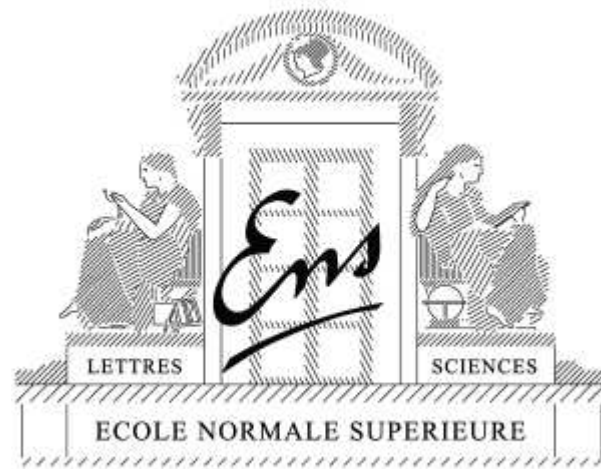


Stochastic gradient methods for machine learning

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Joint work with Eric Moulines, Nicolas Le Roux
and Mark Schmidt - January 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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- **Examples:** computer vision, bioinformatics, signal processing
- **Ideal running-time complexity:** $O(pn + kn)$
- **Going back to simple methods**
 - Stochastic gradient methods (Robbins and Monro, 1951)
 - Mixing statistics and optimization
 - It is possible to improve on the sublinear convergence rate?

Outline

- **Introduction**

- Supervised machine learning and convex optimization
- Beyond the separation of statistics and optimization

- **Stochastic approximation algorithms** (Bach and Moulines, 2011)

- Stochastic gradient and averaging
- **Strongly convex vs. non-strongly convex**

- **Going beyond stochastic gradient** (Le Roux, Schmidt, and Bach, 2012)

- More than a single pass through the data
- **Linear (exponential) convergence rate for strongly convex functions**

Supervised machine learning

- **Data:** n observations $(x_i, y_i) \in \mathcal{X} \times \mathcal{Y}$, $i = 1, \dots, n$, **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

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

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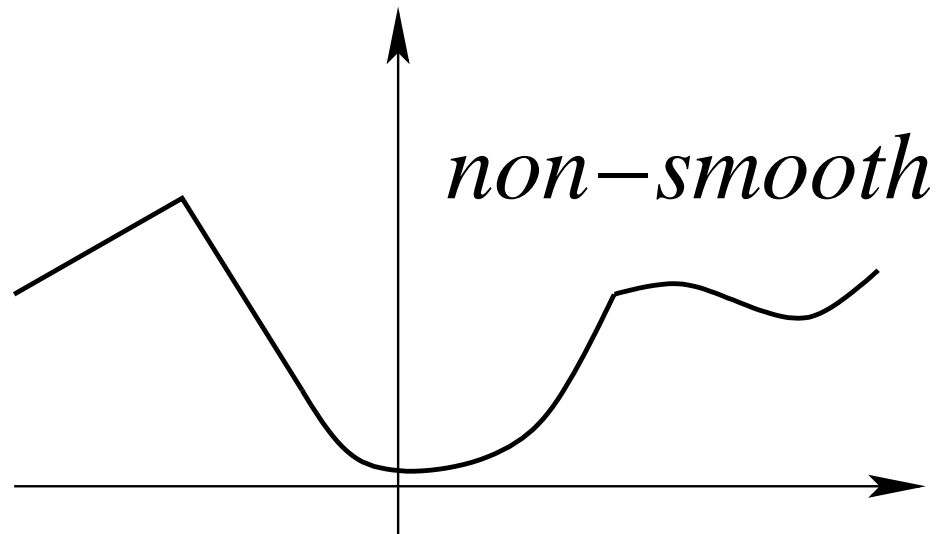
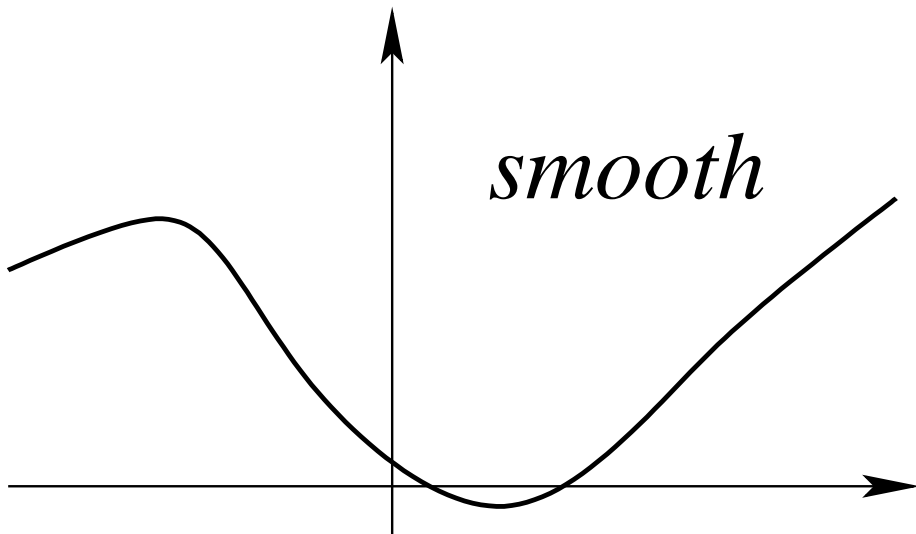
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- **Two fundamental questions:** (1) computing $\hat{\theta}$ and (2) analyzing $\hat{\theta}$
 - **May be tackled simultaneously**

Smoothness and strong convexity

- A function $g : \mathbb{R}^p \rightarrow \mathbb{R}$ is **L -smooth** if and only if it is differentiable and its gradient is L -Lipschitz-continuous

$$\forall \theta_1, \theta_2 \in \mathbb{R}^p, \|g'(\theta_1) - g'(\theta_2)\| \leq L \|\theta_1 - \theta_2\|$$

- If g is twice differentiable: $\forall \theta \in \mathbb{R}^p, g''(\theta) \preceq L \cdot Id$



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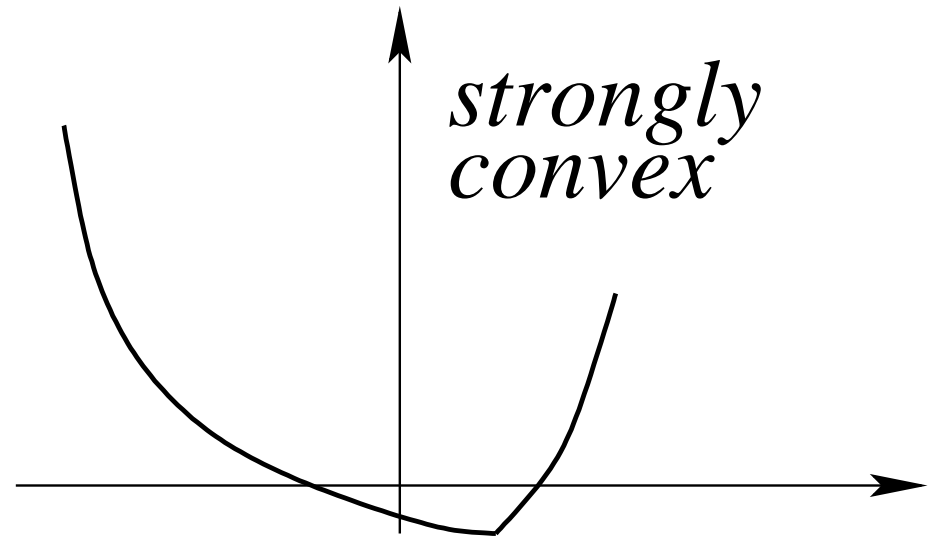
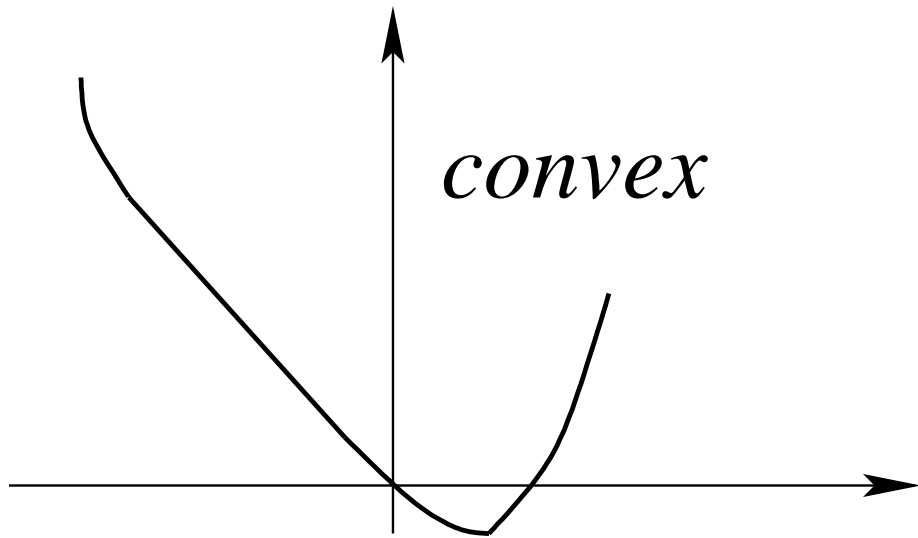
- with $g(\theta) = \frac{1}{n} \sum_{i=1}^n \ell(y_i, \theta^\top \Phi(x_i))$
- Hessian \approx covariance matrix $\frac{1}{n} \sum_{i=1}^n \Phi(x_i) \Phi(x_i)^\top$
- **Bounded data**

Smoothness and strong convexity

- A function $g : \mathbb{R}^p \rightarrow \mathbb{R}$ is μ -strongly convex if and only if

$$\forall \theta_1, \theta_2 \in \mathbb{R}^p, g(\theta_1) \geq g(\theta_2) + \langle g'(\theta_2), \theta_1 - \theta_2 \rangle + \frac{\mu}{2} \|\theta_1 - \theta_2\|^2$$

- Equivalent definition: $\theta \mapsto g(\theta) - \frac{\mu}{2} \|\theta\|^2$ is convex
- If g is twice differentiable: $\forall \theta \in \mathbb{R}^p, g''(\theta) \succeq \mu \cdot Id$



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- If g is twice differentiable: $\forall \theta \in \mathbb{R}^p, \quad g''(\theta) \succeq \mu \cdot Id$
- **Machine learning**
 - with $g(\theta) = \frac{1}{n} \sum_{i=1}^n \ell(y_i, \theta^\top \Phi(x_i))$
 - Hessian \approx covariance matrix $\frac{1}{n} \sum_{i=1}^n \Phi(x_i) \Phi(x_i)^\top$
 - **Data with invertible covariance matrix** (low correlation/dimension)
 - ... or with added regularization by $\frac{\mu}{2} \|\theta\|^2$

Stochastic approximation

- **Goal:** Minimizing a function f defined on a Hilbert space \mathcal{H}
 - given only unbiased estimates $f'_n(\theta_n)$ of its gradients $f'(\theta_n)$ at certain points $\theta_n \in \mathcal{H}$
- **Stochastic approximation**
 - Observation of $f'_n(\theta_n) = f'(\theta_n) + \varepsilon_n$, with $\varepsilon_n =$ i.i.d. noise
 - Non-convex problems

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 - Non-convex problems
- **Machine learning - statistics**
 - **loss for a single pair of observations:** $f_n(\theta) = \ell(y_n, \theta^\top \Phi(x_n))$
 - $f(\theta) = \mathbb{E} f_n(\theta) = \mathbb{E} \ell(y_n, \theta^\top \Phi(x_n)) =$ **generalization error**
 - Expected gradient: $f'(\theta) = \mathbb{E} f'_n(\theta) = \mathbb{E} \{ \ell'(y_n, \theta^\top \Phi(x_n)) \Phi(x_n) \}$

Convex smooth stochastic approximation

- Key properties of f and/or f_n
 - Smoothness: f_n L -smooth
 - Strong convexity: f μ -strongly convex

Convex smooth stochastic approximation

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 - **Smoothness:** f_n L -smooth
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- **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:

$$\gamma_n = Cn^{-\alpha}$$

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- **Desirable practical behavior**

- Applicable (at least) to least-squares and logistic regression
- Robustness to (potentially unknown) constants (L, μ)
- Adaptivity to difficulty of the problem (e.g., strong convexity)

Convex stochastic approximation

Related work

- **Machine learning/optimization**

- Known minimax rates of convergence (Nemirovski and Yudin, 1983; Agarwal et al., 2010)
 - **Strongly convex: $O(n^{-1})$**
 - **Non-strongly convex: $O(n^{-1/2})$**
- Achieved with and/or without averaging (up to log terms)
- Non-asymptotic analysis (high-probability bounds)
- Online setting and regret bounds
- 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)

Convex stochastic approximation

Related work

- **Stochastic approximation**

- Asymptotic analysis
- Non convex case with strong convexity around the optimum
- $\gamma_n = Cn^{-\alpha}$ with $\alpha = 1$ is not robust to the choice of C
- $\alpha \in (1/2, 1)$ is robust **with averaging**
- Broadie et al. (2009); Kushner and Yin (2003); Kul'chitskiĭ and Mozhgovoĭ (1991); Fabian (1968)
- Polyak and Juditsky (1992); Ruppert (1988)

Problem set-up - General assumptions

- **Unbiased gradient estimates:**

- $f_n(\theta)$ is of the form $h(z_n, \theta)$, where z_n is an i.i.d. sequence
- e.g., $f_n(\theta) = h(z_n, \theta) = \ell(y_n, \theta^\top \Phi(x_n))$ with $z_n = (x_n, y_n)$
- NB: can be generalized

- **Variance of estimates:** There exists $\sigma^2 \geq 0$ such that for all $n \geq 1$, $\mathbb{E}(\|f'_n(\theta^*) - f'(\theta^*)\|^2) \leq \sigma^2$, where θ^* is a global minimizer of f

Problem set-up - Smoothness/convexity assumptions

- **Smoothness of f_n** : For each $n \geq 1$, the function f_n is a.s. convex, differentiable with L -Lipschitz-continuous gradient f'_n :
 - Bounded data

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 - Bounded data
- **Strong convexity of f** : The function f is strongly convex with respect to the norm $\|\cdot\|$, with convexity constant $\mu > 0$:
 - Invertible population covariance matrix
 - or regularization by $\frac{\mu}{2}\|\theta\|^2$

Summary of new results (Bach and Moulines, 2011)

- Stochastic gradient descent with learning rate $\gamma_n = Cn^{-\alpha}$
- **Strongly convex smooth objective functions**
 - Old: $O(n^{-1})$ rate achieved **without** averaging for $\alpha = 1$
 - New: $O(n^{-1})$ rate achieved **with** averaging for $\alpha \in [1/2, 1]$
 - Non-asymptotic analysis with explicit constants
 - Forgetting of initial conditions
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 - Robustness to the choice of C
- **Proof technique**
 - Derive deterministic recursion for $\delta_n = \mathbb{E}\|\theta_n - \theta^*\|^2$
$$\delta_n \leq (1 - 2\mu\gamma_n + 2L^2\gamma_n^2)\delta_{n-1} + 2\sigma^2\gamma_n^2$$
 - Mimic SA proof techniques in a non-asymptotic way

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 - Non-asymptotic analysis with explicit constants
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 - Robustness to the choice of C
- **Convergence rates** for $\mathbb{E}\|\theta_n - \theta^*\|^2$ and $\mathbb{E}\|\bar{\theta}_n - \theta^*\|^2$
 - no averaging: $O\left(\frac{\sigma^2\gamma_n}{\mu}\right) + O(e^{-\mu n\gamma_n})\|\theta_0 - \theta^*\|^2$
 - averaging: $\frac{\text{tr } H(\theta^*)^{-1}}{n} + \mu^{-1}O(n^{-2\alpha} + n^{-2+\alpha}) + O\left(\frac{\|\theta_0 - \theta^*\|^2}{\mu^2 n^2}\right)$

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- **Non-strongly convex smooth objective functions**
 - Old: $O(n^{-1/2})$ rate achieved **with** averaging for $\alpha = 1/2$
 - New: $O(\max\{n^{1/2-3\alpha/2}, n^{-\alpha/2}, n^{\alpha-1}\})$ rate achieved **without** averaging for $\alpha \in [1/3, 1]$
- **Take-home message**
 - Use $\alpha = 1/2$ with averaging to be adaptive to strong convexity

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

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
- **Future/current work - open problems**
 - High-probability through all moments $\mathbb{E}\|\theta_n - \theta^*\|^{2d}$
 - Analysis for logistic regression using self-concordance (Bach, 2010)
 - Including a non-differentiable term (Xiao, 2010; Lan, 2010)
 - Non-random errors (Schmidt, Le Roux, and Bach, 2011)
 - Line search for stochastic gradient
 - Non-parametric stochastic approximation
 - Online estimation of uncertainty
 - 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 $\mathbb{E}_z h(\theta, z) = \mathbb{E}_{(x,y)} \ell(y, \theta^\top \Phi(x))$

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- **Machine learning practice**

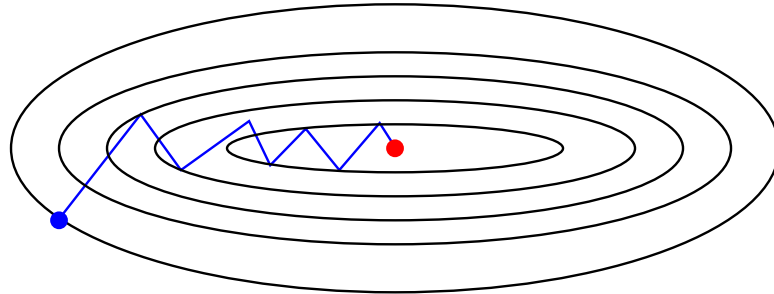
- Finite data set (z_1, \dots, z_n)
- Multiple passes
- Minimizes **training** cost $\frac{1}{n} \sum_{i=1}^n h(\theta, z_i) = \frac{1}{n} \sum_{i=1}^n \ell(y_i, \theta^\top \Phi(x_i))$
- Need to regularize (e.g., by the ℓ_2 -norm) to avoid overfitting

Stochastic vs. deterministic methods

- 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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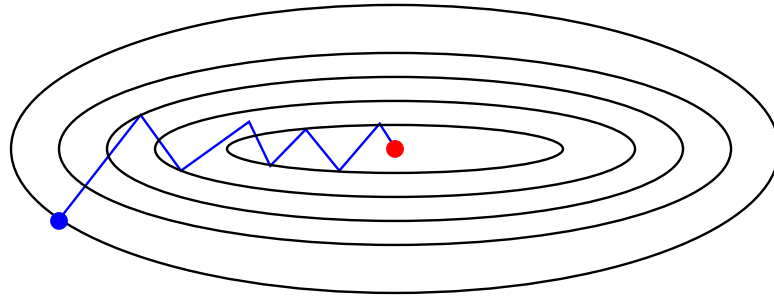


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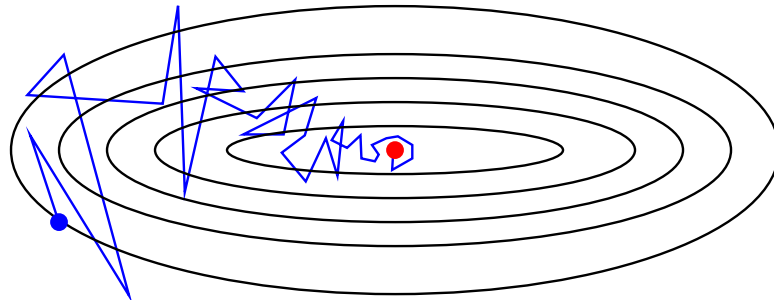
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 - Linear (e.g., exponential) convergence rate
 - Iteration complexity is linear in n
- **Stochastic** gradient descent: $\theta_t = \theta_{t-1} - \gamma_t f'_{i(t)}(\theta_{t-1})$
 - Sampling with replacement: $i(t)$ random element of $\{1, \dots, n\}$
 - Convergence rate in $O(1/t)$
 - Iteration complexity is independent of n

Stochastic vs. deterministic methods

- 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: $\theta_t = \theta_{t-1} - \gamma_t g'(\theta_{t-1}) = \theta_{t-1} - \frac{\gamma_t}{n} \sum_{i=1}^n f'_i(\theta_{t-1})$

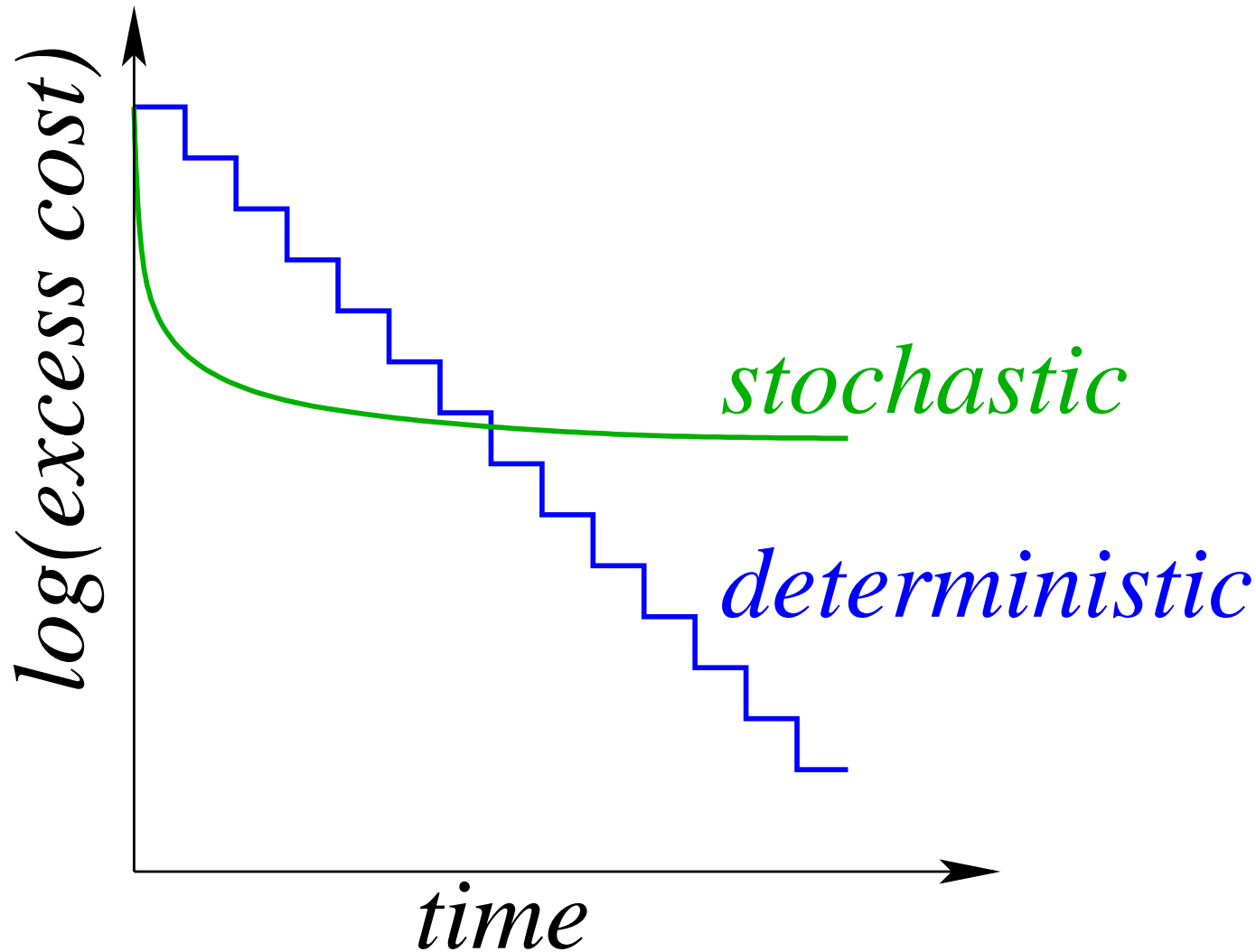


- **Stochastic** gradient descent: $\theta_t = \theta_{t-1} - \gamma_t f'_{i(t)}(\theta_{t-1})$



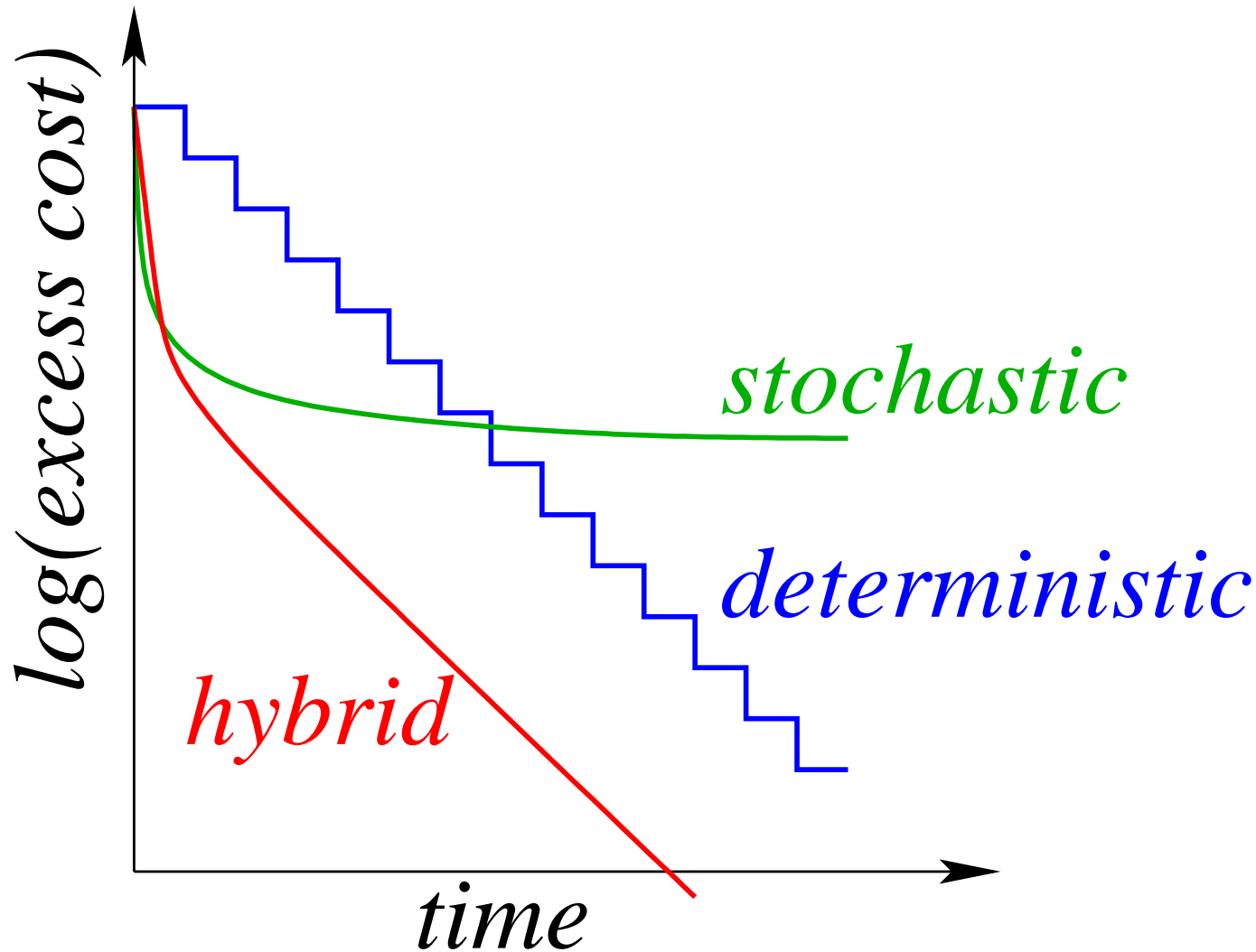
Stochastic vs. deterministic methods

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



Stochastic vs. deterministic methods

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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 , $i = 1, \dots, n$
 - Random selection $i(t) \in \{1, \dots, n\}$ with replacement
 - Iteration: $\theta_t = \theta_{t-1} - \frac{\gamma_t}{n} \sum_{i=1}^n y_i^t$ with $y_i^t = \begin{cases} f'_i(\theta_{t-1}) & \text{if } i = i(t) \\ y_i^{t-1} & \text{otherwise} \end{cases}$

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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 - I

- Assume each f_i is L -smooth and $\hat{f} = \frac{1}{n} \sum_{i=1}^n f_i$ is μ -strongly convex

- **Constant step size** $\gamma_t = \frac{1}{2nL}$:

$$\mathbb{E}[\|\theta_t - \theta^*\|^2] \leq \left(1 - \frac{\mu}{8Ln}\right)^t \left[3\|\theta_0 - \theta^*\|^2 + \frac{9\sigma^2}{4L^2}\right]$$

- Linear rate with iteration cost independent of n ...
- ... but, same behavior as batch gradient and IAG (cyclic version)
- **Proof technique**
 - Designing a quadratic Lyapunov function for a n -th order non-linear stochastic dynamical system

Stochastic average gradient

Convergence analysis - II

- Assume each f_i is L -smooth and $\hat{f} = \frac{1}{n} \sum_{i=1}^n f_i$ is μ -strongly convex
- **Constant step size** $\gamma_t = \frac{1}{2n\mu}$, if $\frac{\mu}{L} \geq \frac{8}{n}$

$$\mathbb{E}[\hat{f}(\theta_t) - \hat{f}(\theta^*)] \leq C \left(1 - \frac{1}{8n}\right)^t$$

with $C = \left[\frac{16L}{3n} \|\theta_0 - \theta^*\|^2 + \frac{4\sigma^2}{3n\mu} \left(8 \log \left(1 + \frac{\mu n}{4L} \right) + 1 \right) \right]$

- Linear rate with iteration cost independent of n
- Linear convergence rate “independent” of the condition number
- After each pass through the data, constant error reduction

Rate of convergence comparison

- Assume that $L = 100$, $\mu = .01$, and $n = 80000$

- Full gradient method has rate

$$\left(1 - \frac{\mu}{L}\right) = 0.9999$$

- Accelerated gradient method has rate

$$\left(1 - \sqrt{\frac{\mu}{L}}\right) = 0.9900$$

- Running n iterations of SAG for the same cost has rate

$$\left(1 - \frac{1}{8n}\right)^n = 0.8825$$

- *Fastest possible* first-order method has rate

$$\left(\frac{\sqrt{L} - \sqrt{\mu}}{\sqrt{L} + \sqrt{\mu}}\right)^2 = 0.9608$$

- **Beating two lower bounds** (with additional assumptions)

- (1) stochastic gradient and (2) full gradient

Stochastic average gradient

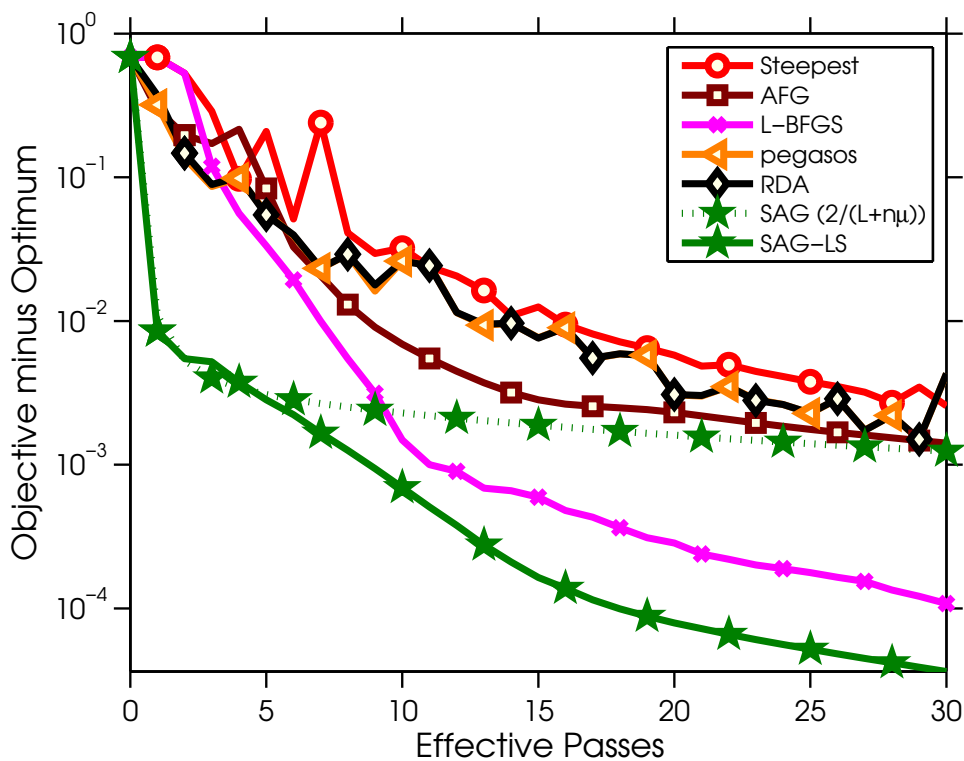
Implementation details and extensions

- The algorithm can use **sparsity** in the features to reduce the storage and iteration cost
- **Grouping functions together** can further reduce the memory requirement
- We have obtained good performance when L is not known with a **heuristic line-search**
- Algorithm allows **non-uniform sampling**
- Possibility of making **proximal, coordinate-wise, and Newton-like** variants

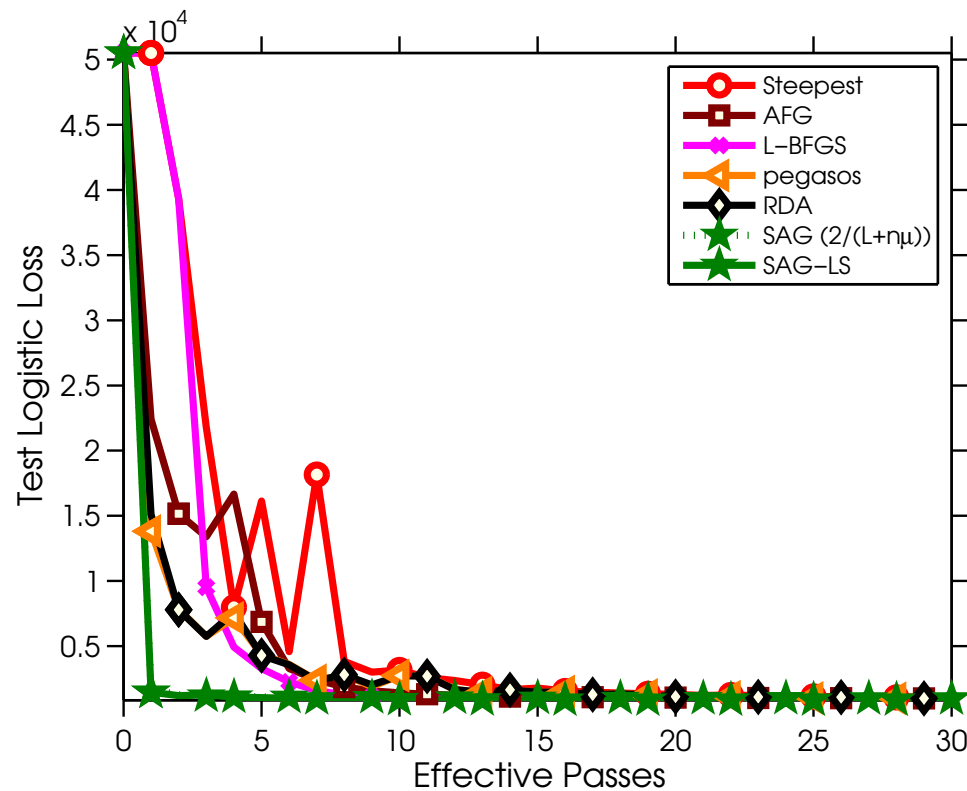
Stochastic average gradient

Simulation experiments

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



Training cost

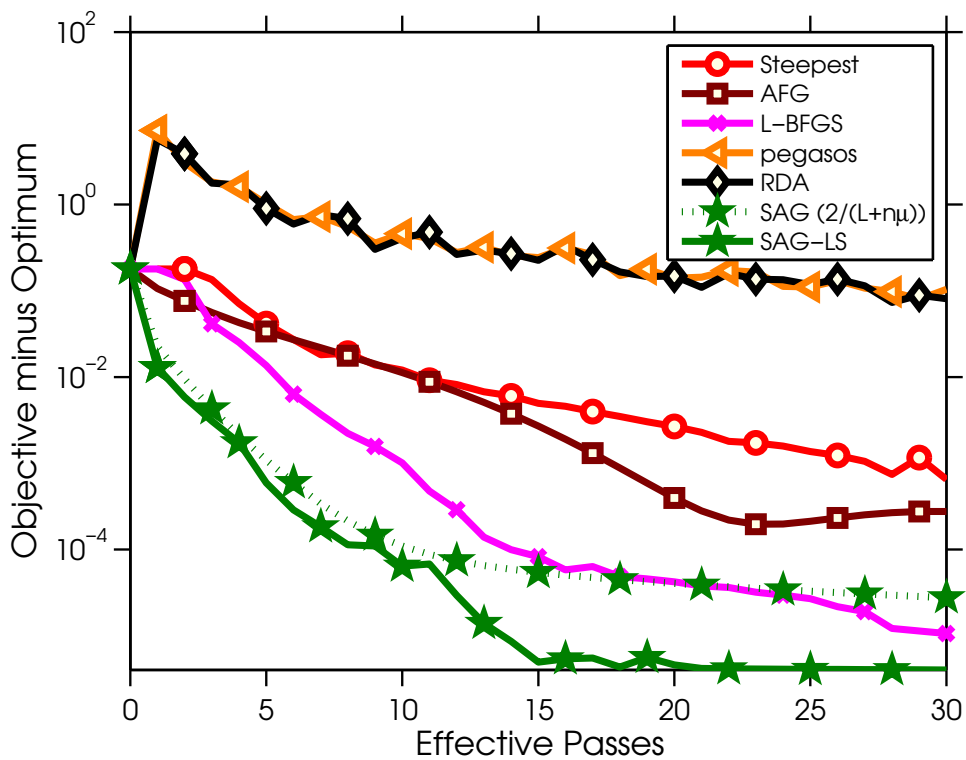


Testing cost

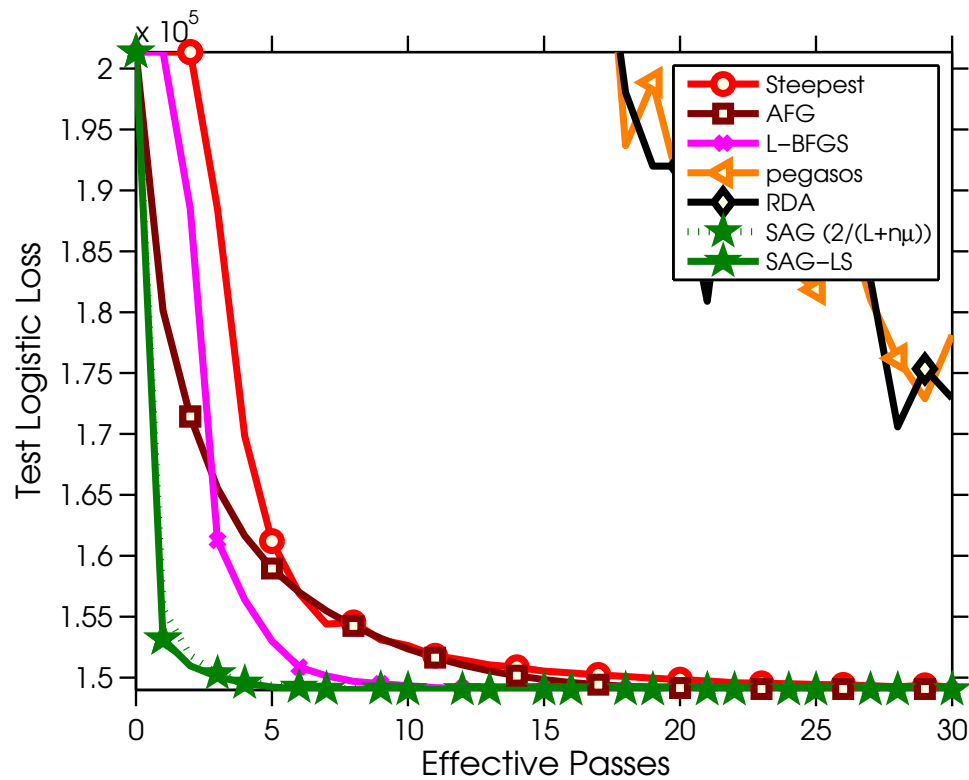
Stochastic average gradient

Simulation experiments

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



Training cost



Testing cost

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
 - Beyond machine learning

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
 - Beyond machine learning
- **Future/current work - open problems**
 - Including a non-differentiable term
 - Line search
 - Using second-order information or non-uniform sampling
 - **Going beyond finite training sets (bound on testing cost)**
 - **Non strongly-convex case**

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