

# The science of networks

- (1) Real world networks and their universal properties
- (2) Mathematical framework
- (3) Networks of configuration type
- (4) Preferential attachment networks

# Real world networks and their universal properties

Networks consists of a large, but finite, number of **nodes** connected by **links**.  
In the modern world, networks are **ubiquitous**:

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Scientists who have studied these networks have made a number of surprising discoveries, based on **statistical**, **numerical** and **nonrigorous analytical** arguments. Our aim is to put these claims in a rigorous mathematical framework and to verify or refine them in this framework.

# Real world networks and their universal properties

A crucial claim often made is that these networks are **scale-free**. In the *language of network science* this means that

$$\frac{\text{\#nodes with degree } k}{\text{\#nodes in the network}} \approx k^{-\tau},$$

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Some estimated power-law exponents:

- an **e-mail** network at the University of Kiel:  $\tau \approx 1.81$ ,
- the **world-wide web**:  $\tau \approx 2.1$ ,
- the **internet**:  $\tau \approx 2.2$ ,
- the **movie actor network**:  $\tau \approx 2.3$ ,
- the **collaboration graph** in mathematics:  $\tau \approx 2.4$ ,
- a network of **sexual relationships** in Sweden:  $\tau \approx 3.3$ .

# Real world networks and their universal properties

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The typical distances observed in networks are often very small. In the *language of network science* networks are called **small worlds** if the distance of two randomly chosen nodes in the giant component is of order  $\log N$ , and they are called **ultrasmall** if it is of order  $\log \log N$ , where  $N$  is the total number of nodes.

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## Claim 3

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There are a lot of further claims in the network sciences literature, typically about **processes on the networks** but we stop here and start with our journey looking at mathematical models and results.

# Real world networks and their universal properties

A small selection of references:

- [Albert, Jeong, Barabasi](#). Error and attack tolerance of complex networks. Nature 406, 378-382 (2000)
- [Albert, Barabasi](#). Statistical mechanics of complex networks. Rev. Mod. Phys. 74, 4797 (2002)
- [Cohen, Havlin](#). Scale-free networks are ultrasmall. Phys. Rev. Lett. 90, 058701 (2003)
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# Mathematical framework

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The general approach is to define a **sequence  $(\mathcal{G}_N)$  of random graphs** with  $N$  vertices and study **asymptotic** properties as  $N$  goes to infinity. In this framework we can give rigorous definitions of the main notions of network science.

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The general approach is to define a **sequence** ( $\mathcal{G}_N$ ) of **random graphs** with  $N$  vertices and study **asymptotic** properties as  $N$  goes to infinity. In this framework we can give rigorous definitions of the main notions of network science.

We always assume that the vertices of  $\mathcal{G}_N$  are labelled as  $1, \dots, N$  and define the **empirical degree distribution** of ( $\mathcal{G}_N$ ) as ( $X_N(k): k = 0, 1, \dots$ ) where

$$X_N(k) = \frac{1}{N} \sum_{i=1}^N \mathbf{1}\{\text{degree of vertex } i = k\}.$$

We call ( $\mathcal{G}_N$ ) **scale-free** with **power-law exponent**  $\tau$  if

$$\lim_{N \rightarrow \infty} X_N(k) = \mu(k) \quad \text{in probability,}$$

for some nonrandom probability vector ( $\mu(k): k = 0, 1, \dots$ ) and

$$\lim_{k \rightarrow \infty} \frac{\log \mu(k)}{\log k} = -\tau.$$

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Let  $\mathcal{C}_N \subset \mathcal{G}_N$  be the **largest connected component** of the network. We say that  $(\mathcal{G}_N)$  has a **giant component** if

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Given  $\mathcal{G}_N$  and a deletion parameter  $q$  we obtain the **percolated network**  $\mathcal{G}_N(q)$  by removing every edge of  $\mathcal{G}_N$  independently with probability  $q$ . We say the network is **robust** if, for every  $0 < q < 1$  the network  $(\mathcal{G}_N(q))$  has a giant component.

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Given  $\mathcal{G}_N$  we let  $d(\cdot, \cdot)$  be the **graph distance** of two vertices, i.e. the length of the shortest path between them. Picking two vertices  $V, W \in \mathcal{C}_N$  independently, uniformly from  $\mathcal{C}_N$ , we say the network is **ultrasmall** if

$$\lim_{N \rightarrow \infty} \frac{d(V, W)}{\log \log N} = c > 0 \quad \text{in probability.}$$

# Networks of configuration type

The most studied network model in the mathematical literature is the sparse **Erdős-Rényi graph** . Given a parameter  $p > 0$  we obtain  $\mathcal{G}_N$  by putting an edge between any two vertices independently with probability  $p/N$ .

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In this model the degree of a vertex is binomially distributed with parameters  $N - 1$  and  $p/N$ , so that by the **law of small numbers**

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In particular these networks are **not scale free** as the asymptotic degree distribution has light tails. It is not a suitable model for 'real' networks.

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Take  $D_1, D_2, \dots$  iid random variables with

$$\mathbb{P}\{D_1 > x\} = x^{1-\tau}(c + o(1)) \quad \text{as } x \uparrow \infty$$

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Given  $D_1, \dots, D_N$  we construct the **network**  $\mathcal{G}_N$  as follows:

- To any vertex  $n \leq N$  we attach  $D_n$  **half-edges** or **stubs**.
- We start by **matching** a stub with a (uniformly) randomly chosen other stub, and continue matching every unpaired stub with a remaining randomly chosen stub until all (or all but one) stubs are matched.
- Any matched pair of stubs are **connected** to form an edge.

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Obviously the resulting network can have self-loops and multiple edges. However, if  $\tau > 2$ , the network has **power law exponent**  $\tau$  even if self-loops and multiple edges are removed.

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We describe a model introduced by **Norros and Reittu** under the name **conditionally Poissonian random graph**. It is based on drawing an iid fitness sequence  $\Lambda_1, \Lambda_2, \dots$  with

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Conditional on this sequence, the network is constructed as follows:

- $\mathcal{G}_1$  consists of a single vertex and no edges,
- given  $\mathcal{G}_N$  we **insert** one new vertex and, independently for any  $n \leq N$  introduce a random number of edges between the new vertex and  $n$  according to a **Poisson distribution** with mean

$$\frac{\Lambda_n \Lambda_{N+1}}{\sum_{k=1}^{N+1} \Lambda_k},$$

- we further **remove** each edge in  $\mathcal{G}_N$  independently with probability

$$1 - \frac{\sum_{k=1}^N \Lambda_k}{\sum_{k=1}^{N+1} \Lambda_k},$$

and thus obtain  $\mathcal{G}_{N+1}$ .

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The **preferential attachment paradigm** claims to offer a simple and credible explanation for the occurrence of scale-free networks. At the same time it gives rise to a **very nice class of network models**, which can still be studied rigorously, although they are more complex than the configuration type models.

# Networks of configuration type

A small selection of references:

- [Aiello, Chung, Lu](#). A random graph model for power law graphs. *Experiment. Math.* 10, 53-66 (2001)
- [Britton, Deijfen, Martin-Löf](#). Generating simple random graphs with prescribed degree distribution. *J. Statist. Phys.* 124, 1377–1397 (2006).
- [van der Hofstad, Hooghiemstra](#). Universality for distances in power-law random graphs. *J. Math. Phys.* 49, 125-209 (2008).
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- [Norros, Reittu](#). Attack resistance of power-law random graphs. *Internet Math.* 5, 251-266 (2008)

# Preferential attachment networks

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Growing networks are built by [adding vertices successively](#). When a new vertex is introduced, attachment to vertices with [higher degree is preferred](#), following the principle that [the rich get richer](#). Roughly speaking, a new vertex is connected by edges to a fixed or random number of existing nodes with a probability [proportional to a nondecreasing function  \$f\$  of their degree](#). The function  $f$ , which regulates the strength of the preferential attachment is called the [attachment rule](#).

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We first discuss a version of the model where new vertices are connected to a [fixed number  \$m \geq 2\$](#)  of old vertices. Here the attachment rule is affine, more precisely there exist  $\delta > -m$  such that  $f(k) = k + \delta$ . The case  $\delta = 0$  is studied extensively in the work of [Bollobas and Riordan](#).

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- $\mathcal{G}_1$  consists of a single vertex with  $m$  self loops.
- In each further step, given  $\mathcal{G}_N$ , we **insert one new vertex** and then successively **insert  $m$  edges** connecting the new vertex to vertex  $n \leq N$  with probability

$$\sim (\text{degree of vertex } n) + \delta$$

or to itself with probability

$$\sim (\text{current degree}) + \frac{\delta}{m}.$$

# Preferential attachment networks

## Theorem

Denoting by

$$p_k = \left(2 + \frac{\delta}{m}\right) \frac{\Gamma(k + \delta)\Gamma(m + 2 + \delta + \frac{\delta}{m})}{\Gamma(m + \delta)\Gamma(k + 3 + \delta + \frac{\delta}{m})} \text{ for } k \geq m$$

we have

$$\lim_{N \uparrow \infty} X_N(k) = p_k \quad \text{for all } k, \text{ in probability.}$$

In particular, the network is **scale-free** with power-law exponent

$$\tau = 3 + \frac{\delta}{m}.$$

This was first proved for  $\delta = 0$  by **Bollobas, Riordan, Spencer** and **Tusnady**.

# Preferential attachment networks

Our focus in this course will be on a **preferential attachment network with variable outdegree** introduced by **Dereich and Mörters**. The variability of the outdegree will be used to maximise independence in the network. This makes the model easier to study than the model with fixed outdegree. An immediate advantage is that **nonlinear attachment rules** can be handled.

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Our focus in this course will be on a **preferential attachment network with variable outdegree** introduced by **Dereich and Mörters**.

We fix a concave function  $f: \mathbb{N} \cup \{0\} \rightarrow (0, \infty)$  with  $f(0) \leq 1$  and

$$\Delta f(k) := f(k+1) - f(k) < 1 \quad \text{for all } k \in \mathbb{N}.$$

At time  $N = 1$ , we have a single vertex (labeled 1). In each time step  $N \rightarrow N + 1$  we

- add a **new vertex** labeled  $N + 1$ , and
- for each  $n \leq N$  independently introduce an **oriented edge** from the new vertex  $N + 1$  to the old vertex  $n$  with probability

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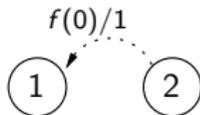
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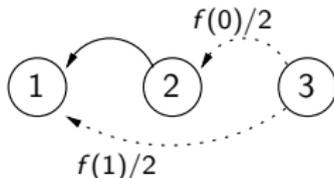
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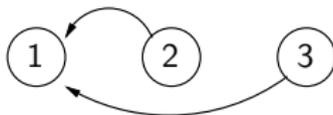
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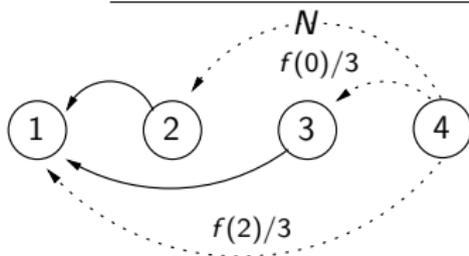
We fix a concave function  $f: \mathbb{N} \cup \{0\} \rightarrow (0, \infty)$  with  $f(0) \leq 1$  and

$$\Delta f(k) := f(k+1) - f(k) < 1 \quad \text{for all } k \in \mathbb{N}.$$

At time  $N = 1$ , we have a single vertex (labeled 1). In each time step  $N \rightarrow N + 1$  we

- add a **new vertex** labeled  $N + 1$ , and
- for each  $n \leq N$  independently introduce an **oriented edge** from the new vertex  $N + 1$  to the old vertex  $n$  with probability

$$\frac{f(\text{indegree of } n \text{ at time } N)}{N}$$



**Example:**

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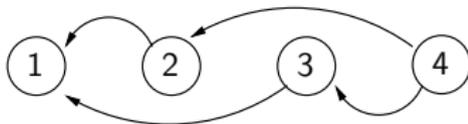
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The empirical indegree distribution of  $\mathcal{G}_N$  is given by

$$X_N^{in}(k) = \frac{1}{N} \sum_{i=1}^N \mathbf{1}\{\text{indegree of vertex } i = k\}.$$

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## Theorem 1

Denoting by

$$\mu(k) = \frac{1}{1 + f(k)} \prod_{l=0}^{k-1} \frac{f(l)}{1 + f(l)},$$

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The limit

$$\gamma := \lim_{k \uparrow \infty} \frac{f(k)}{k} = \inf_{n \geq 1} \Delta f(n)$$

exists by concavity and, by Theorem 1, under the assumption that  $\gamma > 0$  the network is **scale-free** with power-law exponent

$$\tau = \frac{1 + \gamma}{\gamma}.$$

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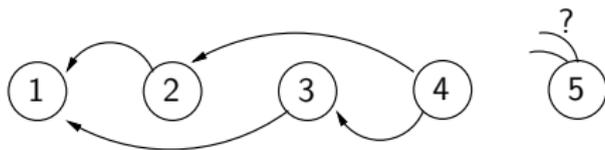
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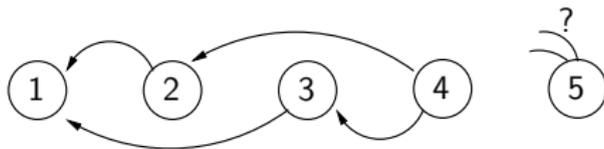
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If  $f(k) \sim k^\alpha$  for  $0 \leq \alpha < 1$ , then  $\log \mu(k) \sim -\frac{1}{1-\alpha} k^{1-\alpha}$ .

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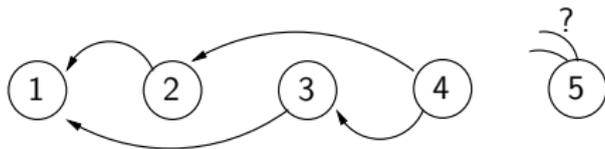


## Theorem 2

The **conditional distribution of the outdegree** of the vertices with label  $N + 1$ , given the graph at time  $N$ , converges almost surely in the variational topology to the **Poisson distribution** with parameter

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The outdegree distribution is therefore **light-tailed** and does not interfere with the power-law exponent.

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A small selection of references:

- **Barabasi, Albert.** Emergence of scaling in random networks. *Science* 286, 509–512 (1999)
- **Bollobas, Riordan, Spencer, Tusnady.** The degree sequence of a scale-free random graph process. *Random Structures Algorithms* 18, 279–290 (2001).
- **Dereich, Mörters.** Random networks with sublinear preferential attachment: Degree evolutions. *Electronic Journal of Probability* 14, 1222-1267 (2009).
- **Krapivsky, Redner.** Organization of growing random networks. *Phys. Rev. E* 63, 066123 (2001).