# Spectral Clustering

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### Quadratic Form

- Let **A** be some  $n \times n$  matrix.
- What is Ax? What's the type of the output? What may x represent?
  - Some numeric assignment on all nodes in G.
  - E.g., what if  $x_i \in \{0,1\}$ ?  $x_i \in [0,1]$ ?  $x_i \in \Re$ ?
- What is x<sup>T</sup>Ax? What's the type of the output? Why it is called a qudratic form?

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- What is x<sup>T</sup>Ax? What's the type of the output? Why it is called a qudratic form?
  - $\mathbf{x}^{\top} \mathbf{A} \mathbf{x} = \sum_{i,j} A_{ij} \cdot (x_i x_j)$

## Unnormalized Graph Laplacian

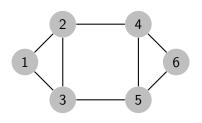
- Let **A** is the adjacency matrix of a "normal" (unweighted) undirected graph G.  $\mathbb V$  are the vertices of G and  $\mathbb E$  are the edges of G
  - An edge between  $v_i$  and  $v_j$  is modelled as (i,j) and (j,i), i.e.,  $A_{ij} = A_{ji} = 1$ .
  - $A_{ii} = \underline{\hspace{1cm}}?$
- What about  $\mathbf{x}^{\top}\mathbf{I}\mathbf{x}$ ?
- What about  $\mathbf{x}^{\top}\mathbf{D}\mathbf{x}$ , where  $\mathbf{D} = \mathrm{Diag}(d_1, d_2, \dots, d_n)$  and  $d_i = deg(v_i)$ ?
- What about  $\mathbf{x}^{\top} \mathbf{A} \mathbf{x}$ ?
- Now what about  $2(\mathbf{x}^{\top}\mathbf{D}\mathbf{x} \mathbf{x}^{\top}\mathbf{A}\mathbf{x})$ ?

### Example

$$\mathbf{x}^{ op} \mathbf{L} \mathbf{x} = rac{1}{2} \cdot \sum_{\mathbf{e}_{ij} \in \mathbb{E}} (x_i - x_j)^2$$
 , where  $\mathbf{L} = \mathbf{D} - \mathbf{A}$ .

ullet  $\ell_2$  differences between assignments on the two ends of an edge, summed over all edges.

## Example



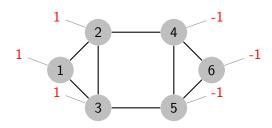
	$n_1$	$n_2$	n <sub>3</sub>	n <sub>4</sub>	<i>n</i> <sub>5</sub>	n <sub>6</sub>
$n_1$						
$n_2$						
<i>n</i> <sub>3</sub>						
$n_4$						
n <sub>5</sub>						
<i>n</i> <sub>6</sub>						

- 1 is the one vector.
- L1 = (NB:  $L^{\top} = L$ )  $\Longrightarrow$   $\lambda_1 = 0, \nu_1 = 1$

$$\implies \lambda_1 = 0, v_1 = 1$$

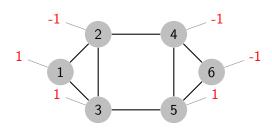
 $\mathbf{o} \mathbf{x}^{\mathsf{T}} \mathbf{L} \mathbf{x} =$ 

# Binary x induces a Clustering /1



- x =
- $\mathbf{o} \mathbf{x}^{\mathsf{T}} \mathbf{L} \mathbf{x} =$

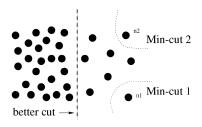
# Binary x induces a Clustering /2



- x =
- $\bullet$   $\mathbf{x}^{\top}\mathbf{L}\mathbf{x} =$
- $\bullet \ \mathbf{x}^{\top}\mathbf{x} =$

### Min Cut vs. Normalized Cut

- Min cuts are not always desirable.
  - Biased towards cutting small sets of isolated nodes.



- Cut:  $cut(A, B) = \sum_{v_i \in A, v_i \in B} w_{ij}$ .
- Normalized cut:

$$ncut(A, B) = \frac{cut(A, B)}{vol(A)} + \frac{cut(A, B)}{vol(B)},$$

where 
$$vol(A) = \sum_{v_i \in A} d_i) = \sum_{v_i \in A, v_j \in \mathbb{V}} w_{i,j}$$
.

### Connection to L

$$ncut(A, B) = cut(A, B) \left(\frac{1}{vol(A)} + \frac{1}{vol(B)}\right)$$

- Let  $x_i = \frac{1}{vol(A)}$  if  $v_i \in A$ , and  $= \frac{-1}{vol(B)}$  otherwise.
- $\mathbf{x}^{\top} \mathbf{L} \mathbf{x} = \sum_{e} w_{ij} (x_i x_j)^2 = 0 + \sum_{v_i \in A, v_j \in B} \left( \frac{1}{vol(A)} + \frac{1}{vol(B)} \right)^2$
- $\mathbf{x}^{\top} \mathbf{D} \mathbf{x} = (\mathbf{x}^{\top} D) \mathbf{x} = \sum_{E} d_{i} x_{i}^{2} = \sum_{v_{i} \in A} \frac{d_{i}}{vol(A)^{2}} + \sum_{v_{j} \in B} \frac{d_{j}}{vol(B)^{2}} = \frac{1}{vol(A)} + \frac{1}{vol(B)}$

$$ncut(A, B) = \frac{\mathbf{x}^{\top} \mathbf{L} \mathbf{x}}{\mathbf{x}^{\top} \mathbf{D} \mathbf{x}}$$

## Relaxation and Optimization

Minimize 
$$ncut(A, B) = \frac{\mathbf{x}^{\top} \mathbf{L} \mathbf{x}}{\mathbf{x}^{\top} \mathbf{D} \mathbf{x}}$$
 Subject to  $x_i \in \left\{ \frac{1}{vol(A)}, \frac{-1}{vol(B)} \right\}$ 

- NP-hard to optimize under the discrete constraint.
- Relaxation: grow the feasible region of x and find the minimum value within the enlarged region.
  - allow x to be a real vector?
  - Yes, but too large.
  - This gives the constraint:  $\mathbf{x}^{\mathsf{T}}\mathbf{D}\mathbf{1} = 0$  or equivalently  $\mathbf{x}^{\mathsf{T}}\mathbf{D} \perp \mathbf{1}$  (You can verify this by plugging in any discrete vectors)
- Solution: the second smallest eigenvector of the generalized eigen value problem  $\mathbf{L}\mathbf{x} = \lambda \mathbf{D}\mathbf{x}$ .
- Normalized Laplacian:

$$\mathbf{L}' = \mathbf{D}^{-\frac{1}{2}}(\mathbf{D} - \mathbf{W})\mathbf{D}^{-\frac{1}{2}} = \mathbf{I} - \mathbf{D}^{-\frac{1}{2}}\mathbf{W}\mathbf{D}^{-\frac{1}{2}}$$

## Spectral Clustering Algorithm Framework

- Algorithm SC\_recursive\_bin\_cut(data, k)
  - Construct the weighted graph G
  - Construct the special graph laplacian *L* for *G*.
  - Compute the smallest non-zero eigenvector for L. This is the new representation of vertices in a new 1-dimensional space (i.e., embedding).
  - Cluster the vertices in the embedding space according to the objective function.
  - For each cluster, recursively call the algorithm if more clusters are needed

# Spectral Clustering Algorithm Framework

- Algorithm SC\_k\_way\_cut(data, k)
  - ullet Construct the weighted graph G
  - Construct the special graph laplacian *L* for *G*.
  - Compute the smallest t non-zero eigenvector for L. This is the new representation of vertices in a new t-dimensional space (i.e., embedding).
  - Cluster the vertices in the embedding space using another clustering algorithm (e.g., k-means)

### Notes on the Algorithms

- How to construct the weighted graph if only n objects are given?
  - Be based on the similarity or distance among objects.
  - E.g.,  $w_{ij} = \exp(\frac{\|f(o_i) f(o_j)\|}{2\sigma^2})$  where f(o) is the feature vector of object o. One can also induce a sparse graph if one caps the raw weights by a threshold.
- Which Laplacian to use?
  - Unnormalized graph laplacian  $\mathbf{L} = \mathbf{D} \mathbf{W}$ .
  - Normalized graph laplacian  $\mathbf{L} = \mathbf{D}^{-\frac{1}{2}}(\mathbf{D} \mathbf{W})\mathbf{D}^{-\frac{1}{2}}$ .

## Comments on Spectral Clustering

#### Pros:

- Usually better quality than other methods.
- Can be thought of (non-linear) dimensionality reduction or embedding.
- Freedom to construct a (sparse) G to preserve local similarity/connectivity.
- Only requires some similarity measure.
- Could be more efficient than k-means for high-dimensional sparse vectors (esp. if k-means is not fully optimized for such case).

### Cons:

- Still need to determine k
- Assumes clusters are of similar sizes.
- Does not scale well with large datasets; but more scalable variants exist.
- One of the relaxation of the original NP-hard problem may not be the tightest relaxation.