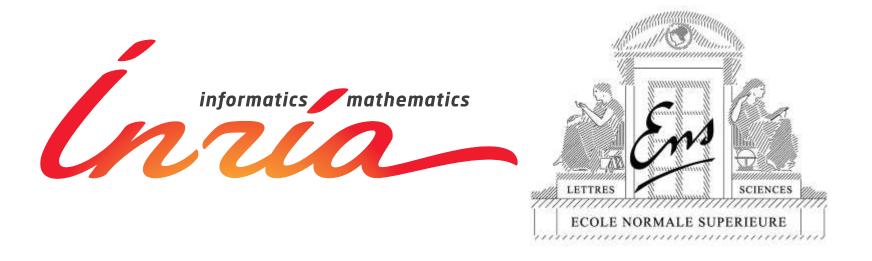
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Stochastic gradient methods for machine learning

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Joint work with Eric Moulines, Nicolas Le Roux and Mark Schmidt - April 2013

Context Machine learning for "big data"

- Large-scale machine learning: large p, large n, large k
 - -p: dimension of each observation (input)
 - -k: number of tasks (dimension of outputs)
 - -n: number of observations
- Examples: computer vision, bioinformatics, signal processing
- Ideal running-time complexity: O(pn + kn)

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- Ideal running-time complexity: O(pn + kn)
- Going back to simple methods
 - Stochastic gradient methods (Robbins and Monro, 1951)
 - Mixing statistics and optimization

Outline

Introduction

- Supervised machine learning and convex optimization
- Stochastic approximation algorithms (Bach and Moulines, 2011; Bach, 2013)
 - Stochastic gradient and averaging
 - Strongly convex vs. non-strongly convex
 - Adaptivity
- Going beyond stochastic gradient (Le Roux, Schmidt, and Bach, 2012, 2013)
 - More than a single pass through the data
 - Linear (exponential) convergence rate for strongly convex functions

Supervised machine learning

- Data: n observations $(x_i, y_i) \in \mathcal{X} \times \mathcal{Y}$, $i = 1, \ldots, n$, i.i.d.
- Prediction as a linear function $\theta^{\top}\Phi(x)$ of features $\Phi(x) \in \mathcal{F} = \mathbb{R}^p$
- (regularized) empirical risk minimization: find $\hat{\theta}$ solution of

$$\min_{\theta \in \mathcal{F}} \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, \theta^{\top} \Phi(x_i)) + \mu \Omega(\theta)$$

convex data fitting term + regularizer

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convex data fitting term + regularizer

- Empirical risk: $\hat{f}(\theta) = \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, \theta^{\top} \Phi(x_i))$ training cost
- Expected risk: $f(\theta) = \mathbb{E}_{(x,y)} \ell(y, \theta^{\top} \Phi(x))$ testing cost
- Two fundamental questions: (1) computing $\hat{\theta}$ and (2) analyzing $\hat{\theta}$

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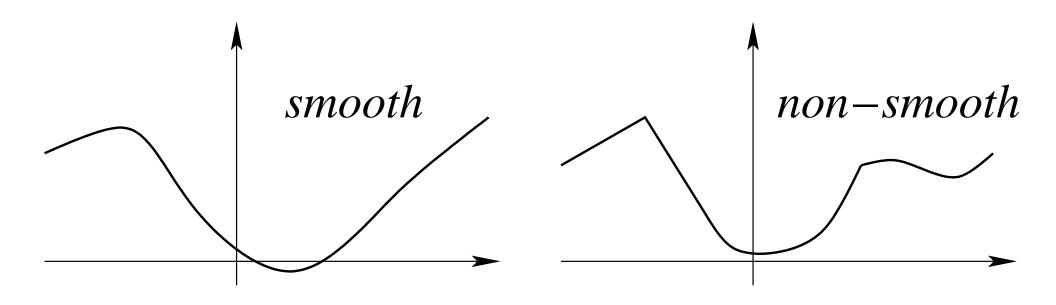
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- Two fundamental questions: (1) computing $\hat{\theta}$ and (2) analyzing $\hat{\theta}$
 - May be tackled simultaneously

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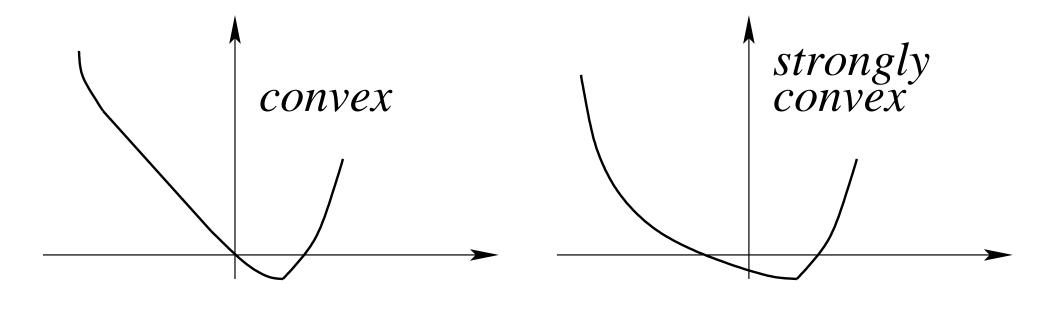
Machine learning

- with $g(\theta) = \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, \theta^{\top} \Phi(x_i))$
- Hessian \approx covariance matrix $\frac{1}{n} \sum_{i=1}^{n} \Phi(x_i) \Phi(x_i)^{\top}$
- Bounded data

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 - with $g(\theta) = \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, \theta^{\top} \Phi(x_i))$
 - Hessian \approx covariance matrix $\frac{1}{n} \sum_{i=1}^{n} \Phi(x_i) \Phi(x_i)^{\top}$
 - Data with invertible covariance matrix (low correlation/dimension)
 - ... or with added regularization by $\frac{\mu}{2} \|\theta\|^2$

Iterative methods for minimizing smooth functions

- **Assumption**: g convex and smooth on $\mathcal{F} = \mathbb{R}^p$
- Gradient descent: $\theta_t = \theta_{t-1} \gamma_t g'(\theta_{t-1})$
 - O(1/t) convergence rate for convex functions
 - $O(e^{-\rho t})$ convergence rate for strongly convex functions
- Newton method: $\theta_t = \theta_{t-1} g''(\theta_{t-1})^{-1}g'(\theta_{t-1})$
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- Key insights from Bottou and Bousquet (2008)
 - 1. In machine learning, no need to optimize below statistical error
 - 2. In machine learning, cost functions are averages

⇒ Stochastic approximation

Stochastic approximation

- ullet Goal: Minimizing a function f defined on $\mathcal{F}=\mathbb{R}^p$
 - given only unbiased estimates $f_n'(\theta_n)$ of its gradients $f'(\theta_n)$ at certain points $\theta_n \in \mathcal{F}$

Stochastic approximation

- Observation of $f'_n(\theta_n) = f'(\theta_n) + \varepsilon_n$, with $\varepsilon_n = \text{i.i.d.}$ noise
- Non-convex problems

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Machine learning - statistics

– loss for a single pair of observations: $|f_n(\theta) = \ell(y_n, \theta^{\top} \Phi(x_n))|$

$$f_n(\theta) = \ell(y_n, \theta^{\top} \Phi(x_n))$$

- $-f(\theta) = \mathbb{E} f_n(\theta) = \mathbb{E} \ell(y_n, \theta^{\top} \Phi(x_n)) =$ generalization error
- Expected gradient: $f'(\theta) = \mathbb{E} f'_n(\theta) = \mathbb{E} \left\{ \ell'(y_n, \theta^\top \Phi(x_n)) \Phi(x_n) \right\}$

Convex smooth stochastic approximation

• **Key assumption**: smoothness and/or strongly convexity

Convex smooth stochastic approximation

- **Key assumption**: smoothness and/or strongly convexity
- **Key algorithm:** stochastic gradient descent (a.k.a. Robbins-Monro)

$$\theta_n = \theta_{n-1} - \gamma_n f'_n(\theta_{n-1})$$

- Polyak-Ruppert averaging: $\bar{\theta}_n = \frac{1}{n} \sum_{k=0}^{n-1} \theta_k$
- Which learning rate sequence γ_n ? Classical setting: $| \gamma_n = Cn^{-\alpha} |$

$$\gamma_n = C n^{-\alpha}$$

- Known global minimax rates of convergence (Nemirovski and Yudin, 1983; Agarwal et al., 2010)
 - Strongly convex: $O(n^{-1})$ Attained by averaged stochastic gradient descent with $\gamma_n \propto (\mu n)^{-1}$
 - Non-strongly convex: $O(n^{-1/2})$ Attained by averaged stochastic gradient descent with $\gamma_n \propto n^{-1/2}$

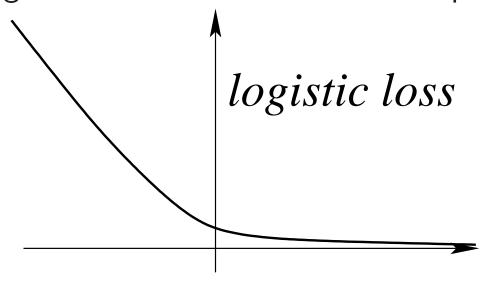
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- Bottou and Le Cun (2005); Bottou and Bousquet (2008); Hazan et al. (2007); Shalev-Shwartz and Srebro (2008); Shalev-Shwartz et al. (2007, 2009); Xiao (2010); Duchi and Singer (2009); Nesterov and Vial (2008); Nemirovski et al. (2009)

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- Asymptotic analysis of averaging (Polyak and Juditsky, 1992;
 Ruppert, 1988)
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 - A single algorithm with global convergence rate?

Adaptive algorithm for logistic regression

- Logistic regression: $(x_n, y_n) \in \mathbb{R}^p \times \{-1, 1\}$
 - Single data point: $f_n(\theta) = \log(1 + \exp(-y_n \theta^{\top} x_n))$
 - Generalization error: $f(\theta) = \mathbb{E}f_n(\theta)$
- Cannot be strongly convex ⇒ local strong convexity
 - unless restricted to $|\theta^{\top}x_n| \leqslant M$
 - μ = lowest eigenvalue of the Hessian at the optimum $f''(\theta_*)$



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 - unless restricted to $|\theta^{\top}x_n| \leqslant M$
 - μ = lowest eigenvalue of the Hessian at the optimum $f''(\theta_*)$
- n steps of averaged SGD with constant step-size $1/(2R^2\sqrt{n})$
 - with R = radius of data (Bach, 2013):

$$\mathbb{E}f(\bar{\theta}_n) - f(\theta_*) \leqslant \min\left\{\frac{1}{\sqrt{n}}, \frac{R^2}{n\mu}\right\} \left(15 + 5R\|\theta_0 - \theta_*\|\right)^4$$

- Proof based on generalized self-concordance (Bach, 2010)

Conclusions / Extensions Stochastic approximation for machine learning

- Mixing convex optimization and statistics
 - Non-asymptotic analysis through moment computations
 - Averaging with longer steps is (more) robust and adaptive

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Mixing convex optimization and statistics

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• Future/current work - open problems

- High-probability through all moments $\mathbb{E}\|\theta_n-\theta_*\|^{2d}$
- Non-random errors (Schmidt, Le Roux, and Bach, 2011)
- Line search for stochastic gradient
- Non-parametric stochastic approximation
- Going beyond a single pass through the data

Going beyond a single pass over the data

Stochastic approximation

- Assumes infinite data stream
- Observations are used only once
- Directly minimizes testing cost $\mathbb{E}_{(x,y)} \ell(y, \theta^{\top} \Phi(x))$

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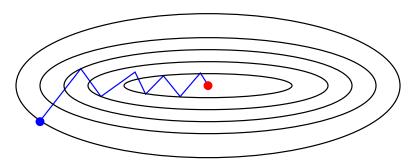
Machine learning practice

- Finite data set $(x_1, y_1, \dots, x_n, y_n)$
- Multiple passes
- Minimizes training cost $\frac{1}{n} \sum_{i=1}^{n} \ell(y_i, \theta^{\top} \Phi(x_i))$
- Need to regularize (e.g., by the ℓ_2 -norm) to avoid overfitting

- Minimizing $g(\theta) = \frac{1}{n} \sum_{i=1}^{n} f_i(\theta)$ with $f_i(\theta) = \ell(y_i, \theta^{\top} \Phi(x_i)) + \mu \Omega(\theta)$
- Batch gradient descent: $\theta_t = \theta_{t-1} \gamma_t g'(\theta_{t-1}) = \theta_{t-1} \frac{\gamma_t}{n} \sum_{i=1}^{n} f_i'(\theta_{t-1})$
 - Linear (e.g., exponential) convergence rate
 - Iteration complexity is linear in n

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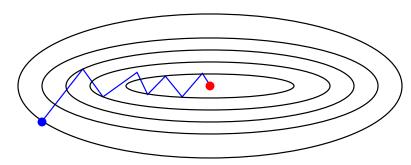
损失函数

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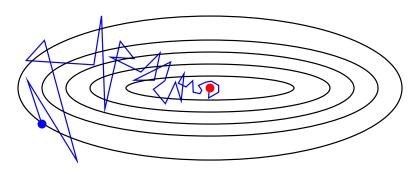
- Stochastic gradient descent: $\theta_t = \theta_{t-1} \gamma_t f'_{i(t)}(\theta_{t-1})$
 - Sampling with replacement: i(t) random element of $\{1,\ldots,n\}$
 - Convergence rate in O(1/t)
 - Iteration complexity is independent of n

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

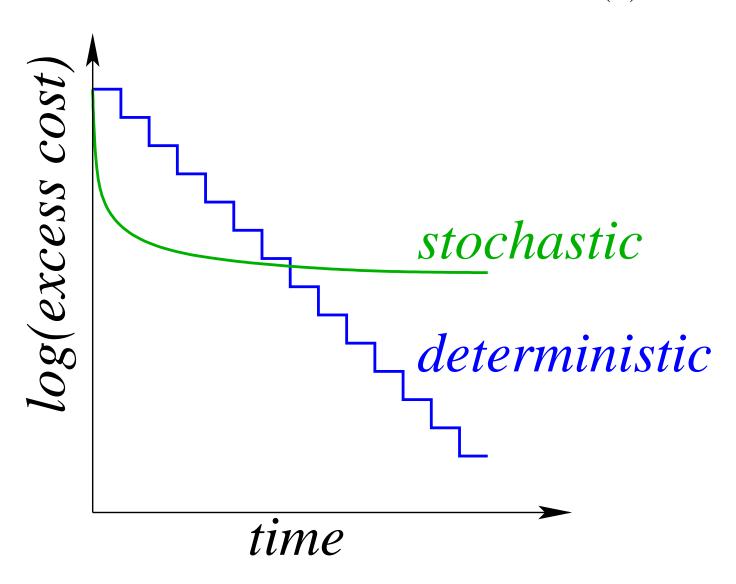
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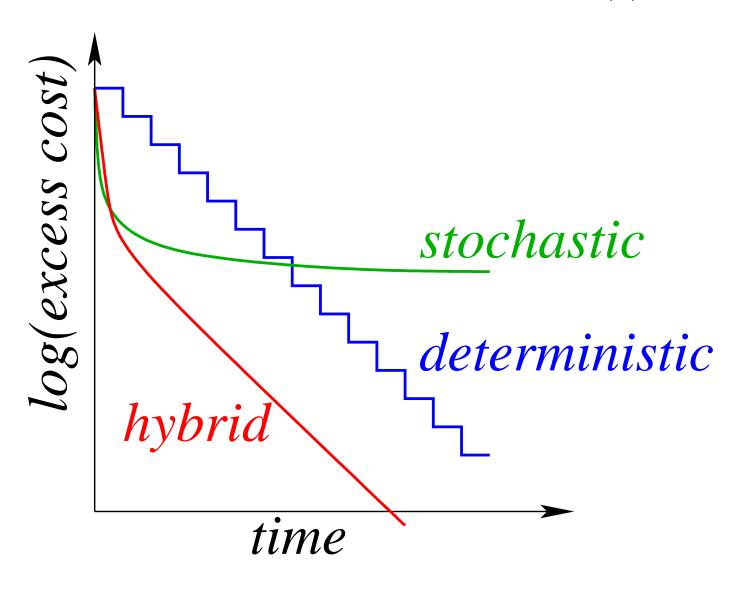
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• Goal = best of both worlds: linear rate with O(1) iteration cost



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Accelerating gradient methods - Related work

Nesterov acceleration

- Nesterov (1983, 2004)
- Better linear rate but still O(n) iteration cost
- Hybrid methods, incremental average gradient, increasing batch size
 - Bertsekas (1997); Blatt et al. (2008); Friedlander and Schmidt (2011)
 - Linear rate, but iterations make full passes through the data.

Accelerating gradient methods - Related work

- Momentum, gradient/iterate averaging, stochastic version of accelerated batch gradient methods
 - Polyak and Juditsky (1992); Tseng (1998); Sunehag et al. (2009);
 Ghadimi and Lan (2010); Xiao (2010)
 - Can improve constants, but still have sublinear O(1/t) rate
- Constant step-size stochastic gradient (SG), accelerated SG
 - Kesten (1958); Delyon and Juditsky (1993); Solodov (1998); Nedic and Bertsekas (2000)
 - Linear convergence, but only up to a fixed tolerance.
- Stochastic methods in the dual
 - Shalev-Shwartz and Zhang (2012)
 - Linear rate but limited choice for the f_i 's

Stochastic average gradient (Le Roux, Schmidt, and Bach, 2012)

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

$$- \text{ Iteration: } \theta_t = \theta_{t-1} - \frac{\underline{\gamma_t}}{\underline{n}} \sum_{\underline{i=1}}^n \underline{y_{\underline{i}}^t \text{ with } y_i^t} = \underbrace{\begin{cases} \underline{f_i'(\theta_{t-1})} & \text{if } i = i(t) \\ \underline{y_{\underline{i}}^{t-1}} & \text{otherwise} \end{cases}}_{}$$

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- Stochastic version of incremental average gradient (Blatt et al., 2008)
- Extra memory requirement
 - Supervised machine learning
 - If $f_i(\theta) = \ell_i(y_i, \Phi(x_i)^\top \theta)$, then $f_i'(\theta) = \ell_i'(y_i, \Phi(x_i)^\top \theta) \Phi(x_i)$
 - Only need to store n real numbers

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 (with potentially $\mu = 0$)
- constant step size $\gamma_t = 1/(16L)$
- initialization with one pass of averaged SGD

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- initialization with one pass of averaged SGD
- Strongly convex case (Le Roux et al., 2012, 2013)

$$\mathbb{E}\left[g(\theta_t) - g(\theta_*)\right] \leqslant \left(\frac{8\sigma^2}{n} + \frac{4L\|\theta_0 - \theta_*\|^2}{n}\right) \exp\left(-t \min\left\{\frac{1}{8n}, \frac{\mu}{16L}\right\}\right)$$

- Linear (exponential) convergence rate with O(1) iteration cost
- After one pass, reduction of cost by $\exp\left(-\min\left\{\frac{1}{8},\frac{n\mu}{16L}\right\}\right)$

Stochastic average gradient - Convergence analysis

Assumptions

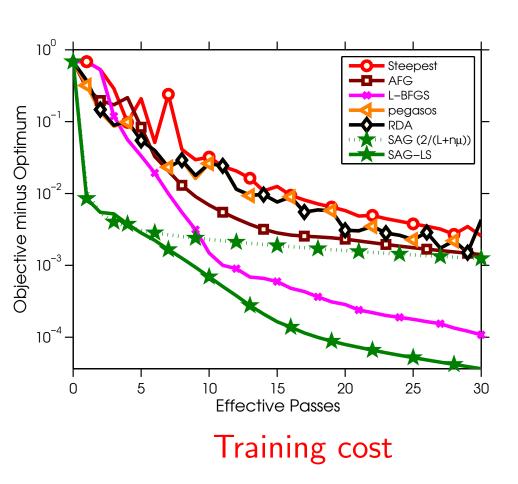
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- Non-strongly convex case (Le Roux et al., 2013)

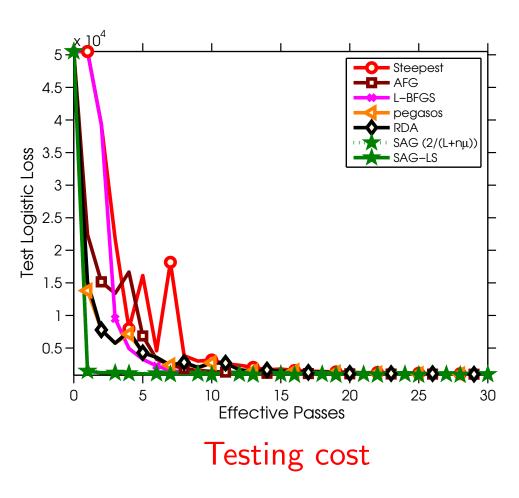
$$\mathbb{E}\left[g(\theta_t) - g(\theta_*)\right] \leqslant 48 \frac{\sigma^2 + L\|\theta_0 - \theta_*\|^2}{\sqrt{n}} \frac{n}{k}$$

- Improvement over regular batch and stochastic gradient
- Adaptivity to potentially hidden strong convexity

Stochastic average gradient Simulation experiments

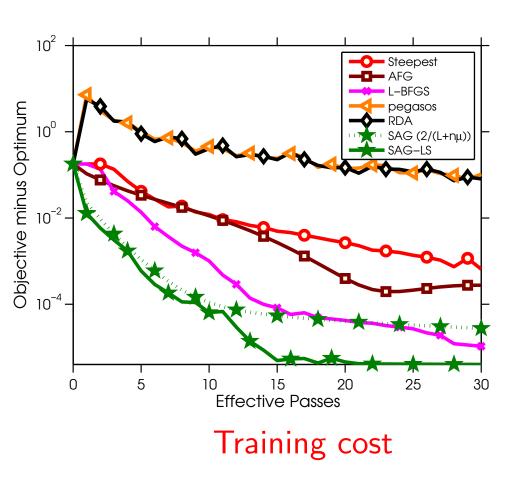
- protein dataset (n = 145751, p = 74)
- Dataset split in two (training/testing)

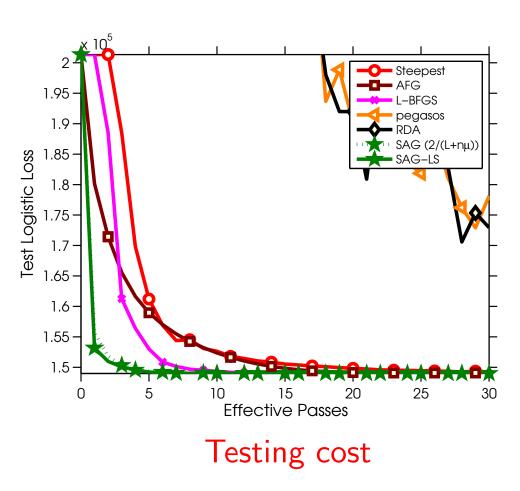




Stochastic average gradient Simulation experiments

- cover type dataset (n = 581012, p = 54)
- Dataset split in two (training/testing)





Conclusions / Extensions Stochastic average gradient

- Going beyond a single pass through the data
 - Keep memory of all gradients for finite training sets
 - Linear convergence rate with O(1) iteration complexity
 - Randomization leads to easier analysis and faster rates

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Future/current work - open problems

- Including a non-differentiable term
- Line search
- Using second-order information or non-uniform sampling
- Distributed optimization
- Going beyond finite training sets (bound on testing cost)

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