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Least Squares Learning

Approximately solve

$$\min_{w \in \mathcal{H}} \mathcal{L}(w), \qquad \mathcal{L}(w) = \frac{1}{2} \mathbb{E}[(Y - \langle w, X \rangle)^2].$$

 \mathcal{H} : real separable Hilbert space

Define

$$\Sigma = \mathbb{E}[X \otimes X], \quad h = \mathbb{E}[XY].$$

Optimal solution w_* satisfies normal equation:

 $\Sigma w_* = h$.

Mini-Batch SGD Recursion

Let $t = 0, ..., T, w_0 = 0$ and

$$w_{t+1} = w_t - \frac{\gamma_t}{b} \sum_{i=b(t-1)+1}^{bt} (\langle w_t, x_{j_i} \rangle - y_{j_i}) x_{j_i},$$

 $j_1, ..., j_{bT} \sim i.i.d Unif[n].$

Tail AveSGD: Tail-length L = 1, ..., T $\bar{w}_L := \frac{1}{I} \sum_{t=T-L+1}^{T} w_t$ Unif AveSGD: L = T

Why Tail-Averaging ? (Part I)



 \odot too small step sizes \Rightarrow slow convergence

 \square larger step sizes \Rightarrow improved convergence + noisy trajectories Unif AveSGD: allows large/ constant step sizes since it reduces the variance of SGD

Tail AveSGD: sufficiently "long" tail preserve this benefit

Assumption I: Regularity

For some $r \geq 0$ we assume $w_* \in \operatorname{Ran}(\Sigma^r)$. Note: $\operatorname{Ran}(\Sigma^0) = \mathcal{H}$ and $\operatorname{Ran}(\Sigma^r) \subseteq \mathcal{H}$

Beating SGD Saturation with Tail-Averaging and Minibatching

Main Theorem: Excess Risk of Tail AveSGD

Define effective dimension

 $\mathcal{N}(1/\gamma L) := \mathrm{Trace}[(\Sigma + 1/(\gamma L))^{-1}\Sigma].$ Let $1 \le L \le T$. Assume $\gamma \kappa^2 < 1/4$. Then

 $\mathbb{E}[\mathcal{L}(\bar{w}_{L}) - \mathcal{L}(w_{*})] \lesssim \operatorname{Approx}_{L}(\Sigma, w_{*}) + \frac{1}{2}$ for n sufficiently large.

Saturation

Let Approx_I (Σ, w_*) denote the Approximation Error. Unif AveSGD: Approx_T(Σ , w_*) $\approx (1/\gamma T)^{2\min(r,1/2)+1}$

Tail AveSGD: Approx_L(Σ, w_*) $\approx (1/\gamma L)^{2r+1}$



Unif AveSGD

Assumption II: Capacity

For some $\nu \in (0, 1]$ we assume $\mathcal{N}(1/\gamma L) \leq (\gamma L)^{\nu}$.

Corollary: Learning Rate

The excess risk of the (tail)-averaged SGD iterate satisfies

 $\mathbb{E}[\mathcal{L}(\bar{w}_{L}) - \mathcal{L}(w_{*})] \lesssim n^{-\frac{2r+1}{2r+1+\nu}}$ for each of the following choices:

one pass: $b_n \simeq 1$, $L_n \simeq n$, $\gamma_n \simeq n^{-\frac{2r+\nu}{2r+1+\nu}}$ one pass: $b_n \simeq n^{\frac{2r+\nu}{2r+1+\nu}}$, $L_n \simeq n^{\frac{1}{2r+1+\nu}}$, $\gamma_n \simeq 1$ $\bigcirc \bigcirc \left(n^{\frac{1}{2r+1+\nu}}\right) \text{ passes: } b_n \simeq n, \ L_n \simeq n^{\frac{1}{2r+1+\nu}}, \ \gamma_n \simeq 1.$

$$\frac{\mathcal{N}(1/\gamma L)}{n} + \frac{\gamma \operatorname{Trace}[\Sigma^{\nu}]}{b(\gamma L)^{1-\nu}}$$

Tail AveSGD

Experiment: Saturation



Excess risk as a function of regularity r with uniform and tail averaging.

Unif AveSGD: starts to lag behind its tail-averaged counterpart for larger values of r exceeding 1/2, flattening out.

Tail AveSGD: continues to improve for large values of r, confirming that this algorithm can indeed massively benefit from favorable structural properties of the data.



Unif AveSGD

Single pass performance as a function of the stepsize γ and the minibatch-size b.

Performance: remains largely constant as $\gamma \cdot b$ remains constant for both algorithms, until a critical threshold stepsize is reached.

Tail AveSGD: permits the use of larger minibatch sizes, allowing for more efficient parallelization.

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Tail AveSGD