

Introduction to Machine Learning

Spectral Clustering

Mingchen Gao

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1 Spectral Clustering

- An alternate approach to clustering
- Let the data be a set of N points

$$\mathbf{X} = \mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_N$$

- Let \mathbf{S} be a $N \times N$ **similarity matrix**

$$S_{ij} = \text{sim}(\mathbf{x}_i, \mathbf{x}_j)$$

- $\text{sim}(\cdot)$ is a similarity function
- Construct a weighted undirected graph from \mathbf{S} with adjacency matrix,
 \mathbf{W}

$$W_{ij} = \begin{cases} \text{sim}(\mathbf{x}_i, \mathbf{x}_j) & \text{if } \mathbf{x}_i \text{ is nearest neighbor of } \mathbf{x}_j \\ 0 & \text{otherwise} \end{cases}$$

- Can use more than 1 nearest neighbors to construct the graph
- Clustering \mathbf{X} into K clusters is equivalent to finding K cuts in the graph \mathbf{W}
 - A_1, A_2, \dots, A_K
- Possible objective function

$$cut(A_1, A_2, \dots, A_K) \triangleq \frac{1}{2} \sum_{k=1}^K W(A_k, \bar{A}_k)$$

- where \bar{A}_k denotes the nodes in the graph which are **not in** A_k and

$$W(A, B) \triangleq \sum_{i \in A, j \in B} W_{ij}$$

Normalized Min-cut Problem

$$normcut(A_1, A_2, \dots, A_K) \triangleq \frac{1}{2} \sum_{k=1}^K \frac{W(A_k, \bar{A}_k)}{vol(A_k)}$$

where $vol(A) \triangleq \sum_{i \in A} d_i$, d_i is the weighted degree of the node i

- Equivalent to solving a 0-1 knapsack problem
- Find N binary vectors, \mathbf{c}_i of length K such that $c_{ik} = 1$ only if point i belongs to cluster k
- If we relax constraints to allow c_{ik} to be real-valued, the problem becomes an eigenvector problem
 - Hence the name: **spectral clustering**

1.1 Graph Laplacian

$$\mathbf{L} \triangleq \mathbf{D} - \mathbf{W}$$

- \mathbf{D} is a diagonal matrix with degree of corresponding node as the diagonal value

Properties of Laplacian Matrix

1. Each row sums to 0
2. $\mathbf{1}$ is an eigen vector with eigen value equal to 0
 - This means that $\mathbf{L}\mathbf{1} = \mathbf{0}\mathbf{1}$.
3. Symmetric and positive semi-definite
4. Has N non-negative real-valued eigenvalues
5. If the graph (\mathbf{W}) has K connected components, then \mathbf{L} has K eigenvectors spanned by $\mathbf{1}_{\mathbf{A}_1}, \dots, \mathbf{1}_{\mathbf{A}_K}$ with 0 eigenvalue.

To see why \mathbf{L} is positive semi-definite:

$$\begin{aligned}\mathbf{xLx}^\top &= \mathbf{xDx}^\top - \mathbf{xWx}^\top \\ &= \sum_i d_i x_i^2 - \sum_i \sum_j x_i x_j w_{ij} \\ &= \frac{1}{2} \left(\sum_i d_i x_i^2 - 2 \sum_i \sum_j x_i x_j w_{ij} + \sum_j d_j x_j^2 \right) \\ &= \frac{1}{2} \sum_i \sum_j w_{ij} (x_i - x_j)^2\end{aligned}$$

which is $\geq 0 \forall \mathbf{x}$.

1.2 Spectral Clustering Algorithm

Observation

- In practice, \mathbf{W} might not have K exactly isolated connected components
- By *perturbation theory*, the smallest eigenvectors of \mathbf{L} will be close to the ideal indicator functions

Algorithm

- Compute first (smallest) K eigen vectors of \mathbf{L}
- Let \mathbf{U} be the $N \times K$ matrix with eigenvectors as the columns
- Perform kMeans clustering on the rows of \mathbf{U}

References

Murphy Book Chapter 21.5

References