Outline for today

- Methods for learning the net parameters:
 - Stochastic gradient descent (SGD)
 - Polyak and Nesterov momentum
 - AdaGrad and variants toward Adam
- Simplified version of AdaGrad with scalar, adaptive, stepsize
- ► Convergence dependence of scalar AdaGrad on initial stepsize

Stochastic gradient descent (SGD)

Given a loss function $\mathcal{L}(\theta; X, Y)$, gradient descent is given by

$$\theta^{(k+1)} = \theta^{(k)} - \eta \cdot \mathsf{grad}_{\theta} \mathcal{L}(\theta, X, Y)$$

where η is referred to as the stepsize, or in deep learning the "learning rate."

Recall, we typically have a loss function which is the sum of n individual loss functions, independent for each data point:

$$\mathcal{L}(\theta; X, Y) = n^{-1} \sum_{\mu=1}^{n} I(\theta; x_{\mu}, y_{\mu})$$

For $n\gg 1$ gradient descent is computationally too costly and instead one can break appart the n loss functions into "mini-batches" and repeatedly solve

$$\theta^{(k+1)} = \theta^{(k)} - \eta |\Lambda_k|^{-1} \operatorname{grad}_{\theta} \sum_{\mu \in \Lambda_k} I(\theta; x_{\mu}, y_{\mu}).$$

This is referred to as stochastic gradient descent as typically Λ_k is chosen in some randomized method, usually as a partition of [n] and a sequence of Λ_k which cover [n] is referred to as an "epoch."

Stochastic gradient descent: challenges and benefits

$$\theta^{(k+1)} = \theta^{(k)} - \eta |\Lambda_k|^{-1} \operatorname{grad}_{\theta} \sum_{\mu \in \Lambda_k} I(\theta; x_{\mu}, y_{\mu}).$$

- SGD is preferable for large n as it reduces the per iteration computational cost dependence on n to instead depend on $|\Lambda_k|$ which can be set by the user as opposed to n which is given by the data set.
- SGD, and gradient descent, require selection of a learning rate (stepsize) which in deep learning is typically selected using some costly trial and error heuristics.
- ▶ The learning rate is typically chosen adaptively in a way that satisfies $\sum_{k=1}^{\infty} \eta_k = \infty$ and $\sum_{k=1}^{\infty} \eta_k^2 < \infty$; in particular as $\eta_k \sim k^{-1}$.
- ► The optimal selection of learning weight, and selection of Λ , depends on the unknown local Lipschitz constant $\|\operatorname{grad} I(\theta_1; x_\mu, y_\mu) \operatorname{grad} I(\theta_2; x_\mu, y_\mu)\| \le L_\mu \|\theta_1 \theta_2\|.$

SGD improvements: momentum

There are many improvements of stochastic gradient descent typically used in practise for deep learning; particularly popular is Polyak momentum:

$$\theta^{(k+1)} = \theta^{(k)} + \beta(\theta^{(k)} - \theta^{(k-1)}) - \alpha \cdot \operatorname{grad}_{\theta} \mathcal{L}\left(\theta^{(k)}\right)$$

or Nesterov's accelerated gradient:

$$\begin{array}{rcl} \hat{\theta}^k & = & \theta^{(k)} + \beta(\theta^{(k)} - \theta^{(k-1)}) \\ \theta^{(k+1)} & = & \hat{\theta}^{(k)} - \alpha \cdot \mathrm{grad}_{\theta} \mathcal{L}\left(\hat{\theta}^{(k)}\right) \end{array}$$

These acceleration methods give substantial improvements in the linear convergence rate for convex problems; linear convergence rates are: Normal GD $\frac{\kappa-1}{\kappa+1}$, Polyak $\frac{\sqrt{\kappa}-1}{\sqrt{\kappa}+1}$ and NAG $\sqrt{\frac{\sqrt{\kappa}-1}{\sqrt{\kappa}}}$.

SGD improv. : Adaptive sub-gradients (Duchi et al. 11'1)

A standard method for improving the convergence rate of line-search methods is preconditioning, Adaptive sub-gradients (AdaGrad) is such a method.

Let $g^{(k)}(\theta^{(k)}) =: \operatorname{grad}_{\theta} \mathcal{L}(\theta^{(k)})$ be the gradient of the training loss function at iteration k, then an efficient method for preconditioning is via a diagonal matrix B with entries

$$B_k(i,i) = \left(\sum_{j=1}^k \left(g^{(j)}(\theta^{(j)})(i)\right)^2\right)^{1/2}$$

which is the diagonal of the square-root of the sum of prior gradient outer-products. AdaGrad is the gradient descent method

$$\theta^{(k+1)} = \theta^{(k)} - \eta |\Lambda_k|^{-1} (B^{(k)} + \epsilon I)^{-1} \mathsf{grad}_{\theta} \sum_{\mu \in \Lambda_k} I(\theta; \mathsf{x}_{\mu}, \mathsf{y}_{\mu}).$$

 $\epsilon l > 0$ added to avoid poor scaling of small values of $B^{(k)}(i,i)$.

1 http://jmlr.org/papers/volume12/duchi11a/duchi11a.pdf

AdaGrad improvements: RMSProp and AdaDelta

AdaGrad preconditions with the inverse of

$$B_k(i,i) = \left(\sum_{j=1}^k \left(g^{(j)}(\theta^{(j)})(i)\right)^2\right)^{1/2}.$$

RMSProp (Hinton) gives more weight to the current gradient

$$B_k^{RMS}(i,i) = \gamma B_{k-1}^{RMS}(i,i) + (1-\gamma) \left(g^{(k)}(\theta^{(k)})(i)\right)^2$$

for some $\gamma \in [0,1]$ and updates as

$$\theta^{(k+1)} = \theta^{(k)} - \eta |\Lambda_k|^{-1} (B^{(k)} + \epsilon I)^{-1/2} \operatorname{grad}_{\theta} \sum_{\mu \in \Lambda_k} I(\theta; x_{\mu}, y_{\mu}).$$

AdaDelta (Zeiler 12'²) extends AdaGrad using a similar preconditioned as B_k^{RMS} , but also estimates the stepsize using an average difference in $\theta^{(k)} - \theta^{(k-1)}$.

²https://arxiv.org/abs/1212.5701

Adaptive moment estimation (Adam) (Kingma et al. 15'3)

Algorithm 1: Adam, our proposed algorithm for stochastic optimization. See section 2 for details, and for a slightly more efficient (but less clear) order of computation. g_t^2 indicates the elementwise square $g_t \odot g_t$. Good default settings for the tested machine learning problems are $\alpha = 0.001$, $\beta_1 = 0.9, \beta_2 = 0.999$ and $\epsilon = 10^{-8}$. All operations on vectors are element-wise. With β_1^t and β_2^t we denote β_1 and β_2 to the power t.

```
Require: \alpha: Stepsize
Require: \beta_1, \beta_2 \in [0, 1): Exponential decay rates for the moment estimates
Require: f(\theta): Stochastic objective function with parameters \theta
Require: \theta_0: Initial parameter vector
   m_0 \leftarrow 0 (Initialize 1st moment vector)
   v_0 \leftarrow 0 (Initialize 2<sup>nd</sup> moment vector)
   t \leftarrow 0 (Initialize timestep)
   while \theta_t not converged do
       t \leftarrow t + 1
       g_t \leftarrow \nabla_{\theta} f_t(\theta_{t-1}) (Get gradients w.r.t. stochastic objective at timestep t)
       \begin{array}{l} m_t \leftarrow \beta_1 \cdot m_{t-1} + (1-\beta_1) \cdot g_t \text{ (Update biased first moment estimate)} \\ v_t \leftarrow \beta_2 \cdot v_{t-1} + (1-\beta_2) \cdot g_t^2 \text{ (Update biased second raw moment estimate)} \end{array}
       \widehat{m}_t \leftarrow m_t/(1-\beta_1^t) (Compute bias-corrected first moment estimate)
       \hat{v}_t \leftarrow v_t/(1-\beta_2^t) (Compute bias-corrected second raw moment estimate)
       \theta_t \leftarrow \theta_{t-1} - \alpha \cdot \widehat{m}_t / (\sqrt{\widehat{v}_t} + \epsilon) (Update parameters)
   end while
   return \theta_t (Resulting parameters)
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³https://arxiv.org/pdf/1412.6980.pdf

Adaptive moment estimation (Adam) (Kingma et al. 15^{'4})

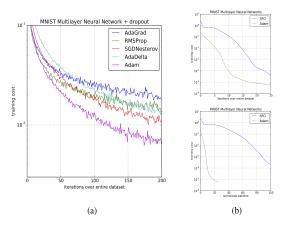


Figure 2: Training of multilayer neural networks on MNIST images. (a) Neural networks using dropout stochastic regularization. (b) Neural networks with deterministic cost function. We compare with the sum-of-functions (SFO) optimizer (Sohl-Dickstein et al., 2014)

⁴https://arxiv.org/pdf/1412.6980.pdf

Adaptive moment estimation (Adam) (Kingma et al. 15'5)

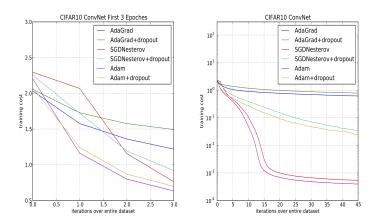


Figure 3: Convolutional neural networks training cost. (left) Training cost for the first three epochs. (right) Training cost over 45 epochs. CIFAR-10 with c64-c64-c128-1000 architecture.

⁵https://arxiv.org/pdf/1412.6980.pdf

AdaGrad: as adaptive stepsize rule (Ward et al. 18'6)

Let $g^{(k)}(\theta^{(k)}) =: \operatorname{grad}_{\theta} \mathcal{L}(\theta^{(k)})$ be the gradient of the training loss function at iteration k, then an efficient method for preconditioning is via a diagonal matrix B with entries

$$B_k(i,i) = \left(\sum_{j=1}^k \left(g^{(j)}(\theta^{(j)})(i)\right)^2\right)^{1/2}$$

which is the diagonal of the square-root of the sum of prior gradient outer-products. AdaGrad is the gradient descent method

$$\theta^{(k+1)} = \theta^{(k)} - \eta |\Lambda_k|^{-1} (B^{(k)} + \epsilon I)^{-1} \mathsf{grad}_{\theta} \sum_{\mu \in \Lambda_k} I(\theta; \mathsf{x}_{\mu}, \mathsf{y}_{\mu}).$$

A simplified version, focusing on the per iteration (as opposed to per index) update is to let $B_k = b_k I$ where $b_{k+1}^2 = b_k^2 + \|g^{(k)}\|_2^2$.

⁶https://arxiv.org/pdf/1806.01811.pdf

AdaGrad: scalar update in batch setting (Ward et al. 18'7)

Given a loss function $\mathcal{L}(\theta; X, Y)$ consider the scalar AdaGrad update algorithm: Initialize with $\theta^{(0)}$ and $b_0 > 0$

$$\begin{array}{rcl} b_k^2 & = & b_{b-1}^2 + \|\mathrm{grad}_{\theta}\mathcal{L}(\theta^{(k)})\|_2^2 \\ \theta^{(k)} & = & \theta^{(k-1)} - b_k^{-1}\mathrm{grad}_{\theta}\mathcal{L}(\theta^{(k)}) \end{array}$$

Consider $\mathcal{L}(\theta) \in C_L^1$, that is L is the smallest value for which $\|\|\operatorname{grad}_{\theta}\mathcal{L}(\theta_1) - \operatorname{grad}_{\theta}\mathcal{L}(\theta_2)\|_2 \le L\|\theta_1 - \theta_2\|$ for all θ_1, θ_2 , then scalar batch AdaGrad satisfies $\min_{k=1,\ldots,T-1}\|\operatorname{grad}_{\theta}\mathcal{L}(\theta^{(k)})\|_2^2 \le \epsilon$ for either

$$T = 1 + \left\lceil 2\epsilon^{-1}\mathcal{L}(\theta^{(0)})(b_0 + 2\mathcal{L}(\theta^{(0)})) \right\rceil \quad \text{if} \quad b_0 \ge L, \quad \text{or}$$

$$= 1 + \left\lceil \epsilon^{-1} \left(L^2 - b_0^2 + 4(\mathcal{L}(\theta^{(0)}) + (3/4 + \log{(L/b_0)})L)^2 \right) \right\rceil$$
if $b_0 < L$. In contrast, if b_k is held as a fixed constant b , then if $b < L/2$ gradient descent can diverge, while if $b \ge L$ then
$$T = 2b\epsilon^{-1}\mathcal{L}(\theta^{(0)}).$$

⁷https://arxiv.org/pdf/1806.01811.pdf

AdaGrad: scalar proof ingredients (Ward et al. 18'8)

The convergence complexity for AdaGrad in the scalar batch setting follows from the following properties for any non-negative values a_1,\ldots,a_T with $a_1>0$, (with a_k taking the place of $\|\mathrm{grad}_{\theta}\mathcal{L}(\theta^{(k)})\|_2^2$)

$$\sum_{\ell=1}^{T} \frac{a_{\ell}}{\sum_{i=1}^{\ell} a_{i}} \leq \log \left(\sum_{i=1}^{T} a_{i} \right) + 1 \quad \text{and} \quad \sum_{\ell=1}^{T} \frac{a_{\ell}}{\sqrt{\sum_{i=1}^{\ell} a_{i}}} \leq 2 \sqrt{\sum_{i=1}^{T} a_{i}}.$$

Also, for any fixed $\epsilon \in (0,1]$ and $L,b_0>0$, the iterates $b_{k+1}^2=b_k^2+a_k$ has the property that after $N=\lceil \epsilon^{-1}(L^2-b_o^2)\rceil+1$ iterations either $\min_{k=0}^{N-1}a_k\leq \epsilon$ or $b_N\geq L$. Lastly, letting k_0 be the first iterate such that $b_{k_0}\geq L$, then for all $k\geq k_0$ $b_k\leq b_{k_0-1}+2\mathcal{L}(\theta^{(k_0-1)})$ (bounded above) and $\mathcal{L}(\theta^{(k_0-1)})\leq \frac{L}{2}\left(1+2\log(b_{k_0-1}/b_0)\right)$ (not diverged).

⁸https://arxiv.org/pdf/1806.01811.pdf

AdaGrad: scalar stochastic (Ward et al. 18'9)

Let $g^{(k)}$ be an unbiased estimator of the gradient $\operatorname{grad}_{\theta}\mathcal{L}(\theta^{(k)})$ of the training loss function at iteration k; that is $\mathbb{E}(g^{(k)}) = \operatorname{grad}_{\theta}\mathcal{L}(\theta^{(k)})$. Moveover, let there be a uniform bound $\mathbb{E}(\|g^{(k)}\|_2^2) \leq c_g^2$. Then consider the stochastic scalar AdaGrad update as

$$b_k^2 = b_{b-1}^2 + ||g^{(k)}||_2^2$$

$$\theta^{(k)} = \theta^{(k-1)} - b_k^{-1} g^{(k)}.$$

Unlike in the batch version of AdaGrad where b_k converges to a fixed stepsize, stochastic AdaGrad converges roughly at the rate $b_k \approx c_g \, k^{1/2}$. Morevover Ward et al. showed that

$$\min_{\ell=0,\dots,N-1} \left(\mathbb{E} \| \mathsf{grad}_{\theta} \mathcal{L}(\theta^{(k)}) \|^{4/3} \right)^{3/2} \leq \mathcal{O} \left(\frac{b_0 + c_g}{\mathsf{N}} + \frac{c_g}{\mathsf{N}^{1/2}} \right) \log(\mathsf{N} c_g^2/b_0^2)$$

⁹https://arxiv.org/pdf/1806.01811.pdf

AdaGrad: batch MNIST (Ward et al. 18'10)

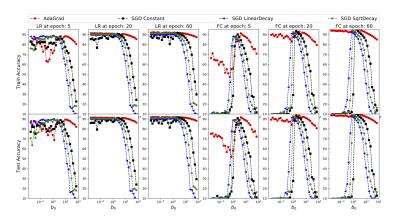


Figure 1: Batch setting on MNIST. Top (bottom) row are plots of train (test) accuracy with respect to the initialization b_0 . The left 6 figures are for logistic regression (LR) with snapshots at epoch 5, 20 and 60 in the 1st, 2nd and 3rd column respectively. The right 6 figures are for two fully connected layers (FC) with snapshots at epoch 5, 20 and 60 in the 4th, 5th and 6th column.

¹⁰https://arxiv.org/pdf/1806.01811.pdf

AdaGrad: stochastic MNIST (Ward et al. 18'11)

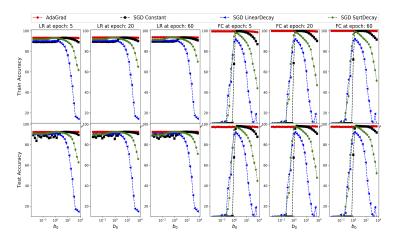


Figure 2: Stochastic setting on MNIST. Left 6 figures by logistic regression and right 6 figures by two fully connected layer. Note that the scale of y-axis change. See Figure 1 for reading instruction.

¹¹https://arxiv.org/pdf/1806.01811.pdf

AdaGrad: stochastic CIFAR10 (Ward et al. 18'12)

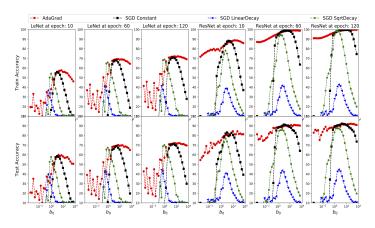


Figure 3: Stochastic setting on CIFAR10. Left 6 figures by LeNet and right 6 figures by ResNet. Note that the epoch (see title) is different from previous figures and no momentum is used. See Figure 1 for reading instruction.

¹²https://arxiv.org/pdf/1806.01811.pdf