## Community Structure in Networks

Thanks to Jure Leskovec, Stanford and Panayiotis Tsaparas, Univ. of loannina for slides

## Agenda

- Network Communities
- Community Detection
- Method 1: Girvan-Newman
- Method 2: Modularity Optimization
- Community Detection
- Graph Cuts
- Spectral Clustering
- Network Profiling
- Communities: Issues and Questions

Network Communities

## Networks \& Communities

- We often think of networks "looking" like this:

- What lead to such a conceptual picture?


## Networks: Flow of Information

- How information flows through the network?
- What structurally distinct roles do nodes play?
- What roles do different links (short vs. long) play?
- How people find out about new jobs?
- Mark Granovetter, part of his PhD in 1960s
- People find the information through personal contacts
- But: Contacts were often acquaintances rather than close friends
- This is surprising: One would expect your friends to help you out more than casual acquaintances
- Why is it that acquaintances are most helpful?


## Granovetter's Explanation

- Granovetter makes a connection between social and structural role of an edge
- First point: Structure
- Structurally embedded edges are socially strong
- Long-range edges spanning different parts of the network are socially weak
- Second point: Information
- Long-range edges allow you to gather information from different parts of the network and get a job
- Structurally embedded edges are heavily redundant in terms of information access



## Conceptual Picture of Networks

- Granovetter's theory leads to the following conceptual picture of networks



## Network Communities

- Granovetter's theory
suggest that networks are composed of tightly connected sets of nodes


Communities, clusters, groups, modules

- Sets of nodes with lots of connections inside and few to outside (the rest of the network)


## Finding Network Communities

- How to automatically find such densely connected groups of nodes?
- Ideally such automatically detected clusters would then correspond to real groups
- For example:


Communities, clusters, groups, modules

## Social Network Data



- Zachary's Karate club network:
- Observe social ties and rivalries in a university karate club
- During his observation, conflicts led the group to split
- Split could be explained by a minimum cut in the network


## NCAA Football Network



Nodes: Teams
Edges: Games played

## NCAA Football Network



## NCAA conferences

O Mid American
O Big East

- Atlantic Coast

O SEC
O Conference USA

- Big 12

O Western Athletic
$\bigcirc$ Pacific 10
O Mountain West

- Big 10

O Sun Belt
$\bigcirc$ Independents

Nodes: Teams
Edges: Games played

## Facebook Ego-network



## Facebook Ego-network



## Protein-Protein Interactions



Nodes: Proteins
Edges: Interactions

## Protein-Protein Interactions



Nodes: Proteins
Edges: Interactions

## Community Detection

## How to find communities?



We will work with undirected (unweighted) networks

## Method 1: Strength of Weak Ties

- Edge betweenness: Number of shortest paths passing over the edge
- Intuition:


Edge strengths (call volume) in a real network


Edge betweenness in a real network

## Method 1: Girvan-Newman

- Divisive hierarchical clustering based on the notion of edge betweenness:
Number of shortest paths passing through the edge
- Girvan-Newman Algorithm:
- Undirected unweighted networks
- Repeat until no edges are left:
- Calculate betweenness of edges
- Remove edges with highest betweenness
- Connected components are communities
- Gives a hierarchical decomposition of the network


## Girvan-Newman: Example



Need to re-compute betweenness at every step

## Girvan-Newman: Example

Step 1:



Step 3:

(7)

(5)

Step 2:

-

Hierarchical network decomposition:


## Girvan-Newman: Results



Communities in physics collaborations

## Girvan-Newman: Results

- Zachary’s Karate club: Hierarchical decomposition



## We need to resolve 2 questions

1. How to compute betweenness?
2. How to select the number of clusters?


## How to Compute Betweenness?

- Want to compute betweenness of paths starting at node $\boldsymbol{A}$

- Breadth first search starting from $A$ :



## How to Compute Betweenness?

- Count the number of shortest paths from $A$ to all other nodes of the network:



## How to Compute Betweenness?

- Compute betweenness by working up the tree: If there are multiple paths count them fractionally

The algorithm:
-Add edge flows:
-- node flow =
$1+\sum$ child edges
-- split the flow up
based on the parent value

- Repeat the BFS procedure for each starting node $U$



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## We need to resolve 2 questions

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2. How to select the number of clusters?


## Network Communities

- Communities: sets of tightly connected nodes
- Define: Modularity Q
- A measure of how well a network is partitioned into communities

- Given a partitioning of the network into groups $\boldsymbol{S} \in \boldsymbol{S}$ :
$Q \propto \sum_{s \in S}[(\#$ edges within group $s)-$ $\underbrace{\text { (expected \# edges within group } s \text { )] }}$

Need a null model!

## Null Model: Configuration Model

- Given real $\boldsymbol{G}$ on $\boldsymbol{n}$ nodes and $\boldsymbol{m}$ edges, construct rewired network $\boldsymbol{G}^{\prime}$
- Same degree distribution but random connections
- Consider $\boldsymbol{G}^{\prime}$ as a multigraph

- The expected number of edges between nodes $\boldsymbol{i}$ and $\boldsymbol{j}$ of degrees $\boldsymbol{k}_{\boldsymbol{i}}$ and $\boldsymbol{k}_{\boldsymbol{j}}$ equals to: $\boldsymbol{k}_{\boldsymbol{i}} \cdot \frac{\boldsymbol{k}_{\boldsymbol{j}}}{2 \boldsymbol{m}}=\frac{\boldsymbol{k}_{\boldsymbol{i}} \boldsymbol{k}_{\boldsymbol{j}}}{2 \boldsymbol{m}}$
- The expected number of edges in (multigraph) $\mathbf{G}^{\prime}$ :

$$
\begin{aligned}
& =\frac{1}{2} \sum_{i \in N} \sum_{j \in N} \frac{k_{i} k_{j}}{2 m}=\frac{1}{2} \cdot \frac{1}{2 m} \sum_{i \in N} k_{i}\left(\sum_{j \in N} k_{j}\right)= \\
& =\frac{1}{4 m} 2 m \cdot 2 m=m
\end{aligned}
$$

$$
\begin{aligned}
& \text { Note: } \\
& \sum_{u \in N} k_{u}=2 m
\end{aligned}
$$

## Modularity

- Modularity of partitioning S of graph G:
" $\mathbf{Q} \propto \sum_{s \in S}$ [ (\# edges within group $s$ ) (expected \# edges within group $s$ )]
- $\boldsymbol{Q}(\boldsymbol{G}, S)=\underbrace{\frac{1}{2 m}} \sum_{s \in S} \sum_{i \in s} \sum_{j \in s}\left(A_{i j}-\frac{k_{i} k_{j}}{2 m}\right)$ Normalizing cost.: $-1<\mathrm{Q}<1$
- Modularity values take range [-1,1]
- It is positive if the number of edges within groups exceeds the expected number
- 0.3-0.7<Q means significant community structure


## Modularity: Number of clusters

- Modularity is useful for selecting the number of clusters:


Why not optimize Modularity directly?

Modularity Optimization

## Method 2: Modularity Optimization

- Let's split the graph into 2 communities!
- Want to directly optimize modularity!
$-\max _{S} Q(G, S)=\frac{1}{2 m} \sum_{s \in S} \sum_{i \in S} \sum_{j \in s}\left(A_{i j}-\frac{k_{i} k_{j}}{2 m}\right)$
- Community membership vector s:
- $s_{i}=1$ if node $i$ is in community 1
-1 if node $\boldsymbol{i}$ is in community $\mathbf{- 1}$

$$
\frac{s_{i} s_{j}+1}{2}=\begin{aligned}
& 1 . . \text { if } \mathrm{s}_{\mathrm{i}}=\mathrm{s}_{\mathrm{j}} \\
& 0 . . \text { else }
\end{aligned}
$$

- $Q(G, s)=\frac{1}{2 m} \sum_{i \in N} \sum_{j \in N}\left(A_{i j}-\frac{k_{i} k_{j}}{2 m}\right) \frac{\left(s_{i} s_{j}+1\right)}{2}$

$$
=\frac{1}{4 m} \sum_{i, j \in N}\left(A_{i j}-\frac{k_{i} k_{j}}{2 m}\right) s_{i} s_{j}
$$

## Modularity Matrix

Define:

- Modularity matrix: $B_{i j}=A_{i j}-\frac{k_{i} k_{j}}{2 m}$

$$
\begin{aligned}
& \text { Note: each row/col of B } \\
& \text { sums to } 0: \sum_{j} A_{i j}=\boldsymbol{k}_{i}, \\
& \sum_{j} \frac{k_{i} \boldsymbol{k}_{j}}{2 m}=\boldsymbol{k}_{i} \sum_{j} \frac{k_{j}}{2 m}=\boldsymbol{k}_{i}
\end{aligned}
$$

- Membership: $s=\{-1,+1\}$
- Then: $Q(G, s)=\frac{1}{4 m} \sum_{i \in N} \sum_{j \in N}\left(A_{i j}-\frac{k_{i} k_{j}}{2 m}\right) s_{i} s_{j}$

$$
\begin{aligned}
& =\frac{1}{4 m} \sum_{i, j \in N} B_{i j} s_{i} s_{j} \\
& =\frac{1}{4 m} \sum_{i} s_{i} \underbrace{\sum_{j} B_{i j} s_{j}}_{=B_{i} \cdot s}=\frac{1}{4 m} s^{T} B s
\end{aligned}
$$

- Task: Find $\mathbf{s} \in\{-\mathbf{1}, \mathbf{+ 1}\}^{n}$ that maximizes $\mathbf{Q}(\mathbf{G}, \mathbf{s})$


## Quick Review of Linear Algebra

- Symmetric matrix A
- That is positive semi-definite:

$$
\left[\begin{array}{ccc}
a_{11} & \ldots & a_{1 n} \\
\vdots & & \vdots \\
a_{n 1} & \ldots & a_{n n}
\end{array}\right]\left[\begin{array}{c}
x_{1} \\
\vdots \\
x_{n}
\end{array}\right]=\lambda\left[\begin{array}{c}
x_{1} \\
\vdots \\
x_{n}
\end{array}\right]
$$

$\boldsymbol{A}=\boldsymbol{U} \cdot \boldsymbol{U}^{\boldsymbol{T}}$

- Then solutions $\lambda, \boldsymbol{x}$ to equation $\boldsymbol{A} \cdot \boldsymbol{x}=\lambda \cdot \boldsymbol{x}$ :
- Eigenvectors $\boldsymbol{x}_{\boldsymbol{i}}$ ordered by the magnitude of their corresponding eigenvalues $\lambda_{i}\left(\lambda_{1} \leq \lambda_{2} \ldots \leq \lambda_{n}\right)$
- $\boldsymbol{x}_{\boldsymbol{i}}$ are orthonormal (orthogonal and unit length)
- $\boldsymbol{x}_{\boldsymbol{i}}$ form a coordinate system (basis)
- If $\boldsymbol{A}$ is positive-semidefinite: $\lambda_{i} \geq 0$ (and they always exist)
- Eigen Decomposition theorem: Can rewrite matrix $\boldsymbol{A}$ in terms of its eigenvectors and eigenvalues: $\boldsymbol{A}=$ $\sum_{i} x_{i} \cdot \lambda_{i} \cdot x_{i}^{T}$


## Modularity Optimization

- Rewrite: $Q(G, s)=\frac{1}{4 m} s^{\mathrm{T}} B s$ in terms of its eigenvectors and eigenvalues:
$=\mathrm{s}^{\mathrm{T}}\left[\sum_{i=1}^{n} x_{i} \lambda_{i} x_{i}^{T}\right] s=\sum_{i=1}^{n} s^{T} x_{i} \lambda_{i} x_{i}^{T} s=\sum_{i=1}^{n}\left(s^{T} \mathrm{X}_{i}\right)^{2} \lambda_{i}$
- So, if there would be no other constraints on $s$ then to maximize $Q$, we make $s=x_{n}$
- Why? Because $\lambda_{n} \geq \lambda_{n-1} \geq \cdots$
- Remember $\boldsymbol{s}$ has fixed length!
- Assigns all weight in the sum to $\lambda_{\boldsymbol{n}}$ (largest eigenvalue) - All other $\boldsymbol{s}^{\boldsymbol{T}} \boldsymbol{x}_{\boldsymbol{i}}$ terms are zero because of orthonormality


## Finding the vector $s$

- Let's consider only the first term in the summation (because $\lambda_{\boldsymbol{n}}$ is the largest): $\max _{s} Q(G, s)=\sum_{i=1}^{n}\left(s^{T} x_{i}\right)^{2} \lambda_{i} \approx\left(s^{T} x_{n}\right)^{2} \lambda_{n}$
- Let's maximize: $\sum_{j=1}^{n} s_{j} \cdot x_{n, j}$ where $s_{j} \in\{-1,+1\}$
- To do this, we set:
$s_{j}= \begin{cases}+1 & \left.\text { if } x_{n, j} \geq 0 \text { ( } j-\text { th coordinate of } x_{n} \geq 0\right) \\ -1 & \left.\text { if } x_{n, j}<0 \text { ( } j-\text { th coordinate of } x_{n}<0\right)\end{cases}$
- Continue the bisection hierarchically


## Summary: Modularity Optimization

- Fast Modularity Optimization Algorithm:
- Find leading eigenvector $\boldsymbol{x}_{\boldsymbol{n}}$ of modularity matrix B
- Divide the nodes by the signs of the elements of $\boldsymbol{x}_{\boldsymbol{n}}$
- Repeat hierarchically until:
- If a proposed split does not cause modularity to increase, declare community indivisible and do not split it
- If all communities are indivisible, stop
- How to find $\boldsymbol{x}_{\boldsymbol{n}}$ ? Power method!
- Start with random $v^{(0)}$, repeat :
- When converged $\left(v^{(t)} \approx v^{(t+1)}\right)$, set $\boldsymbol{x}_{n}=\boldsymbol{v}^{(t)}$

$$
v^{(t+1)}=\frac{B v^{(t)}}{\left\|B v^{(t)}\right\|}
$$

## Summary: Modularity

- Girvan-Newman:
" Based on the "strength of weak ties"
- Remove edge of highest betweenness
- Modularity:
- Overall quality of the partitioning of a graph
- Use to determine the number of communities
- Fast modularity optimization:
- Transform the modularity optimization to a eigenvalue problem

Community Detection: Graph Cuts \& Spectral Clustering

## Agenda

## Graph Partitioning

- Graph Cuts
- Spectral Clustering


## Graph Partitioning

- Undirected graph $\boldsymbol{G}(\boldsymbol{V}, \boldsymbol{E})$ :
- Bi-partitioning task:

- Divide vertices into two disjoint groups $\boldsymbol{A}, \boldsymbol{B}$

- Questions:
" How can we define a "good" partition of $\boldsymbol{G}$ ?
- How can we efficiently identify such a partition?


## Graph Partitioning

- What makes a good partition?
- Maximize the number of within-group connections
- Minimize the number of between-group connections



## Graph Cuts

- Express partitioning objectives as a function of the "edge cut" of the partition
- Cut: Set of edges with only one vertex in a group:

$$
\operatorname{cut}(A, B)=\sum_{i \in A, j \in B} w_{i j}
$$



## Graph Cut Criterion

- Criterion: Minimum-cut
- Minimize weight of connections between groups $\arg \min _{\mathrm{A}, \mathrm{B}} \operatorname{cut}(A, B)$
- Degenerate case:

- Problem:
- Only considers external cluster connections
- Does not consider internal cluster connectivity


## Graph Bisection

- Since the minimum cut does not always yield good results we need extra constraints to make the problem meaningful
- Graph Bisection
- Partition the graph into two equal sets of nodes
- Kernighan-Lin algorithm
- Start with random equal partitions
- Swap nodes to improve some quality metric (e.g., cut, modularity, etc)


## Ratio Cut

## Criterion: Ratio-cut

Normalize cut by the size of the groups

$$
\text { Ratio-cut }=\frac{\operatorname{Cut}(\mathrm{U}, \mathrm{~V}-\mathrm{U})}{|U|}+\frac{\operatorname{Cut}(\mathrm{U}, \mathrm{~V}-\mathrm{U})}{|V-U|}
$$

## Normalized Cut

## Criterion: Normalized-cut

Connectivity between groups relative to the density of each group

Normalized-cut $=\frac{\operatorname{Cut}(\mathrm{U}, \mathrm{V}-\mathrm{U})}{\operatorname{Vol}(\mathrm{U})}+\frac{\operatorname{Cut}(\mathrm{U}, \mathrm{V}-\mathrm{U})}{\operatorname{Vol}(V-U)}$
$\operatorname{vol}(U)$ : total weight of the edges with at least one endpoint in $U: \operatorname{vol}(U)=\sum_{i \in U} d_{i}$

Why use these criteria?

- Produce more balanced partitions


## An Example



Red is Min-Cut
Ratio-Cut $($ Red $)=\frac{1}{1}+\frac{1}{8}=\frac{9}{8}$
Ratio-Cut(Green) $=\frac{2}{5}+\frac{2}{4}=\frac{18}{20}$
Normalized-Cut(Red) $=\frac{1}{1}+\frac{1}{27}=\frac{28}{27}$
Normalized-Cut(Green) $=\frac{2}{12}+\frac{2}{16}=\frac{14}{48}$

Minimizing Normalizedcut is even better for Green due to density

## Another Example



Which of the three cuts has the best (min, normalized, ratio) cut?

## Graph Cut Criteria

- Criterion: Conductance [Shi-Malik, '97]
- Connectivity between groups relative to the density of each group

$$
\phi(A, B)=\frac{\operatorname{cut}(A, B)}{\min (\operatorname{vol}(A), \operatorname{vol}(B))}
$$

$\operatorname{vol}(\boldsymbol{A})$ : total weight of the edges with at least one endpoint in $A: \operatorname{vol}(A)=\sum_{i \in A} \boldsymbol{k}_{\boldsymbol{i}}$
$■$ Why use this criterion?

- Produces more balanced partitions
- How do we efficiently find a good partition?
- Problem: Computing optimal cut is NP-hard


## Graph Cuts

Ratio-cut and normalized-cut can be reformulated in matrix format and solved using spectral clustering

# Spectral Clustering for Graph Partitioning 

## Spectral Clustering Algorithms

- Three basic stages:
- 1) Pre-processing
- Construct a matrix representation of the graph
- 2) Decomposition
- Compute eigenvalues and eigenvectors of the matrix
- Map each point to a lower-dimensional representation based on one or more eigenvectors
- 3) Grouping
- Assign points to two or more clusters, based on the new representation
- But first, let's define the problem


## Spectral Graph Partitioning

- $\boldsymbol{A}$ : adjacency matrix of undirected $\mathbf{G}$
- $\boldsymbol{A}_{i j}=\mathbf{1}$ if $(\boldsymbol{i}, \boldsymbol{j})$ is an edge, else $\mathbf{0}$
$-\boldsymbol{x}$ is a vector in $\mathfrak{R}^{n}$ with components ( $\boldsymbol{x}_{\boldsymbol{1}}, \ldots, \boldsymbol{x}_{\boldsymbol{n}}$ )
- Think of it as a label/value of each node of $\boldsymbol{G}$
- What is the meaning of $A \cdot x$ ?

$$
\left[\begin{array}{ccc}
a_{11} & \ldots & a_{1 n} \\
\vdots & & \vdots \\
a_{n 1} & \ldots & a_{n n}
\end{array}\right]\left[\begin{array}{c}
x_{1} \\
\vdots \\
x_{n}
\end{array}\right]=\left[\begin{array}{c}
y_{1} \\
\vdots \\
y_{n}
\end{array}\right] \quad y_{i}=\sum_{j=1}^{n} A_{i j} x_{j}=\sum_{(i, j) \in E} x_{j}
$$

- Entry $y_{i}$ is a sum of labels $x_{j}$ of neighbors of $i$


## Spectral Graph Theory

$$
\left[\begin{array}{ccc}
a_{11} & \ldots & a_{1 n} \\
\vdots & & \vdots
\end{array}\right]\left[\begin{array}{c}
x_{1} \\
\vdots
\end{array}\right]=\lambda\left[\begin{array}{c}
x_{1} \\
\vdots
\end{array}\right] \quad \boldsymbol{A} \cdot \boldsymbol{x}=\boldsymbol{\lambda} \cdot \boldsymbol{x}
$$

- Spectral Graph Theory:
" Analyze the "spectrum" of matrix representing $\boldsymbol{G}$
- Spectrum: Eigenvectors $x_{i}$ of a graph, ordered by the magnitude (strength) of their corresponding eigenvalues $\lambda_{i}: \Lambda=\left\{\lambda_{1}, \lambda_{2}, \ldots, \lambda_{n}\right\} \lambda_{1} \leq \lambda_{2} \leq \ldots \leq \lambda_{n}$

Note: We sort $\lambda_{i}$ in ascending (not descending) order!

- Spectral clustering: use the eigenvectors of $A$ or graphs derived by it (mostly graph Laplacian)


## Matrix Representations

- Adjacency matrix (A):
- $n \times n$ matrix
- $\boldsymbol{A}=\left[a_{i j}\right], a_{i j}=1$ if edge between node $\boldsymbol{i}$ and $\boldsymbol{j}$

- Important properties:

|  | 1 | 2 | 3 | 4 | 5 | 6 |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| 1 | 0 | 1 | 1 | 0 | 1 | 0 |
| 2 | 1 | 0 | 1 | 0 | 0 | 0 |
| 3 | 1 | 1 | 0 | 1 | 0 | 0 |
| 4 | 0 | 0 | 1 | 0 | 1 | 1 |
| 5 | 1 | 0 | 0 | 1 | 0 | 1 |
| 6 | 0 | 0 | 0 | 1 | 1 | 0 |

- Symmetric matrix
- Eigenvectors are real and orthogonal


## Matrix Representations

- Degree matrix (D):
- $\boldsymbol{n} \times \boldsymbol{n}$ diagonal matrix
- $D=\left[d_{i i}\right], d_{i i}=$ degree of node $i$



## Matrix Representations

- Laplacian matrix (L):
- $\boldsymbol{n} \times \boldsymbol{n}$ symmetric matrix


|  | 1 | 2 | 3 | 4 | 5 | 6 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | 3 | -1 | -1 | 0 | -1 | 0 |
| 2 | -1 | 2 | -1 | 0 | 0 | 0 |
| 3 | -1 | -1 | 3 | -1 | 0 | 0 |
| 4 | 0 | 0 | -1 | 3 | -1 | -1 |
| 5 | -1 | 0 | 0 | -1 | 3 | -1 |
| 6 | 0 | 0 | 0 | -1 | -1 | 2 |

$$
L=D-A
$$

- Laplacian matrix Limportant properties:
- Eigenvalues are non-negative real numbers
- Eigenvectors are real and orthogonal


## Example: Eigenvalues \& Eigenvectors



| Eigenvalue | 0 | 1 | 3 | 3 | 4 | 5 |
| :--- | :--- | ---: | ---: | ---: | ---: | ---: |
| Eigenvector | 1 | 1 | -5 | -1 | -1 | -1 |
|  | 1 | 2 | 4 | -2 | 1 | 0 |
|  | 1 | 1 | 1 | 3 | -1 | 1 |
|  | 1 | -1 | -5 | -1 | 1 | 1 |
|  | 1 | -2 | 4 | -2 | -1 | 0 |
|  | 1 | -1 | 1 | 3 | 1 | -1 |

## Spectral Clustering Algorithms

- Three basic stages:
- 1) Pre-processing
- Construct a matrix representation of the graph
- 2) Decomposition
- Compute eigenvalues and eigenvectors of the matrix
- Map each point to a lower-dimensional representation based on one or more eigenvectors
- 3) Grouping
- Assign points to two or more clusters, based on the new representation


## Spectral Partitioning Algorithm

- 1) Pre-processing:
- Build Laplacian matrix $L$ of the graph

- 2) 

Decomposition:

- Find eigenvalues $\lambda$ and eigenvectors $\boldsymbol{x}$ of the matrix $L$
- Map vertices to corresponding components of $\boldsymbol{\lambda}_{2}$



How do we now find the clusters?

## Spectral Partitioning

- 3) Grouping:
- Sort components of reduced 1-dimensional vector
- Identify clusters by splitting the sorted vector in two
- How to choose a splitting point?
- Naïve approaches:
- Split at 0 or median value
- More expensive approaches:
- Attempt to minimize normalized cut in 1-dimension (sweep over ordering of nodes induced by the eigenvector)

| 1 | 0.3 |
| :---: | :---: |
| 2 | 0.6 |
| 3 | 0.3 |
| 4 | -0.3 |
| 5 | -0.3 |
| 6 | -0.6 |

Split at 0:
Cluster A: Positive points
Cluster B: Negative points

| 1 | 0.3 |
| :--- | :--- |
| 2 | 0.6 |
| 3 | 0.3 |$\quad$| 4 | -0.3 |
| :---: | :---: |
| 5 | -0.3 |
| 6 | -0.6 |



## Example: Spectral Partitioning




## Example: Spectral Partitioning



Components of $x_{2}$


Rank in $\mathbf{x}_{\mathbf{2}}$

## Example: Spectral Partitioning





## k-Way Spectral Clustering

- How do we partition a graph into $k$ clusters?
- Two basic approaches:
- Recursive bi-partitioning [Hagen et al., '92]
- Recursively apply bi-partitioning algorithm in a hierarchical divisive manner
- Disadvantages: Inefficient, unstable
- Cluster multiple eigenvectors [Shi-Malik, '00]
- Build a reduced space from multiple eigenvectors
- Commonly used in recent papers
- A preferable approach...


## Recursive Bi-partitioning



## Cluster Multiple Eigenvectors

- Use several of the eigenvectors to partition the graph
- If we use m eigenvectors, and set a threshold for each, we can get a partition into $2^{m}$ groups, each group consisting of the nodes that are above or below threshold for each of the eigenvectors, in a particular pattern.


## Example



| Eigenvalue | 0 | 1 | 3 | 3 | 4 | 5 |
| :--- | :--- | ---: | ---: | ---: | ---: | ---: |
| Eigenvector | 1 | 1 | -5 | -1 | -1 | -1 |
|  | 1 | 2 | 4 | -2 | 1 | 0 |
|  | 1 | 1 | 1 | 3 | -1 | 1 |
|  | 1 | -1 | -5 | -1 | 1 | 1 |
|  | 1 | -2 | 4 | -2 | -1 | 0 |
|  | 1 | -1 | 1 | 3 | 1 | -1 |

If we use both the $\mathbf{2}^{\text {nd }}$ and $3^{\text {rd }}$ eigenvectors:

- nodes 2 and $\mathbf{3}$ (positive in both)
- nodes 5 and 6 (negative in $2^{\text {nd }}$, positive in $3^{\text {rd }}$ )
- nodes 1 and 4 alone

Note that while each eigenvector tries to produce a minimum-sized cut, successive eigenvectors have to satisfy more and more constraints => the cuts progressively worse.

## Why use multiple eigenvectors?

- Approximates the optimal cut [Shi-Malik, '00]
- Can be used to approximate optimal $k$-way normalized cut
- Emphasizes cohesive clusters
- Increases the unevenness in the distribution of the data
- Associations between similar points are amplified, associations between dissimilar points are attenuated
" The data begins to "approximate a clustering"
- Well-separated space
- Transforms data to a new "embedded space", consisting of $\boldsymbol{k}$ orthogonal basis vectors
- Multiple eigenvectors prevent instability due to information loss


## Many Other Partitioning Methods

- METIS:
- Heuristic but works really well in practice
- http://glaros.dtc.umn.edu/gkhome/views/metis
- Graclus:
- Based on kernel k-means
- http://www.cs.utexas.edu/users/dml/Software/graclus.html
- Louvain:
- Based on Modularity optimization
- http://perso.uclouvain.be/vincent.blondel/research/louvain.html
- Clique percolation method:
- For finding overlapping clusters
- http://angel.elte.hu/cfinder/


## How to Profile Network Communities?

## Network and Communities

- How should we think about large scale organization of clusters in networks?
- Finding: Community Structure



## Community Score

- How community-like is a set of nodes?
- A good cluster $S$ has
- Many edges internally
- Few edges pointing outside
- What's a good metric:

Conductance
$\phi(S)=\frac{|\{(i, j) \in E ; i \in S, j \notin S\}|}{\sum_{s \in S} d_{s}}$


Small conductance corresponds to good clusters
(Note $|\mathrm{S}|<|\mathrm{V}| / 2)$

## Network Community Profile Plot

- Define:

Network community profile (NCP) plot Plot the score of best community of size $k$

$$
\Phi(k)=\min _{c} \phi(S)
$$

$$
S \subset V,|S|=k
$$




## How to (Really) Compute NCP?

dblp-lars


## NCP Plot: Meshes

- Meshes, grids, dense random graphs:



California road network

## NCP plot: Network Science

- Collaborations between scientists in networks
[Newman, 2005]



Community size, log k
Dips in the conductance graph correspond to the "good" clusters we can visually detect

## Large Networks: Very Different

## Typical example: General Relativity collaborations ( $\mathrm{n}=4,158, \mathrm{~m}=13,422$ )


[Internet Mathematics 'og]

## More NCP Plots of Networks


(a) LIVEJOURNALO1


(b) MESSENGER-DE
 k (number of nodes in the cluster)
(e) WEB-GOOGLE


k (number of nodes in the cluster)
(f) AmAZONALL
-- Rewired graph
-- Real graph

## NCP: LiveJournal (n=5m, m=42m)




## Communities:

Issues and Questions

## What is Cluster Analysis?

Finding groups of objects such that the objects in a group will be similar (or related) to one another and different from (or unrelated to) the objects in other groups


## Clusters Can Be Ambiguous



How many clusters?


Six Clusters


Two Clusters



Four Clusters

## Communities: Issues and Questions

- Some issues with community detection:
- Many different formalizations of clustering objective functions
- Objectives are NP-hard to optimize exactly
- Methods can find clusters that are systematically "biased"
- Questions:
- How well do algorithms optimize objectives?
- What clusters do different methods find?


## Many Different Objective Functions

- Single-criterion:
- Modularity: $m-E(m)$
- Edges cut: $c$
- Multi-criterion:
- Conductance: $c /(2 m+c)$
- Expansion: c/n
- Density: $1-m / n^{2}$

$n$ : nodes in S
$m$ : edges in S
$c$ : edges pointing outside S
- CutRatio: $c / n(N-n)$
- Normalized Cut: $c /(2 m+c)+c / 2(M-m)+c$
- Flake-ODF: frac. of nodes with more than $1 / 2$ edges pointing outside $S$


## Many Classes of Algorithms

Many algorithms to implicitly or explicitly optimize objectives and extract communities:

- Heuristics:
- Girvan-Newman, Modularity optimization: popular heuristics
- Metis: multi-resolution heuristic [Karypis-Kumar '98]
- Theoretical approximation algorithms:
- Spectral partitioning

500 node communities from Spectral:


500 node communities from Metis:


## Properties of Clusters (2)



- Metis gives sets with better conductance
- Spectral gives tighter and more well-rounded sets



## Single-criterion Objectives



Observations:

- All measures are monotonic
- Modularity
- prefers large clusters
- Ignores small clusters


## Multi-criterion Objectives



## All qualitatively similar <br> Observations:

- Conductance, Expansion, Normcut, Cut-ratio are similar
- Flake-ODF prefers larger clusters
- Density is bad
- Cut-ratio has high variance

Normalized Cut Maximum ODF

Avg ODF Flake ODF

