



# HIERARCHICAL BLOCK MODELS

Lecture 7, Fall 2014  
CSCI 5352, Network Analysis and Models

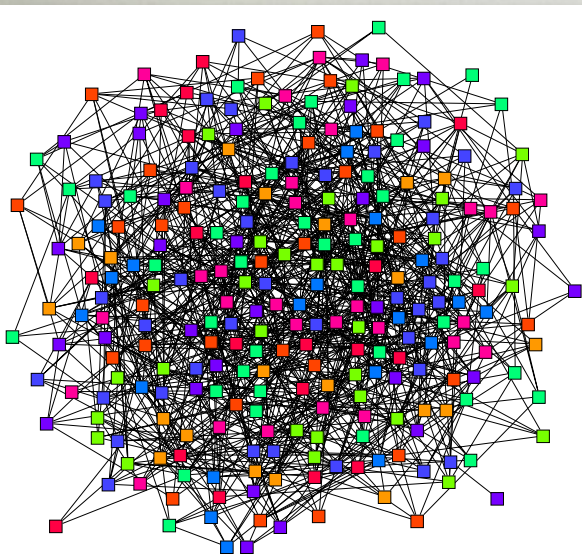
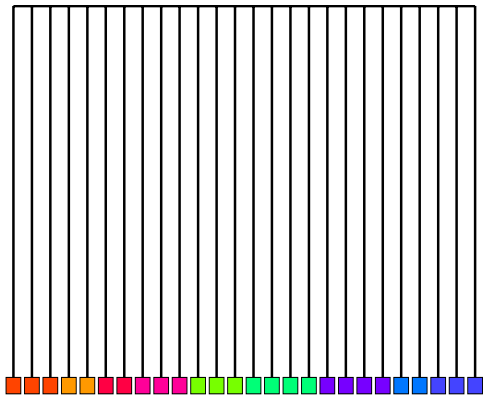
Prof. Aaron Clauset  
University of Colorado, Boulder

# ADVANCED GENERATIVE MODELS

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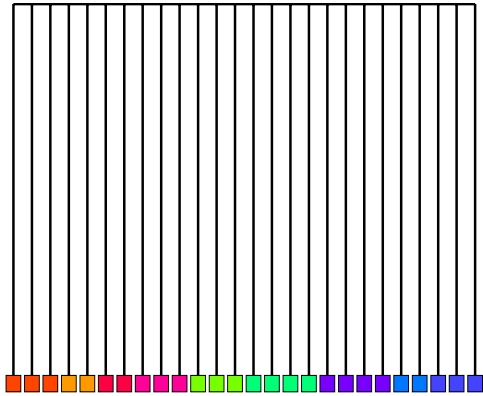
1. hierarchical random graph model example
2. examples of tasks generative models can do
3. application to some real-world data sets

**no structure**

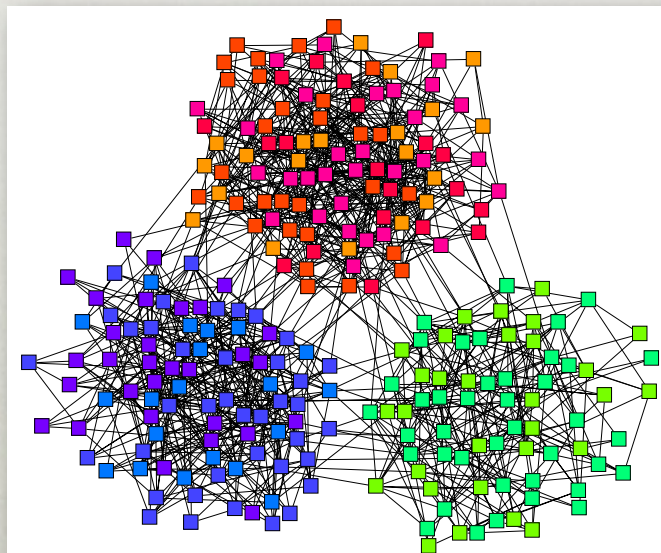
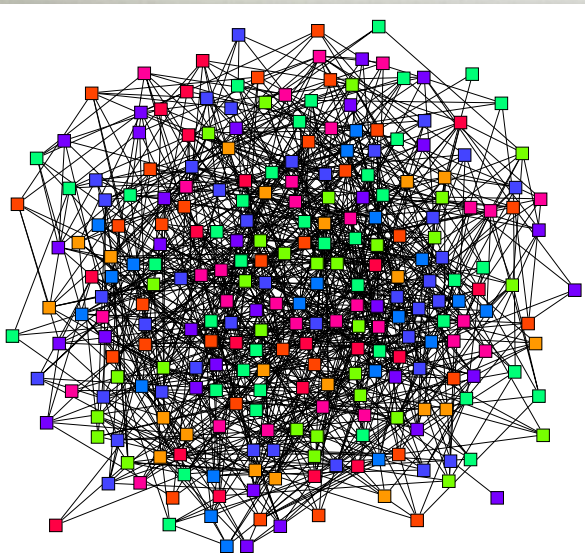
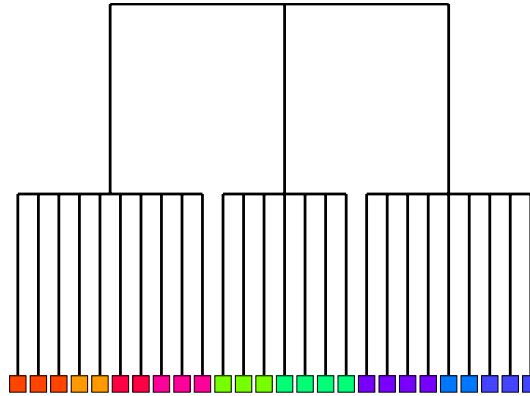




**no structure**

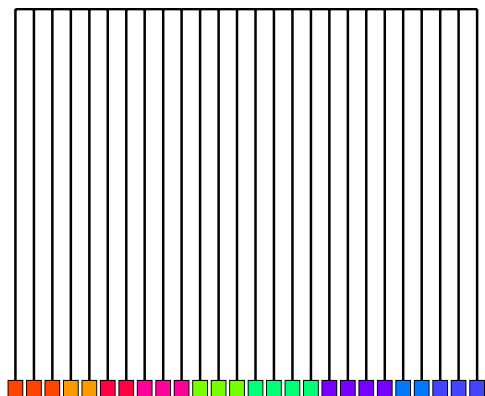


**block structure**

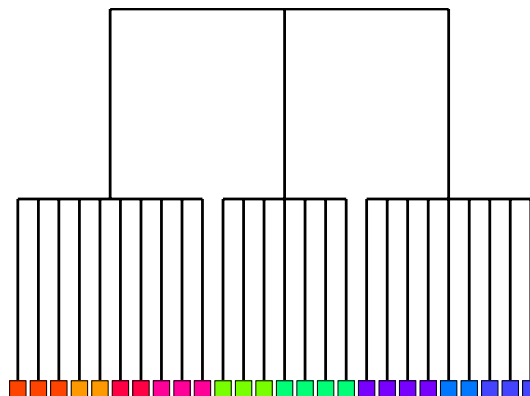


**one scale**

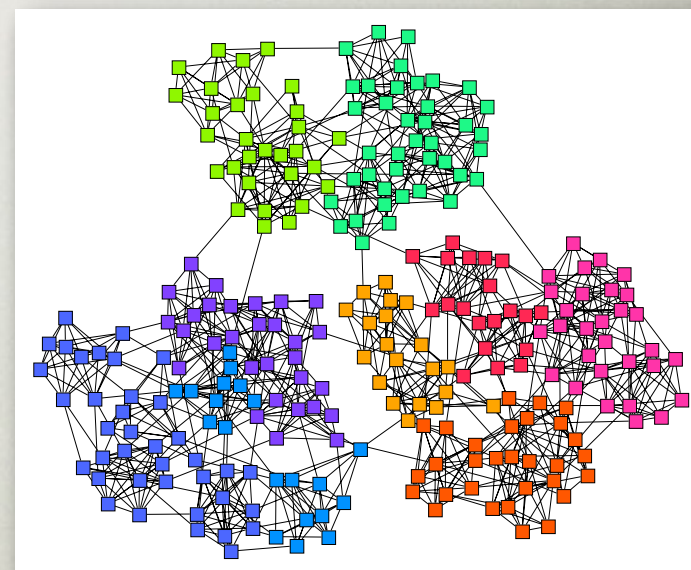
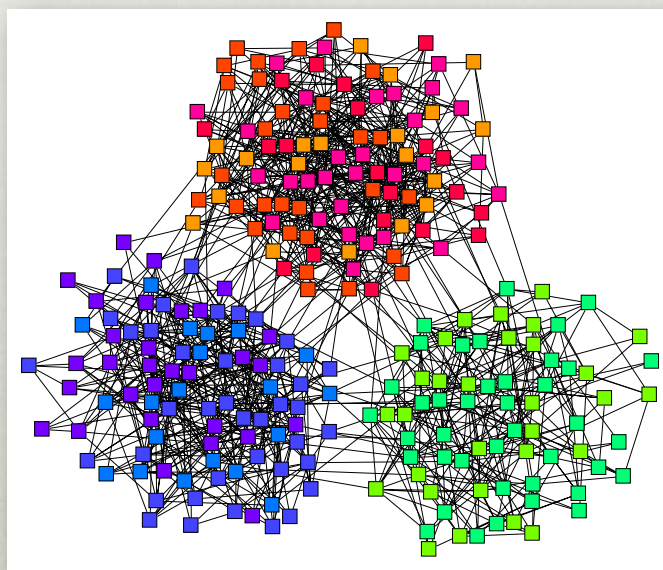
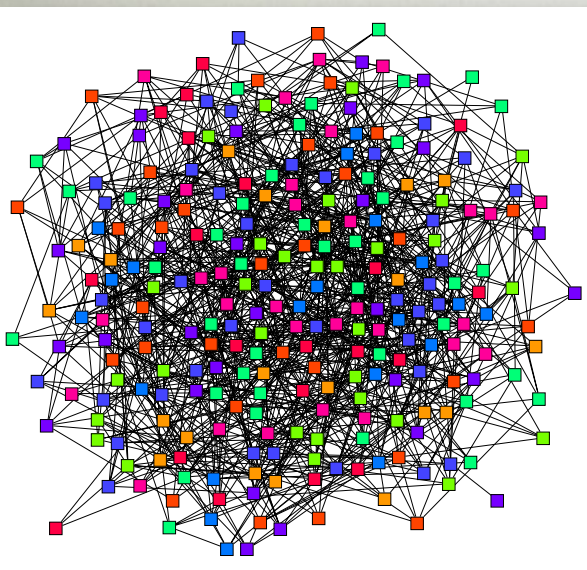
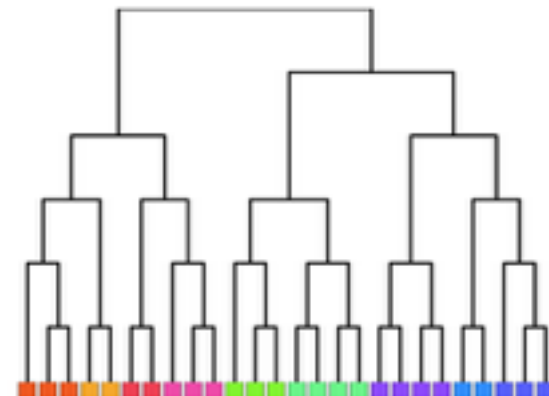
**no structure**



**block structure**



**hierarchical structure**

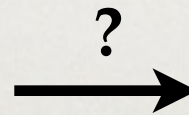
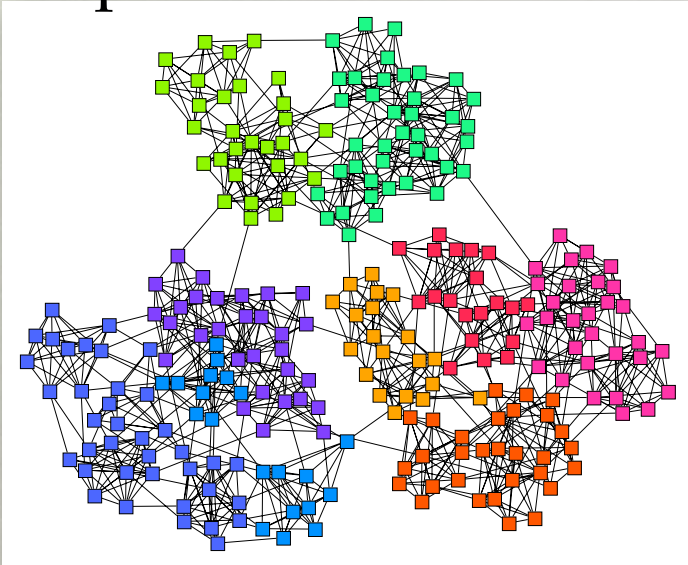


**one scale**

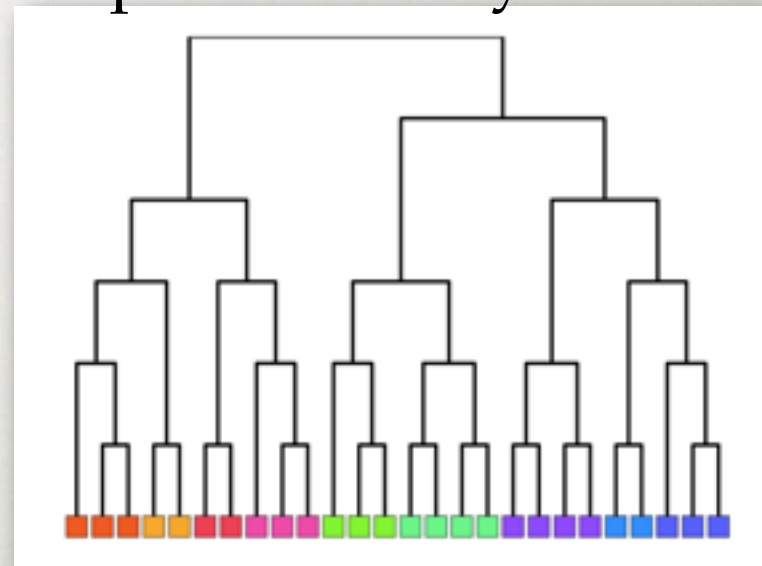
**multi-scale**

how can we extract a network's hierarchy?

step 1: network data



step 3: hierarchy





# GENERATIVE MODELS

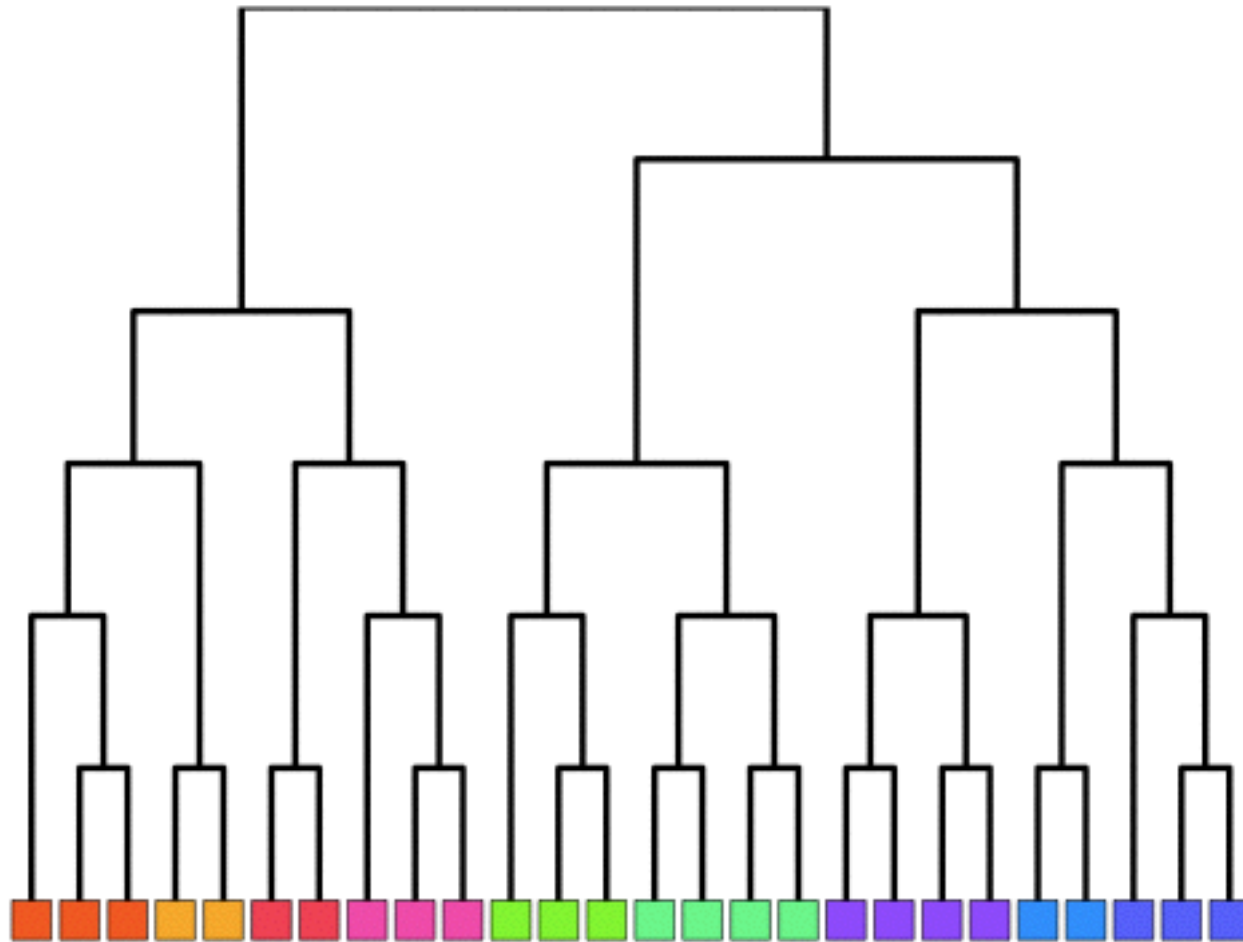
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1. write down hierarchical stochastic block model
2. estimate / learn model from data
3. evaluate model goodness-of-fit
4. evaluate model predictions

# A MODEL OF HIERARCHY

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$\mathcal{D}$

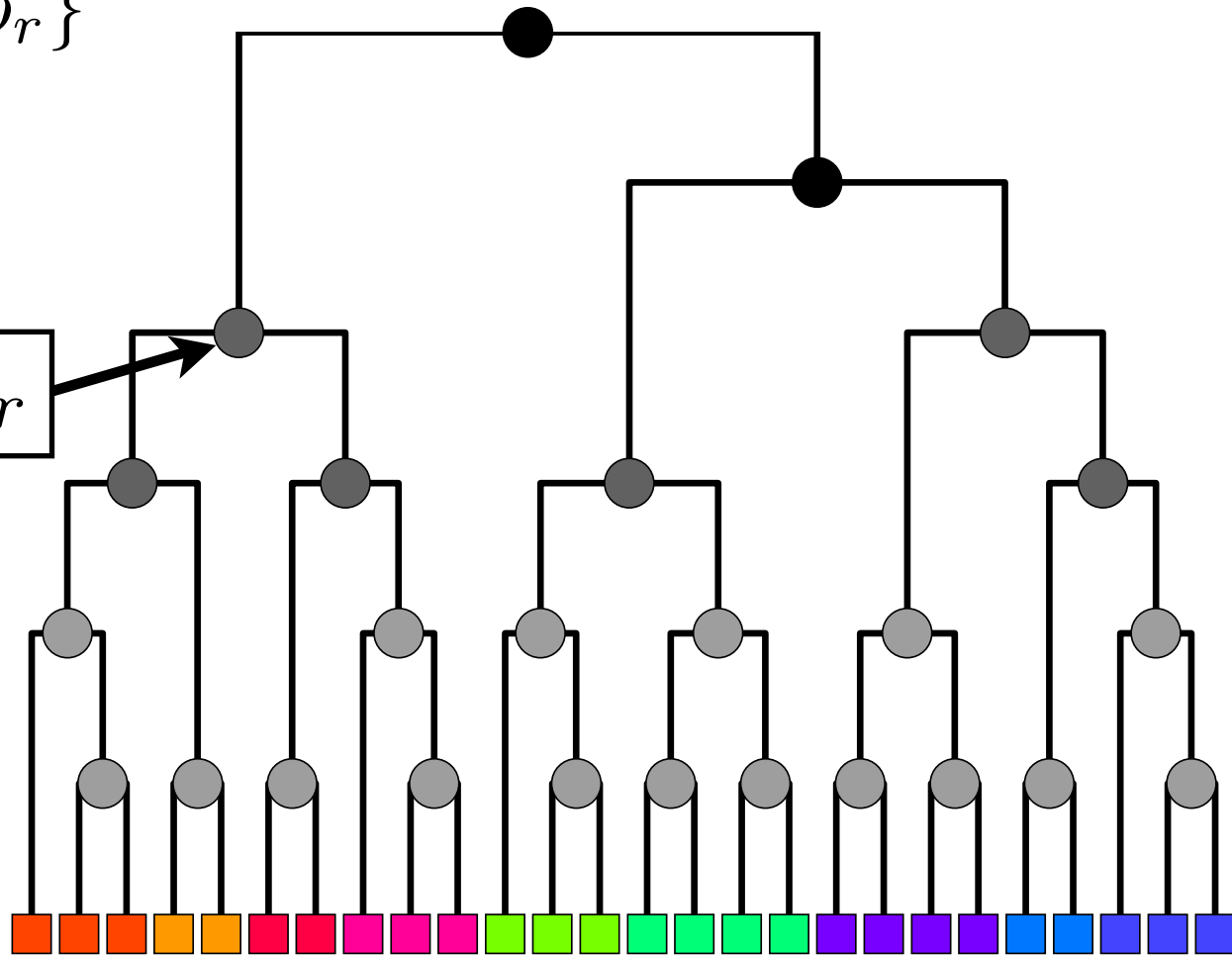




# A MODEL OF HIERARCHY

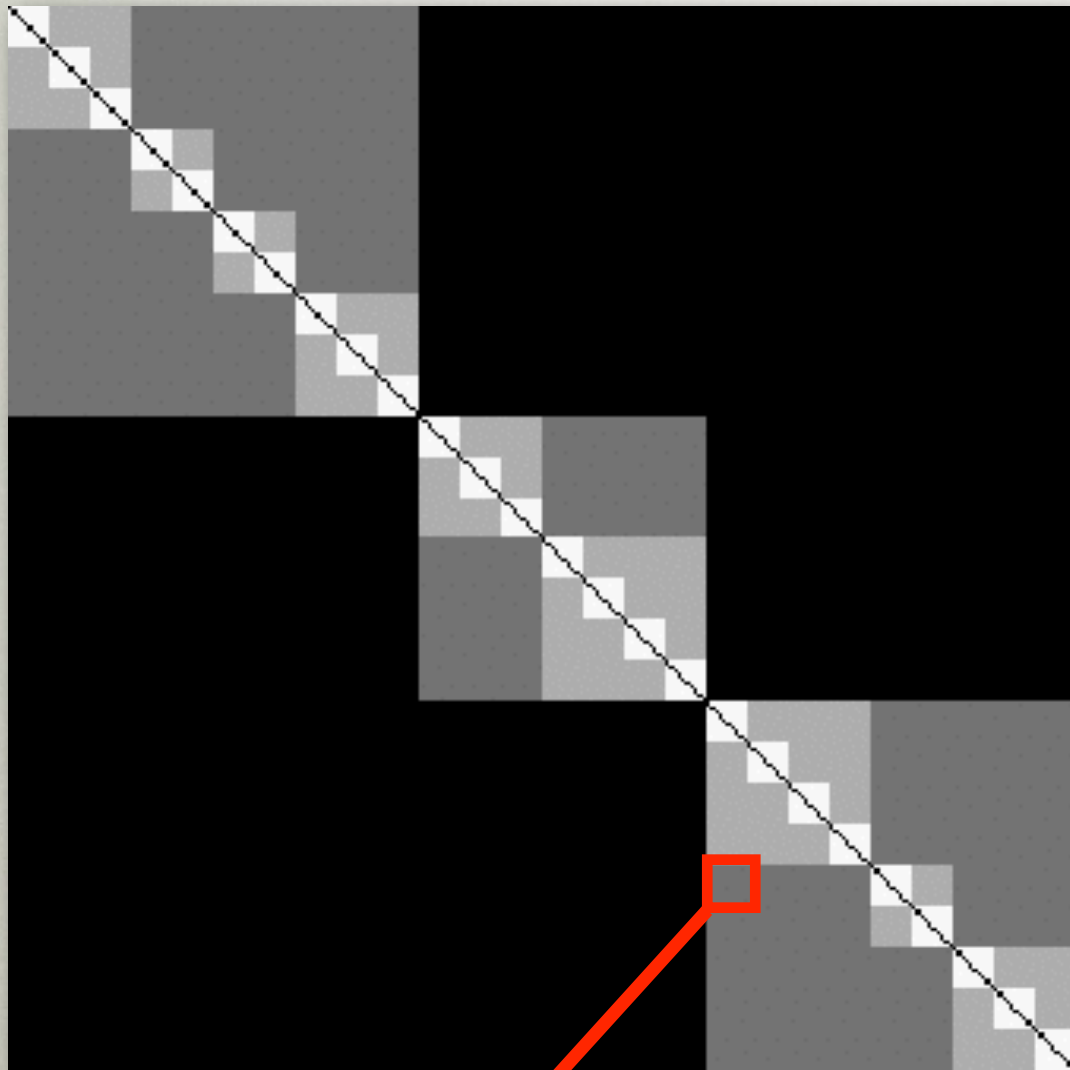
$\mathcal{D}, \{p_r\}$

probability  $p_r$



assortative modules

“inhomogeneous” random graph

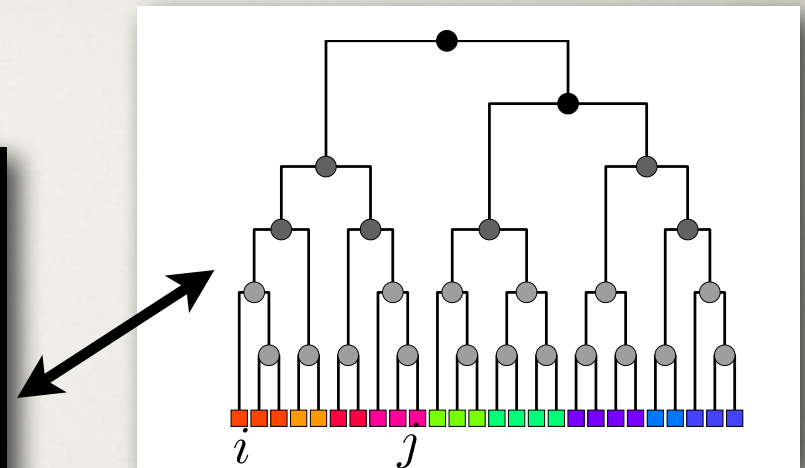


$j$

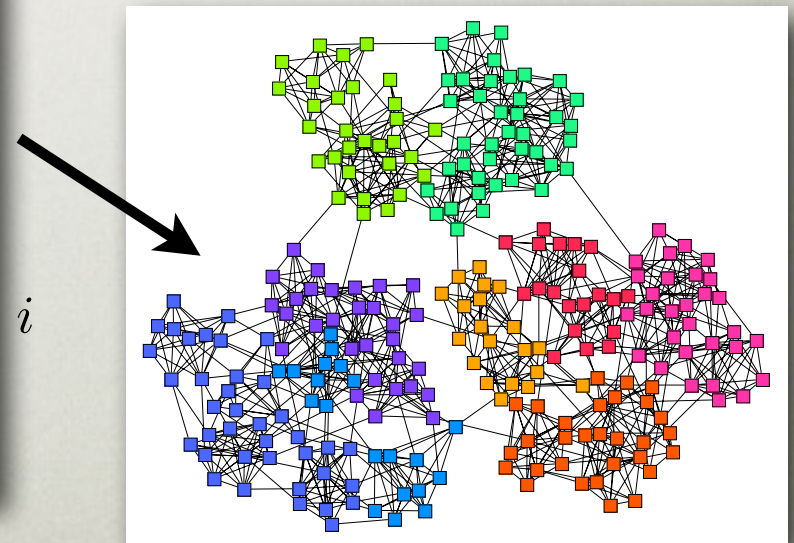
$$\Pr(i, j \text{ connected}) = p_r$$

$$= p_{(\text{lowest common ancestor of } i, j)}$$

model



instance



# HIERARCHICAL RANDOM GRAPH

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## *advantages*

- explicit hierarchical structure
- flexible ( $2n$  parameters)
- captures structure at all scales
- captures mixtures of assortativity, disassortativity
- like SBM, decomposes adjacencies into "bundles" (each a random bipartite graph)
- learnable directly from data
- nice interpretable structure

## *disadvantages*

- flexible ( $2n$  parameters)
- computationally slow
- can overfit to degree structure (like SBM)



# FITTING THE MODEL

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- **likelihood function**  $\mathcal{L} = \Pr(\text{ data } | \text{ model } )$   
(  $\mathcal{L}$  scores **quality** of model)
- **sample all good models**  
via Markov chain Monte Carlo\*  
over all dendrograms
- **technical details in**

Clauset, Moore and Newman, *Nature* **453**, 98-101 (2008) and

Clauset, Moore and Newman, *ICML* (2006)

\* other sampling or optimization methods possible

# LIKELIHOOD FUNCTION

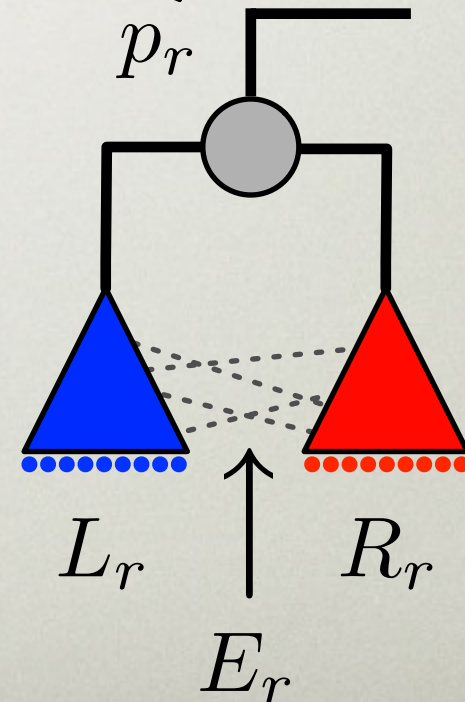
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$$\mathcal{L}(\mathcal{D}, \{p_r\}) = \prod_r p_r^{E_r} (1 - p_r)^{L_r R_r - E_r}$$

$L_r$  = number nodes in left subtree

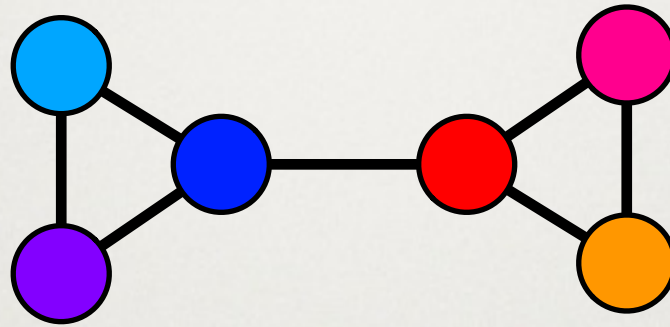
$R_r$  = number nodes in right subtree

$E_r$  = number edges with  $r$  as lowest common ancestor



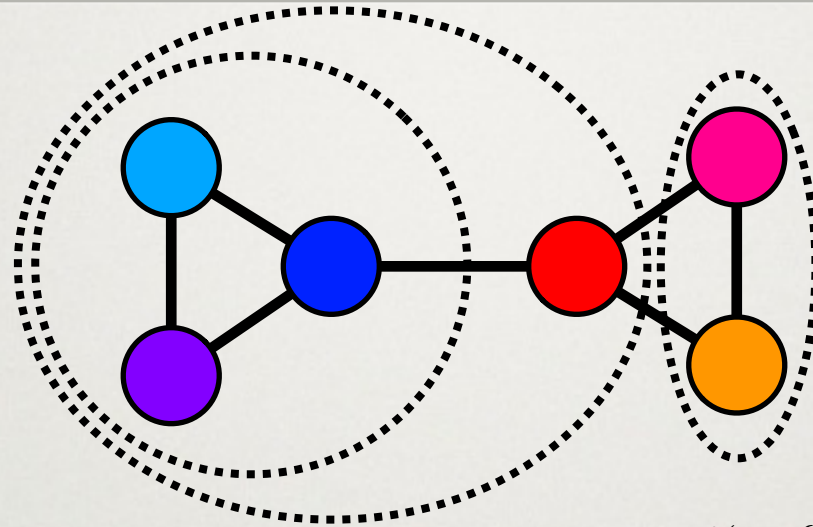
# EXAMPLE

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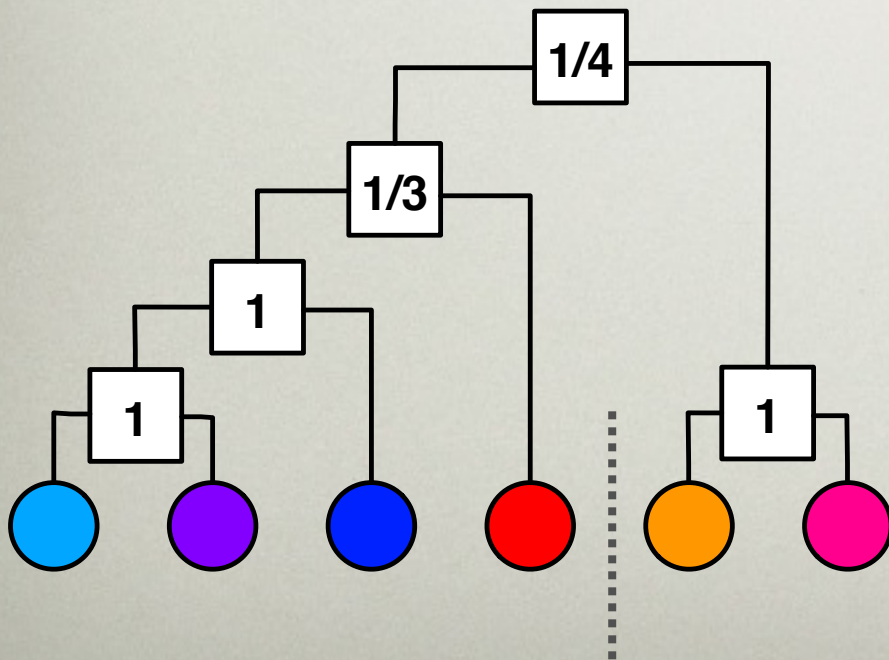




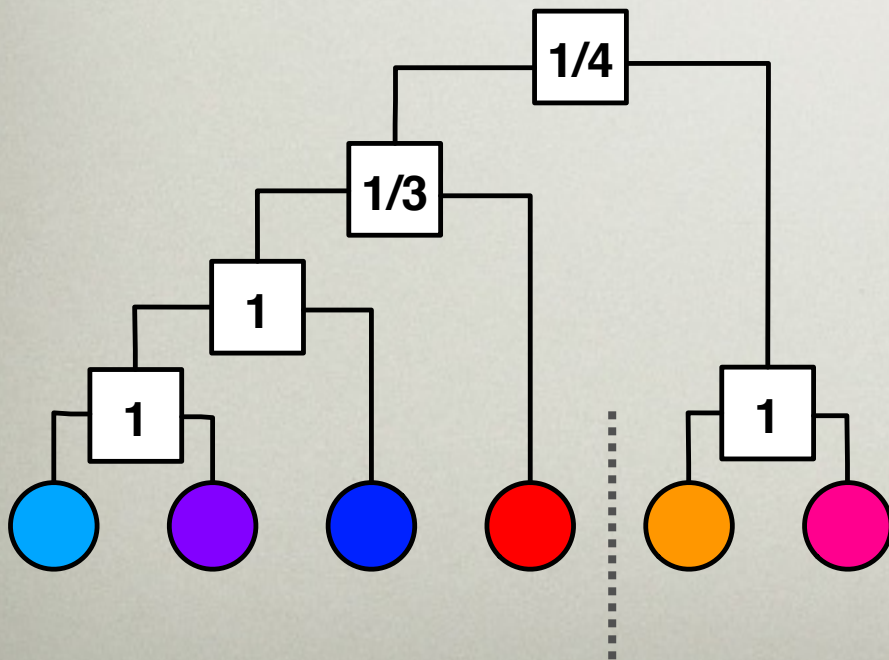
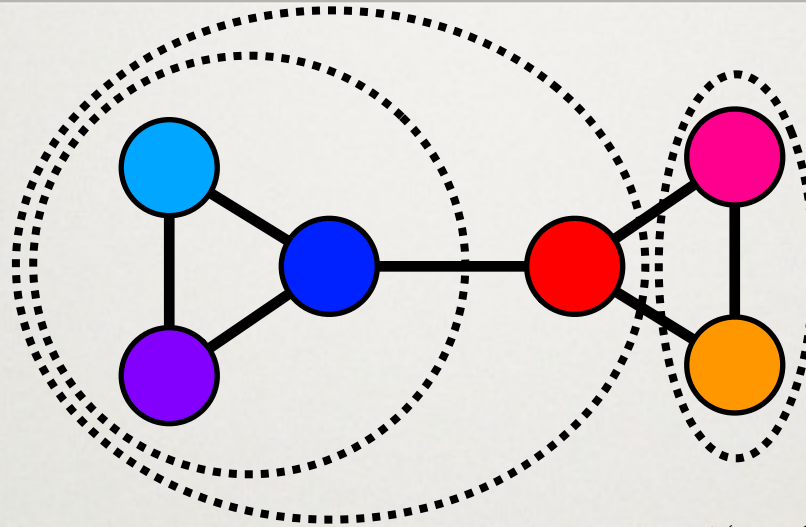
# BAD DENDROGRAM



$$\mathcal{L}(\mathcal{D}, \{p_r\}) = \prod_r p_r^{E_r} (1 - p_r)^{L_r R_r - E_r}$$



# BAD DENDROGRAM



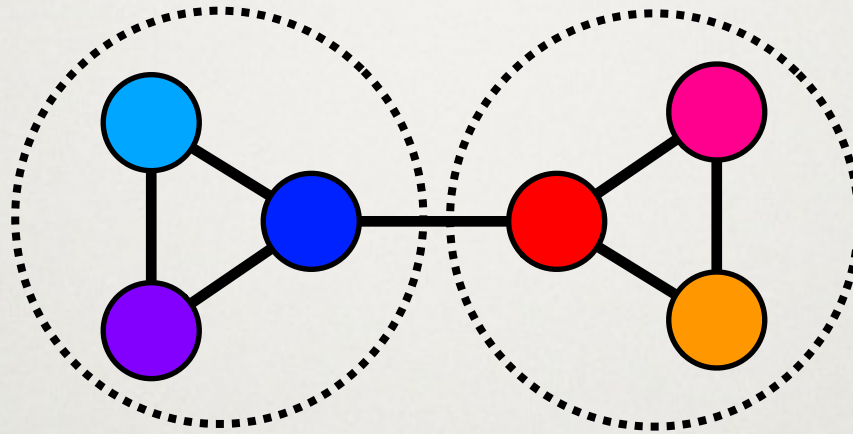
$$\mathcal{L}(\mathcal{D}, \{p_r\}) = \prod_r p_r^{E_r} (1 - p_r)^{L_r R_r - E_r}$$

$$\mathcal{L} = \left[ \left( \frac{1}{3} \right)^1 \left( \frac{2}{3} \right)^2 \right] \cdot \left[ \left( \frac{1}{4} \right)^2 \left( \frac{3}{4} \right)^6 \right]$$

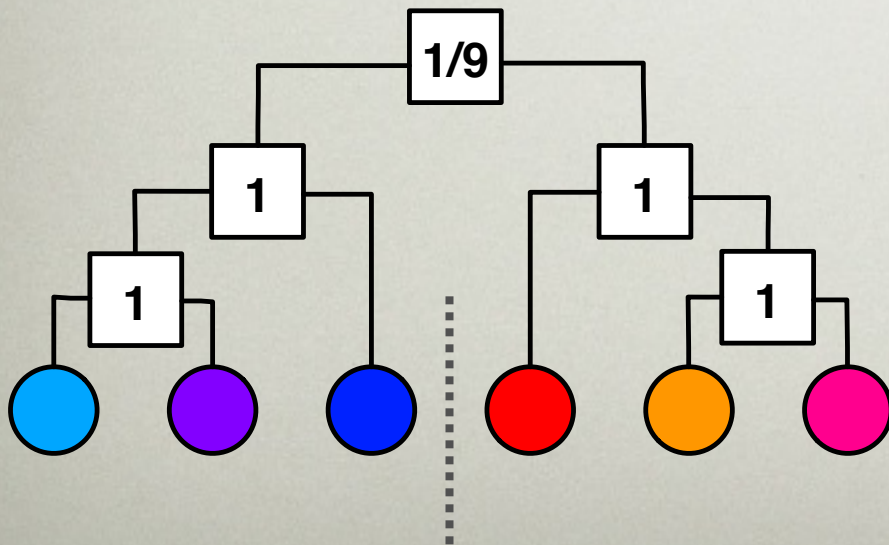
$$\mathcal{L} = 0.0016$$

# GOOD DENDROGRAM

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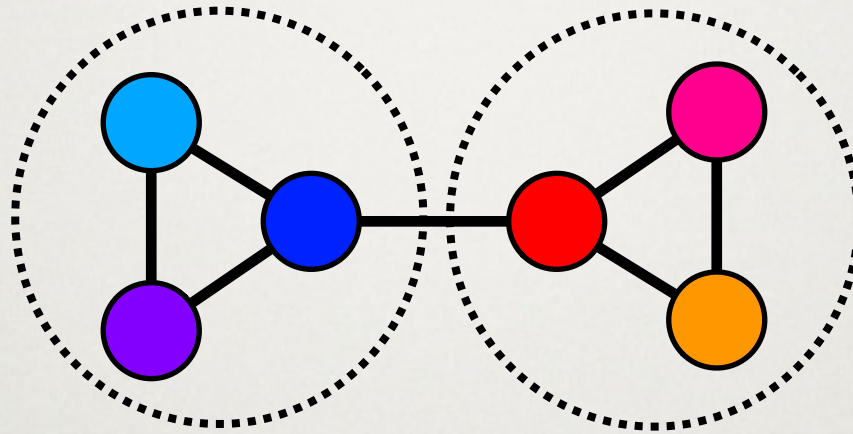


$$\mathcal{L}(\mathcal{D}, \{p_r\}) = \prod_r p_r^{E_r} (1 - p_r)^{L_r R_r - E_r}$$





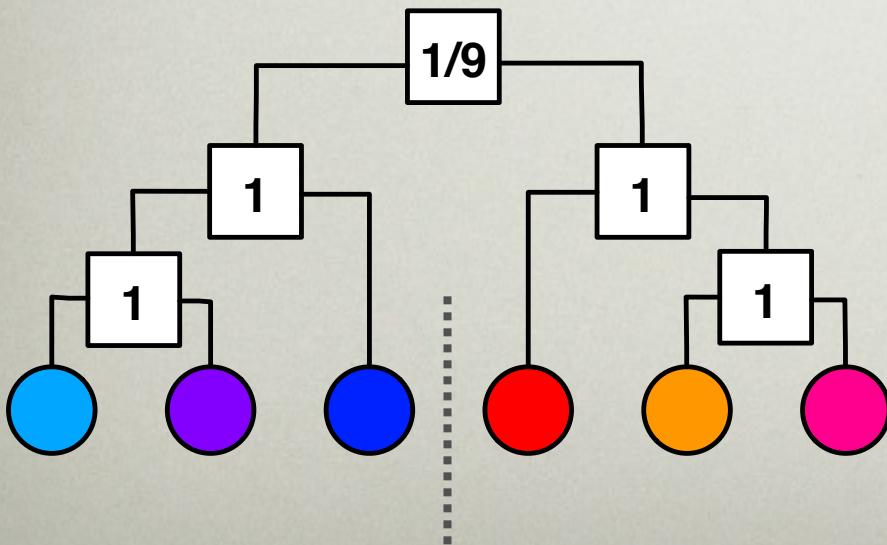
# GOOD DENDROGRAM



$$\mathcal{L}(\mathcal{D}, \{p_r\}) = \prod_r p_r^{E_r} (1 - p_r)^{L_r R_r - E_r}$$

$$\mathcal{L} = \left[ \left( \frac{1}{9} \right)^1 \left( \frac{8}{9} \right)^8 \right]$$

$$\mathcal{L} = 0.0433$$



# MARKOV CHAIN MONTE CARLO (MCMC)

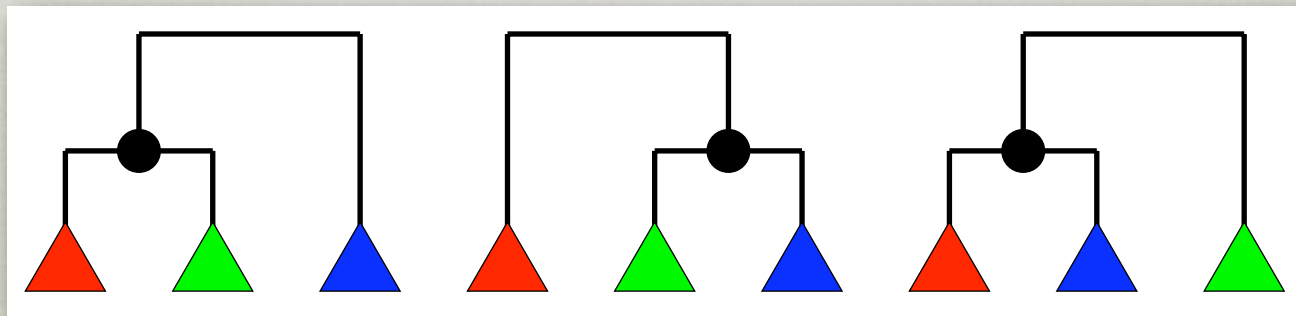
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Given  $\mathcal{D}$ , choose random internal node

Choose random reconfiguration of subtrees [ergodicity]

Recompute probabilities  $\{p_r\}$  and likelihood  $\mathcal{L}$

Sampling states according to their likelihood [detailed balance]



three subtree configurations  
(up to relabeling)

# SOME APPLICATIONS

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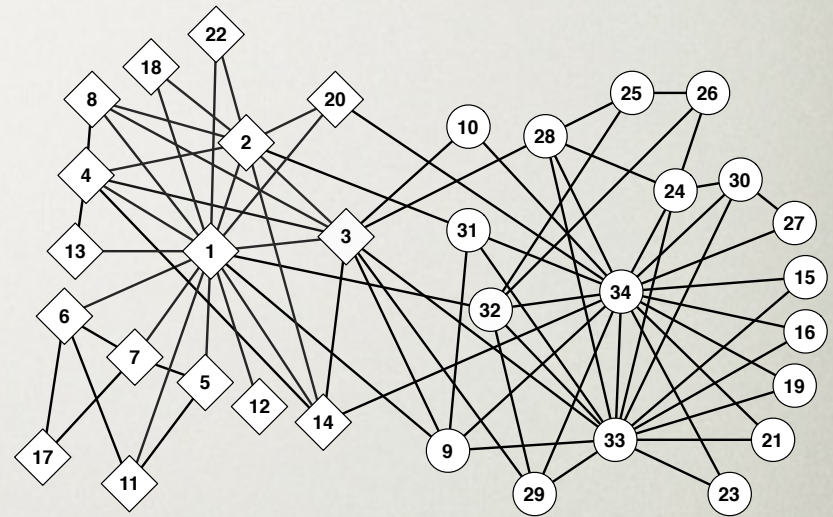
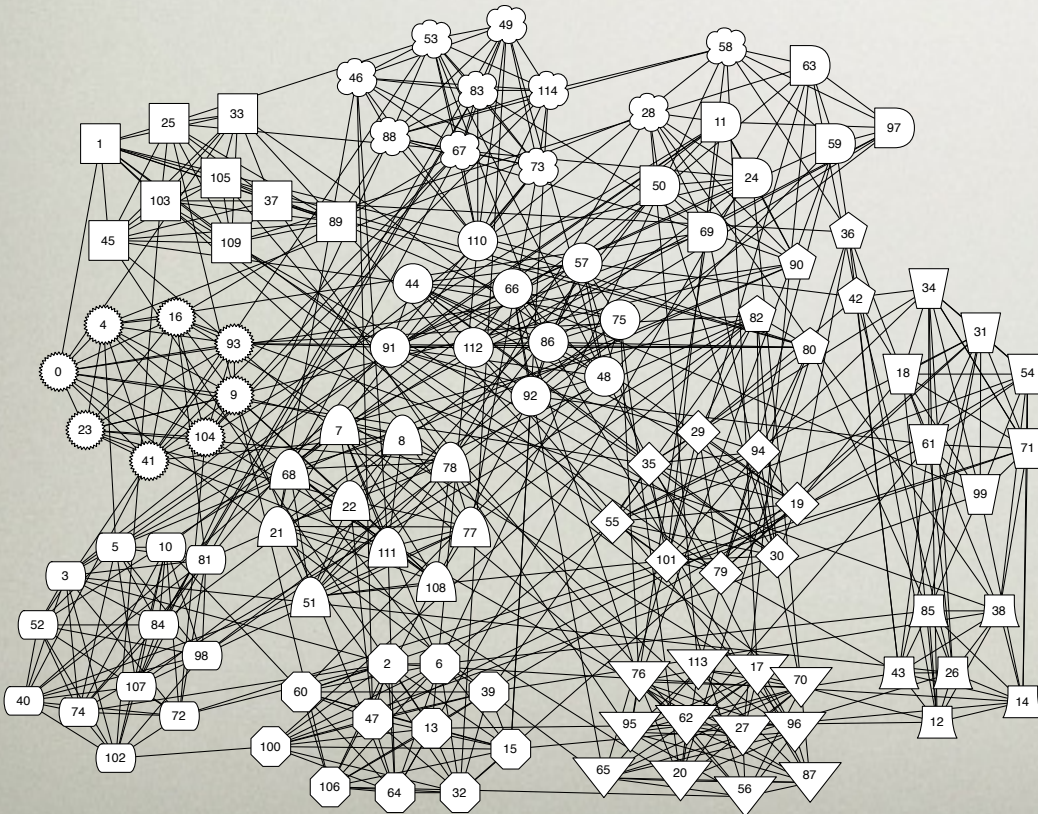


# TWO CASE STUDIES

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NCAA Schedule 2000

$n = 115$     $m = 613$



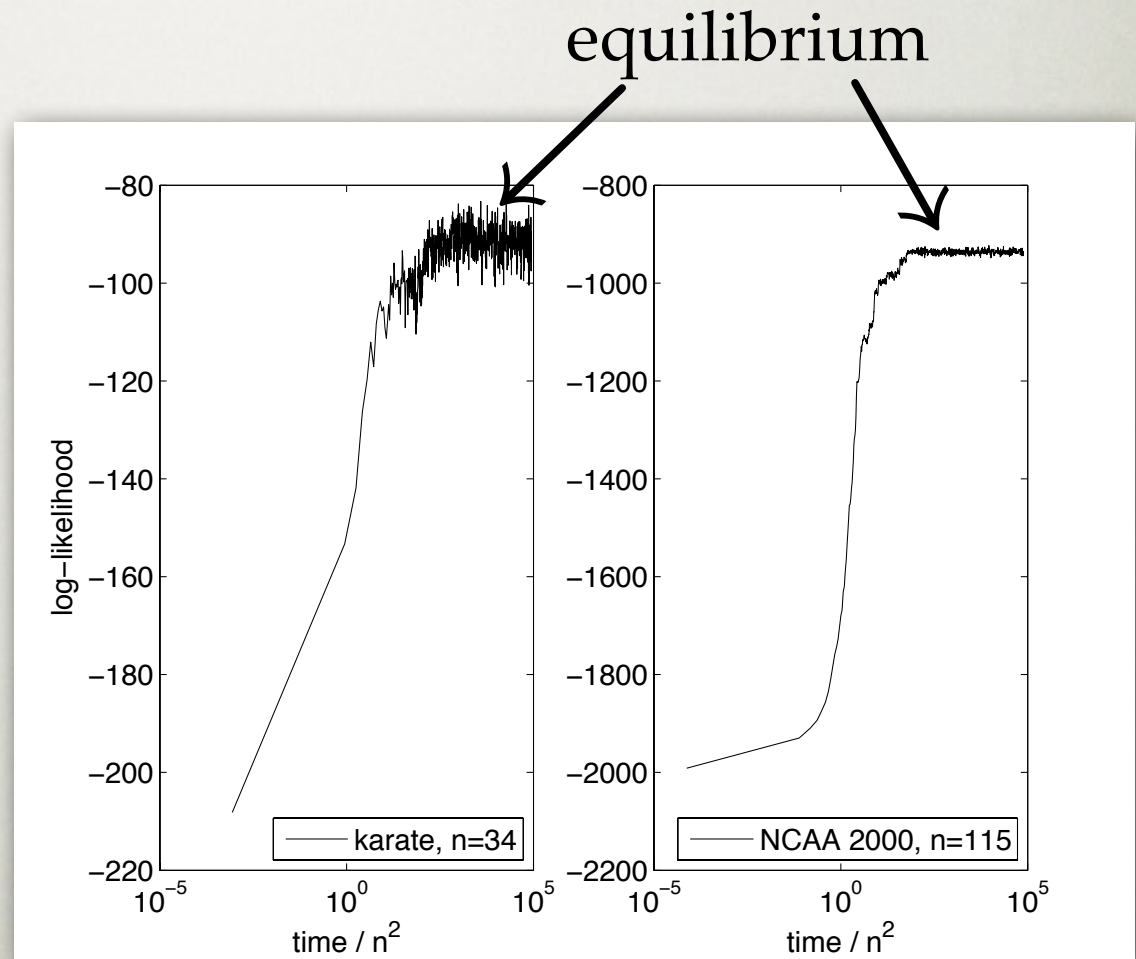
Zachary's Karate Club

$n = 34$     $m = 78$

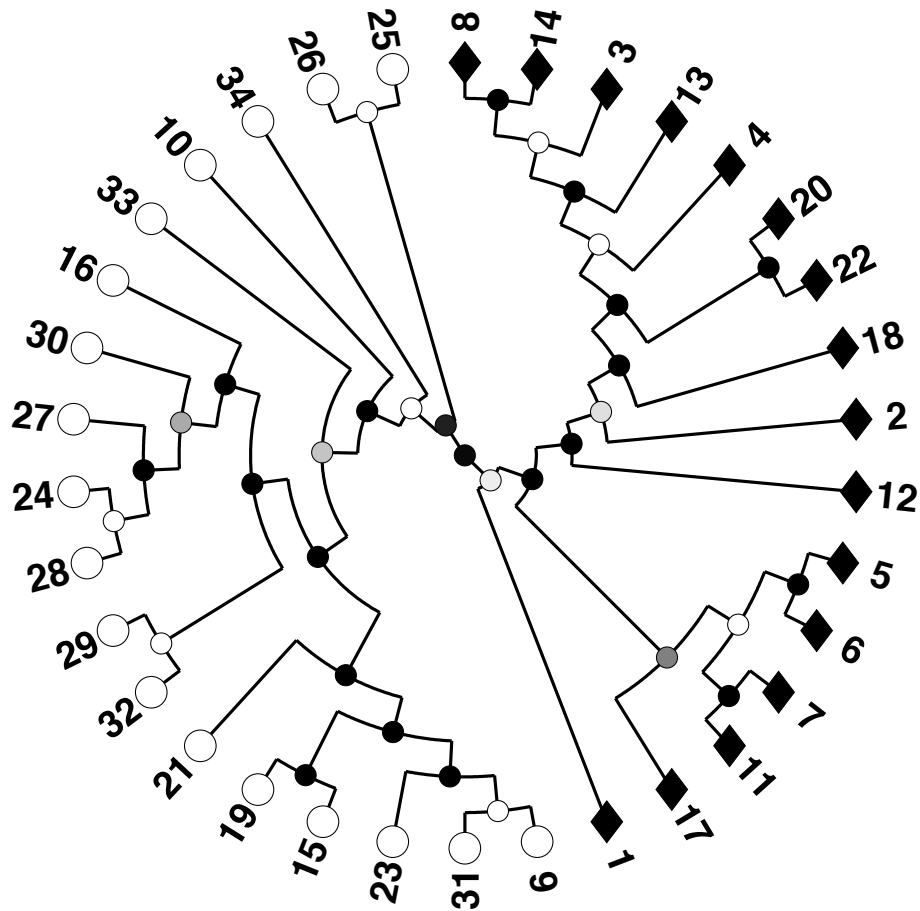
# MIXING TIMES

MCMC mixes  
relatively quickly

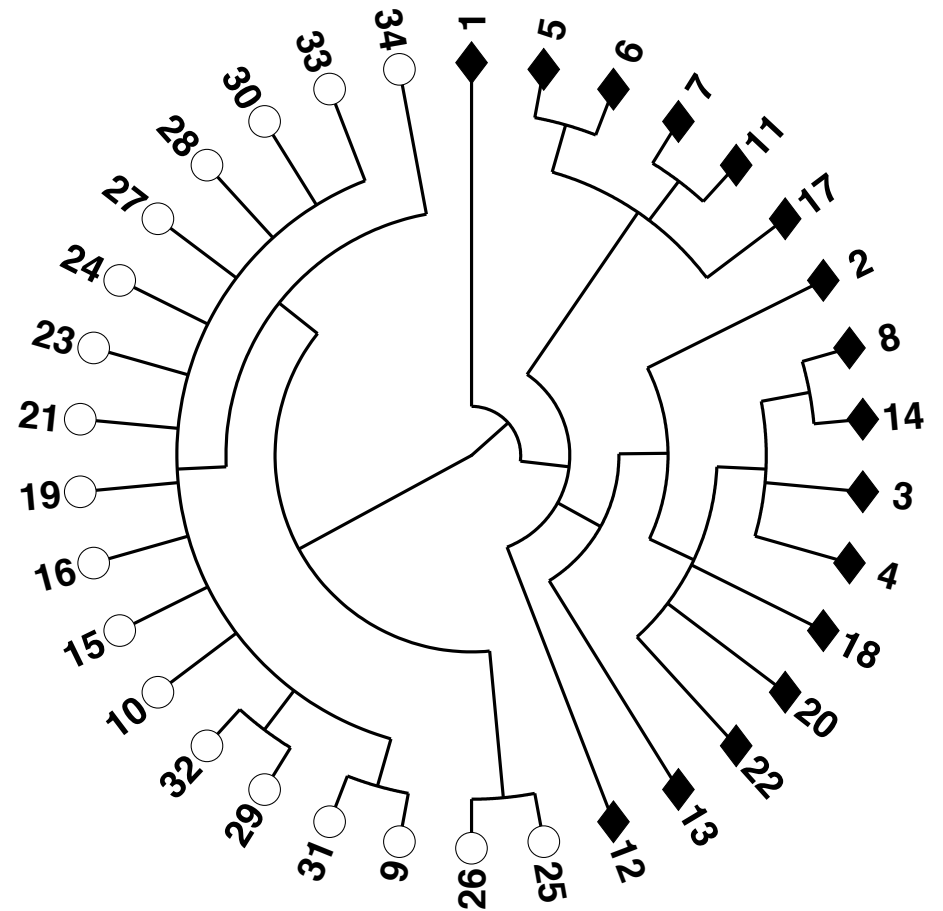
Equilibrium in  
 $\sim O(n^2)$  steps



# HIERARCHIES



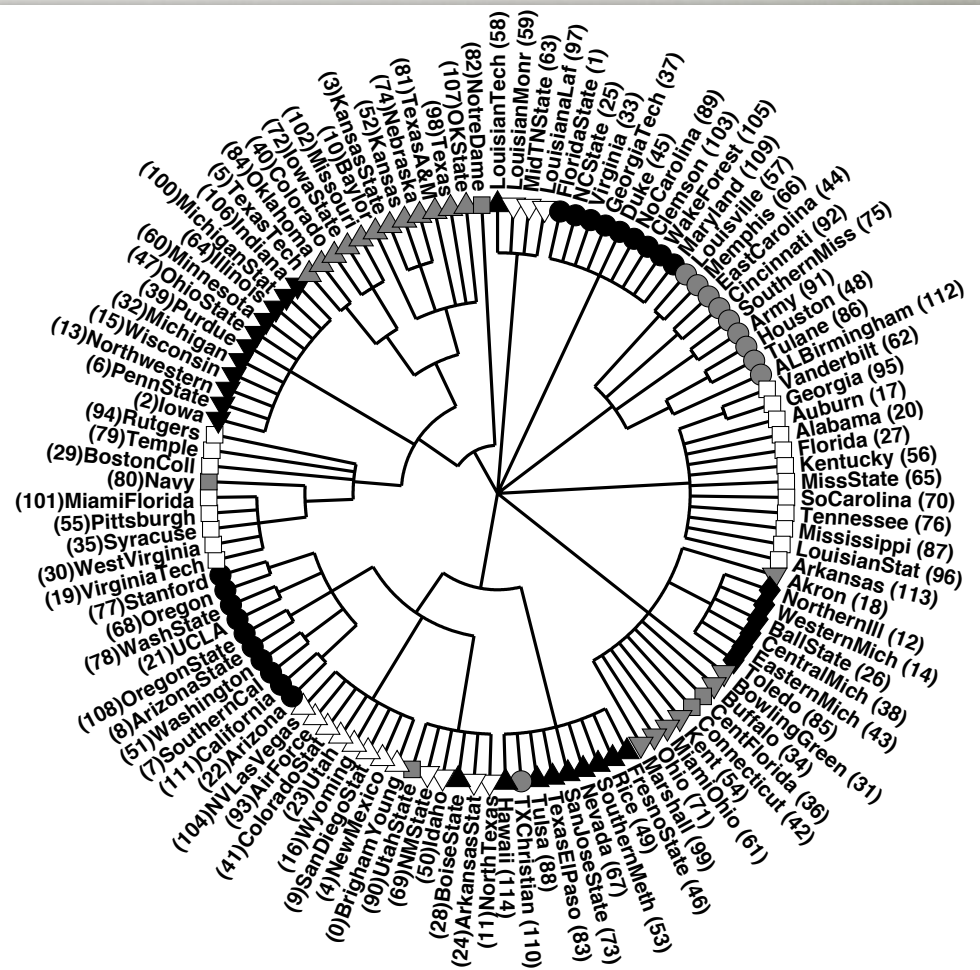
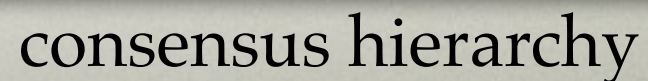
point estimate



consensus hierarchy



# point estimate



# EDGE ANNOTATIONS

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## Average likelihood of edge existing

- For each edge  $(i, j)$  in  $G$ , compute average associated parameter  $\langle \theta_r \rangle_{(i,j)}$  over sampled models
- $\langle \theta_r \rangle_{(i,j)}$  is edge annotation (weight)



# VERTEX ANNOTATIONS

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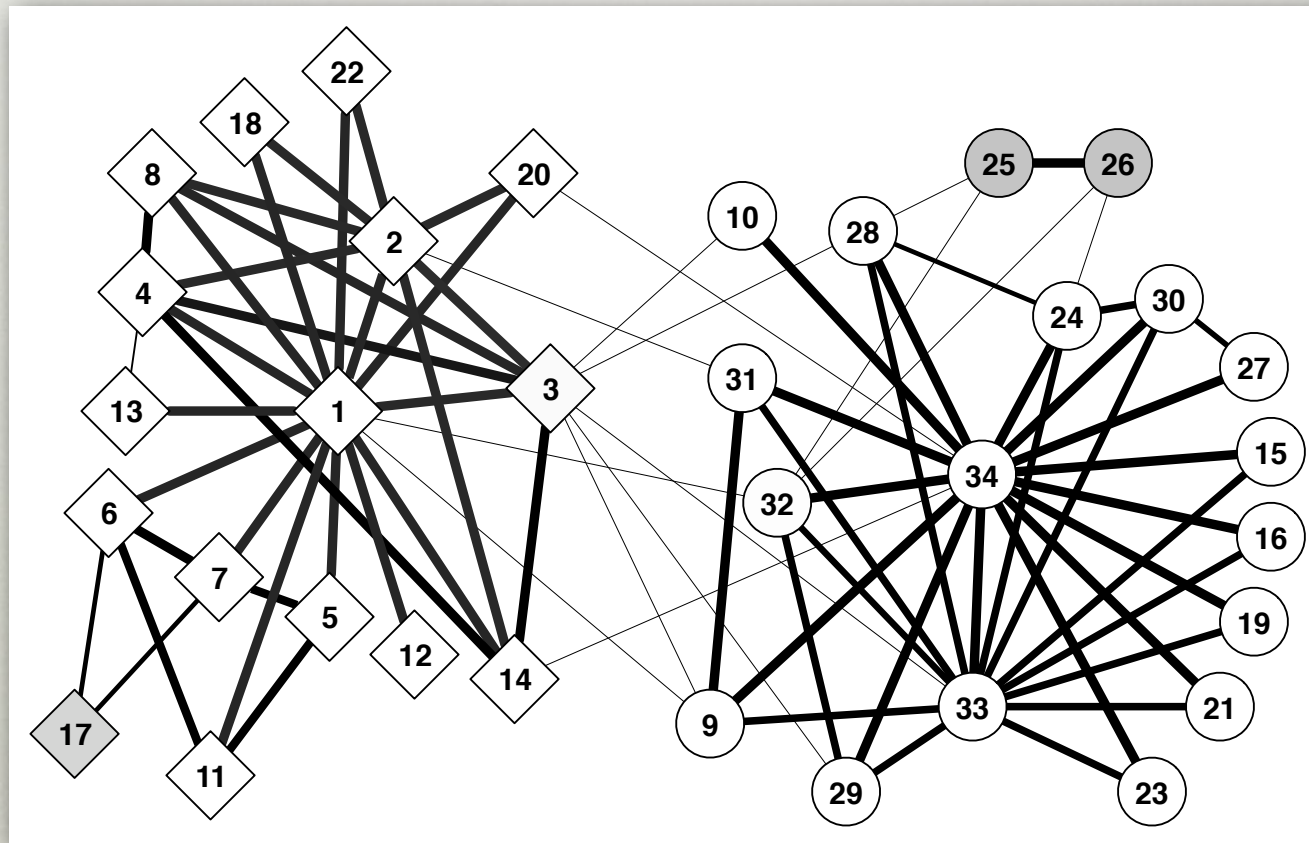
## Group-affiliation strengths

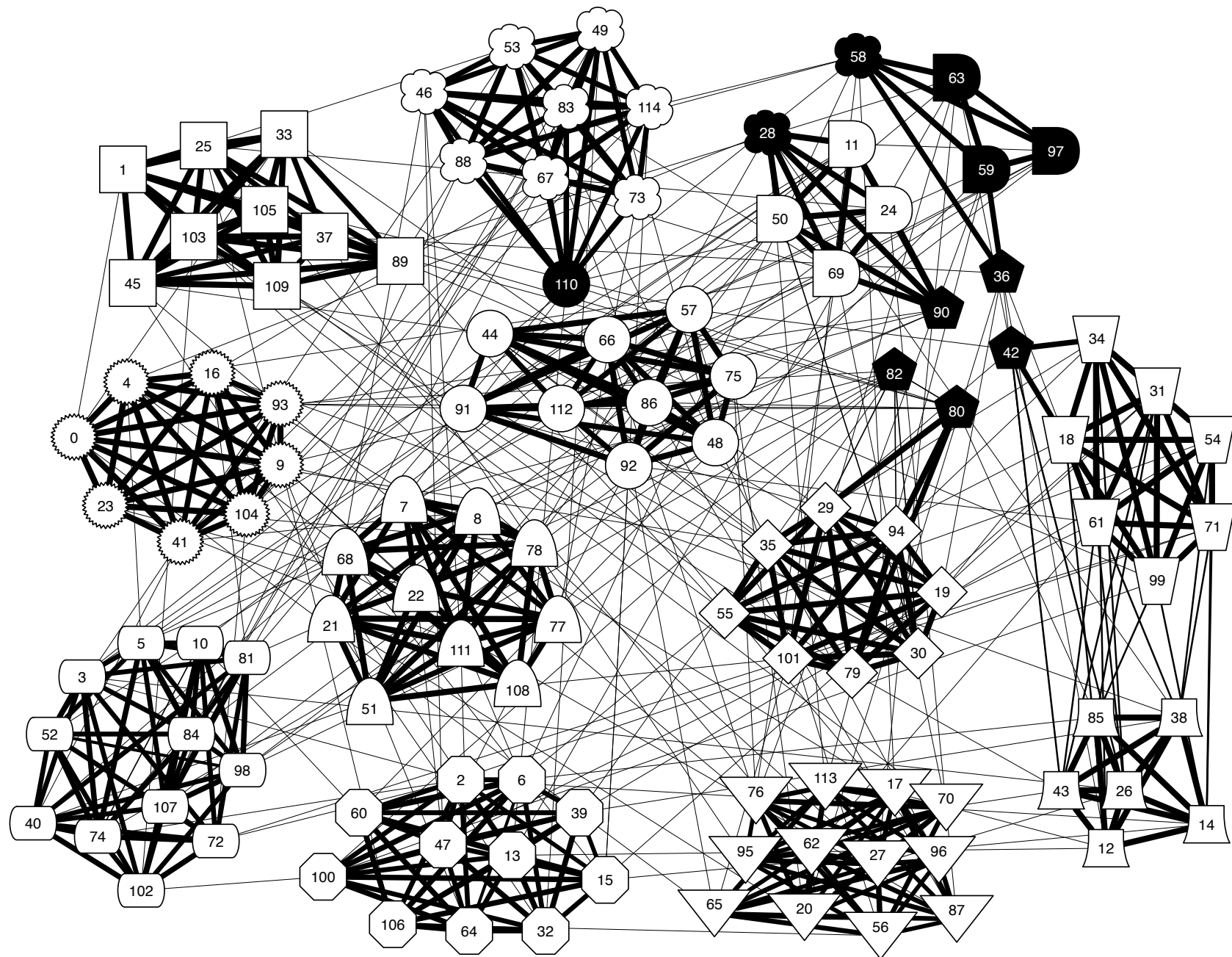
- If each vertex has known group label
- Ask, how often does vertex  $i$  appear in a subtree with majority of its fellows?
- Frequency is vertex annotation (strength)



# EDGE, NOTE ANNOTATIONS

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# MODEL CHECKING

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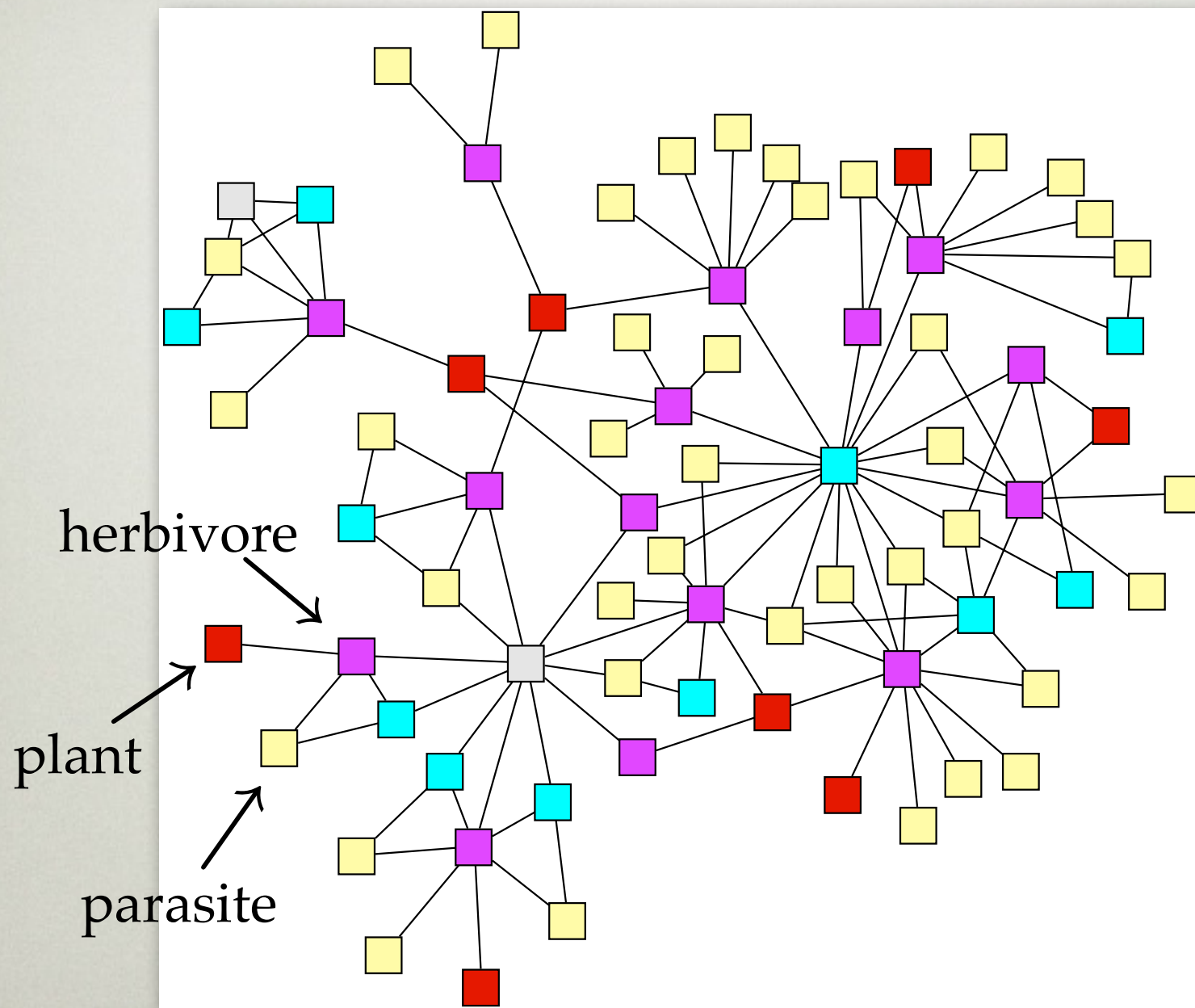


# MODEL CHECKING

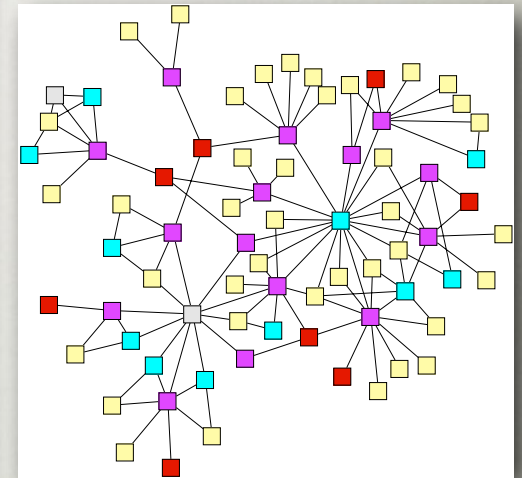
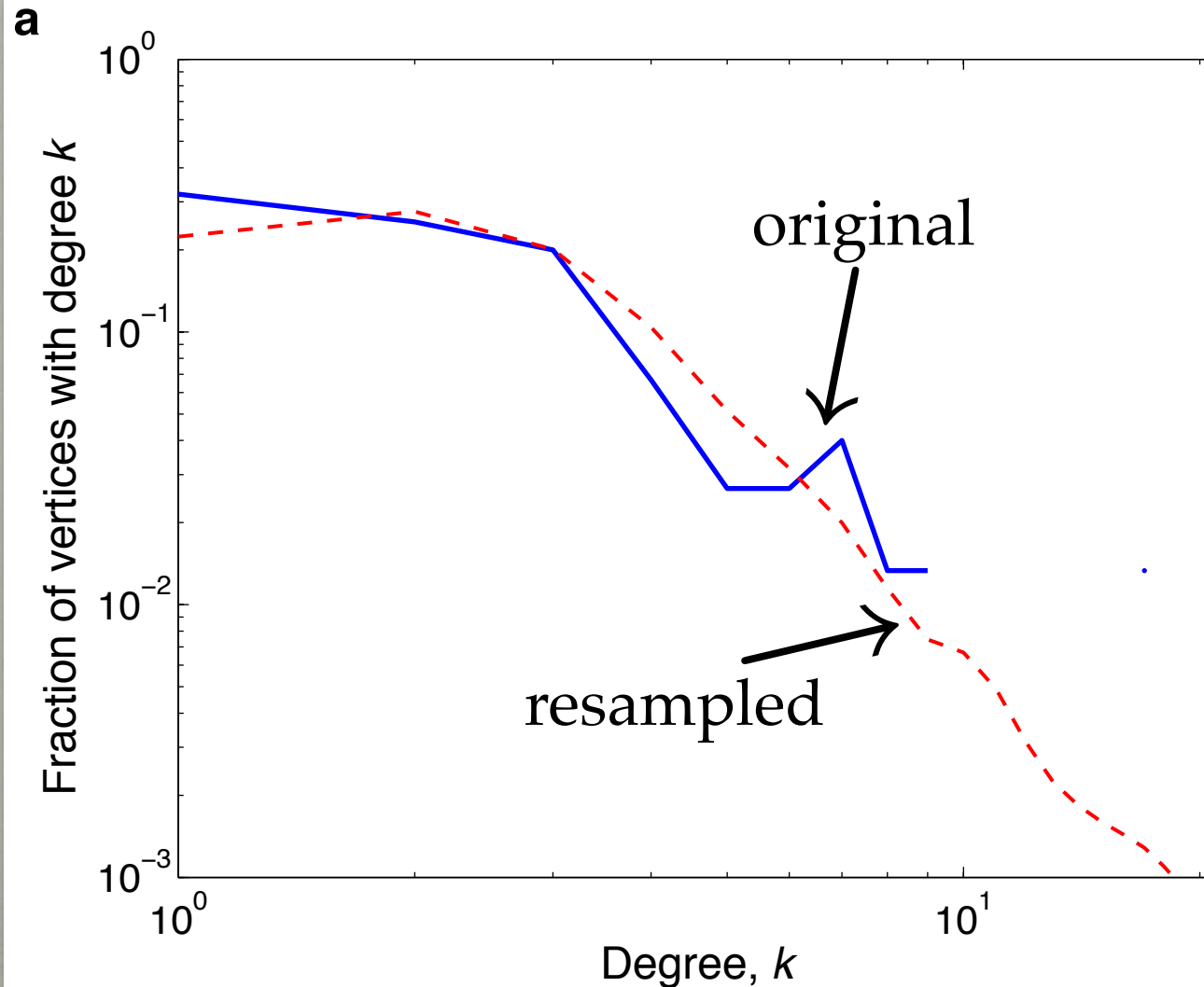
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- Given graph  $G$
- run MCMC to equilibrium
- then, for each sampled  $\mathcal{D}$ , draw a **resampled** graph  $G'$  from ensemble

**Checking the model (goodness-of-fit):  
do resampled graphs look like original?**

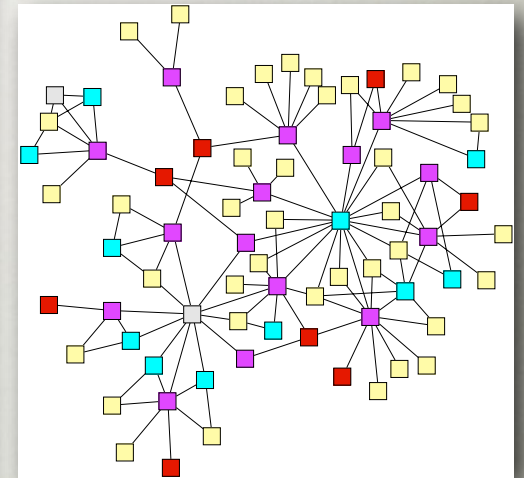
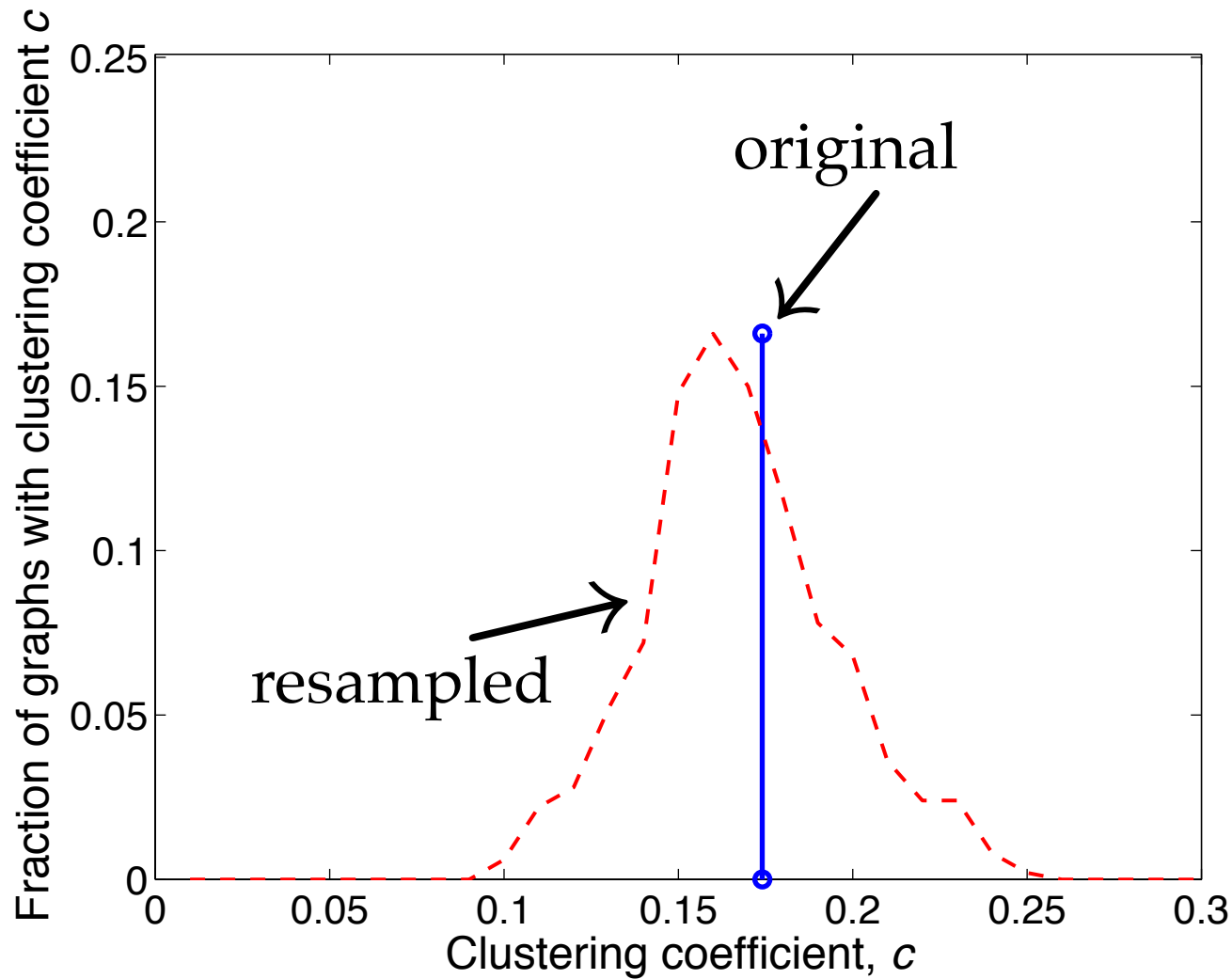


# DEGREE DISTRIBUTION

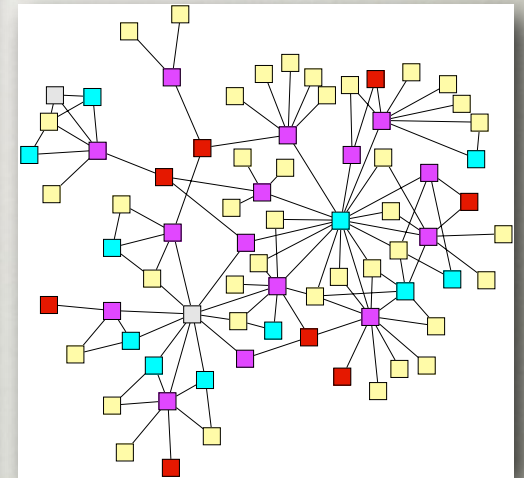
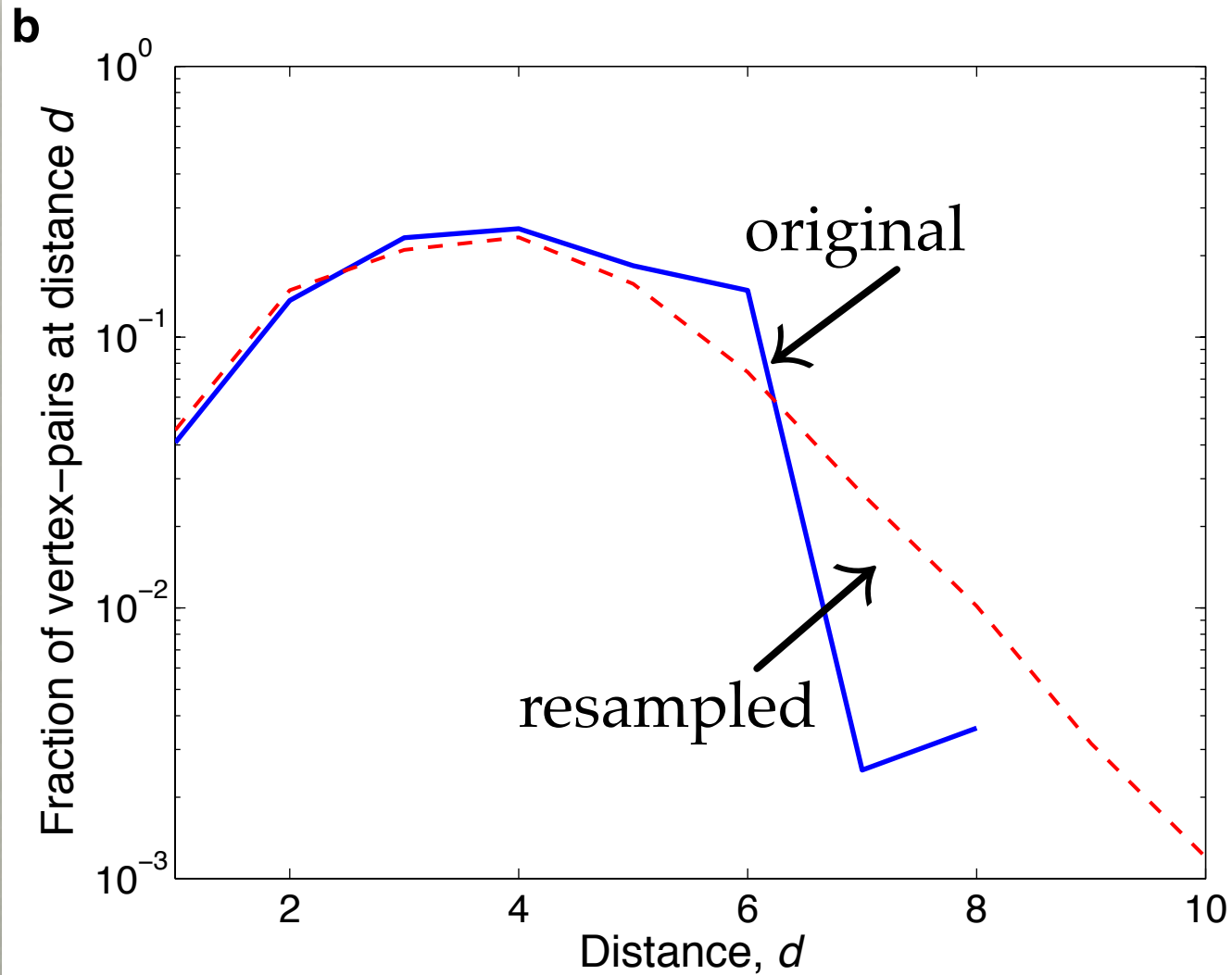




# CLUSTERING COEFFICIENT



# DISTANCE DISTRIBUTION



# PREDICTING MISSING LINKS

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many networks partially known, noisy

- social nets, food webs, protein interactions, etc.

can hierarchies predict their **missing links**?

previous approaches

- Liben-Nowell & Kleinberg (2003)
- Goldberg & Roth (2003)
- Szilágyi et al. (2005)
- many more now



# ACCURACY IS HARD

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- remove  $k$  edges from  $G$
- how easy to guess a missing link?

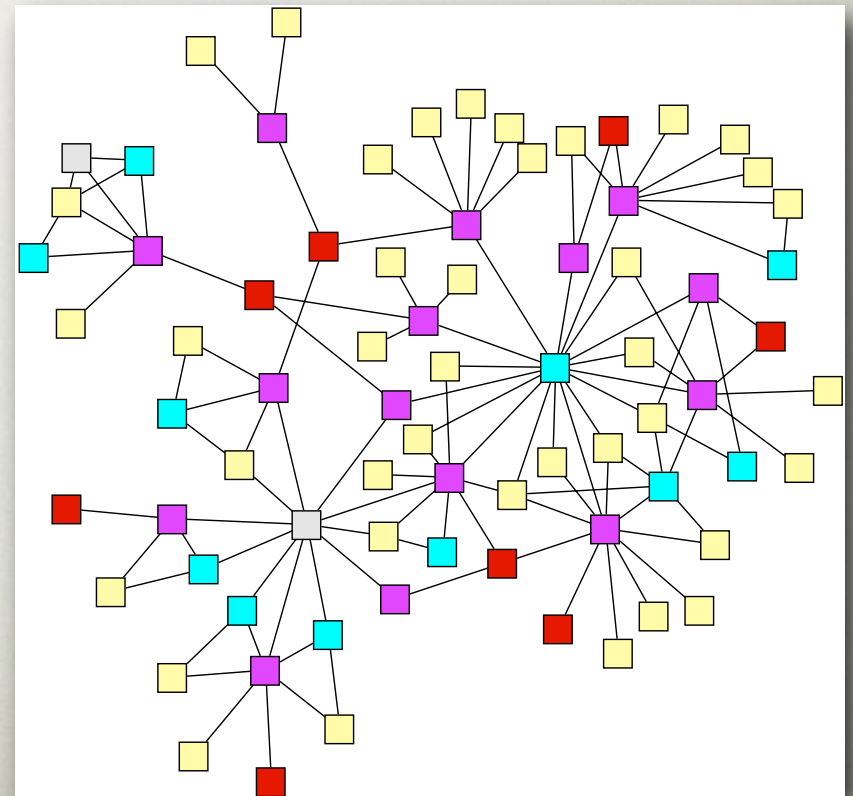
$$p_{\text{guess}} \approx \frac{k}{n^2 - m + k}$$
$$= O(n^{-2})$$

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$$n = 75$$

$$m = 113$$

$$p_{\text{guess}} = k / (2662 + k)$$



# GENERATIVE MODEL APPROACH

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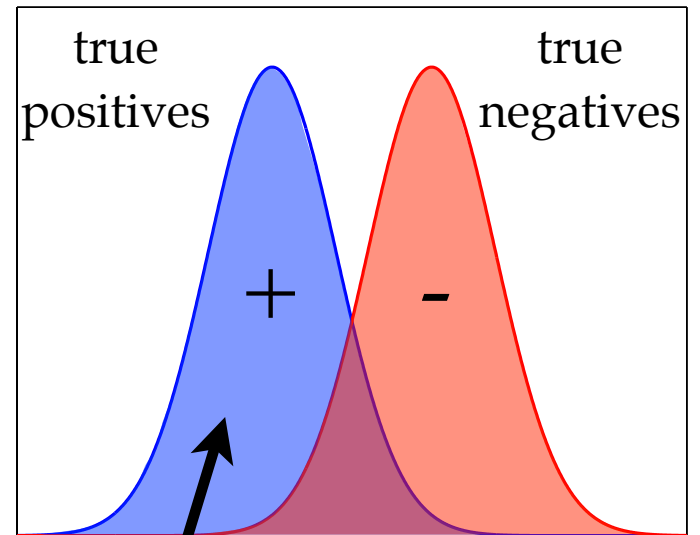
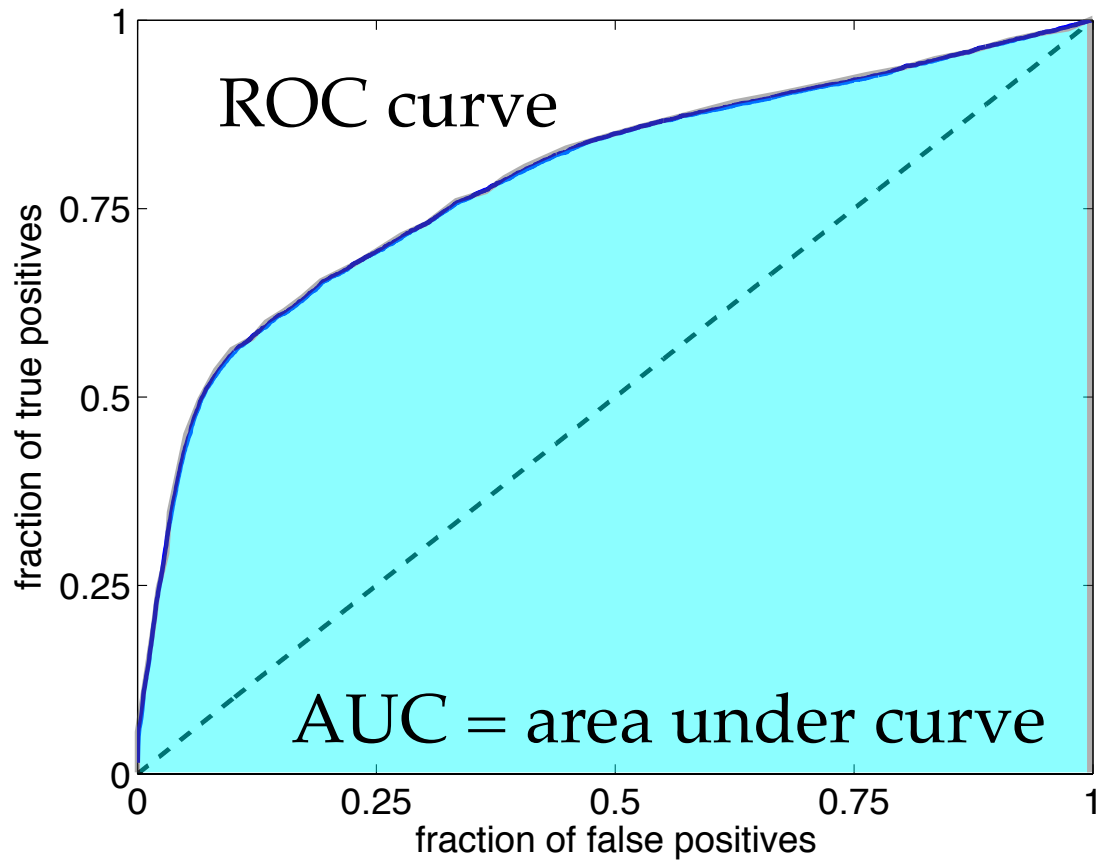
- Given incomplete graph  $G$
- run MCMC to equilibrium
- then, over sampled  $\mathcal{D}$ , compute average  $\langle p_r \rangle$  for links  $(i, j) \notin G$
- predict links with high  $\langle p_r \rangle$  values are missing

**Test via leave- $k$ -out cross-validation**

perfect accuracy:  $\text{AUC} = 1$

no better than chance:  $\text{AUC} = 1/2$

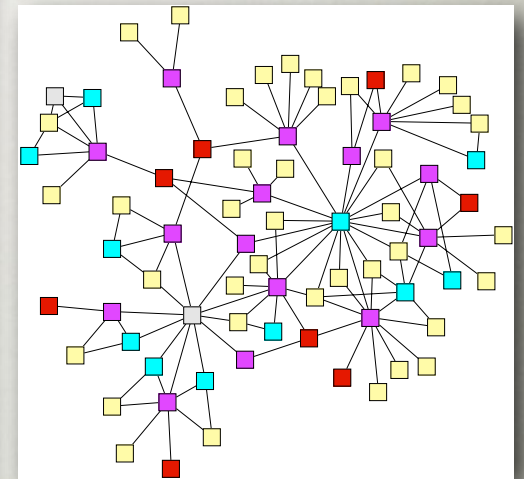
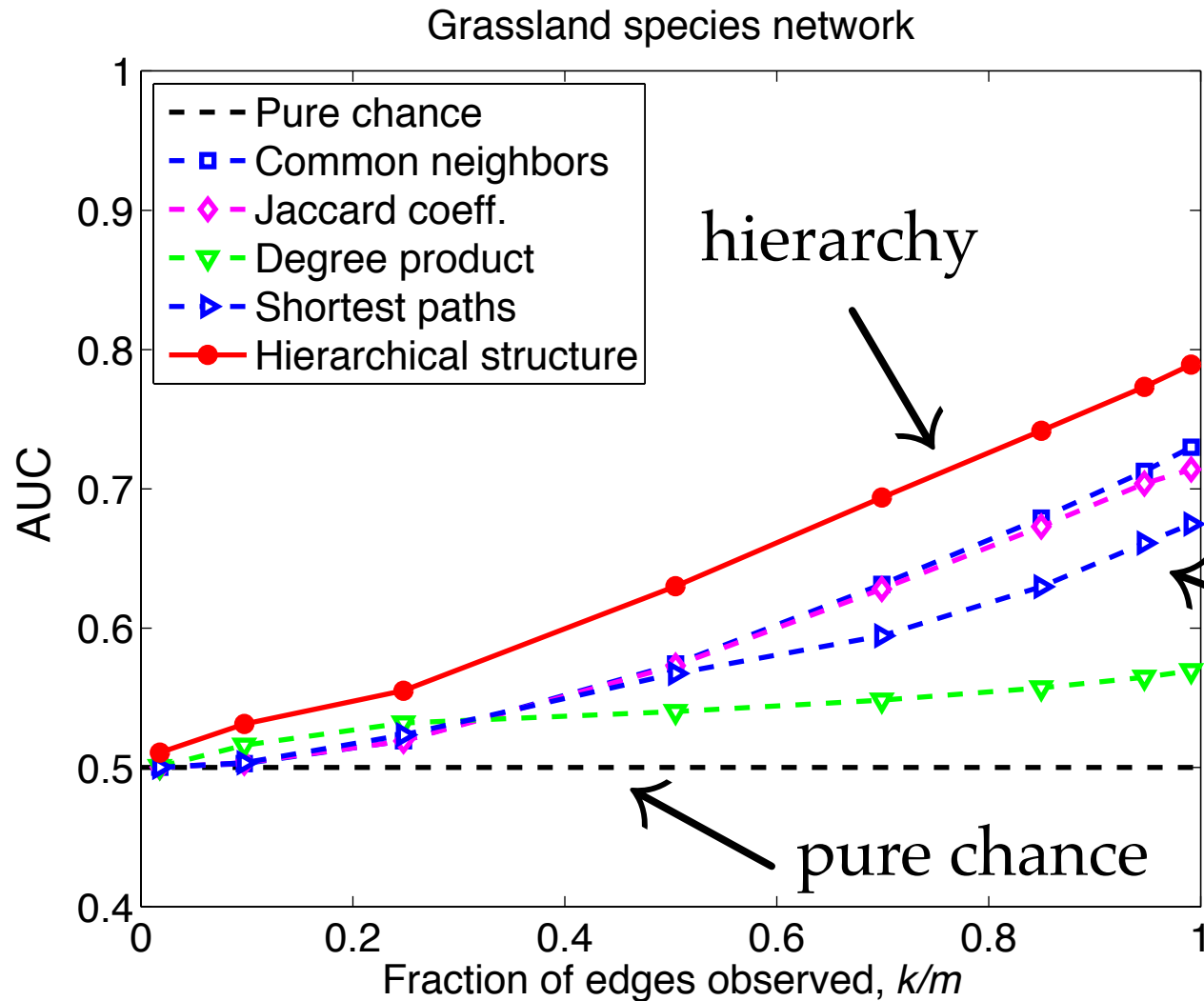
# SCORING THE PREDICTIONS



AUC =  
Pr( distinguish  
+ from - )

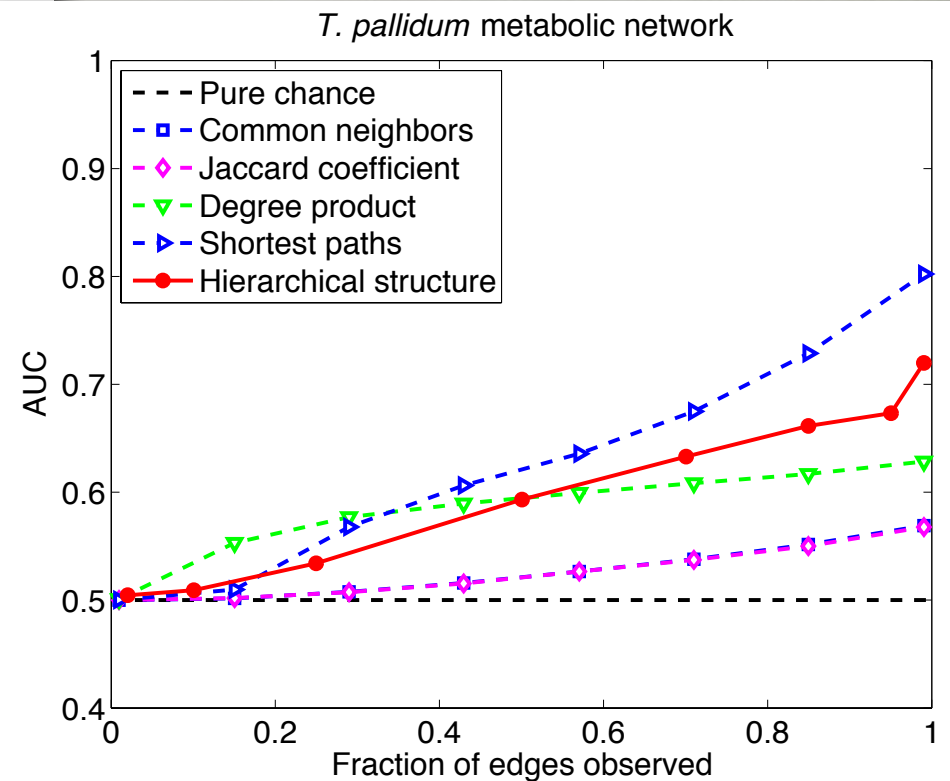
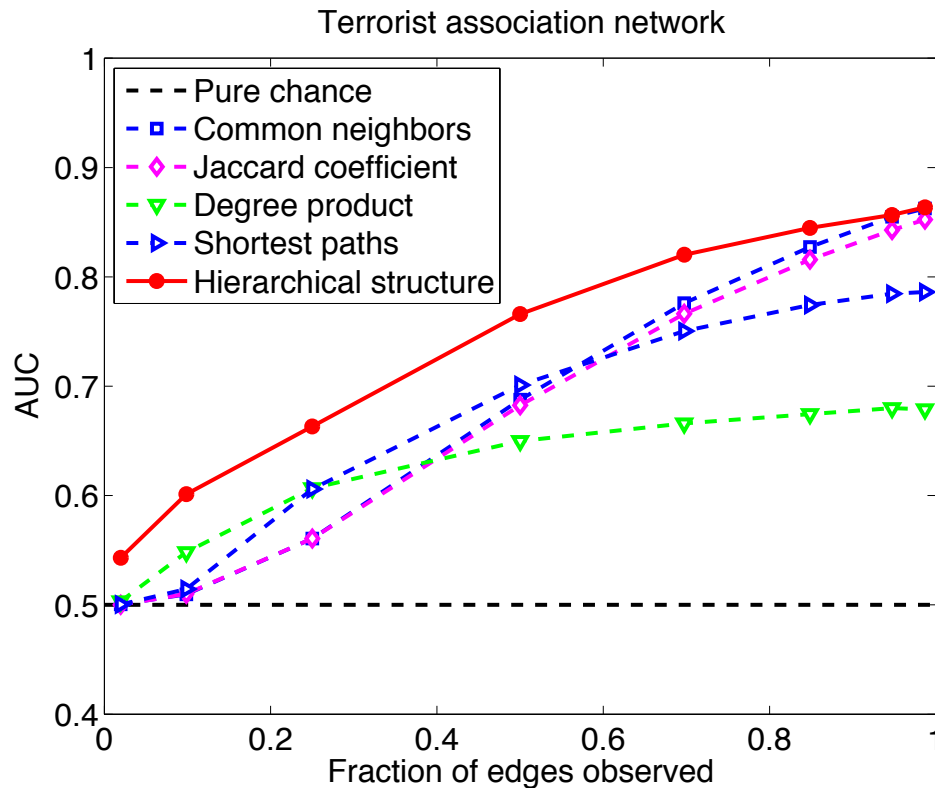


# PERFORMANCE 1



simple predictors

# PERFORMANCE 2



# SOME FINAL THOUGHTS

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- what processes create these hierarchical structures?
- scaling up the running time from  $O(n^2)$  ?
- active learning
- generalization to weighted, directed edges
- generalization to non-Poisson distributions