

PERCOLATION PROCESSES ON DYNAMICALLY GROWN GRAPHS

DUKE MATHEMATICS SENIOR THESIS

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Abstract

We develop the theory of cluster growth near criticality for a class of “two-choice rules” for dynamically grown graphs. We use scaling theory to compute critical exponents for any two-choice rule, and we show special cases in which we can solve for these exponents explicitly. Finally, we compare our results with the corresponding results for the Erdős-Rényi rule, the simplest two-choice rule for which more explicit calculations are possible. We derive several of its important properties, then show that a large subset of two-choice rules - bounded size rules - behave like Erdős-Rényi near criticality.

1 INTRODUCTION

We begin with an empty graph on n vertices, and at each step we add one edge chosen at random. If each step corresponds to time $1/n$, then as $n \rightarrow \infty$ the graph at time t converges to an Erdős-Rényi graph in which the connection probability between any two edges is $2t/n$. A basic result on Erdős-Rényi graphs [7] is that there is a critical time $t_c = 1/2$ at which point a giant component with order n points emerges, which is referred to as “percolation”.

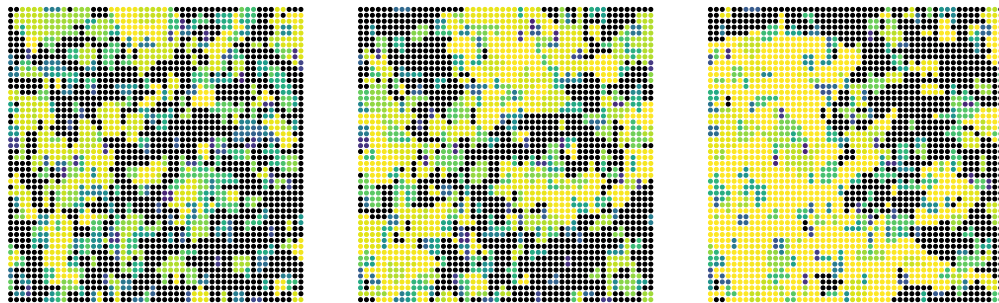


Figure 1: The above is an approximation of the Erdős-Rényi process using 2500 vertices at times $t = \frac{2}{5}, \frac{1}{2},$ and $\frac{3}{5}$, respectively. Black vertices are isolated; otherwise, vertices of the same color belong to the same cluster. Cluster color corresponds to cluster size, with dark clusters being small and light clusters large. Note that after time $\frac{1}{2}$, a single cluster has begun to dominate the graph.

The properties of the Erdős-Rényi model are well understood, so it serves as a natural reference point when studying other models. For example, if $\theta(t)$ is the fraction of vertices in the giant component (after we have let $n \rightarrow \infty$), then $\theta(t) \rightarrow 0$ as $t \searrow t_c$. This shows that the phase transition is continuous. Additionally, $\theta(t) \sim (t - t_c)^\beta$, where $\beta = 1$. This gives us a “critical exponent” β that describes the behavior of the process near t_c .

It is then natural to ask which generalizations of Erdős-Rényi graphs also have these properties. At a Fields Institute workshop in 2000, Dimitris Achlioptas suggested a class of variants of the Erdős-Rényi model, now dubbed “Achlioptas processes”. One starts with an empty graph, and at each step two edges (v_1, v_2) and (v_3, v_4) are randomly chosen from the set of all possible edges. Only one of them is added to the graph, though, according to some rule that depends on the cluster sizes of the current graph.

In 2001, Bohman and Frieze were the first to describe an Achlioptas process that delayed percolation [4]. At each step, the first edge is added if both v_1 and v_2 were isolated vertices, and the second edge is added otherwise. They showed that there was a constant $c_0 > 0.535$ such that the largest component at time c_0 has size bounded by $(\log n)^{O(1)}$. This rule is an example of a “bounded size rule”, in which all components of size greater than some constant K are treated the same.

If κ_i is the cluster size of v_i , then two possible rules are the sum rule and product rule. In the sum rule, the first edge is added if $\kappa_1 + \kappa_2 < \kappa_3 + \kappa_4$, and the second edge is added otherwise. In the product rule, the first edge is added if $\kappa_1 \kappa_2 < \kappa_3 \kappa_4$, and the second edge is added otherwise. In 2009, Achlioptas, D’Souza, and Spencer [1] used numerical experiments to claim that the product rule exhibited “explosive percolation” by causing a discontinuous phase transition at the critical time; however, in 2010, da Costa, et al. [5] found that the product rule’s phase transition was actually continuous by comparing it to a strictly more explosive rule with continuous phase transition; the da Costa rule with parameter m selected two groups of m vertices at random and joined the vertices from each group with the smallest cluster size (in this work, we will classify a rule like this as a “minimizing two-choice rule”). Their arguments relied on scaling theory, and our results in turn rely heavily on their ideas.

In 2011, Riordan and Warnke showed that results for the da Costa rule could be greatly generalized even without the use of scaling theory: the phase transition at criticality is continuous for all ℓ -vertex rules [8], which are a generalization of Achlioptas processes in which ℓ vertices are sampled independently at random and one or more edges are necessarily added between them only if each vertex is in a distinct cluster.

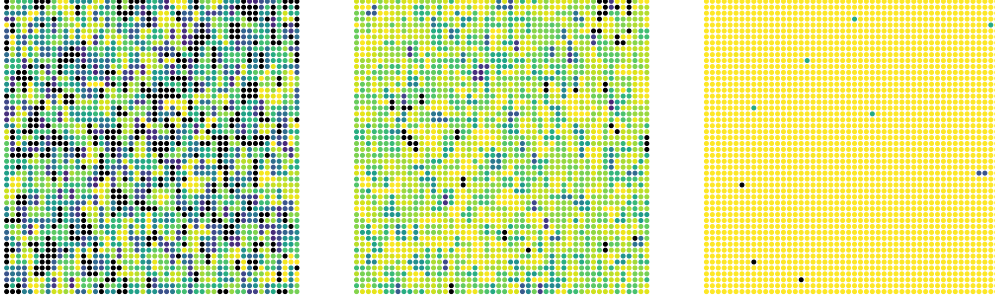


Figure 2: The above is an approximation of the da Costa rule with $m = 5$ at times $t = \frac{1}{2}$, $\frac{3}{4}$, and 1, respectively. Here we see explosive cluster growth: by time $\frac{3}{4}$, there are many large clusters cluttering the graph, and by time $t = 1$, they have all merged into a single cluster that contains all but a handful of vertices.

In 2013, Bhamidi, Budhirja, and Wang showed that the behavior of the Bohman-Frieze process near the critical point was the same as in the Erdős-Rényi case [3]. Explicitly, they showed that at time $t_c + r/n^{1/3}$ for $-\infty < r < \infty$, the system converges to the “multiplicative coalescent”, in which clusters of size x and y merge at rate xy . Aldous had already showed a corresponding result for Erdős-Rényi graphs [2], so this hinted at a deeper connection between the Bohman-Frieze and

Erdős-Rényi processes. Through far from trivial, this result should not be surprising. After some time there will be very few isolated vertices, so the process will start to add edges *almost* uniformly at random.

Then in 2017, Riordan and Warnke showed that *all* bounded size Achlioptas processes share (in a strong sense) all the features of the Erdős-Rényi phase transition [9]. One unfortunate consequence of working in such generality is that Riordan and Warnke’s work is largely based on contradiction, and thus it does not provide much quantitative information about the sizes of the graph clusters near criticality. Under the additional assumption of “scaling behavior” that is supported by numerical evidence (see da Costa et al. [6]), though, it is possible to calculate further critical exponents for these processes.

The focus of this work will then be to generalize existing methods of calculating critical exponents, determining the classes of rules for which this is possible, and studying the limiting behavior of the critical exponents as their corresponding rules’ parameters tend to extreme values.

2 TWO-CHOICE RULES

To begin, we will need to define some basic terms that will be used over and over again. Let S be the fraction of vertices in the dominating cluster as $n \rightarrow \infty$. If x_i is a vertex, then we denote its absolute cluster size by κ_i . Denote the probability that the minimum of m independently and randomly sampled vertices is s by

$$Q_m(s, t) := \mathbb{P}(\min\{\kappa_1, \dots, \kappa_m\} = s \text{ at time } t).$$

For notational simplicity we will drop the time t from the notation. Consider the sum $\langle 1 \rangle_m := \sum_s Q_m(s)$, where the sum is over $s < \infty$. Since S is the probability that a randomly chosen vertex belongs to an infinite cluster, we can write this sum as

$$\langle 1 \rangle_m = \sum_s Q_m(s) = 1 - S^m. \quad (1)$$

Since $S = 0$ for $t \leq t_c$ by definition, $\langle 1 \rangle_m = 1$ for all $t \leq t_c$. Since they frequently show up in common examples, we give $m = 1$ and $m = 2$ shorthands:

$$P := Q_1, \quad Q := Q_2.$$

We also define

$$\langle s^k \rangle_\phi := \sum_s s^k \phi(s)$$

for any real-valued function ϕ . In the special case $\phi = Q_m$, we will use the shorthand

$$\langle s^k \rangle_m := \langle s^k \rangle_{Q_m}.$$

We will use $\langle \cdot \rangle_P$ and $\langle \cdot \rangle_Q$ instead of $\langle \cdot \rangle_1$ and $\langle \cdot \rangle_2$, respectively. Finally, we will use the notation $a(t) \sim b(t)$ to denote that $a(t)$ and $b(t)$ are asymptotically similar in the following sense: if t approaches some critical value t^* (which will be clear from context) such that $a(t), b(t) \rightarrow \infty$ as $t \rightarrow t^*$, then

$$a(t) \sim b(t) \iff \frac{a(t)}{b(t)} \rightarrow C \text{ as } t \rightarrow t^*$$

for some nonzero constant C . It is straightforward to show that if $a(t) \sim b(t)$, then for any nonzero constants λ, μ , we have $a(t) \sim \lambda a(t) \sim a(t) + \mu \sim b(t)$.

With these definitions in hand, we can turn our attention to the main results. We will be discussing rules that add a single edge every $t = 1/n$ units of time, gotten by randomly sampling two groups of vertices independently from the graph, then choosing an endpoint vertex from each group. The following definition is a straightforward generalization of this idea, but we will only concern ourselves in this work with the case $\ell = 2$.

Definition 1. Define a rule \mathcal{R} as follows:

- Every $t = 1/n$ units of time, choose ℓ groups of vertices $\mathcal{V}_1, \dots, \mathcal{V}_\ell$ (of potentially different sizes) by sampling vertices independent and at random from the graph.
- For each $i \in \{1, \dots, \ell\}$, follow some rule \mathcal{F}_i to choose a vertex x_i with cluster size κ_i from group \mathcal{V}_i , subject to the condition that \mathcal{F}_i induces a function $\phi_i(s) = \mathbb{P}(\kappa_i = s)$ that does not depend on any other ϕ_j for $j \neq i$.

We call \mathcal{R} an ℓ -**choice rule**.

If $\phi_i = Q_{m_i}$ for each i , then \mathcal{R} is **minimizing**. We can similarly define **maximizing** rules. If each ϕ_i is the same, we say \mathcal{R} is **symmetric**. We would like to restrict the vertex selection processes in each group as little as possible in order to get a more general theory, but for the concrete examples mentioned above, we can obtain more detailed results.

2.1 THE SCALING ASSUMPTION AND CRITICAL EXPONENTS

Most of the results in this work follow from the assumption that near the critical time t_c , the function P has the following form

$$P(s, t) = s^{1-\tau} f(s\delta^{1/\sigma}), \quad (2)$$

where τ and σ are constants, f is a scaling function, $\delta := t - t_c$, and s is large. The following theorem gives relations between the exponents τ and $1/\sigma$ if some regularity conditions hold for the scaling function f .

Theorem 1. Suppose a rule \mathcal{R} has a scaling function f such that

1. $\lim_{x \rightarrow \infty} x^{2-\tau} f(x)$ is finite; and
2. $\int_0^\infty x^{2-\tau} f'(x) dx$ is finite.

Then there are **critical exponents**

$$\begin{aligned} \beta &= (\tau - 2)/\sigma, \\ \gamma_m &= (m(2 - \tau) + 1)/\sigma. \end{aligned}$$

such that $S \sim \delta^\beta$ and $\langle s^k \rangle_m \sim \delta^{-\gamma_m - (k-1)/\sigma}$.

Proof. We begin proving the first result. Since $S(t) = 1 - \sum_s P(s, t) = \sum_s (P(s, t_c) - P(s, t))$, we can replace the sum with integration and substitute (2) to get

$$S \approx \int_0^\infty s^{1-\tau} (f(0) - f(s\delta^{1/\sigma})) ds.$$

We make the change of variable $s = x\delta^{-1/\sigma}$ to get

$$S = \delta^{(\tau-2)/\sigma} \int_0^\infty x^{1-\tau} (f(0) - f(x)) dx.$$

Integrating by parts gives

$$S = \frac{\delta^{(\tau-2)/\sigma}}{\tau-2} \left[-x^{2-\tau} (f(0) - f(x)) \right]_{x=0}^{x=\infty} - \int_0^\infty x^{2-\tau} f'(x) dx.$$

So by our assumptions on f , we have $S \sim \delta^\beta$, where $\beta = (\tau - 2)/\sigma$. The derivation of γ_m is similar. \square

We will primarily be concerned with determining the critical exponents $1/\sigma$, τ , and γ_m . More generally, we can also consider exponents γ_ϕ for any real-valued function ϕ : γ_ϕ is the unique real number such that $\langle s \rangle_\phi \sim \delta^{-\gamma_\phi}$, if it exists. If so, we can generalize the previous theorem to extend its results to this case.

Theorem 2. *Suppose $\phi(s)$ can be written in the form $\psi(P(s))$, and suppose further that there is a γ_ϕ such that $\langle s \rangle_\phi \sim \delta^{-\gamma_\phi}$. Then*

$$\langle s^k \rangle_\phi \sim \delta^{-\gamma_\phi - (k-1)/\sigma}$$

for all k .

Proof. The integral definition of $\langle s^k \rangle_\phi$ and the scaling assumption give

$$\begin{aligned} \langle s^k \rangle_\phi &= \int_0^\infty s^k \psi(P(s)) ds \\ &= \int_0^\infty \psi(s^{1-\tau} f(s^{1/\sigma})) ds. \end{aligned}$$

We make the change of variable $s = x\delta^{1/\sigma}$ to get

$$\langle s^k \rangle_\phi = \delta^{-(k+1)/\sigma} \int_0^\infty x^k \psi(x^{1-\tau} \delta^{(1-\tau)/\sigma} f(x)) dx.$$

Let $g_k(x) := x^k \delta^{-(k+1)/\sigma} \psi(x^{1-\tau} \delta^{(1-\tau)/\sigma} f(x))$, then

$$\begin{aligned} \langle s^k \rangle_\phi &= \int_0^\infty g_k(x) dx = \delta^{-1/\sigma} \int_0^\infty x g_{k-1}(x) dx \\ &= \delta^{-(k-1)/\sigma} \int_0^\infty x^{k-1} g_1(x) dx. \end{aligned}$$

By our assumption that γ_ϕ exists, the above integral is $\sim \delta^{-\gamma_\phi}$. Thus $\langle s^k \rangle_\phi \sim \delta^{-\gamma_\phi - (k-1)/\sigma}$ for all $k > 0$. \square

As a consequence, we can determine γ_ϕ from $1/\sigma$ using any k . Thus we can always choose whichever one makes computations easiest. From now on, we will assume that all rules we work with satisfy the conditions of Theorem 2.

2.2 INDUCED ORDER MAPS

An important tool used throughout this work is the *induced order map*, which encodes how quickly the giant component is growing at criticality. We use the term *scaling form* to denote the function of δ that a particular value takes on as a result of the scaling assumption (2). For example, the scaling form of S is δ^β , a direct result of Theorem 1.

Definition 2. For ζ a real-valued function of S , the **induced order map** of ζ is the unique map $F_\zeta : \mathbb{R} \rightarrow \mathbb{R}$ such that

$$\zeta(S) \sim \delta^{F_\zeta(\beta)},$$

i.e. $F_\zeta(\beta)$ is the order of the scaling form of $\zeta(S)$. The **i -th standard induced order map** F_i of an ℓ -choice rule is the induced order map of $1 - \langle 1 \rangle_{\phi_i}$, the probability that the vertex selected by ϕ_i belongs to an infinite cluster.

Example 1. Consider the map $f(S) = S^a$ for some constant a . In scaling form, $f(S) \sim \delta^{a\beta}$, so the induced order map F_f is multiplication by a . Alternatively, consider the map $g(S) = bS$ for some constant b . In scaling form, $g(S) \sim b\delta^a$, so the induced order map F_g is the identity. This is expected, as g does not change the asymptotic behavior of the growth of S .

Example 2. Consider the two-choice rule with $\phi_1 = \phi_2 = Q_m$. Both standard induced order maps are the same: the unique map F such that $F(\beta)$ is the order of the scaling form of $1 - \langle 1 \rangle_m$. Recall the identity $1 - \langle 1 \rangle_m = S^m$, which means the scaling form is $\delta^{m\beta}$, so $F(\beta) = m\beta$. Note that $F(\beta) = \beta$ for the Erdős-Rényi rule ($m = 1$), i.e. $1 - \langle 1 \rangle_P = S \sim \delta^\beta$.

We will see later on that the quantity $1 - \langle 1 \rangle_\phi$ occurs quite naturally when describing the growth of the giant component near criticality, which motivates why we use it to define the standard induced order map. On a different note, the following theorem ties together critical exponents and induced order maps, allowing us to translate between the two whenever necessary.

Proposition 1. For any two-choice rule \mathcal{R} ,

$$\gamma_{\phi_i} = \frac{1}{\sigma} - F_i(\beta),$$

where F_i is the i -th standard induced order map of \mathcal{R} .

Proof. γ_{ϕ_i} is defined to be the real number satisfying $\langle s \rangle_\phi \sim \delta^{-\gamma_{\phi_i}}$. But by Theorem 2 and the definition of standard induced order maps, $\langle 1 \rangle_{\phi_i} \sim \delta^{-\gamma_{\phi_i} + 1/\sigma} = \delta^{F_i(\beta)}$. Thus $\gamma_{\phi_i} = 1/\sigma - F_i(\beta)$. \square

In the following sections, we will derive scaling relations for general two-choice rules in terms of standard induced order maps, then use these to analyze minimizing rules and several other examples.

2.3 SCALING RELATIONS

Suppose \mathcal{R} is a general two-choice rule, then, since the two sets are chosen independently,

$$\partial_t P(s) = s \sum_{u+v=s} \phi_1(u)\phi_2(v) - s\phi_1(s) - s\phi_2(s). \quad (3)$$

The summation term represents two components merging into a new component of size s , and the last two terms each represent a component of size s joining with another component (this also has an integral form, which we will use when it is convenient). We can use this to calculate the growth rate of the giant component, which allows us to prove an important computational theorem.

Lemma 1. For any two-choice rule \mathcal{R} , the fraction of vertices in the giant component S satisfies

$$\partial_t S = \langle s \rangle_{\phi_1} (1 - \langle 1 \rangle_{\phi_2}) + \langle s \rangle_{\phi_2} (1 - \langle 1 \rangle_{\phi_1}).$$

Proof. Using the identity $\sum_s P(s) = 1 - S$, we calculate

$$\begin{aligned}
\partial_t S &= - \sum_s \partial_t P(s) \\
&= - \sum_s s \sum_{u+v=s} \phi_1(u) \phi_2(v) + \sum_s s \phi_1(s) \sum_s + s \phi_2(s) \\
&= - \sum_u \sum_v (u+v) \phi_1(u) \phi_2(v) + \langle s \rangle_{\phi_1} + \langle s \rangle_{\phi_2} \\
&= - \sum_u u \phi_1(u) \sum_v \phi_2(v) - \sum_u \phi_1(u) \sum_v v \phi_2(v) + \langle s \rangle_{\phi_1} + \langle s \rangle_{\phi_2} \\
&= - \langle s \rangle_{\phi_1} \langle 1 \rangle_{\phi_2} - \langle 1 \rangle_{\phi_1} \langle s \rangle_{\phi_2} + \langle s \rangle_{\phi_1} + \langle s \rangle_{\phi_2} \\
&= \langle s \rangle_{\phi_1} (1 - \langle 1 \rangle_{\phi_2}) + \langle s \rangle_{\phi_2} (1 - \langle 1 \rangle_{\phi_1}).
\end{aligned}$$

□

Theorem 3. For any two-choice rule \mathcal{R} , there are nonnegative functions ζ_1 and ζ_2 such that

$$\partial_t S = \langle s \rangle_{\phi_1} \zeta_2(S) + \langle s \rangle_{\phi_2} \zeta_1(S).$$

Furthermore, the two terms above have the same order $F_1(\beta) + F_2(\beta) - 1/\sigma$ when the expression is put into scaling form, where F_i is the i -th standard induced order map of \mathcal{R} .

Proof. By Lemma 1, we know $\partial_t S = \langle s \rangle_{\phi_1} (1 - \langle 1 \rangle_{\phi_2}) + \langle s \rangle_{\phi_2} (1 - \langle 1 \rangle_{\phi_1})$. Now for both i , consider the probability $\zeta_i(S)$ that a vertex chosen from group i belongs to an infinite cluster. This satisfies the relation $\langle 1 \rangle_{\phi_i} = 1 - \zeta_i(S)$, which allows us to recover the desired form of $\partial_t S$. Then by Proposition 1, both terms have order $F_1(\beta) + F_2(\beta) - 1/\sigma$. □

To determine scaling relations for two-choice rules, we will need one last result, a consequence of Theorem 3.

Corollary 1. For any two-choice rule \mathcal{R} , the average cluster size $\langle s \rangle_P$ satisfies

$$\partial_t \langle s \rangle_P = 2 \langle s \rangle_{\phi_1} \langle s \rangle_{\phi_2} - \langle s^2 \rangle_{\phi_1} \zeta_2(S) - \langle s^2 \rangle_{\phi_2} \zeta_1(S),$$

where ζ_1 and ζ_2 are the functions from Theorem 3.

Proof. Recall from the proof of Theorem 3 that there are functions ζ_1 and ζ_2 satisfying the relation $\langle 1 \rangle_{\phi_i} = 1 - \zeta_i(S)$. We can then explicitly compute $\partial_t \langle s \rangle_P$.

$$\begin{aligned}
\partial_t \langle s \rangle_P &= \sum_s s \partial_t P(s) \\
&= \sum_s s^2 \sum_{u+v=s} \phi_1(u) \phi_2(v) - \sum_s s^2 \phi_1(s) - \sum_s s^2 \phi_2(s) \\
&= \sum_u \sum_v (u+v)^2 \phi_1(u) \phi_2(v) - \langle s^2 \rangle_{\phi_1} - \langle s^2 \rangle_{\phi_2} \\
&= \langle s^2 \rangle_{\phi_1} \langle 1 \rangle_{\phi_2} + 2 \langle s \rangle_{\phi_1} \langle s \rangle_{\phi_2} + \langle 1 \rangle_{\phi_1} \langle s^2 \rangle_{\phi_2} - \langle s^2 \rangle_{\phi_1} - \langle s^2 \rangle_{\phi_2} \\
&= 2 \langle s \rangle_{\phi_1} \langle s \rangle_{\phi_2} + \langle s^2 \rangle_{\phi_1} (\langle 1 \rangle_{\phi_2} - 1) + \langle s^2 \rangle_{\phi_2} (\langle 1 \rangle_{\phi_1} - 1) \\
&= 2 \langle s \rangle_{\phi_1} \langle s \rangle_{\phi_2} - \langle s^2 \rangle_{\phi_1} \zeta_2(S) - \langle s^2 \rangle_{\phi_2} \zeta_1(S).
\end{aligned}$$

□

Theorem 4. *For any two-choice rule with standard induced order maps F_i ,*

$$\begin{aligned}\gamma_{\phi_1} &= F_2(\beta) - \beta + 1, \\ \gamma_{\phi_2} &= F_1(\beta) - \beta + 1, \\ \gamma_P &= F_1(\beta) + F_2(\beta) - 2\beta + 1, \\ \frac{1}{\sigma} &= F_1(\beta) + F_2(\beta) - \beta + 1, \\ \tau &= \frac{\beta}{F_1(\beta) + F_2(\beta) - \beta + 1} + 2.\end{aligned}$$

Proof. Since $\partial_t S \sim \delta^{\beta-1}$, Theorem 3 shows

$$\frac{1}{\sigma} = F_1(\beta) + F_2(\beta) - \beta + 1.$$

Since $\partial_t S = \langle s \rangle_{\phi_1} \zeta_2(S) + \langle s \rangle_{\phi_2} \zeta_1(S)$, the two dominating terms near criticality give us the system

$$\beta - 1 = -\gamma_{\phi_1} + F_2(\beta) = -\gamma_{\phi_2} + F_1(\beta).$$

This system implies

$$\begin{aligned}\gamma_{\phi_1} &= F_2(\beta) - \beta + 1, \\ \gamma_{\phi_2} &= F_1(\beta) - \beta + 1.\end{aligned}$$

By Corollary 1, we know $\partial_t \langle s \rangle_P = 2\langle s \rangle_{\phi_1} \langle s \rangle_{\phi_2} - \langle s^2 \rangle_{\phi_1} \zeta_2(S) - \langle s^2 \rangle_{\phi_2} \zeta_1(S)$. The three terms in this expression all have order $F_1(\beta) + F_2(\beta) - 2/\sigma$, which gives the system

$$-\gamma_P - 1 = -\gamma_{\phi_1} - \gamma_{\phi_2} = -\gamma_{\phi_1} - \frac{1}{\sigma} + F_2(\beta) = -\gamma_{\phi_2} - \frac{1}{\sigma} + F_1(\beta).$$

Using our previously calculated expressions for γ_{ϕ_1} and γ_{ϕ_2} this system gives us

$$\gamma_P = F_1(\beta) + F_2(\beta) - 2\beta + 1.$$

Finally, using the identity $\beta = (\tau - 2)/\sigma$ from Theorem 1, we get

$$\tau = \frac{\beta}{F_1(\beta) + F_2(\beta) - \beta + 1} + 2.$$

□

Suppose we are instead working with a minimizing two-choice rule, i.e. $\phi_1 = Q_a$ and $\phi_2 = Q_b$ for two positive integers a and b . In this case, we have simpler forms for the scaling relations.

Corollary 2. *For minimizing two-choice rules, the scaling relations from Theorem 4 become*

$$\begin{aligned}\gamma_a &= 1 + (b - 1)\beta, \\ \gamma_b &= 1 + (a - 1)\beta, \\ \gamma_P &= 1 + (a + b - 2)\beta, \\ \frac{1}{\sigma} &= 1 + (a + b - 1)\beta, \\ \tau &= \frac{\beta}{1 + (a + b - 1)\beta} + 2.\end{aligned}$$

Proof. The relation $\langle 1 \rangle_m = 1 - S^m$ holds for all m , so the induced order map for Q_m is $\beta \mapsto m\beta$. The result then follows from Theorem 4. \square

Example 3. Da Costa et al. [6] studied the two-choice rule with $\phi_1 = \phi_2 = Q_m$ for some m . In this case, Corollary 2 gives critical exponents

$$\begin{aligned}\gamma_m &= 1 + (m-1)\beta, \\ \gamma_P &= 1 + (2m-2)\beta, \\ \frac{1}{\sigma} &= 1 + (2m-1)\beta, \\ \tau &= \frac{\beta}{1 + (2m-1)\beta} + 2,\end{aligned}$$

which agree with those derived in [6].

2.4 LIMITING BEHAVIOR

One might expect that as a and b grow, we can delay the emergence of the giant component since we have more control over which vertices are connected. The following proposition makes this notion precise for minimizing two-choice rules in terms of induced order maps.

Proposition 2. *Suppose \mathcal{R} is a minimizing two-choice rule with standard induced order maps $F_1(\beta) = a\beta$ and $F_2(\beta) = b\beta$. If $a\beta \rightarrow 0$ and $b\beta \rightarrow 0$ as $a, b \rightarrow \infty$, then the critical exponents for \mathcal{R} have limits*

$$\gamma_{\phi_1} = \gamma_{\phi_2} = \gamma_P = \frac{1}{\sigma} = 1, \quad \tau = 2.$$

Proof. If $a\beta, b\beta \rightarrow 0$, then $\beta \rightarrow 0$ necessarily. This is then a straightforward computation using Corollary 2. \square

Of course, not all rules will satisfy this condition, as not all rules attempt to delay the critical time in the first place. It is an interesting question to determine broad families of rules for which $F_i(\beta) \rightarrow 0$ for both i as $a, b \rightarrow \infty$. We conjecture that this applies for all minimizing two-choice rules.

Conjecture 1. *Suppose \mathcal{R} is a minimizing two-choice rule with group sizes a and b . Then $a\beta \rightarrow 0$ and $b\beta \rightarrow 0$ as $a, b \rightarrow \infty$, i.e. the limits in Proposition 2 apply.*

2.5 CLUSTER SIZE VARIANCE

The variance of the cluster size (at a fixed time) is, by definition, $\text{Var}_i(s) = \langle s^2 \rangle_{\phi_i} - \langle s \rangle_{\phi_i}^2$. For two-choice rules, we can use our scaling relations to put this solely in terms of β and the induced order maps, at least near criticality.

$$\begin{aligned}\text{Var}_1(s) &= \delta^{-[\gamma_{\phi_1} + 1/\sigma]} - \delta^{-2\gamma_{\phi_1}} \\ &= \delta^{-[F_1(\beta) + 2F_2(\beta) - 2\beta + 2]} + \delta^{-[2F_2(\beta) - 2\beta + 2]}.\end{aligned}$$

Similarly,

$$\text{Var}_2(s) = \delta^{-[2F_1(\beta) + F_2(\beta) - 2\beta + 2]} + \delta^{-[2F_1(\beta) - 2\beta + 2]}.$$

Then if $F_i(\beta) \geq 0$ for both i , this is dominated by the first term. So near t_c , we have

$$\begin{aligned}\text{Var}_1(s) &\approx \delta^{-[F_1(\beta)+2F_2(\beta)-2\beta+2]}, \\ \text{Var}_2(s) &\approx \delta^{-[2F_1(\beta)+F_2(\beta)-2\beta+2]}.\end{aligned}$$

If $F_i(\beta) \rightarrow 0$ for both i , then we get the asymptotic behavior $\text{Var}_i(s) \rightarrow \delta^{-2}$. If $F_i(\beta) \leq 0$ for either i , then a similar argument with more cases gives the same limit.

Example 4. If \mathcal{R} is a minimizing two-choice rule, then $F_1(\beta) = a\beta$ and $F_2(\beta) = b\beta$. Then

$$\begin{aligned}\text{Var}_1(s) &= \delta^{(2-a-2b)\beta-2}, \\ \text{Var}_2(s) &= \delta^{(2-2a-b)\beta-2}.\end{aligned}$$

If Conjecture 1 is true, then $\text{Var}_i(s) \rightarrow \delta^{-2}$ for both i .

We can consider what happens in the limit for certain rules. If $F_i(\beta) \rightarrow 0$ as for both i as $a, b \rightarrow \infty$, then by the proof of Proposition 2, $\beta \rightarrow 0$ as well. Then a simple computation yields

$$\text{Var}_i(s) = \delta^{-2}$$

for both i . Thus if Conjecture 1 holds, this is true for minimizing two-choice rules.

If a rule is simple enough, we need not consider a limiting case of it to determine the variance of its cluster size distribution. For example, we know the Erdős-Rényi rule has $\beta = 1$ (see Section 3.1) and $F_i(\beta) = \beta = 1$ for both i . Thus for this simple rule, we get $\text{Var}_i(s) = \delta^{-3}$ for both i .

3 ERDŐS-RÉNYI

The Erdős-Rényi rule is the simplest of all the two-choice rules. For each group, simply pick one vertex at random to be the group representative. Thus an equivalent way of defining the Erdős-Rényi rule is to add one of the $\binom{n}{2}$ possible edges in the graph at random at each time step.

This rule also has strong connections with the usual Erdős-Rényi random graphs. In particular, we can view the graph created by the Erdős-Rényi rule at time t/n (so t edges have been added) as an Erdős-Rényi graph with connection probability $\frac{2t}{n(n-1)}$. Thus in addition to our scaling theory, we can apply the theory of Erdős-Rényi random graphs to study this rule. We will denote Erdős-Rényi random graphs on n vertices and with connection probability p_c as $G(n, p_c)$.

Before deriving our first results, we should note that the Erdős-Rényi rule is both minimizing and symmetric, so plugging in $a = b = 1$ into the scaling relations from Corollary 2 gives

$$\gamma_P = 1, \quad \frac{1}{\sigma} = 1 + \beta, \quad \tau = \frac{\beta}{1 + \beta} + 2. \quad (4)$$

With such a simple rule, we can in fact determine much more. In particular, we can determine β , which in turn fixes the other critical exponents. We will show $\beta = 1$ (see Theorem 5 below), which subsequently implies $\sigma = 1/2$ and $\tau = 5/2$.

3.1 FINDING β FOR ERDŐS-RÉNYI

Lemma 2. *For the Erdős-Rényi rule, the critical time is $1/2$.*

Proof. This is a result of translating between the Erdős-Rényi rule and Erdős-Rényi random graphs. It is well known [10] that percolation in Erdős-Rényi random graphs occurs when the connection probability is $1/(n-1)$, which corresponds to using the Erdős-Rényi rule up until time $1/2$ by our earlier discussion. \square

Note that since the critical time for the Erdős-Rényi rule is $t_c = t/n = 1/2$, this corresponds to adding $t = n/2$ edges to the graph. To find β then, we can apply the theory of branching processes to study how connected the static graphs $G(n, t/n)$ are for $t < n/2$, as branching processes arise quite naturally in them.

Theorem 5. *For the Erdős-Rényi rule, $\beta = 1$.*

Proof. Suppose $t < n/2$, i.e. we are in the subcritical regime, and consider the Erdős-Rényi random graph $G(n, t/n)$. This has connection probability proportional to $G(n, 2t/(n(n-1)))$ by a factor of $2/(n-1)$, so as $n \rightarrow \infty$, this corresponds to a slowed down version of the Erdős-Rényi rule with critical time $t_c = 1$.

Let v be a randomly chosen vertex and let Z_m be the number of vertices at distance m from v . When $m = o(\log n)$, the cluster containing x is a tree with high probability, which makes Z_m into a branching process where each individual has a $\text{Poisson}(t)$ number of offspring.

The generating function for this distribution is

$$\begin{aligned}\phi(z) &= \sum_{k=0}^{\infty} r_k z^k \\ &= \sum_{k=0}^{\infty} \frac{t^k e^{-t}}{k!} z^k \\ &= e^{-t(1-z)}.\end{aligned}$$

We care about this in the limit as $t \rightarrow n/2$ and as $n \rightarrow \infty$, so we can suppose $t > 1$. A standard result on branching processes is when $t > 1$, the probability that the process dies out is given by the smallest $z \in [0, 1]$ such that $z = \phi(z)$. This means the probability that the process *never* dies out is $\theta = 1 - z$.

Letting $\mu := \phi''(1)$ and $z = 1 - \theta$ and expanding $\phi(1 - \theta)$ in power series near 1 gives

$$\begin{aligned}z &= \phi(z) \\ 1 - \theta &= \phi(1 - \theta) \\ 1 - \theta &\approx 1 - t\theta + \frac{\mu\theta^2}{2} \\ \theta &\approx \frac{2}{\mu}(t - 1).\end{aligned}$$

Thus $\theta \sim t - 1$. But in this slowed down version of Erdős-Rényi, the critical time is 1, so this is in the form $\theta \sim \delta$. By definition, β is the exponent such that $\theta \sim \delta^\beta$, so $\beta = 1$. \square

3.2 THE SCALING WINDOW

A critical assumption of this work was the scaling assumption, and for Erdős-Rényi, we can find an explicit bound on the size of the *scaling window*, the region where the scaling assumption is accurate.

Theorem 6. *For Erdős-Rényi, the scaling assumption holds for clusters of size s after t_c when $\delta = \Theta(s^{-1/2})$.*

Proof. We use the following two relationships, which [6] derives for symmetric minimizing two-choice rules when $t < t_c$ and s is large.

- $P(x) = \delta^{(\tau-1)/\sigma} \tilde{f}(x\delta^{1/\sigma})$;
- $\tilde{f}(x) \propto x^\lambda \exp(-Cx^{1+\log_2 m})$, where $\lambda = (1 + \log_2 m) \left(1 + \frac{1}{4m-2}\right) - \frac{2m}{2m-1}$;

where m is the group size. In the case of Erdős-Rényi, we know $m = 1$, $\tau = 5/2$, and $1/\sigma = 2$, so these reduce to

$$P(x) = \tilde{C}\delta^2 x^{-1/2} \exp(-Cx\delta^2),$$

where \tilde{C} is the constant of proportionality from \tilde{f} . To clean up notation, we use the shorthand $\mathcal{E}_x := \exp(-Cx\delta^2)$. Also note that $\delta = t - t_c$ when $t > t_c$, so $\partial_t \delta = 1$. With these observations, we can begin the computation. When s is large, Equation (3) becomes

$$\begin{aligned} \partial_t P(s) &= \frac{s}{2} \int_0^s P(u)P(s-u) du - sP(s) \\ \partial_t \left\{ \tilde{C}\delta^2 s^{-1/2} \mathcal{E}_s \right\} &= \frac{s}{2} \int_0^s \tilde{C}^2 \delta^4 (su - u^2)^{-1/2} \mathcal{E}_u \mathcal{E}_{s-u} du - \tilde{C}\delta^2 s^{1/2} \mathcal{E}_s \\ \tilde{C}\delta \mathcal{E}_s \left[2s^{-1/2} - 2C\delta^2 s^{1/2} \right] &= \tilde{C}\delta \mathcal{E}_s \left[\frac{s}{2} \int_0^s \tilde{C}\delta^3 (su - u^2)^{-1/2} du - \delta s^{1/2} \right] \\ 2s^{-1/2} - 2C\delta^2 s^{1/2} &= \frac{s}{2} \tilde{C}\delta^3 \int_0^s (su - u^2)^{-1/2} du - \delta s^{1/2}. \end{aligned}$$

Consider the integral above. It is straightforward to check that $2 \sin^{-1}(\sqrt{u/s})$ is an antiderivative of $\frac{1}{\sqrt{su-u^2}}$ when s is positive, so this integral evaluates to

$$\begin{aligned} \int_0^s \frac{1}{\sqrt{su-u^2}} du &= 2 (\sin^{-1}(1) - \sin^{-1}(0)) \\ &= \pi. \end{aligned}$$

Thus our earlier equation becomes

$$2s^{-1/2} - 2C\delta^2 s^{1/2} = \frac{1}{2} \tilde{C}\delta^3 \pi s - \delta s^{1/2}.$$

Note that the first term on the left hand side goes to zero as $s \rightarrow \infty$, so in the limit this becomes

$$\left(\frac{1}{2} \tilde{C}\pi s \right) \delta^2 + \left(2Cs^{1/2} \right) \delta - s^{1/2} = 0.$$

This quadratic has roots

$$\delta = \frac{-2Cs^{1/2} \pm \sqrt{4C^2 s + 2\tilde{C}\pi s^{3/2}}}{\tilde{C}\pi s}.$$

The only valid root is the positive one since $\delta > 0$ in the postcritical regime. This root is $\Theta(s^{-1/2})$ when s is large. \square

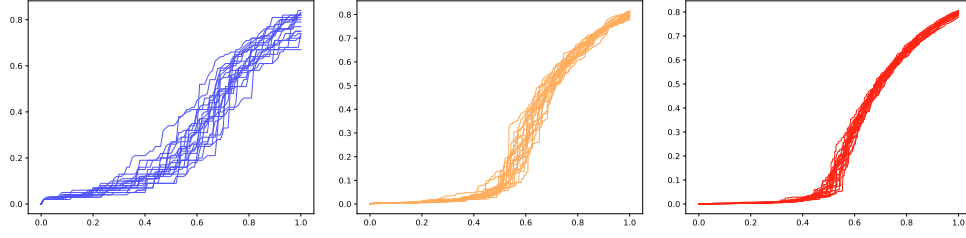


Figure 3: The above plots show the fraction of vertices in the giant component for 20 Erdős-Rényi processes with $n = 100, 1000$, and 2500 vertices, respectively. The fact that $\beta = 1$, i.e. $\theta \sim \delta$, means that the slopes of these lines at $t = 1/2$ become linear in the limit as $n \rightarrow \infty$. Furthermore, Theorem 6 implies that as $n \rightarrow \infty$, the linear portion of the plot becomes smaller, which is somewhat evident from the above plots.

3.3 BOUNDED SIZE RULES

Since the Erdős-Rényi rule is so well understood, it is useful to know how any arbitrary rule relates to it. In particular, we can focus on rules that behave exactly like Erdős-Rényi, as characterized by critical exponents. These rules might have different subcritical and supercritical behavior, but as the giant component emerges, the rule produces similar graphs as Erdős-Rényi.

Bounded size rules are a very general category of rule that eventually start acting like Erdős-Rényi in the following sense: fix a constant $K \in \mathbb{N}$, then \mathcal{R} is a bounded size rule if it treats all sampled vertices with $\kappa_i > K$ identically. In fact, [9] shows that bounded size rules share all the key features of Erdős-Rényi at criticality in a certain precise sense.

Intuitively, it makes sense that a bounded size rule eventually become similar to Erdős-Rényi. After enough clusters have been joined, the number of clusters of size $\leq K$ will be so small that it will not have much of an effect on the growth of the giant component. The key to noticing the complexity here is actually twofold:

1. For any $k \leq K$, the number of clusters of size k can be *nonzero* at percolation (see Proposition 3 below);
2. After all clusters of size $\leq K$ have stopped having a real effect on the giant component, we are in a position much different than that of Erdős-Rényi. Instead of lots of isolated nodes, we have lots of finite size clusters.

We can use a simple bounded size two-choice rule to examine these problems, which we do using a generalization of the well known Bohman-Frieze rule:

1. At each step, pick m vertices. These are the group 1 vertices.
2. If any of the m vertices are isolated, pick the first such one as the group 1 representative. If not, sample one more random vertex and pick it no matter what.
3. Repeat this for group 2, and connect the 2 group representatives with an edge.

For this particular rule, we can explicitly track how the number of isolated clusters is changing through time; we will see that a nonzero proportion of our graph will be made up of isolated nodes at criticality.

Proposition 3. *For the Bohman-Frieze rule with $m = 1$, a nonzero proportion of the vertices in the graph are isolated at percolation.*

Proof. We can explicitly write out the probability $\phi(s)$ that a representative's cluster size is s :

$$\phi(s) = \begin{cases} 1 - (1 - P_1)^{m+1} & s = 1, \\ (1 - P_1)^m P_s & s > 1. \end{cases}$$

Let $m = 1$, then Equation (3) gives

$$\begin{aligned} \partial_t P_1 &= -2\phi(1) \\ &= -2 + 2(1 - P_1)^{m+1} \\ &= -2 + 2(1 - P_1)^2 \\ &= 2P_1^2 - 4P_1. \end{aligned}$$

Solving this differential equation and using the initial condition $P_1(0) = 1$ gives us the solution

$$P_1(t) = \frac{2}{e^{4t} + 1}.$$

We know that percolation occurs at $t = 1$ at the very latest by [8]. But plugging in $t = 1$ gives $P_1(1) = \frac{2}{e^4 + 1} \approx 0.024$, so at criticality, a strictly positive proportion of our nodes are still isolated. \square

As it turns out, this does not significantly affect cluster growth for bounded size rules, which we can formalize in terms of critical exponents (compare the following theorem with (4)). This is a known result [9], but using our theory of induced order maps, the proof is computational instead of by contradiction, as in [9].

Theorem 7. *Every bounded size two-choice rule has critical exponents*

$$\gamma_P = 1, \quad \frac{1}{\sigma} = 1 + \beta, \quad \tau = \frac{\beta}{1 + \beta} + 2.$$

Proof. Suppose we have a bounded size two-choice rule that treats all clusters of size $> K$ the same. Without loss of generality, suppose this rule is also symmetric with $\phi(s) = \mathbb{P}(\kappa = s)$ (this argument really only applies to the representative selection process for one of the two vertex groups, so it readily extends to the case of non-symmetric rules). Finally, suppose each vertex group contains m vertices.

We will also need a bit of new notation. Let $N \in \{0, \dots, m\}$ denote the number of vertices in our group whose cluster size is $\leq K$, and let $P_{\leq K} = P_1 + \dots + P_K$. Then for $s > K$, we can partition $\phi(s)$ and use the fact that it acts exactly like Erdős-Rényi if all m of the vertices in the group come from clusters of size $> K$.

$$\begin{aligned} \phi(s) &= \mathbb{P}(\kappa = s \mid N = 0) \mathbb{P}(N = 0) + \mathbb{P}(\kappa = s \mid N > 0) \mathbb{P}(N > 0) \\ &= \frac{P_s}{S + \sum_{r>K} P_r} (1 - P_{\leq K})^m + \mathbb{P}(\kappa = s \mid N > 0) \mathbb{P}(N > 0). \end{aligned}$$

But $1 - P_{\leq K} = S + \sum_{r>K} P_r$, so this becomes

$$\phi(s) = P_s (1 - P_{\leq K})^{m-1} + \mathbb{P}(\kappa = s \mid N > 0) \mathbb{P}(N > 0).$$

For $s \leq K$, we can use this same partition to get a similar expression.

$$\begin{aligned} \phi(s) &= \mathbb{P}(\kappa = s \mid N = 0) \mathbb{P}(N = 0) + \mathbb{P}(\kappa = s \mid N > 0) \mathbb{P}(N > 0) \\ &= 0 + \mathbb{P}(\kappa = s \mid N > 0) \mathbb{P}(N > 0). \end{aligned}$$

This allows us to sum over all $\phi(s)$ and thus calculate $\langle 1 \rangle_\phi$.

$$\begin{aligned}
\sum_s \phi(s) &= \sum_{s>K} \phi(s) + \sum_{s \leq K} \phi(s) \\
&= \left[\sum_{s>K} (1 - P_{\leq K})^{m-1} \right] + \left[\sum_s \mathbb{P}(\kappa = s \mid N > 0) \mathbb{P}(N > 0) \right] \\
&= \left[(1 - P_{\leq K})^{m-1} (1 - S - P_{\leq K}) \right] + \left[(1 - (1 - P_{\leq K})^m) \sum_s \mathbb{P}(\kappa = s \mid N > 0) \right] \\
&= (1 - P_{\leq K})^{m-1} [1 - S - \xi + (\xi - 1)P_{\leq K}] + \xi,
\end{aligned}$$

where $\xi = \sum_s \mathbb{P}(\kappa = s \mid N > 0)$. In scaling form, the first term above will be dominated by the *smallest* critical exponent since δ is approaching 0. After multiplying everything out, the term with the lowest exponent will be S (with critical exponent β).

We will now induct on the group size m to show that $\xi \sim 1$. When $m = 1$,

$$\sum_s \mathbb{P}(\kappa = s \mid N > 0) = \sum_s \mathbb{P}(\kappa = s \mid N = 1) = 1.$$

Then assuming $\sum_s \mathbb{P}(\kappa = s \mid N > 0) \sim 1$ when the group size is $m - 1$,

$$\begin{aligned}
\sum_s \mathbb{P}(\kappa = s \mid N > 0) &= \sum_s \mathbb{P}(\kappa = s \mid 0 < N < m) \mathbb{P}(0 < N < m \mid N > 0) \\
&\quad + \sum_s \mathbb{P}(\kappa = s \mid N = m) \mathbb{P}(N = m \mid N > 0) \\
&= \mathbb{P}(0 < N < m \mid N > 0) \sum_s \mathbb{P}(\kappa = s \mid 0 < N < m) \\
&\quad + \mathbb{P}(N = m \mid N > 0) \sum_s \mathbb{P}(\kappa = s \mid N = m).
\end{aligned}$$

The first sum above is ~ 1 by our inductive hypothesis, and the second sum is exactly equal to 1. Thus

$$\sum_s \mathbb{P}(\kappa = s \mid N > 0) \sim \mathbb{P}(0 < N < m \mid N > 0) + \mathbb{P}(N = m \mid N > 0) = 1,$$

as desired. Thus the scaling form of $\langle 1 \rangle_\phi = \sum_s \phi(s)$ is

$$\langle 1 \rangle_\phi \sim \delta^\beta.$$

By definition, this fixes the standard induced order map of this rule. Note that for Erdős-Rényi, $\langle 1 \rangle_P = S \sim \delta^\beta$ as well, so both rules share the same standard induced order map. Since our computation of the critical exponents of two-choice rules relied only this map, the desired result follows. \square

A crucial relation used in the previous theorem was the fact that

$$\xi = \sum_s \mathbb{P}(\kappa = s \mid N > 0)$$

is dominated by other terms as we approach criticality. For certain bounded size two-choice rules, more can be said of this term. For example, the Bohman-Frieze rule is a bounded size rule with threshold $K = 1$, and $\xi_{BF} = 1$ since it always prioritizes connecting isolated vertices. It is already intuitively clear that the Bohman-Frieze rule cannot prioritize isolated clusters (more generally, the clusters below its size threshold) any further, but this classification gives that statement rigorous meaning.

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