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EPL, **83** (2008) 68003

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Modular synchronization in complex networks with a gauge Kuramoto model

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received 12 June 2008; accepted in final form 6 August 2008
published online 12 September 2008

PACS 89.75.-k – Complex systems
PACS 89.65.-s – Social and economic systems

Abstract – We modify the Kuramoto model for synchronization on complex networks by introducing a gauge term that depends on the edge betweenness centrality (BC). The gauge term introduces additional phase difference between two vertices from 0 to π as the BC on the edge between them increases from the minimum to the maximum in the network. When the network has a modular structure, the model generates the phase synchronization within each module, however, not over the entire system. Based on this feature, we can distinguish modules in complex networks, with relatively little computational time of $\mathcal{O}(NL)$, where N and L are the number of vertices and edges in the system, respectively. We also examine the synchronization of the modified Kuramoto model and compare it with that of the original Kuramoto model in several complex networks.

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Complex networks have drawn considerable attention from diverse disciplines such as sociology, information science, physics, biology and so on [1]. Many complex networks in real world contain modules within them, which form in a self-organized way to achieve the efficiency functionally or regionally. Such modular systems can exhibit collective synchronized patterns within each module, not forming the global synchronization [2] as can be found in the cortex of neural network [3] or different synchronization transition behaviors depending on the patterns of inter-module connections [4].

In this letter, we study the modular synchronization pattern generated from a modified Kuramoto equation (KE), which we call the gauge KE,

$$\frac{d\phi_i(t)}{dt} = \Omega_i - J \sum_{j=1}^N a_{ij} \sin(\phi_i(t) - \phi_j(t) - \eta g(b_{ij})). \quad (1)$$

Here, ϕ_i is the phase of vertex i , Ω_i is the natural frequency of vertex i selected from the Gaussian distribution $e^{-\Omega^2/2}/\sqrt{2\pi}$, J is the overall coupling constant and a_{ij} is the (i, j) -th component of the adjacency matrix, which is one when the vertices i and j are connected, and

zero otherwise. η is a control parameter. The extra phase term $g(b_{ij})$, we call the gauge term below, is defined as

$$g(b_{ij}) = \frac{b_{ij} - b_{\min}}{b_{\max} - b_{\min}} \pi, \quad (2)$$

where b_{\min} and b_{\max} are the minimum and the maximum edge betweenness centrality (BC) [5] or load [6], respectively, in the system. Here, the edge BC or load is the amount of effective traffic passing through a given edge when every pair of vertices sends and receives a unit packet that travels along the shortest path between them. Then the gauge term $g(b_{ij})$ is in the range from 0 to π depending on the BC of edge. When $\eta = 0$, the gauge KE recovers the standard KE [7] which becomes fully synchronized when J is sufficiently large. The KE with the extra phase of the form $\sin(\phi_i - \phi_j - c)$ ($c = \text{constant}$) was studied first in [8]. The effect of the extra phase is to destroy the synchronization. Intuitively, one expect that the BCs on intra-module links are smaller than those on inter-module. Thus, each module can be synchronized, while the entire system is not. Moreover, the gauge term induces an effective coupling that can be negative at the edges connecting different modules. Due to this negative coupling, the average phase of each module may have velocity different from

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each other. Using this property, the gauge KE can be used for module identification in complex networks.

The module identification in the context of synchronization has been studied [9,10]. These studies are inspired by the so-called dynamic clustering (DC) approach that individual oscillators have different levels of synchronization time owing to the heterogeneity of degree in network. Since vertices within modules are densely connected, they are synchronized more earlier than those between modules. Using this idea, the hierarchical structure can be detected by monitoring the temporal evolution of synchronization [9]. To identify the modules, however, the information of characteristic time at each hierarchical level is needed, which may be obtained from the spectrum of the Laplacian matrix of the system. Boccaletti *et al.* [10] introduced another model, in which the coupling strength of the KE depends on the BC as $b_{ij}^{\alpha(t)}$, where $\alpha(t)$ is negative. Thus, the coupling strengths across the module-connecting edges are weaker than those within module. $\alpha(t)$ is then tuned to detect the modules. In both methods, one needs to control the parameters such as time and $\alpha(t)$. However, our method based on eq. (1) with $\eta = 1$ does not contain any control parameter, so that we can identify the modules without any prerequisite information.

We begin to study the synchronization pattern generated from eq. (1). Firstly, we apply the gauge KE to an *ad hoc* network [11] with a modular structure. The network is composed of $N = 128$ vertices and $L = 1024$ edges. Those vertices are grouped to four modules, each of which is of equal size. And edges are connected with probability p_{in} for pairs of nodes belonging to the same module whereas pairs belonging to different modules have edges with probability p_{out} . By controlling the parameter p_{in} and p_{out} we can obtain a fraction of inter-modular edges, $z_{\text{out}}/\langle k \rangle$ as we want, where z_{out} is the mean degree of inter-modular edges and $\langle k \rangle = 2L/N$ is the mean degree. This *ad hoc* network has been used as a benchmark for module identification algorithms in previous studies [12].

We measure the order parameter defined as

$$\mathcal{M}_{\text{tot}} \equiv \left\langle \left| \frac{1}{N} \sum_{j=1}^N e^{i\phi_j} \right| \right\rangle, \quad (3)$$

where $\langle \dots \rangle$ denotes the time and ensemble average. The order parameter is measured in the steady state. When $\eta = 0$, the order parameter saturates to 1 for large J , however, as η is increased toward 1, it saturates at lower values as shown in fig. 1(a). This behavior indicates that the network is not synchronized globally. To check if the synchronization forms within each module, the local order parameter, defined as $\mathcal{M}_\alpha \equiv \langle |\sum_{j=1}^{N_\alpha} e^{i\phi_j} / N_\alpha| \rangle$, is measured, where α is the module index, N_α is the number of vertices within the module α and the sum is over vertices within the module. We find that indeed the order parameter \mathcal{M}_{mod} reaches 1 for large J as shown in fig. 1(b), indicating that the oscillators within the module are synchronized. We examine the average phase of each

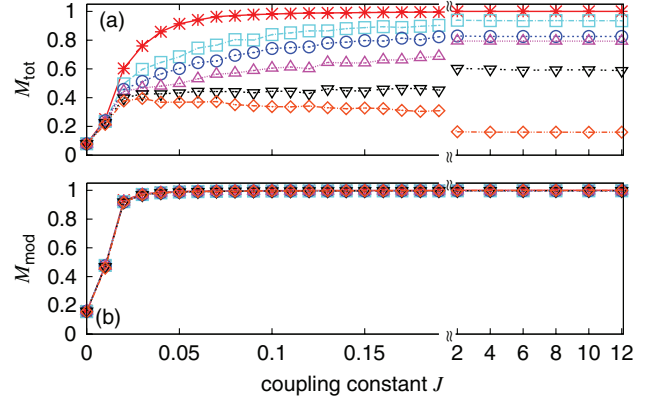


Fig. 1: (Color online) The order parameter defined over the entire network (a) and within a module (b) vs. the coupling constant J for the *ad hoc* network in case of $z_{\text{out}}/\langle k \rangle = 0.05$. Data are for $\eta = 0.0, 0.6, 0.7, 0.8, 0.9$ and 1.0 from the top in (a). The same symbols are used for (b), but data for different η collapse onto the single curve.

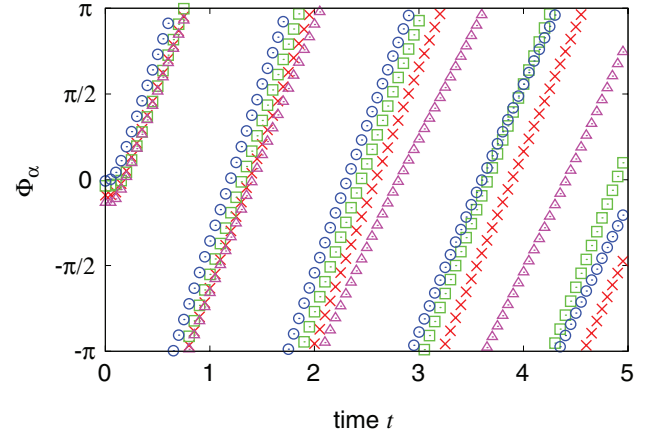


Fig. 2: (Color online) The time evolution of average phases of the four modules, distinguished by different symbols, for the *ad hoc* network with $z_{\text{out}}/\langle k \rangle = 0.05$ when $\eta = 1.0$ and $J = 2.0$.

module as a function of time. As shown in fig. 2, the modules are distinguishable by different average phases and average phase velocities.

The stability of synchronization of the model (1) is examined. Assuming the fully synchronized state of the form $\phi_i^* = \phi_i^0 + \Omega t$, and linearizing eq. (1), we get $\dot{\xi}_i(t) = -J \sum_j G_{ij} \xi_j(t)$, where $\xi_i(t) = \phi_i(t) - \phi_i^*$, $G_{ij} = (\sum_k a_{ik} \omega_{ik}) \delta_{ij} - a_{ij} \omega_{ij}$ and $\omega_{ij} = \cos(\phi_i^0 - \phi_j^0 - \eta g(b_{ij}))$. $\lambda_1 = 0$ is the trivial eigenvalue of G and the sign of other eigenvalues determines the stability of the fully synchronized state. Due to the negative element of the coupling matrix G , its eigenvalues can be negative, and then the Lyapunov exponent in the linear stability analysis can be as well. In that case, the synchronization is no longer stable. We obtain ω_{ij} from $\cos(\phi_i(t) - \phi_j(t) - \eta g(b_{ij}))$ at an arbitrary but sufficiently large t and trace out the eigenvalues for the *ad hoc* network having $z_{\text{out}}/\langle k \rangle = 0.05$

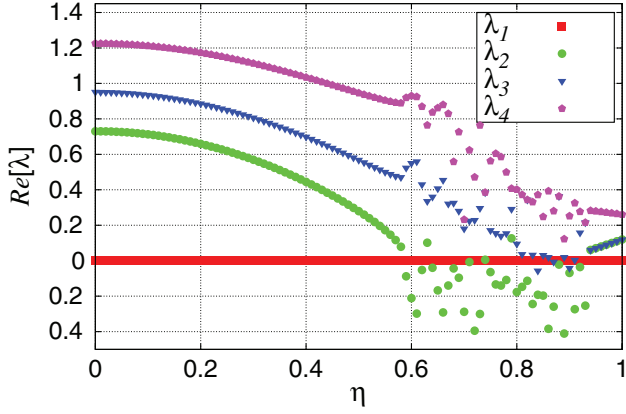


Fig. 3: (Color online) The first 4 eigenvalues, $\lambda_1 = 0$, λ_2 , λ_3 and λ_4 , of G_{ij} vs. the parameter η for the *ad hoc* network in case of $z_{\text{out}}/\langle k \rangle = 0.05$ and $J = 2.0$. Data beyond $\eta_c \approx 0.59$ depend sensitively on time t where ω_{ij} is obtained.

and plot the first 3 non-zero eigenvalues vs. η in fig. 3. λ_2 is positive at $\eta = 0$ and decreases to zero as η increases from 0 to $\eta_c \approx 0.59$. And increasing η further above η_c drives the system to unstable state. For $0 \leq \eta < \eta_c$, the order parameter \mathcal{M}_{tot} is almost 1 in the steady state, whereas \mathcal{M}_{tot} has a smaller constant value for $\eta > \eta_c$. In many cases, they actually oscillates in time before the time average due to disparate group velocities of the modules as shown in fig. 2. The curve fitting of λ_2 in the vicinity of $\eta = \eta_c$ shows $\lambda_2 \propto (\eta_c - \eta)^{1/2}$. The square-root singularity of λ_2 near the stability edge is the signature of the saddle-node bifurcation [13].

We introduce how to identify modules with the gauge KE. To this end, we take the following steps:

- i) We apply the gauge KE (1) to all oscillators with a sufficiently large coupling constant J . The phases $\{\phi_i(t)\}$ of each oscillator are obtained in the steady state.
- ii) We measure the phase similarity defined as $C_{ij} = \langle [1 + \cos(\phi_i(t) - \phi_j(t))]/2 \rangle$ for each connected pair of oscillators (i, j) . The brackets are the average over different times, natural frequencies $\{\Omega_i\}$, and initial random phases $\{\phi_i(0)\}$.
- iii) From the empty state, where all edges are absent, we add edges (i, j) one by one that are chosen following the descending order of C_{ij} . Clusters after the step iii) are regarded as modules. The edges that existed originally, but not connected yet until the step iii) are regarded as inter-modular edges.
- iv) We repeat the step iii) until the modularity of the system becomes maximum. The modularity Q is defined as

$$Q = \sum_{\alpha} e_{\alpha\alpha} - a_{\alpha}^2, \quad (4)$$

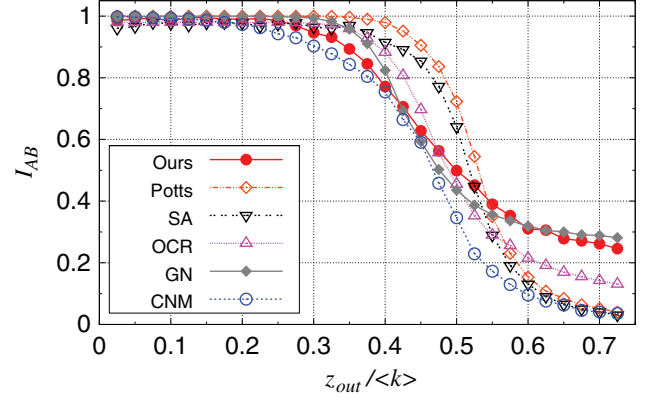


Fig. 4: (Color online) The mutual information vs. $z_{\text{out}}/\langle k \rangle$, the fraction of inter-modular edges per mean degree for the *ad hoc* network. See the text for abbreviations.

where $a_{\alpha} = \sum_{\beta} e_{\alpha\beta}$, and $e_{\alpha\beta}$ is the fraction of edges that connect the vertices belonging to the modules α and β [11].

To test the performance of our algorithm, we measure the mutual information on several networks, defined as

$$I(A, B) = \frac{-2 \sum_{i=1}^M \sum_{j=1}^{M'} \log \left(\frac{N_i^j}{N_i N^j} \right)}{\sum_{i=1}^M N_i \log \left(\frac{N_i}{N} \right) + \sum_{j=1}^{M'} N^j \log \left(\frac{N^j}{N} \right)}, \quad (5)$$

where $M = 4$ is the number of preassigned modules and M' is the number of detected modules. N_i^j is the number of vertices belonging to the i -th preassigned and the j -th detected modules, $N_i = \sum_j N_i^j$ and $N^j = \sum_i N_i^j$ [12].

Figure 4 shows the mutual information measured on the *ad hoc* network as a function of $z_{\text{out}}/\langle k \rangle$ for several module-detecting algorithms. The preliminary result of this comparison was presented in a proceeding paper [14]. The performance of our algorithm is not better than those of the Potts model and the simulated annealing (SA) [15,16]. Even though they are better in performance, if we count for their long computation time, then ours may be useful practically. The performance of opinion-changing rate model (OCR) algorithms [10] is somewhat better, however, it requires an extra task of parameter tuning, so that ours is easier to implement. Since our algorithm shares with the Girvan-Newman (GN) algorithm [17] the idea of clustering based on BC, the performances of the two algorithms are close to each other. However, since ours calculates the BC on each edge only once, whereas the GN algorithm does it repeatedly for each disconnected cluster, the computational time can be reduced drastically from $\mathcal{O}(NL^2)$ to $\mathcal{O}(NL)$. The performance of our algorithm is better than that of the Clauset-Newman-Moore (CNM) algorithm [18], which runs in $\mathcal{O}(N \ln^2 N)$ for sparse graphs.

Secondly, we apply our algorithm to the hierarchical network proposed by Ravasz and Barabási [19]. When the number of levels is two, the modules are well selected in a

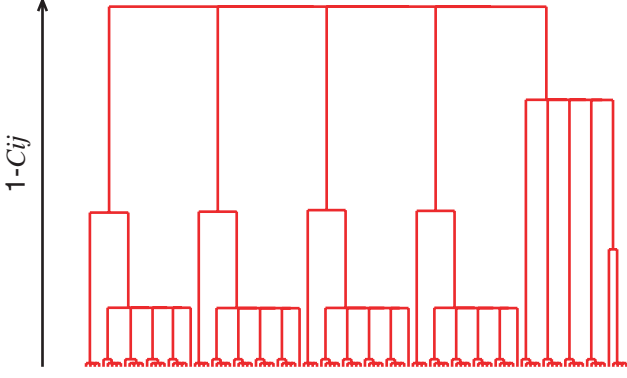


Fig. 5: (Color online) The dendrogram based on the phase similarity between connected pairs of vertices for the hierarchical network with three levels.

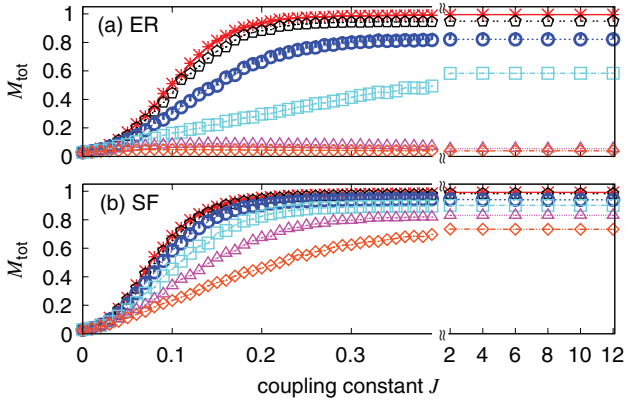


Fig. 6: (Color online) The order parameter *vs.* the coupling constant *J* for the ER (a) and the SF network with the degree exponent 3.5 (b). The data are for the cases of $\eta = 0, 0.2, 0.4, 0.6, 0.8$ and 1.0 from the top.

similar way as in fig. 3 of ref. [9]. For the three level case, the dendrogram constructed by our method is shown in fig. 5. Here, the hub at the second level is grouped with one of the four identical modules connected to it in the second level.

Thirdly, we apply the gauge KE to Erdős-Rényi (ER) random networks and scale-free (SF) networks with no modular structure to see the dependence of the network structure. The SF network is generated using the static model with degree exponent 3.5 [6]. The order parameter (3) behaves differently for the two networks. For the ER network, the saturated value of the order parameter decreases from 1 to 0 as η increases from 0 to 1 (fig. 6(a)). However, for the SF network, the order parameter does not decrease to 0, but ≈ 0.7 even if η reaches 1 (fig. 6(b)).

The mechanism of this difference is complex. To understand it, we measure the phase difference $\Delta\phi$ across an edge connected to the hub for the case $\eta = 1$. Typical patterns are as follows: For the ER network, the phase difference changes with time running from $-\pi$ to π as shown in fig. 7(a). For the SF network, its pattern shows two different patterns depending on the degree of the other

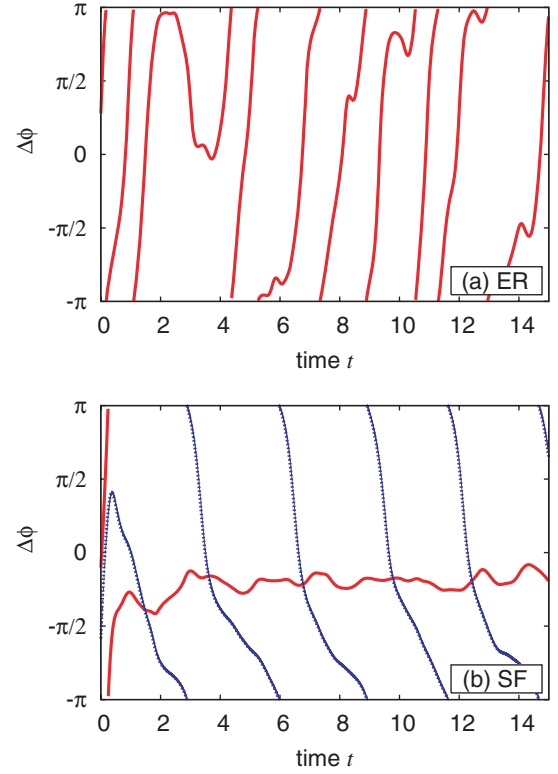


Fig. 7: (Color online) Typical patterns of the phase difference across an edge connected to the hub for the ER network (a) and the SF network (b). In (b), the solid (dotted) curve is likely to occur across an edge between the hub and a node with large (small) degree.

node. It either oscillates as for the ER network (dotted curve in fig. 7(b)) or stays around a smaller value in short intervals as shown in fig. 7(b) (solid curve). As a consequence of this partially oscillating behavior of phase difference for the case of SF network, the order parameter can remain finite as $\mathcal{M}_{\text{tot}} \approx 0.7$. We find that the different patterns in $\Delta\phi$ for the ER and the SF network are closely related to the fraction of edges that have the gauge term larger than $\pi/2$. This fraction is about 15% and 1% for the ER network and SF network, respectively. Note that the BC distribution follows an exponential function (a power law) for the ER (SF) network [6], so that the fraction of edges with BC larger than $\pi/2$ is larger for the ER network than for the SF network. When the gauge term is larger than $\pi/2$, the edge acts as an anti-ferromagnetic coupling between the two nodes. The anti-ferromagnetic coupling induces a frustration in synchronization. As the fraction of the anti-ferromagnetic coupling increases, the destruction of phase and frequency synchronization is more likely to occur. Thus, the feature in figs. 6 and 7 can emerge.

In summary, we have introduced a gauge KE in which the gauge term depends on the edge BC. The gauge term drives the phase difference between the two vertices of an edge from 0 to π as the BC across the edge increases. As a result, the phase difference of two oscillators belonging

to different modules is large, however, it is small across the edges within modules. Thus, the model generates the phase synchronization within each module, however, it does not globally. Measuring the phase similarity between two connected oscillators, we constructed the dendrogram and identified the modules. Such module-detecting method works efficiently.

This work was supported by KOSEF grant Acceleration Research (CNRC) (No. R17-2007-073-01001-0) in SNU and the Korea Research Foundation Grant funded by the Korean Government (MOEHRD, Basic Research Promotion Fund) (KRF-2007-355-C00030) in KIST.

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