A Generic Quasi-Newton Algorithm for Faster Gradient-Based Optimization

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An alternate title: Acceleration by Smoothing

Collaborators



Publications and pre-prints

H. Lin, J. Mairal and Z. Harchaoui. A Generic Quasi-Newton Algorithm for Faster Gradient-Based Optimization. arXiv:1610.00960. 2017

C. Paquette, H. Lin, D. Drusvyatskiy, J. Mairal, Z. Harchaoui. Catalyst Acceleration for Gradient-Based Non-Convex Optimization. *arXiv:1703.10993.* 2017

H. Lin, J. Mairal and Z. Harchaoui. A Universal Catalyst for First-Order Optimization. *Adv. NIPS* 2015.

Minimizing large finite sums

Consider the minimization of a large sum of convex functions

$$\min_{x\in\mathbb{R}^d}\left\{f(x)\triangleq\frac{1}{n}\sum_{i=1}^n f_i(x)+\psi(x)\right\},\,$$

where each f_i is *L*-smooth and convex and ψ is a convex regularization penalty but not necessarily differentiable.

Motivation

	Composite	Finite sum	Exploit "curvature"
First-order methods	 ✓ 		
Quasi-Newton			

[Nesterov, 2013, Wright et al., 2009, Beck and Teboulle, 2009, Chambolle and Pock, 2011, Combettes and Wajs, 2005],...

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[Schmidt et al., 2017, Xiao and Zhang, 2014, Defazio et al., 2014a,b, Shalev-Shwartz and Zhang, 2012, Mairal, 2015, Zhang and Xiao, 2015]

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Expected number of gradients ∇f_i to compute to guarantee $f(x_k) - f^* \leq \varepsilon$, when the objective f is μ -strongly convex: • accelerated proximal gradient: $O\left(n\sqrt{\frac{L_f}{\mu}}\log\left(\frac{1}{\varepsilon}\right)\right)$; • incremental gradient methods: $O\left(\left(n + \frac{L}{\mu}\right)\log\left(\frac{1}{\varepsilon}\right)\right)$.

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[Byrd et al., 2015, Lee et al., 2012, Scheinberg and Tang, 2016, Yu et al., 2008, Ghadimi et al., 2015, Stella et al., 2016],...

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Quasi-Newton		×	✓

[Byrd et al., 2016, Gower et al., 2016]

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Motivation



QuickeNing: main idea (an old one)

Idea: Smooth the function and then apply Quasi-Newton.

• The strategy appears in early work about variable metric bundle methods. [Chen and Fukushima, 1999, Fukushima and Qi, 1996, Mifflin, 1996, Fuentes, Malick, and Lemaréchal, 2012, Burke and Qian, 2000] ...

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The Moreau-Yosida envelope

Given $f : \mathbb{R}^d \to \mathbb{R}$ a convex function, the Moreau-Yosida envelope of f is the function $F : \mathbb{R}^d \to \mathbb{R}$ defined as

$$F(x) = \min_{w \in \mathbb{R}^d} \left\{ f(w) + \frac{\kappa}{2} \|w - x\|^2 \right\}.$$

The proximal operator p(x) is the unique minimizer of the problem.

The Moreau-Yosida regularization

$$F(x) = \min_{w \in \mathbb{R}^d} \left\{ f(w) + \frac{\kappa}{2} \|w - x\|^2 \right\}.$$

Basic properties [see Lemaréchal and Sagastizábal, 1997]

• Minimizing f and F is equivalent in the sense that

$$\min_{x\in\mathbb{R}^d}F(x)=\min_{x\in\mathbb{R}^d}f(x),$$

and the solution set of the two problems coincide with each other.

• F is continuously differentiable even when f is not and

$$\nabla F(x) = \kappa(x - p(x)).$$

In addition, ∇F is Lipschitz continuous with parameter $L_F = \kappa$.

• If f is μ -strongly convex then F is also strongly convex with parameter $\mu_F = \frac{\mu\kappa}{\mu+\kappa}$.

The Moreau-Yosida regularization

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F enjoys nice properties: smoothness, (strong) convexity and we can control its condition number $1 + \kappa/\mu$.

A fresh look at Catalyst [Lin et al., 2015]



A fresh look at the proximal point algorithm

A naive approach consists of minimizing the smoothed objective F instead of f with a method designed for smooth optimization.

Consider indeed

$$x_{k+1} = x_k - \frac{1}{\kappa} \nabla F(x_k).$$

By rewriting the gradient $\nabla F(x_k)$ as $\kappa(x_k - p(x_k))$, we obtain

$$x_{k+1} = p(x_k) = \operatorname*{arg\,min}_{w \in \mathbb{R}^p} \left\{ f(w) + \frac{\kappa}{2} \|w - x_k\|^2 \right\}.$$

This is exactly the **proximal point algorithm** [Martinet, 1970, Rockafellar, 1976].

A fresh look at the accelerated proximal point algorithm

Consider now

$$x_{k+1} = y_k - rac{1}{\kappa}
abla F(y_k)$$
 and $y_{k+1} = x_{k+1} + eta_{k+1}(x_{k+1} - x_k),$

where β_{k+1} is a Nesterov-like extrapolation parameter. We may now rewrite the update using the value of ∇F , which gives:

$$x_{k+1} = p(y_k)$$
 and $y_{k+1} = x_{k+1} + \beta_{k+1}(x_{k+1} - x_k)$

This is the accelerated proximal point algorithm of Güler [1992].

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Remarks

- *F* may be better conditioned than *f* when $1 + \kappa/\mu \le L/\mu$;
- Computing $p(y_k)$ has a cost!

A fresh look at Catalyst [Lin, Mairal, and Harchaoui, 2015]

Catalyst is a particular accelerated proximal point algorithm with inexact gradients [Güler, 1992].

$$x_{k+1} \approx p(y_k)$$
 and $y_{k+1} = x_{k+1} + \beta_{k+1}(x_{k+1} - x_k)$

The quantity x_{k+1} is obtained by using an optimization method M for approximately solving:

$$x_{k+1} pprox rgmin_{w \in \mathbb{R}^p} \left\{ f(w) + rac{\kappa}{2} \|w - y_k\|^2
ight\},$$

Catalyst provides Nesterov's acceleration to ${\mathcal M}$ with...

- restart strategies for solving the sub-problems;
- global complexity analysis resulting in theoretical acceleration;
- optimal balancing between outer and inner computations.

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see also [Frostig et al., 2015]
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Limited-Memory BFGS (L-BFGS)

Pros

• one of the largest practical success of smooth optimization.

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Cons

- worst-case convergence rates for strongly-convex functions are linear, but much worse than the gradient descent method.
- proximal variants typically requires solving many times

$$\min_{x\in\mathbb{R}^d}\frac{1}{2}(x-z)B_k(z-z)+\psi(x).$$

• no guarantee of approximating the Hessian.

QuickeNing

Main recipe

- L-BFGS applied to the **smoothed objective** *F* with **inexact gradients** [see Friedlander and Schmidt, 2012].
- inexact gradients are obtained by solving sub-problems using a first-order optimization method *M*;
- ideally, \mathcal{M} is able to adapt to the problem structure (finite sum, composite regularization).
- replace L-BFGS steps by proximal point steps if no sufficient decrease is estimated ⇒ no line search on F;

Obtaining inexact gradients

Algorithm Procedure ApproxGradient

input Current point x in \mathbb{R}^d ; smoothing parameter $\kappa > 0$.

1: Compute the approximate mapping using an optimization method \mathcal{M} :

$$z pprox rgmin_{w \in \mathbb{R}^d} \left\{ h(w) \stackrel{\scriptscriptstyle \Delta}{=} f(w) + rac{\kappa}{2} \|w - x\|^2
ight\},$$

2: Estimate the gradient $\nabla F(x)$

$$g=\kappa(x-z).$$

output approximate gradient estimate g, objective value $F_a \triangleq h(z)$, proximal mapping z.

Algorithm QuickeNing

input x_0 in \mathbb{R}^p ; number of iterations K; $\kappa > 0$; minimization algorithm \mathcal{M} .

- 1: Initialization: $(g_0, F_0, z_0) = \text{ApproxGradient}(x_0, \mathcal{M}); B_0 = \kappa I.$
- 2: for k = 0, ..., K 1 do
- 3: Perform the Quasi-Newton step

$$egin{aligned} & x_{ ext{test}} = x_k - B_k^{-1} g_k \ & (g_{ ext{test}}, F_{ ext{test}}, z_{ ext{test}}) = ext{ApproxGradient}\left(x_{ ext{test}}, \mathcal{M}
ight) \,. \end{aligned}$$

4: **if**
$$F_{\text{test}} \leq F_k - \frac{1}{2\kappa} ||g_k||^2$$
, **then**

5:
$$(x_{k+1}, g_{k+1}, F_{k+1}, z_{k+1}) = (x_{\text{test}}, g_{\text{test}}, F_{\text{test}}, z_{\text{test}}).$$

- 6: **else**
- 7: Update the current iterate with the last proximal mapping:

$$egin{aligned} & x_{k+1} = z_k = x_k - (1/\kappa)g_k \ & (g_{k+1}, \mathcal{F}_{k+1}, z_{k+1}) = \operatorname{\mathsf{ApproxGradient}}\left(x_{k+1}, \mathcal{M}
ight) \,. \end{aligned}$$

8: end if

9: update
$$B_{k+1} = L\text{-BFGS}(B_k, x_{k+1} - x_k, g_{k+1} - g_k)$$
.

10: end for

output last proximal mapping $z_{\mathcal{K}}$ (solution).

Algorithm QuickeNing

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The main characters:

- the sequence $(x_k)_{k\geq 0}$ that minimizes F;
- the sequence $(z_k)_{k\geq 0}$ produced by \mathcal{M} that minimizes f;
- the gradient approximations $g_k \approx \nabla F(x_k)$;
- the function value approximations $F_k \approx F(x_k)$;
- an L-BFGS update with inexact gradients;
- an approximate sufficient descent condition.

10: **епа тог**

output last proximal mapping z_K (solution).

Requirements on $\mathcal M$ and restarts

$\mathsf{Method}\ \mathcal{M}$

• Say a sub-problem consists of minimizing h; we want \mathcal{M} to produce a sequence of iterates $(w_t)_{t\geq 0}$ with linear convergence rate

$$h(w_t) - h^* \leq C_{\mathcal{M}}(1 - \tau_{\mathcal{M}})^t (h(w_0) - h^*).$$

Restarts

• When f is smooth, we initialize $w_0 = x$ when solving

$$\min_{w\in\mathbb{R}^d}\left\{f(w)+\frac{\kappa}{2}\|w-x\|^2\right\}.$$

• When $f = f_0 + \psi$ is composite, we use the initialization

$$w_0 = \operatorname*{arg\,min}_{w \in \mathbb{R}^d} \left\{ f_0(x) + \langle \nabla f_0(x), w - x \rangle + \frac{L + \kappa}{2} \|w - x\|^2 + \psi(w) \right\}$$

When do we stop the method \mathcal{M} ?

Three strategies to balance outer and inner computations

- (a) use a **pre-defined sequence** $(\varepsilon_k)_{k\geq 0}$ and stop the optimization method \mathcal{M} when the approximate proximal mapping is ε_k -accurate.
- (b) define an **adaptive stopping criterion** that depends on quantities that are available at iteration *k*.
- (c) use a **pre-defined budget** T_M of iterations of the method M for solving each sub-problem.

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- (c) use a **pre-defined budget** T_M of iterations of the method M for solving each sub-problem.

Remarks

- (a) is the less practical strategy.
- (b) is simpler to use and conservative (compatible with theory).
- (c) requires $T_{\mathcal{M}}$ to be large enough in theory. The aggressive strategy $T_{\mathcal{M}} = n$ for an incremental method is extremely simple to use and effective in practice.

When do we stop the method \mathcal{M} ?

Three strategies for μ -strongly convex objectives f

(a) use a **pre-defined sequence** $(\varepsilon_k)_{k\geq 0}$ and stop the optimization method \mathcal{M} when the approximate proximal mapping is ε_k -accurate.

$$arepsilon_k = rac{1}{2}C(1-
ho)^{k+1} \quad ext{with} \quad C \geq f(x_0) - f^* \; ext{ and } \;
ho = rac{\mu}{4(\mu+\kappa)}.$$

(b) For minimizing $h(w) = f(w) + (\kappa/2) ||w - x||^2$, stop when

$$h(w_t)-h^{\star}\leq \frac{\kappa}{36}\|w_t-x\|^2.$$

(c) use a **pre-defined budget** T_M of iterations of the method M for solving each sub-problem with

$$T_{\mathcal{M}} = rac{1}{ au_{\mathcal{M}}} \log \left(19 C_{\mathcal{M}} rac{L + \kappa}{\kappa}
ight)$$
. (be more aggressive in practice)

Remarks and worst-case global complexity

Composite objectives and sparsity

Consider a composite problem with a sparse solution (e.g., $\psi = \ell_1$). The method produces two sequences $(x_k)_{k\geq 0}$ and $(z_k)_{k\geq 0}$;

- $F(x_k) \rightarrow F^*$, minimizes the smoothed objective \Rightarrow no sparsity;
- f(z_k) → f^{*}, minimizes the true objective ⇒ the iterates may be sparse if *M* handles composite optimization problems;

Global complexity

The number of iterations of \mathcal{M} to guarantee $f(z_k) - f^* \leq \varepsilon$ is at most

• $\tilde{O}(\frac{\mu+\kappa}{\tau_{M}\mu}\log(1/\varepsilon))$ for μ -strongly convex problems.

•
$$\tilde{O}(\frac{\kappa R^2}{\tau_{\mathcal{M}}\varepsilon})$$
 for convex problems.

Global Complexity and choice of κ

Example for gradient descent

With the right step-size, we have $\tau_M = (\mu + \kappa)/(L + \kappa)$ and the complexity for $\mu > 0$ becomes

$$ilde{O}\left(rac{L+\kappa}{\mu}\log(1/arepsilon)
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Example for SVRG for minimizing the sum of *n* functions $\tau_{\mathcal{M}} = \min(1/n, (\mu + \kappa)/(L + \kappa))$ and the complexity for $\mu > 0$ is

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QuickeNing does not provide any theoretical acceleration, but it does not degrade significantly the worst-case performance of \mathcal{M} (unlike L-BFGS vs gradient descent).

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Then, how to choose κ ? (i) assume that L-BFGS steps do as well as Nesterov. (ii) **choose** κ **as in Catalyst**.

Experiments: formulations

• ℓ_2 -regularized Logistic Regression:

$$\min_{x \in \mathbb{R}^d} \quad \frac{1}{n} \sum_{i=1}^n \log\left(1 + \exp(-b_i a_i^T x)\right) + \frac{\mu}{2} \|x\|^2,$$

• ℓ_1 -regularized Linear Regression (LASSO):

$$\min_{x\in\mathbb{R}^d} \quad \frac{1}{2n}\sum_{i=1}^n (b_i - a_i^T x)^2 + \lambda \|x\|_1,$$

• $\ell_1 - \ell_2^2$ -regularized Linear Regression (Elastic-Net):

$$\min_{x\in\mathbb{R}^d} \quad \frac{1}{2n}\sum_{i=1}^n (b_i - a_i^T x)^2 + \lambda \|x\|_1 + \frac{\mu}{2} \|x\|^2,$$

Experiments: Datasets

We consider four standard machine learning datasets with different characteristics in terms of size and dimension

name	covtype	alpha	real-sim	rcv1
n	581012	250 000	72 309	781 265
d	54	500	20 958	47 152

- we simulate the ill-conditioned regime $\mu = 1/(100n)$;
- λ for the Lasso leads to about 10% non-zero coefficients.

Experiments: QuickeNing-SVRG

We consider the methods

- SVRG: the Prox-SVRG algorithm of Xiao and Zhang [2014].
- Catalyst-SVRG: Catalyst applied to SVRG;
- L-BFGS (for smooth objectives): Mark Schmidt's implementation.
- **QuickeNing-SVRG1**: QuickeNing with aggressive strategy (c): one pass over the data in the inner loop.
- QuickeNing-SVRG2: strategy (b), compatible with theory.

We produce 12 figures (3 formulations, 4 datasets).

Experiments: QuickeNing-SVRG (log scale)



- QuickeNing-SVRG1 ≥ SVRG, QuickeNing-SVRG2;
- QuickeNing-SVRG2
 SVRG;
- QuickeNing-SVRG1 ≥ Catalyst-SVRG in 10/12 cases.

Experiments: QuickeNing-ISTA

We consider the methods

- ISTA: the proximal gradient descent method with line search.
- FISTA: the accelerated ISTA of Beck and Teboulle [2009].
- L-BFGS (for smooth objectives): Mark Schmidt's implementation.
- **QuickeNing-ISTA1**: QuickeNing with aggressive strategy (c): one pass over the data in the inner loop.
- QuickeNing-ISTA2: strategy (b), compatible with theory.

Experiments: QuickeNing-ISTA (log scale)



- L-BFGS (for smooth f) is slightly better than QuickeNing-ISTA1;
- QuickeNing-ISTA \geq or \gg FISTA in 11/12 cases.
- QuickeNing-ISTA1 \geq QuickeNing-ISTA2.

Experiments: Influence of κ



- κ₀ is the parameter (same as in Catalyst) used in all experiments;
- QuickeNing slows down when using κ > κ₀;
- here, for SVRG, QuickeNing is robust to small values of $\kappa!$

Experiments: Influence of /



- I = 100 in all previous experiments;
- I = 5 seems to be a reasonable choice in many cases, especially for sparse problems.

Conclusions and perspectives

- A simple generic Quasi-Newton method for composite functions, with simple sub-problems, and complexity guarantees.
- We also have a variant for dual approaches.
- Does not solve the gap between theory and practice for L-BFGS.

Perspectives

- QuickeNing-BCD, QuickeNing-SAG,SAGA,SDCA...
- Other types of smoothing? ⇒ Links with recent Quasi-Newton methods applied to other envelopes [Stella et al., 2016].

Outer-loop convergence analysis

Lemma: approximate descent property

$$F(x_{k+1}) \leq f(z_k) \leq F(x_k) - \frac{1}{4\kappa} \|\nabla F(x_k)\|_2^2 + 2\varepsilon_k.$$

Then, ε_k should be smaller than $\frac{1}{4\kappa} \|\nabla F(x_k)\|_2^2$, and indeed

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Then, ε_k should be smaller than $\frac{1}{4\kappa} \|\nabla F(x_k)\|_2^2$, and indeed Proposition: convergence with impractical ε_k and $\mu > 0$ If $\varepsilon_k \leq \frac{1}{16\kappa} \|\nabla F(x_k)\|_2^2$, define $\rho = \frac{\mu}{4(\mu+\kappa)}$, then $F(x_{k+1}) - F^* \leq f(z_k) - f^* \leq (1-\rho)^{k+1} (f(x_0) - f^*).$

Unfortunately, $\|\nabla F(x_k)\|$ is unknown.

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Unfortunately, $\|\nabla F(x_k)\|$ is unknown.

Lemma: convergence with adaptive ε_k and $\mu > 0$ If $\varepsilon_k \leq \frac{1}{36\kappa} \|g_k\|^2$, then $\varepsilon_k \leq \frac{1}{16} \|\nabla F(x_k)\|^2$.

This is strategy (b). g_k is known and easy to compute.

Inner-loop complexity analysis

Restart for L-smooth functions

For minimizing *h*, initialize the method \mathcal{M} with $w_0 = x$. Then,

$$h(w_0) - h^* \leq \frac{L+\kappa}{2\kappa^2} \|\nabla F(x)\|^2.$$
(1)

Proof.

We have the optimality condition $\nabla f(w^*) + \kappa(w^* - x) = 0$. As a result,

$$\begin{split} h(w_0) - h^* \\ &= f(x) - \left(f(w^*) + \frac{\kappa}{2} \| w^* - x \|^2 \right) \\ &\leq f(w^*) + \langle \nabla f(w^*), x - w^* \rangle + \frac{L}{2} \| x - w^* \|^2 - \left(f(w^*) + \frac{\kappa}{2} \| w^* - x \|^2 \right) \\ &= \frac{L + \kappa}{2} \| w^* - x \|^2 = \frac{L + \kappa}{2\kappa^2} \| \nabla F(x) \|^2. \end{split}$$

Presentation borrowed from Mark Schmidt, NIPS OPT 2010

• Quasi-Newton methods work with the parameter and gradient differences between successive iterations:

$$s_k \triangleq x_{k+1} - x_k, \quad y_k \triangleq \nabla f(x_{k+1}) - \nabla f(x_k).$$

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 Since B_{k+1} is not unique, the Broyden-Fletcher-Goldfarb-Shanno (BFGS) method chooses the symmetric matrix whose difference with B_k is minimal:

$$B_{k+1} = B_k - \frac{B_k s_k s_k B_k}{s_k B_k s_k} + \frac{y_k y_k^{\top}}{y_k^{\top} s_k}.$$

Presentation borrowed from Mark Schmidt, NIPS OPT 2010

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- We can solve a linear system involving these updates when B₀ is diagonal in O(dl) [Nocedal, 1980].

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