

Learning from Networks

Significance and Random Graphs

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Motivation

Assume you computed some feature of a graph, e.g. the clustering coefficient.

Now you want to understand if it provides an *interesting* information.

[Published: 04 June 1998](#)

Collective dynamics of ‘small-world’ networks

[Duncan J. Watts](#) ✉ & [Steven H. Strogatz](#)

[Nature](#) **393**, 440–442 (1998) | [Cite this article](#)

Table 1 Empirical examples of small-world networks

	L_{actual}	C_{actual}
Film actors	3.65	0.79
Power grid	18.7	0.080
<i>C. elegans</i>	2.65	0.28

Ideas?



One Solution

Compare your observation with the one obtained from a *random network*.

Intuition: you will understand if what you measured is not explained by the characteristics of a random network \Rightarrow it is interesting!

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Table 1 Empirical examples of small-world networks

	L_{actual}	L_{random}	C_{actual}	C_{random}
Film actors	3.65	2.99	0.79	0.00027
Power grid	18.7	12.4	0.080	0.005
<i>C. elegans</i>	2.65	2.25	0.28	0.05

Simple Examples

Comparison with Random Networks

We need to describe:

- how do we **compare** the observation in the real network with the observation from a random network?
- how do we **compute** the measure on a random network?

Comparison with Random Networks

Scenario:

- you have measured the clustering coefficient for your network;
- you can compute the clustering coefficient for a random network

What would you do to understand if the clustering coefficient of your network is interesting?

Assessing Significance

We are interested in verifying (actually, *falsifying*) whether simple characteristics of the graph explain our measured feature \mathcal{F} (e.g., the clustering coefficient).

We use a framework corresponding to statistical hypothesis testing.

Note: we present it for our scenario, not the most general version.

valore di feature \mathcal{F} osservato in rete reale

We assume the null hypothesis H_0 that (f) well conforms with the distribution observed in a random graph

random graph: graph taken uniformly at random among all graphs (with $|V|$ vertices and) with the required simple characteristics.

Example


Given a graph $G = (V, E)$, we compute its clustering coefficient
 $cc(G)$

Question: is $cc(G)$ explained by the number $|E|$ of edges of G ?

null hypothesis H_0 : $cc(G)$ well conforms with the distribution of
 $cc(G)$ in a random graph


random graph: graph taken uniformly at random among all graphs with $|V|$ vertices and $|E|$ edges (or where the expected number of edges is $|E|$).

G :


$$\Rightarrow cc(G) = \frac{2}{\binom{6}{3}} = \frac{2}{20} = 0.1$$

random graph on 6 nodi e 7 lati
↓

G' :


$$\Rightarrow cc(G') = \frac{1}{\binom{6}{3}} = \frac{1}{20} = 0.05$$

anche G è random graph

Measures of Significance

How do we assess whether the null hypothesis H_0 is true or not?

Note: given that we are considering random graphs, our observation is (almost) always obtainable in at least one or few random graphs

Needed: way to quantitatively assess how likely it is that the value f of feature \mathcal{F} arises when H_0 is true.

Commonly used measures:

- z -score
- p -value

z-score

Let f be the measure of the feature \mathcal{F} of interest.

Let $X_{\mathcal{F}}$ be the random variable corresponding to the value of the feature \mathcal{F} when the null hypothesis H_0 is true, i.e. in a random graph

Definition

The z-score of H_0 is:

$$\frac{f - \mathbb{E}[X_{\mathcal{F}}]}{\sigma[X_{\mathcal{F}}]}$$

Notes:

- the expectation $\mathbb{E}[X_{\mathcal{F}}]$ and standard deviation $\sigma[X_{\mathcal{F}}]$ are w.r.t. the distribution of random graphs given by the null hypothesis H_0
- the z-score can be positive or negative

p -value

Measures the probability of observing a value for feature \mathcal{F} at least as extreme as f when the null hypothesis H_0 is true.

Meaning of “at least as extreme” depends on the feature and the specific situation.

Example:

Given the clustering coefficient $cc(G)$, “at least as extreme” could be:

- $\geq cc(G)$
- or $\leq cc(G)$

p -value (continue)

Definition

The p -value $p(H_0)$ of H_0 is

$$p(H_0) = \mathbb{P}[X_{\mathcal{F}} \text{ is at least as extreme as } f | H_0 \text{ is true}]$$

Notes:

- if we are interested in understanding if the value of f is higher than expected (when H_0 is true):

$$p(H_0) = \mathbb{P}[X_{\mathcal{F}} \geq f | H_0 \text{ is true}]$$

- if we are interested in understanding if the value of f is lower than expected (when H_0 is true):

$$p(H_0) = \mathbb{P}[X_{\mathcal{F}} \leq f | H_0 \text{ is true}]$$

- if $p(H_0)$ is small it is unlikely that the observed value f is obtained when H_0 is true

Example

$$G: \begin{array}{c} \circ \\ \diagup \quad \diagdown \\ \circ \quad \circ \end{array} \Rightarrow cc(G) = \frac{1}{\binom{4}{3}} = \frac{1}{4} = \frac{1}{4}$$

random graph con $|V|=4$, $|E|=4 \Rightarrow$ prendo dist. uniforme

$$X_G = cc(\text{grafo preso da dist.}) \Rightarrow \mathbb{E}[X_G] = 0.2, \sigma[X_G] \approx 0.72 \Rightarrow$$

$$\Rightarrow z\text{-score} = \frac{0.25 - 0.2}{0.72} \approx 0.07$$

$$\Pr[X_G \geq 0.25] = 0.8$$

$$\Pr[X_G \leq 0.25] = 0.5$$

Small Digression

Usually the p -value is used to *reject* a null hypothesis H_0 using the following rule:

Rejection rule: given a value $\alpha \in (0, 1)$, reject H_0 if $p(H_0) \leq \alpha$.

Rejecting H_0 : flagging feature f as *significant*

The following (easy to prove) proposition provides the ground for the rejection rule above.

Proposition

If the rejection rule above is used, then $\mathbb{P}[H_0 \text{ rejected} | H_0 \text{ true}] \leq \alpha$.

Small Digression (continue)

Note that in statistical hypothesis testing two errors are possible:

- H_0 is true *but* H_0 is rejected (type-I error or false positive)
- H_0 is not true *but* H_0 is not rejected (type-II error or false negative)

The previous proposition shows that the **rejection rule** provides guarantees on false positives.

Recap

Up to know: how to compare the observation in the real network with the observation from random networks.

Still missing: what is the *actual* definition of random graph?

From before: graph taken uniformly at random among all graphs (with $|V|$ vertices and) with the required simple characteristics.

In practice?

Erdős-Rényi Random Graphs

Idea: the simple characteristics that is preserved is the number of edges in the graph.


Definition

A random graph from the $G(n, m)$ model is a graph chosen uniformly at random among all graphs with n vertices and m edges.

Note: the vertices are implicitly considered labelled

Introduced by Paul Erdős and Alfréd Rényi in 1959.

Example

G :  $\Rightarrow |V|=4, |E|=4$

$|G(n,m)| = \binom{6}{4} = 15 \Rightarrow \text{probability } 1/15$



...

Erdős-Rényi-Gilbert Random Graphs

Similar to the previous model, but the number of edges is preserved only *in expectation*.

Definition

A random graph from the $G(n, p)$ model is a graph with n vertices and where each edge appears with probability p independently of all other events.

Introduced by Paul Erdős, Alfréd Rényi, and Edgar Gilbert in 1959.

Example

grafo di prima: $G(n, p)$, $|V|=4$, $p \in (0, 1)$

$|G(n, p)| = 2^6 = 64$ (6 archi da cui scegliere)

che valore dare a p per fare confronti?

↓
voglio $\mathbb{E}(m \text{ in grafo da } G(n, p)) = m$
↓
dist. bin

Erdős-Rényi-Gilbert Random Graphs (continue)

Proposition

Let $m(n, p)$ be the number of edges of a random graph from $G(n, p)$. Then

$$\mathbb{E}[m(n, p)] = \binom{n}{2} p$$

$G(n, m)$ vs $G(n, p)$

Informally: if $p = \frac{m}{\binom{n}{2}}$, then $G(n, m)$ and $G(n, p)$ are relatively similar models.

Note:

- the two models are not identical, but several properties are the same
- $G(n, p)$ is usually preferred because it allows for easier analytical results.

Example

What is the expected clustering coefficient of a random graph from $G(n, p)$?

Erdős-Rényi-Gilbert Random Graphs (continue)

However, not all properties can be easily computed analytically!

e.g.: *presence of triangles is not independent \Rightarrow difficult to calculate*

Note: depending on the analysis/property/measure, we need

- the *expectation* and *standard deviation* (e.g., for *z*-score)
- the *distribution* (e.g., for *p*-value)

Example

What Now?



Monte-Carlo Approach

Let f be the measure of the feature \mathcal{F} of interest on your real network.

If you can generate a random graph, then you use the following Monte-Carlo approach:

- 1 generate P instances G_1, G_2, \dots, G_P of a random graph
- 2 compute the measure of interest on the P instances:
 $\mathcal{F}(G_1), \mathcal{F}(G_2), \dots, \mathcal{F}(G_P)$
- 3 use $\mathcal{F}(G_1), \mathcal{F}(G_2), \dots, \mathcal{F}(G_P)$ to estimate the z -value (with estimated expectation and standard deviation) or p -value for f

Notes:

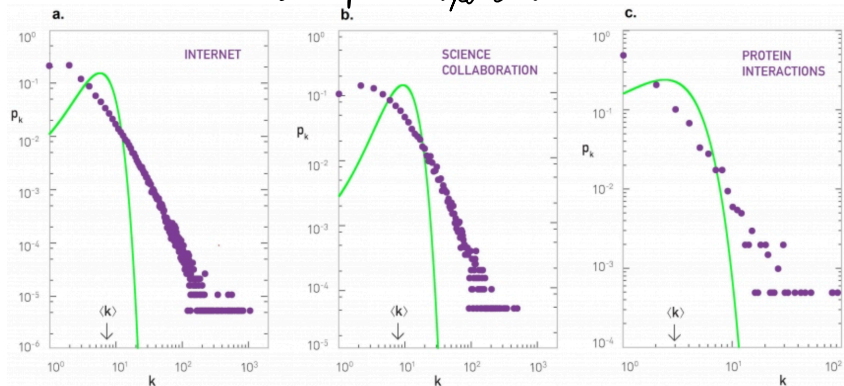
- can be very expensive computationally!
- it is embarrassingly parallel

Example

Back to $G(n, p)$

While useful, $G(n, p)$ does not resemble real world networks.

anche altre feature importanti



Random Graphs with Given Degree Sequence

Problem: given a graph G , you want to generate a random graph where each vertex has the same degree as in G .

More formally: let

- $V = \{1, \dots, n\}$
- $d_i = \deg(i)$ the degree of vertex i in G

Given: d_i for $i = 1, 2, \dots, n$

Goal: generate a random graph $G' = (V, E')$ where the degree of vertex i is d_i .

Ideas?

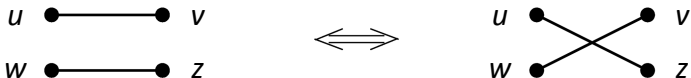


Random Graphs with Given Degree Sequence (continue)

Idea:

- start from G
- swap pairs of edges in G while preserving the degree of each vertex
- repeat the previous steps many times

Edge swap operation:



Random Graphs with Given Degree Sequence (continue)

Algorithm RandomGraphFixedDegrees(G, k)

Input: graph $G = (V, E)$; $k \in \mathbb{N}^+$

Output: random graph G' where each vertex has the same degree as in G

```
 $E' \leftarrow E$ ;  $G' = (V, E')$ ;  
for  $i \leftarrow 1$  to  $k$  do  
    sample edges  $(u, v)$  and  $(w, z)$  from  $E'$  uniformly at  
    random;  
    if  $(u, z) \notin E'$  and  $(w, v) \notin E'$  then  
         $E' \leftarrow E' \setminus \{(u, v), (w, z)\} \cup \{(u, z), (w, v)\}$ ;  
return  $G'$ ;
```

Example

Interpretation

Interpretation (continue)

Analysis

What we want: the output of `RandomGraphFixedDegrees(G, k)` is a graph *chosen uniformly at random* among *all the graphs* with the same degree sequence as G .

Equivalent: every graph with the same degree sequence as G is produced in output by `RandomGraphFixedDegrees(G, k)` with the same probability.

Question 1: can `RandomGraphFixedDegrees(G, k)` produce in output every graph with the same degree sequence as G ?

Proposition

Every graph with the same degree sequence as G can be obtained by a sequence of edge swap operation.

Proved in Ryser (1957), *Combinatorial properties of matrices of zeros and ones*, Canadian J. Math.

Question 1: can `RandomGraphFixedDegrees(G, k)` produce in output every graph with the same degree sequence as G ?

Answer: yes.

Analysis (continue)

Question 2: does every graph with the same degree sequence as G have the same probability to be produced in output by RandomGraphFixedDegrees?

Answer: yes, if k is large enough

Intuition

A Different, but Not Correct, Algorithm

Algorithm RandomGraphFixedDegrees(G, k)

Input: graph $G = (V, E)$; $k \in \mathbb{N}^+$

Output: random graph G' where each vertex has the same degree as in G

$E' \leftarrow E$; $G' = (V, E')$;

for $i \leftarrow 1$ **to** k **do**

 sample edges (u, v) and (w, z) from E , with $(u, z) \notin E$ and $(w, v) \notin E$, uniformly at random;

$E \leftarrow E \setminus \{(u, v), (w, z)\} \cup \{(u, z), (w, v)\}$;

return G' ;

Intuition

Analysis (continue)

Question 2: does every graph with the same degree sequence as G have the same probability to be produced in output by RandomGraphFixedDegrees?

Answer: yes, if k is large enough

Question 3: how large should k be?

Answer: empirically, $k \in \Omega(|E|)$ (e.g., $k = 100|E|$)

Related Random Graph: The Chung-Lu Model

Idea: instead of having the *exact* degree sequence, each node has (approximately) the same degree as in G *in expectation*

How?

Let $\text{deg}(u)$ be the degree of $u \in V$ in G , and $|E| = m$

For each pair (u, v) , let $p_{u,v} = \frac{\text{deg}(u)\text{deg}(v)}{2m}$.

Random graph G' : every edge (u, v) appears with probability $p_{u,v}$ independently of all other events.

Introduced by Chung and Lu (2002), *The average distances in random graphs with given expected degrees*, PNAS.

Analysis

Proposition

Let d_u the degree of u in random graph from the Chung-Lu model.
Then

$$\mathbb{E}[d_u] = \deg(u) \left(1 - \frac{\deg(u)}{2m} \right)$$

Notes

- $\mathbb{E}[d_u] \approx \text{deg}(u)$
- some nodes u may have degree fairly different from $\text{deg}(u)$ (e.g., low degree nodes)
- the model is (fairly) amenable to analytical results
- it is the underlying model for a commonly used heuristic algorithm to find *communities* in networks