

Statistical properties of sampled networks

Sang Hoon Lee,^{*} Pan-Jun Kim,[†] and Hawoong Jeong[‡]

Department of Physics, Korea Advanced Institute of Science and Technology, Daejeon 305-701, Korea

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We study the statistical properties of the sampled scale-free networks, deeply related to the proper identification of various real-world networks. We exploit three methods of sampling and investigate the topological properties such as degree and betweenness centrality distribution, average path length, assortativity, and clustering coefficient of sampled networks compared with those of original networks. It is found that the quantities related to those properties in sampled networks appear to be estimated quite differently for each sampling method. We explain why such a biased estimation of quantities would emerge from the sampling procedure and give appropriate criteria for each sampling method to prevent the quantities from being overestimated or underestimated.

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I. INTRODUCTION

Recently, a huge amount of research on complex networks has been achieved in interdisciplinary fields including mathematics, statistical physics, computer science, sociology, biology, etc. [1–3]. Complex networks are ubiquitous in the real world, e.g., there are technological networks such as the Internet [4], biological networks such as protein interaction networks [5], and social networks such as scientific collaboration networks [6]. Various models to explain the observed properties of those real networks have been introduced and studied by both numerical and analytic approaches. Relatively fewer works, however, have been done about possible error or bias in collecting data and identifying real networks in a practical sense, and most works deal exclusively with either social networks or the Internet [7–17].

For instance, a survey of relationships among participants has to be conducted to construct a social network, but the collected network data may be incomplete or erroneous since a survey usually targets only a partial sample of a whole population [9]. The topology of the Internet is inferred by aggregating paths or *traceroutes* [18], which also reveals only a part of the whole Internet [10–13]. In biology, protein-protein interaction networks are identified by seeking contextual or cellular functions mostly within specific functional modules [5]. Identification of such networks by experiments also has a fundamental limit naturally. Thus, all these networks identified are *sampled* networks from complete structures. In addition, if the size of an entire network is too large to measure some quantities such as betweenness centrality [19,20] due to time complexity, inevitably a sampling process is necessary.

So far models of networks have been designed based on features observed in real networks, such as the small-world effect [21] and the power-law degree distribution [22,23]. But what if those *observed* characteristics from the *sampled*

networks are considerably different from the original structures of the real networks? It has been shown that the sampled networks based on the traceroute sampling method may have significantly different topological properties from the original network in some cases [10–13]. Effects of missing data in social networks are discussed in Ref. [9], in which it was shown that some problems in conceiving social networks can cause incompleteness of data and lead to misestimation of quantities like mean node degree, clustering coefficient, assortativity, etc. At this point, bias in such quantities needs to be considered in a more general sense.

In a statistical sense, inference from a sample provides fairly reasonable estimation of a whole population if a large number of objects are selected randomly enough to be representative in the population. This naive criterion, however, cannot be applied directly to sampling networks, since there are two different elements, i.e., nodes and links in a network. A degree distribution of nodes is, for example, a statistic of a network, but the degree is not an independent characteristic of each node. Nodes are literally connected to one another, by the other kind of components called links from which a degree is defined. Similarly, other properties of a network also heavily depend on the way that nodes and links are interwoven. There could be several different ways of sampling networks due to the two interrelated elements (nodes and links), and each method may give distinctive features with respect to such properties.

There has been a large amount of work on random breakdowns or intentional attack on complex networks, considered as the exact reverse process of sampling, in the physics community [24–27]. The analytic methods in that work, therefore, can be also applied to the sampling problem. In this paper, we adopt three basic methods of sampling networks and investigate the effect of each method on measuring several well-known network quantities such as degree distribution, average path length, betweenness centrality distribution [19,20], assortativity [28], and clustering coefficient [21]. Observed bias of such quantities is explained, and we provide appropriate criteria for choosing sampling methods to measure the quantities more accurately. Some typical real networks as well as the Barabási-Albert model [22] are sampled for this analysis. More general sampling processes

^{*}Electronic address: lshlj@stat.kaist.ac.kr

[†]Electronic address: pj@stat.kaist.ac.kr

[‡]Electronic address: hjeong@kaist.ac.kr

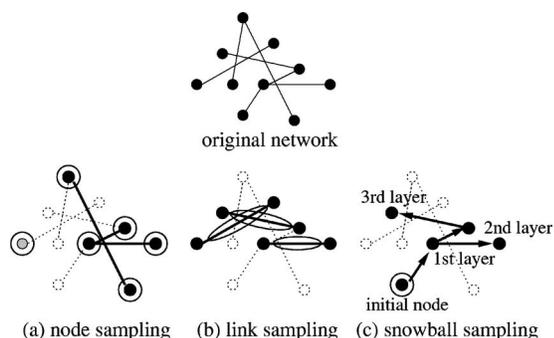


FIG. 1. Three kinds of sampling method. (a) Node sampling: Select the circled nodes, keep three links among them, and the isolated node is removed. (b) Link sampling: Select the three circled links and six nodes attached to them. (c) Snowball sampling: Starting from the circled node, select nodes and links attached to them by tracing links.

used to identify real networks may consist of some combinations of methods presented here or variations of them, but we can infer by using the results from the basic methods.

II. SAMPLING METHODS AND NETWORKS

We introduce three kinds of sampling method called node sampling, link sampling, and snowball sampling. In node sampling, a certain number of nodes are randomly chosen and links among them are kept. The sampling fraction in this method is defined as the ratio of the number of chosen nodes (including isolated nodes that will be removed later) to that of all the nodes in the original network. As in Fig. 1(a), isolated nodes are neglected for convenience, although they are fully predictable, so the number of nodes in a sampled network is a little bit less than that of selected nodes. We observe the dependence of the number of chosen links on that of nodes, since it is related to the average degrees and average path length of sampled networks, discussed later on. Suppose the fraction of number of selected nodes is α and that of links among them is β . Then it is found that $\beta \sim \alpha^2$ if we pick nodes randomly, since the maximum number of (undirected) links possible for n selected nodes are $\binom{n}{2} = n(n-1)/2 \sim n^2$ [29].

In link sampling, a certain number of links are randomly selected and nodes attached to them are kept, as in Fig. 1(b). In snowball sampling [30,31], we first choose a single node and all the nodes directly linked to it are picked. Then all the nodes connected to those picked in the last step are selected, and this process is continued until the desired number of nodes are sampled. The set of nodes selected in the n th step is denoted as the n th layer, in the same sense of “radius” for ego-centered networks in Ref. [31]. See Fig. 1(c) for illustration. To control the number of nodes in the sampled network, a necessary number of nodes are randomly chosen from the last layer. Similar to the cluster-growing method used to calculate the fractal dimension of percolation clusters in Ref. [32], the snowball sampling method tends to pick hubs (nodes with many links) in short step due to high connectivity of them. So whether the initial node is a hub or not

TABLE I. The numbers of nodes n and links l for each real network.

Network	n	l	Ref.
PIN	5077	16 449	[5,33]
Internet AS	10 515	21 455	[34]
arxiv.org	49 983	245 300	[6]

does not make a noticeable difference in characterizing the sampled network.

For numerical analysis of the sampled networks, we use Barabási-Albert (BA) scale-free network as an example of model networks, which follows the power-law degree distribution $p(k) \sim k^{-3}$, with 30 000 nodes and $m_0 = m = 4$ [22]. We also consider three real-world networks from various fields, including protein interaction network (PIN) [5,33], the Internet at the autonomous systems (AS) level [34], and e-print archive coauthorship network (arxiv.org) [6]. The numbers of nodes and links for each network are in Table I. Although results from other homogeneous networks are also discussed in Sec. IV, most of networks considered in this work are undirected and scale-free networks following power-law degree distribution, $p(k) \sim k^{-\gamma}$, where $2 < \gamma \leq 3$.

III. CHARACTERISTICS OF SAMPLED NETWORKS

A. Degree distribution and average path length

A degree of a node is defined as the number of links attached to the node. Many real networks are shown to have a power-law degree distribution $p(k) \sim k^{-\gamma}$ [1–3], including the networks considered in this paper. We found that in general degree distributions of sampled networks from the four networks obtained by all three methods follow the power-law as well as those of the original networks. The exponents of degree distribution γ (degree exponent) are extracted using maximum likelihood estimate given by the formula [35]

$$\gamma = 1 + n \left[\sum_{i=1}^n \ln \frac{k_i}{k_{\min}} \right]^{-1}, \quad (1)$$

where n is the number of elements in a set $\{k_i\}$ whose elements follow the power-law distribution $p(k) \sim k^{-\gamma}$, and k_{\min} is the smallest element for which the power-law behavior holds. Figure 2 shows the change of the degree exponent for the sampled networks from each network obtained by numerical simulation for each method as we change the sampling fraction α .

For node sampling, we fix the number of sampled nodes and select nodes randomly. In this case, the new degree distribution $p'(k)$ of the sampled network is expressed as

$$p'(k) = \sum_{k_0=k}^{n-1} p(k_0) \binom{k_0}{k} \binom{n-k_0-1}{n_s-k-1} \bigg/ \binom{n-1}{n_s-1}, \quad (2)$$

where $p(k)$ is the degree distribution of the original network, n is the number of nodes in the original network, and $n_s = \alpha n$ is the fixed number of sampled nodes. In the case that

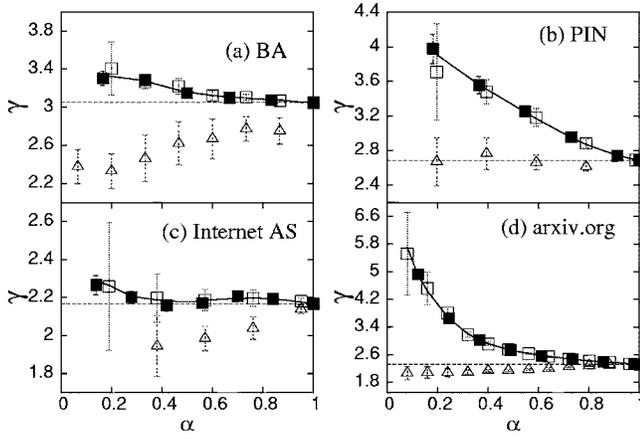


FIG. 2. Changes of degree exponent γ for each network's sampled networks according to the sampling fraction α , averaged over ten independent realizations. Empty squares (\square) stand for node sampling, filled squares (\blacksquare) for link sampling, and empty triangles (\triangle) for snowball sampling. The horizontal dashed lines are the values for the original exponent of each network, and the solid lines represent the values obtained by Eq. (5).

the number of nodes in sampled networks is not fixed but only the probability α with which individual nodes are selected is given [14,17], Eq. (2) should be written as

$$p'(k) = \sum_{n_s=k+1}^n \sum_{k_0=k}^{n-1} p(k_0) f(n_s) \binom{k_0}{k} \frac{\binom{n-k_0-1}{n_s-k-1}}{\binom{n-1}{n_s-1}}, \quad (3)$$

where the probability that n_s number of nodes are chosen is $f(n_s) = \binom{n}{n_s} \alpha^{n_s} (1-\alpha)^{n-n_s}$. If the number of nodes is fixed, $f(n_s) = \delta(n_s - an)$ and Eq. (3) becomes Eq. (2) with $n_s = an$. Even if the number of nodes is not fixed, when the system size is large enough to use the approximation $f(n_s) \approx \delta(n_s - an)$, we can safely use Eq. (2). Equation (2) can be further reduced by $n!/(n-m)! \approx n^m$ for $n \gg m$. Suppose $n, n_s \gg k_0, k$. Then

$$\begin{aligned} \frac{\binom{n-k_0-1}{n_s-k-1}}{\binom{n-1}{n_s-1}} &\approx n_s^k (n-n_s)^{k_0-k} / n^{k_0} \\ &= \left(\frac{n_s}{n}\right)^k \left(1 - \frac{n_s}{n}\right)^{k_0-k} \\ &= \alpha^k (1-\alpha)^{k_0-k}, \end{aligned} \quad (4)$$

which leads to the formula previously used in Refs. [14,17,25]

$$p'(k) = \sum_{k_0=k}^{\infty} p(k_0) \binom{k_0}{k} \alpha^k (1-\alpha)^{k_0-k}. \quad (5)$$

The sizes of all the four networks studied in this paper are larger than 5000, and we have checked that Eqs. (2) and (5) give practically the same values of $p'(k)$ and are indistinguishable in the graphs. For much smaller networks, on the other hand, Eq. (2) actually predicts the degree distribution

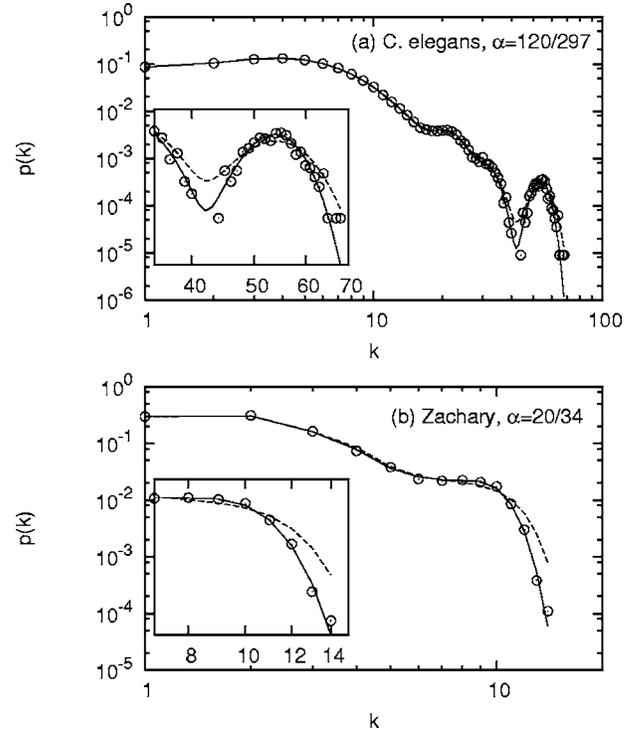


FIG. 3. Degree distribution for sampled networks of (a) *C. elegans* neural network with $\alpha=120/297$ and (b) Zachary karate club network with $\alpha=20/34$, obtained from the node sampling. Empty circles are simulation results from 1000 sampling processes. Solid lines correspond to Eq. (2) and dashed lines to Eq. (5). Insets show the part of large degrees, where the difference between two formulae is prominent, for each graph.

of sampled networks better than Eq. (5). In Fig. 3, we compare the simulation results for two small networks, the nematode *C. elegans* neural network [37] with 297 nodes and 2359 links and the Zachary karate club network [38] with 34 nodes and 77 links, with those two equations by substituting the original degree distribution $p(k) = n_k/n$, where n_k is the number of nodes with degree k . The figure clearly shows that Eq. (2) is more accurate.

The above equations turn out to be applied to the link sampling with the same sampling fraction α as well. Here we can use the technique in Ref. [36] to solve the bond percolation or epidemic model. Suppose a node, which originally had k_0 links before sampling, comes to have k links. Because the random link sampling chooses links uniformly, the probability of the node having k out of k_0 links is $p(k|k_0) = \binom{k_0}{k} \alpha^k (1-\alpha)^{k_0-k}$. Consequently the probability that a node in the sampled network has degree k from all the possible original degree k_0 is $p'(k) = \sum_{k_0=k}^{\infty} p(k_0) p(k|k_0)$, which leads us back to Eq. (5). The fact that those two sorts of sampling are described by the same equation is also supported by Fig. 2 showing the similar degree exponent changes for both node and link sampling.

As Stumpf *et al.* point out in Refs. [14,17], Eq. (5) for a power-law degree distribution $p(k) \sim k^{-\gamma}$ yields deviation of $p'(k)$ from the original power-law form for quite small sampling fraction α . For moderate values of α , however, the

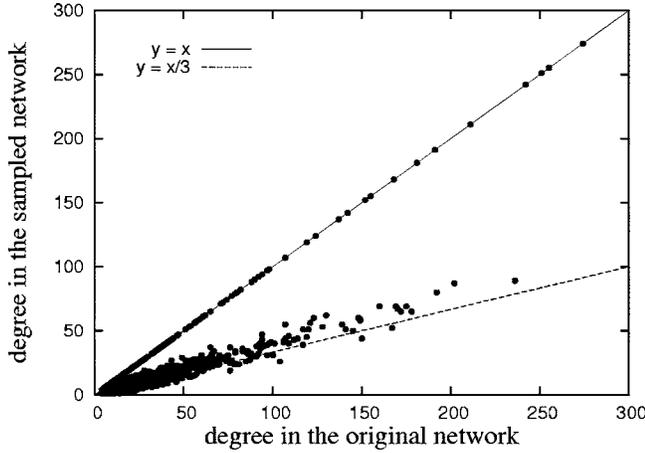


FIG. 4. Change of nodes' degree in BA network for snowball sampling. The sampling fraction is 10 000/30 000.

deviation is not significant and we observe that the tangent of $p'(k)$ in the log-log plot actually becomes steeper from Eq. (5), consistent with our numerical observation about the node and link sampling as shown in Fig. 2. To extract the degree exponents from Eq. (5), first we calculate the degree distribution of the original networks by $p(k)=n_k/n$. Substituting that $p(k)$ into Eq. (5), we obtain the degree distribution $p'(k)$ of the sampled networks corresponding to a given sampling fraction α . The degree exponents from those $p'(k)$ in Fig. 2 show good agreement with the values from numerical simulation for both node and link sampling cases.

On the contrary, it is found that a degree exponent decreases for snowball sampling as we decrease the sampling fraction. By the definition of snowball sampling, hubs are more likely to be selected by this method. Furthermore, once a hub is picked, every node connected to the hub is selected in the next step unless it belongs to the previous layer. This characteristic of snowball sampling tends to *conserve* the degrees of easily selected hubs, which leads to the decrease of degree exponents by holding the “tail” of the power-law degree distribution. Figure 4 shows the degrees in a sampled network obtained by snowball sampling, and the nodes with large degree on the $y=x$ line clearly indicates a tendency to choose hubs and conserve their degrees. Therefore, the snowball sampling underestimates the degree exponent. In Ref. [11], they show that the traceroute sampling can underestimate the degree exponent of a scale-free network by undersampling the low-degree nodes relative to the high-degree ones. In spite of the difference between the snowball and traceroute sampling, both of these methods overrepresent hubs and have the same “crawling” character used to identify the nodes. We infer that the decrease of degree exponents for both sampling methods is caused by these similar features.

We also check two closely related quantities, namely the average degree and the average path length (APL) in the sampled networks. APL is the average of shortest paths between all the pairs of nodes in a network, often used as a measure of network efficiency. In Fig. 5, we present APL of the *giant component* in the sampled networks obtained by the numerical simulation. For snowball sampling, APL decreases according to the decreased system size of sampled networks.

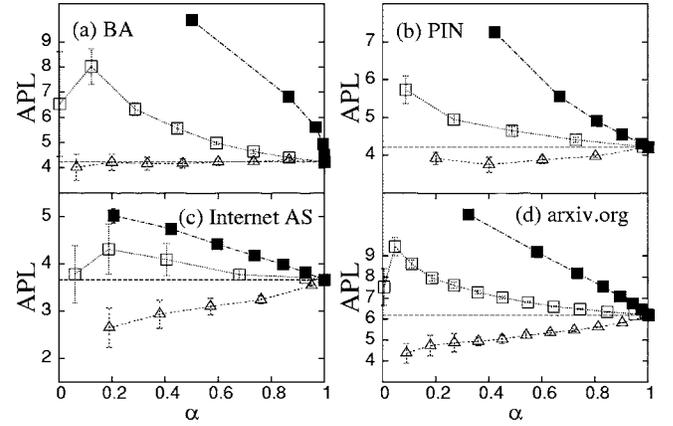


FIG. 5. Changes of APL for each network's sampled networks according to the ratio ξ of the size of giant component in the sampled networks to that of the original ones, averaged over ten independent realizations. Empty squares (\square) stand for node sampling, filled squares (\blacksquare) for link sampling, and empty triangles (\triangle) for snowball sampling. The horizontal dashed lines are the values for the original APL of each network, and the other lines are guides to the eyes.

On the other hand, for node and link sampling, the APL of a sampled network is larger than that of the original network for not-so-small sampling fraction, even though the size of the sampled network itself is smaller than the original one. As presented earlier, for node sampling, the number of links is proportional to the square of the number of the nodes, which leads to $\langle k \rangle = 2l/n \propto n$, where l and n are the numbers of links and nodes in a sampled network, respectively. This suggests that the average degree in a sampled network decreases as the sampling fraction becomes smaller. Obviously, for a given network, APL decreases as the average degree increases. The diminishment in the average degree, therefore, seems to have a stronger effect on APL than the overall system size in this case. Similar behavior of the average degree and APL is observed for link sampling, but in this case it seems that the “treelike” structure of sampled networks, related to the clustering coefficient discussed later, is responsible for that behavior.

B. Betweenness centrality distribution

Betweenness centrality (BC or load), which measures the centrality of a node by the traffic flow in a network, of node b is defined as

$$g_b = \sum_{i \neq j} \frac{C_b(i,j)}{C(i,j)}, \quad (6)$$

where $C(i,j)$ is the number of all the shortest pathways between a pair of nodes (i,j) and $C_b(i,j)$ is that of the shortest pathways running through a node b [20]. It is known that the BC distribution follows a power-law $p(g) \sim g^{-\eta}$ for scale-free networks [19,20].

Similar to the degree distribution, the BC distribution of sampled networks also follows power-law well as do the original networks. Figure 6 shows the change of the BC ex-

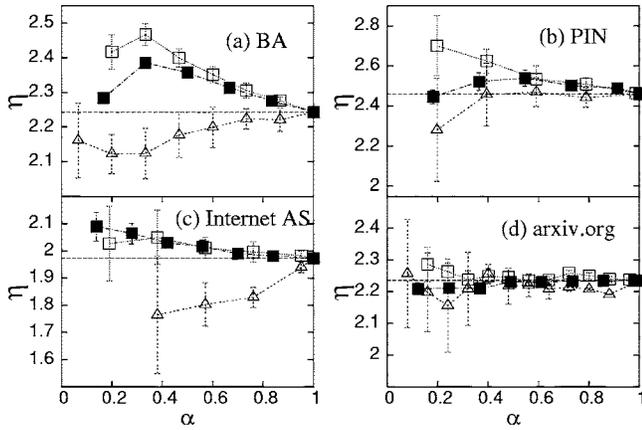


FIG. 6. Changes of BC exponent η for each network's sampled networks according to the sampling fraction α , averaged over ten realizations. Empty squares (\square) stand for node sampling, filled squares (\blacksquare) for link sampling, and empty triangles (\triangle) for snowball sampling. The horizontal dashed lines are the values for the original exponent of each network, and the other lines are guides to the eyes.

ponent, also obtained by Eq. (1), for each network and each sampling method. Similar to the degree exponent case, in general, BC exponents increase for node and link sampling and decrease for snowball sampling as the sampling fraction gets lower. Figure 6 bears a resemblance to Fig. 2 except for the case of arxiv.org, for which the BC exponent seems to be conserved for all the sampling methods. The correlation between degree and BC of nodes [39], shown in Fig. 7, could explain the same direction of changes of degree and BC exponents. For assortative networks such as arxiv.org here, however, it is known that the degree-BC correlation is not clear [40], which explains the different behavior in Fig. 6(d). Therefore, at least empirically, we expect overestimation of a BC exponent by node and link sampling and underestimation by snowball sampling.

C. Assortativity

The assortativity r , which measures the correlation between degrees of node linked to each other, is defined as the

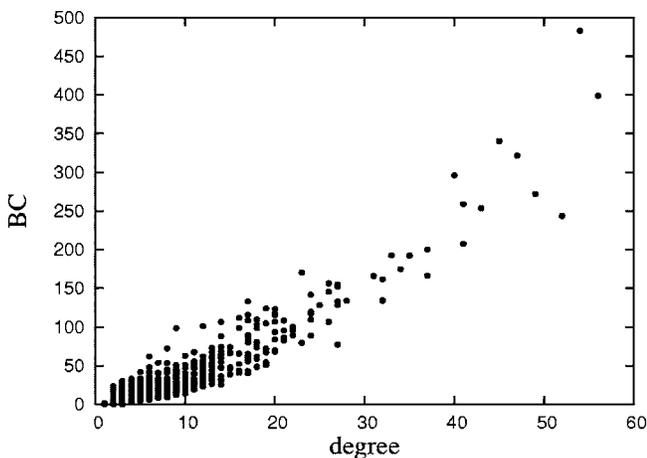


FIG. 7. Degree and BC of nodes in a sampled network of BA network by node sampling. The sampling fraction is 10 000/30 000. The value of BC is rescaled by the number of nodes.

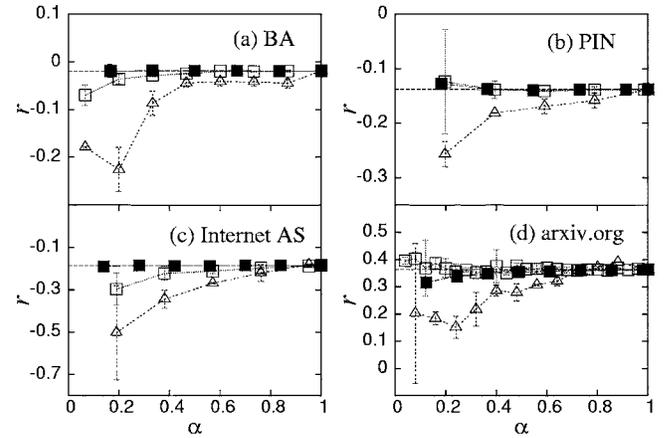


FIG. 8. Changes of assortativity r for each network's sampled networks according to the sampling fraction α , averaged over ten realizations. Empty squares (\square) stand for node sampling, filled squares (\blacksquare) for link sampling, and empty triangles (\triangle) for snowball sampling. The horizontal dashed lines are the values for the original assortativity of each network, and the other lines are guides to the eyes.

Pearson correlation coefficient of degrees between pairs of nodes [28]. Positive values of r stand for the positive degree-degree correlation which means that nodes with large degrees tend to be connected to one another. Most social networks have this positive degree correlation (*assortative* mixing), including the arxiv.org network considered in this paper. On the other hand, most biological and technological networks show negative degree correlation (*disassortative* mixing), including PIN and Internet AS network here. If there is no degree correlation among nodes (neutral), as in the case of BA model, the value of r is in the vicinity of 0.

The change of assortativity for each network and each method is shown in Fig. 8. For node and link sampling, no noticeable changes of assortativity in the sampled networks are observed. Random choice of nodes or links appears to conserve assortativity well for these two methods. Sampled networks from snowball sampling, however, are shown to be more disassortative than the original networks. This pattern is common no matter whether the original network is assortative (arxiv.org), disassortative (PIN and Internet AS), or neutral (BA).

There is another way to check the degree correlation, which is measuring the quantity $\langle k_{nn}(k) \rangle = \sum_k k' p(k'|k)$, i.e., the average degree of nearest neighbors of nodes with degree k [41]. Assortative mixing is represented by a positive slope of the $\langle k_{nn}(k) \rangle$ graph, while the others by horizontal (neutral) or a negative slope (disassortative). Figure 9 shows the changes of these slopes for $\langle k_{nn}(k) \rangle$ graphs of the sampled networks from two kinds of original networks by snowball sampling. The slope decreases, i.e., moves toward the negative value as the sampling fraction gets lower for both disassortative Internet AS and assortative arxiv.org.

We suggest that the more disassortative nature of sampled networks compared with the original ones is due to the last layer of snowball sampling method. In contrast to the conserved structure of the inner layers, a considerable number of

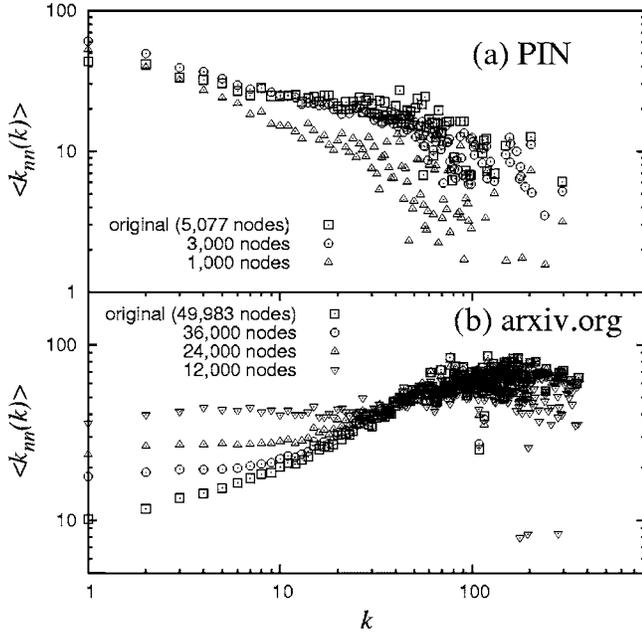


FIG. 9. $\langle k_{nn}(k) \rangle$ distribution for sampled networks of (a) PIN, (b) arxiv.org by snowball sampling.

links are lost for the nodes in the last layer. Meanwhile, hubs are likely to be selected for snowball sampling. This separation of “core” and “periphery” part is seen in Fig. 4, and the connections between hubs and nodes of the last layer can reduce the value of assortativity. The simulation shows that a sampled network containing the entire last layer is more disassortative than the one where only parts of the last layer are kept, which supports the hypothesis that the effect of the last layer induces disassortative mixing. Therefore, we have to be careful when measuring the assortativity for the network from the snowball sampling.

D. Clustering coefficient

The clustering coefficient C_i of node i is the ratio of the total number y of the links connecting its nearest neighbors to the total number of all possible links between all these nearest neighbors [3],

$$C_i = \frac{2y}{k_i(k_i - 1)}, \quad (7)$$

where k_i is the degree of node i . The clustering coefficient of a network is the average of this value over all the nodes $C = \sum_i C_i / n$, where n is the number of nodes. Most real networks have much larger value of clustering coefficient than model networks such as ER or BA network due to, e.g., the community or modular structure.

In Fig. 10, we show the change of clustering coefficient for each original network and each sampling method. For node and snowball sampling, there is a little change of clustering coefficient depending on networks. On the other hand, link sampling prominently reduces the clustering coefficient. This effect is obvious since the random omission of links, the reverse process of link sampling, “opens up triangles fast” as

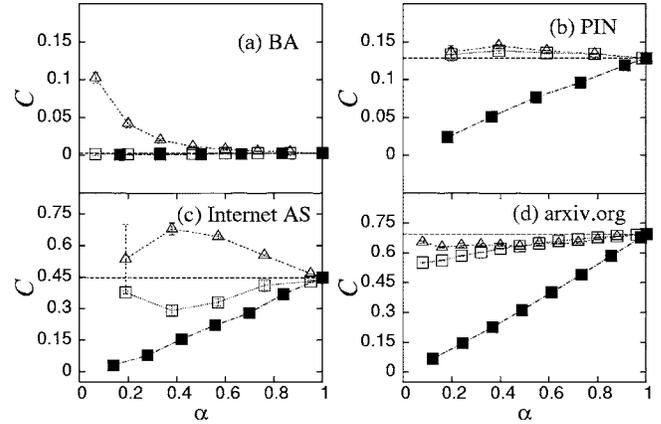


FIG. 10. Changes of clustering coefficient C for each network’s sampled networks according to the sampling fraction α , averaged over ten realizations. Empty squares (\square) stand for node sampling, filled squares (\blacksquare) for link sampling, and empty triangles (\triangle) for snowball sampling. The horizontal dashed lines are the values for the original clustering coefficient of each network, and the other lines are guides to the eyes.

stated in Ref. [9]. The link sampling, therefore, underestimates clustering coefficient of a network.

IV. DISCUSSION AND CONCLUSIONS

In this paper, we have studied the changes of well-known quantities in complex networks for randomly sampled networks. Three kinds of sampling methods are applied, and three representative real-world networks, along with the BA model, are used as the original networks for numerical investigation. We have measured four typical quantities in sampled networks, which shows some characteristic patterns in changes of the quantities for each sampling method. Based on properties of sampling methods, possible explanations for such changes as well as the mathematical analysis are provided. We have also analyzed other networks than the scale-free ones such as Erdős-Rényi random network [42] and the growing network without the preferential attachment [22], and the results show that the form of the degree distribution is conserved for the node and link sampling in those cases, consistent with the previous work [17].

Table II summarizes the results. To check the generality of the results, we also investigated the randomized version of

TABLE II. The changes of quantities in networks by each sampling method. As the sampling fraction gets lower (\Downarrow at the very right of each sampling method indicates this), \Uparrow stands for increase, \Downarrow for decrease, $=$ for the same, and \Downarrow for depending on networks.

	Degree Exponent γ	BC Exponent η	Assortativity r	Clustering Coefficient C
Node \Downarrow	\Uparrow	\Uparrow	$=$	\Downarrow
Link \Downarrow	\Uparrow	\Uparrow	$=$	\Downarrow
Snowball \Downarrow	\Downarrow	\Downarrow	\Downarrow	\Downarrow

each network in a similar fashion. The randomized networks were constructed by shuffling the links while conserving only the degree distribution [28]. We found the same results with the original networks. The results in Table II, therefore, seem to hold for scale-free networks in general and provide criteria for sampling method when some specific quantity is supposed to be investigated by the sampling. From another viewpoint, bias of some quantities can be predicted if a specific sampling method used to identify a network is known. If we are interested in the assortativity of a network, for example, node or link sampling can give fairly accurate values. For a clustering coefficient, on the other hand, the link sampling method should be avoided.

Sampling problems should be taken into account for real network research, but not much work has been done so far. Exploration of other characteristics of complex networks or

using other sampling methods, rigorous analytic approaches, and establishing solid principles by more systematic investigation could all be important research topics for the future. We hope this work can make a contribution to this direction of research.

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