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Effecient Sensor Position Selection Using Graph Signal Sampling Theory

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I. INTRODUCTION

Sensor Placement Problem

- ullet Selects F sensor locations from N locations to optimize performance
- Often has to handle *large number* of sensors *distributed nonuniformly*
- Many conventional researches [1, 2]
- Assume that the spatial phenomena are modeled as a Gaussian process
- Place the sensors at the most informative locations

Graph Signal Processing

- Can efficiently analyze complex, irregular, and high-dimensional data
- Relationship between data points is represented as edges in a graph This Paper
- · Proposes a sensor selection approach based on the sampling theory for graph signals [4, 5, 6]
- · Shows that the conventional approaches can be viewed as graph vertex domain operations

Graph Signals

- Graph: $\mathcal{G}=\{\mathcal{V},\mathcal{E}\}$ where \mathcal{V} and \mathcal{E} are sets of nodes and links
- Adjacency Matrix: A(m, n) = weight of link between node m and n
- Diagonal Degree Matrix: $D(m,m) = \sum_n A(m,n)$
- Normalized Graph Laplacian Matrix (GLM): $\mathcal{L} = \mathbf{I} \mathbf{D}^{-1/2} \mathbf{A} \mathbf{D}^{-1/2}$
 - \rightarrow eigenvalues: $\{\lambda_i\}_{i=0...N-1} = \sigma(\mathcal{L}) = spectrum of graph$ eigenvectors: $\{\boldsymbol{u}_{\lambda_i}\}_{i=0...N-1}$
- $\mathbf{X}_{\mathcal{A}\mathcal{B}}$: matrix extracting \mathcal{A} rows and \mathcal{B} columns from \mathbf{X}
 - $oldsymbol{x}_{\mathcal{A}}:$ restriction of $oldsymbol{x}$ to its components indexed by \mathcal{A}

III. SENSOR PLACEMENT PROBLEM

Problem Setting

- Selecting $|\mathcal{S}| = F$ sensors from $|\mathcal{V}| = N$ possible locations
- ullet Signal f has the Gaussian joint zero-mean distribution:

$$p(\boldsymbol{f}) = \frac{1}{(2\pi)^{\frac{N}{2}}|\mathbf{K}|} \exp\left(-\frac{1}{2}\boldsymbol{f}^T\mathbf{K}^{-1}\boldsymbol{f}\right)$$
 K: covariance matrix

Entropy Criterion [1]

★ Sensors are selected so that the uncertainty of a measurement with respect to previous measurements is maximized

$$\mathcal{S}^* = \underset{\mathcal{S} \subset \mathcal{V}: |\mathcal{S}| = k}{\operatorname{arg \, min}} H(f_{\mathcal{S}^c}|f_{\mathcal{S}}) = \underset{\mathcal{S} \subset \mathcal{V}: |\mathcal{S}| = k}{\operatorname{arg \, max}} H(f_{\mathcal{S}}) \quad \mathcal{S}^c = \mathcal{V} \setminus \mathcal{S}$$

• Conditional entropy: $H(f(y)|\mathbf{f}_{\mathcal{S}}) = \frac{1}{2}\log(2\pi e(\mathcal{K}(y,y) - \mathbf{K}_{y\mathcal{S}}\mathbf{K}_{\mathcal{S}\mathcal{S}}^{-1}\mathbf{K}_{\mathcal{S}y}))$

Mutual Information Criterion [2]

★ Selects the locations that more significantly reduce the uncertainty of the rest of the space

$$\mathcal{S}^* = \underset{\mathcal{S} \subset \mathcal{V}: |\mathcal{S}| = k}{\operatorname{arg max}} H(f_{\mathcal{S}^c}) - H(f_{\mathcal{S}^c}|f_{\mathcal{S}}) := \operatorname{MI}(\mathcal{S}) \quad \overline{\mathcal{S}} = \mathcal{V} \setminus (\mathcal{S} \cup \mathcal{Y})$$

IV. GRAPH SIGNAL FOR SENSOR SELECTION

• Sensors and observed signals are viewed as nodes and graph signals How to determine the connection of nodes

Assumption: the random signals have following distributions:

$$p(\mathbf{f}) \propto \exp\left(-\sum_{i}\sum_{j}A(i,j)(f(i)-f(j))^{2}-\delta\sum_{i}f(i)^{2}\right)$$

= $\exp\left(-\mathbf{f}^{T}(\mathbf{L}+\delta\mathbf{I})\mathbf{f}\right)$ δ : parameter

★ Graph Laplacian matrix is obtained from the covariance matrix

$$\mathbf{L} = \mathbf{K}^{-1} - \delta \mathbf{I}$$

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- [5] I. Shomorony and A. S. Avestimehr, "Sampling large data on graphs," in *Proc. GlobalSIP'14*, 2014, pp. 933–936. [6] A.Gadde and A.Ortega, "A probabilistic interpretation of sampling theory of graph signals," in *Proc. ICASSP'15*, 2015, pp. 3257–3261.

V. SELECTION USING GRAPH SAMPLING THEORY

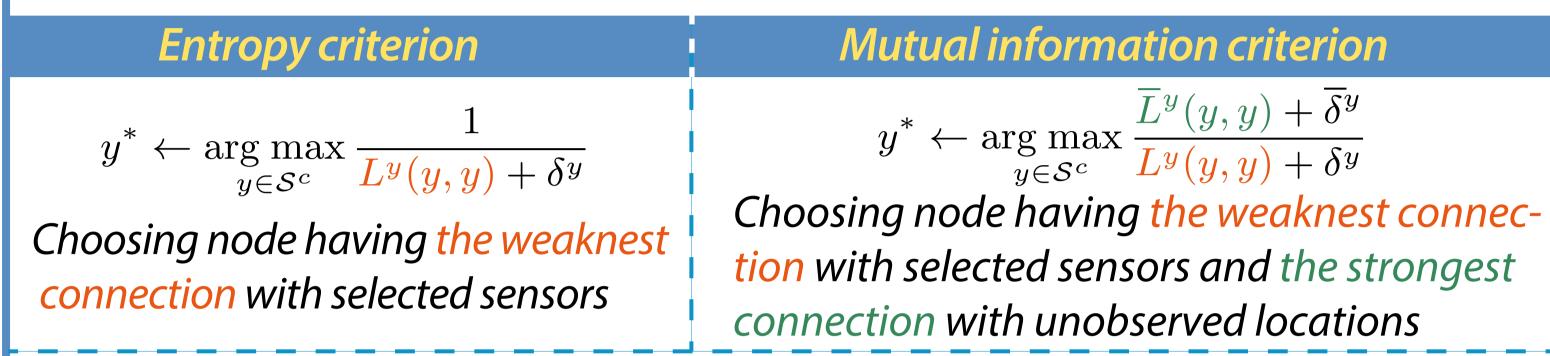
Chooses sensors so as to maximize the cut-off frequency in the graph spectral domain: $S^ = \arg \max \omega_c(S)$

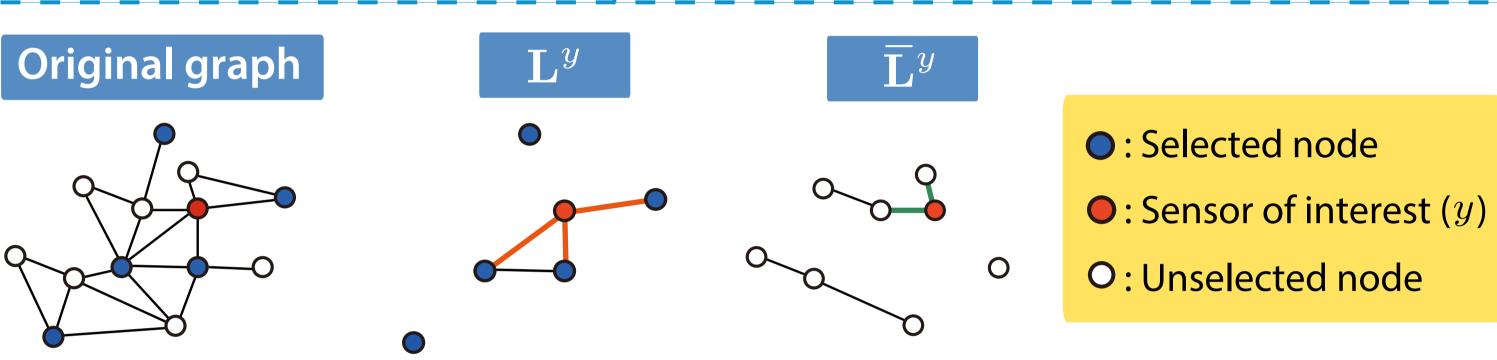
Cut-off frequncy $\omega_c(\mathcal{S})$: the maximum frequency of a signal that can be perfectly recoverd from the samples on the subset ${\cal S}$

• For optimization, we use three methods based on; eigenvalue [3] (EV), singlar value [4] (SVD), and standard basis [5] (SV)

Relationships with Existing Methods

- Representing the entropy and MI criteria by using graph Laplacian matrix
 - Using the relationship [6]: $\mathbf{L}_{\mathcal{S}^c} + \delta \mathbf{I} = (\mathbf{K}_{\mathcal{S}^c} \mathbf{K}_{\mathcal{S}^c\mathcal{S}}(\mathbf{K}_{\mathcal{S}})^{-1}\mathbf{K}_{\mathcal{S}^c\mathcal{S}}^T)^{-1}$





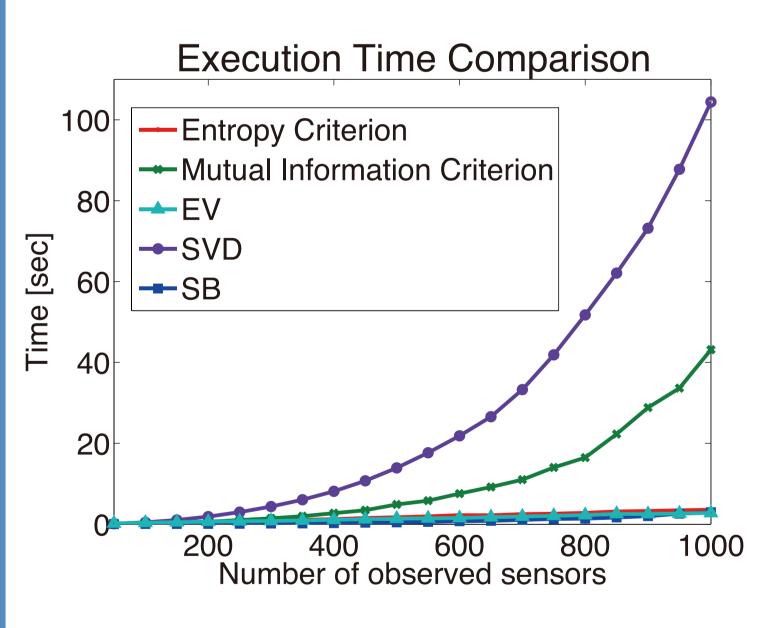
 $L^y(y,y)$ = degree of yth node in graph having nodes $S \cup y$ and edges in $S \cup y$ $\overline{L}^y(y,y) = \text{degree of } y \text{th node in graph having unselected nodes } \mathcal{S}^c \text{ and edges in } \mathcal{S}^c$

Conventional approach can be viewed as graph vertex domain operations, whereas our approach can be viewed as graph spectral domain approach

VI. EXPERIMENTAL RESULTS

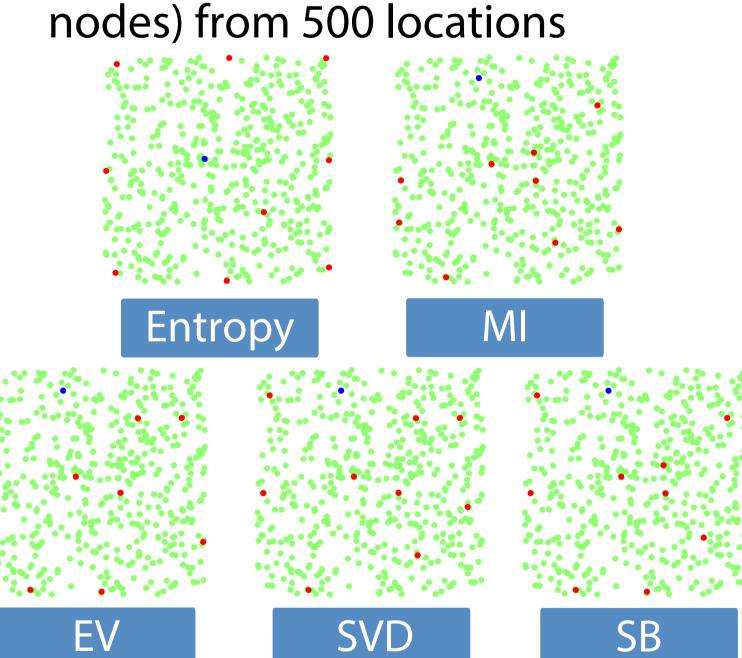
1. Execution Time

• $|\mathcal{S}| = |\mathcal{V}|/10$ sensors are selected with various number of \mathcal{V}



2. Choosing Sensor Positions

 10 selected locations (red and blue nodes) from 500 locations



3. Predicting Values on Unobserved Locations

- Test signals are randomly generated according to the GP model and are corrupted by the additive white Gaussian noise
- ullet Original signals are reconstruct only from the signals on ${\mathcal S}$
- Estimated signal is obtained by method in [3]

Performance Comparison (Average of 500 Tested Signals): SNR [dB]

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$ \mathcal{S} $	10	20	40	60	80	100
Entropy	-2.77	-2.17	0.28	2.02	2.13	5.75
MI	3.44	5.88	7.57	8.75	9.38	10.25
EV	3.52	6.46	7.74	8.53	9.44	10.37
SVD	3.83	5.98	7.55	8.88	9.14	10.29
SB	3.61	6.34	7.86	8.21	8.67	10.31

VII. CONCLUSION

- Optimal sensor selection method based on the graph sampling theorem has been proposed
- All the proposed methods achieved better performance
- As a future work, we will further investigate the theoretical issues and better reconstruction algorithms