Optimal Statistical Rates for Decentralised Non-Parametric Regression with Linear Speed-Up

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- Desiderata for Distributed Machine Learning —

Suppose n agents solving a machine learning problem. Properties of an ideal distributed algorithm:

Statistics: retain optimal statistical precision

Runtime: speed-up over single-agent due to parallel computing: ideally factor n

Communication: fixed cost per agent per step, ideally independent of n

Consensus Optimisation for Decentralised Learning -

Network of agents G = (V, E), each with data points sampled i.i.d $(x_{i,v}, y_{i,v})$ perform linear regression. Therefore each agent wishes to minimise with coefficient $\boldsymbol{\omega} \in \mathbb{R}^d$

$$F(\boldsymbol{\omega}) = \frac{1}{\mathbf{n}} \sum_{\mathbf{v} \in \mathbf{V}} \frac{1}{\mathbf{m}} \sum_{\mathbf{i} \in [\mathbf{m}]} (\langle \boldsymbol{\omega}, x_{i,v} \rangle - y_{i,v})^2 = \underbrace{\frac{1}{\mathbf{n}} \sum_{\mathbf{v} \in \mathbf{V}} F_v(\boldsymbol{\omega})}_{\mathbf{v} \in \mathbf{V}},$$
Empirical Loss for Agent v
Consensus Optimisation

where F_v is function held by agent v.

Consensus Optimisation

Suppose F_v arbitrary and each agent wants to minimise $\frac{1}{n} \sum_{\mathbf{v} \in \mathbf{V}} F_v$.

- \rightarrow Agents alternate: local gradient descent steps on F_v and local averaging on network [4]
 - ✓ Low communication cost per agent for sparse graphs
- Robust/Decentralised as no single node responsible for disseminating information
- Performance depends on network topology [2, 5].

Graphs with smaller spectral gap benefit less from decentralisation

In machine learning F_v are often not arbitrary

In our case, concentration states for $v \in V$

$$F_{v}(\boldsymbol{\omega}) = \frac{1}{\mathbf{m}} \sum_{\mathbf{i} \in [\mathbf{m}]} (\langle \boldsymbol{\omega}, x_{i,v} \rangle - y_{i,v})^{2} \xrightarrow{\mathbf{m} \to \infty} \underbrace{\int_{X \times Y} (\langle \boldsymbol{\omega}, x \rangle - y)^{2} d\rho(x,y)}_{\text{Test Risk}} =: \mathcal{E}(\boldsymbol{\omega})$$

where ρ is the distribution of the data points.

 \mathbb{V} In large data scenario, all functions are converging to the same quantity $\to \mathrm{Test} \, \mathrm{Risk}$.

Main Question:

Can concentration speed-up consensus learning for any network topology?

Non-parametric Statistical Assumptions -

Use tools from non-parametric regression [1].

Predictor minimising Test Risk over set of linear predictors $x \to \langle \omega, \cdot \rangle$ denoted f_H .

Noise: Exists $M \in (0, \infty)$, $\nu \in (1, \infty)$ so for any $\ell \in \mathbb{N}$ we have $\int_V y^{2\ell} d\rho(x|y) \leq \nu \ell! M^{\ell}$.

Difficulty of estimation problem:

For $f \in L^2(H, \rho_X)$ let $\mathcal{L}_{\rho}(f) = \int_X \langle x \cdot \rangle f(x) d\rho_X(x)$. Exists r > 0 such that $\|\mathcal{L}_{\rho}^{-r} f_H\| < \infty$. **Spectrum of covariance operator:**

Exists $\gamma \in (0,1], c_{\gamma} > 0$ such that $\text{Tr}(\mathcal{L}_{\rho}(\mathcal{L}_{\rho} + \lambda)^{-1}) \leq c_{\gamma}\lambda^{-\gamma}$ for all $\lambda > 0$.

Distributed Gradient Descent

Consider a simple consensus optimisation algorithm [4]:

Communication Matrix: $\mathbf{P} \in \mathbb{R}^{n \times n}$ symmetric doubly stochastic matrix support on the graph

$$\mathbf{P} = \mathbf{P}^{\top}$$
, $\mathbf{P}\mathbf{1} = \mathbf{1}$ and for $v \neq w$ $\underbrace{\mathbf{P}_{\mathbf{v}\mathbf{w}} \neq \mathbf{0} \text{ only if } (\mathbf{v}, \mathbf{w}) \in \mathbf{E}}_{\mathbf{Sparsity pattern matches network}}$

Network Dependence: Let σ_2 be second largest eigenvalue in magnitude for **P**.

Scaling $O((1 - \sigma_2)^{-1}) = n^2$ (Cycle), n (Grid/Random Geo.), and 1 (Complete/Expander). **Algorithm**: Initalised $w_{1,v} = 0$ for $v \in V$, iterates updated for all $v \in V$

$$\boldsymbol{\omega_{t+1,v}} = \underbrace{\sum_{\mathbf{w} \in \mathbf{V}} \mathbf{P_{vw}}}_{\text{Local Communication}} \left(\boldsymbol{\omega_{t,w}} - \eta \frac{1}{m} \sum_{i=1}^{m} (\langle \boldsymbol{\omega_{t,w}}, x_{i,w} \rangle_{H} - y_{i,w}) x_{i,w} \right)$$
Local Gradient Descent

Implict Regularisation:

Model Complexity controlled through early stopping (Iterations t) and step size η .

Optimal Statistical Rate with Implicit Regularisation –

Let aforementioned assumptions hold with $r \ge 1/2$ and $2r + \gamma \ge 2$. Fix

$$t = \underbrace{(\mathbf{nm})^{\mathbf{1}/(\mathbf{2r} + \gamma)}}_{\mathbf{Single-Machine Iterations}} \times \begin{cases} \left(\frac{(nm)^{2r/(2r + \gamma)}}{\mathbf{m}(1 - \sigma_2)^{\gamma}}\right)^{1/\gamma} \vee 1 & \text{if } m \geq n^{2r/\gamma} \\ \frac{(nm)^{r/(2r + \gamma)}}{\sqrt{\mathbf{m}}(1 - \sigma_2)} & \text{otherwise} \end{cases}$$

 $\eta = \frac{\kappa^{-2}(nm)^{1/(2r+\gamma)}}{t}$ and let $m \ge n^{\frac{2r+2+\gamma}{2r+\gamma-2}}$ then $\forall v \in V$:

 $\mathbf{E}[\mathcal{E}(\boldsymbol{\omega_{t+1,v}})] - \inf_{\omega} \mathcal{E}(\omega) \lesssim (\mathbf{nm})^{-2\mathbf{r}/(2\mathbf{r}+\gamma)}$

Time Model

Gradient computation costs 1 unit of time.

Communication delay costs τ units of time, for some $\tau > 0$.

Distributed time per iteration =

Local Gradient Computation Communicating/Aggreagating Neighbours Information

Speed-Up defined as

$$Speed-Up = \frac{Single \ Machine \ Run \ time}{Distributed \ Run \ time} = \frac{Single \ Machine \ Iterations}{Distributed \ Iterations} \times \underbrace{\frac{nm}{m+\tau+Deg(P)}}_{Ratio \ of \ Time \ Per \ Iteration}$$

— Main Result: Speed-Up with Consensus Methods Utilising Concentration —

Iterations required decreasing in number of samples m up to a point

$$\underbrace{\mathbf{m} \geq \frac{n^{2r/\gamma}}{(\mathbf{1} - \sigma_{\mathbf{2}})^{2\mathbf{r} + \gamma}} \vee n^{\frac{2r + 2 + \gamma}{2r + \gamma - 2}}}_{\text{Sufficiently Many Samples}} \implies \text{Distributed Iterations} = \underbrace{\mathbf{Single Machine Iterations}}_{\text{Speed-Up}}$$

$$\Longrightarrow \text{Speed-Up}^{\tau + \text{Deg}(P) = O(m)} O(n)$$

therefore linear speed-up for any network topology.

	Cycle	Grid	R. Geom.	Complete	Expander
Speed-Up	O(n)	O(n)	O(n)	O(n)	O(n)
Communication	O(1)	O(1)	O(1)	O(n)	O(1)

Speed-Up with Single-Step Consensus Methods -

Single-Step methods typically require iterations to scale with inverse spectral gap e.g. [2, 5]

Distributed Iterations = Single Machine Iterations
$$\times (\mathbf{1} - \sigma_{\mathbf{2}})$$

$$\downarrow \downarrow$$
Speed-Up = $\frac{nm}{m + \tau + \mathrm{Deg}(P)} (\mathbf{1} - \sigma_{\mathbf{2}})^{\tau + \mathrm{Deg}(P) = O(m)} O(n(\mathbf{1} - \sigma_{\mathbf{2}}))$

therefore linear speed-up restricted to well connected topologies.

	Cycle	Grid	R. Geom.	Complete	Expander
Speed-Up	O(1/n)	O(1)	O(1)	O(n)	O(n)
Communication	O(1)	O(1)	O(1)	O(n)	O(1)

Speed-Up with Multi-Step Consensus Methods —

Multi-Step methods perform multiple communication steps between gradient descent steps [5],

Distributed Iterations = Single Machine Iterations

But communication rounds scales with inverse spectral gap.

	Cycle	Grid	R. Geom.	Complete	Expander
Speed-Up	O(n)	O(n)	O(n)	O(n)	O(n)
Communication	O(n)	$O(\sqrt{n})$	$O(\sqrt{n})$	O(n)	O(1)

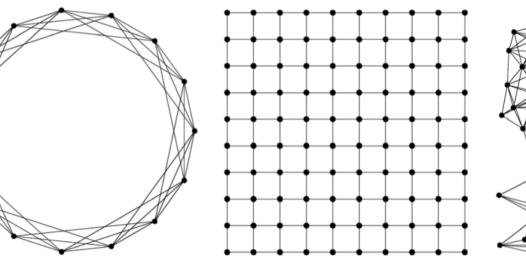
Therefore trade off between **Speed-Up** and **Communication Cost**.

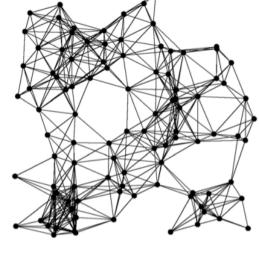


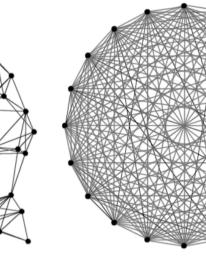


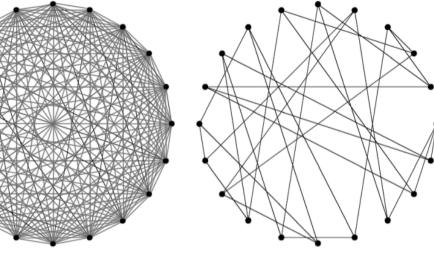












Detailed Error Decomposition

The Test Error is decomposed as follows when $m \ge n^{2r/\gamma}$

$$\textbf{Test Error} \lesssim \underbrace{(\eta t)^{-2r}}_{\textbf{Bias}} + \underbrace{\frac{(\eta t/(nm)^{1/(2r+\gamma)})^2}{(nm)^{2r/(2r+\gamma)}}}_{\textbf{Sample Variance}} + \underbrace{\frac{\eta^{\gamma}}{\textbf{m}(\textbf{1}-\sigma_{\textbf{2}})^{\gamma}}}_{\textbf{Population Network Error}} + \underbrace{\frac{(\eta t)^{\gamma+2}}{m^2}}_{\textbf{Residual Network Error}}$$

Bias and **Sample Variance** align with Gradient Descent with nm samples [3]

 \implies Fix $\eta t = (nm)^{1/(2r+\gamma)}$.

Population and **Residual Network Error**: arise due from averaging steps with the matrix **P**.

Population Network Error ———— Residual Network Error ———

Follows standard network term [2]

Decreasing with step size η Depends on spectral gap $(1 - \sigma_2)$

Due to concentration, decreasing with m

Small by picking $\eta = \left(\frac{\mathbf{m}(\mathbf{1} - \sigma_{\mathbf{2}})^{\gamma}}{(nm)^{2r/(2r+\gamma)}}\right)^{1/\gamma} \vee 1$.

Higher order term

From empirical covariance multiplying the iterates at each iteration

Utilise concentration and contraction to

 \implies require $m \ge n^{\frac{2r+2+\gamma}{2r+\gamma-2}}$.

- Proof Sketch: Population Network Error —

Let \mathcal{T}_{ρ} be conjugate of \mathcal{L}_{ρ} and for $k \geq 1$, let $\mathbf{N}_{\mathbf{k}}$ be r.v. with zero mean concentrating to zero such that $\|(\mathcal{T}_{\rho} + \lambda \mathbf{I})^{-1/2} \mathbf{N}_{\mathbf{k}}\| \lesssim 1/\sqrt{\lambda^{\gamma} \mathbf{m}}$ w.h.p. Then, with mixing time $t^* \simeq (1 - \sigma_2)^{-1}$

$$\begin{split} \mathbf{E}[(\mathbf{Pop.\ Net.\ Error})] &\leq \mathbf{E}\Big[\Big(\sum_{k=1}^{t} \sigma_{\mathbf{2}}^{\mathbf{t}-\mathbf{k}+\mathbf{1}} \eta \| \mathcal{T}_{\rho}^{1/2} (I - \eta \mathcal{T}_{\rho})^{t-k} \mathbf{N}_{\mathbf{k}} \|\Big)^{2}\Big] \\ &\lesssim \frac{1}{\mathbf{m} \lambda^{\gamma}} \Big(\sum_{k=t-\mathbf{t}^{\star}}^{t} \sigma_{\mathbf{2}}^{\mathbf{t}-\mathbf{k}+\mathbf{1}} \eta \underbrace{\|\mathcal{T}_{\rho}^{1/2} (I - \eta \mathcal{T}_{\rho})^{t-k} (\mathcal{T}_{\rho} + \lambda I)^{1/2}\|}_{\mathbf{Contraction}} \Big)^{2} \lesssim \frac{\log(t^{\star}) + \lambda \eta \mathbf{t}^{\star}}{\mathbf{m} \lambda^{\gamma}} \\ &\overset{\mathbf{t}^{\star} \ \mathbf{Poorly\ Mixed\ terms}}{} &\overset{\mathbf{t}^{\star} \ \mathbf{Poorly\ Mixed\ terms}}{} \end{split}$$

Optimised by picking $\lambda = (\eta \mathbf{t}^*)^{-1}$.

Future Work -

- Non-parametric setting: extending analysis to non-attainable case $r \le 1/2$ and tighter analysis on Residual Network Error.
- General Loss Function: Squared loss yields bias/variance decomposition, concentration likely hold for more general losses.
- Statistics/Communication trade off with sparse/randomised gossip: linear speed-up independent of topology \rightarrow agents randomly gossip at each iteration.
- Stochastic Gradient Descent and mini-batches: random subset of data at each iteration [3].

References -

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