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# Quantum communication capacity transition of complex quantum networks

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# Quantum communication capacity transition of complex quantum networks

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Quantum network is the key to enable distributed quantum information processing. As the single-link communication rate decays exponentially with the distance, to enable reliable end-to-end quantum communication, the number of nodes need to grow with the network scale. For highly connected networks, we identify a threshold transition in the capacity as the density of network nodes increases—below a critical density, the rate is almost zero, while above the threshold the rate increases linearly with the density. Surprisingly, above the threshold the typical communication capacity between two nodes is independent of the distance between them, due to multi-path routing enabled by the quantum network. In contrast, for less connected networks such as scale-free networks, the end-to-end capacity saturates to constants as the number of nodes increase, and always decays with the distance. Our results are based on capacity evaluations, therefore the minimum density requirement for an appreciable capacity applies to any general protocols of quantum networks.

### I. INTRODUCTION

Quantum information (QI) science has brought advantages in various applications [1–4]. To unleash the full power of QI processing in distributed tasks [5, 6], a quantum network (QN) [7–11] aiming at entanglement distribution and QI transmission is the key.

The Internet is mainly built upon fiber networks, with photons as the information carrier. Similarly, photons as the only known "flying qubits" will likely be the information carrier in a QN. In both cases, channel loss is the major challenge to communication. Therefore, networking protocols that make use of intermediate nodes or repeaters are important for both. Unlike classical information, quantum information cannot be simply cloned and amplified, and therefore increasing the number of nodes, even repeater nodes [12–22], are costly. In this regard, a key question for designing a QN is to understand the trade-off between the density of nodes and the entanglement distribution rate: how many nodes are necessary to guarantee reliable QI transmission between multiple users in a fixed region?

The answer not only depends on the overall distances between the users, but also on the topology of the QN to be built [23]. As it is likely that well-developed classical fiber networks can be adopted as the base of QNs, Ref. [24] developed a model for QN on the the probabilistic transmission of single photons and took a classical network science approach to study its connectivity by the giant component. However, for QNs exploiting quantum technologies such as quantum error correction [25] and non-classical state generation [26, 27], the semi-classical approach has a limited implication. In particular, Ref. [24]'s critical density highly depends on the repetition of each channel uses and thus blurs the essential constraints. More recently, Refs. [28, 29] considered effects from repeater nodes. As the results rely on specific protocols, the fundamental limits of the tradeoff remains unclear. We address the same question with a full quantum information approach based on the fundamental limits [30–34], and obtain a minimum density requirement that generally applies to any protocols.

As the exact architecture and protocols of QNs are unclear, we take the information-theoretical approach and evaluate the end-to-end capacity [33] of QI transmission. To account for different possibilities of the future QNs, we consider typical types of network models [40], based on the Waxman networks [36, 37], Erdős-Rényi model and scale-free networks [38, 39]. Our results provide an upper bound to characterize the quantum capacity of QNs and the analysis applies to all kinds of quantum communication. In Waxman and Erdős-Rényi QNs, the ensemble-averaged capacity abruptly transits from almost zero to nonzero values at a critical density of nodes. Above the threshold, it grows with the density linearly, at a rate depending on the statistical properties of the QN. Surprisingly, in this region the end-to-end capacity typically does not depend on the distance between the two end nodes, due to the multi-path routing enabled by the coordination of the entire QN. In scale-free QNs, the ensemble-averaged capacity saturates to a constant depending on the scale of the network as the density of nodes increases, due to the limited connectivity in the network that prevents efficient multi-path routing.

#### **II. MODEL OF QNS**

The skeleton of a QN can be described by a graph, with vertices  $\mathcal{G}$  being the network nodes and edges  $\mathcal{E}$ representing the transmission links [41]. As nodes are located geographically, we can assign a 2-D coordinate

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 $\boldsymbol{x}$  to each node. The transmission link along each edge  $E_{\boldsymbol{x},\boldsymbol{x}'}$  is modeled as a bosonic pure loss channel, with a transmissivity  $\eta(\boldsymbol{x},\boldsymbol{x}') = 10^{-\gamma D(\boldsymbol{x},\boldsymbol{x}')}$  for fiber length  $D(\boldsymbol{x},\boldsymbol{x}')$  at a state-of-the-art rate  $\gamma = 0.02$  per kilometer (km). For simplicity, we assume that for each edge, the fiber length  $D(\boldsymbol{x},\boldsymbol{x}')$  and the geographical distance  $\|\boldsymbol{x} - \boldsymbol{x}'\|_2$  are identical.

With the transmission links on each edge defined, one needs to specify the graph structure—the coordinates and connections of the vertices—to specify the QN. Without loss of generality, we choose the coordinates  $\boldsymbol{x}$  of the N nodes uniformly random in a square  $\Omega_R \equiv [-R, R] \times [-R, R]$ , with an area of  $|\Omega_R| = 4R^2$ .

In the random Waxman model [36, 37], each pair of nodes is connected with a probability  $\Pi(x, x') =$  $e^{-D(\boldsymbol{x},\boldsymbol{x}')/\alpha L}$  decaying exponentially with the distance. Here  $L = 2\sqrt{2}R$  is the maximum possible distance in a square; the constant  $\alpha$  controls the typical fiber length and is fixed so that  $\alpha L = 226$ km to model the U.S. fiber-optics networks [37]. It is worthy to point out that Ref. [24] adopted the same Waxman QNs. In the scale-free model [39], the network is built up dynamically: when each node x is being added, it is connected to *m* nodes out of all the previous added nodes. The probability of node x' being connected to node x is proportional to the current degree  $D_q(\mathbf{x}')$  and inversely proportional to the distance  $D(\boldsymbol{x}, \boldsymbol{x}')$ , i.e.,  $\Pi(\boldsymbol{x}, \boldsymbol{x}') \propto$  $D_{q}(\boldsymbol{x})/D(\boldsymbol{x},\boldsymbol{x}')$ , in contrast to the Waxman model's exponential decay with distance.

To obtain a direct impression, we visualize the two models in Fig. 2 (a) and Fig. 4 (a) respectively. Immediate differences in the connectivity can be seen, e.g. by comparing Fig. 2 (a2) and Fig. 4 (a1): for the same N = 1585 nodes in a region of scale  $R \simeq 800$  km, the Waxman model is much more connected and homogeneous, while the scale-free model is less connected and heterogeneous. These differences can be captured by their statistical properties. As shown in Fig. 1(a)(b), the Waxman QN model has a Poisson degree distribution and the average degree grows with the number of nodes linearly [45]; while the scale-free QN model has a long-tailed power-law degree distribution and a bounded average of 2m. It is also worthy mentioning that the Waxman model has a percolation phase transition (see Appendix A), where the percentage of the giant component of the graph increases sharply from close to zero to unity as the density  $\rho = N/|\Omega_R|$  increases above a critical value of  $\rho_G \simeq 7 \times 10^{-6}$ . However, we show that this necessary condition is far from being sufficient.

While we base our QN models on the Internet, a QN will be majorly different from Internet. In particular, classical repeaters [42] are not counted as network nodes in the study of Internet [40], as they are universally deployed and cheap. In contrast, quantum repeaters are nontrivial and therefore directly considered as network nodes in this study. In this regard, the Waxman model's exponential decay of long direct links will be more likely for QNs. However, our goal is *not* to determine which

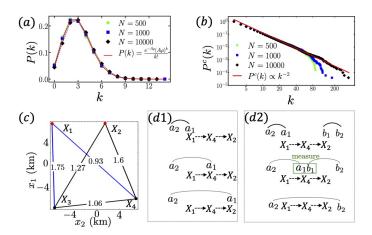


Figure 1. (a) The degree distribution of Waxman network, with density  $\rho = 10^{-5}$ , fits well with a Poisson distribution (red curve) with  $A = 3.0 \times 10^5$ . (b) The cumulative degree distribution of the scale-free model, with density  $\rho = 10^{-5}$ , fits well with a power-law (red curve). (c) A fournode QN, with the axes as the geographical coordinates. The blue color indicates a cut between  $X_1$  and  $X_2$  and the number on each edge equals the edge capacity in Eq. (1). For the cut indicated by blue edges, the cut capacity of Eq. (2)  $\mathcal{C}(\mathbb{U}_{x,x'}) = 1.75 \pm 0.93 = 2.68$ , which turns out to be the minimum cut. (d1-d2) Examples of entanglement distribution protocols. (d1) Direct communication strategy with potential error correction. (d2) Entanglement-swap strategy. After an entanglement swap measurement on  $a_1b_1$ , nodes  $X_1$  and  $X_2$ can share an entangled state in  $a_2b_2$ .

model can better represent a QN, an emerging technology, but to characterize each model in terms of quantum communications.

# III. PROTOCOLS AND CAPACITY FORMULA

To distribute entanglement between two nodes  $X_1$  and  $X_2$  in a QN, the nodes can transmit quantum states between all links and perform two-way classical communication in combination of local operations at each node. To begin with, let's consider an instance of a four-node network in Fig. 1(c). In a single-path routing strategy, one can choose a path from  $X_1$  to  $X_2$  (e.g.  $X_1 - X_4 - X_2$ ,  $X_1 - X_3 - X_2$ , or  $X_1 - X_3 - X_4 - X_2$ ) and utilize all the channels along the path once to distribute the entanglement. With the path fixed, one can either perform direct communication or adopt entanglement swap [43], as shown in Fig. 1(d). A more efficient approach is to adopt multi-path routing. For example, nodes  $X_1$  and  $X_2$  in Fig. 1(c) can utilize multiple non-overlapping paths simultaneously  $(X_1 - X_3 - X_2 \text{ and } X_1 - X_4 - X_2)$  and achieve a better performance.

As protocols vary, to obtain universal results, we consider the ultimate achievable entanglement distribution rate among *all* protocols [32, 33]. In contrast to classi-

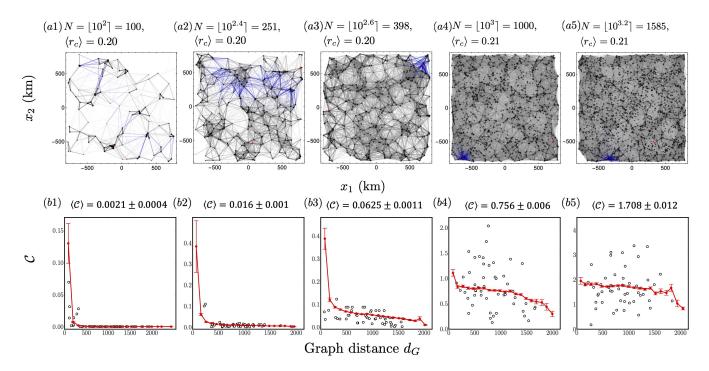


Figure 2. Waxman QNs,  $\alpha = 0.1$  ( $R \simeq 800$  km). (a1)-(a5) Visualizations with different number of nodes N. The darkness and opacity of the color of the nodes and edges indicate the relative amplitude of the capacity (darker means larger). The blue edges are the minimum cut solution to random pairs of end nodes indicated by the red dots. (b1)-(b5) The end-to-end capacity  $\mathcal{C}(\boldsymbol{x}, \boldsymbol{x}')$  between random nodes  $\boldsymbol{x}, \boldsymbol{x}'$  for QNs with fixed  $N, \alpha$ . The x-axis is the graph distance between two nodes  $d_G(\boldsymbol{x}, \boldsymbol{x}')$ , in terms of the shortest fiber path. The scattered circles are 50 random pairs in a single QN sample and the red dashed lines indicate the average obtained from 5000 random data.

cal communication [2, 44], QI transmission rate for each edge is fundamentally limited by the channel loss to be

$$\mathcal{C}_E\left(E_{\boldsymbol{x},\boldsymbol{x}'}\right) = -\log_2\left(1-\eta\right) = -\log_2\left(1-10^{-\gamma D(\boldsymbol{x},\boldsymbol{x}')}\right),\tag{1}$$

regardless of the energy, where  $\eta = 10^{-\gamma D(\boldsymbol{x}, \boldsymbol{x}')}$  is the channel loss [32]. To characterize the importance of a single node, we define the node capacity  $C_N(\boldsymbol{x}) = \sum_{\boldsymbol{x}' \in \mathcal{N}(\boldsymbol{x})} C_E(E_{\boldsymbol{x}, \boldsymbol{x}'})$ , as the sum of the edge capacities.

Consider the graph with edge capacities  $\{C_E(E_{\boldsymbol{x},\boldsymbol{x}'})\}$  as the weights (e.g. Fig. 1(c)), the problem of solving the end-to-end capacity is reduced to solving the minimum cut [33]. Let's first introduce a cut  $\mathbb{U}_{\boldsymbol{x},\boldsymbol{x}'}$  between two nodes  $\boldsymbol{x}$  and  $\boldsymbol{x}'$  as the set of edges such that their deletion will disconnect the two nodes. For example, in Fig. 1(c), the blue part indicates a cut for A and B. Then the capacity between end nodes  $\boldsymbol{x}$  and  $\boldsymbol{x}'$  is given by the "edge connectivity" between them [33]

$$\mathcal{C}(\boldsymbol{x}, \boldsymbol{x}') = \min_{\mathbb{U}_{\boldsymbol{x}, \boldsymbol{x}'}} \mathcal{C}_{U}(\mathbb{U}_{\boldsymbol{x}, \boldsymbol{x}'}) \equiv \min_{\mathbb{U}_{\boldsymbol{x}, \boldsymbol{x}'}} \sum_{E_{\boldsymbol{y}, \boldsymbol{y}'} \in \mathbb{U}_{\boldsymbol{x}, \boldsymbol{x}'}} \mathcal{C}_{E}(E_{\boldsymbol{y}, \boldsymbol{y}'}).$$
(2)

To obtain further insights, we derive an upper bound of the end-to-end capacity by the node capacities of the two end nodes,  $C(\boldsymbol{x}, \boldsymbol{x}') \leq \min \{C_N(\boldsymbol{x}), C_N(\boldsymbol{x}')\}$ , as one can always choose the cut that consists of all edges connected to one of the end nodes.

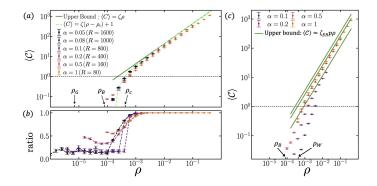


Figure 3. (a) Average end-to-end capacity  $\langle \mathcal{C} \rangle$  vs. nodes density  $\rho$ . The dark green solid line and the light green dashed line give the upper bounds  $\langle \mathcal{C} (\boldsymbol{x}) \rangle \simeq \zeta \rho$  and its shifted fitting  $\langle \mathcal{C} \rangle = \zeta(\rho - \rho_c) + 1$  respectively. The arrows indicate critical densities for the birth of giant connected component  $(\rho_G)$ , for the prediction of Ref. [24]  $(\rho_B)$  and  $\rho_c \simeq 4.25 \times 10^{-4}$  is when  $\langle \mathcal{C} \rangle = 1$ . (b) The average of the ratio of end-node edges inside the minimum cut. It shares the same legend as in (a). (c) Average end-to-end capacity  $\langle \mathcal{C} \rangle$  of Erdős-Rényi model vs node density  $\rho$ . The green lines from top to bottom correspond to the asymptotic upper bound  $\langle \mathcal{C} \rangle = \zeta_{ER} \rho \rho$  for  $\alpha = 1, 0.5, 0.2$ .

We take a statistical approach and evaluate the average end-to-end capacity  $\langle \mathcal{C}(\boldsymbol{x}, \boldsymbol{x}') \rangle$  in an ensemble of network models, where the average is over the choices of the end

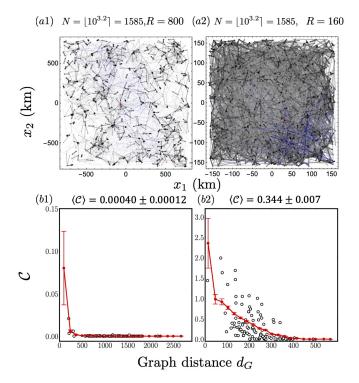


Figure 4. Scale-free QNs with N = 1585 nodes, with similar arrangements and ensemble-averaging to Fig. 2. (a1)-(a2) Visualizations of the QN model with different scales R = 800, 160 km, with fixed N = 1585 nodes. (b1)-(b2) The corresponding end-to-end capacity.

nodes  $\boldsymbol{x}, \, \boldsymbol{x}'$  and the random realization of the network, with fixed numbers of nodes N and scale R. In this regard,

$$\langle \mathcal{C} (\boldsymbol{x}, \boldsymbol{x}') \rangle \leq \langle \min \{ \mathcal{C}_N (\boldsymbol{x}), \mathcal{C}_N (\boldsymbol{x}') \} \rangle \leq \langle \mathcal{C}_N (\boldsymbol{x}) \rangle, \quad (3)$$

upper bounded by the ensemble-averaged node capacity. Compared to the edge connectivity approach based on probabilistic single-photon transmission in Ref. [24], our quantum capacity approach applies to all protocols and reveals essential features of a network.

# IV. RATE TRANSITION OF WAXMAN QNS

To study Waxman QNs, we first fix the scale  $R \simeq 800$  km and vary the number of nodes N. In Fig. 2 (b), we plot the end-to-end capacity  $C(\boldsymbol{x}, \boldsymbol{x}')$  of random pairs vs. the graph distance  $d_G(\boldsymbol{x}, \boldsymbol{x}')$  (the shortest path length) between them. When the number of nodes is small (e.g. Fig. 2 (b1)), the capacity decays with the graph distance drastically; while surprisingly, when the number of nodes becomes larger (e.g. Fig. 2 (b2)), the capacity is almost independent of the graph distance [35]. This is due to the effect of multi-path routing—the number of possible paths increases significantly with distance when the nodes are dense.

To systematically evaluate the transition in the endto-end capacity, we evaluate the ensemble-averaged capacity  $\langle \mathcal{C}(\boldsymbol{x}, \boldsymbol{x}') \rangle$  for different values of R and N. We expect the density of nodes  $\rho$  to be the crucial parameter. Indeed, we can show that when R is large, the ensemble-averaged node capacity  $\langle \mathcal{C}_N(\boldsymbol{x}) \rangle \simeq \zeta \rho$ , as the upper bound in Ineq. (3), is linear in density  $\rho$  with the coefficient  $\zeta \simeq 4358$  (see Appendix C).

In Fig. 3 (a), we plot the average capacity vs the node density  $\rho$  for different system size R. Overall, for a fixed density  $\rho$ , the capacity  $\langle C \rangle$  converges as the scale Rincreases. When the density is small, the capacity is mostly close to zero (see Appendix B); As the density increases, we see a sudden transition from almost zero capacity to o(1) capacity at a critical density. The transition happens at around  $\langle C \rangle \sim 1$  corresponding to a density  $\rho_c \simeq 4.25 \times 10^{-4}$ , which is much larger than the giant component transition  $\rho_G \simeq 7 \times 10^{-6}$  and the result  $\rho_B \simeq 6.82 \times 10^{-5}$  from Ref. [24].

After this transition, the average capacity increases linearly with node density  $\rho$ , approaching the upper bound  $\zeta\rho$  (dark green line). The reason of the convergence can be observed from Fig. 2 (a): when the connectivity is high, the minimum cut (blue edges) becomes a cut formed by all the edges connecting to one of the end points. To be more quantitative, we calculate the ratio of the edges in the minimum cut that contain at least one end node. As shown in Fig. 3 (b), the ratio transits from close to zero to unity at the same time as the end-to-end capacities approach the upper bounds. In fact, we find that a shifted upper bound  $\zeta(\rho - \rho_c) + 1$  fits the overall numerical results well, as shown by the green dashed line in Fig. 3(a).

Note that Ref. [24]'s critical density depends on the protocol parameters—e.g. the number of repetition  $n_p$  for each link; therefore the value of their critical density is not an essential characterization of the QN. Their results have to obey the constraint in our paper, as any protocol has its rate bounded by the capacity. We can confirm as follows: as they consider  $n_p = 1000$  repeated use of each channel to successfully establish one single Bell pair, the end-to-end capacity per channel use in their protocol is merely  $10^{-3}$  for density  $\rho = \rho_B$ , which is in fact within the vanishing capacity region in our results.

## V. RATE SATURATION OF SCALE-FREE QNS

Now we switch the focus to scale-free QNs (see Fig. 4). Similarly, we evaluate the end-to-end capacity for the same set of choices of R and N. In Fig. 5 (a), the ensemble-averaged capacity  $\langle \mathcal{C}(\boldsymbol{x}, \boldsymbol{x}') \rangle$  grows as N increases and saturates to a constant dependent on the scale R of the network. This is due to the limited degree of scale-free networks, which constrains the upper bound of the node capacity to be bounded by a constant  $\propto m$  and dependent on R (see Appendix E). As we can see in Fig. 5 (b), the ratio of edges of end points being

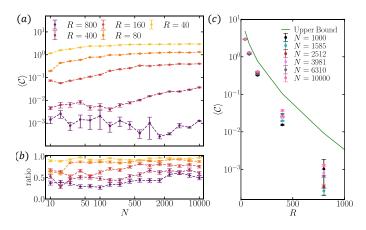


Figure 5. Scale-free model. (a) Average end-to-end capacity  $\langle \mathcal{C} \rangle$  vs. the number of nodes N for various scales R's. (b) The average of the ratio of end-node edges inside the minimum cut. (c) Capacity vs. the scale of the QN. The orange curve is the upper bound  $\langle \mathcal{C} (\boldsymbol{x}) \rangle$  from numerical integration (see Appendix E).

in the minimum cut is now determined by the network scale, and gets close to unity when the network is small. Indeed, in Fig. 5 (c) we see the gap between the saturated capacity and the upper bound from node capacity is small for small R, while larger with R increasing. Overall, the capacity decays with R exponentially, even when the number of nodes is large.

In additional to the saturation of capacity, the graphdistance-independence of the capacity is absent for scalefree QNs. In Fig. 4 (b), regardless of the capacity being large or small, there is a sharp decrease of the end-to-end capacity as the graph distance increases, in contrast to the Waxman QNs in Fig. 2 (b). This is due to the lack of multi-path routing, constrained by the connectivity of the scale-free networks. Indeed, we can find the average clustering coefficient  $\langle r_c \rangle$  decaying with the system size, instead of saturating to constants with number of nodes in the Waxman case (see Appendix A).

## VI. RATE TRANSITION OF ERDŐS-RÉNYI QNS

We also extend our analyses to the Erdős-Rényi model, a network model with uniform edge connection probability p. To compare with the Waxman model, we match the number of edges in Erdős-Rényi model to the Waxman model with same  $\alpha$  and N, via choosing a proper p. The corresponding degree distribution is binomial (see Appendix A).

We evaluate the transition of average end-to-end capacity  $\langle \mathcal{C} \rangle$  with node density in Fig. 3(c), and identify similar trend to the Waxman model: when  $\rho$  is large,  $\langle \mathcal{C} \rangle$  grows linearly with  $\rho$ ; while when  $\rho$  is small, there is still a sharp decrease in the capacity. While in the Waxman model, the capacity  $\langle \mathcal{C} \rangle$  agrees among differin Eq. (3).  $\langle C_N(\boldsymbol{x}) \rangle \simeq \zeta_{ER} p \rho$ , where  $\zeta_{ER} \simeq 5137.9$  (see Appendix D). We can directly see the dependence of  $\langle C \rangle$  on connection probability p from the asymptotic upper bound and we show them in Fig. 3(c).

#### VII. CONCLUSION AND DISCUSSIONS

In this paper, we examine the end-to-end quantum communication capacity in Waxman, Erdős-Rényi QNs and scale-free QNs. Our results provide guidance on the design of QN infrastructure, as the capacity places an achievable upper bound on rates of quantum communication protocols.

In particular, our results suggest that when the connectivity of the QN is high (like in the Waxman case), multi-path routing will enable reliable quantum communication. On the practical side, considering that quantum repeaters might be as costly and expensive as user nodes, this indicates that at a moderate metropolitan scale where users are dense and direct links are possible, it might be better to simply build more direct links between the users and utilize the multi-path routing for reliable quantum communication.

Our results are based on network capacity results and therefore reveals essential property of a QN, independent on the protocol. We reveal more detailed properties of QNs, other than the simple connectivity properties in Ref. [24]. Our results address the entanglement generation capacity, which is the most relevant quantity in a QN. In particular, our results allow unlimited two-way classical communication (via an underlying classical network) as assistance in the entanglement generation process. Ref. [24] limits the protocols to be at a single photon level, and is strongly dependent on the specific protocol parameters to generate entanglement. The density of nodes to guarantee reliable communication would depend on the exact meaning of reliable communication, however, a network above the threshold we identified is preferable as the capacity starts to become distanceindependent.

#### ACKNOWLEDGMENTS

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#### Appendix A: Basic properties networks.

As shown in Fig. 6, in the Waxman model, the average degree  $\langle k \rangle$  of the nodes increases with the number of nodes N linearly, at a rate depending on the scale  $\alpha$ ; In the scale-free network, the average degree saturates to  $\langle k \rangle = (2N - 1 - m)m/N \simeq 2m$  as the number of nodes N increases. Here m is the number of edges brought by the addition of each single node.

We plot the degree distribution of Erdős Rényi model and mean degree in Fig. 7.

The Waxman model has a giant component transition as the density of nodes  $\rho$  increases. As shown in Fig. 8(a), the ratio of the size of the largest connected component  $N_G$  over the total number of nodes N increases from close to zero to unity abruptly at a density of  $\rho_G \simeq 7 \times 10^{-6}$ . The transition becomes sharper as the number of nodes increase.

To understand the connectivity of the networks, we plot the clustering coefficient's dependence network parameters. For a single node, the single-node local clustering coefficient  $r_c(\boldsymbol{x}) = t/[k(k-1)/2]$  identifies the existence of connections between its k neighbors  $\mathcal{N}(\boldsymbol{x})$ . Here t is number of triangles that is attached to the node  $\boldsymbol{x}$ . We can define the graph clustering coefficient  $\langle r_c \rangle$  by averaging over all nodes. For Waxman networks,  $\langle r_c \rangle$  con-

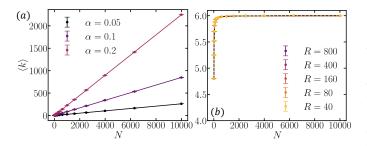


Figure 6. Average degree of (a) Waxman and (b) Yook models and its dependence on the number of nodes N. (a) Solid lines gives linear fitting results of  $\langle k \rangle = A\rho$  where  $\rho$  is the density of nodes. (b) Dashed lines show the theory curve  $\langle k \rangle = (2N - 1 - m)m/N$ .

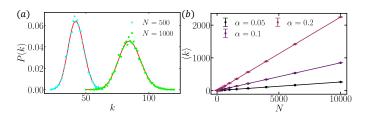


Figure 7. Degree distribution and mean degree of Erdős Rényi model. (a) Degree distribution of Erdős Rényi model with  $\alpha = 0.1$ . The red ccurves represent the analytical expression for it. (b) Mean degree of Erdős Rényi model with different  $\alpha$ .

verging to a constant dependent on  $\alpha$  as the number of nodes increases, as shown in Fig. 8(b). While for the scale-free networks,  $\langle r_c \rangle$  decays to zero as the number of nodes N increases, as shown in Fig. 8(c).

# Appendix B: Additional data for the end-to-end capacity

We provide additional data of the numerical calculations. First, we show the distribution of the end-to-end capacity between random pairs of nodes in each ensemble of networks. Fig. 9 shows the Waxman case, corresponding to Fig. 2(b1)-(b5); while Fig. 10 shows the scale-free case, corresponding to Fig. 4(b1)(b2) of the main paper. The average of the data utilized here gives the red curves in the corresponding plots of the main paper, which are also shown as red curves in these plots.

Next, we present an in-depth analyses of Fig. 3 in the main paper. Fig. 11(a) shows each curve of capacity vs. number of nodes for different scales individually, without collapsing everything in plotting with density. In the main paper, we do not show the long tails, as these tails are mainly due to rare cases of random pairs of nodes lying very close to each other. Indeed, if we plot the median instead of the mean, as shown in Fig. 12, these long tails are not present and we see a clear sharp drop. To avoid burying the main take-away in such technical details, we do not present the entire data in the main paper. Here we also evaluated the exact upper bound from Eq. (C1) for each curve, which converges to the asymptotic results shown in the main paper (see Fig. 13 for details of the convergence). In Fig. 11 (b), we calculate the critical number  $N_c$  for  $\langle C \rangle = 1$ , which is much larger than the giant component transition point  $N_G$  or the results from Ref. [24]. We can also solve  $\langle \mathcal{C}(\boldsymbol{x}) \rangle = 1$  in Eq. (C1) to obtain a lower bound estimate on  $N_c$ , which works well when R is large as shown in Fig. 11(c). In Fig. 11(d), we plot the capacity in linear scale to show the deviations between the actual capacities and the upper bounds in more detail. The major reason for the deviation at large R and high density is due to the second inequality of Ineq. (3) of the main paper, which we also print here

$$\langle \mathcal{C}(\boldsymbol{x}, \boldsymbol{x}') \rangle \leq \langle \min \{ \mathcal{C}(\boldsymbol{x}), \mathcal{C}(\boldsymbol{x}') \} \rangle \leq \langle \mathcal{C}(\boldsymbol{x}) \rangle, \quad (B1)$$

as interchanging the order of ensemble averaging and minimization is not tight.

#### Appendix C: Derivation of the asymptotic results for Waxman model

Due to the independence between the edges between nodes, we have

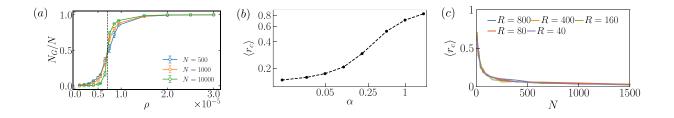


Figure 8. (a) The relative size  $N_G/N$  of the largest component in the Waxman QN model vs. the density of nodes,  $\alpha, L$  are both determined by N and density. To obtain the average, we sampled 10, 10, 5 graphs for  $N = 500, 1000, 10^4$  separately. The dashed vertical line at a density  $\sim 7 \times 10^{-6}$  indicates the transition point. (b) Clustering coefficients vs  $\alpha$  for Waxman QN model, in the large number of nodes  $N \gg 1$  limit. (c) Clustering coefficients of the scale-free QN model.

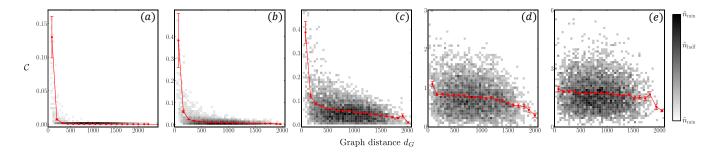


Figure 9. More details on Fig. 2(b1)-(b5). We have used the same numbering of the subplots for consistency. The gray scale PDF represents the statistical distribution (plotted in nonlinear scale  $\sqrt{\tilde{n}}$  for visualization) of end-to-end capacity over 5000 random pairs of end nodes (50 pairs from each of the 100 random QNs). The red lines are the average end-to-end capacity in each of the distance window. We sort the 5000 samples according to the graph distances from small to large and divided them into 20 groups of 250 points accordingly. We take the average of the capacity and graph distance in each group and obtain a data point.

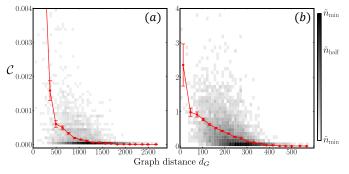


Figure 10. More details on Fig. 4(b1)(b2) in the main paper. We have used the same numbering of the subplots for consistency. The gray scale PDF represents the statistical distribution (plotted in nonlinear scale  $\sqrt{\tilde{n}}$  for visualization) of end-to-end capacity over 5000 random pairs of end nodes (50 pairs from each of the 100 random QNs). The red lines are the average end-to-end capacity in each of the distance window. We sort the 5000 samples according to the graph distances from small to large and divided them into 20 groups of 250 points accordingly. We take the average of the capacity and graph distance in each group and obtain a data point.

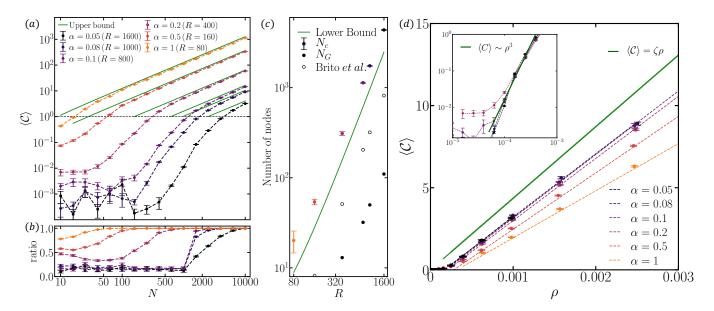


Figure 11. (a) Average end-to-end capacity  $\langle C \rangle$  vs. number of nodes N for various  $\alpha$ 's. The system size  $R \simeq 80/\alpha$  km. We see a critical drop at small N when  $\alpha$  is not too large, indicated by the dashed lines going down to zero. The green solid lines gives the upper bounds in Eq. (C1). (b) The average of the ratio of end-node edges inside the minimum cut. It shares the same legend as in (a). (c) We plot the critical number of nodes vs  $\alpha$ . The green curve indicates a lower bound from solving  $\langle \mathcal{C}(\boldsymbol{x}) \rangle = 1$  in Eq. (C1). In comparison, we plotted the results predicted from Ref. [24] (black open circles) and the critical number of nodes  $N_G$  for the appearance of giant components. (d) Average end-to-end capacity  $\langle C \rangle$  vs. density of nodes  $\rho$  for various  $\alpha$ 's in a linear scale. The system size  $R \simeq 80/\alpha$  km. The green line represents the asymptotic upper bounds  $\langle \mathcal{C} \rangle \simeq \zeta \rho$ and the dashed lines with same colors as dots shows the linear fitting in the range  $\langle \mathcal{C} \rangle > 0.1$ . The inset is the dependence of average end-to-end capacity with density in the range  $\rho \in (10^{-5}, 10^{-3})$ . The dark blue solid line presents a power-law relation as  $\langle \mathcal{C} \rangle \sim \rho^3$ .

$$\left\langle \mathcal{C}\left(\boldsymbol{x}\right)\right\rangle = \frac{(N-1)}{|\Omega_{R}|^{2}} \int_{\Omega_{R}} d^{2}\boldsymbol{x} \int_{\Omega_{R}} d^{2}\boldsymbol{x}' \Pi\left(\boldsymbol{x}, \boldsymbol{x}'\right) \mathcal{C}\left(E_{\boldsymbol{x}, \boldsymbol{x}'}\right)$$
(C1)

$$= -\frac{(N-1)}{|\Omega_R|^2} \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_R} d^2 \boldsymbol{x}' e^{-D(\boldsymbol{x}, \boldsymbol{x}')/\alpha L} \log_2\left(1 - 10^{-\gamma D(\boldsymbol{x}, \boldsymbol{x}')}\right)$$
(C2)

$$= -\frac{(N-1)}{|\Omega_R|^2} \left[ \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_\infty} d^2 \boldsymbol{x}' e^{-D(\boldsymbol{x}, \boldsymbol{x}')/\alpha L} \log_2 \left( 1 - 10^{-\gamma D(\boldsymbol{x}, \boldsymbol{x}')} \right) + O(R) \right]$$
(C3)

$$= -\frac{(N-1)}{|\Omega_R|} \int_{\Omega_\infty} d^2 \mathbf{x}' e^{-D(\mathbf{x},\mathbf{x}')/\alpha L} \log_2\left(1 - 10^{-\gamma D(\mathbf{x},\mathbf{x}')}\right) + O(NR^{-3})$$
(C4)

$$= -\frac{(N-1)\pi}{2R^2} \int_0^\infty r \ dr \ e^{-r/\alpha L} \log_2\left(1 - 10^{-\gamma r}\right) + O(NR^{-3}) \tag{C5}$$

$$= -\frac{(N-1)\pi}{2R^2} \int_0^\infty r \, dr \, e^{-r/\alpha L} \log_2\left(1 - 10^{-\gamma r}\right) + O(NR^{-3}) \tag{C6}$$

$$= -2\pi\rho \int_0^\infty r \, dr \, e^{-r/\alpha L} \log_2\left(1 - 10^{-\gamma r}\right) + O(NR^{-3}) + O(R^{-2}). \tag{C7}$$

Inputting  $\alpha L = 226$  and  $\gamma = 0.02$  we have the asymptotic expansion of

$$\langle \mathcal{C} \left( \boldsymbol{x} \right) \rangle = \zeta \rho, \tag{C8}$$
  
$$\zeta = -2\pi \int^{\infty} dr \ r e^{-r/226} \log_2 \left( 1 - 10^{-0.02r} \right) \simeq 4357.9.$$

In Fig. 13, we compare the asymptotic results with the exact numerical integration in Eq. (C1). A good convergence towards the asymptotic result is found with the increasing scale R.

$$T = -2\pi \int_0^\infty dr \ r e^{-r/226} \log_2\left(1 - 10^{-0.02r}\right) \simeq 4357.9.$$
(C9)

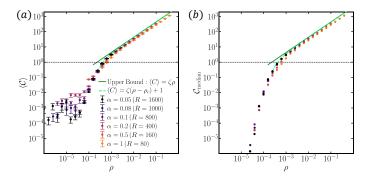


Figure 12. (a) Average end-to-end capacity  $\langle \mathcal{C} \rangle$  vs. density of nodes  $\rho$  for various  $\alpha$ 's. The system size  $R \simeq 80/\alpha$  km. We see a critical drop at small N when  $\alpha$  is not too large, indicated by the dashed lines going down to zero. The green solid lines gives the upper bounds in Eq. (C1). (b) The median end-to-end capacity  $\langle \mathcal{C} \rangle$  vs. density of nodes  $\rho$  for various  $\alpha$ 's. It shares the same legend as in (a). Instead of a long tail, we see clear sharp drop around the transition point.

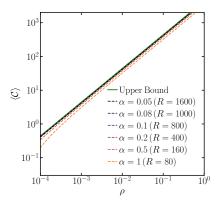


Figure 13. Comparison of the exact upper bound in Eq. (C1) and its asymptotic limit  $\langle C \rangle \simeq \zeta \rho$  for the Waxman model. We plot those upper bound by dashed lines with different  $\alpha$  and asymptotic limit in orange line.

#### Appendix D: Derivation of node capacity in Erdős Rényi models

$$\begin{aligned} \langle \mathcal{C} \left( \boldsymbol{x} \right) \rangle &= \frac{(N-1)}{|\Omega_R|^2} \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_R} d^2 \boldsymbol{x}' \Pi \left( \boldsymbol{x}, \boldsymbol{x}' \right) \mathcal{C}_E \left( E_{\boldsymbol{x}, \boldsymbol{x}'} \right) \\ &= -\frac{(N-1)p}{|\Omega_R|^2} \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_R} d^2 \boldsymbol{x}' \log_2 \left( 1 - 10^{-\gamma D(\boldsymbol{x}, \boldsymbol{x}')} \right) \\ &= -\frac{(N-1)p}{|\Omega_R|} \int_{\Omega_\infty} d^2 \boldsymbol{x}' \log_2 \left( 1 - 10^{-\gamma D(\boldsymbol{x}, \boldsymbol{x}')} \right) + O(NR^{-3}) \\ &= -\frac{(N-1)p\pi}{2R^2} \int_0^\infty r \ dr \log_2 \left( 1 - 10^{-\gamma r} \right) + O(NR^{-3}) \\ &= \zeta_{ER} p\rho + O(NR^{-3}) + O(R^{-2}), \end{aligned}$$

#### Appendix E: Derivation of the asymptotic results for scale-free model

Considering the on average 2m neighbours as independent, the ensemble-averaged node capacity is

$$\left\langle \mathcal{C}\left(\boldsymbol{x}\right)\right\rangle = \frac{2m}{|\Omega_{R}|^{2}} \int_{\Omega_{R}} d^{2}\boldsymbol{x} \int_{\Omega_{R}} d^{2}\boldsymbol{x}' \left\langle \Pi\left(\boldsymbol{x},\boldsymbol{x}'\right) \mathcal{C}\left(E_{\boldsymbol{x},\boldsymbol{x}'}\right)\right\rangle$$

$$= \frac{2m}{A} \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_R} d^2 \boldsymbol{x}' \left\langle \frac{D_g(\boldsymbol{x}')}{D(\boldsymbol{x}, \boldsymbol{x}')} \mathcal{C}(E_{\boldsymbol{x}, \boldsymbol{x}'}) \right\rangle$$
(E2)

where the normalization constant

2

$$\mathbf{A} = \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_R} d^2 \boldsymbol{x}' \left\langle \frac{D_g(\boldsymbol{x}')}{D(\boldsymbol{x}, \boldsymbol{x}')} \right\rangle.$$
(E3)

The  $\langle \cdot \rangle$  inside the integral now denotes average over the degree distribution of neighbours, conditioned on the neighbour being at  $\mathbf{x}'$ . We can approximate the distribution of the degree as independent of the distance to node  $\mathbf{x}$ , then  $\langle D_g(\mathbf{x}') f(\mathbf{x}, \mathbf{x}') \rangle = \langle D \rangle f(\mathbf{x}, \mathbf{x}')$ , where  $\langle D \rangle$  is a constant and  $f(\mathbf{x}, \mathbf{x}')$  is an arbitrary function of  $\mathbf{x}, \mathbf{x}'$ . We can cancel out the constant and equivalently calculate

$$\left\langle \mathcal{C}\left(\boldsymbol{x}\right)\right\rangle = \frac{2m}{A'} \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_R} d^2 \boldsymbol{x}' \left\langle \frac{1}{D\left(\boldsymbol{x}, \boldsymbol{x}'\right)} \mathcal{C}\left(E_{\boldsymbol{x}, \boldsymbol{x}'}\right)\right\rangle$$
(E4)

$$A' = \int_{\Omega_R} d^2 \boldsymbol{x} \int_{\Omega_R} d^2 \boldsymbol{x}' \left\langle \frac{1}{D(\boldsymbol{x}, \boldsymbol{x}')} \right\rangle.$$
(E5)

The above integral can be numerically calculated. It is clear that  $\langle \mathcal{C}(\boldsymbol{x}) \rangle$  does not grow with the number of nodes N, as m is now a constant.

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