MAT 271: Applied & Computational Harmonic Analysis Lecture 20: Multiscale Basis Dictionaries on Graphs and Networks

Naoki Saito (with help from Jeff Irion)

Department of Mathematics University of California, Davis

March 13, 2014

- Introductory Remarks
- Motivations: Why Graphs?
- Background
 - Basic Graph Theory Terminology
 - Graph Laplacians
 - Graph Partitioning via Spectral Clustering
- Multiscale Basis Dictionaries
 - Hierarchical Graph Laplacian Eigen Transform (HGLET)
 - Generalized Haar-Walsh Transform (GHWT)
- Best-Basis Algorithm for HGLET & GHWT
- 6 Approximation Experiments
- Summary and Future Work

- For much more details of this part of lecture, please check my course website on "Harmonic Analysis on Graphs & Networks": http://www.math.ucdavis.edu/~saito/courses/HarmGraph/ as well as my two short articles with Jeff Irion at http://www.math.ucdavis.edu/~saito/publications/.
- We rely on the so-called graph Laplacians to construct our multiscale basis dictionaries. Some good references on graph Laplacian eigenvalues are:
 - R. B. Bapat: *Graphs and Matrices*, Universitext, Springer, 2010.
 - A. E. Brouwer & W. H. Haemers: Spectra of Graphs, Springer, 2012.
 - F. R. K. Chung: Spectral Graph Theory, Amer. Math. Soc., 1997.
 - D. Cvetković, P. Rowlinson, & S. Simić: An Introduction to the Theory of Graph Spectra, Vol. 75, London Mathematical Society Student Texts, Cambridge Univ. Press, 2010.
 - D. Spielman: "Spectral graph theory," in Combinatorial Scientific Computing (O. Schenk, ed.), Chap. 18, pp. 495–524, CRC Press, 2012.
- As for the graph Laplacian eigenfunctions, there are not too many books (although there may be many papers); one of the good books is
 - T. Bıyıkoğlu, J. Leydold, & P. F. Stadler, Laplacian Eigenvectors of Graphs, Lecture Notes in Mathematics, vol. 1915, Springer, 2007.

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- More and more data are collected in a distributed and irregular manner; they are not organized such as familiar digital signals and images sampled on regular lattices. Examples include:
 - Data from sensor networks
 - Data from social networks, webpages, ...
 - Data from biological networks
 - . . .
- It is quite important to analyze:
 - Topology of graphs/networks (e.g., how nodes are connected, etc.)
 - Data measured on nodes (e.g., a node = a sensor, then what is an edge?)

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- Hence, we need to lift such tools for unorganized and irregularly-sampled datasets including those represented by graphs and networks.
- Moreover, constructing a graph from a usual signal or image and analyzing it can also be very useful! E.g., Nonlocal means image denoising of Buades-Coll-Morel.

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An Example of Sensor Networks

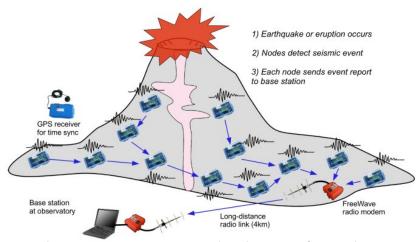


Figure : Volcano monitoring sensor network architecture of Harvard Sensor Networks Lab

An Example of Social Networks

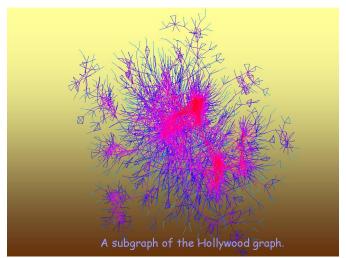


Figure: Through the courtesy of Prof. Fan Chung, UC San Diego

An Example of Biological Networks

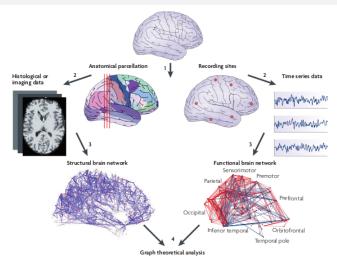
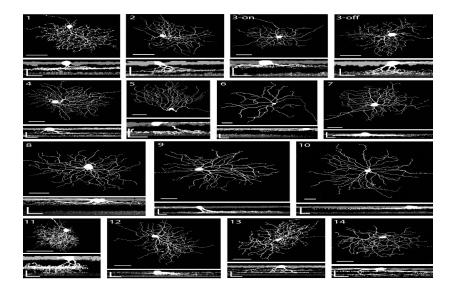
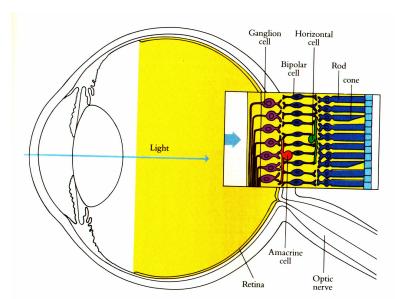


Figure: From E. Bullmore and O. Sporns, *Nature Reviews Neuroscience*, vol. 10, pp.186–198, Mar. 2009.

Another Biological Example: Retinal Ganglion Cells

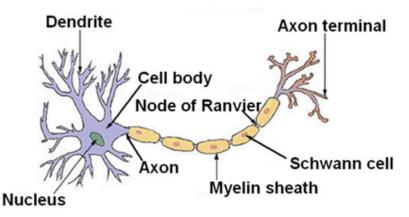


Retinal Ganglion Cells (D. Hubel: Eye, Brain, & Vision, '95)

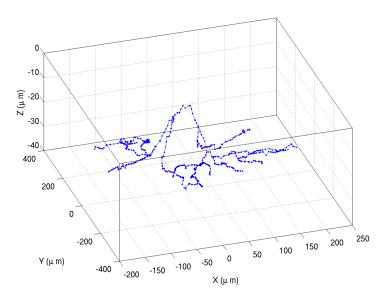


A Typical Neuron (from Wikipedia)

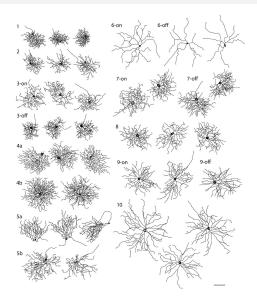
Structure of a Typical Neuron



Mouse's RGC as a Graph



Clustering using Features Derived by Neurolucida®



- Construct a graph each of whose vertices represents $k \times k$ patch of a given image (k may be 3,5,..., etc.) So each vertex represents a point in \mathbb{R}^{k^2} .
- Connect every pair of vertices with the weight $W_{ij} = \exp(-\| \mathrm{patch}_i \mathrm{patch}_j \|^2 / \epsilon^2)$ with appropriately chosen scale parameter $\epsilon > 0$.
- Compute the weighted average of the center pixel of each patch using the normalized weights $W_{ij}/\sum_{\ell}W_{i\ell}$. More precisely, the average of the center of the ith patch, $\overline{c}_i = \sum_j W_{ij}c_j/\sum_{\ell}W_{i\ell}$.
- See also an interesting work by Daitch-Kelner-Spielman: "Fitting a Graph to Vector Data," *Proc. 26th Intern. Conf. Machine Learning*, 2009.

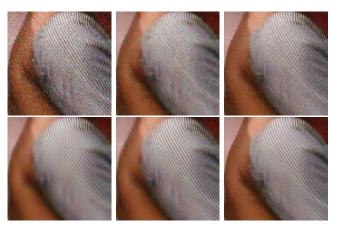
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From: A. Buades, B. Coll, and J.-M. Morel, *SIAM Review*, vol. 52, no. 1, pp. 113–147, 2010.

Noisy Image; Total Variation Denoising; Neighborhood Filter



Trans. Inv. Wavelets; Empirical Wiener; Nonlocal Means

Wavelets

- Have been quite successful on regular domains
- Have been extended to irregular domains ⇒ "2nd Generation Wavelets"

For example

- Hammond, Vandergheynst, and Gribonval (2011): wavelets via spectral graph theory
- Coifman and Maggioni (2006): diffusion wavelets
 Bremer et al. (2006): diffusion wavelet packets
- **Key difficulty:** The notion of *frequency* is ill-defined on graphs ⇒ The Fourier transform is not properly defined on graphs
- Common strategy: Develop wavelet-like multiscale transforms
- **Key Idea:** Use of the graph Laplacian eigenvectors as the substitution of the Fourier basis

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Goals

- Develop and implement multiscale transforms for data on graphs and networks; in particular, build *multiscale basis dictionaries* on graphs.
- Investigate their usefulness for a variety of applications including approximation, denoising, classification, and regression on graphs.
- In this lecture, we will focus on how to construct such dictionaries on graphs and demonstrate their usefulness for data approximation on graphs.

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- $V = V(G) = \{v_1, \dots, v_N\}$ is the set of vertices.
- $E = E(G) = \{e_1, \dots, e_{N'}\}$ is the set of edges, where $e_k = (v_i, v_j)$ represents an edge (or line segment) connecting between adjacent vertices v_i, v_j for some $1 \le i, j \le N$.
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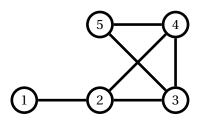




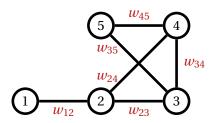




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Note that there are many ways to define w_{ij} .

For example, for unweighted graphs, we typically use

$$w_{ij} \coloneqq \begin{cases} 1 & \text{if } v_i \sim v_j \text{ (i.e., } v_i \text{ and } v_j \text{ are adjacent);} \\ 0 & \text{otherwise.} \end{cases}$$

This is often referred to as the adjacency matrix and denoted by A(G).

For weighted graphs, w_{ij} should reflect the similarity (or affinity) of information at v_i and v_j , e.g., if $v_i \sim v_j$, then

$$w_{ij} := 1/\operatorname{dist}(v_i, v_j)$$
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Graph Laplacians

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- sorted eigenvalues $0 = \lambda_0 < \lambda_1 \le \cdots \le \lambda_{N-1}$
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Why Graph Laplacians?

• Let $f \in \mathbb{R}^N$. Then

$$Lf(v_i) = d_{v_i}f(v_i) - \sum_{j \neq i} w_{ij}f(v_j).$$

This is a generalization of the finite difference approximation to the Laplace operator.

- After all, sines (cosines) are the eigenfunctions of the Laplacian on the rectangular domain with Dirichlet (Neumann) boundary conditions.
- Hence, the expansion of data measured at the vertices w.r.t. the eigenvectors of a graph Laplacian can be viewed as Fourier (or spectral) analysis of the data on that graph.

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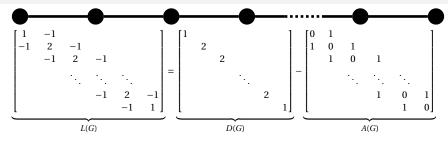
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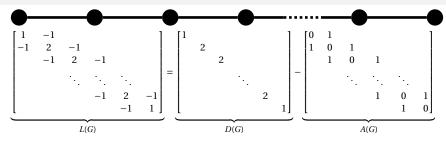
A Simple Yet Important Example: A Path Graph



The eigenvectors of this matrix are exactly the *DCT Type II* basis vectors used for the JPEG image compression standard! (See e.g., Strang, SIAM Review, 1999).

- $\lambda_k = 2 2\cos(\pi k/N) = 4\sin^2(\pi k/2N), k = 0, 1, ..., N 1.$
- $\phi_k(\ell) = \sqrt{2/N} \cos\left(\pi k(\ell + \frac{1}{2})/N\right), \ k, \ell = 0, 1, ..., N-1.$
- λ (eigenvalue) is a monotonic function w.r.t. k (frequency). However, for general graphs, λ does not have a simple relationship with k.

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$$(X, X^c) := \frac{\text{cut}(X, X^c)}{|X|} + \frac{\text{cut}(X, X^c)}{|X^c|},$$

$$\operatorname{cut}(X, X^c) := \sum_{v_i \in X; v_j \in X^c} W_{ij}$$

 Dividing by the number of nodes ensures that the partitions are of roughly the same size ⇒ we do not simply cleave a small number of nodes

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Let us reformulate the RatioCut minimization problem.

① Define $\mathbf{f} \in \mathbb{R}^N$ as

$$f_i := \begin{cases} \sqrt{\frac{|X^c|}{|X|}} & \text{if } v_i \in X \\ -\sqrt{\frac{|X|}{|X^c|}} & \text{if } v_i \in X^c \end{cases}$$

2 The RatioCut problem can be reformulated as

 $\min_{X\subset V} oldsymbol{f}^{\mathsf{T}} L oldsymbol{f}$ subject to $oldsymbol{f}$ defined as above

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$$f^{\mathsf{T}}Lf = \frac{1}{2} \sum_{\substack{i,j=1 \ v_j \in X}}^{N} W_{ij} (f_i - f_j)^2$$

$$= \frac{1}{2} \sum_{\substack{v_i \in X^c \ v_j \in X^c}} W_{ij} \left(\sqrt{\frac{|X^c|}{|X|}} + \sqrt{\frac{|X|}{|X^c|}} \right)^2$$

$$+ \frac{1}{2} \sum_{\substack{v_i \in X^c \ v_j \in X}} W_{ij} \left(-\sqrt{\frac{|X^c|}{|X|}} - \sqrt{\frac{|X|}{|X^c|}} \right)^2$$

$$= \operatorname{cut}(X, X^c) \left(\frac{|X^c|}{|X|} + \frac{|X|}{|X^c|} + 2 \right)$$

$$= \operatorname{cut}(X, X^c) \left(\frac{|X| + |X^c|}{|X|} + \frac{|X| + |X^c|}{|X^c|} \right)$$

$$= |V| \operatorname{RatioCut}(X, X^c)$$

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Unfortunately, this problem is NP hard... Relax!

A couple things to note about f:

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$$f \perp 1 \Leftrightarrow \sum f_i = 0$$

$$\sum_{i=1}^{N} f_i = \sum_{v_i \in X} \sqrt{\frac{|X^c|}{|X|}} - \sum_{v_i \in X^c} \sqrt{\frac{|X|}{|X^c|}}$$
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• If we relax our previous definition of f and simply require that (i) $f \perp 1$ and (ii) $||f|| = \sqrt{N}$, then we get the relaxed minimization problem¹:

$$\min_{\boldsymbol{f} \in \mathbb{R}^N} \boldsymbol{f}^{\mathsf{T}} L \boldsymbol{f}$$
 subject to $\boldsymbol{f} \perp \boldsymbol{1}, \ \|\boldsymbol{f}\| = \sqrt{N}$

- By the Rayleigh-Ritz Theorem, the solution is given by ϕ_1 (scaled as necessary), where ϕ_1 is the eigenvector corresponding to the second smallest eigenvalue of L.
- ϕ_1 is known as the Fiedler vector and is often used to partition a graph into two subsets.
- von Luxburg recommends the use of the random-walk version of the Laplacian matrix, $L_{\rm rw}:=I-D^{-1}W$, over the usual Laplacian matrix L which leads to the NCut and the generalized eigenvalue problem: $L\phi=\lambda D\phi$.

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Definition (Weak Nodal Domain)

A positive (or negative) weak nodal domain of f on V(G) is a maximal connected induced subgraph of G on vertices $v \in V$ with $f(v) \ge 0$ (or $f(v) \le 0$) that contains at least one nonzero vertex. The number of weak nodal domains of f is denoted by $\mathfrak{W}(f)$.

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Corollary (Fiedler (1975))

If G is connected, then $\mathfrak{W}(\phi_1) = 2$.

Example of Graph Partitioning

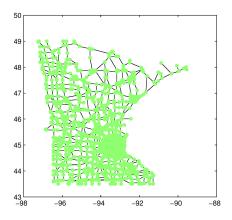


Figure: The MN road network

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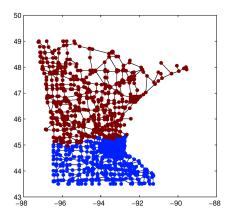


Figure : The MN road network partitioned via the Fiedler vector of L

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Our transforms involve 2 main steps:

- Recursively partition the graph
 - These steps can be performed concurrently, or we can fully partition the graph and then generate a set of bases
- 2 Using the regions on each level of the graph partitioning, generate a set of orthonormal bases for the graph

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Hierarchical Graph Laplacian Eigen Transform (HGLET)

Now we present a novel transform that can be viewed as a generalization of the *block Discrete Cosine Transform*. We refer to this transform as the *Hierarchical Graph Laplacian Eigen Transform (HGLET)*.

The algorithm proceeds as follows...

- Generate an orthonormal basis for the entire graph \Rightarrow Laplacian eigenvectors (Notation is $\phi_{k,l}^j$ with j=0)
- ullet Partition the graph using the Fiedler vector $oldsymbol{\phi}_{k,1}^{j}$
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- For an unweighted path graph, this yields a dictionary of the block DCT-II
- Similar to wavelet packet or local cosine dictionaries in that it generates an overcomplete basis from which we can select a basis useful for the task at hand ⇒ best-basis algorithm, local discriminant basis algorithm, . . .
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Related Work

The following work also proposed a similar strategy to construct a multiscale basis dictionary, i.e., *local cosine dictionary on a graph*:

A. D. Szlam, M. Maggioni, R. R. Coifman, and J. C. Bremer, Jr., "Diffusion-driven multiscale analysis on manifolds and graphs: top-down and bottom-up constructions," in Wavelets XI (M. Papadakis et al. eds.), Proc. SPIE 5914, Paper # 59141D, 2005.

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However, in our opinion, the generalization of the folding/unfolding operations (originally used in the construction of the local cosine transforms on a regular domain) to the graph setting may be harmful. We believe that such operations are not necessary for most tasks in practice. If one needs smoother and overlapping basis vectors, then a better partitioning scheme other than the folding/unfolding operations is called for.

Computational Complexity: HGLET

	Computational	Run Time
	Complexity	for MN ¹
HGLET (redundant)	$O(N^3)$	67 sec

¹Computations performed on a personal laptop (4.00 GB RAM, 2.26 GHz), N = 2640 and nnz (\mathbb{W}) = 6604.

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Generalized Haar-Walsh Transform (GHWT)

HGLET is a generalization of the block DCT, and it generates basis vectors that are *smooth* on their support.

The Generalized Haar-Walsh Transform (GHWT) is a generalization of the classical Haar and Walsh-Hadamard Transforms, and it generates basis vectors that are *piecewise-constant on their support*.

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The algorithm proceeds as follows...

① Generate a full recursive partitioning of the graph \Rightarrow Fiedler vectors

- ② Generate an orthonormal basis for level j_{max} (the finest level) \Rightarrow scaling vectors on the single-node regions
 - ullet As with HGLET, the notation is $oldsymbol{\psi}_{k,l}^J$
- ③ Using the basis for level j_{max} , generate an orthonormal basis for level $j_{\text{max}} 1 \Rightarrow scaling$ and Haar-like vectors
- Repeat... Using the basis for level j, generate an orthonormal basis for level j-1 ⇒ scaling, Haar-like, and Walsh-like vectors
- Select an orthonormal basis from this collection of orthonormal bases

- Generate a full recursive partitioning of the graph ⇒ Fiedler vectors
- ② Generate an orthonormal basis for level j_{max} (the finest level) \Rightarrow scaling vectors on the single-node regions
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$$\left[\begin{array}{c} \pmb{\psi}_{0,0}^{j_{\mathsf{max}}} \end{array}\right] \quad \left[\begin{array}{c} \pmb{\psi}_{1,0}^{j_{\mathsf{max}}} \end{array}\right] \quad \left[\begin{array}{c} \pmb{\psi}_{2,0}^{j_{\mathsf{max}}} \end{array}\right] \quad \left[\begin{array}{c} \pmb{\psi}_{3,0}^{j_{\mathsf{max}}} \end{array}\right] \quad \cdots \quad \left[\begin{array}{c} \pmb{\psi}_{K/\mathsf{max}}^{j_{\mathsf{max}}} \end{array}\right] \quad \left[\begin{array}{c} \pmb{\psi}_{K/\mathsf{max}=1.0}^{j_{\mathsf{max}}} \end{array}\right]$$

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$$\left[\begin{array}{c} \pmb{\psi}_{0,0}^{j_{\mathsf{max}}-1} & \pmb{\psi}_{0,1}^{j_{\mathsf{max}}-1} \end{array} \right] \quad \left[\begin{array}{c} \pmb{\psi}_{1,0}^{j_{\mathsf{max}}-1} & \pmb{\psi}_{1,1}^{j_{\mathsf{max}}-1} \end{array} \right] \quad \cdots \quad \left[\begin{array}{c} \pmb{\psi}_{K^{j_{\mathsf{max}}-1}-1,0}^{j_{\mathsf{max}}-1} & \pmb{\psi}_{K^{j_{\mathsf{max}}-1}-1,1}^{j_{\mathsf{max}}-1} \end{array} \right]$$

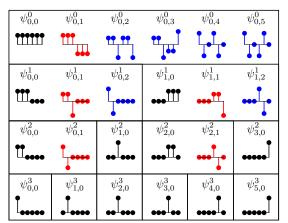
$$\left[\begin{array}{c} \pmb{\psi}_{0,0}^{j_{\mathsf{max}}} \end{array} \right] \quad \left[\begin{array}{c} \pmb{\psi}_{1,0}^{j_{\mathsf{max}}} \end{array} \right] \quad \left[\begin{array}{c} \pmb{\psi}_{2,0}^{j_{\mathsf{max}}} \end{array} \right] \quad \cdots \quad \left[\begin{array}{c} \pmb{\psi}_{K^{j_{\mathsf{max}}-2},0}^{j_{\mathsf{max}}} \end{array} \right] \quad \left[\begin{array}{c} \pmb{\psi}_{K^{j_{\mathsf{max}}-1},0}^{j_{\mathsf{max}}} \end{array} \right]$$

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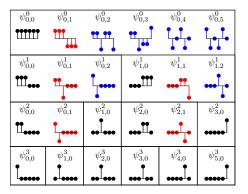


Figure: Default dictionary; i.e., coarse-to-fine

• This reorganization gives us more options for choosing a good basis

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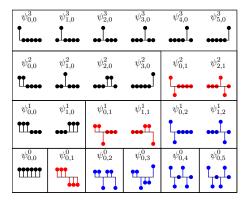


Figure: Reordered & regrouped dictionary; i.e., fine-to-coarse

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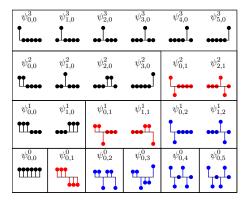
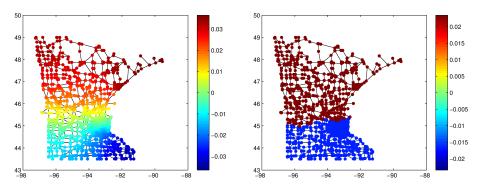


Figure: Reordered & regrouped dictionary; i.e., fine-to-coarse

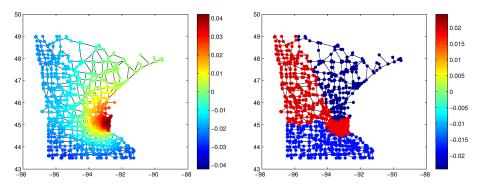
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Here we display some of the basis vectors generated by our HGLET (left) and GHWT (right) schemes on the MN road network. (Note: j=0 is the coarsest scale, j=14 is the finest.)

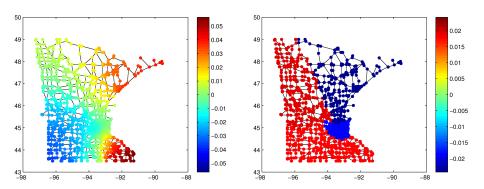
Level j = 0, Region k = 0, l = 1



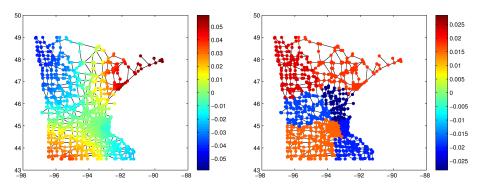
Level
$$j = 0$$
, Region $k = 0$, $l = 2$



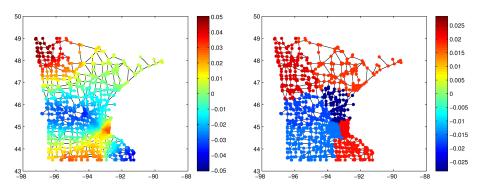
Level
$$j = 0$$
, Region $k = 0$, $l = 3$



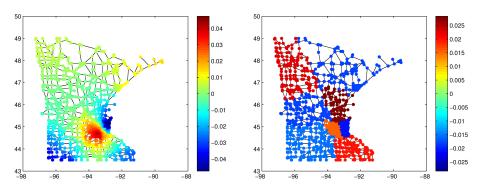
Level
$$j = 0$$
, Region $k = 0$, $l = 4$



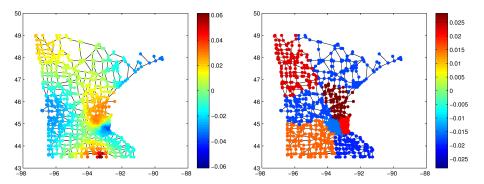
Level
$$j = 0$$
, Region $k = 0$, $l = 5$



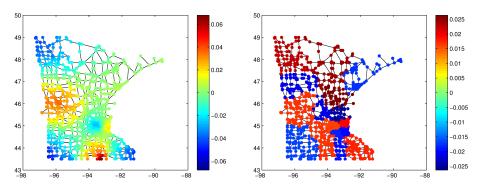
Level
$$j = 0$$
, Region $k = 0$, $l = 6$



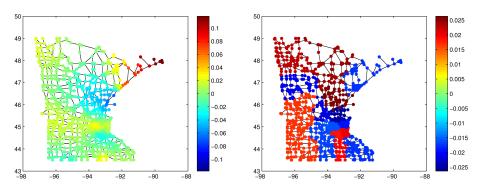
Level
$$j = 0$$
, Region $k = 0$, $l = 7$



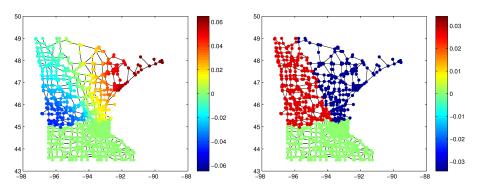
Level
$$j = 0$$
, Region $k = 0$, $l = 8$



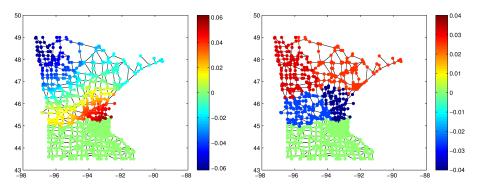
Level
$$j = 0$$
, Region $k = 0$, $l = 9$



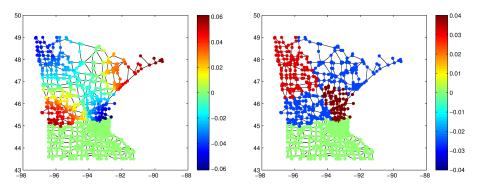
Level
$$j = 1$$
, Region $k = 0$, $l = 1$



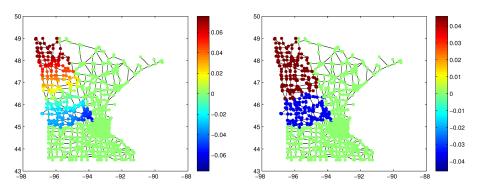
Level
$$j = 1$$
, Region $k = 0$, $l = 2$



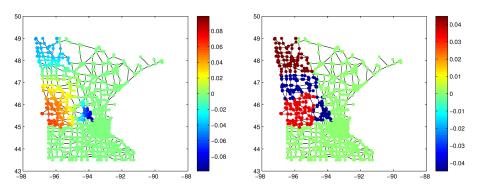
Level
$$j = 1$$
, Region $k = 0$, $l = 3$



Level
$$j = 2$$
, Region $k = 0$, $l = 1$

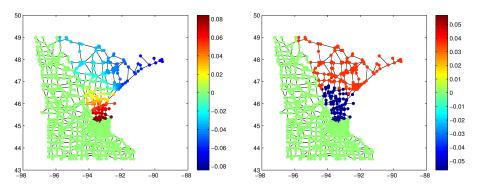


Level
$$j = 2$$
, Region $k = 0$, $l = 2$

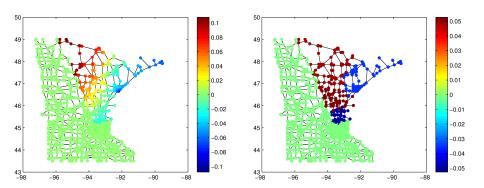


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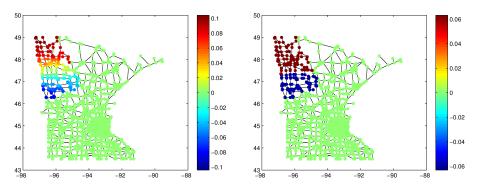
Level j = 2, Region k = 1, l = 1



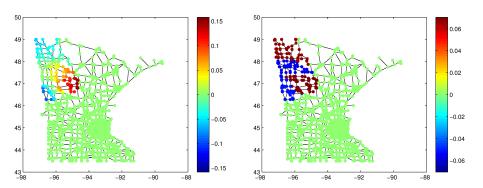
Level
$$j = 2$$
, Region $k = 1$, $l = 2$



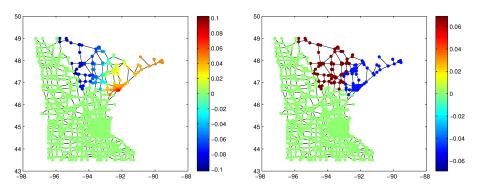
Level
$$j = 3$$
, Region $k = 0$, $l = 1$



Level
$$j = 3$$
, Region $k = 0$, $l = 2$

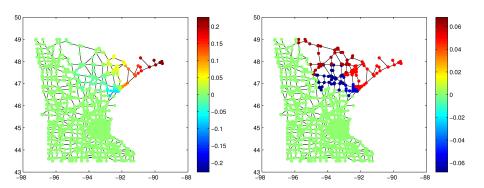


Level
$$j = 3$$
, Region $k = 2$, $l = 1$



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Level j = 3, Region k = 2, l = 2



Computational Complexity: GHWT

	Computational	Run Time
	Complexity	for MN ¹
HGLET (redundant)	$O(N^3)$	67 sec
GHWT (redundant)	$O(N^2)$	10 sec

¹Computations performed on a personal laptop (4.00 GB RAM, 2.26 GHz), N = 2640 and nnz (\mathbb{W}) = 6604.

Related Work

The following articles also discussed the Haar-like transform on graphs and trees, but *not the Walsh-Hadamard transform* on them:

- A. D. Szlam, M. Maggioni, R. R. Coifman, and J. C. Bremer, Jr., "Diffusion-driven multiscale analysis on manifolds and graphs: top-down and bottom-up constructions," in Wavelets XI (M. Papadakis et al. eds.), Proc. SPIE 5914, Paper # 59141D, 2005.
- F. Murtagh, "The Haar wavelet transform of a dendrogram," J. Classification, vol. 24, pp. 3–32, 2007.
- A. Lee, B. Nadler, and L. Wasserman, "Treelets—an adaptive multi-scale basis for sparse unordered data," Ann. Appl. Stat., vol. 2, pp. 435–471, 2008.
- M. Gavish, B. Nadler, and R. Coifman, "Multiscale wavelets on trees, graphs and high dimensional data: Theory and applications to semi supervised learning," in *Proc. 27th Intern. Conf. Machine Learning* (J. Fürnkranz et al. eds.), pp. 367–374, Omnipress, Haifa, 2010.

- Introductory Remarks
- 2 Motivations: Why Graphs?
- Background
 - Basic Graph Theory Terminology
 - Graph Laplacians
 - Graph Partitioning via Spectral Clustering
- Multiscale Basis Dictionaries
 - Hierarchical Graph Laplacian Eigen Transform (HGLET)
 - Generalized Haar-Walsh Transform (GHWT)
- Best-Basis Algorithm for HGLET & GHWT
- 6 Approximation Experiments
- Summary and Future Work

Coifman and Wickerhauser (1992) developed the best-basis algorithm as a means of selecting the basis from a dictionary of wavelet packets that is "best" for approximation/compression.

We generalize this approach, developing and implementing an algorithm for selecting the basis from the dictionary of HGLET / GHWT bases that is "best" for approximation.

As before, we require a cost functional J. For example

$$\mathcal{J}(\mathbf{x}) = \left(\sum_{i=1}^{n} |x_i|^p\right)^{1/p} = \text{norm}(\mathbf{x}, \mathbf{p}) \quad 0$$

• For our approximation experiments in the following pages, we used p=0.1.

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$$\left[\begin{array}{ccccc} \pmb{\phi}^0_{0,0} & \pmb{\phi}^0_{0,1} & \pmb{\phi}^0_{0,2} & \cdots & \pmb{\phi}^0_{0,N^0_0-1} \\ \\ d^0_{0,0} & d^0_{0,1} & d^0_{0,2} & \cdots & d^0_{0,N^0_0-1} \end{array} \right]$$

$$\begin{bmatrix} \boldsymbol{\phi}_{0,0}^1 & \boldsymbol{\phi}_{0,1}^1 & \boldsymbol{\phi}_{0,2}^1 & \cdots & \boldsymbol{\phi}_{0,N_0^1-1}^1 \\ d_{0,0}^1 & d_{0,1}^1 & d_{0,2}^1 & \cdots & d_{0,N_0^1-1}^1 \end{bmatrix} \quad \begin{bmatrix} \boldsymbol{\phi}_{1,0}^1 & \boldsymbol{\phi}_{1,1}^1 & \boldsymbol{\phi}_{1,2}^1 & \cdots & \boldsymbol{\phi}_{1,N_1^1-1}^1 \\ d_{1,0}^1 & d_{1,1}^1 & d_{1,2}^1 & \cdots & d_{1,N_1^1-1}^1 \end{bmatrix}$$

$$\begin{bmatrix} \pmb{\phi}_{0,0}^2 \, \pmb{\phi}_{0,1}^2 \cdots \pmb{\phi}_{0,N_0^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{1,0}^2 \, \pmb{\phi}_{1,1}^2 \cdots \pmb{\phi}_{1,N_1^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{2,0}^2 \, \pmb{\phi}_{2,1}^2 \cdots \pmb{\phi}_{2,N_2^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{3,0}^2 \, \pmb{\phi}_{3,1}^2 \cdots \pmb{\phi}_{3,N_3^2-1}^2 \end{bmatrix} \\ d_{0,0}^2 \, d_{0,1}^2 \cdots \, d_{0,N_0^2-1}^2 \quad d_{1,0}^2 \, d_{1,1}^2 \cdots \, d_{1,N_1^2-1}^2 \quad d_{2,0}^2 \, d_{2,1}^2 \cdots \, d_{2,N_2^2-1}^2 \quad d_{3,0}^2 \, d_{3,1}^2 \cdots \, d_{3,N_3^2-1}^2 \end{bmatrix}$$

$$\begin{bmatrix} \boldsymbol{\phi}_{0,0}^0 & \boldsymbol{\phi}_{0,1}^0 & \boldsymbol{\phi}_{0,2}^0 & \cdots & \boldsymbol{\phi}_{0,N_0^0-1}^0 \end{bmatrix}$$

$$d_{0,0}^0 & d_{0,1}^0 & d_{0,2}^0 & \cdots & d_{0,N_0^0-1}^0$$

$$\begin{bmatrix} \pmb{\phi}_{0,0}^2 \, \pmb{\phi}_{0,1}^2 \cdots \pmb{\phi}_{0,N_0^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{1,0}^2 \, \pmb{\phi}_{1,1}^2 \cdots \pmb{\phi}_{1,N_1^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{2,0}^2 \, \pmb{\phi}_{2,1}^2 \cdots \pmb{\phi}_{2,N_2^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{3,0}^2 \, \pmb{\phi}_{3,1}^2 \cdots \pmb{\phi}_{3,N_3^2-1}^2 \end{bmatrix} \\ d_{0,0}^2 \, d_{0,1}^2 \cdots \, d_{0,N_0^2-1}^2 \quad d_{1,0}^2 \, d_{1,1}^2 \cdots \, d_{1,N_1^2-1}^2 \quad d_{2,0}^2 \, d_{2,1}^2 \cdots \, d_{2,N_2^2-1}^2 \quad d_{3,0}^2 \, d_{3,1}^2 \cdots \, d_{3,N_3^2-1}^2 \end{bmatrix}$$

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$$\begin{bmatrix} \pmb{\phi}_{0,0}^2 \, \pmb{\phi}_{0,1}^2 \cdots \, \pmb{\phi}_{0,N_0^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{1,0}^2 \, \pmb{\phi}_{1,1}^2 \cdots \, \pmb{\phi}_{1,N_1^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{2,0}^2 \, \pmb{\phi}_{2,1}^2 \cdots \, \pmb{\phi}_{2,N_2^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{3,0}^2 \, \pmb{\phi}_{3,1}^2 \cdots \, \pmb{\phi}_{3,N_3^2-1}^2 \end{bmatrix} \\ d_{0,0}^2 \, d_{0,1}^2 \cdots \, d_{0,N_0^2-1}^2 \quad d_{1,0}^2 \, d_{1,1}^2 \cdots \, d_{1,N_1^2-1}^2 \quad d_{2,0}^2 \, d_{2,1}^2 \cdots \, d_{2,N_2^2-1}^2 \quad d_{3,0}^2 \, d_{3,1}^2 \cdots \, d_{3,N_3^2-1}^2 \end{bmatrix}$$

$$\begin{bmatrix} \boldsymbol{\phi}_{0,0}^{0} & \boldsymbol{\phi}_{0,1}^{0} & \boldsymbol{\phi}_{0,2}^{0} & \cdots & \boldsymbol{\phi}_{0,N_{0}^{0}-1}^{0} \end{bmatrix}$$
$$d_{0,0}^{0} & d_{0,1}^{0} & d_{0,2}^{0} & \cdots & d_{0,N_{0}^{0}-1}^{0} \end{bmatrix}$$

$$\begin{bmatrix} \boldsymbol{\phi}_{2,0}^2 \, \boldsymbol{\phi}_{2,1}^2 \cdots \boldsymbol{\phi}_{2,N_2^2-1}^2 \\ d_{2,0}^2 \, d_{2,1}^2 \cdots \, d_{2,N_2^2-1}^2 \end{bmatrix} \quad \begin{bmatrix} \boldsymbol{\phi}_{3,0}^2 \, \boldsymbol{\phi}_{3,1}^2 \cdots \boldsymbol{\phi}_{3,N_3^2-1}^2 \\ d_{2,0}^2 \, d_{2,1}^2 \cdots \, d_{2,N_2^2-1}^2 \\ \end{bmatrix}$$

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$$\begin{bmatrix} \pmb{\phi}_{2,0}^2 \, \pmb{\phi}_{2,1}^2 \cdots \, \pmb{\phi}_{2,N_2^2-1}^2 \\\\ d_{2,0}^2 \, d_{2,1}^2 \cdots \, d_{2,N_2^2-1}^2 \\\\ \end{bmatrix} \quad \begin{bmatrix} \pmb{\phi}_{3,0}^2 \, \pmb{\phi}_{3,1}^2 \cdots \, \pmb{\phi}_{3,N_3^2-1}^2 \\\\ d_{3,0}^2 \, d_{3,1}^2 \cdots \, d_{3,N_3^2-1}^2 \\\\ \end{bmatrix}$$

$$\begin{bmatrix} & \phi_{0,0}^0 & \phi_{0,1}^0 & \phi_{0,2}^0 & \cdots & \phi_{0,N_0^0-1}^0 \\ & d_{0,0}^0 & d_{0,1}^0 & d_{0,2}^0 & \cdots & d_{0,N_0^0-1}^0 \end{bmatrix}$$

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According to cost functional \mathcal{J} , this is the best basis for approximation.

saito@math.ucdavis.edu (UC Davis)

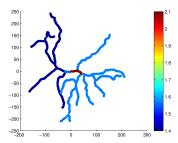
$$\left[\begin{array}{ccccc} \pmb{\phi}_{0,0}^1 & \pmb{\phi}_{0,1}^1 & \pmb{\phi}_{0,2}^1 & \cdots & \pmb{\phi}_{0,N_0^1-1}^1 \\ d_{0,0}^1 & d_{0,1}^1 & d_{0,2}^1 & \cdots & d_{0,N_{\alpha}^1-1}^1 \end{array} \right]$$

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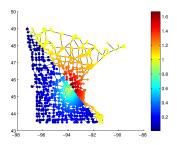
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 With the GHWT bases, we run the best-basis algorithm on both the default (coarse-to-fine) dictionary and the reorganized (fine-to-coarse) dictionary and then compare the cost of the 2 bases to determine the best-basis.

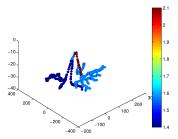
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- 6 Approximation Experiments
- Summary and Future Work



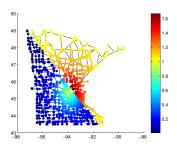
(a) Thickness data on a dendritic tree



(b) A mutilated Gaussian on the MN road network

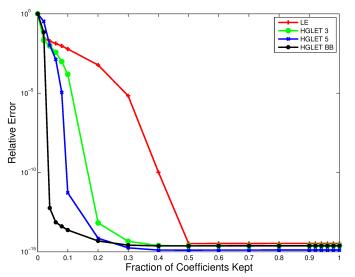


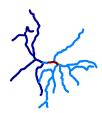
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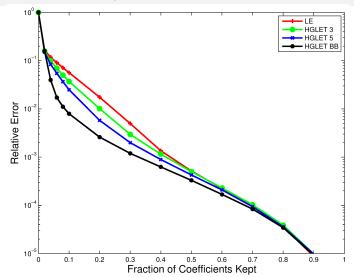
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HGLET on Dendrite (weights = inv. Euclidean dist.)



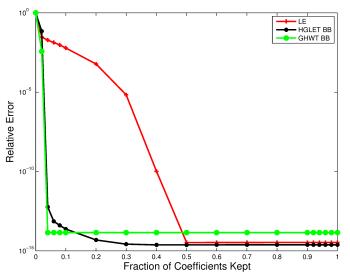


HGLET on MN Mutilated Gaussian (weights = inv. Euclidean dist.)



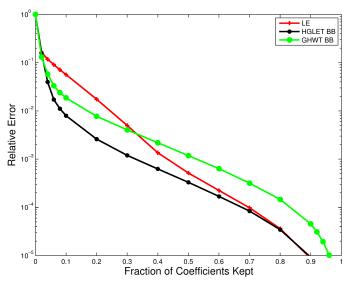


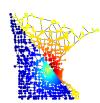
GHWT vs. HGLET on Dendrite





GHWT vs. HGLET on MN Mutilated Gaussian





- From the HGLET plots, we see that HGLET best-basis > HGLET Level 5 > HGLET Level 3 > Laplacian eigenvectors (HGLET Level 0)
- The HGLET best-basis performs the best on the MN Mutilated Gaussian dataset while the GHWT best-basis outperformed the others on the Dendrite dataset
- These performances make a strong case for using localized basis vectors on multiple scales
- Also, these indicate that the smoothness of the basis vectors matters depending on the smoothness inherent in data

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- We developed multiscale basis dictionaries on graphs and networks: HGLET and GHWT. We also developed a corresponding best-basis algorithm.
- The HGLET is a direct generalization of *Hierarchical Block Discrete*
- The GHWT is a generalization of the *Haar Transform* and the
- Both of these transforms allow us to choose an orthonormal basis
- They may also be useful for regularly-sampled signals, e.g., can deal
- Developing a true generalization of smoother wavelet and wavelet

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Future Work

- Perform classification experiments and compare the results using HGLET and GHWT.
- Explore other methods for graph partitioning:
 - Allow for splitting of a region into an arbitrary number of subregions;
 - Consider a bottom-up clustering method, rather than a top-down partitioning method;
 - Incorporate the diffuse interface model and the minimization of the Ginsburg-Landau functional proposed by Bertozzi and Flenner (2012).

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References & Acknowledgments

- http://www.math.ucdavis.edu/~saito/courses/HarmGraph/ contains my course slides and useful information on "Harmonic Analysis on Graphs and Networks"
- Also visit http://www.math.ucdavis.edu/~saito/publications/ for various related publications including:
 - N. Saito: "Data analysis and representation using eigenfunctions of Laplacian on a general domain," Applied & Computational Harmonic Analysis, vol. 25, no. 1, pp. 68-97, 2008.
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