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Degree and Principal Eigenvectors in Complex Networks

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Abstract. The largest eigenvalue λ_1 of the adjacency matrix powerfully characterizes dynamic processes on networks, such as virus spread and synchronization. The minimization of the spectral radius by removing a set of links (or nodes) has been shown to be an NP-complete problem. So far, the best heuristic strategy is to remove links/nodes based on the principal eigenvector corresponding to the largest eigenvalue λ_1 . This motivates us to investigate properties of the principal eigenvector x_1 and its relation with the degree vector. (a) We illustrate and explain why the average $E[x_1]$ decreases with the linear degree correlation coefficient ρ_D in a network with a given degree vector; (b) The difference between the principal eigenvector and the scaled degree vector is proved to be the smallest, when $\lambda_1 = \frac{N_2}{N_1}$, where N_k is the total number walks in the network with k hops; (c) The correlation between the principal eigenvector and the degree vector decreases when the degree correlation ρ_D is decreased.

Keywords: networks; spectral radius; principal eigenvector; degree; assortativity

1 Introduction

Dynamic phenomena occurring on networks are affected by the structure of networks, e.g., the absence of epidemic thresholds in large scale free networks [2][3][6], the effect of the degree correlations on the percolation of networks [8]. The largest eigenvalue $\lambda_1(A)$ of the adjacency matrix A , called the spectral radius of the graph, has been shown to play an important role in dynamic processes on graphs, such as SIS (susceptible-infected-susceptible) virus spread [12] and the Kuramoto type of synchronization process of coupled oscillators [11] on a given network topology. For instance, in a SIS spreading model, the epidemic threshold $\tau_c \simeq \frac{1}{\lambda_1(A)}$ separates two different phases of a dynamic process on a network: if the spreading rate τ is above the threshold, the infection spreads and becomes persistent in time; where $\tau < \tau_c$, the infection dies out exponentially fast [10][12]. In the past decade, researches have focused on how topological changes, such as link (or node) removal, may alter the spectral radius. Milanese *et al.* [7] studied the dynamical importance of the structural perturbation by removing one node or link. Van Mieghem *et al.* [15] have proved that to minimize

the largest eigenvalue by removing a set of links or nodes is a NP-hard problem and have shown that the best strategy so far is based on the components of the principal eigenvector x_1 , which underlines the importance of the principal eigenvalue in characterizing the influence of link/node removal on the spectral radius. Our main objective is to investigate the topological meaning of x_1 , which has been rarely studied. Especially, we aim to understand the relation between x_1 and the degree vector/sequence¹ d , the computationally simplest and mostly studied property of a network.

The degree correlation, also called the assortativity ρ_D is computed as the linear correlation coefficient of the degree of nodes connected by a link. It describes the tendency of network nodes to connect preferentially to other nodes with either similar (when $\rho_D > 0$) or opposite (when $\rho_D < 0$) properties i.e. degree [9]. The assortativity was widely studied after it was realized that the degree distribution alone provides an insufficient characterization of complex networks. Networks with the same degree distribution may still differ significantly in various topological features. Degree-preserving rewiring [13] allows us to either increase or decrease the assortativity of a network without changing the degree of each node. The relation between the principal eigenvector and the degree vector is systematically investigated in networks with various degree distributions and degree correlations.

Section 2 illustrates the importance of the principal eigenvector in characterizing the influence of link/node removal on the spectral radius by two key theories developed in our early work and further simulations. Subsequently, we explore the properties of the principal eigenvector and the relation between the (normalized) degree vector and the principal eigenvector in networks with different degree correlation and with the degree distribution derived from the Erdős-Rényi random graphs² [4], the Bárábasi-Albert graphs³ [1], and real-world networks (see Section 3). Our major contributions are: (a) the average of the components in the principal eigenvector $E[x_1]$ is shown and explained to decrease with the assortativity ρ_D ; (b) the difference between the principal eigenvector and the degree vector is proved to be the smallest, when $\lambda_1 = \frac{N_2}{N_1}$, where N_k is the total number of walks with k hops in a network and (c) the correlation between principal eigenvector and the degree vector decreases as the assortativity ρ_D is decreased. These finds provide essential inspiration on when the degree vector well approximates the principal eigenvector. Finally, we illustrate the possibility to approximate the principal eigenvector based strategy to minimize the largest

¹ The degree vector/sequence is composed of the degree of each node, following the same ordering as the principal eigenvector.

² An Erdős-Rényi random graph can be generated from a set of N nodes by randomly assigning a link with probability p to each pair of nodes.

³ A Bárábasi-Albert graph starts with m nodes. At every time step, we add a new node with m links that connect the new node to m different nodes already present in the graph. The probability that a new node will be connected to node i in step t is proportional to the degree $d_i(t)$ of that node. This is referred to as preferential attachment.

eigenvalue by removing links/nodes by its corresponding degree based strategy (see Section 4), which can be well explained by the findings in early sections.

2 The decrease of the spectral radius

We consider a network as a graph $G = (\mathcal{N}, \mathcal{L})$, where \mathcal{N} is the set of nodes and \mathcal{L} is the set of links. The number of nodes is denoted by $N = |\mathcal{N}|$ and the number of links is represented by $L = |\mathcal{L}|$. The graph G can be represented by the $N \times N$ adjacency matrix A , consisting of elements a_{ij} that are either one or zero depending on whether there is a link between nodes i and j . The eigenvalues of the adjacency matrix are ordered as $\lambda_N \leq \lambda_{N-1} \leq \dots \leq \lambda_1$, where λ_1 is the spectral radius and the corresponding eigenvector x_1 is called the principal eigenvector. Let \mathcal{L}_m (or \mathcal{N}_m) denote the set of the m links (or nodes) that are removed from G , and $G_m(\mathcal{L}) = G \setminus \mathcal{L}_m$ (or $G_m(\mathcal{N}) = G \setminus \mathcal{N}_m$) is the resulting graph after the removal of m links (or nodes) from G . We denote the adjacency matrix of $G_m(\mathcal{L})$ (or $G_m(\mathcal{N})$) by $A_m(\mathcal{L})$ (or $A_m(\mathcal{N})$), which is still a symmetric matrix.

Theorem 1. *For any graph G and graph $G_m(\mathcal{L}) = G \setminus \mathcal{L}_m$, by removing m links from G , it holds that*

$$2 \sum_{l \in \mathcal{L}_m} (w_1)_{l^+} (w_1)_{l^-} \leq \lambda_1(A) - \lambda_1(A_m(\mathcal{L})) \leq 2 \sum_{l \in \mathcal{L}_m} (x_1)_{l^+} (x_1)_{l^-} \quad (1)$$

where x_1 and w_1 are the principal eigenvectors of A and $A_m(\mathcal{L})$ corresponding to the largest eigenvalues $\lambda_1(A)$ and $\lambda_1(A_m(\mathcal{L}))$, respectively, and where a link l joins the nodes l^+ and l^- .

Proof. [15] □

The decrease of the largest eigenvalue $\lambda_1(A) - \lambda_1(A_m(\mathcal{L}))$ tends to be larger if the upper bound $2 \sum_{l \in \mathcal{L}_m} (x_1)_{l^+} (x_1)_{l^-}$ is larger. This motivates the principal eigenvector strategy to minimize the largest eigenvalue: removing the set of links that maximizes $2 \sum_{l \in \mathcal{L}_m} (x_1)_{l^+} (x_1)_{l^-}$. Moreover, when only one link is removed, removing the link with the maximum $(x_1)_{l^+} (x_1)_{l^-}$, maximizes not only the upper bound of (1), but likely the lower bound as well, since w_1 is close to x_1 in this case. This eigenvector strategy performs almost optimally in this situation.

Theorem 2. *For any graph G and graph $G_m(\mathcal{N}) = G \setminus \mathcal{N}_m$, by removing m nodes from G , it holds that*

$$0 \leq \lambda_1(A) - \lambda_1(A_m(\mathcal{N})) \leq 2 \sum_{n \in \mathcal{N}_m} (x_1)_n^2 \lambda_1(A) - \sum_{j \in \mathcal{N}_m} \sum_{i \in \mathcal{N}_m} a_{ij} (x_1)_i (x_1)_j \quad (2)$$

where x_1 is the principal eigenvectors of A corresponding to the largest eigenvalues $\lambda_1(A)$. In particular, if $m = 1$, then

$$0 \leq \lambda_1(A) - \lambda_1(A_1(\mathcal{N})) \leq 2 (x_1)_n^2 \lambda_1(A) \quad (3)$$

where n is the node removed.

Proof. [5] □

Theorem 2 implies that the decrease of spectral radius by removing a node or a set of nodes is strongly related to the principal eigenvector components corresponding to the removed nodes. Motivated by Theorem 2, the eigenvector based one node removal strategy to minimize the largest eigenvalue simply removes the node with the largest principal eigenvector component $(x_1)_n$.

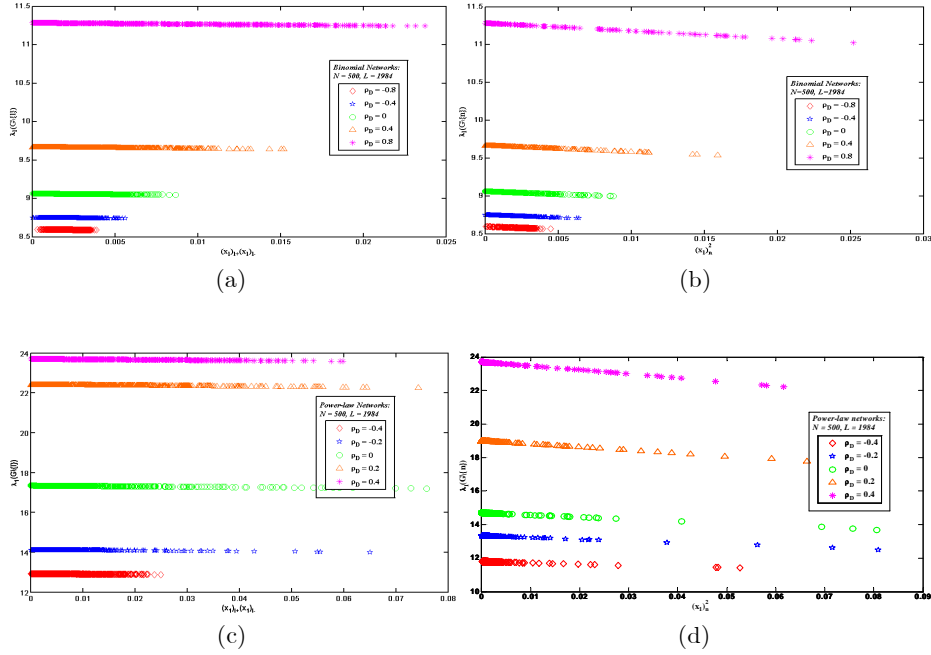


Fig. 1. The spectral radius of graphs by removing a link (or node) as a function of corresponding components in principal eigenvector (a), (b) in Binomial graphs, (c), (d) in power-law graphs.

We perform further simulations to illustrate the importance of the principal eigenvector components in characterizing the influence of the link/node removal on λ_1 . We deduce networks with different assortativities but with a given degree vector, which may follow a binomial or power-law degree distribution. Upon each network, we try all possible one link (or node) removal and examine the largest eigenvalue $\lambda_1(G \setminus (l))$ (or $\lambda_1(G \setminus (n))$) after removing one link (or node) as a function of $(x_1)_{l+}$ $(x_1)_{l-}$ (or $(x_1)_n^2$) corresponding to the link (or node) removed. By the Perron-Frobenius theorem [14], all components of x_1 and w_1 are non-negative (positive if the corresponding graph is connected). Interestingly, $\lambda_1(G \setminus (l))$ (or $\lambda_1(G \setminus (n))$) decreases linearly as a function of increasing $(x_1)_{l+}$ $(x_1)_{l-}$ (or $(x_1)_n^2$),

as shown in Fig. 1. In other words, the spectral radius will be decreased more if the link (or node) removed has a larger $(x_1)_{l+}$ $(x_1)_{l-}$ (or $(x_1)_n^2$).

3 Relation between the principal eigenvector and the degree vector

In view of the importance of the principal eigenvector in characterizing the influence of link/node on the spectral radius, in this section, we explore how the average $E[x_1]$ as well the variance of x_1 changes with the assortativity ρ_D when the degree vector, which may follow the degree distribution of network models or of real-world networks, remains the same. Moreover, we explore the difference and the linear correlation coefficient between the principal eigenvector and the degree vector, the simplest and mostly studies network metric, which as well provides important insights on under which condition the degree vector/sequence well approximates the principal eigenvector.

3.1 Properties of the principal eigenvector

Two types of degree distributions have been so far widely studied: the binomial and power-law degree distribution. The binomial degree distribution is a characteristic of an Erdős-Rényi random graph $G_p(N)$, which has N nodes and any two nodes are connected independently with a probability p . Such a random construction leads to a zero assortativity as proved in [13]. However, the class of graphs $G(N, p)$ with the same binomial degree distribution $\Pr[D_G = k] = \binom{N-1}{k} p^k (1-p)^{N-1-k}$ as Erdős-Rényi random graphs $G_p(N)$ and obtained, for instance, by degree-preserving rewiring feature an assortativity that may vary within a wide range. The power-law degree distribution $\Pr[D = k] = ck^{-\alpha}$, where $c = 1/\sum_{k=1}^{N-1} k^{-\alpha}$ has been widely observed in real-world networks. Similarly, graphs with a given power-law degree distribution, for example, generated by the Barabási-Albert power model [1] can be altered by the degree-preserving rewiring to obtain different assortativity.

We explore the principal eigenvector components (see Figure 2) as well as its average $E[x_1]$ (see Figure 3) in graphs with the same degree distribution (i.e. binomial or power-law) but with different assortativities ρ_D obtained by degree-preserving rewiring. Figure 2 shows that the variance of the principal eigenvector increases with assortativity ρ_D . Furthermore, as shown in Figure 3, $E[x_1]$ decreases with the increase of assortativity ρ_D . Similarly, we consider a set of 11 real-world networks. We apply degree-preserving rewiring to each real-world network to derive network instances with different assortativity. In other words, we derive a class of networks that possess the same degree distribution as a real-world network but different assortativities. Interestingly, we observe the same, $E[x_1]$ decreases with increasing assortativity (see Figure 3(b)).

The decrease of $E[x_1]$ and the increase of the variance of the principal eigenvector components with increasing assortativity can be qualitatively explained

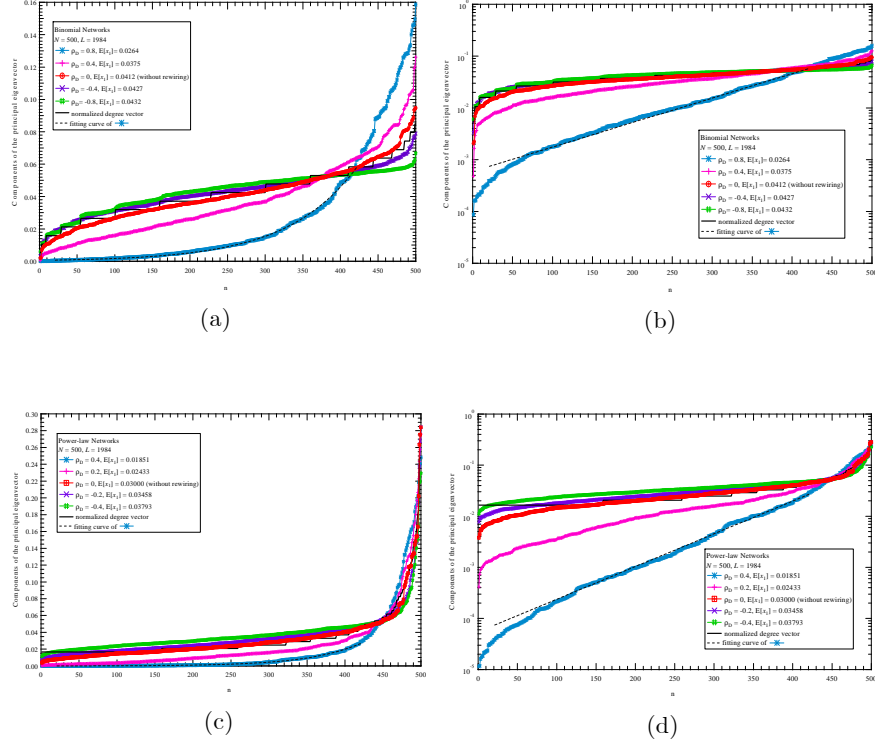


Fig. 2. The components of the principal eigenvector in increasing order. Images (a) (linear) (b) (semilogarithmic) are binomial graphs with different assortativity. Images (c) (linear) (d) (semilogarithmic) are power-law graphs with different assortativity.

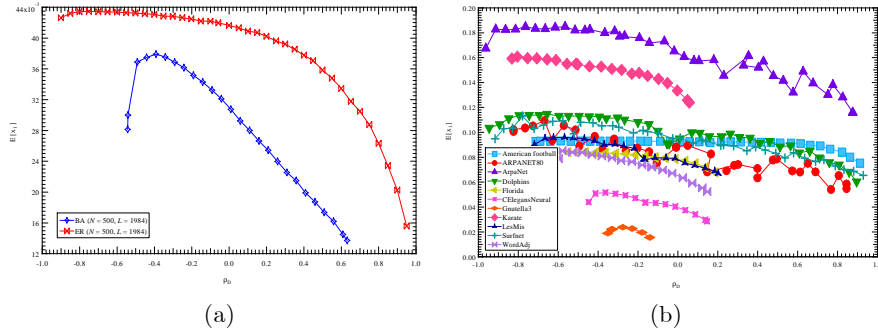


Fig. 3. The average of the components of the principal eigenvector versus the assortativity. (a) in binomial and power-law graphs (b) in network instances derived from real-world networks via degree-preserving rewirings.

as follows. As defined, the principal eigenvector x_1 corresponds to the largest eigenvalue λ_1 follows

$$\lambda_1(x_1)_j = \sum_{q=1}^N a_{jq}(x_1)_q, \quad (4)$$

where $a_{jq} = 1$ if q is a neighbor of node j , or else $a_{jq} = 0$. The j -th component of the principal eigenvector $(x_1)_j$ tends to be large if node j has a large degree (number of neighbors) or if the components corresponding to its neighbors are large. When ρ_D is large, high degree nodes prefer to link with other high degree nodes. In this case, a high degree node possesses a large number of neighbors, whose corresponding eigenvector components are again likely to be large, whereas a low degree node connects to a small number of neighbors, whose corresponding components tend to be small. Both a large variance in degree and a large assortativity ρ_D contribute to a large variance $Var[x_1]$ of the principal eigenvector x_1 . This explains why variance $Var[x_1]$ of x_1 increases with ρ_D and with a given assortativity, the power-law graphs have a larger $Var[x_1]$ than the binomial graphs (see Figure 2). Furthermore, since $Var[X] = E[X^2] - (E[X])^2$ and $x_1^T x_1 = 1$,

$$E[x_1] = \sqrt{\frac{1}{N} - Var[x_1]}, \quad (5)$$

Correspondingly, both a large variance in degree and a large assortativity ρ_D contribute to a small $E[x_1]$ of the principal eigenvector x_1 . Hence, $E[x_1]$ decreases with increasing ρ_D and tends to be smaller when the degree variance is larger. Moreover, considering Eq. (5), we can deduce the upper bound $E[x_1] \leq \frac{1}{\sqrt{N}}$.

Figure 2 compares as well the principal eigenvector x_1 with the normalized degree vector $\bar{d} = \frac{d}{\sqrt{d^T d}}$ in binomial graphs and power-law graphs ($N = 500$, $L = 1984$) with different assortativities. The components of x_1 and \bar{d} are plotted in the order of increasing magnitude. The difference between x_1 and \bar{d} is affected by ρ_D , which will be further explored in the following part.

3.2 Relation between degree vector and principal eigenvector

In this section, we investigate the relation between the principal eigenvector and the degree vector by their difference and linear correlation coefficient. The degree vector has to be first normalized to quantify its difference with the principal eigenvector. We propose two scalings of the degree vector $\bar{d} = \frac{d}{\sqrt{d^T d}}$ and $\tilde{d} = \frac{\alpha}{\lambda_1} d$, where α is a constant. The corresponding *difference vector* between x_1 and the scaled degree vector is $w = x_1 - \frac{d}{\sqrt{d^T d}}$ and $y = x_1 - \frac{\alpha}{\lambda_1} d$, respectively. The overall difference can be quantified by either the relative difference $u^T w$ (or $u^T y$) or the absolute difference $w^T w$ (or $y^T y$), actually, the square sum or the sum of the components in the difference vector respectively. The first scaling of the degree vector $\bar{d} = \frac{d}{\sqrt{d^T d}}$ aims to obtain the same norm for the degree vector and the principal eigenvector: $\sqrt{d^T d} = \sqrt{x_1^T x_1} = 1$. The other $\tilde{d} = \frac{\alpha}{\lambda_1} d$ is motivated by $(x_1)_j = \frac{1}{\lambda_1} \sum_{r=1}^N a_{jr} (x_1)_r \leq \frac{d_j}{\lambda_1}$ and the constant α is determined (see Theorem

3) as the one minimize the absolute difference $y^T y$. Note that both linear scalings of the degree vector will not change the linear correlation coefficient between the principal eigenvector and the degree vector.

Theorem 3. *The absolute difference $w^T w$ (or $y^T y$) between the principal eigenvector and the degree vector is the smallest ($w^T w = 0$ or $y^T y = 0$) when the spectral radius follows $\lambda_1 = \frac{N_2}{N_1}$, where N_k is the total number of k hop walks between any two nodes which can be the same.*

Proof. The absolute difference

$$w^T w = \left(x_1 - \frac{d}{\sqrt{d^T d}}\right)^T \left(x_1 - \frac{d}{\sqrt{d^T d}}\right) = x_1^T x_1 - 2 \frac{d^T x_1}{\sqrt{d^T d}} + \frac{d^T d}{\left(\sqrt{d^T d}\right)^2} = 2 - 2 \frac{d^T x_1}{\sqrt{d^T d}}. \quad (6)$$

Moreover, the generalized form of (4) for the k -th largest eigenvalue λ_k and the corresponding eigenvector x_k follow $(x_k)_j = \frac{1}{\lambda_k} \sum_{r=1}^N a_{jr} (x_k)_r = \alpha \frac{d_j}{\lambda_k} - \frac{1}{\lambda_k} \sum_{r=1}^N a_{jr} (\alpha - (x_k)_r)$, we will determine α so that $y_k = x_k - \frac{\alpha}{\lambda_k} d$ has minimum norm. Hence,

$$y_k^T y_k = \left(x_k - \frac{\alpha}{\lambda_k} d\right)^T \left(x_k - \frac{\alpha}{\lambda_k} d\right) = 1 - 2 \frac{\alpha}{\lambda_k} d^T x_k + \frac{\alpha^2}{\lambda_k^2} d^T d, \quad (7)$$

is minimized with respect to α if $-\frac{2}{\lambda_k} d^T x_k + 2 \frac{\alpha}{\lambda_k^2} d^T d = 0$ or $\frac{\alpha}{\lambda_k} = \frac{d^T x_k}{d^T d}$. Let $y = y_1$, we obtain

$$y^T y = 1 - \frac{(d^T x_1)^2}{d^T d}, \quad (8)$$

using the α derived in the last step. In both Eq. (6) and Eq. (8), $w^T w = 0$ and $y^T y = 0$ if $d^T x_1 = \sqrt{d^T d}$. In other words, when the principal eigenvector is proportion to degree vector, $w = \mathbf{0}$ (or $y = \mathbf{0}$). Since $Ax_1 = \lambda_1 x_1$, $d^T x_k = \lambda_1 u^T x_1$. The condition $d^T x_1 = \sqrt{d^T d}$ implies

$$\lambda_1 u^T x_1 = \sqrt{d^T d} = \sqrt{N_2}$$

where $N_2 = d^T d$. Since $x_1 = \frac{d}{\sqrt{d^T d}}$, and $u^T d = N_1$, Lemma 3 follows. \square

Notice that in some approximate mean-field models for virus spreading [10], $\tau_c \sim \frac{N_1}{N_2} = \frac{1}{\lambda_1}$. Furthermore, $w^T w = 0$ (or $y^T y = 0$) is a special case of $u^T y_k = 0$, when $\lambda_1 = \frac{N_2}{N_1}$.

The relative difference $w^T u = u^T x_1 - \frac{d^T u}{\sqrt{d^T d}}$ ($y^T u$) is zero when the absolute difference is zero. We explore the relative difference in general cases by considering the binomial graphs as an example. The sum of the principal eigenvector $u^T x_1$ and the relative difference $w^T u$ as a function of the assortativity are shown in Figure 4 to follow exactly the same trend, since the degree of each node, thus, $\frac{d^T u}{\sqrt{d^T d}}$ remains the same when we change the assortativity by degree-preserving

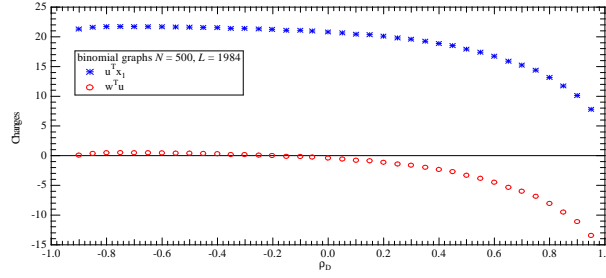


Fig. 4. The difference between the principal eigenvector and degree vector as a function of the assortativity ρ_D .

rewiring. When the assortativity $\rho_D = 0$, the binomial graphs are actually Erdős-Rényi random graphs, for which $\lambda_1 \simeq \frac{N_2}{N_1}$ when the network size is large [14]. Hence, both the absolute and relative difference are zero when the assortativity is around zero. The sum of the principal eigenvector $u^T x_1$ decreases with the assortativity ρ_D , as explained in Subsection 3.1.

3.3 Correlation between the principal eigenvector and the degree vector

Recall that so far the best strategy to minimize the spectral radius by links/nodes removal is based on the principal eigenvector. When the correlation $\rho(x_1, d)$ between the principal eigenvector and the degree vector is positively strong, we may use the degree vector instead of the principal eigenvector to determine which links/nodes to remove, which will be further illustrated in Section 4. Here, we investigate the linear correlation coefficient $\rho(x_1, d)$ between the principal eigenvector and the degree vector as a function of ρ_D . Linear scaling of the degree vector will not change the linear correlation coefficient. Hence, we consider the original degree vector. When the absolute difference between the principal eigenvector and the scaled degree vector is zero, the principal eigenvector is proportion to degree vector. In this case, $\rho(x_1, d) = 1$, which seldom occurs in real-world networks. A strong positive correlation, not necessarily to be one, is already interesting with respect to approximate the eigenvector strategy by the corresponding degree vector strategy in minimizing the spectral radius.

Figure 5(a) depicts that $\rho(x_1, d)$ is mostly positively strong in the Erdős-Rényi random graphs and Barabasi-Albert graphs. However, $\rho(x_1, d)$ decreases dramatically when the assortativity is decreased, actually around the minimal assortativity. Similarly, we derive networks with different assortativities by applying degree preserving rewiring to each of the 11 real-world networks. As in Figure 5(b), We are interested in how $\rho(x_1, d)$ changes with the assortativity ρ_D in real-world networks. Figure 5(b) illustrates that, the correlation $\rho(x_1, d)$ creases as the assortativity is decreased, especially around the minimal assortativity, which is the same as observed in network models. In the simulations

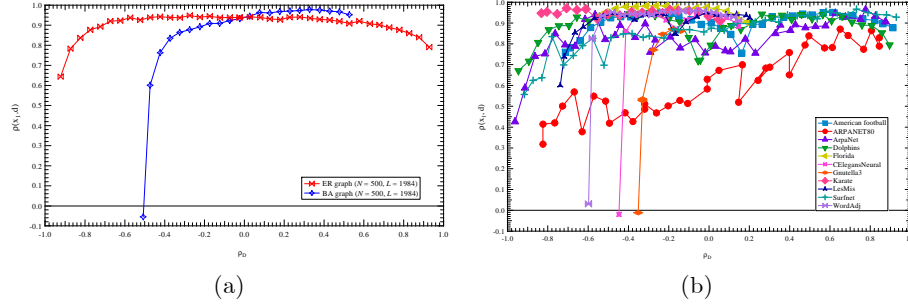


Fig. 5. The linear correlation coefficient between the degree vector and the principal eigenvector as a function of the assortativity (a) in both binomial graphs (red marks and line) and power-law graphs (blue marks and line); (b) in network instances derived from real-world networks via degree-preserving rewirings.

of both network models and real-world networks, the most evident decrease is observed in networks with a power-law degree distribution such as the C. elegans neural network, the Gnutella 3 network and the WordAdj network.

These observations can be explained similarly as we explain the average/variance of the principal eigenvector versus assortativity in Section 3.1. In general, if a node has a large degree, its corresponding principal eigenvector component tends to be large even when the assortativity is zero, due to (4). A large positive assortativity implying large (or small) degree nodes tend to connect to other large (or small) degree nodes, further enforces a large degree node to have more likely a even larger principal eigenvector component compared to a small assortativity. Hence, a negative assortativity will weaken the correlation $\rho(x_1, d)$. Note that the correlation coefficient is not necessarily the maximum at the maximal assortativity as shown in Figure 5, because here we examine the linear correlation coefficient but not the rank correlation.

4 Application: degree *vs.* principal eigenvector strategy in minimizing the spectral radius

In this section, we illustrate the possibility to replace the principal eigenvector strategy by the degree vector in minimizing the spectral radius λ_1 via an example of node removal in power-law networks with different assortativities. As mentioned in Section 2, so far the best node removal strategy removes the node with the largest principal eigenvector component $(x_1)_j$. A widely applied strategy to minimize λ_1 by removing m nodes (a) removes the set of m nodes with the highest component in the principal eigenvector of the original graph. The corresponding degree vector strategy (b) removes the set of m nodes with the highest degree in the original graph. We compare these two strategies in remov-

ing $m \in [1, 200]$ nodes in graphs with positive, zero and negative assortativity (see Fig. 6) but with the same power-law degree distribution as in Fig. 5(a).

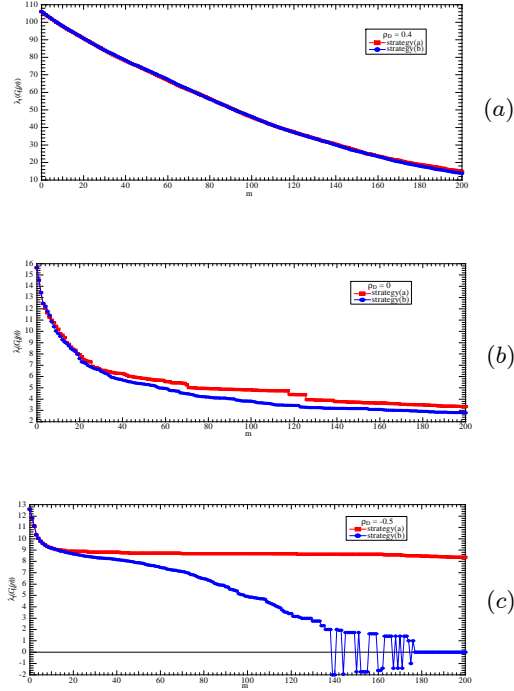


Fig. 6. The decrease of the spectral radius by successively removing m nodes in power-law networks. The square and circle dot dash lines show the decrease of the spectral radius by strategies (a) and (b) separately.

Figure 6 shows that the decreases of λ_1 by removing nodes with strategy (a) and (b) are almost same when ρ_D is large. The eigenvector strategy (a) decreases the spectral radius more thus performs better than the degree vector strategy (b) when the assortativity is small. When the assortativity is large, the degree vector is positively and strongly correlated with the principal eigenvector. In such a case, the degree vector strategy, the simplest to compute, well approximates the principal eigenvector strategy in minimizing the spectral radius.

5 Conclusions

The principal eigenvector is essential in characterizing the influence of link/node on the spectral radius, whereas its topological meaning is far from well understood. This work, via both theoretical analysis and systematic simulations,

contributes to the following aspects: (a) the average $E[x_1]$ (or variance) of the principal eigenvector is shown and explained to decrease (or increase) with the assortativity ρ_D ; (b) the difference between the principal eigenvector and the degree vector is proved to be the smallest, when $\lambda_1 = \frac{N_2}{N_1}$ and (c) we illustrate and explain why the correlation between principal eigenvector and the degree vector decreases as ρ_D is decreased. In general, both a large variance (heterogeneity) in nodal degree and a large degree correlation (homogeneity in connection) contribute to a large average and a small variance of the principal eigenvector and a strong correlation between the degree and the principal eigenvector. As a straightforward application of these finds, we illustrate that when the assortativity is large, we could approximate the well performance principal eigenvector based strategy (to minimize λ_1 by removing links/nodes) by the corresponding degree vector, which is the simplest network property to compute.

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