Signal Processing on Expanding Graphs

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Graphs and graph signals

Graphs (Networks)

- ▶ Captures pair-wise and higher order relationships between entities
- Examples: Social networks, sensor networks, citation networks
- ▶ Irregular domain, unlike time, grid.

Graph Signals

- ▶ Data associated with a graph
- ▶ Node level data: User opinions, temperature readings, paper label
- ▶ Different way to interpret data w.r.t an underlying topology



Graph signal processing

Processing data w.r.t the underlying graph topology for a specific task.

- ▶ Filtering: Topology aware filtering out certain variations in the signal [1]
- ▶ Interpolation: Filling in missing values at certain nodes from other node values [2]
- Classification: Assign a node to a class based on graph signals [3]
- ▶ Topology Identification: Estimate the graph structure (Edges) from graph signals [4]



Rating prediction in Recommender System

- ▶ Nodes: Users in a recommender system with known ratings
- ▶ Graph built from rating similarities between the users
- ▶ Graph signal: All user ratings for a movie
- Task: Predict unobserved ratings (Interpolation) [5], predicting associations between users and items (topology identification) [6]



Expanding graphs



- Example: New users joining a recommender system
- ▶ Received attention from network science via well-known growth models [7]
- ▶ In GSP, majority of works focus on graphs (static and dynamic) with fixed number of nodes [8, 9]
- ▶ Lacking a principled approach from a signal processing POV



Expanding graphs: Challenges

- **Dependence on connectivity**
 - \blacktriangleright Often the connections of new nodes are not known
 - ▶ Inability to use GSP tools to perform rating prediction (Cold start users)



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► Streaming nature of data

- No Batch type data
- \blacktriangleright Need to efficiently re-train as new users arrive over time



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Non-stationarity

- \blacktriangleright New nodes may represent different entities with different distribution of data
- ▶ Signal processing algorithms to adapt to the preferences of different users with different tastes



Overview of this talk

▶ Task aware GSP with filters under uncertainty

 \blacktriangleright Online graph filter design

▶ Online topology identification



Commonly used notation

- ▶ $\mathcal{G} = {\mathcal{V}, \mathcal{E}}$: Graph with node set \mathcal{V} and edge set \mathcal{E}
- ▶ $\mathbf{S} \in \mathbb{R}^{N \times N}$: Shift operator of an N node graph with $S_{ij} \neq 0$ if an edge exists between node i and j; $S_{ij} = 0$ otherwise
- $\blacktriangleright~\mathbf{x} \in \mathbb{R}^{N \times 1}$: Graph signal
- \blacktriangleright A: Adjacency matrix, L: Graph Laplacian matrix
- \blacktriangleright Graphs are undirected without self-loops, unless specified otherwise



Task aware GSP with filters under uncertainty



Graph filters

- ▶ Flexible, parametric, and localized operator for processing signals on graphs.
- \blacktriangleright Sx: Shifts x. each node weighs signals from immediate neighbours
- $\blacktriangleright {\bf S}^k {\bf x}_k$ shifts signal k times over the graph and accumulates information at each node up to its k-hop neighborhood
- \blacktriangleright Combine shifts to get output $\mathbf{y} = \mathbf{H}(\mathbf{S})\mathbf{x} = \sum_{k=0}^{K} h_k \mathbf{S}^k \mathbf{x}$
- ▶ h_k weighs the kth hop, filtering combines them.
- ▶ Used in a variety of applications for graphs of fixed size [10, 1]



Graph filters: Illustration





Stochastic attachment

- ▶ Let $\mathcal{G} = \{\mathcal{V}, \mathcal{E}\}$ with N nodes $\{v_1, \ldots, v_N\}$ with $\mathbf{A} \in \mathbb{R}^{N \times N}$ as adjacency matrix
- ▶ v_+ incoming node forms graph $\mathcal{G}_+ = \{\mathcal{V} \cup v_+, \mathcal{E}_+\}$
- \blacktriangleright However, we do not know how v_+ will connect
- ▶ Let $\mathbf{a}_+ \in \mathbb{R}^N$ be a random vector containing attachment pattern of v_+

$$[\mathbf{a}_{+}]_{i} = \begin{cases} w_{i} & \text{with probability } p_{i} \\ 0 & \text{with probability } (1 - p_{i}) \end{cases}$$
(1)

▶ v_+ attaches to v_i with probability p_i with edge weight w_i



Modeling unknown connectivity

- \blacktriangleright w: vector of edge weights ${\bf p}$: vector of probabilities characterise the attachment behaviour of v_+
- \blacktriangleright The updated adjacency and Laplacian matrices are

$$\mathbf{A}_{+} = \begin{bmatrix} \mathbf{A} & \mathbf{a}_{+} \\ \mathbf{a}_{+}^{\top} & 0 \end{bmatrix}, \ \mathbf{L}_{+} = \begin{bmatrix} \mathbf{L} + \operatorname{diag}(\mathbf{a}_{+}) & -\mathbf{a}_{+} \\ -\mathbf{a}_{+}^{\top} & \mathbf{a}_{+}^{\top} \mathbf{1} \end{bmatrix}$$
(2)

- $\blacktriangleright \text{ Statistics of } \mathbf{a}_+ \colon \mathbb{E}[\mathbf{a}_+] = \mathbf{w} \circ \mathbf{p}, \quad \operatorname{cov}(\mathbf{a}_+) = \mathbf{\Sigma}_+ = \operatorname{diag}(\mathbf{w}^{\circ 2} \circ \mathbf{p} \circ (\mathbf{1} \mathbf{p}))$
- \blacktriangleright What is a good w, p? Our answer: It depends on the task!
- \blacktriangleright Target: Find w and p which learns attachment behviour of new nodes



Processing data with this attachment model

- ▶ Let $\mathbf{x} = [x_1, \dots, N]^\top$ be the existing graph signal on \mathcal{G}
- ▶ We want to predict x_+ at v_+
- \blacktriangleright For this we use a graph filter $\mathbf{h} \in \mathbb{R}^{K+1}$ of order K
- ▶ Goal: Learn **w**, **p** for predicting x_+
- ▶ Done via minimizing a task-aware loss



Key Idea



Key Idea: Learning from how other new users have already interacted with the network with their data



Key Idea: More concrete

Learn empirically from a batch of new nodes, each with its own attachment pattern \mathbf{a}_+ and signal x_+



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Learn empirically from a batch of new nodes, each with its own attachment pattern \mathbf{a}_+ and signal x_+

$$\min_{\mathbf{p},\mathbf{w}} \mathbb{E} \left[f_{\mathcal{T}}(\mathbf{p}, \mathbf{w}, \mathbf{a}_{t+}, \mathbf{x}, x_{+}) \right] + g_{\mathcal{T}}(\mathbf{p}, \mathbf{b}_{t+}) + h_{\mathcal{T}}(\mathbf{w}, \mathbf{a}_{t+})$$
subject to $\mathbf{p} \in \mathcal{P}, \mathbf{w} \in \mathcal{W}$

$$(3)$$

- ► $f_{\mathcal{T}}(\mathbf{p}, \mathbf{w}, \mathbf{a}_{t+}, \mathbf{x}, x_{+})$: Task-aware loss
- ▶ $g_{\mathcal{T}}(\mathbf{p}, \mathbf{b}_{t+})$: Priors on the attachment probabilities
- \blacktriangleright $h_{\mathcal{T}}(\mathbf{w}, \mathbf{a}_{t+})$: Priors on the weights
- ▶ Nonconvex problem, learnt by alternating projected gradient descent
- ▶ Analysis of convergence and perturbation behaviour in [11].



Experimental Setup

We compare with the following baselines:

- Uniformly-at-random: Heuristic, $\mathbf{p} = \frac{1}{N} \mathbf{1}$
- ▶ Degree-biased preferential attachment, Heuristic, $\mathbf{p} = \frac{\mathbf{d}}{\mathbf{1}^\top \mathbf{d}}$
- \blacktriangleright Data-based attachment: Learnt empirically from the \mathbf{a}_+ s during training
- \blacktriangleright Erdos Renyi and Barabasi Albert graphs for synthetic data
- **Movielens 100K** for real world data
- \blacktriangleright All hyperparameters found via cross validation



Synthetic Results



- ▶ Outperforms heuristical attachment models
- \blacktriangleright Typically, ℓ_2 regularizers on ${\bf p}$ and ${\bf w}$ work better
- \blacktriangleright We also found that training both ${\bf p}$ and ${\bf w}$ is more beneficial



Rating Prediction on Movielens 100K data

- Existing 35 nearest neighbor graph of N = 50 users for each movie
- ▶ Trained on individual movie with varying degrees of ratings (personalized)
- ▶ Trained on all movies as well

	Proposed	Attachment Only	Random	Preferential	Mean Prediction
Item 1	0.494	0.537(+8.7)	0.5417(+9.7)	0.527(+6.7)	0.669(+35.4)
Item 48	0.492	0.611(+24.2)	0.53(+7.7)	0.62(+26)	0.55(+11.8)
Item 459	0.462	0.49(+6)	0.52(+12.5)	0.40(+12.5)	1.07(+131)
Item 550	0.512	0.678(+32.4)	0.66(+29)	0.692(+35.2)	0.643(+25.6)
Item 57	0.049	0.057(+16.3)	0.41(+736)	0.20(+308)	0.32(+553)
Item 877	0.99	1.04(+5)	1.07(+8)	1.05(+6)	1.01(+1.72)
All Items	0.799	0.802(+0.38)	0.821(+2.75)	0.820(+2.63)	0.832(+4.1)

- Proposed approach does much better in personalized cases
- ▶ Differences evened out for all items



Overview

- ▶ Learning task-aware stochastic attachment model
- \blacktriangleright We can design graph filters under the same scenarios. More details in [12]
- ▶ Some results using graph convolutional neural networks¹
- ▶ Also possible to learn the attachment model and filters jointly
- ▶ However, all these works consider one incoming node added with replacement for training

Online graph filter design



Sequentially expanding graphs

- ▶ Graphs with growing number of nodes
- ▶ Delay-sensitive applications
- ▶ Time-varying nature of data at incoming node





Scenario

Challenges

 \blacktriangleright Streaming nature of the growing graph and data over it

Lack of connectivity of the incoming nodes (i.e., cold start in recommendation)

Targets

▶ How to adapt the graph filter to the change in topology?

▶ How to design the filter when the connectivity is not known and the graph is growing?



Proposed Solution

▶ Train graph filters online, i.e., update with each incoming node

▶ Use stochastic attachment when node information is not known and update filters online

Adapt the stochastic attachment pattern to predict the attachment over time

▶ Obtain performance bounds w.r.t known attachment patterns



Problem Formulation: Deterministic attachment

▶ Starting graph $\mathcal{G}_0 = \{\mathcal{V}_0, \mathcal{E}_0\}$: N_0 nodes, M_0 edges, adj. matrix \mathbf{A}_0 .

▶ v_1, \ldots, v_T be T sequentially incoming nodes, node v_t attaches to graph \mathcal{G}_{t-1}

▶ Vector $\mathbf{a}_t = [a_1, \dots, a_{N_{t-1}}]^\top \in \mathbb{R}^{N_{t-1}}$ represents directed connectivity of v_t at time t

▶ The expanded adjacency matrix at time $t \ \mathbf{A}_t \in \mathbb{R}^{N_t \times N_t}$ equals

$$\mathbf{A}_t = \begin{bmatrix} \mathbf{A}_{t-1} & \mathbf{0} \\ \mathbf{a}_t^\top & \mathbf{0} \end{bmatrix}$$



Node v_t attaches to $v_i \in \mathcal{V}_{t-1}$ obeying Prob. vector $\mathbf{p}_t \in \mathbb{R}^{N_{t-1}}$ and weights $\mathbf{w}_t \in \mathbb{R}^{N_{t-1}}$



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 $\blacktriangleright \text{ Time-varying statistics of } \mathbf{a}_t: \mathbb{E}\left[\mathbf{a}_t\right] = \mathbf{p}_t \circ \mathbf{w}_t \ ; \ \boldsymbol{\Sigma}_t = \mathsf{diag}(\mathbf{w}_t^{\circ 2} \circ \mathbf{p}_t \circ (\mathbf{1} - \mathbf{p}_t))$



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▶ Parameter of interest: an order K graph filter $\mathbf{h} = [h_0, \dots, h_K]^\top$



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$$\blacktriangleright \text{ Output at new node } [\tilde{\mathbf{y}}_t]_{N_t} := \hat{x}_t = \mathbf{a}_t^\top \sum_{k=1}^K h_k \mathbf{A}_{t-1}^{k-1} \tilde{\mathbf{x}}_t = \mathbf{a}_t^\top \mathbf{A}_{x,t-1} \mathbf{h}$$



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▶ Update $\mathbf{h}(t)$ based on the loss and current estimate $\mathbf{h}(t-1)$.


Online graph filtering

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- ▶ Update $\mathbf{h}(t)$ based on the loss and current estimate $\mathbf{h}(t-1)$.
- $\blacktriangleright \ \mathbf{a}_t$ revealed in the stochastic setting



Deterministic online filtering

► Loss function
$$l_t(\mathbf{h}, x_t) = \frac{1}{2} (\mathbf{a}_t^\top \mathbf{A}_{x,t-1} \mathbf{h} - x_t)^2 + \mu ||\mathbf{h}||_2^2$$
 with $\mu > 0$.

▶ Update: Projected online gradient descent [13]

$$\mathbf{h}(t) = \prod_{\mathcal{H}} (\mathbf{h}(t-1) - \eta \nabla_h l_t(\mathbf{h}, x_t) |_{\mathbf{h}(t-1)})$$

where $\prod_{\mathcal{H}}(\cdot)$ denotes projection on \mathcal{H} .

▶ Computational complexity of order $\mathcal{O}(K(M_t + M_{\max})), M_{\max}$: maximum edges formed by v_t



Stochastic online filtering

 \blacktriangleright Predefined \mathbf{p}_t and \mathbf{w}_t based on stochastic attachment rule and \mathbf{w}_t for all t

 $\blacktriangleright \text{ Loss function } l_t(\mathbf{h}, x_t) = \mathbb{E}\left[\frac{1}{2}(\mathbf{a}_t^\top \mathbf{A}_{x,t-1}\mathbf{h} - x_t)^2\right] + \mu ||\mathbf{h}||_2^2 \text{ with } \mu > 0.$

▷ Computational complexity at time t of order $\mathcal{O}(K(M_t + N_t))$

▶ Potential drawback: Reliance on one type of attachment rule



 \triangleright Update \mathbf{p}_t and \mathbf{w}_t for all t, i.e., adapt to the incoming node behaviour



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Projected online gradient descent but update $\mathbf{h}(t)$, $\mathbf{m}(t)$, and $\mathbf{n}(t)$

▶ Computational complexity at time t of order $\mathcal{O}(K(M_0 + N_t) + N_t(M))$



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$$R_T(\mathbf{h}^*) = \sum_{t=1}^T l_t(\mathbf{h}_t, x_t) - l_t(\mathbf{h}^*, x_t)$$



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Deterministic regret

$$\frac{1}{T}R_T(\mathbf{h}^{\star}) \le \frac{||\mathbf{h}^{\star}||_2^2}{2\eta T} + \frac{\eta}{2}L_d^2$$

Stochastic regret

$$\frac{1}{T}R_{s,T}(\mathbf{h}^{\star}) \leq \frac{1}{T} \left(\sum_{t=1}^{T} w_{h}^{2} Y^{2}(||\mathbf{p}_{t}||_{2}^{2} + M_{max}) + 2Rw_{h} Y \sqrt{||\mathbf{p}_{t}||_{2}^{2} + M_{max}} + w_{h}^{2} Y^{2} \bar{\sigma}_{t}^{2} + L_{d} ||\mathbf{h}^{s}(t-1) - \mathbf{h}^{d}(t-1)|| \right) + \frac{||\mathbf{h}^{\star}||_{2}^{2}}{2\eta} + \frac{\eta}{2} L_{d}^{2} T$$
(4)

Stochastic regret depends on choice of attachment rule



Results

Compare with (1) Batch filter, (2) pre-trained filter, (3) Online kernel-based methods OKL and OMHKL, (4) Prediction Correction-based Online filtering (PC-OGF)

	Synthetic Data						Real Data			
Method	Filter		WMean		Kernel		Movielens100K		COVID	
	NRMSE	Sdev	NRMSE	Sdev	NRMSE	Sdev	NRMSE	Sdev	NRMSE	Sdev
D-OGF (ours)	0.038	0.04	0.02	0.02	0.28	0.05	0.26	0.01	0.21	0.02
S-OGF (ours)	0.26	0.04	0.3	0.06	0.36	0.11	0.28	0.007	0.31	0.02
Ada-OGF (ours)	0.26	0.04	0.33	0.07	0.48	0.21	0.28	0.007	0.26	0.007
PC-OGF (ours)	0.22	0.04	0.26	0.04	0.32	0.06	0.29	0.01	0.26	0.003
Batch	0.05	0.03	0.08	0.04	1.12	0.26	6.7	0.1	0.17	0.03
pre-trained	0.13	0.07	0.10	0.05	0.59	0.36	0.84	0.02	2.5	0.9
OKL	0.24	0.03	0.27	0.05	0.30	0.06	0.27	0.01	0.25	0.02
OMHKL	0.25	0.04	0.4	0.05	0.5	0.2	0.27	0.01	0.25	0.02



Results



Normalized cumulative regret for stochastic online filtering on synthetic data



Overview

▶ We proposed online filtering over graphs that grow sequentially over time. More details in [14]

Extension to growing simplicial complexes in [15]

▶ Provide regret analysis in terms of attachment behaviour of incoming nodes

▶ Numerical results indicate online filters perform collectively better than kernel methods which do not utilize the data, pre-trained filters, and even a batch filter.



Online topology identification



Graph Topology Identification

- ▶ Graph topology is often unknown or unavailable
- ▶ Topology identification concerns learning the topology associated with observed data [4]
- ▶ Typically uses prior information linking the graph with the topology, e.g., Gaussianity [16], smoothness [17], stationarity [18]



Contribution

- Existing works focus on estimating graphs (static or dynamic) of fixed size [19, 20, 21]
- ▶ Instead, we focus on cases where the underlying graph grows steadily in size over time
- Examples: New companies joining financial networks [22]
- **Target** Learn the growing topology





Fundamentals of graph learning

- ▶ $\mathbf{X} \in \mathbb{R}^{N \times T}$: Signals collected over T time instances
- ▶ \mathcal{G} involved via the graph-shift operator $\mathbf{S} \in \mathbb{R}^{N \times N}$ (e.g. \mathbf{A}, \mathbf{L})
- \blacktriangleright Estimate ${\bf S}$ from observed signals ${\bf X} \in \mathbb{R}^{N \times T}$ solving

$$\min_{\mathbf{S}\in\mathcal{S}} \quad \mathcal{L}(\mathbf{S}, \hat{\mathbf{C}}) + \lambda \|\mathbf{S}\|_1 \tag{5}$$

- ▶ Loss \mathcal{L} relates **S** and sample covariance $\hat{\mathbf{C}} = \frac{1}{T} \mathbf{X} \mathbf{X}^{\top}$
- \triangleright ℓ_1 norm promotes sparsity



Expanding graphs

- \triangleright In the expanding graph setting new nodes may arrive every time instant t
- ▶ The size of the graph increases, i.e., $|\mathcal{V}_t| = N_t \leq |\mathcal{V}_{t+1}| = N_t + 1$





Expanding graphs



▶ Sequential arrival of nodes results in the following block structure

$$\mathbf{S}_{t+1} = \begin{bmatrix} [\mathbf{S}_{t+1}]_{\mathcal{V}_t} & [\mathbf{S}_{t+1}]_{\mathcal{V}_t,\mathcal{I}_{t+1}} \\ [\mathbf{S}_{t+1}]_{\mathcal{I}_{t+1},\mathcal{V}_t} & [\mathbf{S}_{t+1}]_{\mathcal{I}_t} \end{bmatrix}, \qquad \mathbf{x}_{t+1} = \begin{bmatrix} [\mathbf{x}_{t+1}]_{\mathcal{V}_t} \\ [\mathbf{x}_{t+1}]_{\mathcal{I}_t} \end{bmatrix}$$

▶ Where \mathcal{I}_t denoting the indexes of incoming nodes at time t



Batch Solution

- \blacktriangleright One option is to learn the new topology \mathbf{S}_t^* every time a new node and its data appears
- We calculate the sample covariance $\hat{\mathbf{C}}_t$ and solve

$$\mathbf{S}_{t}^{*} = \operatorname{argmin}_{\mathbf{S} \in \mathcal{S}_{N_{t}}} \mathcal{L}(\mathbf{S}, \hat{\mathbf{C}}_{t}) + \alpha d([\mathbf{S}]_{\mathcal{V}_{t-1}}, \mathbf{S}_{t-1}^{*}) + \lambda \|\mathbf{S}\|_{1}$$
(6)

- Second term accounts for smooth variation between old nodes
- \blacktriangleright Provide high-quality estimates \mathbf{S}_t^*
- Excessive computational cost, not fit for delay-sensitive applications



Online Algorithm

- **Step 1**: At time $t, \mathbf{x}_t \in \mathbb{R}^{N_t}$ is available
- ▶ Update the covariance matrix as

$$\hat{\mathbf{C}}_t = \mathbf{M}_1 \circ \underline{\hat{\mathbf{C}}}_{t-1}^{(N_t)} + \mathbf{M}_2 \circ \mathbf{x}_t \mathbf{x}_t^{\top}$$
(7)

 \blacktriangleright Mask \mathbf{M}_1 scales importance of past observations, \mathbf{M}_2 that of the latest



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- \blacktriangleright Mask \mathbf{M}_1 scales importance of past observations, \mathbf{M}_2 that of the latest
- **Step 2**: Loss function at time t given by

$$f_t(\mathbf{S}) = \mathcal{L}(\mathbf{S}, \hat{\mathbf{C}}_t) + \alpha d([\mathbf{S}]_{\mathcal{V}_{t-1}}, \hat{\mathbf{S}}_{t-1}) + ||\mathbf{S}||_1$$
(8)

- ► Proximal gradient step: $\check{\mathbf{S}}_t = \Pi_{\mathcal{S}_{N_t}} \left(T_{\eta\lambda} \left(\hat{\underline{\mathbf{S}}}_{t-1}^{(N_t)} \eta \nabla f(\hat{\underline{\mathbf{S}}}_{t-1}^{(N_t)}) \right) \right)$
- ▶ Projection on to constraint set of $N_t \times N_t$ shift operators



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$$\hat{\mathbf{C}}_t = \mathbf{M}_1 \circ \underline{\hat{\mathbf{C}}}_{t-1}^{(N_t)} + \mathbf{M}_2 \circ \mathbf{x}_t \mathbf{x}_t^{\top}$$
(7)

- \blacktriangleright Mask \mathbf{M}_1 scales importance of past observations, \mathbf{M}_2 that of the latest
- **Step 2**: Loss function at time t given by

$$f_t(\mathbf{S}) = \mathcal{L}(\mathbf{S}, \hat{\mathbf{C}}_t) + \alpha d([\mathbf{S}]_{\mathcal{V}_{t-1}}, \hat{\mathbf{S}}_{t-1}) + ||\mathbf{S}||_1$$
(8)

- ► Proximal gradient step: $\check{\mathbf{S}}_t = \Pi_{\mathcal{S}_{N_t}} \left(T_{\eta\lambda} \left(\hat{\underline{\mathbf{S}}}_{t-1}^{(N_t)} \eta \nabla f(\hat{\underline{\mathbf{S}}}_{t-1}^{(N_t)}) \right) \right)$
- ▶ Projection on to constraint set of $N_t \times N_t$ shift operators
- **Step 3**: update estimate GSO as $\hat{\mathbf{S}}_t = h\check{\mathbf{S}}_t + (1-h)\hat{\mathbf{S}}_{t-1}^{(N_t)}$



The Gaussian case

- Assume signals are drawn from Gaussian distribution $\mathbf{x}_t \sim \mathcal{N}(\mathbf{0}, \mathbf{S}_t^{-1})$
- ▶ GSO estimated maximizing the regularized log-likelihood [16]

$$\nabla f_t(\mathbf{S}) = \hat{\mathbf{C}}_t - (\mathbf{S} + \epsilon \mathbf{I})^{-1} + \alpha \nabla d([\mathbf{S}]_{\mathcal{V}_{t-1}}, \hat{\mathbf{S}}_{t-1})$$
$$\Pi_{\mathcal{S}_{N_t}}(\mathbf{S}) = \mathbf{V} \min\left\{ \max\left\{\mathbf{\Lambda}, 0\right\}, \sigma^{1/2} \right\} \mathbf{V}^{\top}$$

- \blacktriangleright Gradient from log-likelihood and projection to PSD matrices set
- ▶ Computational complexity of $\mathcal{O}(N_t^3)$ with over-the-shelf methods



- ▶ We want to know how accurately the online approach can track the growing topology
- ▶ Cumulative regret measures error made over a sequence

$$\sum_{t=1}^{T} \|\hat{\mathbf{S}}_t - \mathbf{S}_t^*\|_F \tag{9}$$

▶ With step size $\eta \leq \epsilon^2$, dynamic cumulative regret is upper bounded by

$$\sum_{t=1}^{T} \|\hat{\mathbf{S}}_{t} - \mathbf{S}_{t}^{*}\|_{F} \le K_{1} + K_{2} \sum_{t=2}^{T} \|\mathbf{S}_{t}^{*} - \underline{\mathbf{S}}_{t-1}^{*(N_{t})}\|_{F}$$
(10)

▶ Bounded by how much the optimal topology changes.



Numerical evaluation

- \blacktriangleright ER graphs growing from 100 to 120 nodes with different frequencies
- ▶ Low-frequency setting: 20 new nodes arrive simultaneously
- ▶ High-frequency setting: nodes arrive in 4 groups of 5 nodes



 \blacktriangleright Recovers from disruptions in both settings, insufficient recovery tme \implies error accumulation



Numerical Evaluation

- ▶ COVID-19 dataset with incidence rates reported by various U.S. states and territories.
- Start with 46 nodes and reports from additional states are available at $t = \{312, 652\}$.



- Error rapidly increases after t = 350 coinciding with a rise in incident rates.
- ▶ Our algorithm adapts to the new topology and the error decreases.



Overview

- ▶ An online graph learning method tailored to expanding graphs
- ▶ Define efficient node-dependent covariance updates
- ▶ Propose an algorithm based on PPG descent to learn from streaming data
- ▶ Bounded dynamic regret



Challenges and further possibilities

Overview

- ▶ A SP-based principled framework for expanding graphs
- \blacktriangleright Stochastic model-based attachments for unknown connectivity
- ▶ Online approaches for SP and Topology identification with some theoretical analysis
- Relatively well compared to known topology



Challenges and further possibilities

Overview

- ▶ A SP-based principled framework for expanding graphs
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Challenges and further possibilities

- $\blacktriangleright\,$ Modeling data on expanding graphs a challenge
- ▶ More refined stochastic attachment models
- ▶ Potential for using physics-driven models involving the spectrum
- \blacktriangleright Extension to expanding higher order networks possible



END



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