

# SPECTRAL CLUSTERING

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Let  $x_1, \dots, x_n$  be objects we wish to cluster.

Spectral clustering refers to the following:

①

- form a \_\_\_\_\_
- construct an  $n \times n$  matrix, called the \_\_\_\_\_, from this graph.
- infer the clusters from the \_\_\_\_\_ or \_\_\_\_\_ decomposition of this matrix.

The ingredients that determine a sim. graph are

- 
- 

Given these two ingredients, the weighted adjacency matrix is

$$w_{ij} = \begin{cases} s_{ij} & \text{if } x_i, x_j \text{ adjacent} \\ 0 & \text{otherwise} \end{cases}$$

Examples

▷  $\epsilon$ -neighborhood graph

- $x_i, x_j$  adjacent  $\iff \|x_i - x_j\| \leq \epsilon$
- $s_{ij} = 1$

$\implies$  locality captured entirely by graph

▷ complete (fully connected) graph

- all  $x_i, x_j$  adjacent

- $s_{ij} = \exp\left\{-\frac{\|x_i - x_j\|^2}{2\sigma^2}\right\}, \sigma^2 > 0$

⇒ locality captured entirely by similarities

▷ l-nearest neighbor graph

- $x_i, x_j$  adjacent  $\iff x_i$  is an l-nearest neighbor of  $x_j$  or vice-versa.

- $s_{ij} = \exp\left\{-\frac{\|x_i - x_j\|^2}{2\sigma^2}\right\}$

▷ mutual  $l$ -nearest neighbor graph

- $x_i, x_j$  adjacent  $\iff x_i$  is an  $l$ -nearest neighbor of  $x_j$  and vice versa

Unfortunately, these similarity graphs have free parameters  $(\epsilon, \sigma, l)$  that can be difficult to tune.

The reason for constructing a similarity graph is that it reduces the problem of clustering to \_\_\_\_\_ :

(c)

Find a partition of the graph such that

- edges between clusters have \_\_\_\_\_ weights
- edges within clusters have \_\_\_\_\_ weights

# Graph Laplacians

## Definitions

• The degree of a vertex  $x_i$  is

①  $d_i :=$

• The degree matrix is the diagonal matrix

$$D :=$$

• The unnormalized graph Laplacian is

$$L :=$$

## Note

$L$  is independent of the self-similarity weights  $w_{ii}$ , because

$$L_{ii}$$

## Proposition 1 (Properties of $L$ )

1) For every  $f \in \mathbb{R}^n$

$$f^T L f = \frac{1}{2} \sum_{i,j=1}^n w_{ij} (f_i - f_j)^2$$

2)  $L$  is symmetric and positive semi-definite

3) The smallest eigenvalue of  $L$  is 0

with corresponding eigenvector  $\underline{1} = [1 \ 1 \ \dots \ 1]^T$

4)  $L$  has  $n$  non-negative, real-valued eigenvalues  $0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$

## Proof 1

$$1) \quad f^T L f = f^T D f - f^T W f$$

$$= \sum_{i=1}^n d_i f_i^2 - \sum_{i,j=1}^n w_{ij} f_i f_j$$

$$= \frac{1}{2} \left( \sum_i d_i f_i^2 - 2 \sum_{i,j} w_{ij} f_i f_j + \sum_j d_j f_j^2 \right)$$

$$= \frac{1}{2} \sum_i \sum_j w_{ij} (f_i^2 - 2 f_i f_j + f_j^2)$$

$$= \frac{1}{2} \sum_{i,j} w_{ij} (f_i - f_j)^2$$

2) Symmetry? follows from symmetry of  $W$

PSD follows from 1) because, for any  $f \in \mathbb{R}^n$ ,

$$\textcircled{E} \quad f^T L f =$$
$$\geq$$

3)

$$L \underline{1} = \begin{bmatrix} d_1 & & & & \\ & d_2 & & & \\ & & \ddots & & \\ & & & \ddots & \\ & & & & d_n \end{bmatrix} \begin{bmatrix} 1 \\ 1 \\ \vdots \\ 1 \end{bmatrix} - \begin{bmatrix} w_{11} & w_{12} & \dots & w_{1n} \\ w_{21} & w_{22} & \dots & w_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ w_{n1} & w_{n2} & \dots & w_{nn} \end{bmatrix} \begin{bmatrix} 1 \\ 1 \\ \vdots \\ 1 \end{bmatrix}$$

=

=

4) Follows because  $L$  is

The unnormalized graph Laplacian encodes many properties of the graph. The following is one such property that is relevant for clustering.

Notation | Let  $A \subseteq \{x_1, \dots, x_n\}$  be a cluster.

Define the indicator vector

$$\underline{1}_A = \begin{bmatrix} f_1 \\ f_2 \\ \vdots \\ f_n \end{bmatrix} \in \mathbb{R}^n$$

where  $f_i = 1$  if  $x_i \in A$  and  $f_i = 0$  if  $x_i \notin A$ .

Proposition | Suppose the graph has connected

(F) components  $A_1, \dots, A_k$ . Then the \_\_\_\_\_  
of  $L$  has dimension \_\_\_\_\_ and is  
spanned by



Recall | The nullspace of  $L$  is

$$N(L) := \{f : Lf = 0\}$$

The proposition is claiming

- $\underline{1}_{A_k} \in N(L)$  for each  $k=1, \dots, K$

- If  $f \in N(L)$ , then

$$f = \sum_{k=1}^K \alpha_k \underline{1}_{A_k}$$

Ⓒ

for some  $\alpha_1, \dots, \alpha_k \in \mathbb{R}$ .

Since  $\underline{1}_{A_1}, \dots, \underline{1}_{A_K}$  are linearly independent,

it follows that

$$\dim(N(L)) =$$

## Proof

- $K=1$ : We have already seen  $\underline{1} \in N(L)$ .  
We need to show that if  $Lf = 0$ ,  
then

$$f = \alpha \cdot \underline{1}$$

for some  $\alpha \in \mathbb{R}$ .

If  $Lf = 0$ , then

$$0 = f^T L f = \sum_{i,j=1}^n w_{ij} (f_i - f_j)^2$$

If  $x_i, x_j$  are adjacent, then

$$w_{ij} > 0 \implies f_i = f_j$$

Since  $K=1$ , any two points can be  
connected by a path

$$\implies f_i = \text{constant}$$

- $K > 1$ : Suppose WLOG that the data are ordered according to connected component. Then  $W$  and  $L$  are block diagonal

$$L = \begin{bmatrix} L_1 & & & 0 \\ & L_2 & & \\ & & \ddots & \\ 0 & & & L_k \end{bmatrix}$$

Note each  $L_k$  is a graph Laplacian on  $A_k$ . Then by the previous case we know

- $L \mathbf{1}_{A_k} = 0$  for each  $k$
- If  $Lf = 0$ , then  $f_i$  is constant on each  $A_k$

$$\Rightarrow f = \sum_{k=1}^k \alpha_k \mathbf{1}_{A_k}$$

□

## Spectral Clustering

How can we use this result to devise a clustering algorithm?

### Ideal case

In the ideal case where there are  $K$  connected components and  $K$  is known, we could proceed as follows:

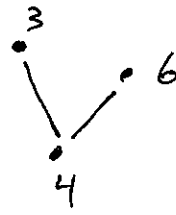
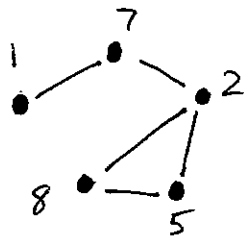
- compute  $L$
- compute a basis  $u_1, \dots, u_k$  for  $N(L)$ .  
 $\hookrightarrow \text{span} \{u_1, \dots, u_k\} = \text{span} \{\mathbf{1}_{A_1}, \dots, \mathbf{1}_{A_k}\}$

- define

$$y_i = (u_{i1}, \dots, u_{ik})$$

- if  $x_i, x_j$  in same components then

Example



$$L = \begin{bmatrix} \times & & & & & & & \\ & \times & & & & & & \\ & & \times & & & & & \\ & & & \times & & & & \\ & & & & \times & & & \\ & & & & & \times & & \\ & & & & & & \times & \\ & & & & & & & \times \end{bmatrix}$$

$$\frac{1}{A_1} = \begin{bmatrix} \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \end{bmatrix}, \quad \frac{1}{A_2} = \begin{bmatrix} \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \end{bmatrix}$$

Use spectral decomposition of  $L$  to find vectors  $u_1, u_2$  that span  $N(L)$ .

Write

$$u_1 = \alpha_1 \frac{1}{A_1} + \beta_1 \frac{1}{A_2}$$
$$u_2 = \alpha_2 \frac{1}{A_1} + \beta_2 \frac{1}{A_2}$$

Set

$$y_i = (u_{1i} \quad u_{2i})$$

Then

$$\begin{bmatrix} y_1 \\ y_2 \\ y_3 \\ y_4 \\ y_5 \\ y_6 \\ y_7 \\ y_8 \end{bmatrix} = \begin{bmatrix} \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \\ \phantom{\times} \end{bmatrix} \Rightarrow \text{determines connected components}$$

In conclusion, we can use the spectral decomposition of the graph Laplacian  $L$  to determine the connected components.

However, there are two problems with this approach

- 1) There are simpler ways of finding the connected components of a graph
- 2) It is very difficult to construct a similarity graph such that its connected components are clusters.

### Nonideal case

More realistically, the components of our similarity graph will not coincide with clusters.

In practice, we should choose the similarity graph such that

number of connected components  $\ll$  desired number of clusters

Then there will be edges connecting points in different clusters, but the weights

⑤ between clusters should be  $\text{---}$   $\text{---}$  the weights within clusters.

Then, if points are indexed according to cluster, we have

$$L = \begin{bmatrix} \boxed{\text{large}} & & \\ & \boxed{\text{large}} & \\ \text{Small} & & \boxed{\text{large}} \\ & \text{Small} & \end{bmatrix} = \begin{bmatrix} \boxed{\text{large}} & & 0 \\ & \boxed{\text{large}} & \\ 0 & & \boxed{\text{large}} \end{bmatrix} + \begin{bmatrix} & & \\ & & \text{small} \\ & & \end{bmatrix}$$



## Perturbation Theory

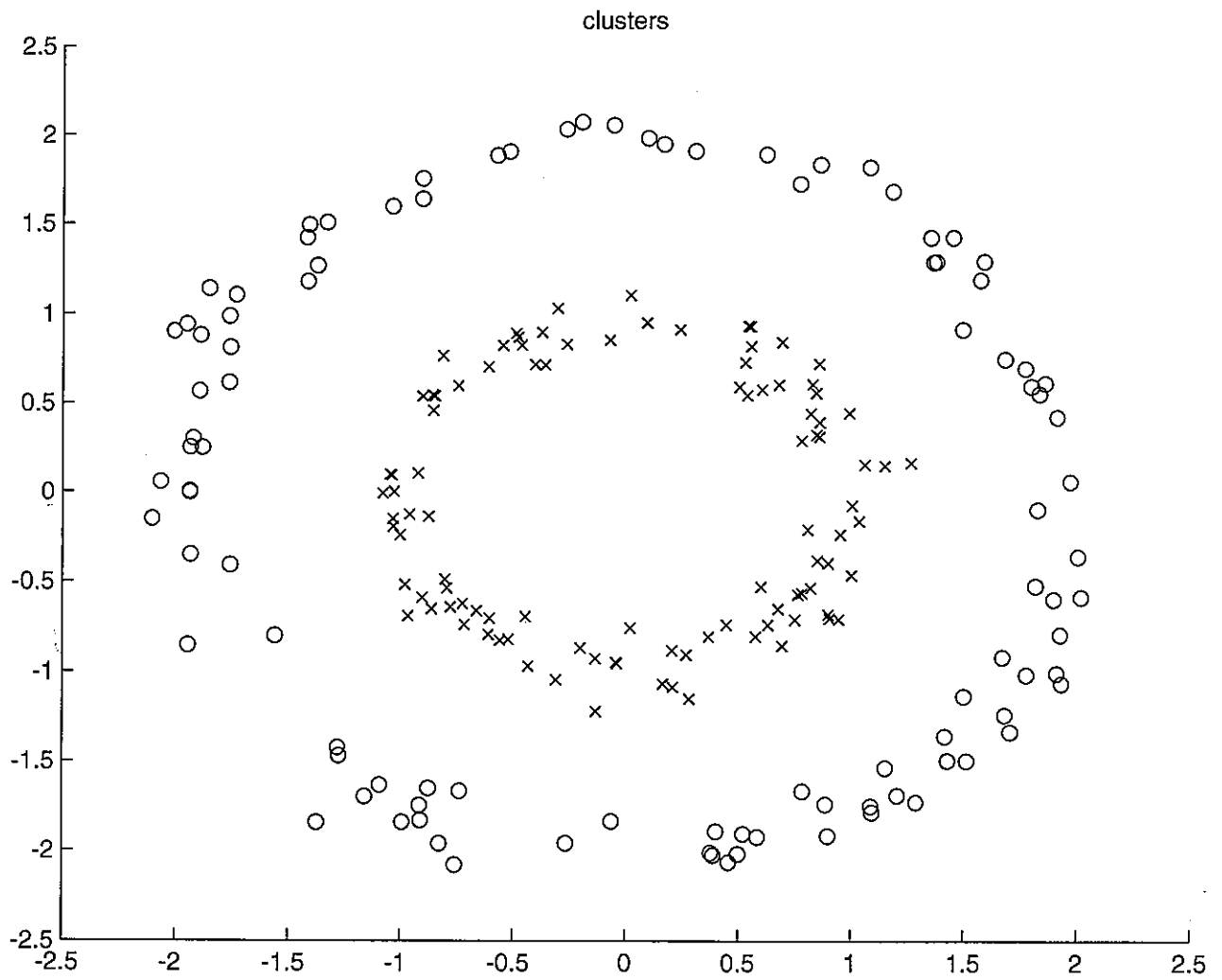
Perturbation theory establishes results that show: if we perturb a matrix by another matrix with small entries, then the eigenvectors and eigenvalues of the matrix are perturbed by a correspondingly small amount.

Recall the nullspace of  $L$  is also the set of eigenvectors with eigenvalue 0. We can use the eigenvectors of  $L$  with the smallest  $K$  eigenvalues as an approximation to the nullspace of an idealized  $L$  based on the true clusters.

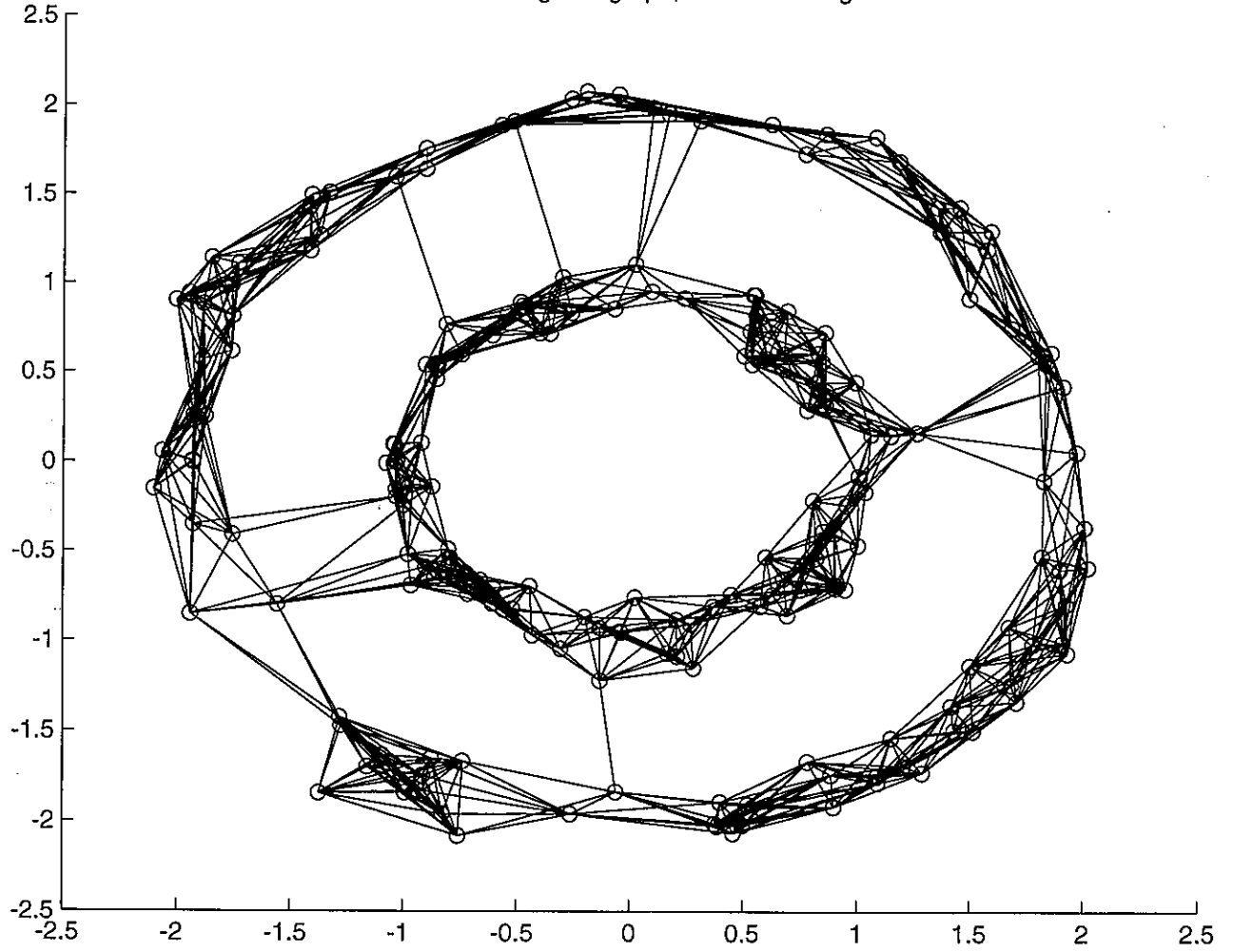


## SPECTRAL CLUSTERING

- Construct similarity graph
- Form graph Laplacian  $L \in \mathbb{R}^{n \times n}$
- Determine the  $K$  smallest eigenvalues of  $L$ ,  
 $0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_K$ , and corresponding  
eigenvectors  $u_1, \dots, u_K \in \mathbb{R}^n$
- Define  $y_i = (u_{1i}, u_{2i}, \dots, u_{Ki})$ ,  $i = 1, \dots, n$
- Cluster  $\{y_i\}_{i=1}^n$  using  $K$ -means clustering  
and assign  $\{x_i\}_{i=1}^n$  to corresponding clusters

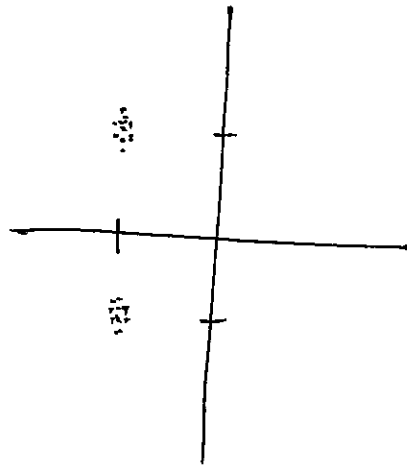


11-nearest neighbor graph, Gaussian weights



-0.070711 -0.07071  
-0.070711 -0.07071  
-0.070711 -0.070709  
-0.070711 -0.07071  
-0.070711 -0.070708  
-0.070711 -0.070708  
-0.070711 -0.070709  
-0.070711 0.071153  
-0.070711 0.071682  
-0.070711 0.070093  
-0.070711 0.071059  
-0.070711 -0.070708  
-0.070711 -0.070709  
-0.070711 0.070098  
-0.070711 -0.07071  
-0.070711 0.071489  
-0.070711 -0.070709  
-0.070711 -0.07071  
-0.070711 0.070098  
-0.070711 -0.070708  
-0.070711 0.070098  
-0.070711 -0.07071  
-0.070711 0.071682  
-0.070711 -0.070709  
-0.070711 -0.07071  
-0.070711 0.070098  
-0.070711 -0.07071  
-0.070711 0.071489  
-0.070711 -0.07071  
-0.070711 -0.070708  
-0.070711 -0.070708  
-0.070711 -0.07071  
-0.070711 -0.070708  
-0.070711 0.071682  
-0.070711 0.070098  
-0.070711 0.070095  
-0.070711 -0.07071  
-0.070711 -0.070708  
-0.070711 -0.07071  
-0.070711 0.070093  
-0.070711 0.071682  
-0.070711 -0.070709  
-0.070711 0.070096  
-0.070711 0.071153  
-0.070711 0.071153  
-0.070711 0.070099  
-0.070711 -0.07071  
-0.070711 0.070093  
-0.070711 0.070093  
-0.070711 0.070098

a few randomly selected  $y_i$ 's



## Normalized Spectral Clustering

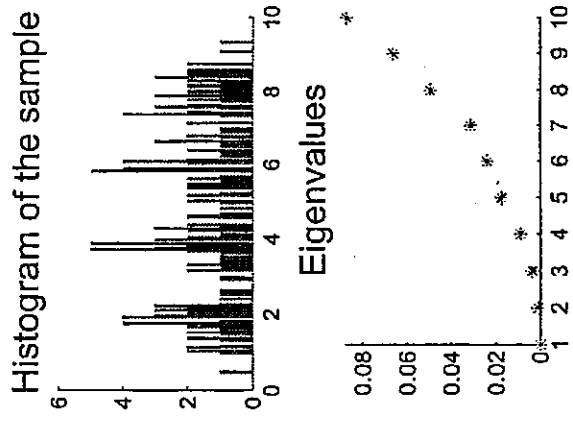
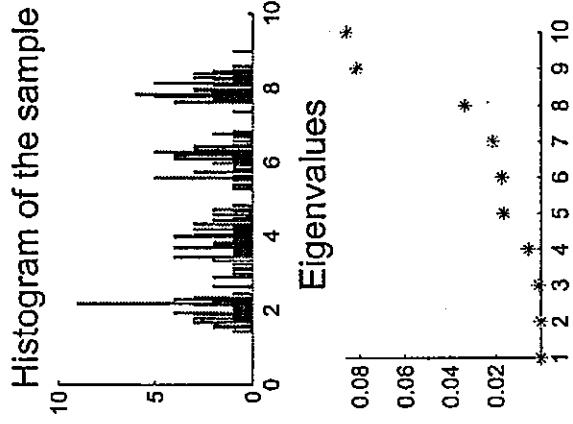
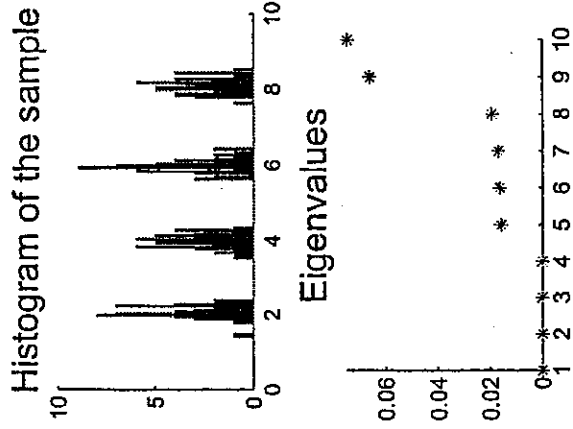
(K) Define the \_\_\_\_\_

$$\tilde{L} :=$$

### Proposition 1

- 1)  $\tilde{L}$  is positive semidefinite with real, nonnegative eigenvalues  $0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$
- 2) 0 is an eigenvalue of  $\tilde{L}$  with corresponding eigenvector  $\underline{1} = [1 \ 1 \ \dots \ 1]^T$ .
- 3) Suppose the graph has connected components  $A_1, \dots, A_K$ . Then the nullspace of  $\tilde{L}$  has dimension K and is spanned by  $\underline{1}_{A_1}, \dots, \underline{1}_{A_K}$

$\implies$  we can substitute  $\tilde{L}$  for  $L$  and get another spectral clustering algorithm.



The first large gap in the spectrum can be used to infer the number of clusters automatically. Results shown are for normalized spectral clustering.

# Graph Cuts

Spectral clustering attempts to find a partition  $A_1, \dots, A_K$  of the similarity graph such that

- (L)
- $w_{ij}$  large if  $x_i, x_j$  in \_\_\_\_\_ cluster
  - $w_{ij}$  small if  $x_i, x_j$  in \_\_\_\_\_ clusters.

using properties of graph Laplacians.

It is also possible to approach this goal directly.

The \_\_\_\_\_ problem is to minimize

$$\text{cut}(A_1, \dots, A_K) := \sum_{i \in A_k, j \in A_l, k \neq l} w_{ij}$$

with respect to  $A_1, \dots, A_K$ , where

$$W(A, B) := \sum_{i \in A, j \in B} w_{ij}$$

The mincut problem is efficiently solvable when  $K=2$ . Unfortunately, it tends to lead to small (and often singleton) clusters.

Therefore researchers have examined modified criteria that favor larger clusters:

- (M)
- Ratio Cut (Hagen and Kahng, 1992)

$$\text{Ratio Cut}(A_1, \dots, A_K) :=$$

where  $|A| :=$

- Normalized Cut (Shi and Malik, 2000)

$$\text{Ncut}(A_1, \dots, A_K) :=$$

=

where  $\text{vol}(A_i) :=$



Unfortunately, minimizing these criteria is NP hard.

(N) Remarkably, however, spectral clustering may be used to solve \_\_\_\_\_ of these problems

In particular,

• unnormalized spectral clustering  $\implies$

• normalized spectral clustering  $\implies$

### Approximating Ratio Cut for $K=2$

We wish to solve

$$\min_A \text{Ratio Cut}(A, \bar{A}) = \min_A \left[ \frac{\text{cut}(A, \bar{A})}{|A|} + \frac{\text{cut}(A, \bar{A})}{|\bar{A}|} \right]$$

Given a subset  $A \subseteq \{1, 2, \dots, n\}$ , define  $f_A := (f_{A1}, \dots, f_{An})^T$  by

$$f_{Ai} := \begin{cases} \sqrt{|\bar{A}|/|A|} & \text{if } i \in A \\ -\sqrt{|A|/|\bar{A}|} & \text{if } i \notin A. \end{cases}$$

Claim  $f_A^T L f_A = n \text{Ratio Cut}(A, \bar{A})$

Proof

$$f_A^T L f_A = \frac{1}{2} \sum_{i,j} w_{ij} (f_{A_i} - f_{A_j})^2$$

$$= \frac{1}{2} \sum_{i \in A, j \in \bar{A}} w_{ij} \left( \sqrt{\frac{|\bar{A}|}{|A|}} + \sqrt{\frac{|A|}{|\bar{A}|}} \right)^2 + \frac{1}{2} \sum_{i \in \bar{A}, j \in A} w_{ij} \left( -\sqrt{\frac{|A|}{|\bar{A}|}} - \sqrt{\frac{|\bar{A}|}{|A|}} \right)^2$$

$$= \frac{1}{2} \text{cut}(A, \bar{A}) \left( \frac{|\bar{A}|}{|A|} + \frac{|A|}{|\bar{A}|} + 2 \right) + \frac{1}{2} \text{cut}(A, \bar{A}) \left( \frac{|A|}{|\bar{A}|} + \frac{|\bar{A}|}{|A|} + 2 \right)$$

$$= \text{cut}(A, \bar{A}) \left( \frac{|\bar{A}|}{|A|} + \frac{|A|}{|\bar{A}|} + 2 \right)$$

$$= \text{cut}(A, \bar{A}) \left( \frac{|A| + |\bar{A}|}{|A|} + \frac{|\bar{A}| + |A|}{|\bar{A}|} \right)$$

$$= \underbrace{(|A| + |\bar{A}|)}_n \cdot \underbrace{\left[ \frac{\text{cut}(A, \bar{A})}{|A|} + \frac{\text{cut}(\bar{A}, A)}{|\bar{A}|} \right]}_{\text{Ratio cut}(A, \bar{A})}$$

Ratio cut(A,  $\bar{A}$ )

Furthermore,  $f_A$  satisfies

$$\textcircled{b} \quad \bullet \quad \underline{1}^T f_A = \sum_{i=1}^n f_{A_i}$$

=

=

$$\bullet \quad \|f_A\|^2 = \sum_{i=1}^n f_{A_i}^2$$

=

=

Therefore, the RatioCut problem can be written as the following optimization problem:

$$\min_A f_A^T L f_A$$

$$\text{s.t.} \quad \underline{1}^T f_A =$$

$$\|f_A\| =$$

If we allow  $f \in \mathbb{R}^n$  we have the following relaxation :

$$\begin{aligned} \min_{f \in \mathbb{R}^n} \quad & f^T L f \\ \text{s.t.} \quad & \mathbf{1}^T f = 0 \\ & \|f\| = \sqrt{n} \end{aligned} \quad \Leftarrow \begin{cases} \text{No longer} \\ \text{require } f = f_A \\ \text{for some } A \end{cases}$$

The solution is

①  $f =$

To recover a solution to the original discrete problem, we can use K-means,  $K=2$ , to cluster the vectors

$$y_i := (\mathbf{1} \quad f_i)$$

Therefore, the approximate solution is given by unnormalized spectral clustering.

A similar analysis applies to  $K > 2$  and to Ncut.

## Final comments

- The  $l$ -nn graph with Gaussian kernel similarity is most common, although the choice of similarity graph is largely an art.
- Another normalized graph Laplacian is

$$\tilde{L} := D^{-\frac{1}{2}} L D^{-\frac{1}{2}}$$

(Ng, Jordan, & Weiss, 2002). It can also be used for spectral clustering but the algorithm requires some modification.

- Which method is preferred?  $\tilde{L}$  is recommended by

U. von Luxburg, "A Tutorial on Spectral Clustering," 2007.

- $b$ -matching is an interesting alternative to nearest neighbor graphs: it ensures that each node has the same number of incident edges (unweighted degree). See Tebara et al, ICML 2009.

• graph Laplacians are also used in  
\_\_\_\_\_ learning.

②

For example, suppose we are in a regression setting with data

$$(x_i, y_i)_{i=1}^m$$

$$(x_i)_{i=m+1}^n$$

To make sure the estimated function doesn't "wiggle" too much, we could minimize

$$\frac{1}{m} \sum_{i=1}^m (y_i - f(x_i))^2 + \frac{\lambda}{2} \underbrace{\sum_{i,j=1}^n w_{ij} (f(x_i) - f(x_j))^2}_{f^T L f}$$

# Key

A. similarity graph, graph Laplacian, eigenvalue/spectral

B. a graph, a similarity matrix  $S = [s_{ij}]_{i,j=1}^n$

C. graph partitioning, low, high

D.  $d_i = \sum_{j=1}^n w_{ij}$ ,  $D = \begin{bmatrix} d_1 & & 0 \\ & \ddots & \\ 0 & & d_n \end{bmatrix}$ ,  $L = D - W$

$$L_{ii} = d_i - w_{ii} = \sum_{j \neq i} w_{ij}$$

$$E. f^T L f = \sum w_{ij} (f_i - f_j)^2 \geq 0$$

$$L \underline{1} = \begin{bmatrix} d_1 \\ \vdots \\ d_n \end{bmatrix} - \begin{bmatrix} \sum w_{ij} \\ \vdots \\ \sum w_{nj} \end{bmatrix} = \begin{bmatrix} 0 \\ \vdots \\ 0 \end{bmatrix} = 0 \cdot \underline{1}$$

L is PSD

F. nullspace (0-eigenspace);  $K$ ;  $\underline{1}_{A_1}, \dots, \underline{1}_{A_K}$

G.  $f = \sum_{k=1}^K \alpha_k \underline{1}_{A_k}$ ,  $\dim(N(L)) = K =$  multiplicity of 0 as an eigenvalue

H.  $y_i = y_j$

I.  $\underline{1}_{A_1} = \begin{bmatrix} 1 \\ 1 \\ 0 \\ 0 \\ 1 \\ 0 \\ 1 \\ 1 \end{bmatrix}$ ,  $\underline{1}_{A_2} = \begin{bmatrix} 0 \\ 0 \\ 1 \\ 1 \\ 0 \\ 1 \\ 0 \\ 0 \end{bmatrix}$ ,  $\begin{bmatrix} y_1 \\ y_2 \\ y_3 \\ y_4 \\ y_5 \\ y_6 \\ y_7 \\ y_8 \end{bmatrix} = \begin{bmatrix} \alpha_1 & \alpha_2 \\ \alpha_1 & \alpha_2 \\ \beta_1 & \beta_2 \\ \beta_1 & \beta_2 \\ \alpha_1 & \alpha_2 \\ \beta_1 & \beta_2 \\ \alpha_1 & \alpha_2 \\ \alpha_1 & \alpha_2 \end{bmatrix}$

J. less than, ideal case, noise/perturbation

K. normalized graph Laplacian  $\tilde{L} = D^{-1}L = I - D^{-1}W$

L. same, different, mincut,  $\text{cut}(A_1, \dots, A_k) = \frac{1}{2} \sum_{k=1}^K W(A_k, \bar{A}_k)$ ,  $W(A, B) = \sum_{i \in A} \sum_{j \in B} w_{ij}$

M. Ratio Cut  $(A_1, \dots, A_k) = \frac{1}{2} \sum_{k=1}^K \frac{W(A_k, \bar{A}_k)}{|A_k|} = \sum_{k=1}^K \frac{\text{cut}(A_k, \bar{A}_k)}{|A_k|}$

$|A_k| = \#$  of nodes in  $A$

$N\text{cut}(A_1, \dots, A_k) = \frac{1}{2} \sum_{k=1}^K \frac{W(A_k, \bar{A}_k)}{\text{vol}(A_k)} = \sum_{k=1}^K \frac{\text{cut}(A_k, \bar{A}_k)}{\text{vol}(A_k)}$

$\text{vol}(A_k) = \sum_{i \in A} \sum_{j \in A} w_{ij}$

N. relaxations, Ratio Cut, Ncut

O.  $\mathbb{1}^T f_A = \sum_{i \in A} \sqrt{\frac{|\bar{A}|}{|A|}} - \sum_{i \in \bar{A}} \sqrt{\frac{|A|}{|\bar{A}|}} = \sqrt{|A| \cdot |\bar{A}|} - \sqrt{|A| \cdot |\bar{A}|} = 0$

$\|f_A\|^2 = \sum_{i \in A} \frac{|\bar{A}|}{|A|} + \sum_{i \in \bar{A}} \frac{|A|}{|\bar{A}|} = |\bar{A}| + |A| = n$

discrete, 0,  $\sqrt{n}$

P. eigenvector of  $L$  corresponding to second smallest eigenvalue

Q. semi-supervised