Stochastic optimization: Beyond stochastic gradients and convexity Part I

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Context

Machine learning for large-scale data

- Large-scale supervised machine learning: large d, large n
 - -d: dimension of each observation (input) or number of parameters
 - -n: number of observations
- **Examples**: computer vision, advertising, bioinformatics, etc.

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Visual object recognition



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- Ideal running-time complexity: O(dn)

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- Large-scale supervised machine learning: large d, large n
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- **Examples**: computer vision, advertising, bioinformatics, etc.
- Ideal running-time complexity: O(dn)
- Going back to simple methods
 - Stochastic gradient methods (Robbins and Monro, 1951)
- Goal: Present recent progress

Outline

1. Introduction/motivation: Supervised machine learning

- Optimization of finite sums
- Existing optimization methods for finite sums

2. Convex finite-sum problems

- Linearly-convergent stochastic gradient method
- SAG, SAGA, SVRG, SDCA, MISO, etc.
- From lazy gradient evaluations to variance reduction

3. Non-convex problems

- 4. Parallel and distributed settings
- 5. Perspectives

References

• Textbooks and tutorials

- Nesterov (2004): Introductory lectures on convex optimization
- Bubeck (2015): Convex Optimization: Algorithms and Complexity
- Bertsekas (2016): Nonlinear programming
- Bottou et al. (2016): *Optimization methods for large-scale machine learning*

• Research papers

- See end of slides
- Slides available at www.ens.fr/~fbach/

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– Linear predictions: $h(x,\theta) = \theta^{\top} \Phi(x)$ with features $\Phi(x) \in \mathbb{R}^d$

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 - Linear predictions: $h(x,\theta) = \theta^{\top} \Phi(x)$ with features $\Phi(x) \in \mathbb{R}^d$
 - Neural networks: $h(x,\theta) = \theta_m^\top \sigma(\theta_{m-1}^\top \sigma(\cdots \theta_2^\top \sigma(\theta_1^\top x)))$



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- Prediction function $h(x, \theta) \in \mathbb{R}$ parameterized by $\theta \in \mathbb{R}^d$
- (regularized) empirical risk minimization: find $\hat{\theta}$ solution of

$$\min_{\theta \in \mathbb{R}^d} \quad \frac{1}{n} \sum_{i=1}^n \quad \ell(y_i, h(x_i, \theta)) \quad + \quad \lambda \Omega(\theta) \qquad = \frac{1}{n} \sum_{i=1}^n f_i(\theta)$$

data fitting term + regularizer

Usual losses

• **Regression**: $y \in \mathbb{R}$

– quadratic loss $\ell(y, h(x, \theta)) = \frac{1}{2}(y - h(x, \theta))^2$

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 - Logistic loss $\ell(y, h(x, \theta)) = \log(1 + \exp(-yh(x, \theta)))$
- Structured prediction
 - Complex outputs y (k classes/labels, graphs, trees, or $\{0,1\}^k$, etc.)
 - Prediction function $h(x, \theta) \in \mathbb{R}^k$
 - Conditional random fields (Lafferty et al., 2001)
 - Max-margin (Taskar et al., 2003; Tsochantaridis et al., 2005)

Supervised machine learning

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data fitting term + regularizer

testing cost

- Optimization: optimization of regularized risk training cost
- Statistics: guarantees on $\mathbb{E}_{p(x,y)}\ell(y,h(x,\theta))$

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$$\forall \theta \in \mathbb{R}^d, | eigenvalues[g''(\theta)] | \leq L$$

• Machine learning

- with
$$g(\theta) = \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, h(x_i, \theta))$$

- Smooth prediction function $\theta \mapsto h(x_i, \theta) + \text{smooth loss}$

• A twice differentiable function $g:\mathbb{R}^d\to\mathbb{R}$ is convex if and only if

$$\forall \theta \in \mathbb{R}^d, \text{ eigenvalues}[g''(\theta)] \ge 0$$



 $\forall \theta \in \mathbb{R}^d$, eigenvalues $[g''(\theta)] \ge \mu$

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 $\forall \theta \in \mathbb{R}^d, \text{ eigenvalues}[g''(\theta)] \geqslant \mu$

– Condition number $\kappa = L/\mu \geqslant 1$



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- Convexity in machine learning
 - With $g(\theta) = \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, h(x_i, \theta))$
 - Convex loss and linear predictions $h(x, \theta) = \theta^{\top} \Phi(x)$

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• Relevance of convex optimization

- Easier design and analysis of algorithms
- Global minimum vs. local minimum vs. stationary points
- Gradient-based algorithms only need convexity for their analysis

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• Strong convexity in machine learning

- With
$$g(\theta) = \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, h(x_i, \theta))$$

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- Invertible covariance matrix $\frac{1}{n} \sum_{i=1}^{n} \Phi(x_i) \Phi(x_i)^{\top} \Rightarrow n \ge d$
- Even when $\mu > 0$, μ may be arbitrarily small!

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- Even when $\mu > 0$, μ may be arbitrarily small!
- Adding regularization by $\frac{\mu}{2} \|\theta\|^2$
 - creates additional bias unless μ is small, but reduces variance
 - Typically $L/\sqrt{n} \geqslant \mu \geqslant L/n$

- Assumption: g convex and L-smooth on \mathbb{R}^d
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 $g(\theta_t) - g(\theta_*) \leqslant O(1/t)$ $g(\theta_t) - g(\theta_*) \leq O((1 - \mu/L)^t) = O(e^{-t(\mu/L)})$ if μ -strongly convex (large $\kappa = L/\mu$) (small $\kappa = L/\mu$)

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- Gradient descent: $\theta_t = \theta_{t-1} \gamma_t g'(\theta_{t-1})$
 - O(1/t) convergence rate for convex functions - $O(e^{-t/\kappa})$ linear if strongly-convex
- Newton method: $\theta_t = \theta_{t-1} g''(\theta_{t-1})^{-1}g'(\theta_{t-1})$
 - $-O(e^{-\rho 2^t})$ quadratic rate

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- Key insights for machine learning (Bottou and Bousquet, 2008)
 - 1. No need to optimize below statistical error
 - 2. Cost functions are averages
 - 3. Testing error is more important than training error

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Stochastic gradient descent (SGD) for finite sums

$$\min_{\theta \in \mathbb{R}^d} g(\theta) = \frac{1}{n} \sum_{i=1}^n f_i(\theta)$$

- Iteration: $\theta_t = \theta_{t-1} \gamma_t f'_{i(t)}(\theta_{t-1})$
 - Sampling with replacement: i(t) random element of $\{1, \ldots, n\}$
 - Polyak-Ruppert averaging: $\bar{\theta}_t = \frac{1}{t+1} \sum_{u=0}^t \theta_u$
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 - Polyak-Ruppert averaging: $\bar{\theta}_t = \frac{1}{t+1} \sum_{u=0}^t \theta_u$
- **Convergence rate** if each f_i is convex *L*-smooth and $g \mu$ -strongly-convex:

$$\mathbb{E}g(\bar{\theta}_t) - g(\theta_*) \leqslant \begin{cases} O(1/\sqrt{t}) & \text{if } \gamma_t = 1/(L\sqrt{t}) \\ O(L/(\mu t)) = O(\kappa/t) & \text{if } \gamma_t = 1/(\mu t) \end{cases}$$

- No adaptivity to strong-convexity in general
- Adaptivity with self-concordance assumption (Bach, 2014)
- Running-time complexity: $O(d \cdot \kappa/\varepsilon)$

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• Minimizing
$$g(\theta) = \frac{1}{n} \sum_{i=1}^{n} f_i(\theta)$$
 with $f_i(\theta) = \ell(y_i, h(x_i, \theta)) + \lambda \Omega(\theta)$

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- Batch gradient descent: $\theta_t = \theta_{t-1} \gamma_t g'(\theta_{t-1}) = \theta_{t-1} \frac{\gamma_t}{n} \sum_{i=1}^n f'_i(\theta_{t-1})$
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 - Iteration complexity is linear in n

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- Stochastic gradient descent: $\theta_t = \theta_{t-1} \gamma_t f'_{i(t)}(\theta_{t-1})$
 - Sampling with replacement: i(t) random element of $\{1,\ldots,n\}$
 - Convergence rate in $O(\kappa/t)$
 - Iteration complexity is independent of n

- Minimizing $g(\theta) = \frac{1}{n} \sum_{i=1}^{n} f_i(\theta)$ with $f_i(\theta) = \ell(y_i, h(x_i, \theta)) + \lambda \Omega(\theta)$
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• Goal = best of both worlds: Linear rate with O(d) iteration cost Simple choice of step size



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• Generic acceleration (Nesterov, 1983, 2004)

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- Still O(nd) iteration cost: complexity = $O(nd \cdot \sqrt{\kappa} \log \frac{1}{\epsilon})$

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 - Similar linear rate but limited choice for the f_i 's
 - Extensions without duality: see Shalev-Shwartz (2016)

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 - Similar linear rate but limited choice for the f_i 's
 - Extensions without duality: see Shalev-Shwartz (2016)
- Stochastic version of accelerated batch gradient methods
 - Tseng (1998); Ghadimi and Lan (2010); Xiao (2010)
 - Can improve constants, but still have sublinear O(1/t) rate

- Stochastic average gradient (SAG) iteration
 - Keep in memory the gradients of all functions f_i , $i = 1, \ldots, n$
 - Random selection $i(t) \in \{1, \ldots, n\}$ with replacement

- Iteration:
$$\theta_t = \theta_{t-1} - \frac{\gamma_t}{n} \sum_{i=1}^n y_i^t$$
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functions
$$g = \frac{1}{n} \sum_{i=1}^{n} f_i$$
 f_1 f_2 f_3 f_4 \cdots f_{n-1} f_n
gradients $\in \mathbb{R}^d$ $\frac{1}{n} \sum_{i=1}^{n} y_i^t$ y_1^t y_2^t y_3^t y_4^t \cdots y_{n-1}^t y_n^t

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functions $g = \frac{1}{n} \sum_{i=1}^{n} f_i$ f_1 f_2 f_3 f_4 \cdots f_{n-1} f_n gradients $\in \mathbb{R}^d$ $\frac{1}{n} \sum_{i=1}^{n} y_i^t$ y_1^t y_2^t y_3^t y_4^t \cdots y_{n-1}^t y_n^t



- Stochastic average gradient (SAG) iteration
 - Keep in memory the gradients of all functions f_i , $i = 1, \ldots, n$
 - Random selection $i(t) \in \{1, \ldots, n\}$ with replacement

- Iteration:
$$\theta_t = \theta_{t-1} - \frac{\gamma_t}{n} \sum_{i=1}^n y_i^t$$
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• Stochastic version of incremental average gradient (Blatt et al., 2008)

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- Stochastic version of incremental average gradient (Blatt et al., 2008)
- Extra memory requirement: n gradients in \mathbb{R}^d in general
- Linear supervised machine learning: only n real numbers

- If
$$f_i(\theta) = \ell(y_i, \Phi(x_i)^\top \theta)$$
, then $f'_i(\theta) = \ell'(y_i, \Phi(x_i)^\top \theta) \Phi(x_i)$

Stochastic average gradient - Convergence analysis

• Assumptions

- Each f_i is L-smooth, $i = 1, \ldots, n$
- $-g = \frac{1}{n} \sum_{i=1}^{n} f_i$ is μ -strongly convex
- constant step size $\gamma_t = 1/(16L)$ no need to know μ

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- constant step size $\gamma_t = 1/(16L)$ no need to know μ
- Strongly convex case (Le Roux et al., 2012; Schmidt et al., 2016)

$$\mathbb{E}\big[g(\theta_t) - g(\theta_*)\big] \leqslant \operatorname{cst} \times \Big(1 - \min\left\{\frac{1}{8n}, \frac{\mu}{16L}\right\}\Big)^t$$

- Linear (exponential) convergence rate with O(d) iteration cost
- After one pass, reduction of cost by $\exp\left(-\min\left\{\frac{1}{8},\frac{n\mu}{16L}\right\}\right)$
- NB: in machine learning, may often restrict to $\mu \ge L/n$ \Rightarrow constant error reduction after each effective pass

Running-time comparisons (strongly-convex)

- Assumptions: $g(\theta) = \frac{1}{n} \sum_{i=1}^{n} f_i(\theta)$
 - Each f_i convex L-smooth and $g \mu$ -strongly convex

| Stochastic gradient descent | $d \times$ | $\frac{L}{\mu}$ | $\times \frac{1}{\varepsilon}$ |
|------------------------------|------------|-------------------------|-------------------------------------|
| Gradient descent | $d \times$ | $n\frac{L}{\mu}$ | $\times \log \frac{1}{\varepsilon}$ |
| Accelerated gradient descent | $d \times$ | $n\sqrt{\frac{L}{\mu}}$ | $\times \log \frac{1}{\varepsilon}$ |
| SAG | $d \times$ | $(n + \frac{L}{\mu})$ | $\times \log \frac{1}{\varepsilon}$ |

- NB-1: for (accelerated) gradient descent, L = smoothness constant of g
- NB-2: with non-uniform sampling, $L = average smoothness constants of all <math>f_i$'s

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- **Beating two lower bounds** (Nemirovski and Yudin, 1983; Nesterov, 2004): with additional assumptions
- (1) stochastic gradient: exponential rate for finite sums
- (2) full gradient: better exponential rate using the sum structure

Running-time comparisons (non-strongly-convex)

- Assumptions: $g(\theta) = \frac{1}{n} \sum_{i=1}^{n} f_i(\theta)$
 - Each f_i convex L-smooth
 - Ill conditioned problems: g may not be strongly-convex ($\mu = 0$)

| Stochastic gradient descent | $d \times$ | $1/\varepsilon^2$ |
|------------------------------|------------|------------------------|
| Gradient descent | $d \times$ | n/arepsilon |
| Accelerated gradient descent | $d \times$ | $n/\sqrt{\varepsilon}$ |
| SAG | $d \times$ | \sqrt{n}/ε |

- Adaptivity to potentially hidden strong convexity
- No need to know the local/global strong-convexity constant

Stochastic average gradient Implementation details and extensions

- Sparsity in the features
 - Just-in-time updates \Rightarrow replace O(d) by number of non zeros
 - See also Leblond, Pedregosa, and Lacoste-Julien (2016)
- Mini-batches
 - Reduces the memory requirement + block access to data
- Line-search
 - Avoids knowing L in advance
- Non-uniform sampling
 - Favors functions with large variations
- See http://www.cs.ubc.ca/~schmidtm/Software/SAG.html

Experimental results (logistic regression)

quantum dataset $(n = 50 \ 000, \ d = 78)$

rcv1 dataset $(n = 697 \ 641, \ d = 47 \ 236)$



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Before non-uniform sampling

protein dataset $(n = 145 \ 751, d = 74)$

sido dataset
$$(n = 12 \ 678, \ d = 4 \ 932)$$



After non-uniform sampling





Linearly convergent stochastic gradient algorithms

• Many related algorithms

- SAG (Le Roux et al., 2012)
- SDCA (Shalev-Shwartz and Zhang, 2013)
- SVRG (Johnson and Zhang, 2013; Zhang et al., 2013)
- MISO (Mairal, 2015)
- Finito (Defazio et al., 2014b)
- SAGA (Defazio, Bach, and Lacoste-Julien, 2014a)

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• Similar rates of convergence and iterations

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- Finito (Defazio et al., 2014b)
- SAGA (Defazio, Bach, and Lacoste-Julien, 2014a)

- Similar rates of convergence and iterations
- Different interpretations and proofs / proof lengths
 - Lazy gradient evaluations
 - Variance reduction

Variance reduction

• **Principle**: reducing variance of sample of X by using a sample from another random variable Y with known expectation

$$Z_{\alpha} = \alpha(X - Y) + \mathbb{E}Y$$

 $-\mathbb{E}Z_{\alpha} = \alpha \mathbb{E}X + (1 - \alpha)\mathbb{E}Y$ $-\operatorname{var}(Z_{\alpha}) = \alpha^{2} [\operatorname{var}(X) + \operatorname{var}(Y) - 2\operatorname{cov}(X, Y)]$ $-\alpha = 1: \text{ no bias, } \alpha < 1: \text{ potential bias (but reduced variance)}$

– Useful if Y positively correlated with X

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• Application to gradient estimation (Johnson and Zhang, 2013; Zhang, Mahdavi, and Jin, 2013)

- SVRG:
$$X = f'_{i(t)}(\theta_{t-1})$$
, $Y = f'_{i(t)}(\tilde{\theta})$, $\alpha = 1$, with $\tilde{\theta}$ stored

 $-\mathbb{E}Y = \frac{1}{n} \sum_{i=1}^{n} f'_{i}(\tilde{\theta}) \text{ full gradient at } \tilde{\theta}, X - Y = f'_{i(t)}(\theta_{t-1}) - f'_{i(t)}(\tilde{\theta})$

Stochastic variance reduced gradient (SVRG) (Johnson and Zhang, 2013; Zhang et al., 2013)

• Initialize $\tilde{\theta} \in \mathbb{R}^d$

• For
$$i_{epoch} = 1$$
 to $\#$ of epochs

- Compute all gradients $f_i'(\tilde{\theta})$ store $g'(\tilde{\theta}) = \frac{1}{n} \sum_{i=1}^n f_i'(\tilde{\theta})$
- Initialize $\theta_0 = \tilde{\theta}$
- For t = 1 to length of epochs

$$\begin{split} \theta_t &= \theta_{t-1} - \gamma \Big[g'(\tilde{\theta}) + \big(f'_{i(t)}(\theta_{t-1}) - f'_{i(t)}(\tilde{\theta}) \big) \Big] \\ \text{Update } \tilde{\theta} &= \theta_t \end{split}$$

- Output: $\tilde{\theta}$
- No need to store gradients two gradient evaluations per inner step
- Two parameters: lengths of epoch + step-size
- Same linear convergence rate as SAG, simpler proof

Interpretation of SAG as variance reduction

• SAG update:
$$\theta_t = \theta_{t-1} - \frac{\gamma}{n} \sum_{i=1}^n y_i^t$$
 with $y_i^t = \begin{cases} f'_i(\theta_{t-1}) & \text{if } i = i(t) \\ y_i^{t-1} & \text{otherwise} \end{cases}$

- Interpretation as lazy gradient evaluations
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 - Biased update (expectation w.r.t. to i(t) not equal to full gradient)

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- SAG update: $\theta_t = \theta_{t-1} \gamma \left[\frac{1}{n} \sum_{i=1}^n y_i^{t-1} + \frac{1}{n} (f'_{i(t)}(\theta_{t-1}) y_{i(t)}^{t-1}) \right]$
 - Biased update (expectation w.r.t. to i(t) not equal to full gradient)
- **SVRG update**: $\theta_t = \theta_{t-1} \gamma \left[\frac{1}{n} \sum_{i=1}^n f'_i(\tilde{\theta}) + \left(f'_{i(t)}(\theta_{t-1}) f'_{i(t)}(\tilde{\theta}) \right) \right]$

- Unbiased update

Interpretation of SAG as variance reduction

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- Unbiased update

- SAGA update: $\theta_t = \theta_{t-1} \gamma_t \left[\frac{1}{n} \sum_{i=1}^n y_i^{t-1} + \left(f'_{i(t)}(\theta_{t-1}) y_{i(t)}^{t-1} \right) \right]$
 - Defazio, Bach, and Lacoste-Julien (2014a)
 - Unbiased update without epochs

SVRG vs. SAGA

• SAGA update:
$$\theta_t = \theta_{t-1} - \gamma_t \left[\frac{1}{n} \sum_{i=1}^n y_i^{t-1} + \left(f'_{i(t)}(\theta_{t-1}) - y_{i(t)}^{t-1} \right) \right]$$

• **SVRG update**:
$$\theta_t = \theta_{t-1} - \gamma \left[\frac{1}{n} \sum_{i=1}^n f'_i(\tilde{\theta}) + \left(f'_{i(t)}(\theta_{t-1}) - f'_{i(t)}(\tilde{\theta}) \right) \right]$$

| | SAGA | SVRG |
|--------------------------------|-----------|---------------------------|
| Storage of gradients | yes | no |
| Epoch-based | no | yes |
| Parameters | step-size | step-size & epoch lengths |
| Gradient evaluations per step | 1 | at least 2 |
| Adaptivity to strong-convexity | yes | no |
| Robustness to ill-conditioning | yes | no |

- See Babanezhad et al. (2015)

$$\frac{1}{n}\sum_{i=1}^{n}f_{i}(\theta)+h(\theta)$$

- f_i smooth and convex
- -h convex, potentially non-smooth

$$\frac{1}{n} \sum_{i=1}^{n} f_i(\theta) + h(\theta)$$

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- Directly extends to variance-reduced gradient techniques
 - Same rates of convergence

Acceleration

• Similar guarantees for finite sums: SAG, SDCA, SVRG (Xiao and Zhang, 2014), SAGA, MISO (Mairal, 2015)

| Gradient descent | $dn\frac{L}{\mu}$ | $\times \log \frac{1}{\varepsilon}$ |
|------------------------------|--------------------------|-------------------------------------|
| Accelerated gradient descent | $dn\sqrt{\frac{L}{\mu}}$ | $\times \log \frac{1}{\varepsilon}$ |
| SAG(A), SVRG, SDCA, MISO | $d(n + \frac{L}{\mu})$ | $\times \log \frac{1}{\varepsilon}$ |

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| SAG(A), SVRG, SDCA, MISO | $d(n + \frac{L}{\mu})$ | $\times \log \frac{1}{\varepsilon}$ |
| Accelerated versions | $d(n + \sqrt{n\frac{L}{\mu}})$ | $\times \log \frac{1}{\varepsilon}$ |

- Acceleration for special algorithms (e.g., Shalev-Shwartz and Zhang, 2014; Nitanda, 2014; Lan, 2015)
- Catalyst (Lin, Mairal, and Harchaoui, 2015)
 - Widely applicable generic acceleration scheme

From training to testing errors

- rcv1 dataset ($n = 697\ 641$, $d = 47\ 236$)
 - NB: IAG, SG-C, ASG with optimal step-sizes in hindsight



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SGD minimizes the testing cost!

- Goal: minimize $f(\theta) = \mathbb{E}_{p(x,y)} \ell(y, \theta^{\top} \Phi(x))$
 - Given n independent samples (x_i, y_i) , i = 1, ..., n from p(x, y)
 - Given a single pass of stochastic gradient descent
 - Bounds on the excess testing cost $\mathbb{E}f(\bar{\theta}_n) \inf_{\theta \in \mathbb{R}^d} f(\theta)$

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 - Optimal for non-smooth losses (Nemirovski and Yudin, 1983)
 - Attained by averaged SGD with decaying step-sizes
- Constant-step-size SGD
 - Linear convergence up to the noise level for strongly-convex problems (Solodov, 1998; Nedic and Bertsekas, 2000)
 - Full convergence and robustness to ill-conditioning?

Robust averaged stochastic gradient (Bach and Moulines, 2013)

- Constant-step-size SGD is convergent for least-squares
 - Convergence rate in ${\cal O}(1/n)$ without any dependence on μ
 - Simple choice of step-size (equal to 1/L)



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 \bullet Convergence in O(1/n) for smooth losses with O(d) online Newton step

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 - Provable and precise rates
 - Improves on two known lower-bounds (by using structure)
 - Several extensions / interpretations / accelerations

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- Extension to saddle-point problems (Balamurugan and Bach, 2016)
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- What's next: non-convexity, parallelization, extensions/perspectives

Postdoc opportunities in downtown Paris



- Machine learning group at INRIA Ecole Normale Supérieure
 - Two postdoc positions (2 years)
 - One junior researcher position (4 years)

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