Research Report - Spectral Clustering

Sourabh Antani supervisor Dr. Yunkai Zhou

Mathematics Department, Southern Methodist University

June 3, 2021

Below are some examples of clustering algorithms used in practice, aside from spectral clustering.

Euclidean distance and spatial density based:

- Euclidean distance and spatial density based:
 - k-means [Macqueen, 1967] [Lloyd 1982]

- Euclidean distance and spatial density based:
 - k-means [Macqueen, 1967] [Lloyd 1982]
 - BIRCH (Balanced Iterative Reducing and Clustering using Hierarchies) [Zhang et al. 1996]

- Euclidean distance and spatial density based:
 - k-means [Macqueen, 1967] [Lloyd 1982]
 - BIRCH (Balanced Iterative Reducing and Clustering using Hierarchies) [Zhang et al. 1996]
 - Mean-shift [Cheng 1995]

- Euclidean distance and spatial density based:
 - k-means [Macqueen, 1967] [Lloyd 1982]
 - BIRCH (Balanced Iterative Reducing and Clustering using Hierarchies) [Zhang et al. 1996]
 - Mean-shift [Cheng 1995]
 - DBSCAN [Ester et al. 1996]

- Euclidean distance and spatial density based:
 - k-means [Macqueen, 1967] [Lloyd 1982]
 - BIRCH (Balanced Iterative Reducing and Clustering using Hierarchies) [Zhang et al. 1996]
 - Mean-shift [Cheng 1995]
 - DBSCAN [Ester et al. 1996]
- Linear Algebra based

- Euclidean distance and spatial density based:
 - k-means [Macqueen, 1967] [Lloyd 1982]
 - BIRCH (Balanced Iterative Reducing and Clustering using Hierarchies) [Zhang et al. 1996]
 - Mean-shift [Cheng 1995]
 - DBSCAN [Ester et al. 1996]
- Linear Algebra based
 - Non-Negative Matrix Factorization [Lawton et al. 1971] [Paatero et al. 1991] [Ding et al. 2005]

- Euclidean distance and spatial density based:
 - k-means [Macqueen, 1967] [Lloyd 1982]
 - BIRCH (Balanced Iterative Reducing and Clustering using Hierarchies) [Zhang et al. 1996]
 - Mean-shift [Cheng 1995]
 - DBSCAN [Ester et al. 1996]
- Linear Algebra based
 - Non-Negative Matrix Factorization [Lawton et al. 1971] [Paatero et al. 1991] [Ding et al. 2005]
 - PCA Based clustering [Zhang et al. 2018]



Graph Laplacian - Construction

• Given an undirected graph, $\mathbf{G} = (\mathbf{V}, \mathbf{E})$ with set of vertices $\mathbf{V} = \{\mathbf{v}_1, \dots, \mathbf{v}_n\}$ and \mathbf{E} set of weighted edges. The weight of the edge between v_i and v_j , denoted by w_{ij} is assigned based on the 'similarity' between the data-points v_i and v_j . For example, if the similarity function chosen is a Gaussian kernel and the data that these vertices represent are scalars x_i and x_j , then $w_{ij} = e^{-|x_i - x_j|^2/2}$

Graph Laplacian - Construction

- Given an undirected graph, $\mathbf{G} = (\mathbf{V}, \mathbf{E})$ with set of vertices $\mathbf{V} = \{\mathbf{v}_1, \dots, \mathbf{v}_n\}$ and \mathbf{E} set of weighted edges. The weight of the edge between v_i and v_j , denoted by w_{ij} is assigned based on the 'similarity' between the data-points v_i and v_j . For example, if the similarity function chosen is a Gaussian kernel and the data that these vertices represent are scalars x_i and x_j , then $w_{ij} = e^{-|x_i x_j|^2/2}$
- Construct the Weighted adjacency matrix $\mathbf{W} = (\mathbf{w_{ij}})_{i,j=1,\dots,n}$ and Degree matrix $\mathbf{D} = diag(d_1,\dots,d_n)$ where $\mathbf{d_i} = \sum_{i=1}^n \mathbf{w_{ij}}$ is called Degree of a vertex v_i .

Graph Laplacian - Construction

- Given an undirected graph, $\mathbf{G} = (\mathbf{V}, \mathbf{E})$ with set of vertices $\mathbf{V} = \{\mathbf{v}_1, \dots, \mathbf{v}_n\}$ and \mathbf{E} set of weighted edges. The weight of the edge between v_i and v_j , denoted by w_{ij} is assigned based on the 'similarity' between the data-points v_i and v_j . For example, if the similarity function chosen is a Gaussian kernel and the data that these vertices represent are scalars x_i and x_j , then $w_{ij} = e^{-|x_i x_j|^2/2}$
- Construct the Weighted adjacency matrix $\mathbf{W}=(\mathbf{w_{ij}})_{i,j=1,\dots,n}$ and Degree matrix $\mathbf{D}=diag(d_1,\dots,d_n)$ where $\mathbf{d_i}=\sum_{j=1}^n\mathbf{w_{ij}}$ is called Degree of a vertex v_i .
- Graph Laplacian is defined as L = D W

Graph Laplacian - Construction

- Given an undirected graph, $\mathbf{G} = (\mathbf{V}, \mathbf{E})$ with set of vertices $\mathbf{V} = \{\mathbf{v}_1, \dots, \mathbf{v}_n\}$ and \mathbf{E} set of weighted edges. The weight of the edge between v_i and v_j , denoted by w_{ij} is assigned based on the 'similarity' between the data-points v_i and v_j . For example, if the similarity function chosen is a Gaussian kernel and the data that these vertices represent are scalars x_i and x_j , then $w_{ij} = e^{-|x_i x_j|^2/2}$
- Construct the Weighted adjacency matrix $\mathbf{W} = (\mathbf{w_{ij}})_{i,j=1,\dots,n}$ and Degree matrix $\mathbf{D} = diag(d_1,\dots,d_n)$ where $\mathbf{d_i} = \sum_{j=1}^n \mathbf{w_{ij}}$ is called Degree of a vertex v_i .
- Graph Laplacian is defined as L = D W
- A fully connected graph can be made sparse by converting it into a k-nearest neighbor or an ϵ -neighborhood graph.

- Unnormalized Graph Laplacian [Mohar, 1991], [Mohar, 1997]
 - L = D W (Symmetric)

- Unnormalized Graph Laplacian [Mohar, 1991], [Mohar, 1997]
 - L = D W (Symmetric)
- Normalized Symmetric Graph Laplacian [Chung, 1997]
 - $L_{sym} = D^{-1/2}LD^{-1/2} = I D^{-1/2}WD^{-1/2}$ (Symmetric)
 - if (λ, u) is an eigenpair of L, then $(\lambda, D^{1/2}u)$ is eigenpair of L_{sym} .

- Unnormalized Graph Laplacian [Mohar, 1991], [Mohar, 1997]
 - L = D W (Symmetric)
- Normalized Symmetric Graph Laplacian [Chung, 1997]
 - $L_{sym} = D^{-1/2}LD^{-1/2} = I D^{-1/2}WD^{-1/2}$ (Symmetric)
 - if (λ, u) is an eigenpair of L, then $(\lambda, D^{1/2}u)$ is eigenpair of L_{sym} .
- Normalized Random Walk Graph Laplacian [Chung 1997]
 - $L_{rw} = D^{-1}L = I D^{-1}W$ (Not-Symmetric)
 - if (λ, u) is an eigenpair of L, then $(\lambda, D^{1/2}u)$ is eigenpair of L_{rw} .

- Unnormalized Graph Laplacian [Mohar, 1991], [Mohar, 1997]
 - L = D W (Symmetric)
- Normalized Symmetric Graph Laplacian [Chung, 1997]
 - $L_{sym} = D^{-1/2}LD^{-1/2} = I D^{-1/2}WD^{-1/2}$ (Symmetric)
 - if (λ,u) is an eigenpair of L , then $(\lambda,D^{1/2}u)$ is eigenpair of L_{sym} .
- Normalized Random Walk Graph Laplacian [Chung 1997]
 - $L_{rw} = D^{-1}L = I D^{-1}W$ (Not-Symmetric)
 - if (λ, u) is an eigenpair of L, then $(\lambda, D^{1/2}u)$ is eigenpair of L_{rw} .
- All Laplacians are Positive-semidefinte, hence have non-Negative Real-Valued Eigenvalues

- Unnormalized Graph Laplacian [Mohar, 1991], [Mohar, 1997]
 - L = D W (Symmetric)
- Normalized Symmetric Graph Laplacian [Chung, 1997]
 - $L_{sym} = D^{-1/2}LD^{-1/2} = I D^{-1/2}WD^{-1/2}$ (Symmetric)
 - if (λ, u) is an eigenpair of L, then $(\lambda, D^{1/2}u)$ is eigenpair of L_{sym} .
- Normalized Random Walk Graph Laplacian [Chung 1997]
 - $L_{rw} = D^{-1}L = I D^{-1}W$ (Not-Symmetric)
 - if (λ, u) is an eigenpair of L, then $(\lambda, D^{1/2}u)$ is eigenpair of L_{rw} .
- All Laplacians are Positive-semidefinte, hence have non-Negative Real-Valued Eigenvalues
- 0 is an eigenvalue with multiplicity equal to number of connected components

Graph Cuts

• The goal of clustering is to maximize the similarity within cluster and minimize the similarity between clusters.

Graph Cuts

- The goal of clustering is to maximize the similarity within cluster and minimize the similarity between clusters.
- Minimizing the between cluster similarity equals to minimizing the objective function

$$\operatorname{\mathsf{cut}}(A_1,\ldots,A_k) := \frac{1}{2} \sum_{i=1}^k \mathit{W}(A_i,\bar{A}_i), \quad \text{where } \mathit{W}(A,B) = \sum_{i \in A, j \in B} \mathit{w}_{ij}, \bar{A} = \mathit{V} \setminus A$$

Graph Cuts

- The goal of clustering is to maximize the similarity within cluster and minimize the similarity between clusters.
- Minimizing the between cluster similarity equals to minimizing the objective function

$$\operatorname{cut}(A_1,\ldots,A_k) := \frac{1}{2} \sum_{i=1}^k W(A_i,\bar{A}_i), \quad \text{where } W(A,B) = \sum_{i \in A,j \in B} w_{ij}, \bar{A} = V \setminus A$$

To maximize the within-cluster similarity, we define two objective functions

$$RatioCut(A_{1},...,A_{k}) = \frac{1}{2} \sum_{i=1}^{k} \frac{W(A_{i},\bar{A}_{i})}{|A_{i}|} = \sum_{i=1}^{k} \frac{cut(A_{i},\bar{A}_{i})}{|A_{i}|} [Hagen, Kahng, 1992]$$

$$Ncut(A_{1},...,A_{k}) = \frac{1}{2} \sum_{i=1}^{k} \frac{W(A_{i},\bar{A}_{i})}{vol(A_{i})} = \sum_{i=1}^{k} \frac{cut(A_{i},\bar{A}_{i})}{vol(A_{i})} [Shi, Malik, 2000]$$

Algorithms - Unnormalized Spectral Clustering - Hagen & Kahng (1992)

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

Construct a similarity graph and its weighted adjacency matrix *W* according to the chosen similarity function

Compute the unnormalized Laplacian L

Compute the first k eigenvectors u_1, \ldots, u_k of L

Let $U \in \mathbb{R}^{n \times k}$ be the matrix containing the vectors u_1, \dots, u_k as columns

For i = 1, ..., n, let $y_i \in \mathbb{R}^k$ be the vector corresponding to the i^{th} row of U

Cluster the points (y_i) , i = 1, ..., n in \mathbb{R}^k with the k-means algorithm into clusters $C_1, ..., C_k$

end

Output: Clusters A_1, \ldots, A_k with $A_i = \{x_j | y_j \in C_i\}$

Algorithms - Normalized Spectral Clustering - Shi & Malik (2000)

This algorithm uses generalized eigenvectors of L, which are the eigenvectors of normalized random-walk Laplacian $L_{rw} = D^{-1}L$

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

Construct a similarity graph and its weighted adjacency matrix *W* according to the chosen similarity function

Compute the unnormalized Laplacian L.

Compute the first k generalized eigenvectors u_1, \ldots, u_k of generalized eigenproblem $Lu = \lambda Du$.

Let $U \in \mathbb{R}^{n \times k}$ be the matrix containing the vectors u_1, \dots, u_k as columns.

For i = 1, ..., n, let $y_i \in \mathbb{R}^k$ be the vector corresponding to the i^{th} row of U.

Cluster the points (y_i) , i = 1, ..., n in \mathbb{R}^k with the k-means algorithm into clusters $C_1, ..., C_k$.

end

Output: Clusters A_1, \ldots, A_k with $A_i = \{x_i | y_i \in C_i\}$



Algorithms - Normalized Spectral Clustering - Ng, Jordan & Weiss (2002)

This algorithm uses the eigenvectors of normalized symmetric Laplacian $L_{sym} = D^{-1/2}LD^{-1/2}$. Note that if $D^{1/2}u$ is an eigenvector of L_{sym} if u is an eigenvector of L hence an additional normalization step is needed.

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

Construct a similarity graph and its weighted adjacency matrix *W* according to the chosen similarity function

Compute the normalized Laplacian L_{sym}

Compute the first k eigenvectors u_1, \ldots, u_k of L_{sym}

Let $U \in \mathbb{R}^{n \times k}$ be the matrix containing the vectors u_1, \dots, u_k as columns

Form the matrix $T \in \mathbb{R}^{n \times k}$ from U by normalizing the norm to 1,

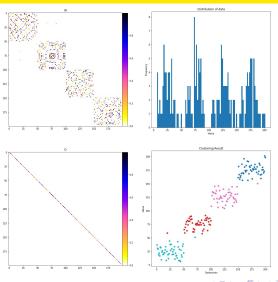
$$t_{ij} = u_{ij}/(\sum_k u_{ik}^2)^{1/2}$$

For $i=1,\ldots,n$, let $y_i\in\mathbb{R}^k$ be the vector corresponding to the i^{th} row of U Cluster the points $(y_i), i=1,\ldots,n$ in \mathbb{R}^k with the k-means algorithm into clusters C_1,\ldots,C_k

end

Output: Clusters A_1, \ldots, A_k with $A_i = \{x_i | y_i \in C_i\}$

Mixture of 4 Gaussians in \mathbb{R}^1 . Similarity function: $e^{-|x_i-x_j|^2/2}$



9/17

1351 Points in \mathbb{R}^2 . Similarity function: $||x - y||_2^{-1}$

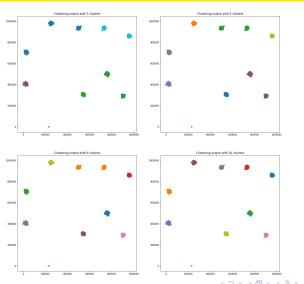
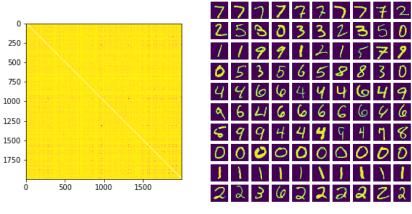


Image Classification. Similarity function: $||\mathbf{v_i} \circ \mathbf{v_j}||_1$

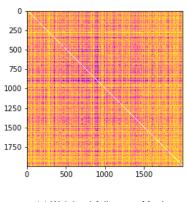


(a) Weighted Adjacency Matrix

(b) Clustering Result

Figure: 2000 images from MNIST data-set were clustered into 10 clusters. (b) shows first 10 images in each cluster with one row per cluster

Image Classification. Similarity function: $||\mathbf{v_i} - \mathbf{v_j}||_2^{-1}$



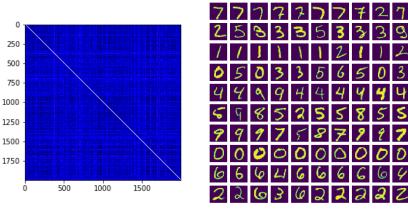




(b) Clustering Result

Figure: 2000 images from MNIST data-set were clustered into 10 clusters. (b) shows first 10 images in each cluster with one row per cluster

Image Classification. Similarity function: number of pixels that either 0 or non-zero in both $v_i \& v_j$



(a) Weighted Adjacency Matrix

(b) Clustering Result

Figure: 2000 images from MNIST data-set were clustered into 10 clusters. (b) shows first 10 images in each cluster with one row per cluster

June 3, 2021

Image Classification. Doubling the number of clusters to capture handwriting style differences

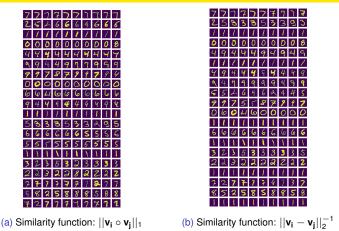


Figure: 2000 images from MNIST data-set were clustered into 20 clusters, showing first 10 images in each cluster with one row per cluster.

June 3, 2021

[Tremblay et al. 2016]

Challenges with Spectral Clustering: Eigenvector computation, Cold-restart, expensive *k*-means due to high dimensionality.

15/17

[Tremblay et al. 2016]

Challenges with Spectral Clustering: Eigenvector computation, Cold-restart, expensive k-means due to high dimensionality.

- Sample $\mathcal{O}(\log(k))$ randomly filtered signals on the graph to serve as feature vectors instead of eigenvectors
 - Estimate kth eigenvalue per [Napoli 2013]
 - Approximate the action of a low-pass filter H that selects the first k eigenvectors on the graph Laplacian
 - Generate $\mathcal{O}(\log(k))$ Gaussian signals, filtered by H to generate feature vectors \mathbf{f}_i

[Tremblay et al. 2016]

Challenges with Spectral Clustering: Eigenvector computation, Cold-restart, expensive *k*-means due to high dimensionality.

- Sample $\mathcal{O}(\log(k))$ randomly filtered signals on the graph to serve as feature vectors instead of eigenvectors
 - Estimate kth eigenvalue per [Napoli 2013]
 - Approximate the action of a low-pass filter H that selects the first k eigenvectors on the graph Laplacian
 - Generate $\mathcal{O}(\log(k))$ Gaussian signals, filtered by H to generate feature vectors \mathbf{f}_i
- Clustering random subset of $\mathcal{O}(k \log(k))$ nodes using random feature vectors

[Tremblay et al. 2016]

Challenges with Spectral Clustering: Eigenvector computation, Cold-restart, expensive *k*-means due to high dimensionality.

- Sample $\mathcal{O}(\log(k))$ randomly filtered signals on the graph to serve as feature vectors instead of eigenvectors
 - Estimate kth eigenvalue per [Napoli 2013]
 - Approximate the action of a low-pass filter H that selects the first k eigenvectors on the graph Laplacian
 - Generate $\mathcal{O}(\log(k))$ Gaussian signals, filtered by H to generate feature vectors \mathbf{f}_i
- Clustering random subset of $\mathcal{O}(k \log(k))$ nodes using random feature vectors
- Infer the cluster label of all N nodes by interpolating.

Future Work

- Accelerated Eigensolvers
 - Accelerating Lanczos
 - Chebychev-Davidson Methods [Zhou & Saad 2010] [Z. Wang, 2015]
 - Taking advantage of structure
- Detecting number of clusters adaptively.
- Hot-restart information from the pass with k clusters to bootstrap clustering process with k + l clusters
- Dimensionality reduction for k-means

Thank You!