

Project 24: Entanglement Percolation

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1 An analytical starter

Task leader(s): Vicentini Gioele

1.1 Introduction: Some concepts to understand the paper

Ideas in the paper are developed in the generating function formalism: $G(x) = \sum_{k=0}^{+\infty} x^k P(k)$. The probability of finding an edge which does not connect to the GCC is u, generated by $H_1(x)$.

As seen in class, by definition this probability can be obtained in the following way: for a generic edge exiting from a node with degree k, there is a probability $1 - p_2$ for it not to be there and a probability $p_2H_1(x)^{k'}q(k')$ that the edge exists and that the neighbors of the connected node are not in the GCC. q(k) is the excess degree distribution $q(k) = kP(k)/\langle k \rangle$ which gives us the probability of having a neighbor of degree k'. Summing over all possible values of k' gives:

$$H_1(x) = (1 - p_2) + \sum_{k'} p_2 H_1(x)^{k'} q(k')$$
(1.1)

$$= (1 - p_2) + \sum_{k'} p_2 H_1(x)^{k'} k' P(k') / \langle k \rangle$$
 (1.2)

$$= (1 - p_2) + p_2 x G_1(H_1(x))$$
(1.3)

where we used the sencond order gen. function $xG_1(x) = \sum_k kP(k) = xG_0'(x)/G_0'(1)$.

This leaves us with $u = H_1(1) = (1 - p_2) + p_2 x G_1(u)$, the usual equation in open form. For the probability of finding a node not connected to the GCC, we proceed similarly: this node has degree k with prob. P(k) and it's not connected to the GCC if all its edges are not: $H_0(1) = \sum_k P(k) H_1(1)^k = G_0(H_1(1))$ and $H_0(x) = x G_0(H_1(x))$. This means that the probability of a node to be in the GCC is $S = 1 - H_0(1)$.

1.2 The contribution of q-swapped cycles

The previous formulas change after the q-swap operations since they change in a relevant way the topology. Assuming node degrees to be uncorrelated:

$$\tilde{H}_1(x) = 1 - p_2 + p_2 \tilde{H}_1(x) P_1(1) + \sum_{q \ge 2} \left[p_2 (1 - \Pi_q) \tilde{H}_1(x)^q P_1(q) + \Pi_q P_1(q) C_q(x) \right]$$
(1.4)

which accounts for (from left to right):

- the probability $1 p_2$ of not having an edge
- the prob. of having an edge to a neighbor of degree 1, for which there is no possible q-swap

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• the prob. (summed over degrees q of the existing neighbor) of not applying q-swap and the neighbor is not in GCC

• the prob. of a neighbor over which we apply the q-swapping operation. $C_q(x)$ generates the prob. of not being in the GCC for the new q-swapped cycle.

In particular, since each edge of the cycle is occupied with prob. p (SCP probability):

$$C_q(x) = qp^{q-1}(1-p)xG_1(\tilde{H}_1(x))^q + p^q xG_1(\tilde{H}_1(x))^q \sum_{l=0}^{q-2} \left[(l+1)p^l xG_1(\tilde{H}_1(x))^q (1-p)^{2-l} \right]$$
(1.5)

note that if an edge connects us to a q-swapped cycle and two or more edges of this cycle are not occupied (SCP probability being p < 1), we have to account in different ways for the cases in which some of the edges are missing (otherwise we count the probability of, for example, having some nodes which do not belong to the GCC but not having a connection to them). if l is the number of occupied edges, this means that l = q has only one disposition of the edges, l = q - 1 has q, l = 0 up to l = q - 2 has l + 1, since l nodes must be connected for us to count properly the probability of not being connected to the GCC. The shape of $C_q(x)$ is connected to this. In the paper they write $(l+1)p^l(1-p)^2 \dots$ instead of the binomial result $(l+1)p^l(1-p)^{2-l} \dots$, but i'm almost sure it's a typo on their side.

For $\tilde{H}_0(x)$, instead, denoting with μ_q the probability for a node of degree q of undergoing q-swap:

$$\tilde{H}_0(x) = P(0) + \tilde{H}_1(x)P(1) + \sum_{q>2} \left[\Pi_q \mu_q P(q) + (1 - \Pi_q \mu_q) P(q) \tilde{H}_1(x)^q \right]$$
(1.6)

which are again the probabilities for degrees 0, 1 which cannot be q-swapped, the prob. of having a q-swap which sets the degree of the node to zero (thus, not in GCC), the prob. of not having a q-swap and not being connected to GCC. In the paper they split the sum in order to obtain $xG_0(H_1(x))$.

2 | Simulations over ER and BA networks, plus a failed attempt on SW

Task leader(s): Vicentini Gioele

For all the networks the strategy is the same: first generate two copies of the network, then apply q-swap of degrees 2,3 to one of the two copies, to then finally study their behavior under bond percolation. The results for ER and BA networks are in chapter 2. BA networks are famous for being robust under bond percolation, so while the effect of 2,3 swaps is still present, it's less relevant. This proves how much the underlying topology of the network is relevant when designing processes apt to enchance specific qualities of the network.

In Figure 2 there's a failed attempt of replicating Figure 3.b of the paper, where percolation and q-swappings are applied on a Small World network with added short-cuts. Due to the limited computational power available, the network studied here is relevantly small, $N \sim 10^3$, which means that the added edges are less than what studied in the paper $(N \sim 10^6)$, although the addition probability is higher (0.5 against 0.25).

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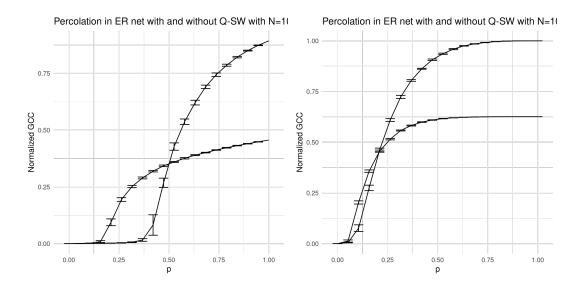


Figure 2.1: Results of simulations of robustness against bond percolation. Left: ER network of 10^4 nodes with avg. degree 2.5. Right: BA network of 10^4 nodes with 3 edges added per timestep. As expected, BA networks are more robust under bond percolation and the 2, 3 swaps are not so relevant.

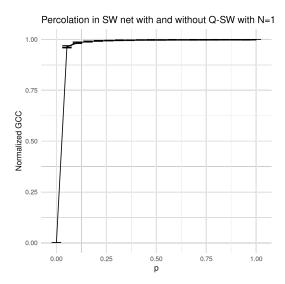


Figure 2.2: Bond percolation after 2, 3 swap on a Watts-Strogats Small World model with 10^3 nodes and shortcut addition probability $\phi = 0.5$. As expected of such highly redundant networks, neither percolation nor q-swaps have relevant effects.