

Structure and dynamics of complex networks

March 26, 2020

Network models

B-A model

Model definition

Simulations

Analytic $p(k)$

C in the B-A model

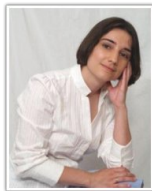
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THE BARABÁSI-ALBERT MODEL

Random graphs vs real networks

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	E-R	W-S
- sparseness, $\langle k \rangle$?	OK	OK
-the small-world effect, $\langle \ell \rangle$?	OK	OK
-large local clustering coeff., $\langle C \rangle$?	NO!	OK
-scale-free $p(k)$?	NO!	NO!

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-scale-free $p(k)$?	NO!	NO!

How to generate a **scale-free** random graph in a simple way?

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- Erdős-Rényi model: fixed N , (static)
- Watts-Strogatz model: fixed N , (static)
- Real networks?

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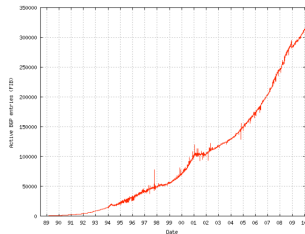
Actor network



Number of movies in IMDb

Herr II, Bruce W., Ke, Weimao, Hardy, Elisha, and Bömer, Katy. (2007) Movies and Actors: Mapping the Internet Movie Database. In Conference Proceedings of 11th Annual Information Visualization International Conference (IV 2007), Zurich, Switzerland, July 4-6, pp. 465-469.

Internet



Growth of the Internet routing table

<http://www.trainingsignaltraining.com/ccna-ipv6>

(from the slides of A.-L. Barabási)

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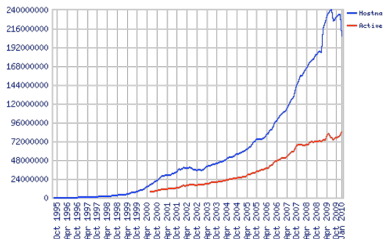
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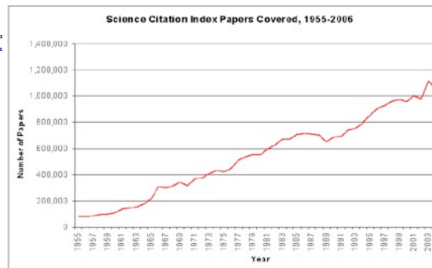
Randomisation

WWW



<http://website101.com/define-ecommerce-web-terms-definitions/>

Scientific Publications



http://www.kk.org/thetechnium/archives/2008/10/the_expansion_o.php

(from the slides of A.-L. Barabási)

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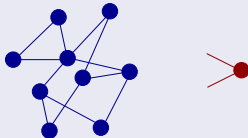
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The Barabási-Albert model

- Growing model: **one new node** in one step with **m new links**.



- The other end of the new links is attached according to the **preferential attachment rule**, i.e., we choose from the 'old' nodes with probabilities proportional to their degree:

$$\mathcal{P}(\text{choosing node } i) \equiv \mathcal{P}_i \sim k_i$$

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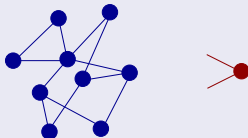
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- Please take a look at the 2 accompanying videos showing a growing B-A network and an E-R graph, where we also add links one by one (in random order).

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- Please take a look at the 2 accompanying videos showing a growing B-A network and an E-R graph, where we also add links one by one (in random order).
- Important differences are apparent already at such small system sizes: we can observe nodes that are likely candidates for becoming HUBS in the long run in the B-A video.

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- Next, take a look at the empirical degree distribution of B-A networks generated using Networkx in the accompanying jupyter notebook.

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- Next, take a look at the empirical degree distribution of B-A networks generated using Networkx in the accompanying jupyter notebook.
- It looks **scale-free!**

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- The scale-free $p(k)$ in the simulation results looks really satisfying.

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- The scale-free $p(k)$ in the simulation results looks really satisfying.
- But can we also derive in more analytic terms that this model leads to a scale-free network?

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- The scale-free $p(k)$ in the simulation results looks really satisfying.
 - But can we also derive in more analytic terms that this model leads to a scale-free network?
- Actually, it can be proven also on pure analytic grounds that the B-A model is scale-free.

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- The scale-free $p(k)$ in the simulation results looks really satisfying.
 - But can we also derive in more analytic terms that this model leads to a scale-free network?
- Actually, it can be proven also on pure analytic grounds that the B-A model is scale-free.
- However, here we are going to discuss a more simple (approximate) analytic derivation of the $p(k)$.

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Main stages of the forthcoming calculation:

- 1) Calculation of the $k_i(t)$, corresponding to the degree of node i at time step t , given that node i appeared as the new node at a given time step t_i .
- 2) Based on the formula obtained for $k_i(t)$, the derivation of $p(k)$.

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OK, so the first step is to study the time evolution of the degree for a given node in this model. Before diving into the formulas, a simple observation:

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OK, so the first step is to study the time evolution of the degree for a given node in this model. Before diving into the formulas, a simple observation:

Due to the model definition, the node degree for any node i in any time step

- can increase if the newcomer node is actually choosing i to connect to,
- or it may remain the same (if the newcomer node is choosing someone else).

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Due to the model definition, the node degree for any node i in any time step

- can increase if the newcomer node is actually choosing i to connect to,
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Let us first take a look at the degree as a function of time in simulation results!

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Time dependent degree

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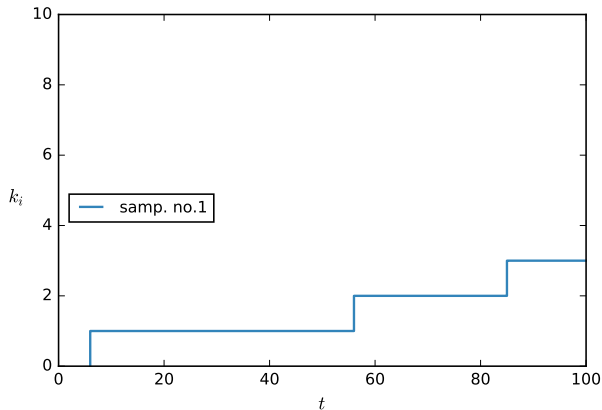
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The degree of a particular node as a function of t in simulations:



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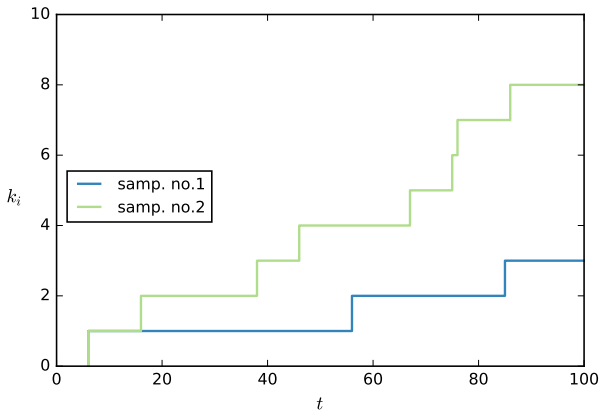
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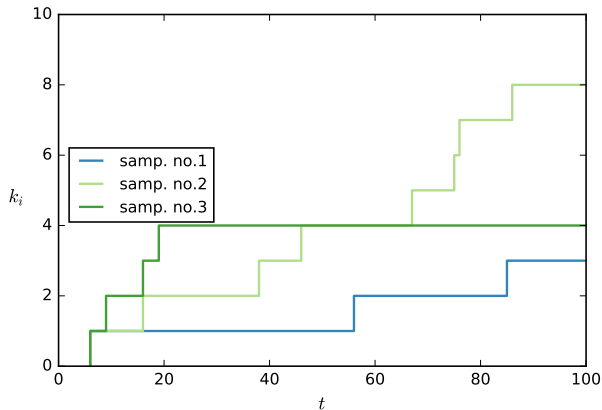
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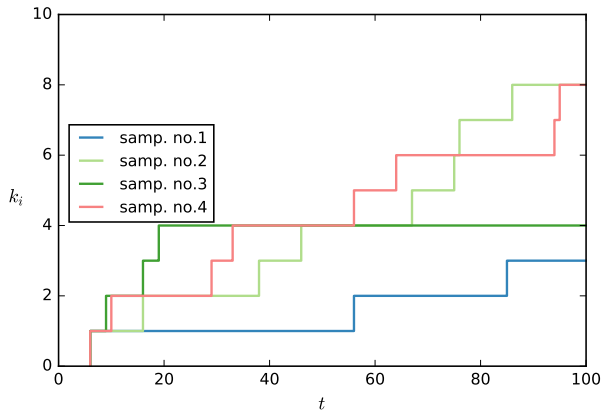
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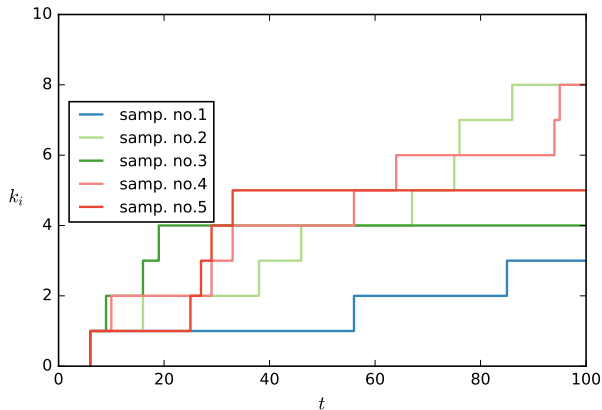
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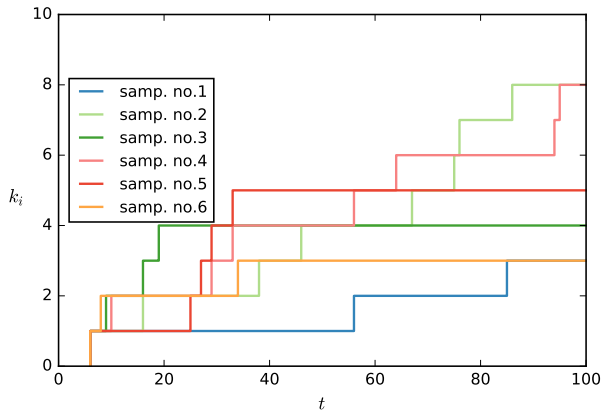
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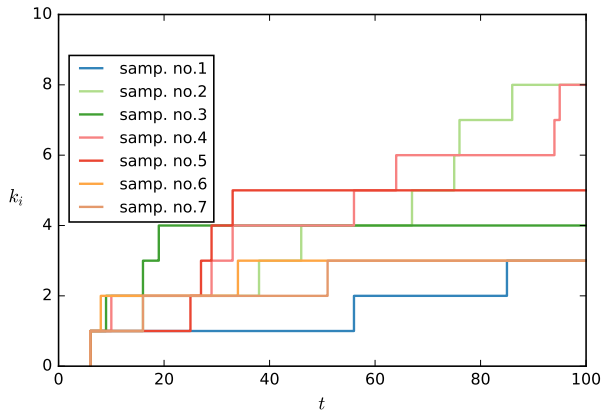
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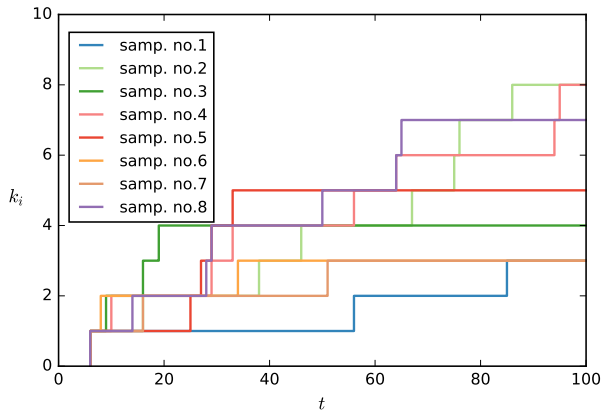
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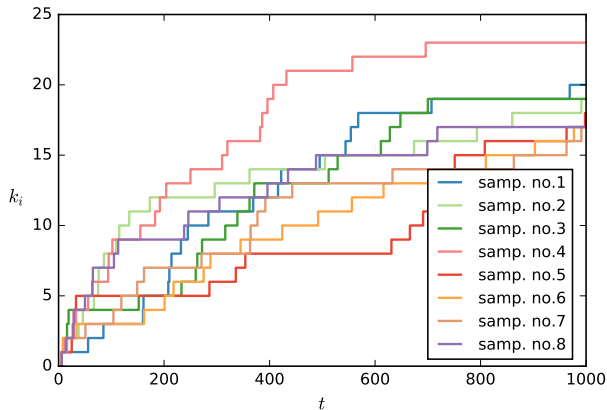
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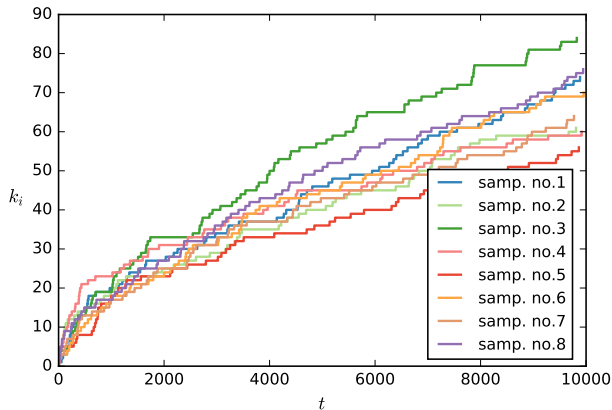
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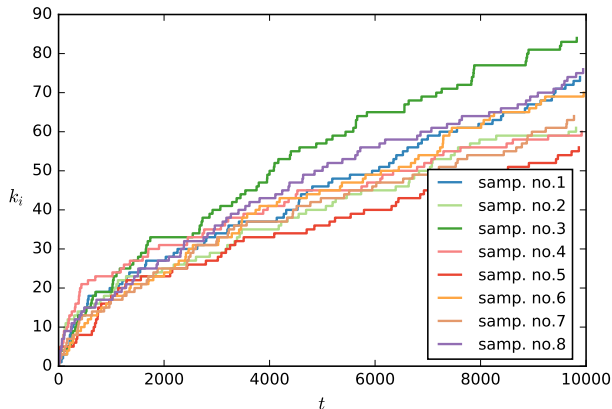
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The degree of a particular node as a function of t in simulations:



We can see a clear trend. Next, we are going to derive an analytic formula for the average $k_i(t)$!

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- For large t the number of nodes and links is

$$\begin{aligned} N &\simeq t, \\ M &\simeq mt. \end{aligned}$$

- The probability of choosing i :

$$\mathcal{P}_i = \frac{k_i}{\sum_j k_j}.$$

→ The approximate change of k_i in a time step:

$$\frac{\Delta k_i}{\Delta t} \simeq m \mathcal{P}_i \quad \rightarrow \quad \frac{\partial k_i}{\partial t} = m \frac{k_i}{\sum_j k_j}.$$

Since the sum of the degrees : $\sum_j k_j = 2M = 2mt$,

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- Solving the diff. eq.:

$$\begin{aligned}\frac{\partial k_i}{\partial t} &= \frac{k_i}{2t} \quad \rightarrow \quad \frac{\partial k_i}{k_i} = \frac{\partial t}{2t} \quad \rightarrow \quad \int \frac{dk_i}{k_i} = \int \frac{dt}{2t} \\ \rightarrow \quad \ln k_i &= \frac{1}{2} \ln t + \text{const.} \quad \rightarrow \quad k_i(t) = ct^{\frac{1}{2}}\end{aligned}$$

- How to determine the constant c ?

At $t = t_i$, (the appearance of node i): $k_i = m$, leading to

$$k_i(t) = m \left(\frac{t}{t_i} \right)^{\frac{1}{2}}.$$

- Thus, the degree of any node is increasing as $k(t) \sim t^\beta$, where $\beta = 1/2$ is the so-called **dynamical exponent**.

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At $t = t_i$, (the appearance of node i): $k_i = m$, leading to

$$k_i(t) = m \left(\frac{t}{t_i} \right)^{\frac{1}{2}}.$$

- Thus, the degree of any node is increasing as $k(t) \sim t^\beta$, where $\beta = 1/2$ is the so-called **dynamical exponent**.

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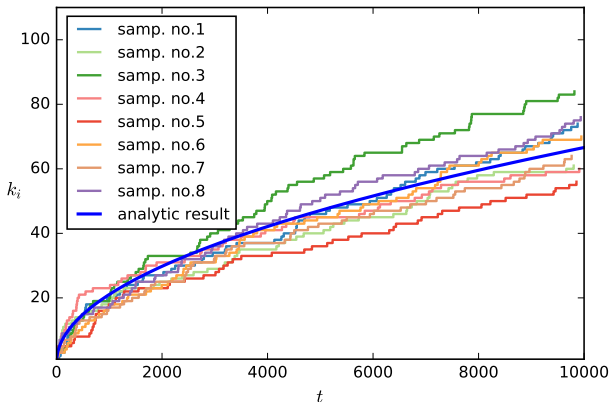
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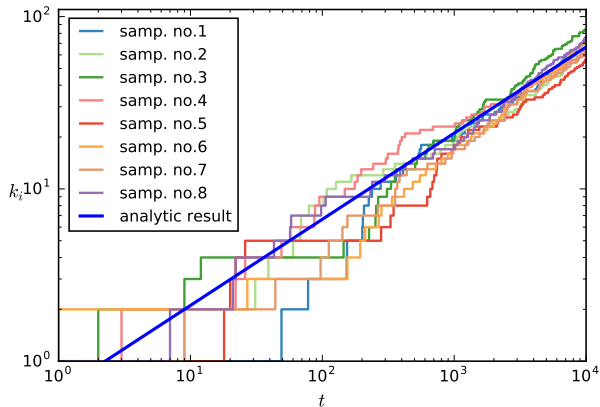
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- How to calculate the degree distribution from $k_i(t)$?

→ The easiest way is by considering the cumulative distribution first.

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- The cumulative degree distribution:

$$P(k) \equiv \mathcal{P}(k_i < k) = \mathcal{P}(m(t/t_i)^{1/2} < k) = \mathcal{P}(t/t_i < (k/m)^2) = \mathcal{P}(t_i/t > (m/k)^2).$$

- The lengths of the time steps:

$$\begin{aligned} P(k) &= 1 - \left(\frac{m}{k}\right)^2 \\ \rightarrow \mathbf{p(k)} &= 2m^2 \mathbf{k}^{-3} \end{aligned}$$

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- The lengths of the time steps:



$$P(k) = 1 - \left(\frac{m}{k}\right)^2$$
$$\rightarrow \mathbf{p(k)} = 2m^2 \mathbf{k}^{-3}$$

- The degree distribution is **SCALE-FREE** with $\gamma = 3$!

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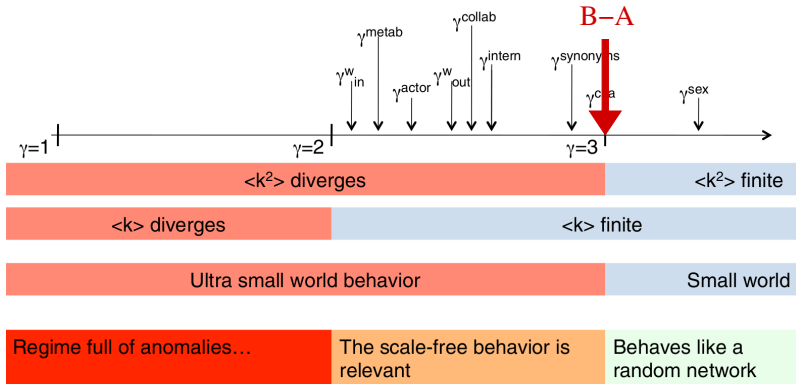
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(from the slides of A.-L. Barabási)

Average distance in scale-free networks

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$$\langle l \rangle \sim \left\{ \begin{array}{ll} \text{const.} & \gamma \leq 2 \\ \frac{\ln \ln N}{\ln(\gamma-1)} & 2 < \gamma < 3 \\ \frac{\ln N}{\ln \ln N} & \gamma = 3 \\ \ln N & \gamma > 3 \end{array} \right\} \begin{array}{l} \text{Ultra Small World} \\ \\ \\ \text{Small World} \end{array}$$

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- What is the **clustering coefficient** in the B-A model?

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- What is the **clustering coefficient** in the B-A model?
- We can calculate it in a rather simple way. As a first step, let us express the probability for nodes i and j , (introduced at time steps t_i and t_j) to be connected!

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- The only moment when they can connect is when the later coming node is actually the new node. Let's assume that i comes first and j comes later.

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- The degree of node i is given by $k_i(t) = m(\frac{t}{t_i})^{1/2}$

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- The degree of node i is given by $k_i(t) = m(\frac{t}{t_i})^{1/2}$
- The probability of a connection to j introduced at $t = t_j$:

$$\mathcal{P}(i-j) = m \frac{k_i}{2mt} = \frac{k_i}{2t} = \frac{m \left(\frac{t_j}{t_i}\right)^{\frac{1}{2}}}{2t_j} = \frac{m}{2} (t_i t_j)^{-\frac{1}{2}}.$$

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- As the next step, let us write down the expected number of links between the neighbours of a node l , (introduced at t_l), at the end of the node generation process $t = N$:

$$\begin{aligned}n_l &= \frac{1}{2} \sum_{t_i=1}^N \sum_{t_j=1}^N \mathcal{P}(l-i) \mathcal{P}(l-j) \mathcal{P}(i-j) \\&\simeq \frac{1}{2} \int_1^N dt_i \int_1^N dt_j \mathcal{P}(l-i) \mathcal{P}(l-j) \mathcal{P}(i-j) \\&= \frac{m^3}{16} \int_1^N dt_i \int_1^N dt_j (t_l t_i)^{-\frac{1}{2}} (t_l t_j)^{-\frac{1}{2}} (t_i t_j)^{-\frac{1}{2}} \\&= \frac{m^3}{16 t_l} \int_1^N dt_i \frac{1}{t_i} \int_1^N dt_j \frac{1}{t_j} = \frac{m^3}{16 t_l} (\ln N)^2.\end{aligned}$$

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- The number of “edge places” between the neighbors of l at the end ($t = N$):

$$\frac{k_l(k_l - 1)}{2} \simeq \frac{k_l^2}{2} = \frac{1}{2} \left[m \left(\frac{t}{t_l} \right)^{\frac{1}{2}} \right]^2 = \frac{m^2 N}{2 t_l}$$

- Thus, the clustering coefficient of l :

$$C_l = \frac{m^3}{16 t_l} (\ln N)^2 \frac{2 t_l}{m^2 N} = \frac{m (\ln N)^2}{8 N}.$$

- Since we choose a general l , and C_l does not depend on l , our result for C in the B-A model:

$$C = \frac{m (\ln N)^2}{8 N}.$$

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$$\langle C \rangle = \frac{m(\ln N)^2}{8N} \rightarrow \text{Is this good or bad?}$$

→ Let's compare it to the Erdős-Rényi model:

- Erdős-Rényi: $\langle C \rangle \simeq \frac{\langle k \rangle}{N} \sim \frac{1}{N}$

- Barabási-Albert: $\langle C \rangle = \frac{m(\ln N)^2}{8N} \sim \frac{(\ln N)^2}{N}$

Slower decay with N , however, no decay at all in real networks:

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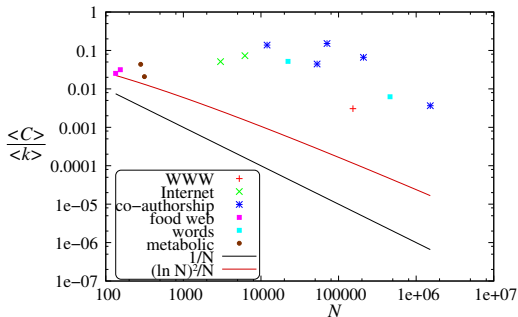
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Slower decay with N , however, no decay at all in real networks:



The Barabási-Albert model

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- **Growth** and **preferential attachment**,
- Dynamical exponent: $k(t) \sim t^\beta$, $\beta = \frac{1}{2}$,
- **SCALE-FREE**: $p(k) \sim k^{-3}$, $\gamma = 3$,
- Small-world: $\langle l \rangle \sim \frac{\ln N}{\ln \ln N}$,
- $C \sim \frac{(\ln N)^2}{N}$.

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Remaining challenges, problems?

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Remaining challenges, problems:

- Original B-A: $\gamma = 3$.



In real networks $2 \leq \gamma \leq 3$.

- Original B-A: oldest nodes have highest degree.



Not always in real systems, e.g., Google.

- Low clustering coefficient.

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Additional fitness

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- A simple solution to the problem of **tunable γ exponent** is given by one of the practice exercise, involving a **uniform additional fitness** beside the degree.

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To enable **newcomers overtake old nodes** in the B-A model we need a further **multiplicative fitness**:

- We introduce a fitness parameter η for the nodes, drawn from some distribution $\rho(\eta)$.
- The probability for node i to acquire new links:

$$\mathcal{P}_i \sim \eta_i k_i \quad \rightarrow \quad \mathcal{P}_i = \frac{\eta_i k_i}{\sum_j \eta_j k_j}$$

→ The time evolution of the degree now depends also on the fitness:

$$k_i(t) \sim t^{\beta(\eta_i)}, \quad \beta(\eta) = c\eta$$

- This also means that now a later coming node with higher fitness (and higher β exponent) can actually overtake older but less fit nodes.

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G. Bianconi and A.-L. Barabási, *Phys. Rev. Lett.* **86**, 5632 (2001)

B-A model with high clustering

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- P. Holme and B. J. Kim: **extra triad formation steps.**
 - Add a new node according to the P.A. rule.
 - After each step, with probability p also form m triangles with new node and randomly chosen neighbours of the nodes the new node was just attached.

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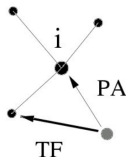
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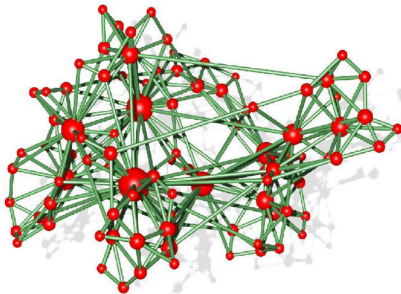
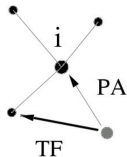
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HK graph with $p = 0.9$

Scale-free model with high clustering

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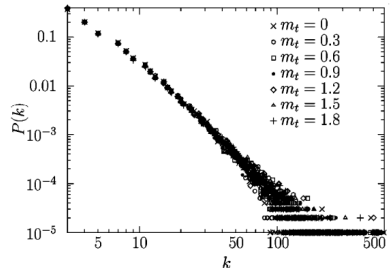
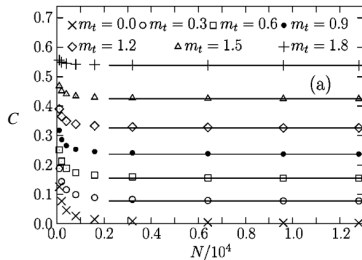
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- Holme-Kim model:



→ tunable clustering coefficient!

(The larger p we set, the larger is $\langle C \rangle$ in the resulting network).

P. Holme and B. J. Kim, *Phys. Rev. E* **65**, 026107 (2002)

Further reading on the Barabási-Albert model (not compulsory):

- Network science book by A.-L. Barabási, chapter 5:
<http://networksciencebook.com/chapter/5>
- Wikipedia: https://en.wikipedia.org/wiki/Barabasi-Albert_model

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THE CONFIGURATION MODEL

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- What if we do not care about the **mechanism** leading to scale-free $p(k)$ and are interested only in the **actual form** of the degree distribution?

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- What if we do not care about the **mechanism** leading to scale-free $p(k)$ and are interested only in the **actual form** of the degree distribution?
- How to generate a random network with arbitrary $p(k)$?

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The configuration model

- The goal is to generate a network with N nodes and a given $p(k)$ corresponding to a random sample from all possible realisations.
- We first draw N times from $p(k)$ to obtain the degree sequence.
- Then the “half links” are joined to connect the nodes with each other.

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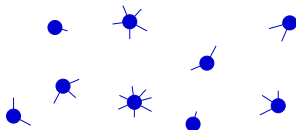
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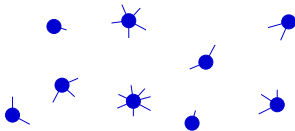
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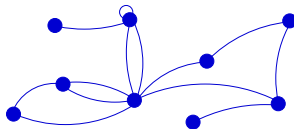
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- If you want a simple graph (without self-connections and multi-edges), connecting the nodes is the hard part...
- How would you start?

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Randomisation

- The usual tactic is to start the connection process at the large degree nodes.

The configuration model

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- The usual tactic is to start the connection process at the large degree nodes.
- However, this way the obtained graph cannot be regarded as completely random sample...

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- The usual tactics is to start the connection process at the large degree nodes.
- However, this way the obtained graph cannot be regarded as completely random sample...

→ Randomisation is needed!

Randomisation

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- If we keep only N and M fixed during the randomisation, we end up with an Erdős-Rényi graph in the end.

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- If we keep only N and M fixed during the randomisation, we end up with an Erdős-Rényi graph in the end.
- The aim is to get rid of any extra correlation coming from the node connection process, but also **preserve the degree of the nodes**.

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- If we keep only N and M fixed during the randomisation, we end up with an Erdős-Rényi graph in the end.
- The aim is to get rid of any extra correlation coming from the node connection process, but also **preserve the degree of the nodes**.
- How can we do that?

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Link randomisation

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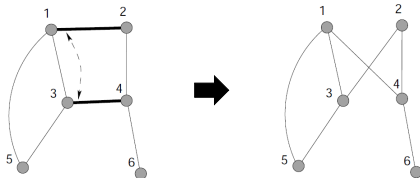
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An intuitive idea is link randomisation:

- in every step, choose 2 links at random,
- and swap one end of the links.
- (Of course, we have to check that we do not introduce self connections or multi edges before actually carrying out the swap).



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Node-randomization

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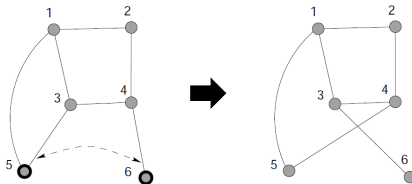
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Another possibility is node randomisation:

- in every step, choose 2 nodes at random,
- and swap one link on the nodes.
- (Again, we have to check that we do not introduce self connections or multi edges before actually carrying out the swap).



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How long do we have to keep on swapping the links?

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How long do we have to keep on swapping the links?

→ A general rule of the thumb is that the number of swaps should be high enough so that on average every link was swapped a few times, i.e., a the number of swaps should be something like 3 or 5 times the number of links.

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Now that we can randomise networks in a way that the node degrees remain the same, comparing any given real network to its configuration model counterpart has become very easy:

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Now that we can randomise networks in a way that the node degrees remain the same, comparing any given real network to its configuration model counterpart has become very easy:

We just **randomise a copy** of the original network (again, the number of carried out rewirings should be a few times the number of links), and the **resulting graph can be viewed as a sample from the configuration model**, where $p(k)$ **is exactly the same** as in the original network!

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Now that we can randomise networks in a way that the node degrees remain the same, comparing any given real network to its configuration model counterpart has become very easy:

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→ The configuration model is also a very important random 'baseline' to which we can compare a given real network.

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- Any other application of randomisation apart from the configuration model?

Measuring significance by randomization

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- Real networks are not completely random nor completely regular...
- To what extent is a given feature (e.g., $\langle \ell \rangle$, C , k_{nn} , etc.) is due to some intrinsic structure, and to what extent is it random?

→ Let's randomise the network!