

recall:

- graph Laplacian $\Rightarrow L = D - A$

- $Lv = \lambda v$
 \swarrow all eigenvectors
 \nwarrow eigenvalues

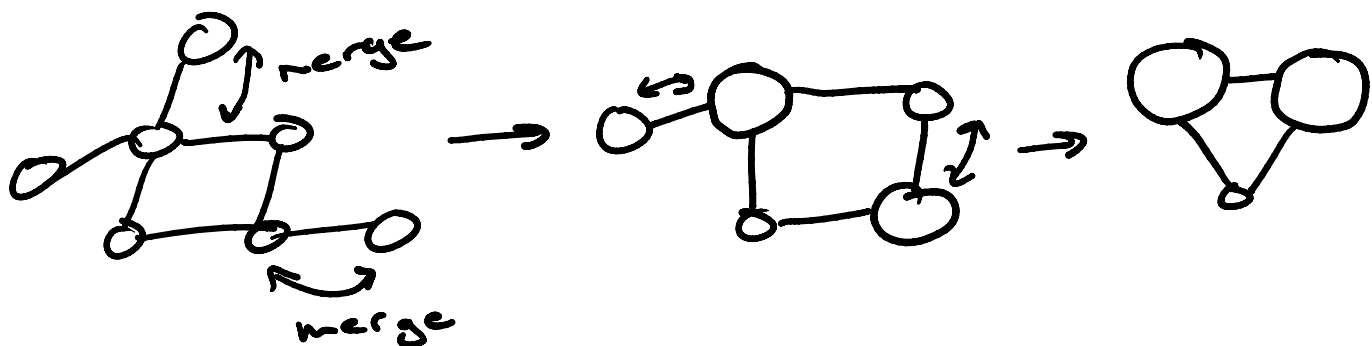
\hookrightarrow we can use these to cut
 our graph into clusters
 (equal sized or not)

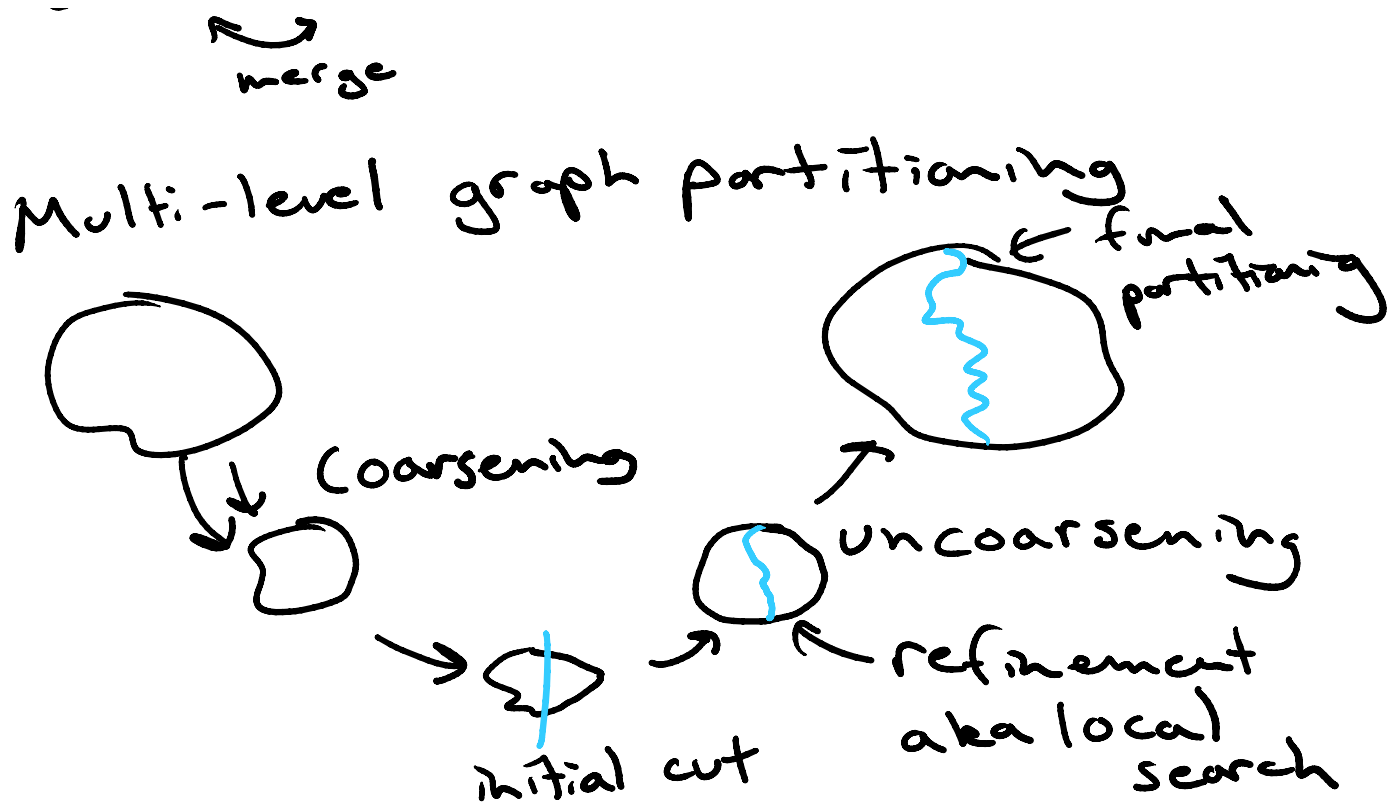
Coarsening

Generally: graph partitioning is hard
 (NP-hard)

What we want to do: transform our
 input to a smaller/easier problem

How? Coarsening





Coarsening: merging vertices via edge contractions to construct a smaller representation of the input

→ many application-specific ways to do this

→ for partitioning, we can preserve spectral properties

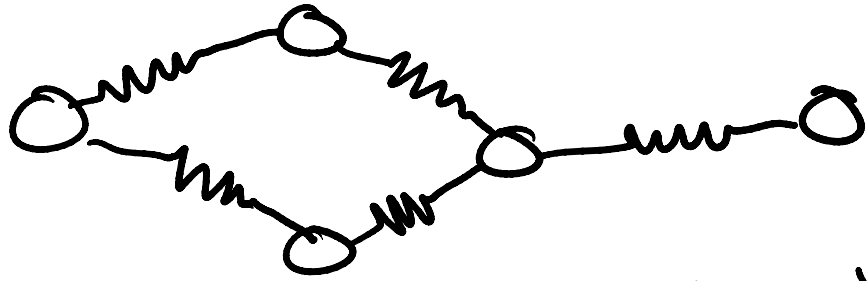
* How? selecting vertices to merge that minimizes change in the graph spectrum
(vertices with similar eigenvector values)

Another application: visualization

A nother application: visualization

Visualization: plotting a graph in
2D (or 3D) space

Common approach: model edges as springs



too close: force push vertizes apart
too far: force pulls vertizes together

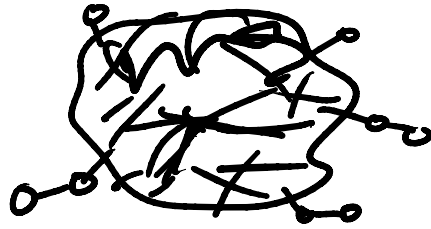
A nother approach: using the spectrum

recall: spectrum is essentially
mapping from discrete to
continuous space

Using two eigenvector values, we
can map a graph into 2D space
by interpreting the eigenvector
values as coordinates

One major issue: large inputs

↳ hairball



Mitigating: can use coarsening
e.g. can coarsen communities
to expose underlying structure

We can use spectral coarsening and
visualization

Coarsen:

Find n pairs of vertices with closest
spectral values and coarsen them

Visualization:

Interpret the two largest eigenvector
values for each vertex as (x, y)
coordinates in 2D space

Random graphs

What is a random graph?

How many $|V|$?

How many $|E|$?

How are edges created?
(attachment probability)

(attachment probability)

Big idea:

- Randomly configured network
- Edges determined randomly

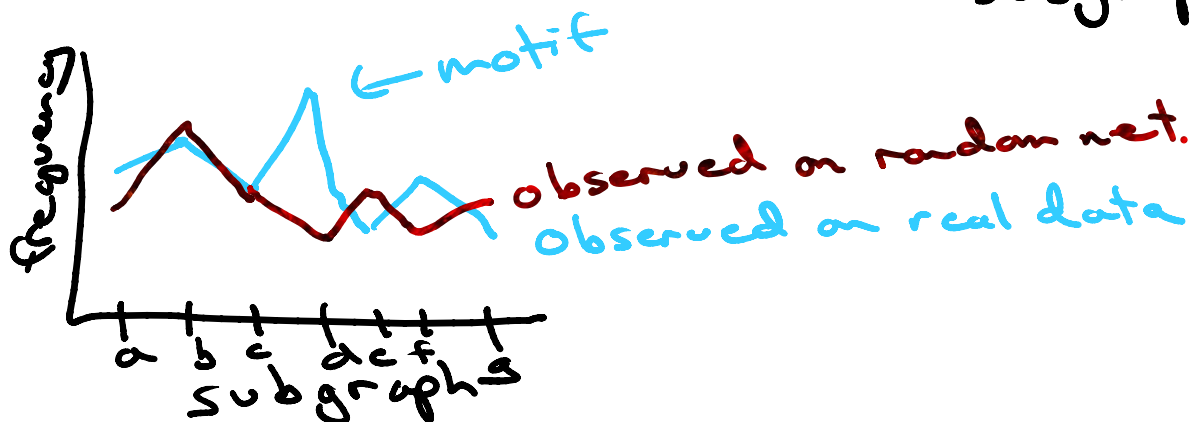
Why do we care:

- Mirror properties of real networks to analytically study
- Use as a "null model" for hypothesis testing or otherwise

E.g. in modularity we compare our observation relative to a randomly configured network with the same degree sequence $(\frac{d_i d_j}{2m})$

E.g. motif finding

motif = a frequently occurring subgraph



^a ³ ^{act}
subgraphs

How do we define a random graph?

Classic model: Erdős-Rényi

$$O.G.: G(n, m) \quad \langle k \rangle = \frac{2m}{n}$$

\uparrow \uparrow
 $|V|$ $|E|$ avg. degree

Issue: generates loopy multigraphs

Newer model: $G(n, p)$ $\langle k \rangle = p(n-1)$

\uparrow
attachment probability

↳ This is a Bernoulli process

Hence, this gives us a binomial distribution for the degree distribution

↳ How many vertices of degree k
(degree k , #vertices w/degree k)

Degree distribution

$$P(k) = \binom{n-1}{k} p^k (1-p)^{(n-1)-k}$$

\uparrow
prob that vertex has degree k

prob that vertex
has degree k

Note: as $n \rightarrow \infty$ and k is fixed

Binomial \rightarrow Poisson

Poisson: $P(k) = \frac{e^{-\langle k \rangle} \langle k \rangle^k}{k!}$ \leftarrow mean value

Why: approximation is good

$n \gg \langle k \rangle$ in real networks

1 fewer parameter

Let's analyze Erdős-Rényi connectivity
in terms of a **GIANT** component

\rightarrow what $\langle k \rangle$ do we need?

$\langle k \rangle = 0 \rightarrow$ completely disconnected

$\langle k \rangle = n-1 \rightarrow$ completely connected

In reality, around $\langle k \rangle = 1$ we observe
emergence of giant component

(see B. section 3.C)

\rightarrow "critical point"

\rightarrow real networks \checkmark

→ critical point

This mirrors real-world networks ✓

What about other real-world properties

- small-world
- low diameter
- Hubbedness
- degree skew

Can we quantify these for an E-K graph?

- consider vertex v
- v has degree of $\langle k \rangle$
- each of v 's neighbors has $\langle k \rangle$ neighbors
- 2-hop neighborhood around v has $\langle k \rangle \langle k \rangle = \langle k \rangle^2$
- 3-hop neighborhood = $\langle k \rangle^3$

$$|N_d(v)| = \langle k \rangle + \langle k \rangle^2 + \langle k \rangle^3 \dots \langle k \rangle^d$$

↑
d-hop neighborhood

$$\approx \frac{\langle k \rangle^{d+1} - 1}{\langle k \rangle - 1}$$

$n = \dots$

To get an estimate for diameter, we can find how many hops in which we would expect to have all vertices

set $|N_d(u)| = n$, solve for d

$$\frac{\langle k \rangle^{d+1} - 1}{\langle k \rangle - 1} = n$$

$$\langle k \rangle^d \approx n$$

$$d \approx \frac{\ln(n)}{\ln(\langle k \rangle)}$$

$$\text{as } n \gg \langle k \rangle \rightarrow d \approx \ln(n)$$

SO IN OTHER WORDS

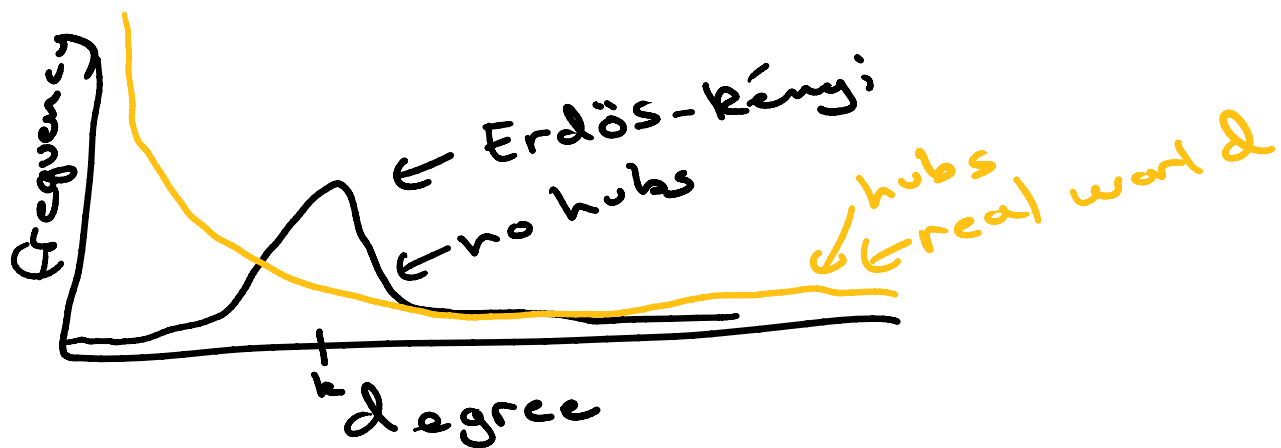
our diameter grows logarithmically
as a function of graph order

↳ that's a small world graph ✓

(and avg. shortest paths grow $\ln(N)$ as well)

Big issue w/ E-R

is the degree distribution



In reality $\langle k \rangle \ll k_{max}$

Introducing: the configuration model

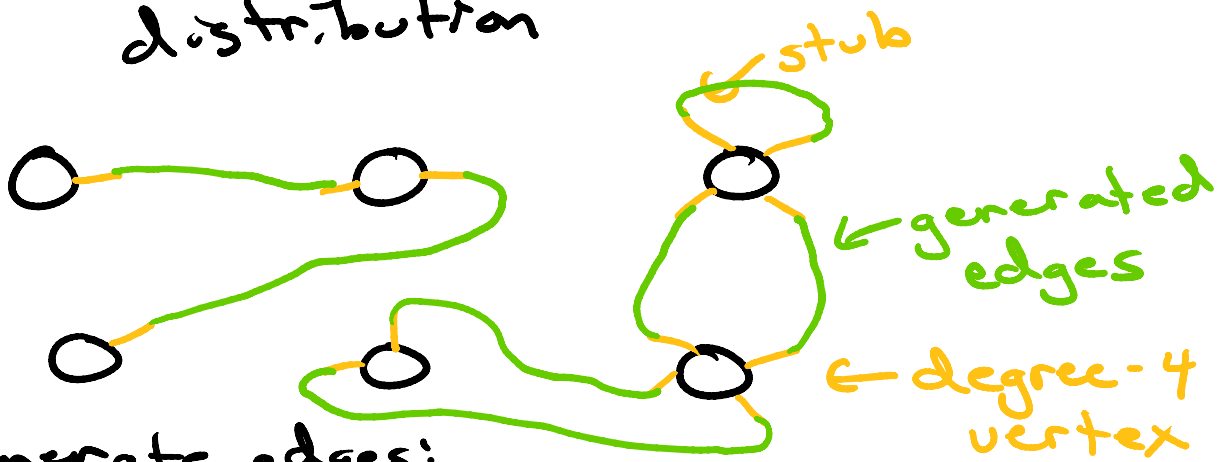
- we can generate a random network with an arbitrary degree distribution

Basic idea

- we have n vertices each with some number of "stubs"
- # of stubs follows the degree distribution

↙ stub

distribution



to generate edges:

- We select two stubs randomly and add an edge between them

What are the attachment probabilities?

consider i j or \cancel{i} \cancel{j}

Note: more likely to select high degree vertex's stubs

Attachment probability is a function of $d(i)$ and $d(j)$ and the sum of the degrees

$$\begin{aligned} \text{prob. of creating edge}(i, j) &= (\text{prob of selecting } i\text{'s stub}) \\ &\quad * (\text{prob of selecting } j\text{'s stub}) \end{aligned}$$

\cdot * (prob of selecting j 's stub)
 \cdot * $Z(i, j), (j, i)$
 \cdot * m (total attempts)

$$\rightarrow P_i P_j Z_m$$

$$= \frac{d(i)}{Z_m} \frac{d(j)}{Z_m} Z_m$$

$$= \frac{d(i) d(j)}{Z_m} \leftarrow \text{as we've seen with modularity}$$

Note: a lot of the same properties hold as with E-R graph, but we also can model skewed distributions and hubs ✓

Issue 1: tough to generate simple graphs

Issue 2: no clustering