating through the server

Let us look at an example of how an ordinary GD becomes distributed.

Algorithm 1 Centralized GD

```
Input: Step size \gamma > 0, starting point x_0 \in \mathbb{R}^d, number iterations K
 1: for k = 0, 1, ..., K - 1 do
        Send x_k to all nodes
                                                                                  by server
 3:
        for m = 1, ..., M in parallel do
             Receive x_k forom server
 4:
                                                                                   by nodes
             Compute gradient \nabla f_m(x_k) at x_k
 5:
                                                                                  by nodes
             Send \nabla f_m(x_k) to server
                                                                                  by nodes
 6:
 7:
        end for
         Receive \nabla f_m(x_k) from nodes
 8:
                                                                                  by server
        Compute \nabla f(x_k) = \frac{1}{M} \sum_{m=1}^{M} \nabla f_m(x_k)
                                                                                  by server
10:
        x_{k+1} = x_k - \gamma \nabla f(x_k)
                                                                                  by server
11: end for
```

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- Q: If we are talking about a federated learning paradigm, what would users want to demand from the learning process?
- Since the data is private and personal, one would like to get its privacy and security.
- The user wants a better result in terms of personalization of learning.
- Q: distributivity is necessary for parallelization, but why can't we achieve full parallelization?
- Communication costs are a waste of time. The problem of communication bottleneck is relevant for all distributed problems.
 There are many ways to fight for fast and effective communication.



6 / 85

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Local updates

The idea – more local computation

- In the basic approach, communications occur every iteration.
- If computing (stochastic) gradients is much cheaper, why not count multiple times between communications.



8 / 85

Parallel SGD/FedAvg/Local SGD

The idea:

• Make local steps (local training):

$$x_{k+1}^m = x_k^m - \gamma \nabla f_m(x_k^m).$$

- Every Tth iteration, forward the current x_k^m to the server. The server averages $x_k = \frac{1}{M} \sum_{m=1}^M x_k^m$, and forwards x_k to the workers. The workers update: $x_k^m = x_k$.
- Centralized distributed SGD is a Local SGD with T=1.
 - Mangasarian O. Parallel Gradient Distribution in Unconstrained Optimization

 McMahan B. et al. Communication Efficient Learning of Doop
 - McMahan B. et al. Communication-Efficient Learning of Deep Networks from Decentralized Data

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How it works

Problems: 1) LSTM on 10 million public posts, 2) CNN on CIFAR-10.

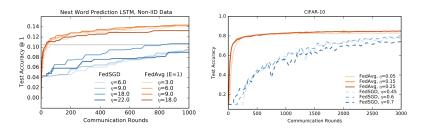


Figure: Comparison of Local SGD (FedAvg) and Centralized Distributed SGD (FedSGD).



McMahan B. et al. Communication-Efficient Learning of Deep Networks from Decentralized Data

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• Problem: logistic regression on a5a LibSVM dataset.

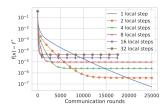


Figure: Dependence of convergence of Local SGD on number of local steps

- Typical convergence of this type of methods: faster in terms of communications, worse quality of ultimate accuracy.
 - Khaled A. et al. Tighter Theory for Local SGD on Identical and Heterogeneous Data

• Q: what causes this effect?

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- In the theoretical estimates of the convergence of the method, this also shows up:

$$\mathcal{O}\left(\frac{\|x^0-x^*\|^2}{\gamma K}+\gamma \cdot \frac{1}{M}\sum_{m=1}^M \|\nabla f_m(x^*)\|^2\right),\,$$

where $\gamma \leq \mathcal{O}\left(\frac{1}{LT}\right)$ – stepsize, K – number of local iterations on each device. The estimation is given for the case of convex and L-smooth f_m .



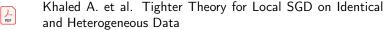
Khaled A. et al. Tighter Theory for Local SGD on Identical and Heterogeneous Data

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• Moreover, the σ_{opt}^2 factor is not eliminated at all.

Glasgow M.R. et al. Sharp bounds for federated averaging (local sgd) and continuous perspective

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Solving the problem

 Q: the problem of the local method is convergence to a neighbourhood. How can it be interpreted and then solved?



Solving the problem

- Q: the problem of the local method is convergence to a neighbourhood. How can it be interpreted and then solved?
- Local task regularization as a defence against local overfitting (FedProx):

$$\tilde{f}_m(x) := f_m(x) + \frac{\lambda}{2} ||x - v||^2,$$

where v – certain reference point.

- Run local iterations not for f_m , but for \tilde{f}_m .
 - PDF L

Li T. et al. Federated Optimization in Heterogeneous Networks

More modifications and generalizations

Using so-called shifts to control bias due to heterogeneity.



Karimireddy S. P. et al. SCAFFOLD: Stochastic Controlled Averaging for Federated Learning



Mishchenko K. et al. ProxSkip: Yes! Local Gradient Steps Provably Lead to Communication Acceleration! Finally!

Using of consensus gossip procedures when centralized communications are unavailable



Beznosikov A. et al. Decentralized Local Stochastic Extra-Gradient for Variational Inequalities

How it works

• Problem: logistic regression on EMNIST (letters).

	Epochs			10% similarity			100% similarity (i.i.d.)		
		Num. of rounds	Speedup	Num. of rounds		Speedup	Num. of rounds		Speedup
SGD	1	317	(1×)	365		(1×)	416		(1×)
SCAFFOL	D1	77 —	(4.1×)	62 -		(5.9×)	60 =		(6.9×)
	5	152	(2.1×)	20 •		(18.2×)	10 •		(41.6×)
	10	286	(1.1×)	16 •		(22.8×)	7 '		$(59.4 \times)$
	20	266	(1.2×)	11 ((33.2×)	4		(104×)
FEDAVG	1	258	(1.2×)	74 -		(4.9×)	83 -		(5×)
	5	428	(0.7×)	34 -		$(10.7 \times)$	10 •		(41.6×)
	10	711	(0.4×)	25 =		(14.6×)	6 -		(69.3×)
	20	1k+	(< 0.3×)	18 •	-1-	$(20.3 \times)$	4	- 1	(104×)
FEDPROX	1	1k+	(< 0.3×)	979	\rightarrow	(0.4×)	459	-	(0.9×)
	5	1k+	(< 0.3×)	794	\rightarrow	(0.5×)	351	-	(1.2×)
	10	1k+	(< 0.3×)	894	\rightarrow	$(0.4 \times)$	308	_	$(1.4 \times)$
	20	1k+	(< 0.3×)	916	\rightarrow	(0.4×)	351	_	(1.2×)

Figure: Comparison of Local SGD (FedAvg), FedProx and SCAFFOLD and Centralized Distributed SGD.



Karimireddy S. P. et al. SCAFFOLD: Stochastic Controlled Averaging for Federated Learning

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What do we want to achieve, anyway?

Lower bounds:

$$K = \Omega\left(\sqrt{\frac{L}{\varepsilon}}\right).$$

- L smoothness constant of f.
- What method will give such estimates?

What do we want to achieve, anyway?

Lower bounds:

$$K = \Omega\left(\sqrt{\frac{L}{\varepsilon}}\right).$$

L – smoothness constant of f.

- What method will give such estimates? Distributed version of Nesterov's accelerated method with 1 local step between communications.
- Note that local methods were invented for stochastic setups. But even here there is no improvement in the general case:
 - Woodworth B. The Min-Max Complexity of Distributed Stochastic Convex Optimization with Intermittent Communication
- But there are settings where localised methods shoot out.

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Data similarity

Distributed learning problem:

$$f(x) = \frac{1}{M} \sum_{m=1}^{M} f_m(x) = \frac{1}{M} \sum_{m=1}^{M} \left[\frac{1}{N} \sum_{i=1}^{N} \ell(x, z_i^m) \right],$$

where z_i^m – data sample (a_i^m, b_i^m) , ℓ – loss of model with weights xon sample z_i^m .

Data similarity

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where z_i^m – data sample (a_i^m, b_i^m) , ℓ – loss of model with weights x on sample z_i^m .

- Suppose we can partition the training data uniformly across devices.
 E.g., if cluster or collaborative computing on open data is used. In fact, we will further understand that it's enough to put a large uniform sample on just one device.
- This gives the similarity of the local loss functions.
- It is asserted that for any x:

$$\|\nabla^2 f_m(x) - \nabla^2 f(x)\| \le \delta.$$

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Theorem (Matrix Hoeffding)

Consider a finite sequence of random square matrices $\{X_i\}_{i=1}^N$. Let the matrices in this sequence be independent, Hermitian and of dimension d. Suppose also that $\mathbb{E}[X_i] = 0$, and $X_i^2 \leq A^2$ is almost surely, where A is a non-random Hermitian matrix. Then with probability 1 - p it is satisfied that

$$\left\|\sum_{i=1}^N X_i\right\| \leq \sqrt{8N\|A^2\| \cdot \ln\left(d/p\right)}.$$



Tropp J. An introduction to matrix concentration inequalities Tropp J. User-friendly tail bounds for sums of random matrices

Similarity parameter

Local loss function:

$$f_m(x) = \frac{1}{N} \sum_{i=1}^N \ell(x, z_i).$$

• ℓ – L-smooth (L-Lipschitz gradient), convex, twice differentiable function (e.g., quadratic or logreg). Then we have $\nabla^2 \ell(x, z_i) \leq LI$ for any x and z_i (here I is a unit matrix).

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- Let us divide all data uniformly over all workers. $X_i = \frac{1}{N} \left[\nabla^2 \ell(x, z_i) \nabla^2 f(x) \right]$. It is easy to check that all conditions of Hoeffding inequality are satisfied for it, in particular, $A^2 = \frac{4L^2}{N^2}I$.

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19 / 85

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Similarity parameter: summary

As a result, we have

$$\|\nabla^2 f_m(x) - \nabla^2 f(x)\| \le \delta \sim \frac{L}{\sqrt{N}}.$$

 Conclusion: the larger the local sample size, the smaller the similarity parameter (hessians are similar to each other).

Method in general terms (not only for similarity)

Consider Mirror Descent:

$$x_{k+1} = \arg\min_{x \in \mathbb{R}^d} \left(\gamma \langle \nabla f(x_k), x \rangle + V(x, x_k) \right),$$

where V(x, y) is the Bregman divergence generated by the strictly convex function $\varphi(x)$:

$$V(x,y) = \varphi(x) - \varphi(y) - \langle \nabla \varphi(y); x - y \rangle.$$

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• Q: Which method do we have if $\varphi(x) = \frac{1}{2} ||x||^2$?

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• Q: Which method do we have if $\varphi(x) = \frac{1}{2} ||x||^2$? Gradient descent.

Convergence in general terms

Definition (Relative smoothness and strong convexity)

Let $\varphi: \mathbb{R}^d \to \mathbb{R}$ is convex and twice differentiable. Let us say that the function f is L_{φ} -smooth and μ_{φ} -strongly convex with respect to φ if for any $x \in \mathbb{R}^d$ the following holds

$$\mu_{\varphi} \nabla^2 \varphi(x) \preceq \nabla^2 f(x) \preceq L_{\varphi} \nabla^2 \varphi(x),$$

or equivalently for any $x, y \in \mathbb{R}^d$

$$\mu_{\varphi}V(x,y) \leq f(x) - f(y) - \langle \nabla f(y); x - y \rangle \leq L_{\varphi}V(x,y).$$



Lu H. et al. Relatively-Smooth Convex Optimization by First-Order Methods, and Applications

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Convergence in general terms: proof

The optimality condition for the Mirror Descent step:

$$\gamma \nabla f(x_k) + \nabla \varphi(x_{k+1}) - \nabla \varphi(x_k) = 0.$$

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From it (here x* – optimal point):

$$\langle \gamma \nabla f(x_k) + \nabla \varphi(x_{k+1}) - \nabla \varphi(x_k), x_{k+1} - x^* \rangle = 0.$$

$$\langle \gamma \nabla f(x_k), x^{k+1} - x^* \rangle = \langle \nabla \varphi(x_k) - \nabla \varphi(x_{k+1}), x^{k+1} - x^* \rangle$$

= $V(x^*, x_k) - V(x^*, x_{k+1}) - V(x_{k+1}, x_k).$

(the last statement is called the Pythagoras' theorem for the Bregman divergence and is verified by the definition)

23 / 85

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Small permutations give:

$$\langle \gamma \nabla f(x_k), x_{k+1} - x_k \rangle + V(x_{k+1}, x_k)$$

$$= V(x^*, x_k) - V(x^*, x_{k+1}) - \langle \gamma \nabla f(x_k), x_k - x^* \rangle.$$

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• Substitute $\gamma = \frac{1}{L_0}$:

$$\langle \nabla f(x_k), x_{k+1} - x_k \rangle + L_{\varphi} V(x_{k+1}, x_k)$$

$$= L_{\varphi} V(x^*, x_k) - L_{\varphi} V(x^*, x_{k+1})$$

$$- \langle \nabla f(x_k), x_k - x^* \rangle.$$

• Substitute $\gamma = \frac{1}{L_{\varphi}}$:

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• Let us use the definition of smoothness with respect to φ c $x = x_{k+1}$, $y = x_k$:

$$f(x_{k+1}) - f(x_k) \leq \langle \nabla f(x_k); x_{k+1} - x_k \rangle + L_{\varphi} V(x_{k+1}, x_k).$$

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• Let us use the definition of smoothness with respect to φ c $x=x_{k+1}$, $y=x_k$:

$$f(x_{k+1}) - f(x_k) \leq \langle \nabla f(x_k); x_{k+1} - x_k \rangle + L_{\varphi} V(x_{k+1}, x_k).$$

Combine the previous two:

$$f(x_{k+1})-f(x_k)\leq L_{\varphi}V(x^*,x_k)-L_{\varphi}V(x^*,x_{k+1})-\langle \nabla f(x_k),x_k-x^*\rangle.$$

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From the previous slide:

$$f(x_{k+1}) - f(x_k) \leq L_{\varphi}V(x^*, x_k) - L_{\varphi}V(x^*, x_{k+1}) - \langle \nabla f(x_k), x_k - x^* \rangle.$$

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Relative strong convexity:

$$\mu_{\varphi}V(x^*,x_k) \leq f(x^*) - f(x_k) - \langle \nabla f(x_k); x^* - x_k \rangle$$

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Relative strong convexity:

$$\mu_{\varphi}V(x^*,x_k) \leq f(x^*) - f(x_k) - \langle \nabla f(x_k); x^* - x_k \rangle$$

Sum the previous two together and shuffle them a bit:

$$f(x_{k+1}) - f(x^*) \le (L_{\varphi} - \mu_{\varphi})V(x^*, x_k) - L_{\varphi}V(x^*, x_{k+1}).$$

From the previous slide:

$$f(x_{k+1}) - f(x_k) \leq L_{\varphi}V(x^*, x_k) - L_{\varphi}V(x^*, x_{k+1}) - \langle \nabla f(x_k), x_k - x^* \rangle.$$

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Sum the previous two together and shuffle them a bit:

$$f(x_{k+1}) - f(x^*) \le (L_{\varphi} - \mu_{\varphi})V(x^*, x_k) - L_{\varphi}V(x^*, x_{k+1}).$$

By virtue of the fact that x* – optimum:

$$V(x^*, x_{k+1}) \leq \left(1 - \frac{\mu_{\varphi}}{L_{\varphi}}\right) V(x^*, x_k).$$

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Theorem (Convergence of Mirror Descent)

Let φ and f satisfy the definition above, then Mirror Descent with step $\gamma = \frac{1}{L_0}$ converges and is satisfied:

$$V(x^*, x_K) \leq \left(1 - \frac{\mu_{\varphi}}{L_{\varphi}}\right)^K V(x^*, x_0).$$

Mirror Descent:

$$x_{k+1} = \arg\min_{x \in \mathbb{R}^d} \left(\gamma \langle \nabla f(x_k), x \rangle + V(x, x_k) \right),$$

where the Bregman divergence of V(x, y) generated by the function $\varphi(x)$ (here we need to require that f_1 is convex):

$$\varphi(x) = f_1(x) + \frac{\delta}{2} ||x||^2.$$

The function f_1 is stored on the server.

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 Q: What is the number of communications that occur in K iterations of Mirror Descent?

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• Q: What is the number of communications that occur in K iterations of Mirror Descent? K of communications (number of ∇f gradient counts), computing arg min requires only computations on the server.

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Algorithm 2 Mirror Descent for the data similarity problem

Input: Stepsize $\gamma > 0$, starting point $x_0 \in \mathbb{R}^d$, number iterations K

- 1: **for** k = 0, 1, ..., K 1 **do**
- Send x_k to all workers
- for m = 1, ..., M in parallel do 3:
- Receive x_k from server 4:
- 5: Compute gradient $\nabla f_m(x_k)$ at x_k
- 6: Send $\nabla f_m(x_k)$ to server
- end for 7:
- Receive $\nabla f_m(x_k)$ from all workers 8:
- Compute $\nabla f(x_k) = \frac{1}{M} \sum_{m=1}^{M} \nabla f_m(x_k)$ 9.
- $x_{k+1} = \operatorname{arg\,min}_{x \in \mathbb{R}^d} \left(\gamma \langle \nabla f(x_k), x \rangle + V(x, x_k) \right)$ 10:
- 11: end for

server

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server

Recall that convergence is defined in terms of constants from the relation:

$$\mu_{\varphi} \nabla^2 \varphi(x) \leq \nabla^2 f(x) \leq L_{\varphi} \nabla^2 \varphi(x),$$

Recall that convergence is defined in terms of constants from the relation:

$$\mu_{\varphi} \nabla^2 \varphi(x) \preceq \nabla^2 f(x) \preceq L_{\varphi} \nabla^2 \varphi(x),$$

In our case.

$$\mu_{\varphi}\left(\delta I + \nabla^2 f_1(x)\right) \leq \nabla^2 f(x) \leq L_{\varphi}\left(\delta I + \nabla^2 f_1(x)\right)$$

 Recall that convergence is defined in terms of constants from the relation:

$$\mu_{\varphi} \nabla^2 \varphi(x) \leq \nabla^2 f(x) \leq L_{\varphi} \nabla^2 \varphi(x),$$

• In our case:

$$\mu_{\varphi}\left(\delta I + \nabla^2 f_1(x)\right) \leq \nabla^2 f(x) \leq L_{\varphi}\left(\delta I + \nabla^2 f_1(x)\right)$$

• Let us find L_{φ} :

$$\|\nabla^{2} f_{1}(x) - \nabla^{2} f(x)\| \leq \delta \Rightarrow \nabla^{2} f(x) - \nabla^{2} f_{1}(x) \leq \delta I$$

$$\Rightarrow \nabla^{2} f(x) \leq \delta I + \nabla^{2} f_{1}(x) \Rightarrow \boxed{L_{\varphi} = 1.}$$

Aleksandr Beznosikov Lecture 29 April 2025 29 / 85

• Let us find μ_{φ} . From the strong convexity of f:

$$\mu I \preceq \nabla^2 f(x) \Rightarrow \delta I \preceq \frac{2\delta}{\mu} \nabla^2 f(x) - \delta I.$$

• Let us find μ_{ω} . From the strong convexity of f:

$$\mu I \preceq \nabla^2 f(x) \Rightarrow \delta I \preceq \frac{2\delta}{\mu} \nabla^2 f(x) - \delta I.$$

• From $\|\nabla^2 f_1(x) - \nabla^2 f(x)\| \le \delta$ we have:

$$\nabla^2 f_1(x) - \nabla^2 f(x) \leq \delta I.$$

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$$\nabla^2 f_1(x) - \nabla^2 f(x) \leq \frac{2\delta}{\mu} \nabla^2 f(x) - \delta I.$$

And we get

$$\nabla^2 f_1(x) + \delta I \leq \frac{2\delta + \mu}{\mu} \nabla^2 f(x) \Rightarrow \boxed{\mu_{\varphi} = \frac{\mu}{2\delta + \mu}}.$$

Aleksandr Beznosikov Lecture 29 April 2025 30 / 85

Theorem (Convergence for the data similarity problem)

Let f be strongly convex, f_1 be convex, and ℓ be smooth, and $\varphi(x)=f_1(x)+\frac{\delta}{2}\|x\|^2$, then Mirror Descent with step $\gamma=1$ converges and is satisfied:

$$V(x^*,x_K) \leq \left(1 - \frac{\mu}{\mu + 2\delta}\right)^K V(x^*,x_0).$$

Theorem (Convergence for the data similarity problem)

Let f be strongly convex, f_1 be convex, and ℓ be smooth, and $\varphi(x)=f_1(x)+\frac{\delta}{2}\|x\|^2$, then Mirror Descent with step $\gamma=1$ converges and is satisfied:

$$V(x^*, x_K) \leq \left(1 - \frac{\mu}{\mu + 2\delta}\right)^K V(x^*, x_0).$$

• This means that if we want to achieve an accuracy ε ($V(x^*, x_K) \sim \varepsilon$), then we need to

$$\mathcal{K} = \mathcal{O}\left(\left[1 + rac{\delta}{\mu}
ight]\lograc{V(x^*, x_0)}{arepsilon}
ight) \ ext{communications}.$$



Hendrikx H. et al. Statistically Preconditioned Accelerated Gradient Method for Distributed Optimization

Aleksandr Beznosikov Lecture 29 April 2025 31 / 85

How it works

Problem: ResNet-18 on CIFAR-10.

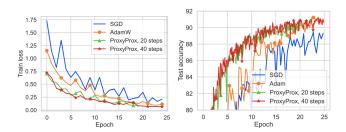


Figure: Comparison of Mirror Descent (ProxyProx) with SOTA optimizers on **non-distributed** problems.



Woodworth B. et al. Two Losses Are Better Than One: Faster Optimization Using a Cheaper Proxy

Research question: acceleration

• Estimate on the number of communications under data similarity:

$$\mathcal{K} = \mathcal{O}\left(\left[1 + \frac{\delta}{\mu}\right] \log \frac{1}{arepsilon}\right).$$

 Estimate on the number of communications for Centralized Distributed Gradient Descent:

$$\mathcal{K} = \mathcal{O}\left(\frac{L}{\mu}\log\frac{1}{arepsilon}
ight).$$

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 Estimate on the number of communications for Centralized Distributed Gradient Descent:

$$\mathcal{K} = \mathcal{O}\left(\frac{L}{\mu}\log\frac{1}{\varepsilon}\right).$$

- Given that δ can be small, we see the improvement.
- But there's also the distributed version of Accelerated Gradient Method that gives estimate:

$$K = \mathcal{O}\left(\sqrt{\frac{L}{\mu}}\log{\frac{1}{\varepsilon}}\right).$$

• It is not clear which is better. Is it possible to accelerate the method for the problem with data similarity?

Another look at Mirror Descent

• Mirror Descent with $\gamma = 1$:

$$x_{k+1} = \arg\min_{x \in \mathbb{R}^d} \left(\left\langle
abla f(x_k), x \right
angle + V(x, x_k) \right),$$

with the Bregman divergence V(x,y), generated by the function $\varphi(x)$:

$$\varphi(x) = f_1(x) + \frac{\delta}{2} ||x||^2.$$

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• Substitute
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:
$$\varphi(x) = f_1(x) + \frac{\delta}{2} ||x||^2.$$
$$x_{k+1} = \arg\min_{x \in \mathbb{R}^d} \left(f_1(x) + \langle \nabla f(x_k) - \nabla f_1(x_k), x \rangle + \frac{\delta}{2} ||x - x_k||^2 \right).$$

Or a little differently:

$$x_{k+1} = \arg\min_{x \in \mathbb{R}^d} \left(f_1(x) + \frac{\delta}{2} \left\| x - \left(x_k - \frac{1}{\delta} (\nabla f(x_k) - \nabla f_1(x_k)) \right) \right\|^2 \right).$$

Aleksandr Beznosikov Lecture 34 / 85

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Q: What method does it look like?

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34 / 85

Aleksandr Beznosikov Lecture 29 April 2025

Proximal Gradient Method for the composite problem $g_1(x) + g_2(x)$:

$$x_{k+1} = \arg\min_{x \in \mathbb{R}^d} \left(\gamma g_2(x) + \frac{1}{2} \|x - (x_k - \gamma g_1(x_k))\|^2 \right).$$

Another look at Mirror Descent

• Proximal Gradient Method for the composite problem $g_1(x) + g_2(x)$:

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• The argmin problem at each iteration can be solved inexactly somehow by a numerical method (e.g., by Gradient Descent or Nesterov's method). The peculiarity of such a method is that ∇g_1 is called much less frequently than ∇g_2 . This kind of algorithms for composite problems that devide oracle complexities are sometimes called slidings.



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- Usually, this kind of algorithms are proposed for composite problems of the form: convex + convex.
- In our case, $g_1 = f f_1$, $g_2 = f_1$. And this is the problem of the form: non-convex + convex = convex.

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Acceleration and optimal algorithm

• Long history:

Reference	Communication complexity	Local gradient complexity	Order	Limitations
DANE [42]	$O\left(\frac{\delta^2}{\mu^2}\log\frac{1}{\varepsilon}\right)$	(2)	1st	quadratic
DiSCO [51]	$\mathcal{O}\left(\sqrt{\frac{\delta}{\mu}}(\log \frac{1}{\varepsilon} + C^2 \Delta F_0)\log \frac{L}{\mu}\right)$	$\mathcal{O}\left(\sqrt{\frac{\delta}{\mu}}(\log \frac{1}{\varepsilon} + C^2 \Delta F_0)\log \frac{L}{\mu}\right)$	2nd	C - self-concordant $^{(3)}$
AIDE [40]	$\mathcal{O}\left(\sqrt{\frac{\delta}{\mu}}\log\frac{1}{\varepsilon}\log\frac{L}{\delta}\right)$	$O\left(\sqrt{\frac{\delta}{\mu}}\sqrt{\frac{L}{\mu}}\log \frac{1}{\varepsilon}\log \frac{L}{\delta}\right)^{(4)}$	1st	quadratic
DANE-LS [50]	$O\left(\frac{\delta}{\mu}\log\frac{1}{\epsilon}\right)$	$O\left(\sqrt{\frac{L}{\mu}} \frac{\delta^{3/2}}{\mu^{3/2}} \log \frac{1}{\epsilon}\right)^{(5)}$	1st/2nd	quadratic (6)
DANE-HB [50]	$O\left(\sqrt{\frac{\delta}{\mu}}\log\frac{1}{\varepsilon}\right)$	$O\left(\sqrt{\frac{L}{\mu}} \frac{\delta}{\mu} \log \frac{1}{\epsilon}\right)^{(5)}$	1st/2nd	quadratic (6)
SONATA [45]	$O\left(\frac{\delta}{\mu}\log\frac{1}{\varepsilon}\right)$	(2)	1st	decentralized
SPAG [21]	$O\left(\sqrt{\frac{L}{\mu}}\log \frac{1}{\varepsilon}\right)^{(1)}$	(2)	1st	M - Lipshitz hessian
DiRegINA [12]	$O\left(\frac{\delta}{\mu}\log\frac{1}{\varepsilon} + \sqrt{\frac{M\delta R_0}{\mu}}\right)$	_(2)	2nd	M -Lipshitz hessian
ACN [1]	$O\left(\sqrt{\frac{\delta}{\mu}}\log\frac{1}{\epsilon} + \sqrt[3]{\frac{M\delta R_0}{\mu}}\right)$	_(2)	2nd	M -Lipshitz hessian
Acc SONATA [46]	$O\left(\sqrt{\frac{\delta}{\mu}}\log\frac{1}{\varepsilon}\log\frac{\delta}{\mu}\right)$	(2)	1st	decentralized
This paper	$O\left(\sqrt{\frac{\delta}{\mu}}\log\frac{1}{\varepsilon}\right)$	$O\left(\sqrt{\frac{L}{\mu}}\log \frac{1}{\varepsilon}\right)$	1st	

Optimal algorithm

For the problem:

$$f(x) = g_1(x) + g_2(x),$$

where $g_1 = f - f_1$ in $g_2 = f_1$.

Algorithm 3 Accelerated Extragradient

Input: Stepsizes γ and θ , momentums α, τ , starting point $x_0 = x_0^f \in \mathbb{R}^d$, number of iterations K

- 1: for $k = 0, 1, 2, \dots, K 1$ do
- $\begin{aligned} x_k^{\mathcal{g}} &= \tau x_k + (1 \tau) x_k^f \\ x_{k+1}^f &\approx \arg\min_{x \in \mathbb{R}^d} \left[\langle \nabla g_1(x_k^{\mathcal{g}}), x x_k^{\mathcal{g}} \rangle + \frac{1}{2\theta} \|x x_k^{\mathcal{g}}\|^2 + g_2(x) \right] \end{aligned}$ 3:
- $x_{k+1} = x_k + \eta \alpha (x_{k+1}^f x_k) \eta \nabla f(x_{k+1}^f)$ 4.
- 5: end for



Kovalev D. et al. Optimal Gradient Sliding and its Application to Distributed Optimization Under Similarity

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- 1 idea Nesterov acceleration.
- 2 idea Sliding.
- 3 idea Extragradient/Tseng's method = method for VIs.
- first two ideas are clear, but the third idea is the key.



How it works

• Problem: logistic regression on different data.

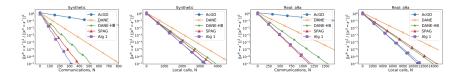


Figure: Comparison of Accelerated ExtraGradient (Alg. 1) with SOTA optimizers for distributed minimization problems under data similarity.

Compression

<u>Unbiased</u> compression (quantization)

Unbiased compression (quantization)

Let us call the stochastic operator Q(x) an unbiased compression (quantization) operator if for any $x \in \mathbb{R}^d$ it is fulfilled:

$$\mathbb{E}[\mathcal{Q}(x)] = x, \quad \mathbb{E}[\|\mathcal{Q}(x)\|_2^2] \le \omega \|x\|_2^2,$$

where $\omega > 1$.

Random sparsification (selection of random components)

Consider a stochastic operator

$$Randk(x) = \frac{d}{k} \sum_{i \in S} [x]_i e_i,$$

where k — some fixed number from the set $\{1,\ldots,d\}$ (the number of components of vector x that we pass; for example, we can choose k=1), S —a random subset of the set $\{1,\ldots,d\}$ of size k (the subset S is chosen randomly and equally likely among all possible subsets of size d), $[\cdot]_i - i$ -th component of the vector, (e_1,\ldots,e_d) —the standard basis in \mathbb{R}^d .



Richtárik P. and Takáč M. Parallel coordinate descent methods for big data optimization

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Richtárik P. and Takáč M. Parallel coordinate descent methods for big data optimization

• Q: why do we need the multiplier $\frac{d}{L}$?

42 / 85

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Richtárik P. and Takáč M. Parallel coordinate descent methods for big data optimization

• Q: why do we need the multiplier $\frac{d}{k}$? For unbiasedness.

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• **Q**: What is ω for random sparsification?

• **Q**: What is ω for random sparsification? $\frac{d}{\nu}$. Each coordinate takes part in Q(x) with probability $\frac{k}{d}$, so

$$\mathbb{E}\left[\|\mathcal{Q}(x)\|^2\right] = \mathbb{E}\left[\sum_{i=1}^d [\mathcal{Q}(x)]_i^2\right]$$
$$= \frac{d^2}{k^2} \left[\sum_{i=1}^d \frac{k}{d} [x]_i^2\right]$$
$$= \frac{d}{k} \|x\|^2.$$

Here $[\cdot]_i$ – *i*th coordinate of the vector.



Three-level ℓ_2 -quantization

Consider the following operator: $[\mathcal{Q}(x)]_i = ||x||_2 \operatorname{sign}(x_i) \xi_i, \ i = 1, \ldots, d,$ where $[\cdot]_i$ is the i-th component of the vector, and ξ_i —a random variable having a Bernoulli distribution with parameter $\frac{|x_i|}{||x||_2}$, i.e.

$$\xi_i = \begin{cases} 1 & \text{with probability } \frac{|x_i|}{\|x\|_2}, \\ 0 & \text{with probability } 1 - \frac{|x_i|}{\|x\|_2}. \end{cases}$$

Thus, if we want to pass a vector $\mathcal{Q}(x)$, we need to pass a vector consisting of zeros and ± 1 , and a real number $\|x\|_2$, with the probability of zeroing a component the greater the component is smaller modulo it. It can be shown that this operator is an unbiased compression with constant $\omega = \sqrt{d}$.

PDF

Alistarh D. et al. QSGD: Communication-Efficient SGD via Gradient Quantization and Encoding

Aleksandr Beznosikov Lecture 29 April 2025 44 / 85

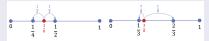
- Q: Is rounding be an unbiased operator?
- Q: Which rounding seems to be the most natural for calculations on a computer?

Natural compression (random rounding to the power of two)

Consider the following operator:

$$[\mathcal{Q}(x)]_i = \begin{cases} \lfloor [x]_i \rfloor_2, & \text{with probabilty } p = \frac{[x]_i - \lceil [x]_i \rceil_2}{\lceil [x]_i \rceil_2 - \lfloor [x]_i \rfloor_2} \\ \lceil [x]_i \rceil_2, & \text{with probability } 1 - p \end{cases}$$

where $[\cdot]_i - i$ -vector component, $[\cdot]_2$ — is the nearest degree of two from the bottom, $[\cdot]_2$ is the nearest degree of two from the top. We round to the two nearest powers of two, the probability of rounding is greater the closer the real number is to the corresponding power of two. It can be shown that $\omega = \frac{9}{8}$.





46 / 85

) PDF Horváth S. et al. Natural compression for distributed deep learning

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Unbiased compression: an idea

 The simplest idea that comes to mind is to use parallel GD, but apply unbiased compression to the gradients sent from the workers to the server.

▷ bv server

by workers

by workers

by workers

Quantized GD (QGD)

Algorithm 1 QGD

Input: step size $\gamma > 0$, starting point $x_0 \in \mathbb{R}^d$, number of iterations K

1: **for**
$$k = 0, 1, ..., K - 1$$
 do

2: Send
$$x_k$$
 to all workers

3: **for**
$$m = 1, ..., M$$
 in parallel **do**

4: Recieve
$$x_k$$
 from server

5: Compute
$$\nabla f_m(x_k)$$
 in point x_k

6: Independently generate
$$g_{k,m} = \mathcal{Q}(\nabla f_m(x_k))$$

7: Send
$$g_{k,m}$$
 to master

10: Compute
$$g_k = \frac{1}{M} \sum_{m=1}^{M} g_{k,m}$$

11:
$$x_{k+1} = x_k - \gamma g_k$$

12: end for Output:
$$x_K$$

7: Send
$$g_{k,m}$$
 to master \triangleright by workers
8: **end for**
9: Recieve $g_{k,m}$ from all workers \triangleright by server
10: Compute $g_k = \frac{1}{M} \sum_{m=1}^{M} g_{k,m}$ \triangleright by server
11: $x_{k+1} = x_k - \gamma g_k$ \triangleright by server

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Theorem (QGD)

Let all local functions f_m be μ -simply convex and have L-Lipschitz gradient, then if $\gamma \leq L^{-1} \left(\frac{2\omega}{M} + 1\right)^{-1}$ then

$$\mathbb{E}\left[\|x_{K}-x^{*}\|^{2}\right] = \mathcal{O}\left((1-\gamma\mu)^{K}\|x_{0}-x^{*}\|^{2}+\gamma\cdot\frac{2\omega}{\mu M^{2}}\sum_{m=1}^{M}\|\nabla f_{m}(x^{*})\|^{2}\right).$$

The selection of γ from the paper was also used to obtain this result:



Stich S. Unified Optimal Analysis of the (Stochastic) Gradient Method

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- Q: what are the problems with this estimation? (recall the convergence estimate GD)

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The selection of γ from the paper was also used to obtain this result:

- Stich S. Unified Optimal Analysis of the (Stochastic) Gradient Method
- Q: what are the problems with this estimation? (recall the convergence estimate GD) Sublinear convergence (depends on the heterogeneity of the data).

49 / 85

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Behaviour in practice:

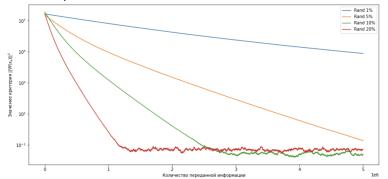


Figure: Behaviour of methods with unbiased compression operator and constant step size

In theory the pitch was chosen cleverly, with constant step the theory predicts exactly the same effect - early plateauing.

Unbiased compression: solving the plateau problem

Method DIANA – QGD with memory:

Algorithm 1 DIANA (sketch)

- 1: Each device m possesses a vector of "memory" $h_0^m = 0$
- 2: The server stores $h_0 = \frac{1}{M} \sum_{m=0}^{M} h_0^m = 0$
- 3: Send a compressed version of the difference to the server $\mathcal{Q}(\nabla f_m(x_k) h_{\nu}^m)$
- 4: Refreshing the memory $h_{k+1}^m = h_k^m + \alpha \mathcal{Q}(\nabla f_m(x_k) h_k^m)$
- 5: The server calculates $g_k = h_k + \frac{1}{M} \sum_{k=1}^{M} \mathcal{Q}(\nabla f_m(x_k) h_k^m)$
- 6: For the update $x_{k+1} = x_k \gamma g_k$
- 7: The server updates $h_{k+1} = h_k + \alpha \frac{1}{M} \sum_{m=1}^{M} \mathcal{Q}(\nabla f_m(x_k) h_k^m)$



Mishchenko K. et al. Distributed Learning with Compressed

Gradient Differences

Lecture

Aleksandr Beznosikov

 Q: What are some other questions about convergence/estimates of convergence?

- Q: What are some other questions about convergence/estimates of convergence? Does it converge better at all?
- Best estimate per number of communications for the unaccelerated method with unbiased compression (DIANA):

$$\mathcal{O}\left(\left[1+rac{\omega}{\mathcal{M}}
ight]rac{L}{\mu}\lograc{1}{arepsilon}
ight).$$

• Estimate on the number of communications for GD:

$$\mathcal{O}\left(\frac{L}{\mu}\log\frac{1}{\varepsilon}\right)$$
.

 In terms of number of communications, compressed methods are inferior to basic methods – this is to be expected (compression fees).
 BUT!

• Compressors compress information β times and it is typical that $\beta > \omega$.

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- Best estimate on the number of information for the unaccelerated method with unbiased compression (DIANA):

$$\mathcal{O}\left(\left[rac{1}{eta} + rac{1}{M}
ight]rac{L}{\mu}\lograc{1}{arepsilon}
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- Compressors compress information β times and it is typical that $\beta > \omega$.
- Best estimate on the number of information for the unaccelerated method with unbiased compression (DIANA):

$$\mathcal{O}\left(\left\lceil \frac{1}{eta} + \frac{1}{M} \right
ceil \frac{L}{\mu} \log \frac{1}{arepsilon}
ight).$$

Estimate on the number of information for GD:

$$\mathcal{O}\left(\frac{L}{\mu}\log\frac{1}{\varepsilon}\right)$$
.

• The unbiased compressor provably improves the number of transmitted information, an improvement factor: $\left[\frac{1}{\beta} + \frac{1}{M}\right]$.

Aleksandr Beznosikov Lecture 29 April 2025 53 / 85

implemented without a "server".

- Often in practice "centralised communications via a server" are
- Architecture with AllGather/AllReduce procedure: some graph of links/communications is given, messages are exchanged according to this graph, including averaging can be organised.



Chan, E. et al. Collective communication: theory, practice, and experience

Centralised communications without a server

Operation	Before	After
Broadcast	$\begin{array}{c c c c c c c c c c c c c c c c c c c $	$\begin{array}{c c c c c c c c c c c c c c c c c c c $
Reduce(- to-one)	$\begin{array}{c c c c c c c c c c c c c c c c c c c $	$\begin{array}{c c c c c c c c c c c c c c c c c c c $
Scatter	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	$\begin{array}{c c c c c c c c c c c c c c c c c c c $
Gather	$\begin{array}{c c c c c c c c c c c c c c c c c c c $	Node 0 Node 1 Node 2 Node 3
Allgather	$\begin{array}{c c c c c c c c c c c c c c c c c c c $	$\begin{array}{c c c c c c c c c c c c c c c c c c c $
Reduce- scatter	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	$\begin{array}{c c c c c c c c c c c c c c c c c c c $
Allreduce	$\begin{array}{c c c c c c c c c c c c c c c c c c c $	$\begin{array}{c c c c c c c c c c c c c c c c c c c $



Aleksandr Beznosikov Lecture 29 April 2025 55 / 85

Ring AllReduce

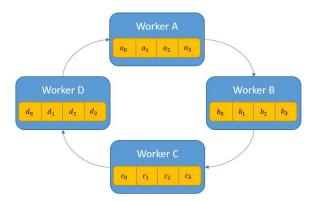


Figure: Picture from here



56 / 85

Ring AllReduce: first step of averaging

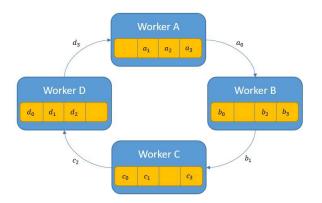


Figure: Picture from here



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Ring AllReduce: second step of averaging

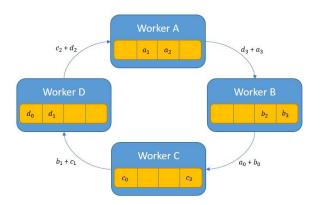


Figure: Picture from here



58 / 85

Ring AllReduce: first step of backprogation

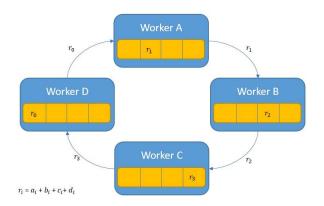


Figure: Картинка отсюда



Aleksandr Beznosikov Lecture

59 / 85

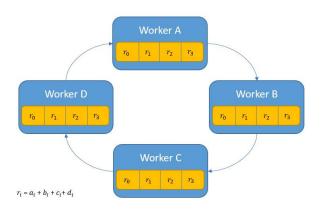


Figure: Картинка отсюда



Quantized GD (QGD) with AllReduce

Algorithm 1 QGD

```
Input: step size \gamma > 0, starting point x_0 \in \mathbb{R}^d, number of iterations K
 1: for k = 0, 1, ..., K - 1 do
```

- for m = 1, ..., M in parallel do 2:
- Compute $\nabla f_m(x_k)$ in x_k 3:
- Indipendetly generate $g_{k,m} = \mathcal{Q}(\nabla f_m(x_k))$ 4:
- Run AllReduce $\{g_{k,m}\}$ and get $g_k = \frac{1}{M} \sum_{m=1}^{M} g_{k,m}$ 5:
- 6: $x_{k+1} = x_k - \gamma g_k$
- end for
- 8: end for
- Output: x_K



Quantized GD (QGD) with AllReduce

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         for m = 1, ..., M in parallel do
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 3:
             Indipendetly generate g_{k,m} = \mathcal{Q}(\nabla f_m(x_k))
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             Run AllReduce \{g_{k,m}\} and get g_k = \frac{1}{M} \sum_{m=1}^{M} g_{k,m}
 5:
 6:
             x_{k+1} = x_k - \gamma g_k
 7:
         end for
 8: end for
Output: X_K
```

• Q: What kind of problems might appear with (for example) Randk?

Quantized GD (QGD) with AllReduce

Algorithm 1 QGD

Output: x_K

```
Input: step size \gamma>0, starting point x_0\in\mathbb{R}^d, number of iterations K

1: for k=0,1,\ldots,K-1 do

2: for m=1,\ldots,M in parallel do

3: Compute \nabla f_m(x_k) in x_k

4: Indipendetly generate g_{k,m}=\mathcal{Q}(\nabla f_m(x_k))

5: Run AllReduce \{g_{k,m}\} and get g_k=\frac{1}{M}\sum_{m=1}^M g_{k,m}

6: x_{k+1}=x_k-\gamma g_k

7: end for

8: end for
```

Q: What kind of problems might appear with (for example) Randk?
 The same non-zero coordinates on different devices can cause collisions.

Aleksandr Beznosikov Lecture 29 April 2025 61 / 85

PermK: do a dependent randomisation

Permutation compressor (dependent RandK)

Suppose that $d \ge n$ and d = qn, where $q \ge 1$ is an integer. Let $\pi = (\pi_1, \dots, \pi_d)$ be a random permutation of $\{1, \dots, d\}$. Then for each $i \in \{1, 2, ..., n\}$ we have the following compression operator

$$Q_i(u) = n \cdot \sum_{j=q(i-1)+1}^{qi} u_{\pi_j} e_{\pi_j}.$$



Szlendak, R. et al. Permutation Compressors for Provably Faster Distributed Nonconvex Optimization

PermK: do a dependent randomisation

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- Szlendak, R. et al. Permutation Compressors for Provably Faster Distributed Nonconvex Optimization
- Friendly to centralised communications without a server.
- In the homogeneous case have the physics of cheap full gradient forwarding.

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62 / 85

Biased compression

Randomisation is good, but there's gap for improvement.

Biased compression

We call the (stochastic) operator C(x) a shifted compression operator if for any $x \in \mathbb{R}^d$ it is fulfilled:

$$\mathbb{E}[\|C(x) - x\|_2^2] \le \left(1 - \frac{1}{\delta}\right) \|x\|_2^2,$$

where $\delta > 0$.

Biased compression: examples

"Greedy" sparsification (selection of the largest modular components)

Consider an operator

$$\mathsf{Top}_k(x) = \sum_{i=d-k+1}^d x_{(i)} e_{(i)},$$

where k — some fixed number from the set $\{1, \ldots, d\}$ (the number of components of vector x that we pass; for example, we can choose k = 1), with the coordinates sorted modulo: $|x_{(1)}| \leq |x_{(2)}| \leq \ldots \leq |x_{(d)}|$, (e_1,\ldots,e_d) — the standard basis in \mathbb{R}^d . It can be shown that this operator is a shifted compression with constant $\delta = \frac{d}{L}k$.



Alistarh D. et al. The convergence of sparsified gradient methods

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Biased compression: examples

- Various examples of compressors (sparsifiers, rounding, etc):
 - Beznosikov A. et al. On Biased Compression for Distributed Learning
- A practical biased compressor based on iterative SVD decomposition: Vogels T. et al. PowerSGD: Practical Low-Rank Gradient Compression for Distributed Optimization



Biased compression: a theorem in the case of 1 node

Theorem (convergence of QGD with shifted compression in case of 1 node)

Let $f \mu$ be strongly convex (or PL) and have L-Lipschitz gradient, then QGD for one node with step $\gamma \leq 1/L$ and with biased compressor with parameter δ converges and is satisfied:

$$f(w_K) - f(w^*) \leq \left(1 - \frac{\gamma \mu}{\delta}\right)^K (f(w_0) - f(w^*)).$$



Beznosikov A. et al. On Biased Compression for Distributed Learning

• Consider the following distributed problem with M=3, d=3 and local functions:

$$\begin{split} f_1(x) &= \langle a, x \rangle^2 + \tfrac{1}{4} \|x\|^2, \ f_2(w) = \langle b, x \rangle^2 + \tfrac{1}{4} \|x\|^2, \ f_3(x) = \langle c, x \rangle^2 + \tfrac{1}{4} \|x\|^2, \\ \text{where } a &= (-3, 2, 2), \ b = (2, -3, 2) \ \text{u} \ c = (2, 2, -3). \end{split}$$

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 where $a = (-3, 2, 2), \ b = (2, -3, 2)$ in $c = (2, 2, -3)$.

- **Q**: where is her optimum? (0,0,0).
- Let the starting point $x_0 = (t, t, t)$ for some t > 0. Then the local gradients are:

$$\nabla f_1(x_0) = \frac{t}{2}(-11, 9, 9), \quad \nabla f_2(x_0) = \frac{t}{2}(9, -11, 9), \quad \nabla f_3(x_0) = \frac{t}{2}(9, 9, -11).$$

 Q: what will the QGD (gradient descent with compressions) step look like if we use *Top1* compression?

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$$f_1(x) = \langle a, x \rangle^2 + \frac{1}{4} \|x\|^2, \ f_2(w) = \langle b, x \rangle^2 + \frac{1}{4} \|x\|^2, \ f_3(x) = \langle c, x \rangle^2 + \frac{1}{4} \|x\|^2,$$
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 Q: what will the QGD (gradient descent with compressions) step look like if we use Top1 compression?

$$x_1 = (t, t, t) + \gamma \cdot \frac{11}{6}(t, t, t) = \left(1 + \frac{11\gamma}{6}\right)x_0.$$

• We move away from the solution geometrically for any $\gamma > 0$.

Aleksandr Beznosikov Lecture 29 April 2025 67 / 85

• Let's try to remember what we didn't pass on in the communication process:

$$e_{1,m} = 0 + \gamma \nabla f_m(x_0) - C(0 + \gamma \nabla f_m(x_0)).$$

Biased compression: error compensation

 Let's try to remember what we didn't pass on in the communication process:

$$e_{1,m} = 0 + \gamma \nabla f_m(x_0) - C(0 + \gamma \nabla f_m(x_0)).$$

And add this to future parcels:

$$C(e_{1,m} + \gamma \nabla f_m(x_1))$$

Biased compression: error compensation

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In an arbitrary iteration, it is written as follows:

Parcel:
$$C(e_{k,m} + \gamma \nabla f_m(x_k))$$
,
 $e_{k+1,m} = e_{k,m} + \gamma \nabla f_m(x_k) - C(e_{k,m} + \gamma \nabla f_m(x_k))$



Biased compression: error compensation

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,
 $e_{k+1,m} = e_{k,m} + \gamma \nabla f_m(x_k) - C(e_{k,m} + \gamma \nabla f_m(x_k))$

This technique is called error compensation (error feedback).

Stich S. et al. Sparsified SGD with memory

Aleksandr Beznosikov Lecture 29 April 2025 68 / 85

by server

by workers

by workers

▷ by workers

by workers

by workers

by server

by server

by server

QGD with error feedback

Algorithm 1 QGD c error feedback

Input: step size $\gamma > 0$, starting point $x_0 \in \mathbb{R}^d$, starting errors $e_{0,m} = 0$ for all m from 1 to M, number of iterations K

- 1: **for** k = 0, 1, ..., K 1 **do**
- 2: Send x_k to all workers
- 3: **for** m = 1, ..., M in parallel **do**
- 4: Recieve x_k from the master
- 5: Compute $\nabla f(w_k)$ in x_k
- 6: Generate $g_{k,m} = C(e_{k,m} + \gamma \nabla f(x_k))$
- 7: Вычислить $e_{k+1,m} = e_{k,m} + \gamma \nabla f_m(x_k) g_{k,m}$
- 8: Send $g_{k,m}$ to master
- $g_{k,m}$ to master
- 9: end for
- 10: Recieve $g_{k,m}$ from all workers
- 11: Compute $g_k = \frac{1}{M} \sum_{m=1}^{M} g_{k,m}$
- 12: $x_{k+1} = x_k g_k$
- 13: end for

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Local updates Compression Dif. priv. Pers.

QGD with error feedback: convergence

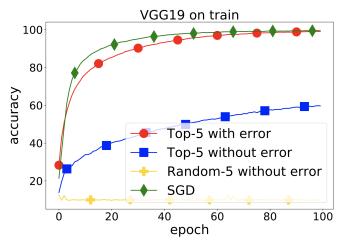


Figure: Accuracy during training of VGG19 on CIFAR10 using different compressors

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QGD c error feedback: convergence

Theorem GD with error feedback

Let all local functions f_m be μ -simply convex and have L-Lipschitz gradient, then if $\gamma \leq \frac{1}{28\delta I}$ then

$$\mathbb{E}\left[f(\tilde{x}_{K}) - f(x^{*})\right] \leq \mathcal{O}\left(\delta L \|x_{0} - x^{*}\|^{2} \exp\left(-\frac{\gamma \mu K}{2}\right)\right) + \gamma \cdot \frac{\delta}{\mu} \cdot \frac{1}{M} \sum_{m=1}^{M} \|\nabla f_{m}(x^{*})\|^{2}\right).$$

Stich S. and Karimireddy S. The error-feedback framework:

- Better rates for SGD with delayed gradients and compressed communication
- Beznosikov A. et al. On Biased Compression for Distributed Learning
- Same problem as QGD the second term in the evaluation should be

aliminated

71 / 85

Biased compression: solving the plateau issue

• Idea like DIANA: memory + compression of difference

Algorithm 1 EF21 (sketch)

- 1: Each device m possesses a vector of "memory" $g_0^m = 0$
- 2: The server stores $h_0 = \frac{1}{M} \sum_{m=1}^{M} h_0^m = 0$
- 3: Send a compressed version of the difference to the server $C(\nabla f_m(w^k) h_k^m)$
- 4: Refreshing the memory $h_{k+1}^m = h_k^m + C(\nabla f_m(x^k) h_k^m)$
- 5: Server compute $h_{k+1} = h_k + \frac{1}{M} \sum_{m=1}^{M} C(\nabla f_m(x^k) h_k^m)$
- 6: Update: $x_{k+1} = x_k \gamma h_{k+1}$



Richtarik P. et al. EF21: A New, Simpler, Theoretically Better, and Practically Faster Error Feedback

Aleksandr Beznosikov Lecture 29 April 2025 72 / 85

 Best estimate on the number of communications for the unaccelerated method with unbiased compression (DIANA):

$$\mathcal{O}\left(\left[1+rac{\omega}{\mathit{M}}
ight]rac{\mathit{L}}{\mu}\lograc{1}{arepsilon}
ight).$$

Best estimate on the number of communications for the unaccelerated method with biased compression (EF-21):

$$\mathcal{O}\left(\left[1+\delta\right]\frac{L}{\mu}\log\frac{1}{arepsilon}
ight).$$

 It has already been discussed that these estimates are worse than for the baseline GD.



73 / 85

Unbiased vs biased

• Compressors compress information β times and it is typical that $\beta \geq \omega$ and $\beta > \delta$.

Unbiased vs biased

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- Best estimate on the number of information for the unaccelerated method with unbiased compression (DIANA):

$$\mathcal{O}\left(\left[\frac{1}{\beta} + \frac{1}{M}\right] \frac{L}{\mu} \log \frac{1}{\varepsilon}\right).$$

 As discussed, the unbiased compressor provably improves the number of transmitted information.

Unbiased vs biased

- Compressors compress information β times and it is typical that $\beta \geq \omega$ and $\beta \geq \delta$.
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$$\mathcal{O}\left(\left[\frac{1}{eta} + \frac{1}{M}\right] \frac{L}{\mu} \log \frac{1}{arepsilon}\right).$$

- As discussed, the unbiased compressor provably improves the number of transmitted information.
- The biased compressor has an estimate:

$$\mathcal{O}\left(\left\lceil \frac{1}{eta} + \frac{\delta}{eta} \right\rceil \frac{L}{\mu} \log \frac{1}{arepsilon} \right).$$

 The biased compressor does not improve the number of transmitted information in the general case. And this is an open question: how to see the theoretical superiority of biased operators. Differential privacy

Why not existing approaches?

Anonymization - Netflix example.

- Anonymization Netflix example.
- Encryption long and computationally expensive.

Differential privacy

- Suppose we are given data size of n where each sample is generated from a domain \mathcal{X} .
- We call two datasets neighboring if they differ by exactly one element.
- Suppose we consider some (possibly stochastic) algorithm \mathcal{A} that, using a given date from \mathcal{X}^d , produces an answer Y (possibly a random variable) from some set E.

Definition

We say that the mechanism \mathcal{A} is (ε, δ) - DP (differential private) if for all neighboring datasets $D, D' \in \mathcal{X}^n$ and for all $E \subseteq \mathcal{Y}$ it follows that

$$\mathbb{P}\{\mathcal{A}(D) \in E\} \leq e^{\varepsilon} \mathbb{P}\{\mathcal{A}(D') \in E\} + \delta.$$

• The essence - change one element in the sample, the algorithm's answer should not change much. I.e. nothing depends on data of one

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77 / 85

Properties of differential privacy

- If $\varepsilon \to 0$, then $e^{\varepsilon} \to 1$. If additionally $\delta \to 0$, we have ideal situation.
- Typically, we want to have $\delta \leq \frac{1}{d}$

Differential privacy mechanisms

Gaussian:

$$\tilde{\mathcal{A}}(D) = \mathcal{A}(D) + Z$$
, with $Z \sim \mathcal{N}(0, \sigma^2)$.

• If we want to have the (ε, δ) - DP algorithm $\tilde{\mathcal{A}}$, we need

$$\sigma = \frac{\Delta \ln \frac{5}{4\delta}}{\varepsilon}.$$

Here we use sensitivity of the algorithm A:

$$\Delta = \max_{\mathsf{neighboring}\ D, D'} \|\mathcal{A}(D) - \mathcal{A}(D')\|$$

Differential privacy mechanisms

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Laplacian:

$$\tilde{A}(D) = A(D) + Z$$
, with $Z \sim \text{Lap}(\alpha)$.

• If we want to have the $(\varepsilon,0)$ - DP algorithm \tilde{A} , we need

$$\alpha = \Delta/\varepsilon$$
.

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Consider

$$\min_{\mathbf{x}\in\mathbb{R}^d}\left[\frac{1}{n}\sum_{i=1}^n\ell(g(\mathbf{x},a_i),b_i))+\lambda\varphi(\mathbf{x})\right].$$

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- For simplicity, we assume that we store all data on single device, but we forward something during training.
- Q: what do we send?



Consider

$$\min_{\mathbf{x}\in\mathbb{R}^d}\left[\frac{1}{n}\sum_{i=1}^n\ell(g(\mathbf{x},a_i),b_i))+\lambda\varphi(\mathbf{x})\right].$$

- For simplicity, we assume that we store all data on single device, but we forward something during training.
- Q: what do we send? Models and gradients.

Model perturbation:

$$\tilde{x}_k = x_k + b$$
 with $b \sim \mathcal{N}(0, \sigma \cdot I)$

• Model perturbation:

$$\tilde{x}_k = x_k + b$$
 with $b \sim \mathcal{N}(0, \sigma \cdot I)$

• Limiting result: if we add noise to x^* (optimal model) and $\sigma^2 \sim \frac{\log \delta^{-1}}{\lambda^2 \varepsilon^2}$, then we have (ε, δ) -DP.



Gradient perturbation:

$$g(x) = \nabla f(x) + b$$
,

Gradient perturbation:

$$g(x) = \nabla f(x) + b$$

but in practice to do in the other way with clipping:

$$g(x) = \min \left\{ \frac{C}{\|\nabla f(x)\|}; 1 \right\} \nabla f(x) + b.$$

Q: why do we need the clipping technique here?

Gradient perturbation:

$$g(x) = \nabla f(x) + b$$

but in practice to do in the other way with clipping:

$$g(x) = \min \left\{ \frac{C}{\|\nabla f(x)\|}; 1 \right\} \nabla f(x) + b.$$

Q: why do we need the clipping technique here? To control gradient and in fact sensitivity in the ball.



• Gradient perturbation:

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Q: why do we need the clipping technique here? To control gradient and in fact sensitivity in the ball.

• Result: if $\sigma^2 \sim \frac{C^2 \log \frac{1}{\delta}}{\varepsilon^2}$, then we have (ε, δ) -DP

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Personalization



4 classical options

What kind of model can be used on local devices?

- Golbal model trained on a big open dataset on the server.
- The same as previous, but with local tuning.
- Global federated learning model.
- The same as previous, but with local tuning.



How to make better?

Change the objective:

• Difference penalization

$$\min_{x_1,...,x_M \in \mathbb{R}^d} \left[\frac{1}{M} \sum_{m=1}^M f_m(x_m) + \frac{\lambda}{2M} \sum_{m=1}^M \|x_m - \bar{x}\|_2^2 \right],$$

where
$$\bar{x} = \frac{1}{M} \sum_{m=1}^{M} x_m$$
.

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$$\min_{x_1,...,x_M \in \mathbb{R}^d} \left[\frac{1}{M} \sum_{m=1}^M f_m(x_m) + \frac{\lambda}{2M} \sum_{m=1}^M \|x_m - \bar{x}\|_2^2 \right],$$

where
$$\bar{x} = \frac{1}{M} \sum_{m=1}^{M} x_m$$
.

• Train the network:

$$\min_{x_1,\ldots,x_M\in\mathbb{R}^d} \min_{W\in\mathcal{C}} \left[\frac{1}{M} \sum_{m=1}^M f_m(x_m) + \frac{\lambda}{2} X^T W X \right],$$

where X is a matrix of x_m , W is a matrix with properties of the communication network.

Aleksandr Beznosikov Lecture 29 April 2025 85 / 85