

Notes on Spectral Clustering

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Talk outline

- Spectral Clustering
 - Distances and similarity graphs
 - Graph Laplacians and their properties
 - Spectral clustering algorithms
 - SC under the hood

Similarity graph

- The objective of a clustering algorithm is partitioning data into groups such that:
 - Points in the **same** group are **similar**
 - Points in **different** groups are **dissimilar**
- **Similarity graph** $G=(V,E)$ *(undirected graph)*
 - Vertices v_i and v_j are connected by a **weighted** edge if their similarity is above a given threshold
 - GOAL: find a partition of the graph such that:
 - edges **within** a group have **high weights**
 - edges **across** different groups have **low weights**

Weighted adjacency matrix

- Let $G(V,E)$ be an undirected graph with vertex set $V = \{v_1, \dots, v_n\}$
- **Weighted adjacency matrix** $W = (w_{ij})_{i,j=1, \dots, n}$
 - $w_{ij} \geq 0$ is the weight of the edge between v_i and v_j
 - $w_{ij} = 0$ means that v_i and v_j are not connected by an edge
 - $w_{ij} = w_{ji}$
- **Degree of a vertex** $v_i \in V$: $d_i = \sum_{j=1..n} w_{ij}$
- **Degree matrix** $D = \text{diag}(d_1, \dots, d_n)$

Different similarity graphs

- ε -neighborhood
 - Connect all points whose pairwise distance is less than ε
- k -nearest neighbors
 - if $v_i \in knn(v_j)$ **OR** $v_j \in knn(v_i)$
 - if $v_i \in knn(v_j)$ **AND** $v_j \in knn(v_i)$ (*mutual knn*)
 - after connecting edges, use similarity as weight
- fully connected
 - all points with similarity $s_{ij} > 0$ are connected
 - To control neighborhoods to be *local*, use a similarity function like the **Gaussian**: $s(x_i, x_j) = \exp(-\|x_i - x_j\|^2 / (2\sigma^2))$

Graph Laplacians

- **Graph Laplacian:**
 - $L = D - W$ (symmetric and positive semi-definite)
- **Properties**
 - Smallest eigenvalue $\lambda_1 = 0$ with eigenvector = $\mathbb{1}$
 - n non-negative, real-valued eigenvalues $0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$
 - the multiplicity k of the eigenvalue 0 of L equals the number of connected components A_1, \dots, A_k in the graph

Spectral Clustering algorithm (1)

Spectral Clustering algorithm

Input: Similarity matrix $S \in \mathbb{R}^{n \times n}$, number k of clusters to construct.

1. Construct a similarity graph as previously described. Let W be its weighted adjacency matrix.
2. Compute the unnormalized Laplacian L
3. Compute the first k eigenvectors u_1, \dots, u_k of L
4. Let $U \in \mathbb{R}^{n \times k}$ be the matrix containing the vectors u_1, \dots, u_k as columns
5. For $i=1, \dots, n$ let $y_i \in \mathbb{R}^k$ be the vector corresponding to the i -th row of U
6. Cluster the points $(y_i)_{i=1, \dots, n}$ in \mathbb{R}^k with the k-means algorithm into clusters C_1, \dots, C_k .

Output: Clusters A_1, \dots, A_k with $A_i = \{j \mid y_j \in C_i\}$.

Normalized Graph Laplacians

- **Normalized graph Laplacians**

- *Symmetric*: $L_{sym} = D^{-1/2} L D^{-1/2} = I - D^{-1/2} W D^{-1/2}$

- *Random Walk*: $L_{rw} = D^{-1} L = I - D^{-1} W$

- **Properties**

- λ is an eigenvalue of L_{rw} with eigenvector u iff λ is an eigenvalue of L_{sym} with eigenvector $w = D^{1/2} u$

- λ is an eigenvalue of L_{rw} with eigenvector u iff λ and u solve the **generalized eigenproblem** $Lu = \lambda Du$

Normalized Graph Laplacians

- **Normalized graph Laplacians**

- *Symmetric:* $L_{sym} = D^{-1/2}LD^{-1/2} = I - D^{-1/2}WD^{-1/2}$

- *Random Walk:* $L_{rw} = D^{-1}L = I - D^{-1}W$

- **Properties (follow)**

- 0 is an eigenvalue of L_{rw} with $\mathbb{1}$ as eigenvector, and an eigenvalue of L_{sym} with eigenvector $D^{1/2}\mathbb{1}$.

- L_{sym} and L_{rw} are positive semi-definite and have n non-negative, real-valued eigenvalues $0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$

- the multiplicity k of the eigenvalue 0 of both L_{sym} and L_{rw} equals the number of connected components A_1, \dots, A_k

Spectral Clustering algorithm (2)

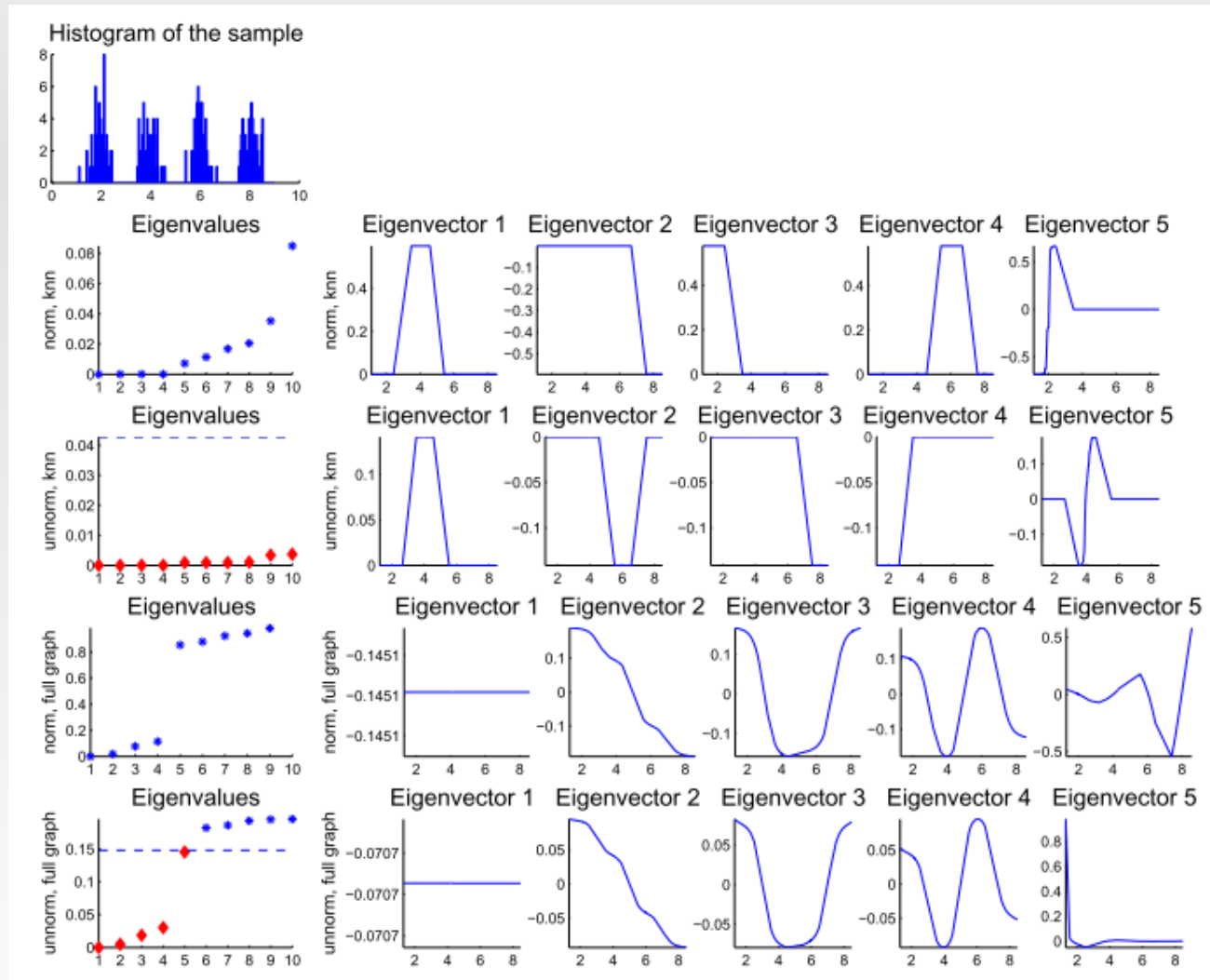
Normalized Spectral Clustering

Input: Similarity matrix $S \in \mathbb{R}^{n \times n}$, number k of clusters to construct.

- L_{rw} :
 3. Compute the first k **generalized** eigenvectors u_1, \dots, u_k of the generalized eigenproblem $Lu = \lambda Du$
- L_{sym} :
 2. Compute the **normalized** Laplacian L_{sym}
 3. Compute the first k eigenvectors u_1, \dots, u_k of L_{sym}
 4. normalize the eigenvectors

Output: Clusters A_1, \dots, A_k with $A_i = \{j \mid y_j \in C_i\}$.

A spectral clustering example



Under the hood

- *0-eigenvalues* in the ideal case
- *parameters* are crucial:
 - k in k nearest neighbors
 - σ in Gaussian kernel
 - k (another one!) in k -means

Random Walk point of view

- Random walk: stochastic process which randomly jumps from one vertex to another
 - Clustering: finding a partition such that a random walk stays long within a cluster and seldom jumps between clusters
- Transition probability $p_{ij} := w_{ij} / d_i$
- Transition matrix: $P = D^{-1}W \quad \Rightarrow \quad L_{rw} = I - P$
 - λ is an eigenvalue of L_{rw} with eigenvector u iff $1-\lambda$ is an eigenvalue of P with eigenvector u

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References

- Von Luxburg, U. (2007). *A tutorial on spectral clustering*. *Statistics and Computing*, 17(4), 395-416. Springer.

Thank you!

Thanks for your attention!

Questions?