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

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## An alternate title: Acceleration by Smoothing

## Collaborators



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Courtney
Paquette

## 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.

## Focus of this work

## 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 $\psi$ is a convex regularization penalty but not necessarily differentiable.

## Motivation

|  | Composite | Finite sum | Exploit "curvature" |
| :---: | :---: | :---: | :---: |
| First-order methods | $\boldsymbol{\swarrow}$ |  |  |
| 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 $\nabla f_{i}$ to compute to guarantee $f\left(x_{k}\right)-f^{\star} \leq \varepsilon$, when the objective $f$ is $\mu$-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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[Byrd et al., 2016, Gower et al., 2016]

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where each $f_{i}$ is L-smooth and convex and $\psi$ is a convex regularization penalty but not necessarily differentiable.

## Motivation

Our goal is to

- accelerate first-order methods with Quasi-Newton heuristics;
- design algorithms that can adapt to composite and finite-sum structures and that can also exploit curvature information.
[Byrd et al., 2016, Gower et al., 2016]


## 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 $f: \mathbb{R}^{d} \rightarrow \mathbb{R}$ a convex function, the Moreau-Yosida envelope of $f$ is the function $F: \mathbb{R}^{d} \rightarrow \mathbb{R}$ 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, $\nabla F$ is Lipschitz continuous with parameter $L_{F}=\kappa$.

- If $f$ is $\mu$-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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In addition, $\nabla F$ is Lipschitz continuous with parameter $L_{F}=\kappa$.
$F$ 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

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

By rewriting the gradient $\nabla F\left(x_{k}\right)$ as $\kappa\left(x_{k}-p\left(x_{k}\right)\right)$, we obtain

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

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

## A fresh look at the accelerated proximal point algorithm

Consider now

$$
x_{k+1}=y_{k}-\frac{1}{\kappa} \nabla F\left(y_{k}\right) \quad \text { and } \quad y_{k+1}=x_{k+1}+\beta_{k+1}\left(x_{k+1}-x_{k}\right),
$$

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

$$
x_{k+1}=p\left(y_{k}\right) \quad \text { and } \quad y_{k+1}=x_{k+1}+\beta_{k+1}\left(x_{k+1}-x_{k}\right)
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This is the accelerated proximal point algorithm of Güler [1992].

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

- $F$ may be better conditioned than $f$ when $1+\kappa / \mu \leq L / \mu$;
- Computing $p\left(y_{k}\right)$ 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].

$$
x_{k+1} \approx p\left(y_{k}\right) \quad \text { and } \quad y_{k+1}=x_{k+1}+\beta_{k+1}\left(x_{k+1}-x_{k}\right)
$$

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

$$
x_{k+1} \approx \underset{w \in \mathbb{R}^{p}}{\arg \min }\left\{f(w)+\frac{\kappa}{2}\left\|w-y_{k}\right\|^{2}\right\},
$$

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.
see also [Frostig et al., 2015]


## Limited-Memory BFGS (L-BFGS)

Pros

- one of the largest practical success of smooth optimization.


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- one of the largest practical success of smooth optimization.


## 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 $\mathcal{M}$;
- 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 $\Rightarrow$ 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 $\mathcal{M}$ :

$$
z \approx \underset{w \in \mathbb{R}^{d}}{\arg \min }\left\{h(w) \triangleq f(w)+\frac{\kappa}{2}\|w-x\|^{2}\right\}
$$

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: $\left(g_{0}, F_{0}, z_{0}\right)=$ ApproxGradient $\left(x_{0}, \mathcal{M}\right) ; B_{0}=\kappa l$.
2: for $k=0, \ldots, K-1$ do
3: $\quad$ Perform the Quasi-Newton step

$$
\begin{aligned}
x_{\text {test }} & =x_{k}-B_{k}^{-1} g_{k} \\
\left(g_{\text {test }}, F_{\text {test }}, z_{\text {test }}\right) & =\text { ApproxGradient }\left(x_{\text {test }}, \mathcal{M}\right) .
\end{aligned}
$$

4: if $F_{\text {test }} \leq F_{k}-\frac{1}{2 \kappa}\left\|g_{k}\right\|^{2}$, then
5: $\quad\left(x_{k+1}, g_{k+1}, F_{k+1}, z_{k+1}\right)=\left(x_{\text {test }}, g_{\text {test }}, F_{\text {test }}, z_{\text {test }}\right)$.
6: else
7: Update the current iterate with the last proximal mapping:

$$
\begin{aligned}
x_{k+1} & =z_{k}=x_{k}-(1 / \kappa) g_{k} \\
\left(g_{k+1}, F_{k+1}, z_{k+1}\right) & =\text { ApproxGradient }\left(x_{k+1}, \mathcal{M}\right) .
\end{aligned}
$$

8: end if
9: update $B_{k+1}=\operatorname{L-BFGS}\left(B_{k}, x_{k+1}-x_{k}, g_{k+1}-g_{k}\right)$.
10: end for
output last proximal mapping $z_{K}$ (solution).

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The main characters:

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

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## Requirements on $\mathcal{M}$ and restarts

## Method $\mathcal{M}$

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

$$
h\left(w_{t}\right)-h^{\star} \leq C_{\mathcal{M}}\left(1-\tau_{\mathcal{M}}\right)^{t}\left(h\left(w_{0}\right)-h^{\star}\right)
$$

## 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

$$
w_{0}=\underset{w \in \mathbb{R}^{d}}{\arg \min }\left\{f_{0}(x)+\left\langle\nabla f_{0}(x), w-x\right\rangle+\frac{L+\kappa}{2}\|w-x\|^{2}+\psi(w)\right\}
$$

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

Three strategies to balance outer and inner computations
(a) use a pre-defined sequence $\left(\varepsilon_{k}\right)_{k \geq 0}$ and stop the optimization method $\mathcal{M}$ when the approximate proximal mapping is $\varepsilon_{k}$-accurate.
(b) define an adaptive stopping criterion that depends on quantities that are available at iteration $k$.
(c) use a pre-defined budget $T_{\mathcal{M}}$ of iterations of the method $\mathcal{M}$ for solving each sub-problem.

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## 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 $\mu$-strongly convex objectives $f$
(a) use a pre-defined sequence $\left(\varepsilon_{k}\right)_{k \geq 0}$ and stop the optimization method $\mathcal{M}$ when the approximate proximal mapping is $\varepsilon_{k}$-accurate.

$$
\varepsilon_{k}=\frac{1}{2} C(1-\rho)^{k+1} \quad \text { with } \quad C \geq f\left(x_{0}\right)-f^{*} \quad \text { and } \quad \rho=\frac{\mu}{4(\mu+\kappa)}
$$

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

$$
h\left(w_{t}\right)-h^{\star} \leq \frac{\kappa}{36}\left\|w_{t}-x\right\|^{2} .
$$

(c) use a pre-defined budget $T_{\mathcal{M}}$ of iterations of the method $\mathcal{M}$ for solving each sub-problem with

$$
T_{\mathcal{M}}=\frac{1}{\tau_{\mathcal{M}}} \log \left(19 C_{\mathcal{M}} \frac{L+\kappa}{\kappa}\right) \cdot \text { (be more aggressive in practice) }
$$

## 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 $\left(x_{k}\right)_{k \geq 0}$ and $\left(z_{k}\right)_{k \geq 0}$;

- $F\left(x_{k}\right) \rightarrow F^{\star}$, minimizes the smoothed objective $\Rightarrow$ no sparsity;
- $f\left(z_{k}\right) \rightarrow f^{\star}$, minimizes the true objective $\Rightarrow$ the iterates may be sparse if $\mathcal{M}$ handles composite optimization problems;

Global complexity
The number of iterations of $\mathcal{M}$ to guarantee $f\left(z_{k}\right)-f^{\star} \leq \varepsilon$ is at most

- $\tilde{O}\left(\frac{\mu+\kappa}{\tau_{\mathcal{M}} \mu} \log (1 / \varepsilon)\right)$ for $\mu$-strongly convex problems.
- $\tilde{O}\left(\frac{\kappa R^{2}}{\tau_{\mathcal{M}} \varepsilon}\right)$ for convex problems.


## Global Complexity and choice of $\kappa$

Example for gradient descent
With the right step-size, we have $\tau_{\mathcal{M}}=(\mu+\kappa) /(L+\kappa)$ and the complexity for $\mu>0$ becomes

$$
\tilde{O}\left(\frac{L+\kappa}{\mu} \log (1 / \varepsilon)\right) .
$$

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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\tilde{O}\left(\max \left(\frac{\mu+\kappa}{\mu} n, \frac{L+\kappa}{\mu}\right) \log (1 / \varepsilon)\right) .
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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 $\kappa$ ?
(i) assume that L-BFGS steps do as well as Nesterov.
(ii) choose $\kappa$ as in Catalyst.

## Experiments: formulations

- $\ell_{2}$-regularized Logistic Regression:

$$
\min _{x \in \mathbb{R}^{d}} \frac{1}{n} \sum_{i=1}^{n} \log \left(1+\exp \left(-b_{i} a_{i}^{T} x\right)\right)+\frac{\mu}{2}\|x\|^{2}
$$

- $\ell_{1}$-regularized Linear Regression (LASSO):

$$
\min _{x \in \mathbb{R}^{d}} \frac{1}{2 n} \sum_{i=1}^{n}\left(b_{i}-a_{i}^{T} x\right)^{2}+\lambda\|x\|_{1}
$$

- $\ell_{1}-\ell_{2}^{2}$-regularized Linear Regression (Elastic-Net):

$$
\min _{x \in \mathbb{R}^{d}} \frac{1}{2 n} \sum_{i=1}^{n}\left(b_{i}-a_{i}^{T} x\right)^{2}+\lambda\|x\|_{1}+\frac{\mu}{2}\|x\|^{2},
$$

## 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 | 250000 | 72309 | 781265 |
| $d$ | 54 | 500 | 20958 | 47152 |

- we simulate the ill-conditioned regime $\mu=1 /(100 n)$;
- $\lambda$ 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 $\geq$ SVRG, QuickeNing-SVRG2;
- QuickeNing-SVRG2 $\geq$ SVRG;
- QuickeNing-SVRG1 $\geq$ 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 $\kappa$



- $\kappa_{0}$ is the parameter (same as in Catalyst) used in all experiments;
- QuickeNing slows down when using $\kappa>\kappa_{0}$;
- 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? $\Rightarrow$ Links with recent Quasi-Newton methods applied to other envelopes [Stella et al., 2016].


## Outer-loop convergence analysis

Lemma: approximate descent property

$$
F\left(x_{k+1}\right) \leq f\left(z_{k}\right) \leq F\left(x_{k}\right)-\frac{1}{4 \kappa}\left\|\nabla F\left(x_{k}\right)\right\|_{2}^{2}+2 \varepsilon_{k} .
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Proposition: convergence with impractical $\varepsilon_{k}$ and $\mu>0$
If $\varepsilon_{k} \leq \frac{1}{16 \kappa}\left\|\nabla F\left(x_{k}\right)\right\|_{2}^{2}$, define $\rho=\frac{\mu}{4(\mu+\kappa)}$, then

$$
F\left(x_{k+1}\right)-F^{*} \leq f\left(z_{k}\right)-f^{*} \leq(1-\rho)^{k+1}\left(f\left(x_{0}\right)-f^{*}\right) .
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Lemma: convergence with adaptive $\varepsilon_{k}$ and $\mu>0$
If $\varepsilon_{k} \leq \frac{1}{36 k}\left\|g_{k}\right\|^{2}$, then $\varepsilon_{k} \leq \frac{1}{16}\left\|\nabla F\left(x_{k}\right)\right\|^{2}$.
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,

$$
\begin{equation*}
h\left(w_{0}\right)-h^{*} \leq \frac{L+\kappa}{2 \kappa^{2}}\|\nabla F(x)\|^{2} \tag{1}
\end{equation*}
$$

## Proof.

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

$$
\begin{aligned}
& h\left(w_{0}\right)-h^{*} \\
& \quad=f(x)-\left(f\left(w^{*}\right)+\frac{\kappa}{2}\left\|w^{*}-x\right\|^{2}\right) \\
& \leq f\left(w^{*}\right)+\left\langle\nabla f\left(w^{*}\right), x-w^{*}\right\rangle+\frac{L}{2}\left\|x-w^{*}\right\|^{2}-\left(f\left(w^{*}\right)+\frac{\kappa}{2}\left\|w^{*}-x\right\|^{2}\right) \\
& \quad=\frac{L+\kappa}{2}\left\|w^{*}-x\right\|^{2}=\frac{L+\kappa}{2 \kappa^{2}}\|\nabla F(x)\|^{2} .
\end{aligned}
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## Quasi-Newton and L-BFGS

## Presentation borrowed from Mark Schmidt, NIPS OPT 2010

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

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s_{k} \triangleq x_{k+1}-x_{k}, \quad y_{k} \triangleq \nabla f\left(x_{k+1}\right)-\nabla f\left(x_{k}\right)
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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:

$$
B_{k+1}=B_{k}-\frac{B_{k} s_{k} s_{k} B_{k}}{s_{k} B_{k} s_{k}}+\frac{y_{k} y_{k}^{\top}}{y_{k}^{\top} s_{k}}
$$

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- We can solve a linear system involving these updates when $B_{0}$ is diagonal in $O(d I)$ [Nocedal, 1980].


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