## Complex Networks Master of Science in Electrical Engineering

Erivelton Geraldo Nepomuceno

Department of Electrical Engineering Federal University of São João del-Rei

April 30, 2019

### 3. Random Graphs

- The term random graph refers to the disordered nature of the arrangement of links between different nodes.
- The systematic study of random graphs was initiated by Erdös and Rényi in the late 1950s.
- We focus our attention on the shape of the degree distributions and on how the average properties of a random graph change as we increase the number of links.

# 3.1 Erdös and Rényi (ER) Models

- A random graph is a graph in which the edges are randomly distributed.
- In the late 1950s, two Hungarian mathematicians, Paul Erdös and Alfréd Rényi came up with a formalism for random graphs that would led to modern graph theory.
- We shall deal with undirected graphs.
- This method is called Erdös and Rényi (ER) random graphs.

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### On random graphs I.

Dedicated to O. Varga, at the occasion of his 50th birthday.

By P. ERDŐS and A. RÉNYI (Budapest).

Let us consider a "random graph"  $\Gamma_{n,N}$  having *n* possible (labelled) vertices and *N* edges; in other words, let us choose at random (with equal probabilities) one of the  $\binom{\binom{n}{2}}{N}$  possible graphs which can be formed from

Figure 1: [1] P. Erdos and A. Rényi, "On random graphs I," Publ Math, vol. 6, pp. 290–297, 1959. GS: 14388.

#### Definition 1 (ER model A: uniform random graphs)

Let  $0 \le K \le M$ , where M = N(N-1)/2. The model, denoted as  $G_{N,K}^{ER}$ , consists in the ensemble of graphs with *N* nodes generated by connecting *K* randomly selected pairs of nodes, uniformly among the *M* possible pairs. Each graph  $G = (\mathcal{N}, \mathcal{L})$  with  $|\mathcal{N}| = N$  and  $K = |\mathcal{L}|$  is assigned the same probability.

• Consider *N* nodes and picks at random the first edge among the *M*possible edges, so that all these edges are equiprobable. Then one chooses the second edge at random with a uniform probability among the remaining M - 1 possible edges, and continues this process until the *K* edges are fixed.

#### Steps to construct a random network<sup>1</sup>

- Start with N isolated nodes.
- Select a node pair and generate a random number between 0 and 1. If the number exceeds *p*, connect the selected node pair with a link, otherwise leave them disconnected.
- **3** Repeat step (2) for each of the N(N-1)/2 node pairs.

<sup>1</sup>Source: Network Science Textbook - Albert-László Barabási. This is a textbook for network science, is freely available under the Creative Commons license.

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- How many different graphs with *N* nodes and *K* edges are there in the ensemble?
- This number is equal to the number of ways we can select *K* objects among *M* possible ones, namely:

 $\frac{M!}{K!(M-K)!}$ 

• This is known as binomial coefficient denoted as

$$C_M^K = \begin{pmatrix} K \\ M \end{pmatrix}$$

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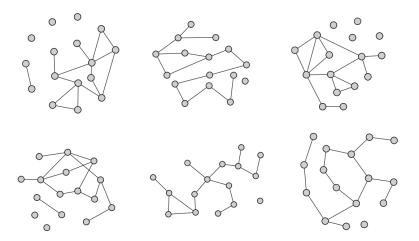


Figure 2: Six different realisations of model A with N = 16 and K = 15.  $C_{120}^{15} \approx 4.73 \times 10^{18}$ .

### Definition 2 (ER model B: binomial random graphs)

Let  $0 \le p \le 1$ . The model, denoted as  $G_{N,p}^{ER}$ , consists in the ensemble of graphs with *N* nodes obtained by connecting each pair of nodes with a probability *p*. The probability  $P_G$  associated with a graph  $G = (\mathcal{N}, \mathcal{L})$  with  $|\mathcal{N}| = N$  and  $|\mathcal{L}| = K$  is  $P_G = p^K (1 - p)^{M-K}$ , where M = N(N-1)/2.

- From  $P_G = p^K (1-p)^{M-K}$ , we can calculate that the probabilities of having an empty or a complete graph with *N* nodes are respectively equal to  $(1-p)^M$  and  $p^M$ .
- However, while all the graphs in the first ensemble have 15 edges, the number of links in the second ensemble.

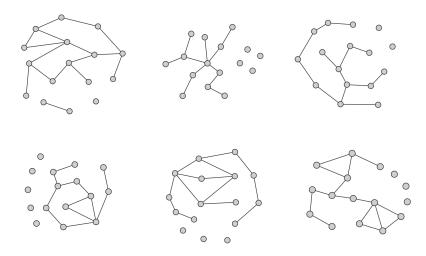


Figure 3: Six different realisations of model *B*, with N = 16 and p = 0.125. We have appropriately chosen the value of *p*, namely pM = 15. Thus, we have  $G_{N,p}^{ER} \equiv G_{N,K}^{ER}$ .

#### Box 1 (Binomial Distribution)

In statistics, the binomial distribution is valid for processes where there are two mutually exclusive possibilities. The binomial distribution is the discrete probability distribution of the number k of successes in a sequence of n independent yes/no experiments, each of which yields success with probability p:

$$Bin(n,p,k) \equiv \binom{n}{k} p^k (1-p)^{n-k}$$

where *k* can take the values 1, 2, 3, ..., n. A simple example is tossing ten coins and counting the number of heads. The distribution of this random number is a binomial distribution with n = 10 and p = 1/2 (if the coin is not biased). As another example, assume 5 per cent of a very large population to be green-eyed. You pick 100 people randomly. The number of green-eyed people you pick is a random variable which follows a binomial distribution with n = 100 and p = 0.05.

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#### Example 1

Suppose a biased coin comes up heads with probability 0.3 when tossed. What is the probability of achieving  $0, 1, \ldots, 6$  heads after six tosses?<sup>*a*</sup>

• 
$$Pr(0 \text{ heads}) = f(0) = Pr(X = 0) = \binom{6}{0} 0.3^0 (1 - 0.3)^{6-0} = 0.117649$$

• 
$$Pr(1 \text{ heads}) = f(1) = Pr(X = 1) = {6 \choose 1} 0.3^1 (1 - 0.3)^{6-1} = 0.302526$$

• 
$$Pr(2 \text{ heads}) = f(2) = Pr(X = 2) = \binom{6}{2}0.3^2(1-0.3)^{6-2} = 0.324135$$

•  $Pr(3 \text{ heads}) = f(3) = Pr(X = 3) = \binom{6}{3}0.3^3(1 - 0.3)^{6-3} = 0.185220$ 

• 
$$Pr(4 \text{ heads}) = f(4) = Pr(X = 4) = \binom{6}{4}0.3^4(1 - 0.3)^{6-4} = 0.059535$$

- $\Pr(5 \text{ heads}) = f(5) = \Pr(X = 5) = \binom{6}{5} 0.3^5 (1 0.3)^{6-5} = 0.010206$
- $Pr(6 \text{ heads}) = f(6) = Pr(X = 6) = \binom{6}{6} 0.3^{6} (1 0.3)^{6-6} = 0.000729$

<sup>a</sup>Source: https://en.wikipedia.org/wiki/Binomial\_distribution

Fluctuations of K in model B is

$$\sigma_K^2 = Mp(1-p).$$

 The ratio between the standard deviation *σ<sub>K</sub>* and the average number of links *K* = *pM* is given by:

$$rac{\sigma_K}{\overline{K}} = \sqrt{rac{(1-p)}{pM}}.$$

• This proves that, in large graphs, the fluctuations in the value of *K* of model *B* can be neglected.

#### Example 2 (Fitting a real network with a random graph)

Suppose you are given a real world network *G* with *N* nodes and adjacency matrix {*aij*} and you want to model it with a binomial random graph ensemble. What is the ensemble of graphs  $G_{N,p}^{ER}$  that best approximates the real network? We can infer the best value of the parameter *p* from maximum likelihood considerations. It is useful to work with the logarithm of  $P_G$ , the so-called log-likelihood  $\mathcal{L}(p)$  that the network *G* belongs to the ensemble:

$$\mathcal{L}(p) = \log P_G(p) = K \log p + [M - K] \log(1 - p)$$

Maximising the log-likelihood with respect to p, i.e. solving the equation  $d\mathcal{L}(p)/dp = 0$ . The solution is p = K/M = 2K/(N(N-1)), where N and K are number of nodes and links in G.

## 3.2 Degree Distribution

- The nice property of random graphs is that they can be studied analytically.
- The degree distribution of ER random graphs can be easily derived analytically in model *B* in the following way.
- If p is the probability that there exists an edge between two generic vertices, the probability that a specific node *i* has degree k<sub>i</sub> equal to k is given by the following expression:

$${
m Prob}_{k_i=k} = C_{N-1}^k 
ho^k (1-
ho)^{N-1-k} \quad 0 \le k \le N-1.$$

 The degree distribution in a random graph is a binomial distribution:

$$p_k = \binom{N-1}{k} p^k (1-p)^{N-1-k} = Bin(N-1, p, k) \quad k = 0, 1, 2, \dots, N-1$$

• Average degree of a randomly chosen node in a random graph:

$$\langle k \rangle = \sum_{k=0}^{N-1} k p_k = p(N-1). \tag{1}$$

The standard deviation around this quantity

$$\sigma_k = \sqrt{p(1-p)(N-1)}$$
(2)

Using the Poisson distribution

$$p_k = \text{Pois}(z, k) = e^{-z} \frac{z^k}{k!}$$
  $k = 0, 1, 2, ...$  (3)

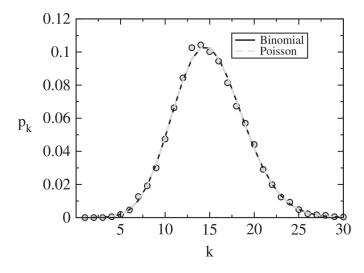


Figure 4: The degree distribution (circles) that results from one realisation of a random graph with N = 10000 nodes and p = 0.0015 is compared to the corresponding binomial (solid line) and Poisson distributions (dashed line).

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### 3.3 Trees, Cycles and Complete Subgraphs

 What are the salient features of a typical random graph with N nodes and K edges?

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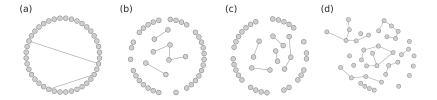


Figure 5: The typical graph of the ensemble  $G_{40,p}^{ER}$  for four different values of p, namely  $p = 2 \times 10^{-3}, 5 \times 10^{-3}, 8 \times 10^{-3}, 3 \times 10^{-2}$ 

• A Random Graph growth. [Video-05].

• GraphStream 1.0 - Dynamics graphs. [Video-06].

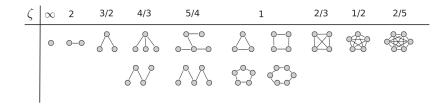


Figure 6: Threshold probabilities  $p(N) \sim N^{-\zeta}$  for the appearance of different subgraphs in a random graph.

### 3.4 Giant Connected Component

 In their works, Erdös and Rényi also discovered a phase transition concerning the order of the largest component in the graph, namely the abrupt appearance of a macroscopic component known as a giant component.

### Definition 3 (Giantcomponent)

A giant component in a graph G is a component containing a number of vertices which increases with the order of G as some positive power of N.

• They proved that a giant component appears at a critical probability function  $p_c(N) = 1/N$ . in model B, or analogously at a critical number of links  $K_c(N) = N/2$  in model A.

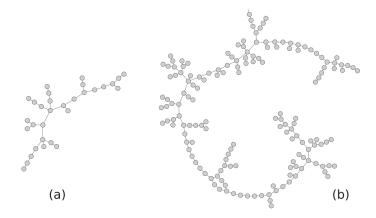


Figure 7: Plot of the largest component  $g_1$  obtained in an ER random graph with N = 1000 nodes and with an average degree respectively below,  $\langle k \rangle = 0.9$  (a), and above,  $\langle k \rangle = 1.1$  (b), the critical value  $\langle k \rangle = 1$ . The largest component  $g_1$  in the first case has  $s_1 = 38$  nodes, while it contains  $s_1 = 202$  nodes, i.e. 20 per cent of the graph nodes, in the second case.

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### Convertion of Vito's file to Pajek format

```
%Write a file.net for Pajek
%Number of Nodes
N = 40;
%Probability
p=8e-3;
fileID = fopen('er.net','w');
fprintf(fileID, '*Vertices %d \n',N);
fprintf(fileID, '*Edges\n');
system(sprintf('er_B %d %f > teste.net',N,p))
A=load('teste.net');
E = [A+1 \text{ ones}(length(A), 1)];
fprintf(fileID,'%d %d %d\n',E');
fclose(fileID);
type er.net
```

### Result

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### **Computational Exercise**

- Use the algorithm in slide 23 to reproduce the results of slide 19.
- Ise the option Fruchterman-Reingold to visualize your graphs.
- Explore the phase transitions properties to find the giant component.
- Develop a code to reproduce the results shown in slide 17.