

Typical distances in ultrasmall random networks

- (1) Ultrasmall networks: a reminder
- (2) Typical distances in configuration networks
- (3) Typical distances in preferential attachment networks
- (4) A model-free approach to lower bounds

Ultrasmall networks: a reminder

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 independently, uniformly from the giant component, we say the network is **ultrasmall** if

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

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Note: We don't expect to gain significant insight into the case $\tau \leq 2$ as our models are undefined or degenerate in this case.

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Note: We expect results about the **diameter** of the giant component to depend considerably on the model details, and therefore to be of less interest.

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Networks with fixed degree sequence use a sequence D_1, D_2, \dots of iid random variables with

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

Given D_1, \dots, D_N we construct the network \mathcal{G}_N by attaching D_i half-edges to the vertex labelled i , and **matching them at random**.

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Theorem (van der Hofstad and Hooghiemstra 2008)

The networks with given fixed degree sequence are **ultrasmall** if and only if

$$2 < \tau < 3.$$

Moreover, for independent, uniformly chosen vertices V and W in the giant component of \mathcal{G}_N , we have

$$\lim_{N \rightarrow \infty} \frac{d(V, W)}{\log \log N} = \frac{2}{-\log(\tau - 2)} \quad \text{in probability.}$$

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Remark: The limit depends only on τ .

Typical distances in configuration networks

Conditionally Poissonian random graphs are based on drawing an iid sequence $\Lambda_1, \Lambda_2, \dots$ of positive fitness values with

$$\mathbb{P}\{\Lambda_1 > x\} = x^{1-\tau}(c + o(1)) \quad \text{as } x \uparrow \infty$$

Conditional on this sequence, we independently connect vertices n, m in \mathcal{G}_N by a Poissonian number of vertices with mean

$$\frac{\Lambda_n \Lambda_m}{\sum_{k=1}^N \Lambda_k}.$$

The conditionally Poissonian random graph is **scale-free** with power-law exponent τ .

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The networks with heavy-tailed fitness distribution are **ultrasmall** if and only if

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Moreover, for independent, uniformly chosen vertices V and W in the giant component of \mathcal{G}_N , we have

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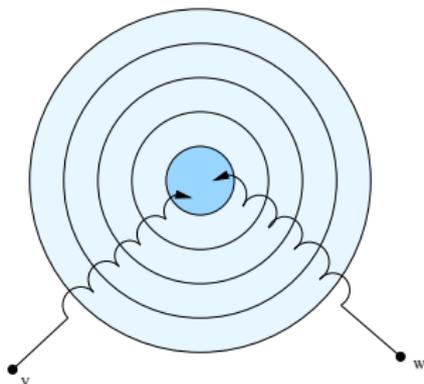
$$\lim_{N \rightarrow \infty} \frac{d(V, W)}{\log \log N} = \frac{2}{-\log(\tau - 2)} \quad \text{in probability.}$$

Remark: The limit is the same as in the first example.

Typical distances in configuration networks

At least heuristically we have some **structural insight** into typical shortest paths in ultrasmall configuration networks:

- typical vertices in the giant component can be connected with a few steps to a **core** of the network;
- within this core there is a hierarchy of **layers** of nodes with increasing connectivity and at the top a small **inner core** of highly connected nodes with very small diameter;
- a shortest path inside the core moves from one layer to the next until the inner core is reached, and then climbing down again until a vertex in the lowest layer of the core is again connected to a typical vertex.



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The j th layer consists of vertices with degree k_j where

$$\log k_j \approx (\tau - 2)^{-j}$$

and there are about

$$\frac{\log \log N}{-\log(\tau - 2)}$$

layers. The graph distance of two randomly chosen vertices in the giant component is therefore

$$(2 + o(1)) \frac{\log \log N}{-\log(\tau - 2)}.$$

Typical distances in preferential attachment networks

We first look at **preferential attachment networks with fixed outdegree**, given by parameters $\delta > -m$ where $m \geq 2$ is an integer.

- \mathcal{G}_1 consists of a single vertex with m self loops.
- 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

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

This network is scale-free with power-law exponent

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

and we expect it to be **ultrasmall** iff $\delta < 0$.

Typical distances in preferential attachment networks

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Theorem

For the preferential attachment model with $\delta = 0$ and independent, uniformly chosen vertices V and W in the giant component of \mathcal{G}_N ,

$$\lim_{N \rightarrow \infty} d(V, W) \frac{\log \log N}{\log N} = 1 \quad \text{in probability.}$$

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For the preferential attachment model with $\delta < 0$ and independent, uniformly chosen vertices V and W in the giant component of \mathcal{G}_N ,

$$\limsup_{N \rightarrow \infty} \frac{d(V, W)}{\log \log N} \leq \frac{4}{-\log(\tau - 2)} \quad \text{in probability.}$$

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Problems:

- Find a [lower bound](#) and thereby verify ultrasmallness if $2 < \tau < 3$.
- Identify the [correct limit](#). Is this limit [universal](#)?
- Find a similar result for preferential attachment networks with [variable](#) outdegree.

A model-free approach to lower bounds

Our first result is based on the following assumption.

Assumption PA(γ)

There exists κ such that, for all pairwise distinct vertices v_0, \dots, v_ℓ in \mathcal{G}_N ,

$$\mathbb{P}\{v_0 \leftrightarrow v_1 \leftrightarrow v_2 \leftrightarrow \dots \leftrightarrow v_\ell\} \leq \prod_{k=1}^{\ell} \kappa (v_{k-1} \wedge v_k)^{-\gamma} (v_{k-1} \vee v_k)^{\gamma-1}.$$

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For preferential attachment models with **fixed or variable outdegree** and power law exponent τ , we can easily verify that

$$\gamma > (\tau - 1)^{-1} \implies \text{Assumption PA}(\gamma).$$

Hence we **expect** networks to be **ultrasmall** if PA(γ) holds for $\frac{1}{2} < \gamma < 1$.

A model-free approach to lower bounds

Theorem 9

Suppose \mathcal{G}_N satisfies **Assumption PA**(γ) for some $\frac{1}{2} < \gamma < 1$. For random vertices V and W chosen independently and uniformly from \mathcal{G}_N , we have

$$d(V, W) \geq 4 \frac{\log \log N}{\log\left(\frac{\gamma}{1-\gamma}\right)} + \mathcal{O}(1) \quad \text{with high probability.}$$

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Corollary 1

The preferential attachment model with fixed outdegree and parameters $\delta > -m$ is **ultrasmall** if and only if $\delta < 0$ or, equivalently $2 < \tau < 3$. Moreover, for independent, uniformly chosen vertices V and W in the giant component of \mathcal{G}_N , we have

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Corollary 2

The preferential attachment model with variable outdegree and attachment rule f is **ultrasmall** if and only if

$$\gamma := \lim_{n \rightarrow \infty} \frac{f(n)}{n} > \frac{1}{2}$$

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We now briefly sketch how assumption **PA**($\gamma + \epsilon$) **can be verified** for the preferential attachment model with variable outdegree.

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For $v < w$, all events $\{v \leftrightarrow w\}$ with different values of v are **independent**. Hence $\mathbb{P}\{v_0 \leftrightarrow \dots \leftrightarrow v_n\}$ can be decomposed into factors of the form $\mathbb{P}\{v \leftrightarrow w\}$ and factors of the form $\mathbb{P}\{u \leftrightarrow v \leftrightarrow w\}$ for $v < u, w$.

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Denoting by $Z[n, N]$ the indegree of vertex n in \mathcal{G}_N we get, for $v < w$,

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For sufficiently large v the increments of f are bounded by $\gamma + \epsilon$, and hence

$$Y_n = f(Z[v, n]) \prod_{i=v}^{n-1} \left(1 + \frac{\gamma + \epsilon}{i}\right)^{-1}$$

defines a **supermartingale**. This implies that

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for a suitable constant $\kappa > 0$, providing the estimate for factors of the **first form**.

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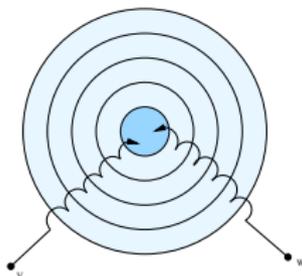
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A similar argument can be used to estimate factors of the **second form**.

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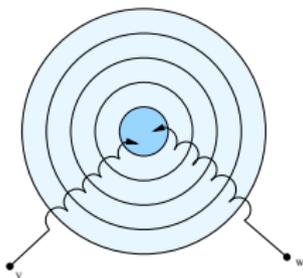
Recall the structural insight into typical shortest paths in ultrasmall configuration networks.



Layers can be identified by vertex degrees, and the j th layer consists of vertices with degree k_j where $\log k_j \approx (\tau - 2)^{-j}$.

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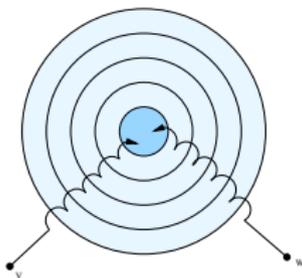


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Therefore a path cannot move directly from one layer to another in one step, but it requires **two steps**: The paths move from one layer to a **young vertex** and from there back into the next higher layer. The distance of two typical vertices is therefore **increased by a factor of two**.

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Both models of configuration type we considered (and more) satisfy $\text{CM}(\gamma)$ for all $\gamma > (\tau - 1)^{-1}$ and we obtain a lower bound from the following theorem.

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Theorem 10

Suppose \mathcal{G}_N satisfies Assumption $\text{CM}(\gamma)$ for some $\frac{1}{2} < \gamma < 1$. For random vertices V and W chosen independently and uniformly from \mathcal{G}_N , we have

$$d(V, W) \geq 2 \frac{\log \log N}{\log\left(\frac{\gamma}{1-\gamma}\right)} + \mathcal{O}(1) \quad \text{with high probability.}$$

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Summary

The results suggest that the ultrasmall networks can be divided into **two universality classes**. In networks of **preferential attachment type** typical vertices have twice the distance compared to networks of **configuration type**. There is also a different structure to shortest paths in the network, with paths alternating between young and old vertices in the case of preferential attachment networks.

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In **real networks**, however, this effect is hard to establish, not least because of the slow growth of $\log \log N$. It also seems that it is often **overruled** by effects not represented in our simple models. For example, in the mathematicians collaboration graph by the effect that the number of authors per paper has increased significantly over the past 50 years, and that mathematicians have a limited period of activity.

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I would however still uphold the claim that this is a good example that **rigorous mathematical analysis** has identified an interesting effect about networks, that can not be identified by simulation or other nonrigorous methods.

A model-free approach to lower bounds

A small selection of references:

- [Bollobas, Riordan](#). The diameter of a scale-free random graph. *Combinatorica* 24, 5–34 (2004)
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