The Power-Oja Method for Decentralized Subspace Estimation/Tracking

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Subspace Estimation

New Challenges

- Tracking and estimation accuracy.
- Distributed system: separate antennas.
- Large scale: massive array.
- Complexity: distribute computations to local processors.



Figure: Two examples: WiFi networks and Radar networks.

Literature Survey

- ▶ The power method [Gv96]:
 - A batch processing method with fast convergence.
 - Non-adaptive, high latency
 - Guarantee for a rank-*p* subspace.
 - Computations can be decentralized [LSM11].
- The Oja's method [OK85]
 - A stochastic gradient decent (SGD) method adaptive for tracking varying statistics.
 - First order method: suffer from slow convergence.
 - unconstrained SGD, no guarantee for a rank-p subspace.
 - Computations can be decentralized [SPK08, SPZ16].
- The key of the decentralization is average consensus [LSM11, SPK08, SPZ16, BDF13]
 - Data are often measured distributively over large networks.
 - Gossip-based consensus algorithms solve multi-agent coordination and optimization problems in a decentralized manner.
 - Their key features are
 - $\checkmark\,$ built-in fault tolerance to intermittent computation/communication.
 - $\checkmark\,$ self reorganization to automatic failure correction.

Problem Statement

We consider a non-stationary stochastic process r(t) ∈ C^N and let T ⊂ {1, 2, ...} be a sampling set. Define the sampled covariance:

$$\hat{R}(\mathcal{T}) := |\mathcal{T}|^{-1} \sum_{s \in \mathcal{T}} \mathbf{r}(s) \mathbf{r}^{H}(s) .$$
(1)

▶ We track top *p*-D subspace by tackling the non-convex, stochastic optimization:

$$\min_{\boldsymbol{U}\in\mathcal{C}^{N\times p}} f_t(\boldsymbol{U}) := \mathbb{E}\left[\|\boldsymbol{r}(t) - \boldsymbol{U}\boldsymbol{U}^{H}\boldsymbol{r}(t)\|^2 \right], \ \forall \ t \ge 1.$$
(2)

• We follow the stochastic approximation to the objective function f(U):

$$\hat{f}(\boldsymbol{U};\mathcal{T}_{\tau}) := \operatorname{Tr}\left(\left(\boldsymbol{U}\boldsymbol{U}^{H}\boldsymbol{U}\boldsymbol{U}^{H}-2\boldsymbol{U}\boldsymbol{U}^{H}\right)\hat{\boldsymbol{R}}(\mathcal{T}_{\tau})\right), \qquad (3)$$

where $\mathcal{T}_{\tau} \subset \{1, 2, ...\}$ is the set of observations made during the τ th batch.

- If $\mathbf{r}(t)$ is stationary for all $t \in \mathcal{T}_{\tau}$, then $\mathbb{E}[\hat{f}(\mathbf{U}; \mathcal{T}_{\tau})] = f_t(\mathbf{U})$.
- When $|\mathcal{T}_{\tau}|$ is large, $\hat{f}(\boldsymbol{U}; \mathcal{T}_{\tau})$ is a good approximation for $f_t(\boldsymbol{U})$.
- No unitary constraint on the subspace U, no guarantee for a rank-p subspace.

Review the Power Method (PM)

The PM works with a whole batch of samples in \mathcal{T}_{τ} .

- **Step 1:** Generate a random vector as an initial point $\tilde{u}^k(1,\tau)$
- ▶ Step 2: For *k* = 1, ..., *p*

$$\begin{split} \tilde{\boldsymbol{u}}^{k}(\ell+1,\tau) &= \hat{\boldsymbol{R}}(\mathcal{T}_{\tau}) \; \tilde{\boldsymbol{u}}^{k}(\ell,\tau) - \sum_{j=1}^{k-1} (\hat{\boldsymbol{u}}^{j}(\tau))^{H} \left(\hat{\boldsymbol{R}}(\mathcal{T}_{\tau}) \; \tilde{\boldsymbol{u}}^{k}(\ell,\tau) \right) \hat{\boldsymbol{u}}^{j}(\tau), \forall \ell = 1, ..., L \\ \hat{\boldsymbol{u}}^{k}(\tau) &:= \tilde{\boldsymbol{u}}^{k}(L,\tau) / \| \tilde{\boldsymbol{u}}^{k}(L,\tau) \| \; . \end{split}$$

▶ Step 3: Output the top-*p* subspace: $\hat{U}_{PM}(\tau) := [\hat{u}^1(\tau) \ \hat{u}^2(\tau) \ \dots \ \hat{u}^p(\tau)].$

We use \hat{U}_0 to initialize the subspace and denote the above power process by

$$\overline{\boldsymbol{U}}_{PM}(\tau) = \mathsf{PM}\big(\{\boldsymbol{r}(\boldsymbol{s})\}_{\boldsymbol{s}\in\mathcal{T}_{\tau}};\,\hat{\boldsymbol{U}}_{0};L\big)\;,\tag{4}$$

Review the Oja's Learning Rule

The Oja's learning rule works with one sample of r(t) at a time.

▶ Let $\hat{U}_{Oja}(t) \in C^{N \times p}$ be an estimate of U(t) at iteration t, we perform the updates:

$$\hat{\boldsymbol{U}}_{Oja}(t+1) = \hat{\boldsymbol{U}}_{Oja}(t) - \gamma_t \nabla \hat{f}(\hat{\boldsymbol{U}}_{Oja}(t); \{t\}) , \qquad (5)$$

with $\nabla \hat{f}(\hat{U}(t), \{t\}) = -2\mathbf{r}(t)\mathbf{r}^{H}(t)\hat{U}(t) + \mathbf{r}(t)\mathbf{r}^{H}(t)\hat{U}(t)\hat{U}^{H}(t)\hat{U}(t) + \hat{U}(t)\hat{U}^{H}(t)\mathbf{r}(t)\mathbf{r}^{H}(t)\hat{U}(t)$.

- Convergence for stationary r(t):
 - When p = 1 and $\gamma_t = c/t$, at a sub-linear rate of O(1/t) [BDF13];
 - If $\sum_t \gamma_t = \infty$, $\sum_t \gamma_t^2 < \infty$, converges almost surely to the principal *p*-dimensional subspace, yet the convergence rate is not given.
- In practice, the Oja's learning rule is often used for non-stationary r(t) with γ_t .

Motivations for the Power-Oja (P-Oja) Method

Observations for PM and Oja:

- For PM, when $\mathbf{r}(t)$ is non-stationary and $|\mathcal{T}_{\tau}| \ll \infty \rightarrow$ a poor approximation for $\hat{\mathbf{R}}(\mathcal{T}_{\tau})$ to the true covariance \rightarrow degraded performance.
- ▶ For Oja, the spectral gap,

$$\sigma_p(\hat{\boldsymbol{R}}(\mathcal{T})) - \sigma_{p+1}(\hat{\boldsymbol{R}}(\mathcal{T}))$$

is an important factor in determining the convergence speed [BDF13].

Our motivations:

- ▶ We want both advantages of the two methods:tracking and estimation accuracy.
- Try to increase the spectral gap.

How to Increase the Spectral Gap

Our approach:

Modify the stochastic approximation of the objective function:

$$\hat{f}_{POja}(\boldsymbol{U};\mathcal{T}) = \operatorname{Tr}\left(\left(\boldsymbol{U}\boldsymbol{U}^{H}\boldsymbol{U}\boldsymbol{U}^{H} - 2\boldsymbol{U}\boldsymbol{U}^{H}\right)\left(\hat{\boldsymbol{R}}(\mathcal{T})\right)^{L}\right).$$
(6)

Apparently, $(\hat{R}(\mathcal{T}))^L$ has a better spectral gap than $\hat{R}(\mathcal{T})$, i.e.,

$$\sigma_{
ho}((\hat{\pmb{R}}(\mathcal{T}))^L) - \sigma_{
ho+1}((\hat{\pmb{R}}(\mathcal{T}))^L) > \sigma_{
ho}(\hat{\pmb{R}}(\mathcal{T})) - \sigma_{
ho+1}(\hat{\pmb{R}}(\mathcal{T})).$$

P-Oja tracks the subspace in a batch by batch manner:

For the \(\tau\) th batch, we have

$$\begin{split} \nabla \hat{f}_{\mathsf{POja}}(\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau);\mathcal{T}_{\tau}) &= -2(\hat{\boldsymbol{R}}(\mathcal{T}_{\tau}))^{L}\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau) + (\hat{\boldsymbol{R}}(\mathcal{T}_{\tau}))^{L}\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau)\hat{\boldsymbol{U}}_{\mathsf{POja}}^{H}(\tau)\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau)\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau)\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau)\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau)\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau). \end{split}$$

• $(\hat{R}(\mathcal{T}_{\tau}))^{L} \hat{U}_{POja}(\tau)$: performing *L* rounds of the power iterations on $\hat{U}_{POja}(\tau)$ and we can approximately calculate it by $\hat{U}_{PM}(\tau) \approx PM(\{r(s)\}_{s \in \mathcal{T}_{\tau}}; \hat{U}_{POja}(\tau); L)$.

The Power-Oja (P-Oja) Method

Finally, the P-Oja method is given by the following iterations:

$$\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau+1) = \hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau) - \gamma_{\tau} \hat{\nabla} \hat{f}_{\mathsf{POja}}(\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau); \mathcal{T}_{\tau}), \qquad (7)$$

where $\hat{\nabla} \hat{f}_{POja}$ is the approximated gradient, evaluated as:

$$\begin{split} \hat{\nabla} \hat{f}_{\mathsf{POja}}(\hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau);\mathcal{T}_{\tau}) &= \overline{\boldsymbol{U}}_{\mathsf{PM}}(\tau) \hat{\boldsymbol{U}}_{\mathsf{POja}}^{\mathsf{H}}(\tau) \hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau) \\ &+ \hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau) \hat{\boldsymbol{U}}_{\mathsf{POja}}^{\mathsf{H}}(\tau) \overline{\boldsymbol{U}}_{\mathsf{PM}}(\tau) - 2 \overline{\boldsymbol{U}}_{\mathsf{PM}}(\tau) \end{split}$$

where $\hat{\boldsymbol{U}}_{PM}(\tau) \approx \mathsf{PM}(\{\boldsymbol{r}(s)\}_{s \in \mathcal{T}_{\tau}}; \hat{\boldsymbol{U}}_{\mathsf{POja}}(\tau); L)$.

Some Remarks for Power-Oja

- The P-Oja method is parametrized by L and T
 - T: controls the variance in the sampled covariance $\hat{R}(\mathcal{T}_{\tau})$;
 - L: the acceleration given by the power method subroutine.
- ▶ P-Oja reduces into the Oja's learning rule when p = 1, T = 1, L = 1.
- If the samples in the batch is sufficient, we are very likely to obtain a rank-p subspace. Recall that no guarantee for a rank-p subspace for Oja.

Decentralization

Preliminaries:

- ▶ We denote the communication network between *M* processor units as an undirected graph G = (V, E) such that $V = \{1, ..., M\}$ and $E \subseteq V \times V$.
- The graph is assumed to be sparse and connected.
- ▶ A doubly stochastic matrix W associated with G, s.t. $[W]_{ij} = 0$ iff $(i, j) \notin E$.
- ► Each processor unit locally processes its subarray's sampling data, and meanwhile exchanges information with its neighbors in *G*.



Figure: Two examples: WiFi networks and Radar networks.

Average Consesus

Stored and computed in processor unit i:

$$\mathbf{r}_i(t) \in \mathbb{C}^{\frac{N}{M}} = [\mathbf{r}_1(t); ...; \mathbf{r}_M(t)]$$

$$\hat{\boldsymbol{u}}_{i}(\ell,\tau) \in \mathbb{C}^{\frac{N}{M}} = [\hat{\boldsymbol{u}}_{1}(\ell,\tau);...;\hat{\boldsymbol{u}}_{M}(\ell,\tau)]$$

•
$$z_i^0 := \boldsymbol{r}_i^H(t) \hat{\boldsymbol{u}}_i(\ell, \tau)$$
 and z_i^0

The power iteration can be expressed as:

$$\sum_{s \in \mathcal{T}_{\tau}} \mathbf{r}(s) \underbrace{\mathbf{r}^{H}(s) \hat{\mathbf{u}}(\ell, \tau)}_{\sum_{i=1}^{\text{centralized:}} Z_{i}^{0}} = \sum_{s \in \mathcal{T}_{\tau}} \mathbf{r}(s) \underbrace{\mathsf{AC}^{i}(\{z_{j}^{0}\}_{j=1}^{M}; K)}_{\substack{\text{decentralized:} \\ z_{i}^{k+1} = \sum_{j=1}^{M} W_{ij} z_{i}^{k}}}$$

with $\|\boldsymbol{z}^{\kappa} - \left(\sum_{i=1}^{M} z_{i}^{0}/M\right) \mathbf{1}\| \leq |\lambda_{2}(\boldsymbol{W})|^{\kappa} \|\boldsymbol{z}_{0} - \left(\sum_{i=1}^{M} z_{i}^{0}/M\right) \mathbf{1}\|$,

- The convergence rate depends on λ_2 (**W**) [DKM⁺10].
- $\lim_{K\to\infty} \sum_{j=1}^{M} z_j^0 = M \cdot AC^i(\{z_j^0\}_{j=1}^M; K)$ at a geometric rate in K [DKM⁺10].

An Illustration of Decentralized Power-Oja

Power Method

$$PM^{i}\left(\{r(s)\}_{s\in\mathcal{T}_{\tau}}, \hat{U}_{POja}(\tau), W, K, L\right) = \sum_{s\in\mathcal{T}_{\tau}} r(s)AC^{i}(\{r_{i}^{H}(t)\hat{u}_{i}(\ell,\tau)\}_{j=1}^{M}; K)$$
Oja's Method

$$\overline{U}_{PM}^{i}(\tau)$$

$$\hat{\nabla}\hat{f}_{POja}^{i}(\hat{U}_{POja}(\tau); \mathcal{T}_{\tau}) = \overline{U}_{PM}^{i}(\tau)(\hat{U}_{POja}^{H}(\tau)\hat{U}_{POja}(\tau))$$

$$+ \hat{U}_{POja}^{i}(\tau)(\hat{U}_{POja}^{H}(\tau)\overline{U}_{PM}(\tau)) - 2\overline{U}_{PM}^{i}(\tau)$$

$$\hat{U}_{POja}^{i}(\tau+1) = \hat{U}_{POja}^{i}(\tau) - \gamma_{\tau}\hat{\nabla}\hat{f}_{POja}^{i}(\hat{U}_{POja}(\tau); \mathcal{T}_{\tau})$$

$$\hat{U}_{POja}^{i}(\tau+1)$$

Key technique in Oja: $\mathbf{A}^{H}\mathbf{B} = \sum_{i=1}^{M} \mathbf{V}_{i} = \sum_{i=1}^{M} \underbrace{(\mathbf{A}^{i})^{H}\mathbf{B}^{i}}_{\mathbf{P} \times \mathbf{P}}$.

Some Remarks for Decentralized Power-Oja

- The message exchanged in the approximate gradient is a $p \times p$ matrix.
- It is crucial for us to choose a proper network topology.

$$\|\boldsymbol{z}^{K}-\bar{\boldsymbol{z}}\boldsymbol{1}\|\leq |\lambda_{2}(\boldsymbol{W})|^{K}\|\boldsymbol{z}_{0}-\bar{\boldsymbol{z}}\boldsymbol{1}\|.$$

- It is more economical to connect the nearby units with a higher probability while the far-apart units with a lower probability.
- ▶ Example: small-world graph with optimal constant weights [XB04]:

$$W = I - rac{2}{\lambda_1(L) + \lambda_{N-1}(L)}L$$

where \boldsymbol{L} is the Laplacian matrix.

Numerical Simulations: parameter settings

- ▶ A massive array with N = 256 antennas grouped to M = 64 subarrays, each equipped with four antennas.
- T = 1500, SNR= 20dB, the power iteration is L = 20.
- ▶ Degree-6 small-world graph with rewiring probability 0.2.
- ▶ $\gamma_t = 5 \times 10^{-4}$ for the Oja's learning rule, $\gamma_t = 0.01t$ with P-Oja for stationary signals and $\gamma_t = 0.04t$ for non-stationary signals.



Figure: Grouping antennas into subarrays with distributed processors for spectrum sensing.

Numerical Simulations: for stationary signals



Figure: The normalized objective value for a constant 2-D signal space.

- ▶ The convergence rate increases as *T* increases.
- ▶ The decentralized performance will approach the centralized one as K increases.
- The P-Oja method converges much faster than the Oja's method, and the decentralized algorithms work well under the chosen graph.

Numerical Simulations: for non-stationary signals



Figure: The normalized objective value for a variant 1-D signal space. The legend is the same as that in Fig. 1, except for the number of gossip iterations is now K = 10. The diamond-marked curve is the NOV for the conventional power method.

- ► As *T* increases, the convergence rate increases.
- The decentralized and centralized methods coincide with each other when K = 10.
- The power method cannot track the change of the covariance.

Conclusions

- ▶ We propose P-Oja method to integrate the Oja's learning rule and power method.
- It exhibits both tracking ability and estimation accuracy.
- ► All the computations are distributed into individual processor units.
- Our simulation results demonstrate that the proposed P-Oja can both track the change of statistic, but converges much faster than the conventional Oja method.

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Thank You

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Question Welcomed!