

## Bootstrap percolation on $G(n, p)$

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Bootstrap percolation on a graph  $G$  is defined as the spread of *activation* or *infection* according to the following rule, with a given threshold  $r \geq 2$ : We start with a set  $\mathcal{A}(0) \subseteq V(G)$  of *active* vertices. Each inactive vertex that has at least  $r$  active neighbours becomes active. This is repeated until no more vertices become active, i.e., when no inactive vertex has  $r$  or more active neighbours.

We are mainly interested in the final size  $A^*$  of the active set, and in particular whether eventually all vertices will be active or not. If they are, we say that the initial set  $\mathcal{A}(0)$  *percolates*. We will study a sequence of graphs of order  $n \rightarrow \infty$ ; we then also say that (a sequence of)  $\mathcal{A}(0)$  *almost percolates* if the number of vertices that remain inactive is  $o(n)$ , i.e., if  $A^* = n - o(n)$ .

Bootstrap percolation has been studied on various graphs, both deterministic and random; one can study either a random initial set or the deterministic problem of choosing an initial set that is optimal in some sense. For example, a classical folklore problem is to find the minimal percolating set in a two-dimensional grid; see Balogh and Pete [3] and Bollobas [5]. (These references also treat higher-dimensional grids.) Some further references for random initial sets on various graphs are Cerf and Manzo [6], Holroyd [7] (grids); Balogh and Bollobás [1] (hypercube); Balogh, Peres and Pete [2] (infinite trees); Balogh and Pittel [4] (random regular graphs).

We here study bootstrap percolation on the Erdős-Rényi random graph  $G_{n,p}$  (which somewhat surprisingly seems to have been neglected so far in this context), with an initial set  $\mathcal{A}(0)$  consisting of a given number  $a$  vertices chosen at random. This was first studied by Vallier [8]; we here present a simple method that allows us to both simplify the proofs and improve the results.

In order to analyze the bootstrap percolation process on  $G_{n,p}$ , we change the time scale and consider at each time step the spread of activation from one vertex only. Choose  $u_1 \in \mathcal{A}(0)$  and give each of its neighbours a *mark*; we then say that  $u_1$  is *used*, and let  $\mathcal{Z}(1) := \{u_1\}$  be the set of used vertices at time 1. We continue recursively: At time  $t + 1$ , choose a vertex  $u_{t+1} \in \mathcal{A}(t) \setminus \mathcal{Z}(t)$  (provided this set is non-empty). We give each neighbour of  $u_{t+1}$  a new mark. Let  $\Delta\mathcal{A}(t+1)$  be the set of inactive vertices with  $r$  marks; these now become active and we let  $\mathcal{A}(t+1) = \mathcal{A}(t) \cup \Delta\mathcal{A}(t+1)$  be the set of active vertices at time  $t$ . We finally set  $\mathcal{Z}(t+1) = \mathcal{Z}(t) \cup \{u_{t+1}\} = \{u_i : i \leq t+1\}$ , the set of used vertices.

The process stops when  $\mathcal{A}(t) \setminus \mathcal{Z}(t) = \emptyset$ , i.e., when all active vertices are used. We denote this time by  $T$ ;

$$(1) \quad T := \min\{t \geq 0 : \mathcal{A}(t) \setminus \mathcal{Z}(t) = \emptyset\}.$$

Thus the final infected set is  $\mathcal{A}(T) = \mathcal{Z}(T)$ , and its size is

$$(2) \quad A^* := |\mathcal{A}(T)| = |\mathcal{Z}(T)| = T.$$

Hence, the set  $\mathcal{A}(0)$  percolates if and only if  $T = n$ , and  $\mathcal{A}(0)$  almost percolates if and only if  $T = n - o(n)$ .

Since  $|\mathcal{Z}(t)| = t$  and  $\mathcal{Z}(t) \subseteq \mathcal{A}(t)$  for  $t = 0, \dots, T$ , we also have, with  $A(t) := |\mathcal{A}(t)|$ , the number of active vertices at time  $t$ ,

$$(3) \quad T = \min\{t \geq 0 : A(t) = t\}.$$

We analyze this process by the standard method of revealing the edges of the graph  $G_{n,p}$  only on a need-to-know basis. We thus begin by choosing  $u_1$  as above and then reveal its neighbours; we then find  $u_2$  and reveal its neighbours, and so on. Let, for  $i \notin \mathcal{Z}(s)$ ,  $I_i(s)$  be the indicator function that there is an edge between the vertices  $u_s$  and  $i$ . This is also the indicator that  $i$  gets a mark at time  $s$ , so if  $M_i(t)$  is the number of marks  $i$  has at time  $t$ , then

$$(4) \quad M_i(t) = \sum_{s=1}^t I_i(s),$$

at least until  $i$  is activated (and what happens later does not matter). Note that if  $i \notin \mathcal{A}(0)$ , then, for every  $t \leq T$ ,  $i \in \mathcal{A}(t)$  if and only if  $M_i(t) \geq r$ .

The crucial feature of this description of the process, which makes the analysis simple, is that the random variables  $I_i(s)$  are i.i.d.  $\text{Be}(p)$ . We have defined  $I_i(s)$  only for  $s \leq T$  and  $i \notin \mathcal{Z}(s)$ , but it is convenient to add further (redundant) variables so that  $I_i(s)$  are defined, and i.i.d., for all  $i \in V_n$  and all  $s \geq 1$ .

Define, for  $i \in V_n \setminus \mathcal{A}(0)$ ,

$$(5) \quad Y_i := \min\{t : M_i(t) \geq r\}.$$

If  $Y_i \leq T$ , then  $Y_i$  is the time vertex  $i$  becomes active, but if  $Y_i > T$ , then  $Y_i$  never becomes active. Thus, for  $t \leq T$ ,

$$\mathcal{A}(t) = \mathcal{A}(0) \cup \{i \notin \mathcal{A}(0) : Y_i \leq t\}.$$

By (4) and (5), each  $Y_i$  has a negative binomial distribution  $\text{NegBin}(r, p)$ ;

$$\mathbb{P}(Y_i = k) = \mathbb{P}(M_i(k-1) = r-1, I_i(k) = 1) = \binom{k-1}{r-1} p^k (1-p)^{r-k};$$

moreover, these random variables  $Y_i$  are i.i.d.

We let, for  $t = 0, 1, 2, \dots$ ,

$$S(t) := |\{i \notin \mathcal{A}(0) : Y_i \leq t\}|,$$

so

$$(6) \quad A(t) = S(t) + A(0) = S(t) + a.$$

By (3), (2) and (6), it suffices to study the stochastic process  $S(t)$ . Note that  $S(t)$  is a sum of  $n - a$  i.i.d. processes  $\mathbf{1}[t \geq Y_i]$ , each of which is 0/1-valued and jumps from 0 to 1 at time  $Y_i$ . The fact that  $S(t)$ , and thus  $A(t)$ , is a sum of i.i.d. processes makes the analysis easy; in particular, for any given  $t$ ,

$$S(t) \sim \text{Bin}(n - a, \mathbb{P}(Y_1 \leq t)).$$

We have, for any given  $t_0$ ,

$$T \geq t_0 \iff \min_{t < t_0} (A(t) - t) > 0 \iff a + \min_{t < t_0} (S(t) - t) > 0 \iff a > -\min_{t < t_0} (S(t) - t).$$

(Note that this is exact; so far no approximation has been done.)

To find the threshold for (almost) percolation, we thus only have to find the minimum  $\min_{t < t_0} (S(t) - t)$  for  $t_0 = n$  or  $t_0$  close to  $n$ . Standard concentration results show that  $S(t) \approx \mathbb{E} S(t)$ , where

$$\mathbb{E} S(t) = (n - a) \mathbb{P}(Y_1 \leq t) = (n - a) \mathbb{P}(M_1(t) \geq r),$$

and explicit results are easily found.

For notational simplicity we state the result for  $r = 2$  only. In this case,  $\mathbb{E} S(t) - t$  has a minimum  $1/(2np^2)$  at  $t = 1/(np^2)$  (asymptotically), and we obtain the following result.

**Theorem 1.** *Let  $r = 2$ , and assume  $n^{-1} \ll p = p(n) \ll n^{-1/2}$ . Then the threshold for (almost) percolation is*

$$a_* := \frac{1}{2np^2}.$$

More precisely, for any fixed  $\delta > 0$ ,

- (i) *If  $|\mathcal{A}(0)| \leq (1 - \epsilon)a_*$ , then whp  $A^* \leq 2|\mathcal{A}(0)|$ .*
- (ii) *If  $|\mathcal{A}(0)| \geq (1 + \epsilon)a_*$ , then whp  $A^* = n - o(n)$ . If further  $np \geq \log n + \log \log n + \omega(n)$  for some  $\omega(n) \rightarrow \infty$ , then whp  $A^* = n$ , so  $\mathcal{A}(0)$  percolates completely.*

Moreover,  $S(t) - \mathbb{E} S(t)$  converges after normalization to a Gaussian process, and it is easy to refine the results above and obtain very precise information on the width of the critical window (which is of the order  $\sqrt{a_*}$ ); we also obtain a Gaussian limit law for the final size  $A^*$  in the subcritical case.

Details will appear.

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