

Stochastic Newton and Cubic Newton Methods with Simple Local Linear-Quadratic Rates

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Problem

We want to solve the **finite-sum optimization** problem

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

- Problem (1) has many applications in machine learning, data science and engineering.
- We focus on the regime when n is **very large**. This is typically the case in big data settings (e.g., massively distributed and federated learning).

Assumptions

- Each f_i is μ -strongly convex ($\mu > 0$), i.e.,
$$\mu I_d \preceq \nabla^2 f_i(x), \forall x \in \mathbb{R}^d,$$
- Each f_i has H -Lipschitz Hessian, i.e.,
$$\|\nabla^2 f_i(x) - \nabla^2 f_i(y)\| \leq H\|x - y\|, \forall x, y \in \mathbb{R}^d.$$

Our Contribution

We develop two new simple and fundamental stochastic second-order methods:

- Stochastic Newton (SN) method,
- Stochastic Cubic Newton (SCN) method.

Our methods have highly desirable properties:

- **Cost of 1 iteration does not depend on n ,**
- **(Local) convergence rate does not depend on the conditioning of the problem.**

Motivation I: The Curse of 1st Order Methods

In this regime, the state of the art methods for solving (1) are variants of **stochastic gradient descent**. However, the performance of all first order methods depends heavily on the **conditioning** of the problem. Various strategies have been proposed to address this problem:

- preconditioning (e.g., data normalization, layer and batch normalization),
- momentum (e.g., Polyak and Nesterov),
- adaptive stepsizes (e.g., Adagrad, ADAM, Barzilai-Borwein, Malitsky-Mishchenko) and line search,
- minibatching and importance sampling.

Some of these methods reduce the effect of conditioning provably, and some are heuristics which often work but sometimes fail. However, **first order methods are inherently incapable to remove the effect of conditioning.**

Newton's Method

It will be useful to recall classical **Newton's method**:

$$\begin{aligned} x^{k+1} &= x^k - \left[\frac{1}{n} \sum_{i=1}^n \nabla^2 f_i(x^k) \right]^{-1} \nabla f(x^k) \\ &= \left[\frac{1}{n} \sum_{i=1}^n \nabla^2 f_i(x^k) \right]^{-1} \left[\frac{1}{n} \sum_{i=1}^n \nabla^2 f_i(x^k) x^k - \nabla f(x^k) \right]. \end{aligned} \quad (2)$$

Fact. If f is μ -strongly convex and its Hessian is Lipschitz continuous, then if $\|x^0 - x^*\|$ is small enough (where $x^* = \arg \min f$), then the iterates (2) of Newton's method converge to x^* **quadratically**: $\|x^{k+1} - x^*\| \leq \frac{1}{2} \|x^k - x^*\|^2$. This means that

$$k \geq \mathcal{O} \left(\log_2 \log_2 \frac{1}{\epsilon} \right) \Rightarrow \|x^k - x^*\| \leq \epsilon.$$

Motivation II: Issues with Existing Stochastic 2nd Order Methods

Because of what we said above, there is **a lot of effort to develop efficient stochastic 2nd order methods**.

Almost every such method has the form

$$x^{k+1} = x^k - (H^k)^{-1} g^k,$$

where $H^k \approx \nabla^2 f(x^k)$ is a stochastic approximation of the Hessian and $g^k \approx \nabla f(x^k)$ is a stochastic gradient approximation of the gradient. Most methods let

$$g^k = \frac{1}{|S_g^k|} \sum_{i \in S_g^k} \nabla f_i(x^k), \quad H^k = \frac{1}{|S_H^k|} \sum_{i \in S_H^k} \nabla^2 f_i(x^k),$$

where S_H^k and S_g^k are suitably chosen random subsets of $\{1, 2, \dots, n\}$. However, all these methods suffer from severe issues:

- they require $\mathcal{O}(\epsilon^{-1})$ or even $\mathcal{O}(\epsilon^{-2})$ samples **in each iteration**, where ϵ is the target accuracy. The number of required samples often exceeds n in theory,
- the resulting convergence rate is **worse** than the rate of first order methods.

New Algorithm: Stochastic Newton

Algorithm 1: Stochastic Newton (SN)

Initialize: Choose $w_1^0, w_2^0, \dots, w_n^0 \in \mathbb{R}^d$ and

$\tau \in \{1, 2, \dots, n\}$

for $k = 0, 1, \dots$ **do**

 Compute Hessian estimator: $H^k = \frac{1}{n} \sum_{i=1}^n \nabla^2 f_i(w_i^k)$;

 Compute gradient estimator: $g^k = \frac{1}{n} \sum_{i=1}^n \nabla f_i(w_i^k)$;

$x^{k+1} = [H^k]^{-1} \left[\frac{1}{n} \sum_{i=1}^n \nabla^2 f_i(w_i^k) w_i^k - g^k \right]$;

 Choose a random set $S^k \subseteq \{1, \dots, n\}$ of cardinality τ ;

 Update

$$w_i^{k+1} = \begin{cases} x^{k+1}, & i \in S^k \\ w_i^k, & i \notin S^k \end{cases}$$

end

Theorem 1 (Stochastic Newton)

The random iterates of SN (Algorithm 1) satisfy the recursion

$$\mathbb{E}_k [\mathcal{W}^{k+1}] \leq \left(1 - \frac{\tau}{n} + \frac{\tau}{n} \left(\frac{H}{2\mu} \right)^2 \right) \mathcal{W}^k,$$

where $\mathcal{W}^k \stackrel{\text{def}}{=} \frac{1}{n} \sum_{i=1}^n \|w_i^k - x^*\|^2$. Furthermore, if $\|w_i^0 - x^*\| \leq \frac{\mu}{H}$ for $i = 1, \dots, n$, then

$$\mathbb{E}_k [\mathcal{W}^{k+1}] \leq \left(1 - \frac{3\tau}{4n} \right) \mathcal{W}^k.$$

New Algorithm: Stochastic Cubic Newton

Motivation: Newton method converges only locally. To fix it, one needs cubic regularization:

$$\Phi^k(x) \stackrel{\text{def}}{=} \langle \nabla f(x^k), x - x^k \rangle + \frac{1}{2} \|x - x^k\|_{\nabla^2 f(x^k)}^2$$

$$x^{k+1} = \operatorname{argmin}_x \left\{ \Phi^k(x) + \frac{M}{6} \|x - x^k\|^3 \right\}.$$

Our algorithm

$$\Psi^k \stackrel{\text{def}}{=} \left\langle \frac{1}{n} \sum_{i=1}^n \nabla f_i(w_i^k), x - x^k \right\rangle + \frac{1}{2n} \sum_{i=1}^n \|x - w_i^k\|_{\nabla^2 f_i(w_i^k)}^2$$

$$x^{k+1} = \operatorname{argmin}_x \left\{ \Psi^k(x) + \frac{M}{6n} \sum_{i=1}^n \|x - w_i^k\|^3 \right\}.$$

Theorem 2 (Stochastic Cubic Newton)

The random iterates of SCN satisfy the recursion

$$\mathbb{E}_k [\mathcal{V}^{k+1}] \leq \left(1 - \frac{\tau}{n} + \frac{\tau}{n} \left(\frac{(M+H)\sqrt{2}}{3\mu^{\frac{3}{2}}} \right)^{3/2} \sqrt{\mathcal{V}^k} \right) \mathcal{V}^k,$$

where $\mathcal{V}^k \stackrel{\text{def}}{=} \frac{1}{n} \sum_{i=1}^n (f(w_i^k) - f(x^*))^{\frac{3}{2}}$. Furthermore, if the vectors w_i^k are close enough to x^* , then

$$\mathbb{E}_k [\mathcal{V}^{k+1}] \leq \left(1 - \frac{\tau}{2n} \right) \mathcal{V}^k.$$

Experiments

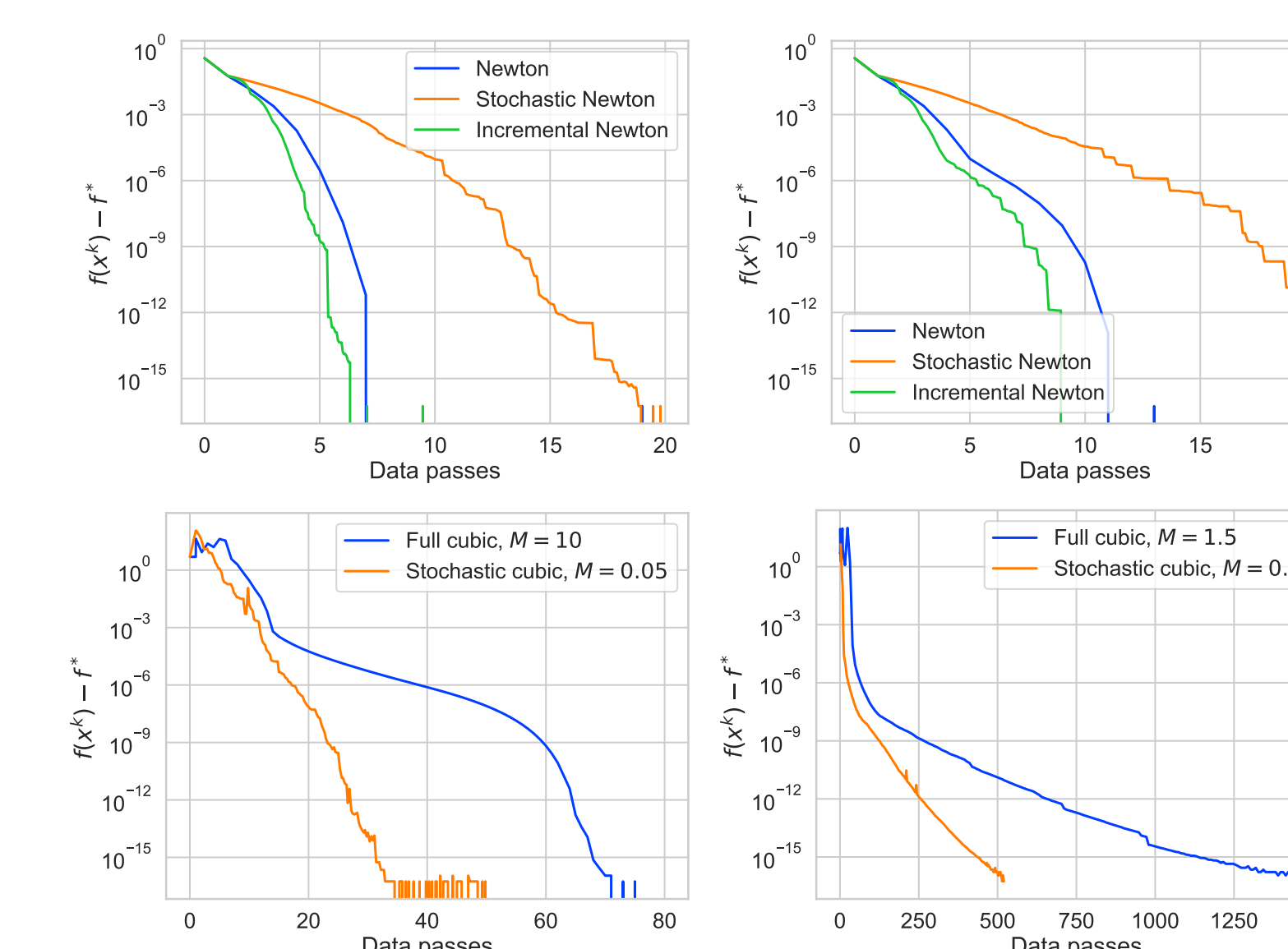


Figure 1: Logistic regression. Top: Newton methods, bottom: cubic Newton methods. Left: $\mu = \frac{1}{100n}$, right: $\mu = \frac{1}{10000n}$.