Variance-Reduced Stochastic Quasi-Newton Methods for Decentralized Learning

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Outline

- Background
- 2 A General Framework of Variance-Reduced Decentralized Stochastic Quasi-Newton Methods
- 3 Two Fully Decentralized Stochastic Quasi-Newton Methods
- Mumerical Experiments

Background

Machine learning over networks

- large-scale learning
- privacy-preserving learning
- decentralized system control







Background

Consider the decentralized convex optimization problem

$$x^* = \arg\min_{x \in \mathbb{R}^d} F(x) \triangleq \frac{1}{n} \sum_{i=1}^n f_i(x), \quad \text{where } f_i(x) \triangleq \frac{1}{m_i} \sum_{l=1}^{m_i} f_{i,l}(x).$$

 $f_{i,l}$ is the *l*-th sample cost on node *i*, assumed to be convex and differentiable.

- We consider the case that m_i is very large.
- An equivalent decentralized formulation is

$$\mathbf{x}^* = \underset{\mathbf{x} = [\mathbf{x}_1; \dots; \mathbf{x}_n] \in \mathbb{R}^{nd}}{\arg \min} f(\mathbf{x}) \triangleq \frac{1}{n} \sum_{i=1}^n f_i\left(\mathbf{x}_i\right), \text{ s.t. } \mathbf{x}_i = \mathbf{x}_j, \ \forall j \in \mathcal{N}_i, \ \forall i.$$

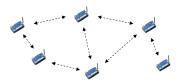


Fig 1. Decentralized network

- Data are distributed across a connected network of n nodes.
- Each node is only allowed to send/receive information to/from its neighboring nodes.
- All the nodes cooperate to obtain a common variable x^* .

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Decentralized Deterministic Algorithms

First-order methods

- penalized gradient methods [Nedic 2009, Yuan 2016]
- EXTRA/PG-EXTRA [Shi 2015a, Shi 2015b]
- primal-dual methods [Shi 2014, Chang 2015]
- gradient tracking [Di Lorenzo 2015, Qu 2017, Nedic 2017, Sun 2019]

Second-order methods

- penalized methods [Mokhtari 2016, Bajovic 2017, Mansoori 2019]
- primal-dual methods [Mokhtari 2016, Zhang 2021]
- approximate Newton-type method [Li 2020]
- cubically-regularized Newton method [Daneshmand 2021]
- adaptive Newton method with a finite-time consensus inner loop [Zhang 2020]

Decentralized Stochastic Algorithms

First-order methods

- stochastic gradient [Chen 2012, Tang 2018, Pu 2019]
- variance-reduced gradient [Mokhtari 2016, Xin 2020, Li 2020, Pu 2021]

Second-order methods

seldom investigated

Can we propose computationally affordable decentralized stochastic second-order methods?

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A General Framework

Assumption 1: The mixing matrix W is nonnegative with $w_{ij} \geq 0$. The weight $w_{ij} = 0$ if and only if $j \notin \mathcal{N}_i$. W is symmetric and doubly stochastic, i.e., $W = W^T$ and $W1_n = 1_n$. The null space of $I_n - W$ is span (1_n) .

• "Averaging" property of the mixing step

$$\|\mathsf{W}\mathsf{x}^k - \mathsf{W}_{\infty}\mathsf{x}^k\| \le \sigma \|\mathsf{x}^k - \mathsf{W}_{\infty}\mathsf{x}^k\|,$$

where $\sigma = \|W - \frac{1}{n} \mathbf{1}_n \mathbf{1}_n^T\|_2 < 1$, $W = W \otimes I_d$ and $W_{\infty} = \frac{\mathbf{1}_n \mathbf{1}_n^T}{n} \otimes I_d$.

ullet σ represents the connectedness of the network.

A General Framework

• Node i updates x_i^{k+1} according to the following decentralized stochastic quasi-Newton step

$$x_i^{k+1} = \sum_{j=1}^n w_{ij} x_j^k - \alpha d_i^k.$$

• In centralized setting, an ideal d_i^{k+1} is the global negative Newton direction

$$\left(\frac{1}{n}\sum_{i=1}^{n}\nabla^{2}f_{i}(\overline{\mathbf{x}}^{k+1})\right)^{-1}\left(\frac{1}{n}\sum_{i=1}^{n}\nabla f_{i}(\overline{\mathbf{x}}^{k+1})\right),\,$$

where $\bar{x}^{k+1} = \frac{1}{n} \sum_{i=1}^{n} x_i^{k+1}$.

- However, computing the global negative Newton direction is expensive: lack of information from the entire network / large m_i .
- ullet Our idea: We update d_i^{k+1} with Hessian inverse approximation H_i^{k+1} and gradient approximation g_i^{k+1} , given by

$$d_i^{k+1} = H_i^{k+1} g_i^{k+1}.$$

We will construct g_i^{k+1} below and H_i^{k+1} in the next part.

Construct g_i^{k+1}

• Node i obtains a corrected stochastic gradient v_i^{k+1} with SVRG, as

$$v_i^{k+1} = \frac{1}{b_i} \sum_{I \in S_i^{k+1}} \left(\nabla f_{i,I}(x_i^{k+1}) - \nabla f_{i,I}(\tau_i^{k+1}) \right) + \nabla f_i(\tau_i^{k+1}),$$

where $S_i^{k+1} \subseteq \{1, \ldots, m_i\}$ with batch size b_i , while $\tau_i^{k+1} = \tau_i^k$ or $\tau_i^{k+1} = x_i^{k+1}$ if mod(k+1, T) = 0. We have $\mathbb{E}[v_i^k] = \nabla f_i(x_i^k)$.

ullet g_i^{k+1} is constructed with a dynamic average consensus (DAC) step

$$g_i^{k+1} = \sum_{i=1}^n w_{ij}g_j^k + v_i^{k+1} - v_i^k,$$

with initialization $g_i^0 = v_i^0 = \nabla f_i(x_i^0)$. We have $\frac{1}{n} \sum_{i=1}^n g_i^k = \frac{1}{n} \sum_{i=1}^n v_i^k, \forall k$.

• With SVRG, $v_i^k \approx \nabla f_i(x_i^k)$. Therefore, with DAC, when x_i^k are almost consensual, $g_i^k \approx \frac{1}{n} \sum_{i=1}^n \nabla f_i(\bar{\mathbf{x}}^k)$.

A General Framework

Proposed general framework

$$\begin{cases} x_{i}^{k+1} = \sum_{j=1}^{n} w_{ij} x_{j}^{k} - \alpha d_{i}^{k}, & x_{i}^{k+1} \\ v_{i}^{k+1} = \frac{1}{b_{i}} \sum_{l \in S_{i}^{k+1}} \left(\nabla f_{i,l}(x_{i}^{k+1}) - \nabla f_{i,l}(\tau_{i}^{k+1}) \right) & y_{i}^{k+1} \approx \nabla f_{i}(x_{i}^{k+1}) \\ + \nabla f_{i}(\tau_{i}^{k+1}), & y_{i}^{k+1} \approx \nabla f_{i}(x_{i}^{k+1}) \\ g_{i}^{k+1} = \sum_{j=1}^{n} w_{ij} g_{j}^{k} + v_{i}^{k+1} - v_{i}^{k}, & H_{i}^{k+1} ? \\ d_{i}^{k+1} = H_{i}^{k+1} g_{i}^{k+1}. & d_{i}^{k+1} \end{cases}$$

Flow on node i

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• In the proposed general framework, H_i^k is constructed locally given g_i^k and x_i^k , without extra sampling or communication.

Linear Convergence of General Framework

Assumption 2: There exist two constants M_1 and M_2 with $0 < M_1 \le M_2 < \infty$ such that

$$M_1I_d \leq H_i^k \leq M_2I_d, \ \forall i=1,\ldots,n, \ \forall k\geq 0.$$

• We will check Assumption 2 in the next part.

Assumption 3: Each local sample cost $f_{i,l}$ is convex and has Lipschitz continuous gradients, i.e.,

$$f_{i,l}(y) \leq f_{i,l}(x) + \nabla f_{i,l}(x)^T (y-x) + \frac{L}{2} ||y-x||^2.$$

Assumption 4: The global cost function F is strongly convex, i.e.,

$$F(y) \ge F(x) + \nabla F(x)^T (y - x) + \frac{\mu}{2} ||y - x||^2.$$

Linear Convergence of General Framework

Theorem 1

Under Assumptions 1–4, if the parameters satisfy

$$\alpha \leq \frac{(1-\sigma^2)^2 \mu \mathit{M}_1}{200 \mathsf{L}^2 \mathit{M}_2^2}, B \leq \frac{1}{160} \min \left\{1, \frac{\zeta (1-\sigma^2)^2}{\gamma^2} \right\}, \, \mathcal{T} \geq \frac{2 \log(280/(\zeta (1-\sigma^2)^2))}{\zeta \widetilde{\alpha}},$$

where $\zeta=\left(\frac{\mu}{L}\right)^2\left(\frac{M_1}{M_2}\right)^2, \gamma=1-\frac{M_1}{M_2},$ and $\widetilde{\alpha}=\frac{M_2^2L^2}{M_1\mu}\alpha$. Then, the proposed general framework converges linearly to the optimal solution, such that

$$\|\boldsymbol{u}^{(t+1)T}\|_{\infty}^{q} \leq 0.9 \|\boldsymbol{u}^{tT}\|_{\infty}^{q}.$$

- ullet Define the non-sampling rate $B=\max_{i\in\{1,\dots,n\}}\left\{rac{m_i-b_i}{(m_i-1)b_i}
 ight\}<1.$
- Define $\kappa_F = L/\mu$ and $\kappa_H = M_2/M_1$.
- The total number of stochastic gradient evaluations is

$$\mathcal{O}\left(\left(\max_{i}\{\textit{\textit{m}}_{i}\}+\frac{\max_{i}\{\textit{\textit{b}}_{i}\}\cdot\kappa_{\textit{F}}^{2}\kappa_{\textit{H}}^{2}\log\frac{\kappa_{\textit{F}}\kappa_{\textit{H}}}{1-\sigma^{2}}}{(1-\sigma^{2})^{2}}\right)\log\frac{1}{\epsilon}\right).$$



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Two Quasi-Newton Methods

• Recall that H_i^k is constructed locally given g_i^k and x_i^k without extra sampling or communication, and is assumed to satisfy

$$M_1I_d \leq H_i^k \leq M_2I_d$$
.

- How to construct Hessian approximation with gradient information?
 - Quasi-Newton methods!

Two Quasi-Newton Methods

• In centralized setting, two well-known quasi-Newton methods are

$$(DFP) \ H^{k+1} = H^k + \frac{s^k (s^k)^T}{(s^k)^T y^k} - \frac{H^k y^k (y^k)^T H^k}{(H^k y^k)^T y^k},$$

$$(BFGS) \ H^{k+1} = \left(I_d - \frac{s^k (y^k)^T}{(s^k)^T y^k}\right) H^k \left(I_d - \frac{y^k (s^k)^T}{(s^k)^T y^k}\right) + \frac{s^k (s^k)^T}{(s^k)^T y^k}.$$

Here, s^k and y^k are defined as

$$s^{k} = x^{k+1} - x^{k}$$
 and $y^{k} = \nabla F(x^{k+1}) - \nabla F(x^{k})$.

- Curvature condition $(s^k)^T y^k > 0$ holds due to strong convexity and thus $H^k > 0$.
- However, g_i^k are noisy due to stochastic gradient noise and disagreement among the nodes, how to preserve $M_1I_d \leq H_i^k \leq M_2I_d$?
 - Damping technique and limited-memory technique!

Damped Regularized Limited-memory DFP: Step I

• To guarantee $\lambda_{\min}(H_i^{k+1}) > 0$, we use **damping technique** and add a regularization

$$H_i^{k+1} = H_i^k + \frac{\hat{s}_i^k (\hat{s}_i^k)^T}{(\hat{s}_i^k)^T \hat{y}_i^k} - \frac{H_i^k \hat{y}_i^k (\hat{y}_i^k)^T H_i^k}{(\hat{y}_i^k)^T H_i^k \hat{y}_i^k} + \rho I_d,$$

where $s_i^k = x_i^{k+1} - x_i^k$, $y_i^k = g_i^{k+1} - g_i^k$, $\hat{s}_i^k = s_i^k - \rho y_i^k$, $\hat{y}_i^k = \theta_i^k y_i^k + (1 - \theta_i^k)(H_i^0 + \epsilon I_d)^{-1} \hat{s}_i^k$,

$$\theta_i^k = \min \left\{ \tilde{\theta}_i^k, \frac{\tilde{L}\|\hat{\mathbf{s}}_i^k\|}{\|y_i^k\|} \right\} \text{ and } \tilde{\theta}_i^k = \left\{ \begin{array}{l} \frac{0.75(\hat{\mathbf{s}}_i^k)^T \left(H_i^0 + \epsilon I_d\right)^{-1} \hat{\mathbf{s}}_i^k}{(\hat{\mathbf{s}}_i^k)^T \left(H_i^0 + \epsilon I_d\right)^{-1} \hat{\mathbf{s}}_i^k - (\hat{\mathbf{s}}_i^k)^T y_i^k}, \\ \text{if } (\hat{\mathbf{s}}_i^k)^T y_i^k \leq 0.25(\hat{\mathbf{s}}_i^k)^T \left(H_i^0 + \epsilon I_d\right)^{-1} \hat{\mathbf{s}}_i^k, \\ 1, \text{ otherwise.} \end{array} \right.$$

- With the corrected \hat{y}_i^k by the damping technique, $(\hat{s}_i^k)^T \hat{y}_i^k > 0$ and thus $H_i^k \succ 0$.
- In addition, the regularization guarantees $\lambda_{\min}(H_i^{k+1}) > \rho$.

Damped Regularized Limited-memory DFP: Step II

- To guarantee $\lambda_{\max}(H_i^{k+1}) < \infty$, we further use **limited-memory technique**.
- ullet Use a fixed moving window of M past variations

$$\{\hat{s}_i^{k+1-M},\hat{s}_i^{k-M},\ldots,\hat{s}_i^k\} \quad \text{and} \quad \{\hat{y}_i^{k+1-M},\hat{y}_i^{k-M},\ldots,\hat{y}_i^k\}.$$

• Recursively use M past variations and update H_i^{k+1} as

$$H_{i}^{k,(t+1)} = H_{i}^{k,(t)} + \frac{\hat{s}_{i}^{p}(\hat{s}_{i}^{p})^{T}}{(\hat{s}_{i}^{p})^{T}\hat{y}_{i}^{p}} - \frac{H_{i}^{k,(t)}\hat{y}_{i}^{p}(\hat{y}_{i}^{p})^{T}H_{i}^{k,(t)}}{(\hat{y}_{i}^{p})^{T}H_{i}^{k,(t)}\hat{y}_{i}^{p}} + \rho I_{d},$$

where $t = 0, \dots, \tilde{M} - 1$, $p = k + 1 - \tilde{M} + t$, and $\tilde{M} = \min\{k + 1, M\}$.

- ullet Restart after $ilde{M}$ iterations with initialization $eta I_d \preceq H_i^{k,(0)} \preceq \mathcal{B} I_d$.
- Computation cost per iteration is $O(Md^2)$, storage is $O(d^2 + Md)$.

Damped Limited-memory BFGS: Step I

• To guarantee $\lambda_{\min}(H_i^{k+1}) > 0$, we use **damping technique**.

$$H_i^{k+1} = \left(I_d - \frac{s_i^k(\hat{y}_i^k)^T}{(s_i^k)^T\hat{y}_i^k}\right)H_i^k\left(I_d - \frac{\hat{y}_i^k(s_i^k)^T}{(s_i^k)^T\hat{y}_i^k}\right) + \frac{s_i^k(s_i^k)^T}{(s_i^k)^T\hat{y}_i^k}.$$

where $\hat{y}_i^k = \theta y_i^k + (1 - \theta)(H_i^0 + \epsilon I)^{-1} s_i^k$, and

$$\theta_{i}^{k} = \min \left\{ \tilde{\theta}_{i}^{k}, \frac{\tilde{L} \| \mathbf{s}_{i}^{k} \|}{\| \mathbf{y}_{i}^{k} \|} \right\}, \ \tilde{\theta}_{i}^{k} = \left\{ \begin{array}{l} \frac{0.75(\mathbf{s}_{i}^{k})^{T} \left(H_{i}^{0} + \epsilon I_{d}\right)^{-1} \mathbf{s}_{i}^{k}}{(\mathbf{s}_{i}^{k})^{T} \left(H_{i}^{0} + \epsilon I_{d}\right)^{-1} \mathbf{s}_{i}^{k} - (\mathbf{s}_{i}^{k})^{T} \mathbf{y}_{i}^{k}}, \\ \text{if } (\mathbf{s}_{i}^{k})^{T} \mathbf{y}_{i}^{k} \leq 0.25(\mathbf{s}_{i}^{k})^{T} \left(H_{i}^{0} + \epsilon I_{d}\right)^{-1} \mathbf{s}_{i}^{k}, \\ 1, \text{ otherwise.} \end{array} \right.$$

• With the corrected \hat{y}_i^k by the damping technique, $(s_i^k)^T \hat{y}_i^k > 0$ and thus $H_i^k \succ 0$.

Damped Limited-memory BFGS: Step II

• To guarantee $\lambda_{\max}(H_i^{k+1}) < \infty$, we further use **limited-memory technique**.

Algorithm 1: Two-loop recursion

- Store M past variations $\{s_i^{k+1-M}, s_i^{k-M}, \dots, s_i^k\}$ and $\{\hat{y}_i^{k+1-M}, \hat{y}_i^{k-M}, \dots, \hat{y}_i^k\}$.
- Instead of generating H_i^{k+1} explicitly, update $H_i^{k+1}g_i^{k+1}$ by two-loop recursion.
- Restart after \tilde{M} iterations with $\beta I_d \leq H_i^{k,(0)} \leq \mathcal{B}I_d$.
- Computation cost per iteration is O(Md), storage is O(Md).

end

Analysis of Proposed DFP and BFGS methods

Theorem 2 (DFP)

The proposed damped regularized limited-memory DFP satisfies

$$M_1I_d \leq H_i^k \leq M_2I_d, \forall i,$$

where
$$M_1 = \rho + (1 + \omega)^{-2M} \left(\frac{1}{\beta} + \frac{1}{4(\mathcal{B} + \epsilon)}\right)^{-1}$$
, $M_2 = \mathcal{B} + M(4\mathcal{B} + 4\epsilon + \rho)$ and $\omega = 4(\mathcal{B} + \epsilon) \left(\tilde{L} + \frac{1}{\beta + \epsilon}\right)$.

Theorem 3 (BFGS)

The proposed damped limited-memory BFGS method satisfies

$$M_1I_d \leq H_i^k \leq M_2I_d$$

where
$$M_1=\left(\frac{1}{eta}+\frac{M\omega^2}{4(\mathcal{B}+\epsilon)}
ight)^{-1}$$
 and $M_2=(1+\omega)^{2M}$ $\left(\mathcal{B}+\frac{1}{\overline{L}(\omega+2)}
ight)$.

Outline

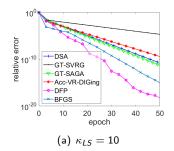
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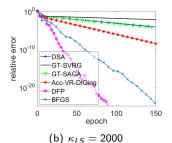
Effects of Condition Number: Synthetic Data

Consider a least-squares problem

$$\min_{x \in \mathbb{R}^d} \frac{1}{2} \sum_{i=1}^n \|A_i x - b_i\|^2.$$

- Define $\kappa_{LS} = \frac{\lambda_{\max}(A^T A)}{\lambda_{\min}(A^T A)}$, where $A = [A_1; \cdots; A_n] \in \mathbb{R}^{nm \times d}$.
- The performance metric: relative error = $\frac{\left\|\mathbf{x}^k \mathbf{x}^*\right\|^2}{n\|\mathbf{x}^0 \mathbf{x}^*\|^2}$.





Comparison with First-order Algorithms: Real Datasets

Consider a logistic regression problem

$$\min_{\mathbf{x} \in \mathbb{R}^d} \frac{\iota}{2} \|\mathbf{x}\|^2 + \frac{1}{n} \sum_{i=1}^n \frac{1}{m_i} \sum_{j=1}^{m_i} \ln \left(1 + \exp \left(-\left(\mathbf{o}_{ij}^T \mathbf{x}\right) \mathbf{p}_{ij}\right)\right),$$

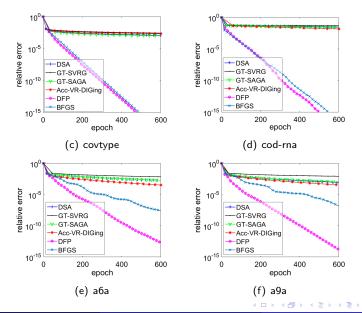
where node i owns m_i training samples $(o_{il}, p_{il}) \in \mathbb{R}^d \times \{-1, +1\}$.

• We normalize each sample such that $\|o_{il}\| = 1, \forall i, l$.

Tabela 1: Datasets used in numerical experiments.

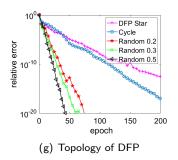
Dataset	# of Samples $(\sum_{i=1}^n m_i)$	# of Features (d)
covtype	40000	54
cod-rna	52000	8
a6a	11220	123
a9a	32560	123
ijcnn1	91700	22

Comparison with First-order Algorithms: Real Datasets



Effects of Topology on ijcnn1

 \bullet of the five graphs are 0.967, 0.950, 0.863, 0.797, and 0.569, respectively.



Decentralized Stochastic Quasi-Newton Methods

Summary

- We propose a general framework of decentralized stochastic quasi-Newton methods, which converges linearly to the optimal solution.
- We specify two fully decentralized stochastic quasi-Newton methods to locally construct Hessian inverse approximations.

Future work

 Improve the theoretical results (with more communication or better initialization).

Thank you!

Two Lemmas

Lemma 4 (DFP)

With the corrected \hat{y}_{i}^{p} by the damping technique, we have

$$0 < \theta_i^p \le 1 \text{ and } (\hat{s}_i^p)^T \hat{y}_i^p \ge 0.25 (\hat{s}_i^p)^T (H_i^{k,(0)} + \epsilon I)^{-1} \hat{s}_i^p.$$

Moreover, H_i^{k+1} keeps positive definite, such that $\lambda_{\min}(H_i^{k+1}) > \rho$.

Lemma 5 (BFGS)

With the corrected \hat{y}_i^p by the damping technique, we have

$$0 < \theta_i^p \le 1 \text{ and } (s_i^p)^T \hat{y}_i^p \ge 0.25 (s_i^p)^T (H_i^{k,(0)} + \epsilon I_d)^{-1} s_i^p.$$

Moreover, H_i^{k+1} keeps positive definite and $\lambda_{\min}(H_i^{k+1}) > 0$.

Effects of Batch Size and Memory Size on ijcnn1

