# Random Graph Models

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#### Introduction

- Random graphs theory is one of the most interesting research areas on the intersection of network science and probability.
- It provides a better understanding of underlying mechanisms that create networks.
- Random graphs are used to benchmark the network's algorithms (e.g. clustering algorithms)
- They are also building blocks of synthetic networks that closely resemble real-world networks.

#### Introduction

- Before we proceed with random models, let us briefly introduce some useful notation.
- While discussing random graphs, we mostly focus on their asymptotical behavior, namely what happens when the number of nodes  $n \to \infty$ .
- We say that an event in a given probability space holds asymptotically almost surely (a.a.s.), if its probability tends to one as n → ∞.

#### Introduction

- For describing the properties of random graphs, we will use asymptotic notation. Given two functions f = f(n) and g = g(n):
  - f(n) = O(g(n)) if there exists an absolute constant c such that  $f(n) \le c g(n)$  for all n.
  - $f(n) = \Omega(g(n))$  if g(n) = O(f(n)).
  - $f(n) = \Theta(g(n))$  if f(n) = O(g(n)) and  $f(n) = \Omega(g(n))$ .
  - f(n) = o(g(n)) or  $f(n) \ll g(n)$  if the limit  $\lim_{n\to\infty} f(n)/g(n) = 0$ .
  - $f(n) = \omega(g(n))$  or  $f(n) \gg g(n)$  if g(n) = o(f(n)).
  - $f(n) \sim g(n)$  if f(n) = (1 + o(1))g(n).

# Binomial Random Graphs

# **Binomial Random Graphs**

- There are three commonly used random graph models:
  - Binomial random graphs
  - Uniform random graphs
  - Erdős-Rényi random graphs
- They are in many cases asymptotically equivalent, thus selection of a specific model depends on its theoretical and computational properties.

# Binomial random graphs

The binomial random graph G(n,p) can be generated by starting with the empty graph on the set of nodes  $[n] = \{1,2,\ldots,n\}$ . For each pair of nodes i,j such that  $1 \leq i < j \leq n$ , we independently introduce an edge i j in G with probability p.

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- G(n, p) is defined as the probability distribution over a family of labeled graphs on n nodes.
- There are  $\binom{N}{m}$  labeled graphs on n nodes and m edges, where  $N=\binom{n}{2}$  is the number of pairs of nodes.
- For a given labeled graph G on n nodes and m edges:

$$\mathbb{P}(G) = p^m (1 - p)^{N - m}$$

# **Uniform Random Graph**

Let  $\Omega$  be the family of all labeled graphs on the set of nodes [n] and exactly m edges, where  $0 \le m \le N$ ,  $N = \binom{n}{2}$ . The uniform random graph G(n,m) assigns to every graph  $G \in \Omega$  the same probability, that is,

$$\mathbb{P}(G) = \frac{1}{|\Omega|} = \binom{N}{m}^{-1}$$

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•  $G(n,p) \approx G(n,m)$ , provided  $m \approx \binom{n}{2}p$ .

# Erdős-Rényi Random Graph

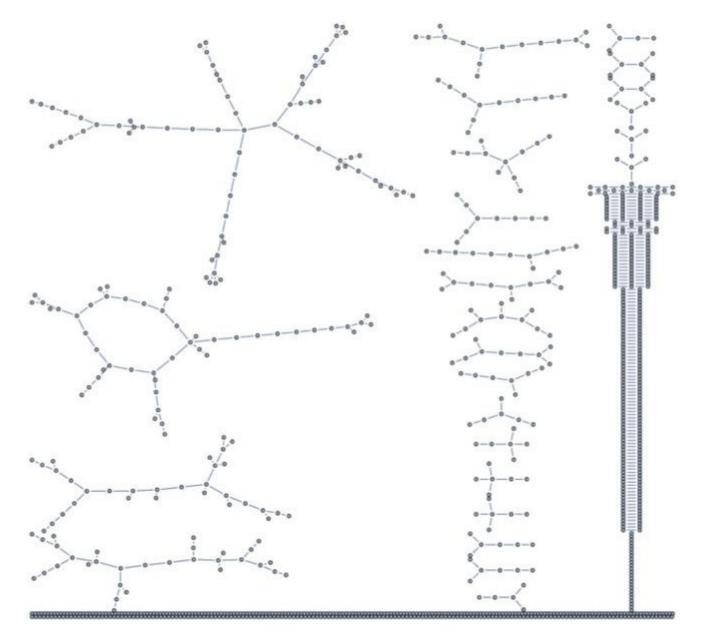
The Erdős-Rényi random graph process is a stochastic process that starts with n labeled nodes and no edges, and at each step adds one new edge chosen uniformly at random from the set of missing edges. Formally, let  $N=\binom{n}{2}$  and let  $e_1,e_2,\ldots,e_N$  be a random permutation of the edges of the complete graph  $K_n$ . The graph process consists of the sequence of random graphs  $(\mathcal{G}(n,m))_{m=0}^N$ , where  $\mathcal{G}(n,m)=([n],E_m)$  and  $Em=\{e_1,e_2,\ldots,e_m\}$ . It is clear that  $\mathcal{G}(n,m)$  is a graph taken uniformly at random from the set of all graphs on n nodes and m edges.

# **Emergence of the giant component**

#### **Subcritical phase:**

$$\langle k \rangle < 1 - \epsilon$$
 for some  $\epsilon > 0$ .

A.a.s. G(n, p) consists of small trees and unicyclic components; the size of the largest component is  $O(\ln n) = o(n)$ .



G(1000, 0.95/1000)

# **Emergence of the giant** component

Critical phase:  $\langle k \rangle \sim 1$ 

The giant component is formed. During that phase, the size of the largest component keeps growing reaching  $\Theta(n^{2/3})$  nodes at precisely  $\langle k \rangle = 1$  a.a.s.

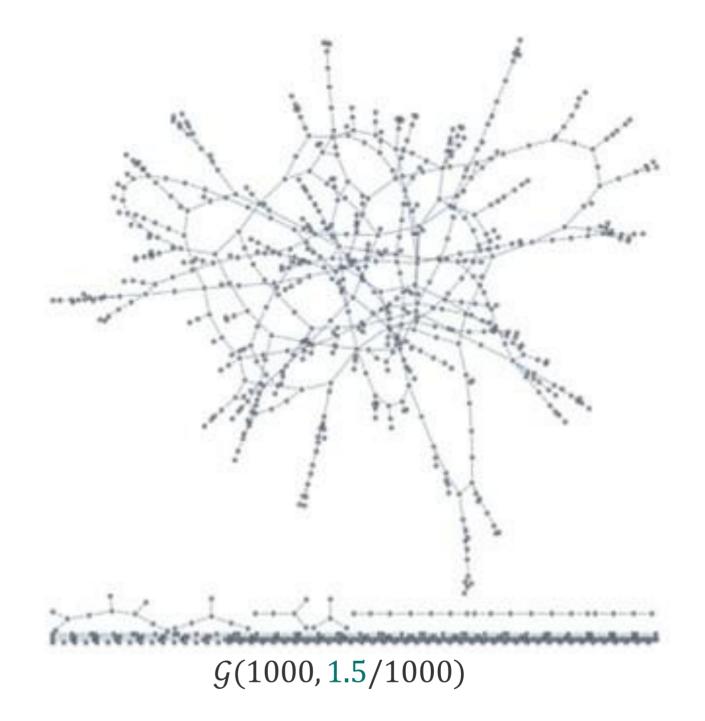
# **Emergence of the giant** component

#### **Supercritical phase:**

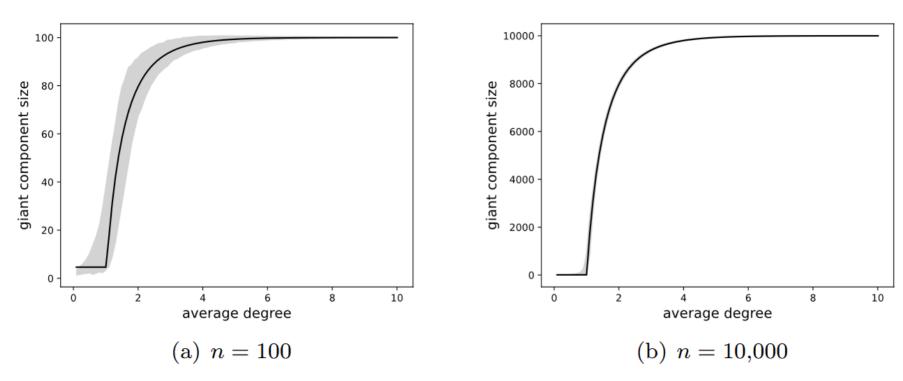
$$\langle k \rangle > 1 + \epsilon$$
 for some  $\epsilon > 0$ .

A.a.s. the size of the giant component is  $(1 + o(1))\beta n$ , where  $\beta + e - \beta \cdot \langle k \rangle = 1$ .

The giant component is unique and the second-largest component is acyclic or unicyclic and has size  $O(\ln n)$ .



# **Binomial Random Graphs**



The order of the giant component: theoretical predictions and empirical results based on 1,000 independent runs for small graphs on (a) n = 100 nodes and larger graphs on (b) n = 10,000 nodes.

# **Connectivity of Binomial Random Graphs**

Assume that

$$p = p(n) = \frac{\ln n + c}{n} \sim \frac{\ln n}{n}$$
 for some  $c \in \mathbb{R}$ 

 then the expected number of isolated nodes is approximately equal to:

$$e^{-c}$$

- Clearly:
  - If  $c \to -\infty$ , then expect isolated nodes.
  - $c \to +\infty$ , then expect the opposite.

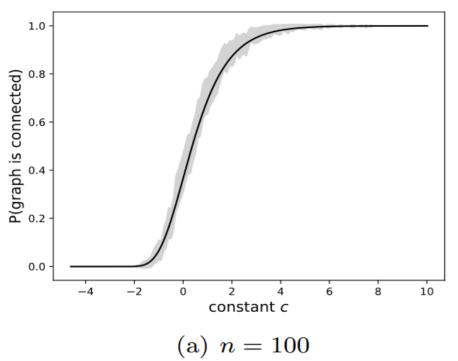
# **Connectivity of Binomial Random Graphs**

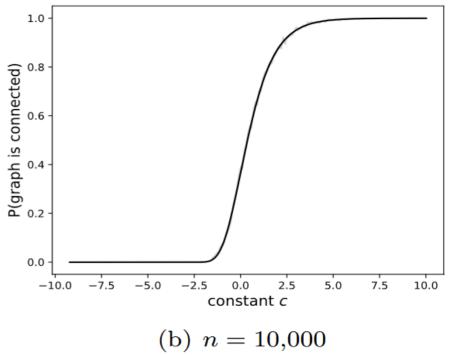
• Once isolated nodes disappear, G(n, p) is connected a.a.s.

$$\mathbb{P}(\mathcal{G}(n,p) \text{ is connected}) \sim \begin{cases} 0 \text{ if } c \to -\infty \\ e^{-c} \text{ if } c \in \mathbb{R} \\ 1 \text{ if } c \to +\infty \end{cases}$$

- For  $any \epsilon > 0$  we have that:
  - If  $pn < (1 \epsilon) \ln n$ , then a.a.s.  $\mathcal{G}(n, p)$  is disconnected.
  - If  $pn > (1 + \epsilon) \ln n$ , then a.a.s.  $\mathcal{G}(n, p)$  is **connected**.

# **Connectivity of Binomial Random Graphs**





he probability that G(n,p) with  $np = \ln n + c$  is connected: theoretical predictions and empirical estimations based on 1,000 independent runs for small graphs on (a) n = 100 nodes and larger graphs on (b) n = 10,000 nodes.

# Degree Distribution of Binomial Random Graphs

• Consider G(n, p) with p = p(n) = c/n for some constant  $c \in \mathbb{R}$  + and large n. For any node  $v \in [n]$ ,

$$\mathbb{E}[deg(v)] = p \cdot (n-1) \sim c$$

- But G(n, p) is not a c-regular graph!
- For any  $\ell \in \mathbb{N} \cup \{0\}$ :

$$\mathbb{P}(\deg(v) = \ell) \sim \frac{c^{\ell}}{\ell!} e^{-c}$$

# **Degree Distribution of Binomial Random Graphs**

• In the limit, the degree distribution of sparse G(n, p) can be approximated by the Poisson distribution, that is,

$$d_{\ell} \sim \frac{c^{\ell}}{\ell!} e^{-c}$$

where c is the asymptotic expected average degree.

• In particular, a.a.s. the maximum degree is  $O(\ln n/\ln \ln n)$ .

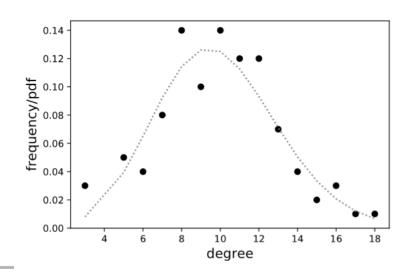
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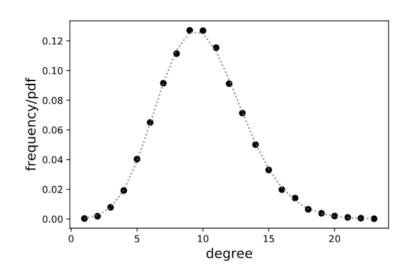
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#### Power-law distribution

- Real-world networks typically do not have the Poisson distribution.
- Typically, degree distribution follows power law:

$$d_{\ell} \approx c * \ell^{\gamma}$$

where  $\gamma > 0$  is a **degree exponent** and c > 0 is some normalizing constant.

Degree distribution can be easily rewritten as:

$$\ln d_{\ell} \approx -\gamma \ln \ell + \ln c$$

(straight line with the slope  $-\gamma$  on a log-log plot).

#### Power-law distribution

• To generate a power-law degree distribution with a given degree exponent  $\gamma$  and minimum degree  $\delta \geq 1$  we assume:

$$d_{\ell} \approx c \int_{\ell}^{\ell+1} x^{-\gamma} dx \approx c * \ell^{\gamma}$$

Note that:

$$1 = \sum_{l=\delta}^{\infty} d_{\ell} = c \int_{\delta}^{\infty} x^{-\gamma} dx = \frac{c(-\delta^{1-\gamma})}{1-\gamma}$$

Finally:

$$d_{\ell} \approx (\gamma - 1)\delta^{\gamma - 1}\ell^{-\gamma}$$

#### Power-law distribution

Average degree:

$$\langle k \rangle = \frac{\gamma - 1}{\gamma - 2} \delta$$

provided that  $\gamma > 2$ .

- Maximum degree:
  - Assuming that the number of nodes of degree at least Δ is close to 1 we get that:

$$\Delta = \delta n^{1/(\gamma - 1)}$$

which is often referred to as the **natural cut-off** of the graph.

Let  $\mathbf{w} = (w_1, w_2, ..., w_n)$  be any vector of n positive real numbers and let  $W = \sum_{i=1}^{n} w_i$ . Random graph  $G(\mathbf{w})$  is generated as follows:

1. Each pair of nodes i, j such that  $1 \le i \le j \le n$  is independently sampled as an edge (or loop if i = j) with probability given by:

$$p_{i,j} \begin{cases} \frac{w_i w_j}{W} & \text{for } i \neq j \\ \frac{w_i^2}{W} & \text{for } i = j \end{cases}$$

• For any any  $i \in [n]$ :

$$\mathbb{E}[\deg(i)] = w_i$$

- The model is also well-studied but the behavior and results are more complex; for example, for any  $\epsilon > 0$ :
  - if  $\langle k \rangle \leq \langle k^2 \rangle / \langle k \rangle < 1 \epsilon$ , then a.a.s. G(w) has no giant,
  - if  $\langle k^2 \rangle / \langle k \rangle \ge \langle k \rangle > 1 + \epsilon$ , then a.a.s. there is one.

Generating power-law graphs with degree exponent γ:

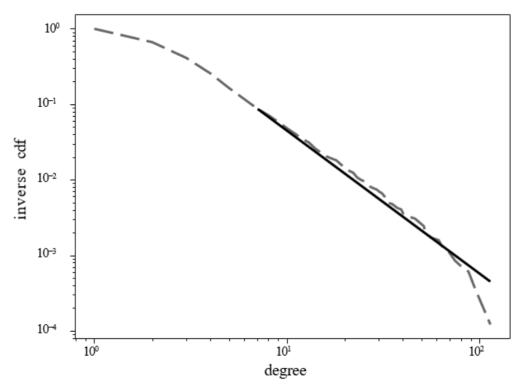
$$w_i = c \cdot (i + i_0 - 1)^{-1/(\gamma - 1)}$$

c=c(n) depends on the minimum (or average) degree  $\delta \geq 1$  and  $i_0=i_0(n)$  depends on the maximum degree  $\Delta$ .

It follows that 
$$c = \delta n^{1/(\gamma-1)}$$
 and  $i_0 = n/\left(\frac{\Delta}{\delta}\right)^{\gamma-1}$  so

$$w_i = \delta \left( \frac{n}{i - 1 + n/(\Delta/\delta)^{\gamma - 1}} \right)^{1/(\gamma - 1)}$$

• It is possible to show that the expected number of nodes of degree k is proportional to  $\Gamma(k-\gamma+1)/\Gamma(k+1)\approx k^{-\gamma}$ , where  $\Gamma(z)=\int_0^\infty x^{z-1}e^{-x}dx$  is the gamma function.



 $\mathcal{G}(w)$  generated on  $n=10{,}000$  nodes using the set of weights with  $\gamma=2.5, \delta=1$ , and  $\Delta=\sqrt{n}=100$ . We got  $\delta'=0$  and  $\Delta'=113$ , respectively. But did we preserve the degree exponent?

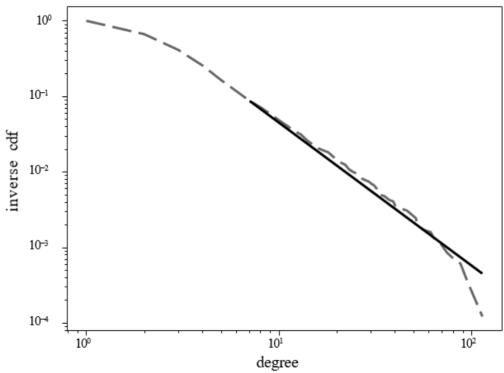
Kolmogorov–Smirnov statistic: focus only on the large degrees.
For a given cutoff for small degrees ℓ ∈ [max{δ, 1}, Δ],

$$\gamma^{\ell} = 1 + \frac{|\{j: \deg(j) \ge \ell|}{\sum_{j: \deg(j) \ge \ell} \ln\left(\frac{\deg(j)}{\ell - 1/2}\right)}$$

 The divergence of the experimental distribution from the theoretical one is defined as:

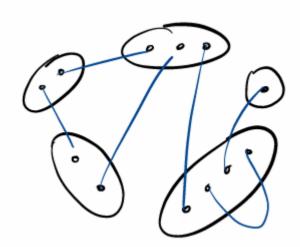
$$D_{\ell} = \max_{k \in [\ell, \Delta]} \left| \frac{|\{j: \deg(j) \ge k|}{|\{j: \deg(j) \ge \ell|} - \frac{\int_{k}^{\infty} x^{-\gamma \ell} dx}{\int_{\ell}^{\infty} x^{-\gamma \ell} dx} \right|$$

• The value of  $\gamma_{\ell}$  that minimizes  $D_{\ell}$  (over all  $\ell \in [\max{\delta, 1}, \Delta]$ ) is used as an estimate  $\gamma'$  of the power-law exponent.



G(w) generated on  $n=10{,}000$  nodes using the set of weights with  $\gamma=2.5, \delta=1$ , and  $\Delta=\sqrt{n}=100$ . We got  $\delta'=0$  and  $\Delta'=113$ , respectively. The fitted line has slope of -1.89 ( $\gamma'=2.89$ ), which was obtained with the procedure we described, with  $\ell'=7$ .

- Chung-Lu model returns a Graph with a degree sequence approximately close to the sequence w.
- Configuration model  $\mathcal{P}_{n,d}$  generates a random graph that strictly follows a given, graphic degree sequence  $d = (\deg(1), \deg(2), ..., \deg(n))$



# Random d-regular graphs

Fix  $d \in \mathbb{N} \cup \{0\}$ . Let  $\Omega$  be the family of all labelled graphs on the set of nodes [n] that are d-regular. The random d-regular graph, denoted by  $\mathcal{G}_{n,d}$ , assigns to every graph  $G \in \Omega$  the same probability, that is,

$$\mathbb{P}(G) = \frac{1}{|\Omega|}$$

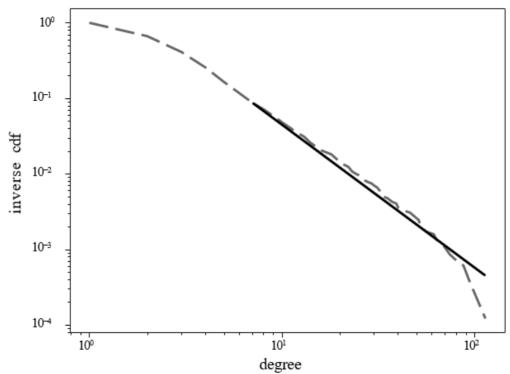
- Since the total volume of any graph is even, n has to be even if d is odd.
- Different mathematical tools are required when d=d(n) grows together with n.
- Generating all d-regular graphs on n nodes is impossible.

# Configuration/pairing model

Consider dn points partitioned into n labelled buckets  $v_1, v_2, \ldots, v_n$  of d points each.

A **pairing** of these points is a perfect matching into dn/2 pairs. Given a random pairing P, we may construct a multigraph  $\mathcal{P}_{n,d}=\mathcal{P}(P)$ , with loops and parallel edges allowed, after contracting buckets into nodes.

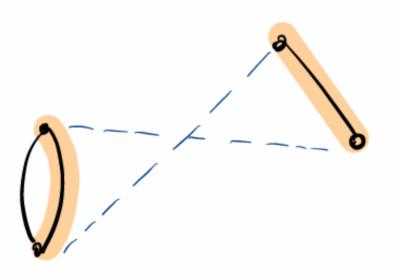
- The restriction of  $\mathcal{P}_{n,d}$  to simple graphs is precisely  $\mathcal{G}_{n,d}$ .
- Practical implication: keep generating  $\mathcal{P}_{n,d}$  (independently) and stop when you get a simple graph; you get  $\mathcal{G}_{n,d}$ .
- Random pairing generates a simple graph with probability asymptotic to  $e^{-(d^2-1)/4}$  depending on d but not on n.
- Practical implication: the expected number of attempts is  $e^{(d^2-1)/4}$  (large for large d but reasonable for small values).
- Theoretical implication: any event holding a.a.s. in  $\mathcal{P}_{n,d}$  also holds a.a.s. in  $\mathcal{G}_{n,d}$ .



 $\mathcal{P}_{n,d}$  is simple with probability  $\sim e^{-(a^2-1)/4}$ Theoretical predictions and empirical results for n=100 and n=10,000.

- Generating  $\mathcal{P}_{n,d}$  is fast but the graph might not be simple!
- Depending on our application, we may:
  - stay with multigraphs;
  - remove all potential loops and parallel edges (erased configuration model); the graph might not match exactly the given sequence d.
  - resample until we get  $\mathcal{G}_{n,d}$  or avoid creating loops/parallel edges and re-start if we get stuck; slow.
  - do **switching**; preserves distribution, typically fast, asymptotically equivalent to  $\mathcal{G}_{n,d}$ .

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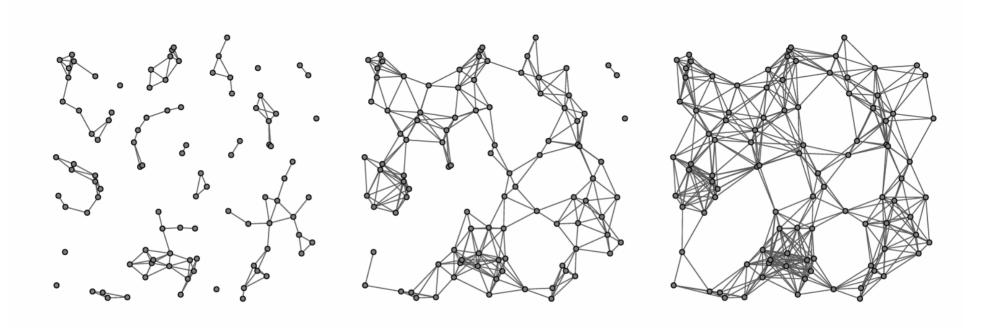


- Often nodes are described by some properties they might impact the existence/strength of an edge.
- While generating a random graph that tries to mimic the real world network, we want to utilize this fact.
- Random Geometric Graphs are a category of models that uses the properties (metadata) to generate a network.

Let  $r \in \mathbb{R} + \cup \{0\}$ . The random geometric graph  $\mathcal{R}\mathcal{G}\mathcal{G}(n,r)$  can be generated by starting with the empty graph on n nodes,  $v_1, v_2, \ldots, v_n$ , that are randomly sampled from the uniform distribution of the underlying space  $[0,1]^d$ . Each pair of nodes  $v_i, v_j$  such that  $1 \le i < j \le n$  is connected by an edge if and only if  $d(v_i, v_j) \le r$ .

- The choice of metric space is important (especially, if d is large)
- For continuous (especially spatial) data one might use the Euclidean distance:

$$d(x,y) = \sqrt{\sum_{i=1}^{d} (x_i - y_i)^2}$$



Instances of RGG(100,r) for  $r \in \{0.1, 0.15, 0.2\}$  on the unit square (d = 2).

Threshold for connectivity is well understood but, for example, the appearance of the giant component is not!