

A unified variance-reduced accelerated gradient method for convex optimization

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Outline



**Convex
finite-sum
optimization**



**Randomized
incremental
gradient
methods**



Our algorithm
Varag

- Convergence results
- Numerical experiments



Future works

Problem of interest: convex finite-sum optimization



$$\psi^* := \min_{x \in X} \left\{ \psi(x) := \frac{1}{m} \sum_{i=1}^m f_i(x) + h(x) \right\}.$$

- Smooth and convex with L_i -Lipschitz continuous gradient over X
- Simple but possibly nonsmooth over X

Let $f(x) := \frac{1}{m} \sum_{i=1}^m f_i(x)$, we assume that f is **possibly strongly convex** with modulus $\mu \geq 0$.

Problem of interest: convex finite-sum optimization



$$\psi^* := \min_{x \in X} \left\{ \psi(x) := \frac{1}{m} \sum_{i=1}^m f_i(x) + h(x) \right\}.$$

- Wide range of applications in machine learning, statistical inference and image processing.
- Take l_2 -regularized logistic regression problem as an example

$$f_i(x) = l_i(x) := \frac{1}{N_i} \sum_{j=1}^{N_i} \log(1 + \exp(-b_j^i a_j^{i T} x)), \quad i = 1, \dots, m, \quad w(x) = R(x) := \frac{1}{2} \|x\|_2^2,$$

- f_i is the loss function based on training data $\{a_j^i, b_j^i\}_j^{N_i}$, or the loss function associated with agent i for a distributed optimization problem.
- Minimization of the empirical risk

$$f_i(x) = l_i(x) := \mathbb{E}_{\xi_i} [\log(1 + \exp(-\xi_i^T x))], \quad i = 1, \dots, m,$$

- f_i given in the form of expectation where ξ_i models the underlying distribution for training dataset i (of agent i for a distributed problem)
- Minimization of the generalized risk

Randomized incremental gradient (RIG) methods



- Derived from SGD and the idea of reducing variance of the gradient estimator
- SVRG[JZ13] exhibits linear rate of convergence $\mathcal{O}\{(m + L/\mu)\log(1/\epsilon)\}$, same result for Prox-SVRG[XZ14], SAGA[DBL14] and SARAH[NLST17] for **strongly convex problems**
 - Update exact gradient \tilde{g} at the outer loop and a gradient of the component function in the inner loop
 - Variance of G_t vanishes as algorithm proceeds
- SVRG++[AY16] obtains $\mathcal{O}\{m\log(1/\epsilon) + L/\epsilon\}$ for smooth convex problems

They are NOT optimal RIG methods!

$$x_t = x_{t-1} - \eta G_t$$

$$\tilde{g} = \nabla f(\tilde{x})$$

$$G_t = \nabla f_{i_t}(x_{t-1}) - \nabla f_{i_t}(\tilde{x}) + \tilde{g}$$

$$y_i^t = \begin{cases} \nabla f_i(x^t), & i = i_t, \\ y_i^{t-1}, & \text{otherwise,} \end{cases}$$

$$G_t = \nabla f_{i_t}(x_t) - y_{i_t}^{t-1} + \frac{1}{m} \sum_{i=1}^m y_i^{t-1}$$

$$G_0 = \nabla f(\tilde{x})$$

$$G_t = \nabla f_{i_t}(x_{t-1}) - \nabla f_{i_t}(x_{t-2}) + G_{t-1}$$

Optimal RIG methods



Accelerated RIG methods: Catalyst[LMH15], RPDG[LZ17], RGEM[LZ18], and Katyusha[A17], etc.

- All exhibit $\mathcal{O}\{(m + \sqrt{mL/\mu})\log(1/\epsilon)\}$ for strongly convex problems
- Except Katyusha^{ns}[A17], none of these methods can be used directly to solve smooth convex problems. They **required perturbation technique**. Katyusha^{ns} is **not advantageous over accelerated gradient method**.
- Except RGEM[LZ18], **none of the optimal methods can solve stochastic finite-sum problems**
- They are **assume the strongly convexity comes from regularizer term $h(x)$**
- **None of them are unified methods that can be adjust to ill-conditioned problem**, e.g., μ is very small.

$$\psi^* := \min_{x \in X} \left\{ \psi(x) := \frac{1}{m} \sum_{i=1}^m f_i(x) + h(x) \right\}.$$

Table 1: Summary of the recent results on accelerated RIG methods

Algorithms	Deterministic smooth strongly convex	Deterministic smooth convex
RPDG[18]	$\mathcal{O}\{(m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon}\}$	$\mathcal{O}\{(m + \sqrt{\frac{mL}{\epsilon}}) \log \frac{1}{\epsilon}\}^1$
Catalyst[20]	$\mathcal{O}\{(m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon}\}^1$	$\mathcal{O}\{(m + \sqrt{\frac{mL}{\epsilon}}) \log^2 \frac{1}{\epsilon}\}^1$
Katyusha[1]	$\mathcal{O}\{(m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon}\}$	$\mathcal{O}\{(m \log \frac{1}{\epsilon} + \sqrt{\frac{mL}{\epsilon}})\}^1$
Katyusha ^{ns} [1]	NA	$\mathcal{O}\{\frac{m}{\sqrt{\epsilon}} + \sqrt{\frac{mL}{\epsilon}}\}$
RGEM[19]	$\mathcal{O}\{(m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon}\}$	NA

The *Varag* algorithm



Algorithm 1 The variance-reduced accelerated gradient (*Varag*) method

Input: $x^0 \in X$, $\{T_s\}$, $\{\gamma_s\}$, $\{\alpha_s\}$, $\{p_s\}$, $\{\theta_t\}$, and a probability distribution $Q = \{q_1, \dots, q_m\}$ on $\{1, \dots, m\}$.

- 1: Set $\tilde{x}^0 = x^0$.
- 2: **for** $s = 1, 2, \dots$ **do**
- 3: Set $\tilde{x} = \tilde{x}^{s-1}$ and $\tilde{g} = \nabla f(\tilde{x})$.
- 4: Set $x_0 = x^{s-1}$, $\bar{x}_0 = \tilde{x}$ and $T = T_s$.
- 5: **for** $t = 1, 2, \dots, T$ **do**
- 6: Pick $i_t \in \{1, \dots, m\}$ randomly according to Q .
- 7: $\underline{x}_t = [(1 + \mu\gamma_s)(1 - \alpha_s - p_s)\bar{x}_{t-1} + \alpha_s x_{t-1} + (1 + \mu\gamma_s)p_s \tilde{x}] / [1 + \mu\gamma_s(1 - \alpha_s)]$.
- 8: $G_t = (\nabla f_{i_t}(\underline{x}_t) - \nabla f_{i_t}(\tilde{x})) / (q_{i_t} m) + \tilde{g}$.
- 9: $\underline{x}_t = \arg \min_{x \in X} \{ \gamma_s [\langle G_t, x \rangle + h(x)] + \mu V(\underline{x}_t, x) \} + V(x_{t-1}, x)$.
- 10: $\bar{x}_t = (1 - \alpha_s - p_s)\bar{x}_{t-1} + \alpha_s \underline{x}_t + p_s \tilde{x}$.
- 11: **end for**
- 12: Set $x^s = x_T$ and $\tilde{x}^s = \sum_{t=1}^T (\theta_t \bar{x}_t) / \sum_{t=1}^T \theta_t$.
- 13: **end for**

- Similar to SVRG algorithmic scheme
- Adopt AC-SA[GL201] in the inner loop
- Allows general distance via prox-function V
- When $\alpha_s = 1, p_s = 0$, *Varag* reduces to non-accelerated method, and achieves $\mathcal{O}\{(m + L/\mu)\log(1/\epsilon)\}$ as SVRG.

Theorem 1 (Smooth finite-sum optimization) Suppose that the probabilities q_i 's are set to $L_i / \sum_{i=1}^m L_i$ for $i = 1, \dots, m$, and weights $\{\theta_t\}$ are set as

$$\theta_t = \begin{cases} \frac{\gamma_s}{\alpha_s}(\alpha_s + p_s) & 1 \leq t \leq T_s - 1 \\ \frac{\gamma_s}{\alpha_s} & t = T_s. \end{cases} \quad (2.2)$$

Moreover, let us denote $s_0 := \lfloor \log m \rfloor + 1$ and set parameters $\{T_s\}$, $\{\gamma_s\}$ and $\{p_s\}$ as

$$T_s = \begin{cases} 2^{s-1}, & s \leq s_0 \\ T_{s_0}, & s > s_0 \end{cases}, \gamma_s = \frac{1}{3L\alpha_s}, \text{ and } p_s = \frac{1}{2}, \text{ with} \quad (2.3)$$

$$\alpha_s = \begin{cases} \frac{1}{2}, & s \leq s_0 \\ \frac{2}{s-s_0+4}, & s > s_0 \end{cases}. \quad (2.4)$$

Then the total number of gradient evaluations of f_i performed by Algorithm 1 to find a stochastic solution of (1) i.e., a point $\bar{x} \in X$ s.t. $\mathbb{E}[\psi(\bar{x}) - \psi^*] \leq \epsilon$, can be bounded by

	Katyusha ^{ns} [1]	$\mathcal{O} \left\{ \frac{m}{\sqrt{\epsilon}} + \sqrt{\frac{mL}{\epsilon}} \right\}$
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$$\bar{N} := \begin{cases} \mathcal{O} \left\{ m \log \frac{D_0}{\epsilon} \right\}, & m \geq D_0/\epsilon, \\ \mathcal{O} \left\{ m \log m + \sqrt{\frac{mD_0}{\epsilon}} \right\}, & m < D_0/\epsilon, \end{cases} \quad (2.5)$$

where D_0 is defined as

$$D_0 := 2[\psi(x^0) - \psi(x)] + 3LV(x^0, x). \quad (2.6)$$

○ **Varag solves smooth problem directly!**

- Doubling epoch length of inner loop
- When the required accuracy ϵ is low and/or the number of components m is large, *Varag* achieves a fast **linear rate of convergence**
- Otherwise, *Varag* achieves an **optimal sublinear rate of convergence**

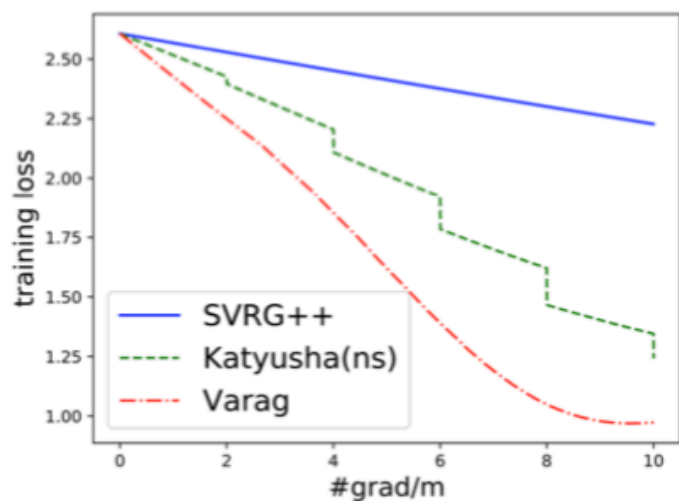
- *Varag* is the first accelerated RIG in the literature to obtain such convergence results by directly solving smooth finite-sum optimization problems.



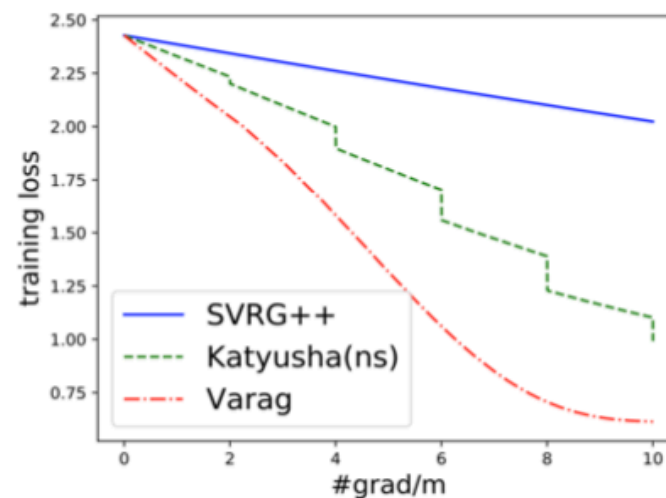
One numerical example – unconstrained logistic models



$$\min_{x \in \mathbb{R}^n} \left\{ \psi(x) := \frac{1}{m} \sum_{i=1}^m f_i(x) \right\} \text{ where } f_i(x) := \log(1 + \exp(-b_i a_i^T x))$$



Diabetes ($m = 1151$), unconstrained logistic



Breast Cancer Wisconsin ($m = 683$), unconstrained logistic

When $\mu \approx 0$ for strongly convex problems...



When the problem is almost not strongly convex, i.e., $\mu \approx 0$, $\sqrt{mL}/\mu \log(1/\epsilon)$ will be dominating and tend to ∞ as μ decreases.

Therefore, these complexity bounds are significantly worse than simply treating the problem as smooth convex problems.

Algorithms	Deterministic smooth strongly convex
RPDG[18]	$\mathcal{O} \left\{ (m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon} \right\}$
Catalyst[20]	$\mathcal{O} \left\{ (m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon} \right\}^1$
Katyusha[1]	$\mathcal{O} \left\{ (m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon} \right\}$
Katyusha ^{ns} [1]	NA
RGEM[19]	$\mathcal{O} \left\{ (m + \sqrt{\frac{mL}{\mu}}) \log \frac{1}{\epsilon} \right\}$

$$\bar{N} := \begin{cases} \mathcal{O} \left\{ m \log \frac{D_0}{\epsilon} \right\}, & m \geq D_0/\epsilon, \\ \mathcal{O} \left\{ m \log m + \sqrt{\frac{mD_0}{\epsilon}} \right\}, & m < D_0/\epsilon, \end{cases}$$

Theorem 2 (A unified result for convex finite-sum optimization) Suppose that the probabilities q_i 's are set to $L_i/\sum_{i=1}^m L_i$ for $i = 1, \dots, m$. Moreover, let us denote $s_0 := \lceil \log m \rceil + 1$ and assume that the weights $\{\theta_t\}$ are set to (2.2) if $1 \leq s \leq s_0$ or $s_0 < s \leq s_0 + \sqrt{\frac{12L}{m\mu}} - 4$, $m < \frac{3L}{4\mu}$. Otherwise, they are set to

$$\theta_t = \begin{cases} \Gamma_{t-1} - (1 - \alpha_s - p_s)\Gamma_t, & 1 \leq t \leq T_s - 1, \\ \Gamma_{t-1}, & t = T_s, \end{cases} \quad (2.7)$$

where $\Gamma_t = (1 + \mu\gamma_s)^t$. If the parameters $\{T_s\}$, $\{\gamma_s\}$ and $\{p_s\}$ set to (2.3) with

$$\alpha_s = \begin{cases} \frac{1}{2}, & s \leq s_0, \\ \max \left\{ \frac{2}{s-s_0+4}, \min \left\{ \sqrt{\frac{m\mu}{3L}}, \frac{1}{2} \right\} \right\}, & s > s_0, \end{cases} \quad (2.8)$$

then the total number of gradient evaluations of f_i performed by Algorithm 1 to find a stochastic ϵ -solution of (1.1) can be bounded by

$$\bar{N} := \begin{cases} \mathcal{O} \left\{ m \log \frac{D_0}{\epsilon} \right\}, & m \geq \frac{D_0}{\epsilon} \text{ or } m \geq \frac{3L}{4\mu}, \\ \mathcal{O} \left\{ m \log m + \sqrt{\frac{mD_0}{\epsilon}} \right\}, & m < \frac{D_0}{\epsilon} \leq \frac{3L}{4\mu}, \\ \mathcal{O} \left\{ m \log m + \sqrt{\frac{mL}{\mu}} \log \frac{D_0/\epsilon}{3L/4\mu} \right\}, & m < \frac{3L}{4\mu} \leq \frac{D_0}{\epsilon}. \end{cases} \quad (2.9)$$

when $f(y) \geq f(x) + \langle \nabla f(x), y - x \rangle + \mu V(x, y), \forall x, y \in X$.

- **Varag is an unified optimal method!**
 - When μ is large enough, *Varag* achieves the optimal linear rate of convergence
 - When μ is relatively small, *Varag* treats the problem as a smooth convex problem.
- *Varag* does not require to know the target accuracy ϵ and the constant D_0 beforehand to obtain optimal convergent rates.
- The unified step-size policy adjusts itself to the value of the condition number
- *Varag* does not assume the strong convexity comes from the regularizer!

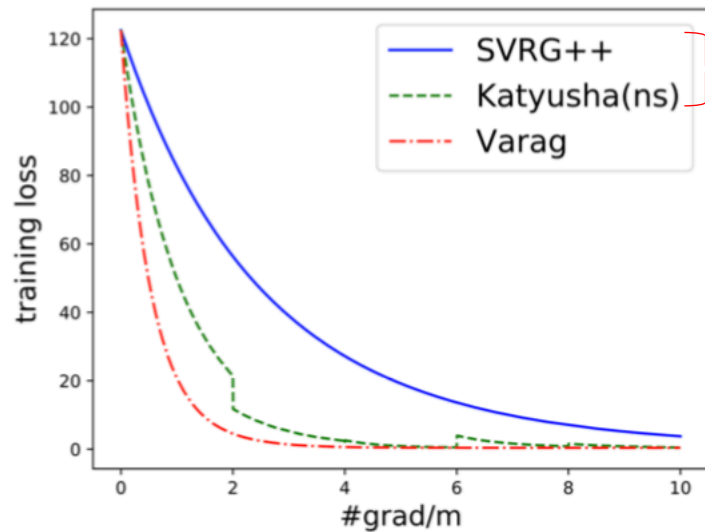
Varag as an unified optimal method



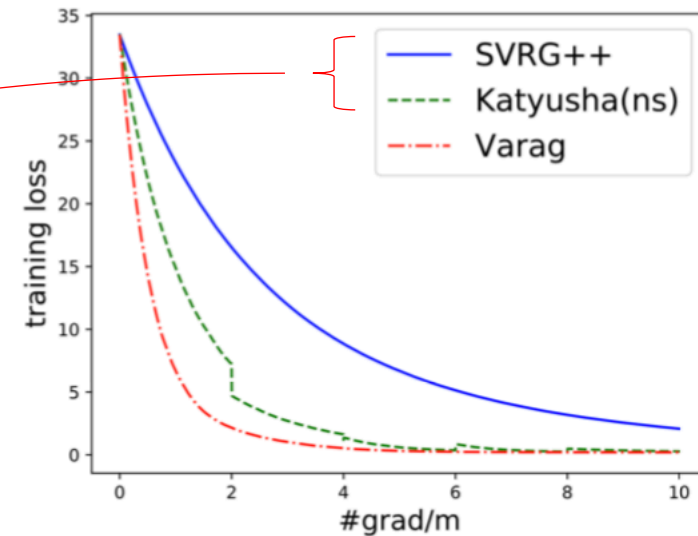
One numerical example – Lasso regression models



$$\min_{x \in \mathbb{R}^n} \{ \psi(x) := \frac{1}{m} \sum_{i=1}^m f_i(x) + h(x) \} \text{ where } f_i(x) := \frac{1}{2} (a_i^T x - b_i)^2, h(x) := \lambda \|x\|_1.$$



Diabetes ($m = 1151$),
Lasso $\lambda = 0.001$



Breast Cancer Wisconsin ($m = 683$),
Lasso $\lambda = 0.001$

Treats it as
smooth convex
problems

$$V(x, X^*) \leq \frac{1}{\bar{\mu}}(\psi(x) - \psi^*), \quad \forall x \in X,$$

Application examples:
 Linear systems,
 quadratic programs,
 linear matrix
 inequalities and
 composite problems,
 etc.

Theorem 3 (Convex finite-sum optimization under error bound) Assume f_i 's are set to $L_i/\sum_{i=1}^m L_i$ for $i = 1, \dots, m$, and θ_t are defined as (2.1) with parameters $\{\gamma_s\}$, $\{p_s\}$ and $\{\alpha_s\}$ as in (2.3) and (2.4), $s = 4 + 4\sqrt{\frac{L}{\bar{\mu}m}}$,

Moreover, if we restart Varag every time it runs s iterations for $k = \log \frac{\psi(x^0) - \psi^*}{\epsilon}$ number of gradient evaluations of f_i to find a stochastic ϵ -solution of (1.1),

$$\bar{N} := k(\sum_s (m + T_s)) = \mathcal{O} \left\{ \left(m + \sqrt{\frac{mL}{\bar{\mu}}} \right) \log \frac{\psi(x^0) - \psi(x^*)}{\epsilon} \right\}. \quad (2.13)$$

Varag is the first randomized method to establish the accelerated linear rate of convergence for solving the above problems!

Generalization of *Varag*

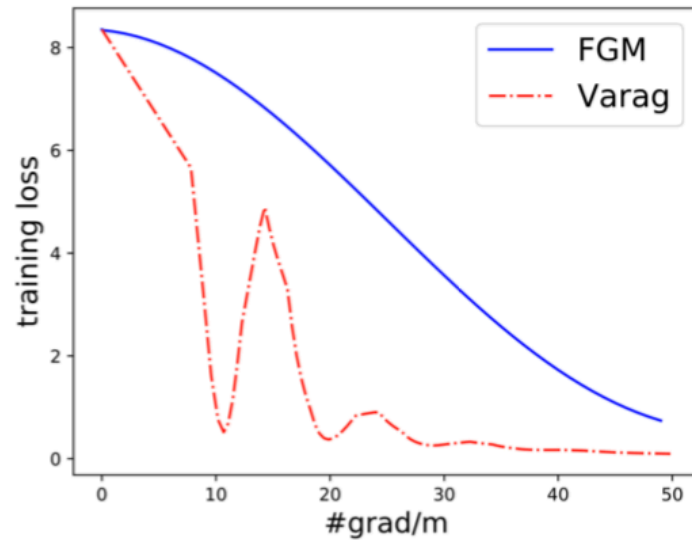


Finite-sum under
 error bound condition

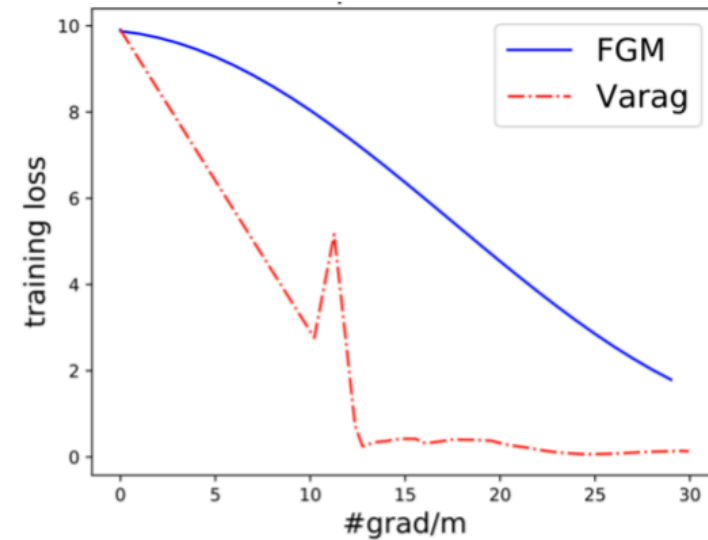
One numerical example – quadratic problems



$$\min_{x \in \mathbb{R}^n} \{ \psi(x) := \frac{1}{m} \sum_{i=1}^m f_i(x) \} \text{ where } f_i(x) := \frac{1}{2} x^T Q_i x + q_i^T x.$$



Diabetes ($m = 1151$)



Parkinsons Telemonitoring ($m = 5875$)

Only noisy gradient information can be accessed via SFO

$$\mathbb{E}_{\xi_j} [G_i(x, \xi_j)] = \nabla f_i(x), \quad i = 1, \dots, m,$$

$$\mathbb{E}_{\xi_j} [\|G_i(x, \xi_j) - \nabla f_i(x)\|_*^2] \leq \sigma^2, \quad i = 1, \dots, m.$$

Algorithm 2 Stochastic accelerated variance-reduced stochastic gradient descent (Stochastic Varag)

This algorithm is the same as Algorithm 1 except that for given batch-size parameters B_s and b_s , Line 3 is replaced by $\tilde{x} = \tilde{x}^{s-1}$ and

$$\tilde{g} = \frac{1}{m} \sum_{i=1}^m \left\{ G_i(\tilde{x}) := \frac{1}{B_s} \sum_{j=1}^{B_s} G_i(\tilde{x}, \xi_j^s) \right\}, \quad (2.16)$$

and Line 8 is replaced by

$$G_t = \frac{1}{q_{i_t} m b_s} \sum_{k=1}^{b_s} (G_{i_t}(\underline{x}_t, \xi_k^s) - G_{i_t}(\tilde{x})) + \tilde{g}. \quad (2.17)$$

Generalization of *Varag*



Stochastic finite-sum

Theorem 4 (Stochastic smooth finite-sum optimization) Assume that θ_t are defined as in (2.2), $C := \sum_{i=1}^m \frac{1}{q_i m^2}$ and the probabilities q_i 's are set to $L_i / \sum_{i=1}^m L_i$ for $i = 1, \dots, m$. Moreover, let us denote $s_0 := \lfloor \log m \rfloor + 1$ and set T_s, α_s, γ_s and p_s as in (2.3) and (2.4). Then the number of calls to the SFO oracle required by Algorithm 2 to find a stochastic ϵ -solution of (1.1) can be bounded by

$$N_{\text{SFO}} = \sum_s (mB_s + T_s b_s) = \begin{cases} \mathcal{O} \left\{ \frac{mC\sigma^2}{L\epsilon} \right\}, & m \geq D_0/\epsilon, \\ \mathcal{O} \left\{ \frac{C\sigma^2 D_0}{L\epsilon^2} \right\}, & m < D_0/\epsilon, \end{cases} \quad (2.18)$$

where D_0 is given in (2.6).

Varag is **the first** to achieve the above complexity results for smooth convex problems!

- RGEM[LZ18] achieves nearly optimal rate $\tilde{\mathcal{O}}\{\sigma^2 / \mu^2 \epsilon\}$ for expected distance between the output and the optimal solution
- Variant of SVRG[KM19] achieves $\mathcal{O}\{m \log m + \sigma^2 / \epsilon\}$ with some specific initial point.

Generalization of *Varag*



Stochastic finite-sum

Future works



- Extend *Varag* to solve nonconvex finite-sum problems
- How to choose stepsize if L and μ are hard to estimate?

A dark blue, irregularly shaped graphic with a splatter effect, containing the text "Thank you!" in white. The graphic has a rough, hand-painted appearance with various shades of blue and white splatters around its edges. The text is centered within the dark blue area.

Thank you!

References

Varag: Lan, G., Li, Z., & Zhou, Y. (2019). *A unified variance-reduced accelerated gradient method for convex optimization*. *arXiv preprint arXiv:1905.12412*. Accepted by NeurIPS 2019.

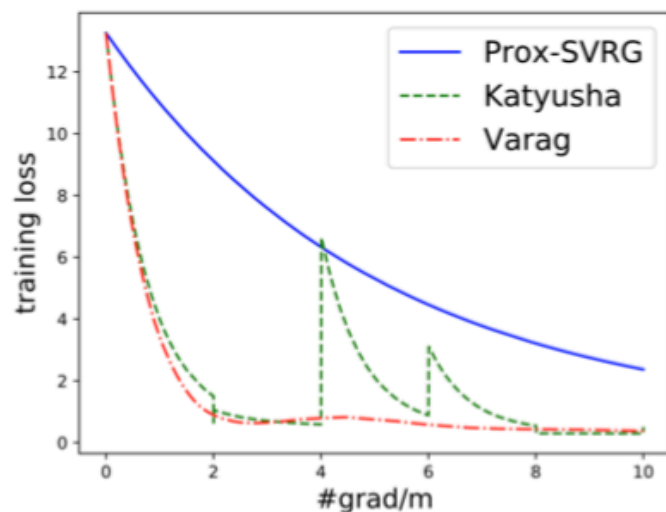
Other RIG methods:

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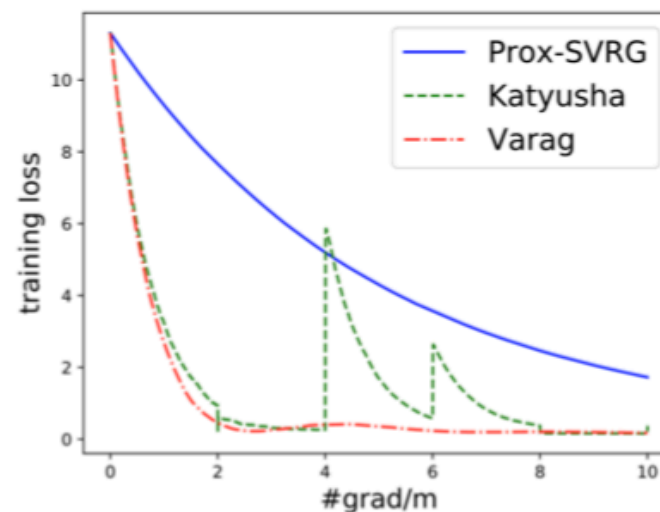
One numerical example – ridge regression models



$$\min_{x \in \mathbb{R}^n} \{ \psi(x) := \frac{1}{m} \sum_{i=1}^m f_i(x) + h(x) \} \text{ where } f_i(x) := \frac{1}{2} (a_i^T x - b_i)^2, h(x) := \lambda \|x\|_2^2.$$



Diabetes ($m = 1151$), ridge $\lambda = 10^{-6}$



Breast-Cancer-Wisconsin ($m = 683$), ridge $\lambda = 10^{-6}$

Varag requires less CPU time per training epoch than Katyusha!