# Stochastic gradient methods for machine learning

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

## Context Machine learning for "big data"

- Large-scale machine learning: large p, large n, large k
  - -p: dimension of each observation (input)
  - -n: number of observations
  - -k: number of tasks (dimension of outputs)
- **Examples**: computer vision, bioinformatics, text processing

# **Search engines - advertising**

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## **Advertising - recommendation**



## **Object recognition**



## **Learning for bioinformatics - Proteins**

- Crucial components of cell life
- Predicting multiple functions and interactions
- Massive data: up to 1 millions for humans!
- Complex data
  - Amino-acid sequence
  - Link with DNA
  - Tri-dimensional molecule



# Context

## Machine learning for "big data"

- Large-scale machine learning: large p, large n, large k
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- **Examples**: computer vision, bioinformatics, text processing
- Ideal running-time complexity: O(pn + kn)

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- Large-scale machine learning: large p, large n, large k
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- **Examples**: computer vision, bioinformatics, text processing
- 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: stochastic approximation algorithms

- Supervised machine learning and convex optimization
- Stochastic gradient and averaging
- Strongly convex vs. non-strongly convex
- Fast convergence through smoothness and constant step-sizes
  - Online Newton steps (Bach and Moulines, 2013)
  - O(1/n) convergence rate for all convex functions
- More than a single pass through the data
  - Stochastic average gradient (Le Roux, Schmidt, and Bach, 2012)
  - 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  $\langle \theta, \Phi(x) \rangle$  of features  $\Phi(x) \in \mathbb{R}^p$
- (regularized) empirical risk minimization: find  $\hat{\theta}$  solution of

$$\min_{\theta \in \mathbb{R}^p} \quad \frac{1}{n} \sum_{i=1}^n \ell(y_i, \langle \theta, \Phi(x_i) \rangle) + \mu \Omega(\theta)$$
  
convex data fitting term + regularizer

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• Empirical risk:  $\hat{f}(\theta) = \frac{1}{n} \sum_{i=1}^{n} \ell(y_i, \langle \theta, \Phi(x_i) \rangle)$  training cost

• Expected risk:  $f(\theta) = \mathbb{E}_{(x,y)} \ell(y, \langle \theta, \Phi(x) \rangle)$  testing cost

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

• A function  $g: \mathbb{R}^p \to \mathbb{R}$  is *L*-smooth if and only if it is twice differentiable and

 $\forall \theta \in \mathbb{R}^p, \ g''(\theta) \preccurlyeq L \cdot \mathrm{Id}$ 



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

• A function  $g: \mathbb{R}^p \to \mathbb{R}$  is  $\mu$ -strongly convex if and only if

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

• If g is twice differentiable:  $\forall \theta \in \mathbb{R}^p, g''(\theta) \succcurlyeq \mu \cdot \mathrm{Id}$ 



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  - Data with invertible covariance matrix (low correlation/dimension)
- Adding regularization by  $\frac{\mu}{2} \|\theta\|^2$

– creates additional bias unless  $\mu$  is small

### Iterative methods for minimizing smooth functions

- Assumption: g convex and smooth on  $\mathbb{R}^p$
- 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:  $\theta_t = \theta_{t-1} g''(\theta_{t-1})^{-1}g'(\theta_{t-1})$ 
  - $O(e^{-\rho 2^t})$  convergence rate

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### • Key insights from Bottou and Bousquet (2008)

In machine learning, no need to optimize below statistical error
 In machine learning, cost functions are averages

 $\Rightarrow$  Stochastic approximation

## **Stochastic approximation**

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

### • 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:

$$f_n(\theta) = \ell(y_n, \langle \theta, \Phi(x_n) \rangle)$$

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

### **Convex 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:  $\bar{\theta}_n = \frac{1}{n+1} \sum_{k=0}^n \theta_k$
- Which learning rate sequence  $\gamma_n$ ? Classical setting:

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

- Known global minimax rates of convergence for non-smooth problems (Nemirovsky and Yudin, 1983; Agarwal et al., 2012)
  - Strongly convex:  $O((\mu n)^{-1})$

Attained by averaged stochastic gradient descent with  $\gamma_n \propto (\mu n)^{-1}$ 

– Non-strongly convex:  $O(n^{-1/2})$ 

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- Non-strongly convex:  $O(n^{-1/2})$ Attained by averaged stochastic gradient descent with  $\gamma_n \propto n^{-1/2}$
- Many contributions in optimization and online learning: 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)
  - All step sizes  $\gamma_n = Cn^{-\alpha}$  with  $\alpha \in (1/2, 1)$  lead to  $O(n^{-1})$  for smooth strongly convex problems

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- A single algorithm for smooth problems with convergence rate O(1/n) in all situations?

### Least-mean-square algorithm

- Least-squares:  $f(\theta) = \frac{1}{2}\mathbb{E}[(y_n \langle \Phi(x_n), \theta \rangle)^2]$  with  $\theta \in \mathbb{R}^p$ 
  - SGD = least-mean-square algorithm (see, e.g., Macchi, 1995)
  - usually studied without averaging and decreasing step-sizes
  - with strong convexity assumption  $\mathbb{E}[\Phi(x_n) \otimes \Phi(x_n)] = H \succcurlyeq \mu \cdot \mathrm{Id}$

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- New analysis for averaging and constant step-size  $\gamma = 1/(4R^2)$ 
  - Assume  $\|\Phi(x_n)\| \leq R$  and  $|y_n \langle \Phi(x_n), \theta_* \rangle| \leq \sigma$  almost surely
  - No assumption regarding lowest eigenvalues of  ${\cal H}$
  - Main result:  $\left| \mathbb{E}f(\bar{\theta}_{n-1}) f(\theta_*) \leqslant \frac{2}{n} \left[ \sigma \sqrt{p} + R \| \theta_0 \theta_* \| \right]^2 \right|$
- Matches statistical lower bound (Tsybakov, 2003)

### Markov chain interpretation of constant step sizes

• LMS recursion for  $f_n(\theta) = \frac{1}{2} (y_n - \langle \Phi(x_n), \theta \rangle)^2$ 

$$\theta_n = \theta_{n-1} - \gamma \big( \langle \Phi(x_n), \theta_{n-1} \rangle - y_n \big) \Phi(x_n)$$

- The sequence  $(\theta_n)_n$  is a homogeneous Markov chain
  - convergence to a stationary distribution  $\pi_{\gamma}$
  - with expectation  $\bar{\theta}_{\gamma} \stackrel{\text{def}}{=} \int \theta \pi_{\gamma}(\mathrm{d}\theta)$

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  - convergence to a stationary distribution  $\pi_{\gamma}$ - with expectation  $\bar{\theta}_{\gamma} \stackrel{\text{def}}{=} \int \theta \pi_{\gamma}(\mathrm{d}\theta)$
- For least-squares,  $\bar{\theta}_{\gamma}=\theta_{*}$ 
  - $\theta_n$  does not converge to  $\theta_*$  but oscillates around it oscillations of order  $\sqrt{\gamma}$
- Ergodic theorem:
  - Averaged iterates converge to  $\bar{\theta}_{\gamma} = \theta_*$  at rate O(1/n)

#### **Simulations - synthetic examples**

• Gaussian distributions - p=20



#### **Simulations - benchmarks**



## **Beyond least-squares - Markov chain interpretation**

- Recursion  $\theta_n = \theta_{n-1} \gamma f'_n(\theta_{n-1})$  also defines a Markov chain
  - Stationary distribution  $\pi_{\gamma}$  such that  $\int f'(\theta) \pi_{\gamma}(\mathrm{d}\theta) = 0$
  - When f' is not linear,  $f'(\int \theta \pi_{\gamma}(\mathrm{d}\theta)) \neq \int f'(\theta) \pi_{\gamma}(\mathrm{d}\theta) = 0$

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  - Stationary distribution  $\pi_{\gamma}$  such that  $\int f'(\theta) \pi_{\gamma}(d\theta) = 0$
  - When f' is not linear,  $f'(\int \theta \pi_{\gamma}(\mathrm{d}\theta)) \neq \int f'(\theta) \pi_{\gamma}(\mathrm{d}\theta) = 0$
- $\theta_n$  oscillates around the wrong value  $\bar{\theta}_{\gamma} \neq \theta_*$

– moreover, 
$$\|\theta_* - \theta_n\| = O_p(\sqrt{\gamma})$$

#### Ergodic theorem

- averaged iterates converge to  $\bar{\theta}_{\gamma} \neq \theta_*$  at rate O(1/n)
- moreover,  $\|\theta_* \bar{\theta}_{\gamma}\| = O(\gamma)$  (Bach, 2013)

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### **Restoring convergence through online Newton steps**

• The Newton step for  $f = \mathbb{E}f_n(\theta) \stackrel{\text{def}}{=} \mathbb{E}[\ell(y_n, \langle \theta, \Phi(x_n) \rangle)]$  at  $\tilde{\theta}$  is equivalent to minimizing the quadratic approximation

$$g(\theta) = f(\tilde{\theta}) + \langle f'(\tilde{\theta}), \theta - \tilde{\theta} \rangle + \frac{1}{2} \langle \theta - \tilde{\theta}, f''(\tilde{\theta})(\theta - \tilde{\theta}) \rangle$$
  
$$= f(\tilde{\theta}) + \langle \mathbb{E}f'_{n}(\tilde{\theta}), \theta - \tilde{\theta} \rangle + \frac{1}{2} \langle \theta - \tilde{\theta}, \mathbb{E}f''_{n}(\tilde{\theta})(\theta - \tilde{\theta}) \rangle$$
  
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• Complexity of least-mean-square recursion for g is O(p)

$$\theta_n = \theta_{n-1} - \gamma \left[ f'_n(\tilde{\theta}) + f''_n(\tilde{\theta})(\theta_{n-1} - \tilde{\theta}) \right]$$

 $-f_n''(\tilde{\theta}) = \ell''(y_n, \langle \tilde{\theta}, \Phi(x_n) \rangle) \Phi(x_n) \otimes \Phi(x_n)$  has rank one

New online Newton step without computing/inverting Hessians

## **Choice of support point for online Newton step**

#### • Two-stage procedure

- (1) Run n/2 iterations of averaged SGD to obtain  $\tilde{ heta}$
- (2) Run n/2 iterations of averaged constant step-size LMS
  - Reminiscent of one-step estimators (see, e.g., Van der Vaart, 2000)
  - Provable convergence rate of O(p/n) for logistic regression
  - Additional assumptions but no strong convexity

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  - Reminiscent of one-step estimators (see, e.g., Van der Vaart, 2000)
  - Provable convergence rate of O(p/n) for logistic regression
  - Additional assumptions but no strong convexity
- Update at each iteration using the current averaged iterate
  - Recursion:  $\theta_n = \theta_{n-1} \gamma \left[ f'_n(\bar{\theta}_{n-1}) + f''_n(\bar{\theta}_{n-1})(\theta_{n-1} \bar{\theta}_{n-1}) \right]$
  - No provable convergence rate (yet) but best practical behavior
  - Note (dis)similarity with regular SGD:  $\theta_n = \theta_{n-1} \gamma f'_n(\theta_{n-1})$

### **Simulations - synthetic examples**

• Gaussian distributions - p=20



#### **Simulations - benchmarks**



## Going beyond a single pass over the data

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

## Going beyond a single pass over the data

#### • Stochastic approximation

- Assumes infinite data stream
- Observations are used only once
- Directly minimizes testing cost  $\mathbb{E}_{(x,y)} \ell(y, \langle \theta, \Phi(x) \rangle)$
- Machine learning practice
  - Finite data set  $(x_1, y_1, \ldots, x_n, y_n)$
  - Multiple passes
  - Minimizes training cost  $\frac{1}{n} \sum_{i=1}^{n} \ell(y_i, \langle \theta, \Phi(x_i) \rangle)$
  - Need to regularize (e.g., by the  $\ell_2\text{-norm})$  to avoid overfitting

• **Goal**: minimize 
$$g(\theta) = \frac{1}{n} \sum_{i=1}^{n} f_i(\theta)$$

• 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 in  $O(e^{-\alpha t})$
  - Iteration complexity is linear in n (with line search)

- 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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  - Linear (e.g., exponential) convergence rate in  $O(e^{-\alpha t})$
  - Iteration complexity is linear in n (with line search)
- 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 (step size selection?)

• Goal = best of both worlds: Linear rate with O(1) iteration cost Robustness to step size



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# 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  $i(t) \in \{1, \ldots, n\}$  with replacement

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

# 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  $i(t) \in \{1, \ldots, n\}$  with replacement

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

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

### **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  $\mu$ -strongly convex (with potentially  $\mu = 0$ )
- constant step size  $\gamma_t = 1/(16L)$
- initialization with one pass of averaged SGD
- Strongly convex case (Le Roux et al., 2012, 2013)

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

- 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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- constant step size  $\gamma_t = 1/(16L)$
- initialization with one pass of averaged SGD
- Non-strongly convex case (Le Roux et al., 2013)

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

- 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**

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



## Conclusions

#### • Constant-step-size averaged stochastic gradient descent

- Reaches convergence rate  ${\cal O}(1/n)$  in all regimes
- Improves on the  $O(1/\sqrt{n})$  lower-bound of non-smooth problems
- Efficient online Newton step for non-quadratic problems

#### • Going beyond a single pass through the data

- Keep memory of all gradients for finite training sets
- Randomization leads to easier analysis and faster rates
- Relationship with Shalev-Shwartz and Zhang (2012); Mairal (2013)

### • Extensions

- Non-differentiable terms, kernels, line-search, parallelization, etc.

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