



Explosive transitions in complex networks' structure and dynamics: Percolation and synchronization



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ABSTRACT

Percolation and synchronization are two phase transitions that have been extensively studied since already long ago. A classic result is that, in the vast majority of cases, these transitions are of the second-order type, i.e. continuous and reversible. Recently, however, explosive phenomena have been reported in complex networks' structure and dynamics, which rather remind first-order (discontinuous and irreversible) transitions. Explosive percolation, which was discovered in 2009, corresponds to an abrupt change in the network's structure, and explosive synchronization (which is concerned, instead, with the abrupt emergence of a collective state in the networks' dynamics) was studied as early as the first models of globally coupled phase oscillators were taken into consideration. The two phenomena have stimulated investigations and debates, attracting attention in many relevant fields. So far, various substantial contributions and progresses (including experimental verifications) have been made, which have provided insights on what structural and dynamical properties are needed for inducing such abrupt transformations, as well as have greatly enhanced our understanding of phase transitions in networked systems. Our intention is to offer here a monographic review on the main-stream literature, with the twofold aim of summarizing the existing results and pointing out possible directions for future research.

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1. Introduction

1.1. Phase transitions

With the term *phase* of a system physicists actually denote one of its states, characterized by a set of physical properties that can be there considered as uniform over a macroscopic length scale.

Now, from the melting of ice into water to ferro-magnetism and from superconductivity to protein folding, nature is daily exhibiting *phase transitions*, i.e. processes in which systems drastically change one or many of their physical properties when a minimal variation in one of their control variables occurs.

In simpler cases, such transitions are described by the variation of the so called *order parameter* (a scalar, or a vector, or even a tensor measuring the degree of order in the system, whose magnitude normally ranges between zero in one phase and nonzero in the other), as the control variable (or the *control parameter*) is gradually and adiabatically changed.

More precisely, let us exemplify the simplest case of a system having only two phases (say phase 1 and phase 2), and passing from one to another with a process involving scalar control and order parameters. To gather a quantitative description and characterization of the transition, the following procedure is usually run after: (i) the system is prepared in one of its two phases (say phase 1), which actually corresponds to set a specific initial value of the control parameter; (ii) the control parameter is then gradually changed (increased, or decreased, depending on the particular case) in very small steps, up to the point at which the other phase is attained; (iii) the corresponding variations of the order parameter are monitored. This procedure can be followed by starting either from phase 1 or from phase 2, producing the curves that describe the *forward* and *backward* transitions, respectively.

Now, a fundamental concept in statistical mechanics is that of *reversibility* of a given process, which, indeed, can be carried out reversibly or irreversibly. Reversibility can be attributed to processes which, after having taken place from an initial system's phase, can be *fully reversed*, i.e. they can be brought back to their original phase via the exactly opposite transition, and with no changes in either the systems or their surroundings. In our case (as well as it happens in many other circumstances) reversibility consists in a full equivalence of the two curves obtained during the forward and backward transitions. When, instead, transitions are irreversible, the two forward and backward curves differ, and a wealth of interesting and relevant events, such as hysteretic loops, occur.

Given the ubiquitous presence of phase transitions in the real world, it is not strange that physicists started, since the very beginning of statistical mechanics, to delve into discovering universal and critical behaviors (or scaling properties in the proximity of the transition points), as well as to pay effort in classifying phase transitions. The first attempt of providing a classification was given by Ehrenfest [1], but more modern classification schemes consider phase transitions as divided into two broad categories: the first-order and the second-order ones.

In thermodynamics, first-order phase transitions are those that involve a latent heat, i.e. those during which the system either absorbs or releases a typically large amount of energy per volume. Normally, they are characterized (at the transition point) by a “mixed-phase regime”, in which some parts of the system have completed the transition and others have not, this way featuring a sort of *coexistence* of the two system's phases. Signatures of such transitions are the abrupt and discontinuous behavior of the order parameter in the proximity of the transition point, as well as the intrinsic irreversibility of the transition, with the presence (in the majority of the cases) of hysteretic loops. At variance, second-order phase transitions (also called continuous phase transitions) are reversible, and correspond to order parameters displaying a continuous behavior near criticality. Examples of the former processes are the melting of ice, or the boiling of water, or the super-cooling and super-heating transitions. Examples of the latter processes are the ferromagnetic transition, the super-conducting transition, and the super-fluid transition.

1.2. Percolation and synchronization

Our main scope is reviewing recent progresses on two classical phase transitions taking place in networked units (percolation and synchronization), as well as on how they can in fact be rendered *explosive*, i.e. *first-order-like*.

Percolation (from the Latin *percolare*, which means “to filter” or “to trickle through”) is a transition involving a drastic change in the network's structure, due to the emergence of a giant connected component whose size is of the order of

the network's size. By means of percolation, a phase of the network where fragmented portions, or clusters, are isolated is changed into a fully connected structure, where nodes can globally interact. The process can be conducted in two ways: either starting from isolated nodes and adding gradually links (what is called *bond percolation*), or sequentially adding nodes to the graph (what is called *site percolation*).

Synchronization (from the Greek σύγ = together and χρόνος = time) is instead a process having to do with network's dynamics. There, dynamical systems adjust some properties of their trajectories (due to their interactions, or to a driving force) so that the network eventually operates in a collective and macroscopically coherent way.

Percolation and synchronization are nowadays classical fields of study in statistical mechanics, and an interested reader can find orientation within the wealth of currently available results in various monographic papers and books [2–8].

A common result is that the vast majority of transitions to percolation and synchronization are of the second-order type, continuous and reversible. However, as soon as networked units with complex architectures of interaction are taken into consideration, two abrupt and explosive phenomena have been recently reported, namely explosive percolation [9–13] and explosive synchronization [14–22], which rather remind first-order like transitions. Explosive percolation (a sudden materialization of a giant connected component in the structure of the network) was discovered in 2009 [9], while explosive synchronization (an abrupt emergence of a collective, synchronized, behavior in network's dynamics) was first described already in 1984 [23]. In literature, the use of the jargon *first-order like* purports that those explosive processes fulfill some of the properties of a first order transition in thermodynamics (as, for instance, discontinuity and irreversibility) and yet they fail featuring *all* the attributes, qualities and traits of a true first-order transition.

Unveiling the main mechanisms at the origin of these two abrupt transitions is of fundamental relevance for a better understanding of networks' structure and dynamics. For instance, explosive percolation is considered to be at the basis of phenomena such as cascading failures, which, on their turn, may have devastating effects (and cause huge economic losses) in various circumstances: electrical blackouts in power grids, financial crises in the network of global financial market, and the spreading of information and rumors through online social networks, such as Facebook or Twitter. On the other hand, synchronization is of primary importance in sustaining basic brain functions such as emotions, complex thoughts, memory, language comprehension, consciousness, etc., while clinical evidence seems to indicate that explosive synchronization is actually the mechanism underlying the transition from normal to pathological brain behavior during epilepsy, one of the world's most prominent brain disorders.

Furthermore, when synchronization occurs through an irreversible transition, one interesting and ingenious application is the construction of *magnetic-like* states of synchronization in networked oscillators: setting the coupling strength inside the hysteresis region, and entraining the phases of the oscillators through a unidirectional coupling with an external pacemaker, the system can be forced to pass from the unsynchronized to the phase-locked configuration, which remains once the pacemaker is switched off, like the case of a permanent magnet.

Given, therefore, their relevance to practical applications, explosive percolation and synchronization have attracted a comprehensive and lasting attention in various fields, and have stimulated a large amount of exciting works and debates. The result is that many substantial contributions and progresses (including experimental verifications) are today available, which provide a rather deep insight on the crucial structural and dynamical mechanisms at the basis of the abruptness of the two transitions.

It is therefore now the time to offer a comprehensive review on the current state of the art on the subject, which would offer a reflected and thought out viewpoint on the many achievements and developments, as well as a weighted and meditated outlook to the still open questions and to the possible directions for future research.

1.3. Outline of the report

After the present introduction, the review is organized as follows. In Section 2, the main results on explosive percolation are summarized and discussed. Starting from the seminal work by Achlioptas, the Section gives a survey on the various processes that have been proposed in order to produce abruptness in the percolation transition, as well as it critically examines the essential physical mechanisms which are at the origin of the explosiveness. The Section ends with accounting for the debate stirred within the scientific community, and reports some rigorous conclusions that have been ultimately made on the nature of the transition in the thermodynamic limit (i.e. for systems whose size tends to infinity).

Section 3 contains a brief excursus on classical synchronization in both identical and non-identical networked systems. An interested reader can definitely find a much more detailed description of the subject in various other books and review articles, like those of Refs. [4–8], whereas Section 3 concentrates on the main ideas and concepts which are then conveniently recalled at the moment of describing those processes where synchronization emerges, instead, abruptly.

In Section 4, we provide a detailed overview of explosive synchronization. Starting from the early evidences of first order transitions in the thermodynamic limit of models of globally coupled phase oscillators (as well as from the seminal work of Ref. [14]), the Section describes the cases where the transition to synchronization is first-order-like, and discusses the generality of explosive synchronization in networked systems with complex architectures and topologies. The Section, furthermore, gives an account of some recently discovered and unveiled coherent states (the Bellerophon states), whose existence is intimately related to the possibility for the system to feature explosive transition to synchronization.

Section 5 is an exhaustive journey on the applications of explosive percolation and synchronization. In particular, applications of explosive percolation are presented in cascading failures of inter-connected multi-layer networks (of

Table 1

Nomenclature and abbreviations used in the manuscript.

Nomenclature		Abbreviations	
A	Adjacency matrix	2D	Two dimensional
k	Node degree	3D	Three dimensional
$\langle k \rangle$	Network average degree	AP	Achlioptas process
$P(k)$	Degree distribution	BA	Barabási-Albert
\mathcal{A}	Degree-degree correlation	BF	Bohman-Frieze
λ_i	Eigenvalues of the adjacency matrix	BFW	Bohman-Frieze-Wormald
E	Number of edges	CA	Cauuthorship
λ	Coupling strength, synchronization control parameter	CM	Configuration Model
N	System size	CS	Complete Synchronization
θ	Oscillator phase	CPT	Continuous Phase Transition
ϕ	Oscillator phase in the rotating frame	DPT	Discontinuous Phase Transition
v	Frequency	dCDGM	da Costa-Dorogovtsev-Goltsev-Mendes
ω	Natural frequency	DLCA	Diffusion-limited cluster aggregation
$g(\omega)$	Natural frequency distribution	EP	Explosive Percolation
ψ	Average phase	ER	Erdős-Rényi
Ω	Average frequency	ES	Explosive Synchronization
r	Kuramoto order parameter	GAP	Generalized Achlioptas process
S	Percolation order parameter	H-PHN	Human protein homology network
s_{\max}	Size of the largest cluster	LCC	Largest Connected Component
s	Size of a network cluster	MC	Minimum Cluster
t	Time, percolation control parameter ($t = E/N$)	MSF	Master Stability Function
		PR	Product Rule
		PT	Phase Transition
		SCA	Spanning Cluster-Avoiding
		SF	Scale-Free
		SP	Suppression Principle
		SR	Sum Rule
		SW	Small-world
		TR	Triangle Rule
		TW	Traveling Wave

relevance in technology and biology), and applications of explosive synchronization are described in power and smart grids, and in brain dynamics. The Section accounts also for experimental realizations of explosive synchronization in controlled laboratory systems.

Finally, Section 6 presents our conclusive remarks and perspective ideas on yet open problems to solve, and on possible research lines for the future of the subject.

For the reference of the reader, we summarize here below (in Table 1) the list of major symbols and abbreviations which are used all throughout the Manuscript.

2. Explosive percolation

2.1. Introduction and historical overview

Percolation transition refers to the emergence of a giant connected cluster (or giant connected component) in a lattice, or a network, when bonds/sites are gradually occupied at random. For simplicity, throughout this review we always consider bond percolation, unless mentioned otherwise.

In bond percolation one starts from N unconnected nodes. At each time step, an edge is added between two nodes selected according to a given rule. The number of edges E added to the system at a certain time step divided by the system size N is the control parameter $t = E/N$ that describes the phase transition (PT). As for the order parameter (usually denoted as $S(t)$) one can take the fraction of nodes belonging to the giant cluster (the largest cluster) in the network. As time increases, more and more edges are formed in the network, which causes the order parameter to increase from zero. In the thermodynamic limit ($N \rightarrow \infty$), $S(t)$ exhibits a PT from zero to $O(1)$ at a critical point t_c .

Classical percolation is a typical geometrical PT, which has been extensively studied since 1940's in many fields, including mathematics, statistical physics, and engineering. Numerous percolation models, such as invasion percolation, first-passage percolation, directed percolation, bootstrap percolation, k -core percolation have been developed, and it has been shown that PT is closely related to various applications, such as conducting materials, fractality of coastlines, turbulence, magnetic models, colloids, growth models, watersheds of landscapes, the spin quantum Hall transition, etc. The reader is addressed to recent reviews in Refs. [24–27] and references therein for an overview of the main results and concepts about classical percolation.

In modern statistical physics, PTs are traditionally classified into two types, namely, the continuous PT (CPT) and the discontinuous PT (DPT), which are also known as the second-order PT and the first-order PT, respectively. CPT exhibits universal characteristics, and can thus be categorized into different universality classes. At variance, DPT is non-universal.

Table 2

Relevant literature concerning AP and GAP models in lattices, random networks, SF networks, and other networks (real-world, directed, and growing networks), including PR, SR, best-of- m rule, dCDGM, and l -vertex rules. Notice that for the best-of- m rule (m -min-cluster rule), both PR and SR can be used and they are no longer distinguishable.

Model	Lattice ($d = 2$)	Lattice ($d \geq 3$)	Other lattices	Random	SF	Others
Product rule	[12,33–35,37–42]	[42]	[43]	[9,12,13,34,39,44–47]	[34,48,49]	[50,51]
Sum rule	[36,40,42]	[42]	[43]	[9,12,13,47]		
Best-of- m	[53–55]	[55]		[13,46,52,56,57]		[51,58]
dCDGM				[11,39,47,64–67]		[68]
l -vertex				[13,46,59–62,69]		

There are many ingredients that determine whether a phase transition occurs in a system via a CPT or DPT, such as, for instance, dimensionality, the type of the dynamical rules, as well as the underlying network's structures.

For decades, most transitions in classical (or ordinary) percolation have been shown to be of CPT (the second-order) type. One example is the prototypical bond percolation in Erdős–Rényi (ER) networks, for which the size of the giant cluster smoothly increases as the control parameter exceeds the percolation threshold. For $t > t_c$, the order parameter $S(t)$ increases with t as $S(t) \propto (t - t_c)^{1/2}$. However, it should be pointed out that there are also rare examples of percolation models that exhibit DPT, such as the bootstrap percolation [28,29], the k -core percolation [28,30,31], and the jamming percolation [32].

In 2009, a competitive percolation process was introduced, later known as the Achlioptas process (AP). AP is essentially different from traditional uncorrelated percolation models. Indeed, at each time step, two edges are first randomly selected as potential candidates. Then, a product rule (PR) is applied: the candidate edge that actually minimizes the product of the sizes of the two clusters containing its end points is eventually established, while the other one is disregarded. An alternative is the sum rule (SR) in which the established link is the one that minimizes the sum of the two clusters' sizes. In AP, a giant cluster emerges after a number of steps that is much smaller than the system size. As a consequence, the order parameter exhibits an extremely abrupt "jump" at the percolation point. At first glance, such a behavior resembles a discontinuous transition and this is the reason why AP was termed "explosive percolation" (EP) [9]. As we will see later on (in Section 2.7), EP can actually be a CPT, or a genuine DPT.

After the initial work by Achlioptas [9], extensive studies on EP have been carried out in various configurations, such as 2D lattices [12,33–42], 3D or high-D lattices [42], Bethe lattices [43], ER networks [9,12,13,34,39,44–47], scale-free (SF) networks [34,48,49], and real-world networks [50,51].

The key factor in AP is the competition (at each time step) between two potential edges before the actual selection of the one which is added to the network. Such a competition mechanism can be straightforwardly extended from two to more edges, leading to the so called "best-of- m " rule, where $m \geq 2$ potential edges are first chosen at each time step. Then, the candidates are evaluated using either PR or SR, and the one minimizing the product or the sum of the two end clusters is eventually established. For this reason, the algorithm is also known as the "min-cluster- m " rule. In the limit case where $m = N$, the "min-cluster- m " rule becomes the smallest cluster model [52], i.e., at each time step the two smallest clusters in the network are identified to connect. So far, the "best-of- m " rule has been studied in 2D lattices [53–55], 3D lattices [55], ER networks [13,46,52,56,57], and real-world networks [51,58].

Other generalized Achlioptas processes (GAP) have been proposed and investigated, including the da Costa–Dorogovtsev–Goltsev–Mendes (dCDGM) model [11] and the l -vertex model [59–61], which has (for $l = 3$) several variants, such as the adjacent edge rule [62], the triangle rule (TR) [59,62], and the clique-3 competition rule [46]. All GAP models have been studied in 2D lattices [41], in ER networks [11,13,39,46,47,59–67], and in real-world networks [68].

The relevant literature regarding AP and GAP models is summarized in Table 2, where the reader can visually find information on which model has been adopted in which specific network's topology.

The fundamental difference between classical percolation and AP is that the latter follows a competitive (or correlated) rule, which involves non-local (or even global) information on the graph. It is well known that introducing long-range interactions during percolation can change the universality class of the PT, or even the transition type. Thus, it is not surprising that a percolation model with non-local competition (like AP) can give rise to abrupt transitions. Following the same motivation, many other models were also investigated, in which various constraints on the occupation of edges were proposed. These models include the Bohman–Frieze–Wormald (BFW) model [13,69–76], the probability model [53,56,77–82], the hybrid model [25,53,83,84], the diffusion-limited cluster aggregation (DLCA) model [10,85], the spanning cluster-avoiding (SCA) model [86,87], the two-species cluster aggregation model [88], the bootstrap model [89], the k -core model [83,90–92], the cascading failure model in inter-dependent networks [93–101], and others [102–105]. The relevant literature regarding all these other models is summarized in Table 3. Like in the case of Table 2, the reader can find orientation on the various adopted models and network structures.

Let us now briefly discuss the mechanisms leading to EP. In classical percolation, nodes are progressively chosen at random to connect. Thus, the probability to establish a specific edge is proportional to the product of the sizes of the clusters where the edge's end points reside. The result is that a giant cluster can quickly form, and this in turn amplifies the probability for its further expansion. It is therefore expected that CPT always occurs under this scheme.

Table 3

Summary of the relevant literature concerning other algorithms leading to EP in lattices, random networks, and complex networks: small-world (SW), SF, and inter-dependent networks, including the BFW model, the probability model, the hybrid model, the DLCA model, the spanning cluster-avoiding (SCA) model, the hierarchical model, the two-species cluster aggregation model (TCA), the bootstrap model, the k -core model, and the cascading failure model.

Model	Lattice ($d = 2$)	Lattice ($d \geq 3$)	Other lattices	Random	SW	SF	Inter-dependent
BFW	[71]	[71]	[71]	[13,69,70,72–75]		[76]	
Probability	[53,77,78,80]	[81]	[81]	[56,77,79,80,82]			
Hybrid	[25,53,83]			[25,83,84]			
DLCA	[10,85]	[85]					
SCA				[86,87]			
TCA				[88]			
Bootstrap				[89]			
k -core	[83,91]	[91]		[83,91,92]			[90]
Cascading failure				[102,105]			[93–101,106]
Others					[103,104]		

In AP and GAP models, instead, competition among potential edges is introduced at the basis of the cluster-merging processes. The key factor leading to EP is here that the competition mechanism systematically suppresses the formation of large components [47,77,78,86], and such a suppression principle generates the necessary “powder keg” (i.e., abundant small-sized clusters) in a specific range before the onset of the transition [52]. Typically, after a number of steps, such small/medium sized clusters merge to each other, and a giant cluster suddenly emerges. Similar suppressive mechanisms have also been pointed out to occur in the DLCA model [10], the spanning cluster-avoiding model [86], and the explosive synchronization model [19].

In BFW models [69,73], another underlying mechanism for EP (namely, the growth by overtaking) was revealed. In these models, the growth of the largest cluster is so severely limited by the edge selection rules that it can only merge with isolated nodes. As a consequence, all significant changes in the size of the giant cluster result from two smaller components merging together and overtaking the previous largest cluster to become the new largest component.

The central issue is whether or not EP is really a discontinuous transition, especially in the thermodynamic limit. Because of the extremely slow convergence to an asymptotic behavior as the system size increases, it is difficult to determine the transition nature of EP with the only help of numerical simulations. Therefore, whether or not EP is a genuinely discontinuous transition became, for some time, the object of a controversial issue [11–13,36,37,45,47,107]. Later, it was eventually established that the EP transition of Ref. [9] is in fact continuous in the thermodynamic limit [11].

The continuous character of EP was initially asserted on the basis of numerical results for a specific model [12], and later on it was further pointed out that EP under AP is continuous, but with a unusual finite size behavior. Namely EP under AP belongs to a new universality class with an extremely small exponent of the order parameter [11]. Numerical results give a power law $S(t) \propto (t - t_c)^\beta$ with $\beta = 0.0555$ (close to $1/18$), for $t_c = 0.9232$. This very small value of the exponent β that leads to a large discrete jump in the order parameter, resembling a discontinuity at the critical point.

Finally, based on numerical investigations regarding the average size of medium-sized cluster [52], Riordan and Warnke provided a mathematical proof that all AP models on ER networks give rise actually to continuous percolation transitions [13]. The proof reveals that the number of clusters participating in the merging process (needed to generate a macroscopic-sized cluster) is not sub-extensive to the system size, and therefore it cannot bring out a DPT in the thermodynamic limit. On the other hand, Ref. [13] also showed that, for a random network, if the number of competitive edges m is allowed to increase with the system size N in a way that $m \rightarrow \infty$ as $N \rightarrow \infty$ [for example, $m \propto \log(\log N)$], a true DPT can occur also in the thermodynamic limit.

It is today accepted that EP under Achlioptas processes are actually CPT (second-order), but with a universality class different from all the others previously observed [12]. The interested reader is addressed to Ref. [25] for a critical and exhaustive review on all critical exponents. While other GAP models (such as minimal cluster model [52] and dCDGM model [11]) turn out to generate explosive (but actually continuous) percolation transitions in finite size systems, it has been discovered that many alternative algorithms lead to genuinely DPT. These latter models include the BFW model [70], the Gaussian model [78], the spanning cluster-avoiding model [86], the hierarchical model [103], and the two-species cluster aggregation model [88].

The next Sections will expand the discussion on all the concepts, ideas and methods summarized so far.

2.2. Explosive percolation in the Achlioptas process

The percolation process introduced by Achlioptas can be described as follows [9]. The starting point is a network with N vertices and no edges, and edges are added one by one. Precisely, at each time step, two potential edges e_1 and e_2 are arbitrarily chosen, each of which is supposed to connect two clusters (i.e. the two end points of each edge are supposed to belong to two distinct clusters). The size of a cluster is defined here as the number of nodes in it. Now, let us denote the sizes of the two end clusters linked by e_1 as s_1 and s'_1 , and the sizes of the two end clusters linked by e_2 by s_2 and s'_2 . In order to select the edge to be actually added to the network, s_1, s'_1, s_2 , and s'_2 are compared among them, by means of the use of certain rules. For instance, one can compare the products $p_1 = s_1 s'_1$ and $p_2 = s_2 s'_2$. If $p_1 < p_2$ ($p_1 > p_2$), then edge e_1 (e_2) is

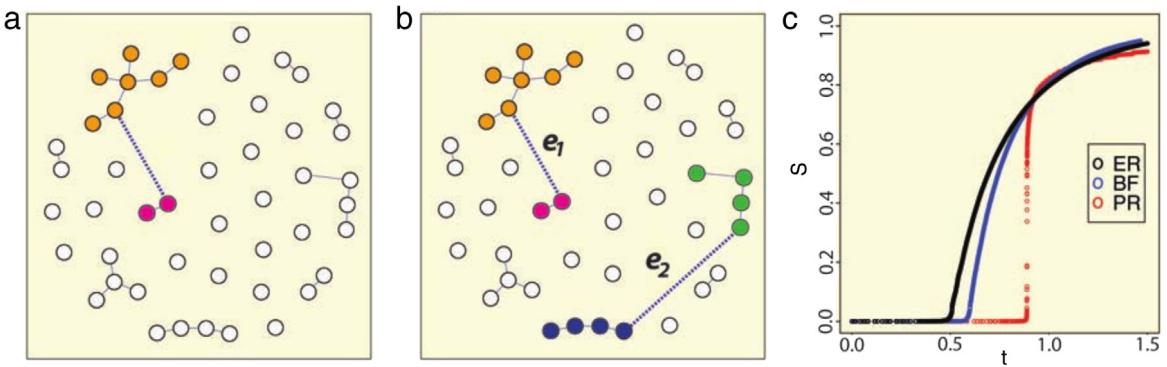


Fig. 1. Comparison between random percolation and AP percolation [65]. (a) Classical ER percolation, where edges are randomly added into the network. (b) AP percolation, where at each step two potential edges compete to be established. (c) Order parameter (the relative size of the giant cluster) vs. control parameter (the number of added edges normalized by system size). For the ER process, PT is continuous (black solid line). For the BF process, it is still continuous with a postponed threshold (blue solid line). For the AP process with the use of the PR rule (red dotted line), EP is observed. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)
Source: From Ref. [9]. Reprinted with permission from the American Association for the Advancement of Science.

established and e_2 (e_1) is discarded. This rule is called “the product rule”, as one always establishes those links minimizing the product of the sizes of the two end clusters. If one instead replaces the “product” by the “sum” of the clusters’ sizes, the resulting rule is called the “sum rule”.

Fig. 1 compares classical percolation and AP, and shows that the latter exhibits two distinguished characteristics. One is that the percolation threshold is significantly postponed (Fig. 1(c)). Another, and most important, is that the order parameter shows an extremely steep “jump” at the percolation point which, at first glance, seemingly resembles a DPT (Fig. 1(c)). In order to characterize the jump, Ref. [9] counts the fraction of vertices added in the network when the giant cluster grows from size at most \sqrt{N} to size at least $N/2$. It is shown that at most $2N^{2/3}$ steps are needed. This implies that, at the phase transition, a constant fraction of the vertices is accumulated into a single giant cluster within a sub-linear number of steps.

It should be emphasized that not all competitive rules lead to EP. Traditionally, the selecting rules in percolation models can be classified into two types. One is the “bounded-size rule”, in which decisions depend only on the sizes of the clusters, and all clusters with size greater than a certain constant K are regarded as the same. For example, $K = 1$ is the Bohman–Frieze (BF) rule, where e_1 is chosen if it joins two isolated vertices, otherwise e_2 is chosen. As shown in Fig. 1(c), the percolation under the BF rule turns out to be still continuous, but with a postponed threshold. In fact, it has been conjectured that the percolation transition is continuous for all bounded-size rules [108].

At variance, an “unbounded-size rule”, like the PR or SR in Ref. [9], treats all clusters of distinct sizes uniquely. Such unbounded-size rules typically lead to interesting transition behaviors, such as shifting the transition threshold and even yielding EP. Not all AP models lead to EP. For example, in the “largest-cluster rule” [9,52], at each step one selects the two largest clusters in the network to connect. Apparently, the process seems to be similar to AP (due to the considered competitive rule for adding edges). However, the percolation turns out to be continuous under this rule.

PR and SR can be straightforwardly extended to the case where at each time step $m > 2$ potential edges are chosen to compete. This leads to the “best-of- m rule” or the “min-cluster- m rule”, which typically leads to EP [52]. Note that when $m = 1$ the “best-of- m rule” recovers the ordinary random percolation, which is continuous. When $m = 2$, it specializes the PR/SR model studied in Ref. [9], which leads to EP. With increasing m , the process becomes more and more competitive. If $m = N$, where N is the total number of nodes, we have global competition among all unoccupied links in the network. This limit situation corresponds to the “smallest-cluster rule” [52], where at each step two smallest clusters in the network are identified to be connected, and global information is used in order to identify the two smallest clusters [46].

2.3. Explosive percolation in generalized Achlioptas processes

Stimulated by the observation that non-local or global competition rules during cluster-merging processes can lead to EP, a number of generalized Achlioptas processes (GAP) have been studied. In the following, we briefly describe two generalizations of the original AP.

2.3.1. The dCDGM model

The model has been studied in Refs. [11,39,47,64–68]. At each step, a pair of clusters are initially randomly selected, and the smaller cluster is kept. Then the same process repeats, and another (smaller) cluster is picked. In order to establish the link, one actually chooses two random nodes (one from each of the chosen clusters) and connects them. The above procedure is sequentially repeated. The dCDGM model provides even more stringent selection of small components during cluster-merging processes than usual AP, since it guarantees that the product of the sizes of the two merging clusters is the

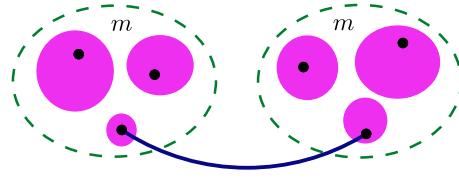


Fig. 2. (Color online). Illustration of the dCDGM algorithm [65]. At each time step, two sets of m nodes are chosen at random. Within each set, the node in the smallest cluster is selected, and the pair of selected nodes are connected.

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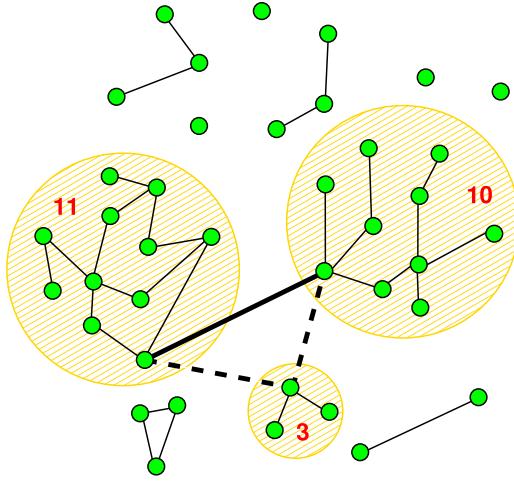


Fig. 3. (Color online). Illustration of the algorithm in a variant of the TR (3-vertex) model [59]. At each step, three vertices are selected and those two vertices that reside in clusters whose size difference is minimal are connected.

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smallest of the four possibilities. This algorithm can be naturally extended to the case of $m > 1$ nodes. For instance, Fig. 2 illustrates this process for $m = 3$ [65].

2.3.2. The l -vertex rule

In this second model, l vertices are randomly chosen at each step, and two of them are connected with an edge using certain rules [59–61,109]. Notice that the resulting percolation process includes the ER model, which is obtained for $l = 2$. Also, Achlioptas processes can be regarded as a special case of the 4-vertex rule. When $l = 3$, the rule has several variants, such as the triangle rule (TR) [62], the adjacent edge rule [62], and the clique-3 competition rule [46]. In all these variants, three distinct vertices are randomly selected at each step. Then, one examines the three possible edges (TR and clique-3 competition) or two adjacent edges in the triangle, and selects the edge that connects the two smallest components.

The bounded-size rule variant occurs when all components have size above the bound K , and one chooses a random edge out of the three; if two components have size above the bound K , one chooses a random edge out of the two adjacent of the smallest component. Furthermore, Nagler et al. have proposed a variant of the TR [59]. There, one first chooses three vertices at random and connects those two vertices that reside in clusters whose size difference is minimal (see Fig. 3). Notice that this latter model does not suppress the growth of the largest cluster, in contrast to many other Achlioptas processes. It is found that the continuous phase transition can have the shape of an incomplete devil's staircase with discontinuous steps in arbitrary vicinity of the transition point [59].

2.4. Explosive percolation under other algorithms

Beyond ordinary percolation, many other models have been developed, where various constraints on the occupation of edges have been considered in order to produce EP. In the following, we briefly review some of these models.

2.4.1. The BFW model

In Ref. [70] a model is introduced (based on the work of Bohman, Frieze, and Wormald (BFW) [110]), where the members of a set of potential edges are considered one at a time, and either added to the network or rejected. Specifically, the algorithm of BFW is as follows. Initially, there are N isolated nodes and all bonds are unoccupied (empty). Therefore, there are N

clusters of unitary size. At each time step, one edge is added to the network. Let u be the total number of selected bonds, t the number of occupied bonds, and k the stage of the process, initially set to $k = 2$. The first bond is occupied at random, such that $t = u = 1$.

Then, at each step u ,

1. One edge is randomly selected out of all edges (or the unoccupied ones);
2. The maximum cluster size l is measured, if the selected edge is occupied;
3. If $l \leq k$, go to 4. Else, if $t/u \geq g(k) = 1/2 + \sqrt{1/(2k)}$, go to 5. Else, increment k by one and go to (3);
4. Occupy the selected edge;
5. Increment u by one.

The procedure is applied iteratively from $t = 0$ (corresponding to a bond occupation fraction $p = t/(2N) = 0$) to $t = 2N$ (where all bonds in the system are occupied, i.e., $p = 1$). Using such a model, Chen et al. have shown that multiple stable giant clusters can coexist, and the percolation transition is strongly discontinuous [70]. The BFW model has been extensively studied in various networks, including lattices and tree-like networks [71], ER networks [13,69,70,72–75], and SF networks [76]. Notice that in Refs. [69,70,73,74,76,110] edges are randomly chosen from all edges (i.e. regardless on whether or not they are occupied), while in Refs. [71,72] edges are only sampled from the set of unoccupied bonds. In Refs. [70–72,76], $g(k) = 1/2 + \sqrt{1/(2k)}$. Later, generalizations of this function have been considered as $g(k) = 1/2 + (2k)^{-\beta}$ [73], and $g(k) = \alpha + (2k)^{-\beta}$ [69,74].

2.4.2. Probability models

In this type of models, an edge between a pair of clusters is occupied with certain probability at each time step. This process differs substantially from AP, since here all unoccupied connections can potentially be chosen at each step, though some of them with a much larger probability [77]. For instance, in cluster aggregation model [79,80], all currently unoccupied edges are considered. Let one denote the sizes of the clusters to be connected by the edge ij as s_i and s_j . Each of these vacant edges is given a weight $(s_i s_j)^\alpha$, and the sum over all the weights is calculated as a normalization constant w . Then the edge ij is occupied with probability $(s_i s_j)^\alpha / w$.

In the Hamiltonian model of Ref. [77], instead, one starts with a network of N vertices without edges, so each vertex initially belongs to a different cluster. First, a simple Hamiltonian H is defined as a function of the cluster sizes and of the number of redundant edges added to the clusters. Then, a new edge b between any pair of vertices not yet connected is added, with a probability proportional to $\exp(-\beta \Delta H_b)$, where ΔH_b is the energy change after adding the edge.

Another model belonging to this class is the largest cluster model (or the Gaussian model) [78]. There, a link is randomly selected among the unoccupied ones. If its occupation would not lead to the formation or growth of the largest cluster, it is always occupied; otherwise, it is occupied with probability $\min\{1, \exp[-\alpha(\frac{s-\bar{s}}{\bar{s}})]\}$. For $\alpha > 0$, this probability suppresses the formation of a cluster significantly larger than the average, inducing a homogenization of cluster sizes, which is important to generate EP.

Such probability models have found applications in 2D lattices [53,77,78,80], 3D and high-D lattices [81], as well as ER networks [56,77,79,80,82].

2.4.3. Hybrid models

Hybrid models are those for which the rule for the occupation of edges is a mixture of ordinary percolation and competitive rules: at each time step, an edge can be chosen randomly with probability p , while it can also be chosen following a given competitive rule with probability $1 - p$. For example, in Refs. [25,84], PR is used; in Ref. [53], the best-of-ten PR is used; in Ref. [83], a mixture of $k = 2$ - and $k = 3$ -core on the random network and a mixture of $k = 3$ -core and jamming percolation models in 2D lattice are studied. Hybrid models have been studied in 2D lattices [25,53,83] and ER networks [25,83,84].

2.4.4. The diffusion-limited cluster aggregation (DLCA) model

In this model, studied in Refs. [10,27,85], N single particles are initially randomly placed in a $L \times L$ square lattice. The particles are assumed to be Brownian, so that the velocity of a cluster is inversely proportional to the square root of its size, i.e. the velocity of an s -sized cluster is given in a general form as $v_s \propto s^{-1/2}$. As a consequence, the larger clusters move considerably more slowly. When two clusters become nearest neighbors, they merge and form a larger cluster. The Brownian motion suppresses the mobility of the largest clusters, impeding their growth, and leading to the discontinuous emergence of a giant cluster. Generalized Brownian motion, for which the velocity is inversely proportional to the size of the cluster to a power η (i.e., $v_s \propto s^\eta$) gives rise to a tri-critical point that separates discontinuous from continuous transitions as a function of η . Namely, as soon as η_c is a tri-critical point, it is shown that when $\eta < \eta_c$ ($\eta > \eta_c$), the PT is discontinuous (continuous) [10].

2.4.5. The spanning cluster-avoiding (SCA) model

The SCA model, studied in Refs. [86,87], considers percolation on an Euclidean lattice, where the emergence of a spanning cluster is concerned. In a spanning cluster, a path of activated links exists that connect sites from one side of the lattice

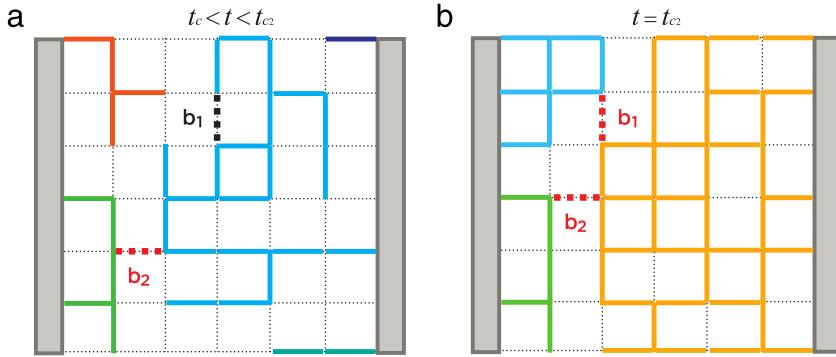


Fig. 4. (Color online). Illustration of the algorithm in the SCA model with $m = 2$ on a square lattice [86]. (a) At each time step, two empty bonds, b_1 and b_2 (shown as dashed lines), are randomly selected. If one of them is a bridge bond (like b_2 in this case), then the non-bridge bond (b_1) is chosen. (b) At $t = t_{c2}$, two bridge bonds can be selected for the first time. Then, one of them is taken randomly and a spanning cluster is formed.

to the other. Inspired by the best-of- m rule, Cho et al. introduced a new global constraint for the dynamics. As shown in Fig. 4, at each time step t , m unoccupied edges are chosen randomly and classified into two types: bridge and non-bridge edges. Bridge (non-bridge) edges are actually those that would (would not) form a spanning cluster when occupied. The SCA model strictly avoids bridge edges to be occupied, and therefore one of the non-bridge edges is randomly selected and occupied. If the m potential edges are all bridge edges, then one of them is randomly chosen and occupied. Once a spanning cluster is formed, no more restrictions are imposed on the occupation of edges. This procedure continues until all edges are occupied. Numerical simulations and theoretical results have shown that the EP in SCA model can be either discontinuous or continuous for $d < d_c = 6$, depending on the number of eligible edges m . For $d \geq d_c$, the transition is shown to be continuous for any finite and fixed value of m . However if m varies with the system size N , a discontinuous transition can also take place for $d \geq d_c$.

2.4.6. The two-species cluster aggregation model

In this model, studied in Ref. [88], N initially isolated nodes are randomly separated into two populations: half of them is colored in black (B) and the other half in white (W), with all nodes in the same cluster keeping the same color. At each time step, the three possible combinations BB, BW, and WW are given selection probabilities $1/(1+2p)$, $p/(1+2p)$, and $p/(1+2p)$, respectively (where p is a model parameter in the range $0 < p \leq 1$). Two clusters are then selected independently of the cluster sizes, and two nodes (one from each cluster) are selected randomly and connected, which causes the two clusters to merge. If the two selected clusters are the same, then two distinct nodes from that cluster are connected. Meanwhile, the colors of all the nodes in the resulting merged cluster have to be updated according to a certain rule. See Fig. 3 in Ref. [88] for an illustration of the algorithm. The cluster aggregation model can exhibit both type-I and -II DPT, as the model parameter changes [88].

2.4.7. The bootstrap models

In the context of EP, some bootstrap models have been proposed [89]. The standard bootstrap percolation process on a lattice assumes that sites are either active or inactive, and the state of a site depends on its neighbors. Initially, each site is activated with a probability p , and inactivated with probability $1 - p$. Every activated site remains in its state, while each inactivated site can become active (and remain active forever) if its k nearest neighbors are active (with $k = 2, 3, \dots$). The procedure is continued until the system reaches the stable configuration which does not change anymore. In the final state of the process, one is concerned on whether or not a spanning giant cluster of active sites exists. Bootstrap percolation can show either continuous or discontinuous transitions, depending on the type of networks (lattice or random), and on the dimensionality of the lattice.

2.4.8. The k -core model

k -core percolation [83,90–92] is a model of correlated percolation beyond the ordinary one. The k -core of a graph is the maximal subgraph for which all vertices have at least k neighbors. It can be found as an asymptotic structure obtained by recursively removing all nodes whose neighbors are less than k . For example, if $k = 2$, all leaves (nodes with only a connection) are removed in a first step. This might turn into some nodes (which originally had 2 or more neighbors) becoming leaves. Then, these latter nodes are removed in a second round, and if this again turns nodes into leaves, they are removed again, and so on. In this way, one ends up either with an empty cluster, or with a cluster where all nodes have $k \geq 2$ neighbors, i.e., the 2-core. It has been proved that k -core percolation on random networks exhibits a CPT for $k \leq 2$, and a DPT for $k \geq 3$ [111]. However, unlike a typical DPT, not one but several diverging length scales exist such that the transition has an unusual nature [112].

2.4.9. Cascading failure model in inter-dependent networks

This type of models aims at studying processes such as cascading-failure dynamics (and related properties, as resilience and robustness) in inter-dependent systems [93–101,106]. For example, the infrastructure network of a country includes both power grids and computer networks. The latter are needed for controlling the power stations, while the former are needed to provide power to the computers. If a node fails in one of the two networks, this can lead to a huge back and forth cascade of failures, as historically happened in many countries. See Fig. 2 of Ref. [106] for an illustration of a cascade of failures in inter-dependent networks which resembles a first-order transition. Several other types of interdependencies were also studied. For example, first-order percolation has been investigated in a single network with two sets of links: the dependency links and the connectivity links [94]. Figure 1 in Ref. [94] illustrates the cascading failures of such a latter case. It should be pointed out that, in this DPT, the order parameter is taken as the fraction of nodes in the mutually connected giant cluster, which is not the same as the standard ones. Inter-dependencies of networks, and robustness of power-grids will be also discussed in Section 5.2.1, from the viewpoint of explosive synchronization.

2.4.10. Other models

Several other models for explosive percolation have been proposed. In the half-restricted ER model, two vertices are connected by an edge at each step, but one of them is arbitrarily chosen and the other is chosen randomly from a restricted set [102]. This variant of random percolation can exhibit a DPT. In the hierarchical model [103], Boettcher et al. studied the effect of an ordinary (uncorrelated) bond-adding rule on a network with a recursive, hierarchical structure. It is shown that the hierarchical adding of SW bonds into a 1D lattice can induce DPT. Furthermore, two works are highly relevant for experimental realizations of EP: in Ref. [45], the formation of single-walled nanotube bundles with uniform diameter is investigated, and by applying the cluster repulsion process to stick percolation, it is shown that the transition becomes explosive. In Ref. [113], the electric breakdown of a substrate was studied, on which highly conducting particles are adsorbed and desorbed with a probability that depends on the local electric field. It is revealed that the electric breakdown due to pollution with metallic powder can become explosive under certain conditions. This model is thus an example of a possible experimental realization exhibiting a true DPT.

2.5. Explosive percolation in 2D lattices

The phase transition corresponding to EP is established to be continuous in the mean-field level, or in higher dimensions. An interesting issue is therefore determining the transition nature for EP in lower dimensions. To answer the question, one needs to understand what differentiates percolation in the cases of higher and lower dimensions. We take here ER networks, and two dimensional (2D) lattices, as illustrative examples. ER networks are, indeed, usually treated as infinite dimensional systems, while 2D lattices are typical low dimensional ones.

One of the main features of complex networks is the SW property, which guarantees a long-range correlation among the nodes. Thus, in the growing process of ER, it is easy to generate a core and then the core gradually grows up. When an AP is applied, the growing of the core is suppressed, and a sharp jump in the growth dynamics occurs for finite-size systems. However, when the size N goes to infinity, the AP rule cannot suppress the growing of the core completely, resulting again in a continuous transition. At variance, for a 2D lattice, the nodes feature only short-range correlations, and therefore it is not clear whether or not an AP would produce a discontinuous transition in these circumstances.

A lot of attention has, indeed, been paid to the case of 2D lattices [33–37,40,42,45,71,86,87]. Another reason is that many interaction networks, whether physical or social, are spatially constrained, and bonds can form only between close neighbors. Therefore, 2D percolation models represent possibly the strongest form of such spatial restrictions.

Both bond percolation and site percolation have been studied in 2D lattices. In the former process, bonds are formed randomly and independently throughout the system, and, at a critical concentration p_c of occupied bonds, a finite fraction of the sites (nodes) are connected together, and percolation takes place through a continuous or second-order transition. The latter process corresponds to consider randomly occupied sites on a given lattice with a probability p , with neighboring occupied sites that merge to form clusters. A remarkable finding here is that, for both bond and site percolation (and together with the existence of the sharp jump in the growth dynamics of the 2D lattices), a hysteretic loop can be observed, which is the signature of an irreversible transition. However, this finding does not mean that all the transitions observed in 2D lattices are discontinuous, and the controversy is still standing on whether the 2D lattices percolation is continuous [36,42] or discontinuous [34,35,37,40,45].

2.5.1. Bond percolation

The AP process in bond percolation of 2D lattices was first considered by Ziff in Ref. [33], which started from a $L \times L$ square lattice with periodic boundary conditions in both directions, and with $n = L^2$ sites. Connected sites are considered as one cluster. As initial sites are fully disconnected, the starting point is a set of n clusters. At each time step, only one bond is added, which always connects different clusters, this way reducing the number of clusters by one.

Let the time t represent the number of bonds successfully added, then the number of clusters N_c in the system is reduced to $n - t$. Let s represent the mass in a cluster, i.e. the number of sites. The key point is that each successful new bond is chosen by the minority product rule (PR) of Ref. [9]. That is, two bonds are randomly chosen at the same time, and only the

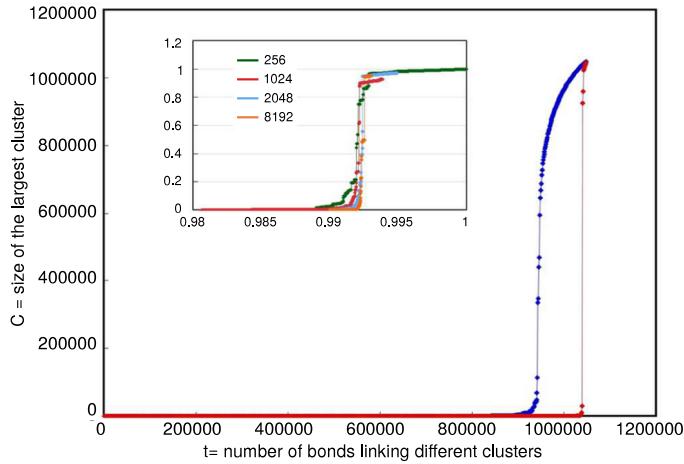


Fig. 5. Regular (blue) and the PR (red) bond percolation on a lattice of size 1024×1024 . Notice the delayed and explosive growth in the PR case. Points are reported every 1024 time steps. The inset zooms on the PR model, for systems of sizes $L = 256, 1024, 2048$, and 8192 (color codes in the legend), with scaled axes C/n vs t/n with $n = L^2$. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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bond that minimizes the product of the masses of the two joined clusters is selected to become occupied, while the other is discarded.

Fig. 5 shows the results obtained in Ref. [33], and reports the maximum cluster size C as a function of t . It is evident that a jump occurs over a small number of time steps, as shown in the inset of Fig. 5 for different system sizes (here the axes are scaled by n). For comparison, regular percolation is also reported in Fig. 5 with a blue curve, pointing to a much smoother transition.

A way to distinguish the difference in the transition between the regular and the PR percolation is by accounting for two different times: the time t_0 at which $C = \sqrt{n}$, and the time at which $t_1 = n/2$ [9]. By calling $\Delta = t_1 - t_0$, Achlioptas, D’Souza, and Spencer found that $\Delta \sim n$ for unbiased growth of ER networks, while $\Delta \sim n^{2/3}$ for the PR growth, and therefore the two transitions are qualitatively different [9]. Ziff measured the same quantities for regular and PR growths on square lattices with different sizes, and also found that the relationship between Δ and n is quite different. For regular percolation, $\Delta/n \sim L^{-36/91}$ was found, while for the PR model $\Delta/n \sim L^{-0.683}$ [33]. Notice that $\Delta/n \rightarrow 0$ as $n \rightarrow \infty$ for both the PR and random growth models. However, it was later discovered that, with a different criterion for the upper end of the gap Δ (say $C = 0.7n$), then one would find that $\Delta/n \rightarrow \text{const.}$ as $n \rightarrow \infty$ for the random growth, while still $\Delta/n \rightarrow 0$ for the PR model [35].

Kim et al. considered the formation of single-walled nano-tube bundles with uniform diameter [45]. Except for the growing process, they also measured the changes in the order parameter during the stick removal process, and found that a hysteresis exists. This result clearly shows that the transition of the stick system with cluster repulsion is discontinuous. AP in such a nano-stick system is defined as follows [45]:

- (I) One first places two sticks α and β at random positions with random orientations.
- (II) Then one lets $\{s_{\alpha_1}, s_{\alpha_2}, \dots, s_{\alpha_n}\} (\{s_{\beta_1}, s_{\beta_2}, \dots, s_{\beta_m}\})$ be the sizes of the clusters which would merge into a bigger one of size $\sum_{k=1}^n s_{\alpha_k} + 1 (\sum_{k=1}^m s_{\beta_k} + 1)$ by the placement of the stick α (β). Here, the size is defined by the number of sticks in the cluster. A PR is adopted by calculating the products:

$$\pi_\alpha = \prod_{i=1}^n s_{\alpha_i}, \quad \pi_\beta = \prod_{j=1}^m s_{\beta_j}. \quad (1)$$

- (III) If $\pi_\alpha \leq \pi_\beta$ ($\pi_\alpha \geq \pi_\beta$) then the stick α (β) is kept and the stick β (α) is removed. Processes (II) and (III) favor the connection between small clusters, which suppresses the growth of larger ones.

An order parameter S is introduced to describe the probability that a stick belongs to a spanning (largest) cluster [45],

$$S = \frac{E_\infty}{E}, \quad (2)$$

where E_∞ is the number of sticks in the spanning cluster, and E is the total number of sticks placed in a given square. In general, EP is marked by the existence of a jump in the order parameter at the critical connecting probability p . For convenience, this probability p is here replaced by the total number of sticks E [45]. Although these two approaches are similar, some differences between the variables p and E have to be noticed. For the former, p is a state variable and is thus

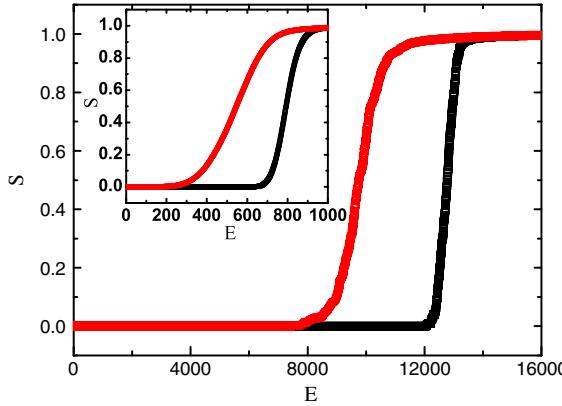


Fig. 6. S vs. E , for $L = 40$. Black (red) solid line indicates S when sticks are added (removed). The two curves manifest the presence of hysteresis. Inset: the same plot but for $L = 10$. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.) Source: Reprinted with permission from Ref. [45].

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independent of the evolutionary process. A small increase of Δp will result in an increase of approximately $\Delta pL^2(L^2 - 1)/2$ new sticks, which is enough to induce a significant increase of E_∞ . For the latter, instead, E is an evolutionary variable and is equivalent to the time t . That is, E will increase one by one, and this will result in a gradual increase of E_∞ . In this sense, we cannot expect a jumping of S in Eq. (2) at the critical point, but only a fast or explosive growing of S .

A first-order like phase transition is usually accompanied with a property of history-dependence, i.e. the existence of a hysteresis loop, because of the presence of meta-stable states. Therefore, one of the most generally accepted (and simplest) methods to verify whether or not an observed transition is first-order like is the existence of a hysteresis loop [114,115]. To measure the hysteresis, one needs to consider both processes of adding and removing sticks. Results are shown in Fig. 6, which clearly manifests the existence of a hysteresis, indicating a first-order like transition.

2.5.2. Site percolation

An alternative mechanism is site percolation [36,37]. It refers to a situation where, initially, all the sites of a square are empty. During the growing process, sites are sequentially occupied, and two occupied neighboring sites are considered to belong to the same cluster. The AP rule is applied as follows. One chooses two candidates at each time step, and investigates which one of them would lead to the smaller clustering. This one is then kept and the other discarded.

Bastas, Kosmidis, and Argyrakis found that this procedure considerably slows down the emergence of the giant component [36]. Their algorithm for the case of the sum rule proceeds as follows: initially, one starts from an empty lattice and randomly occupies a single site. Then, at each time step, one randomly chooses two unoccupied sites, say A and B , and calculates the sizes s_A and s_B of the resulting clusters to which A and B belong, respectively. When $s_A < s_B$ ($s_B < s_A$), site A (B) is permanently occupied, and site B (A) discarded. When $s_A = s_B$, one randomly selects either A or B and permanently occupies it. With S_{max} representing the largest cluster at time t , the process is repeated until the entire lattice is covered.

Fig. 7 shows an illustrative example of the algorithm applied on an $L \times L$ square lattice with periodic boundary conditions, where the white (colored) cells correspond to unoccupied (occupied) sites, and where different colors (red, green, gray, and blue) indicate different clusters. In Fig. 7(a) one randomly selects a pair of unoccupied sites (yellow), say A and B , one at a time, and evaluates the size of the clusters that are formed when A and B are considered, s_A and s_B , respectively. One finds that $s_A = 10$ and $s_B = 14$. Therefore, A is kept (because it leads to the smaller cluster) and B is discarded. The result of this operation is shown in Fig. 7(b).

Similarly to the case of bond percolation, one can introduce here an order parameter S to track the explosive nature of the transition, which is defined as the ratio of the sites that belong to the largest cluster s_{max} to the total number of sites in the lattice,

$$S = \frac{s_{max}}{N}, \quad (3)$$

where $N = t$. By monitoring this quantity, Bastas et al. [36] found an explosive behavior similar to that of Figs. 5 and 6.

Later on, Choi, Yook, and Kim studied the explosive site percolation under AP with a product rule in a 2D square, and also found the existence of a hysteresis loop in the order parameter [37]. The algorithm is similar to the one in Eq. (1) with the word “sticks” replaced by “sites”. The reverse AP process is also the same as in Ref. [45]. The main difference is that $n(m)$ in Eq. (1) is now at most 4, while $n(m)$ for bond percolation is always 2. Choi et al. noticed a qualitative difference in the hysteresis loops between the SR and PR cases. Let $A(L)$ represent the area enclosed by P_{max} for the increasing and decreasing processes on L . The SR shows that $A(L) \rightarrow 0$ as $L \rightarrow \infty$ [36], while the PR shows that $A(L)$ increases when L increases and

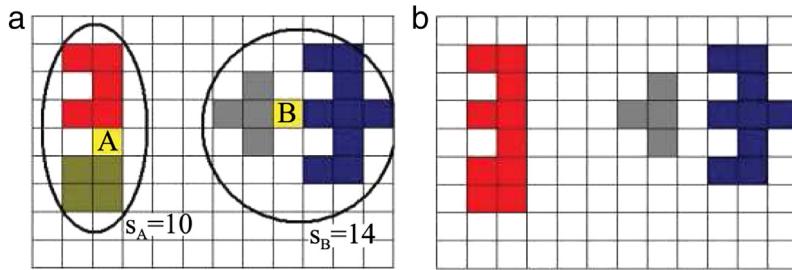


Fig. 7. AP according to the SR for site percolation. White cells correspond to unoccupied sites, while colored cells correspond to occupied sites. Different colors (red, green, gray, and blue) indicate different clusters. (a) One randomly selects two unoccupied sites (yellow), denoted by A and B . The size of the clusters after the addition of A and B , s_A and s_B , respectively, is evaluated. In this example $s_A = 10$ and $s_B = 14$. (b) According to AP, A is kept and B is discarded. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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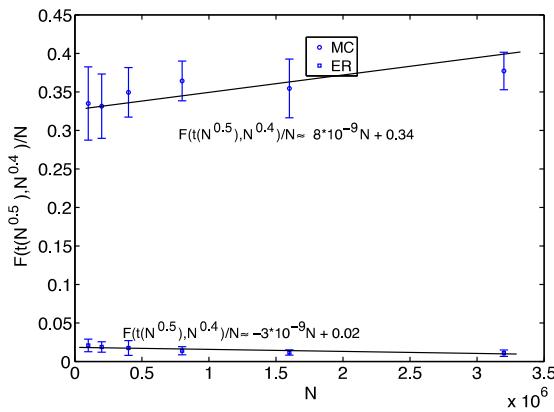


Fig. 8. Size of the powder keg ($\frac{1}{N}F(t(N^{0.5}), N^{0.4})$) as a function of N , for the ER (squares) and AP (circles) sum rules (cluster minimized).

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then saturates to a nonzero value [37]. In a 2D lattice, the PR induces a transition whose nature is completely different from that induced by the SR, as the percolation transition caused by the PR is discontinuous.

2.6. Mechanisms for explosive percolation

The first studies and predictions on EP (with Achlioptas processes on top of ER random networks) were mainly based on numerical simulations [9]. Subsequent studies, however, tried to dwell more deeply upon the generation mechanisms for the observed explosiveness in the transition.

2.6.1. Powder keg

First, Friedman and Landsberg tried to relate EP with a sort of “powder keg” mechanism taking place immediately before the explosive transition in ER-type models [52]. More specifically, before the onset of the transition, a fixed non-zero fraction of nodes in a small-sized cluster constitutes a “powder keg”. Let $F(\tau, a)$ be the number of nodes in clusters of size $\geq a$ after the addition of the τ th edge. In AP processes with sum rules, one finds that the powder keg

$$W = \frac{1}{N}F(t(N^\alpha), N^{1-\beta}) \quad (4)$$

approaches a non-zero constant in the large N limit for some $0 < \alpha < 1$, and $\beta < 1$. In other words, at time $t(N^\alpha)$ (which is the beginning of the phase transition) a fixed non-zero fraction of nodes are contained in small clusters with sizes ranging from $N^{1-\beta}$ to N^α . This set is termed as the powder keg, and has a size approaching a non-zero constant in the large N -limit under the AP rule, while the powder keg is empty in the large N limit for ER rules. Such a conclusion was based on a linear regression of data, comparing the size of the powder keg (estimated numerically) and the system size N , as shown in Fig. 8 for $\alpha = 1/2$ and $\beta = 0.6$.

Although each individual cluster in this set of small-sized clusters contains only a vanishing small fraction of nodes, if taken together these nodes are ignitable and collectively enable an explosive transition. Based on identified powder keg of

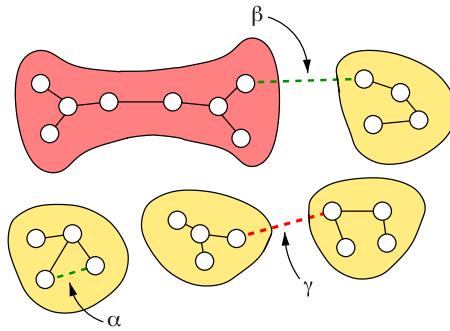


Fig. 9. Three candidate bonds are considered for possible addition at a given step, namely α , β and γ . Among these bonds, α has the smallest probability to be added due to condition (ii) of the main text, β cannot be accepted due to condition (i), and the most probable bond is the γ bond.

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appropriate sizes, a heuristic criterion has been proposed to predict whether or not a given random network will display EP. In particular, it is conjectured that any network model for which the probability of creating a cluster of size a at time t is proportional to $(F(t, a/2)/N)^p$ with $p > 1$ will produce a powder keg. A more simplified criterion based on powder keg has been proposed in Ref. [56], which concludes that, under general conditions, a percolation process will be explosive if the mean number of nodes per cluster diverges at the onset of the phase transition in the thermodynamic limit.

2.6.2. Suppression principle

The original model of AP essentially identifies the dynamics that prevents the creation of a giant cluster by choosing one edge from a given number of randomly selected potential candidates [9]. The following two ingredients are sufficient for obtaining an explosive transition [77]: (i) the size of all growing clusters should be kept approximately the same, and (ii) the merging bonds should dominate with respect to the redundant bonds. The term “merging bonds” indicates those edges that connect vertices in distinct clusters, while “redundant bonds” are edges connecting vertices in the same cluster, as shown in Fig. 9. These two conditions are sufficient for obtaining a first-order like transition in a growth process where bonds are included one by one. More precisely, merging bonds must be introduced with much higher probability than redundant ones.

By defining a Hamiltonian $H(G)$ of a graph G that includes the two ingredients, one can explicitly study the distinct roles of merging and redundant bonds. $H(G)$ reads as

$$H(G) = \sum_{i \in \mathbf{C}} s_i^2 + l_i s_i^\alpha, \quad (5)$$

where the sum is over the entire set of clusters \mathbf{C} , s_i is the number of vertices in cluster i , and l_i is the number of redundant bonds added to the cluster. The parameter α controls the probability of redundant bonds to be added to the system. For small values of α , redundant bonds are favored over merging bonds, and one might expect that a new merging bond will be included only after the addition of all possible redundant bonds. The situation becomes different when α is large, and redundant bonds are not included. By tuning the parameter α in Eq. (5) from small to large values (from $\alpha = 2.5$ to 2.7), a change from a slow continuous growth to a sharp transition is observed, as shown in Fig. 10, which reports the fraction occupied by the largest cluster P_∞ vs. the average connectivity k .

The same idea is also referred to as the suppression principle (SP) [47]. As schematically reported in Fig. 11, one first classifies the types of edge candidate pairs e_1 and e_2 as follows: (i) both e_1 and e_2 connects clusters of potentially different sizes. Clusters of sizes $s_{1a}^{(i)}$ and $s_{1b}^{(i)}$ are connected by the edge e_1 , and clusters of sizes $s_{2a}^{(i)}$ and $s_{2b}^{(i)}$ are connected by the edge e_2 ; (ii) one edge (say, e_1) is an intra-cluster edge (in a cluster of size $s_1^{(i)}$), while the other edge (e_2) is an inter-cluster edge and connects the two clusters of sizes $s_{2a}^{(ii)}$ and $s_{2b}^{(ii)}$; and (iii) both e_1 and e_2 are intra-cluster edges in either the same cluster or two distinct clusters.

Based on the above classification, several variants of dynamic rules have been proposed to determine whether the SP is fulfilled. In a class of models (models of type A) one edge is an intra-cluster edge (cases (ii) or (iii)). In another class (models of type B) when the case (ii) is considered and the intra-cluster edge is selected, so that the cluster size does not increase. When the two potential edges are both intra-cluster edges [case (iii)], one of them is randomly selected. In a third class of models (models of type C), the dynamics proceeds via only inter-cluster connections [case (i)]. Models of type C may be regarded to be the same as models of type B, because the clusters do not grow when the intra-cluster edge is selected. The difference between type B and type C models is that time advances in type B model but not in type C model. Type A models can fail to follow the SP. For instance, when $s_1^{(ii)} = 5$, $s_{2a}^{(ii)} = 3$, and $s_{2b}^{(ii)} = 7$, the edge e_2 is selected in type A models, and therefore, the size of the resulted cluster is 10. On the other hand, if edge e_1 is selected, then none of the clusters would

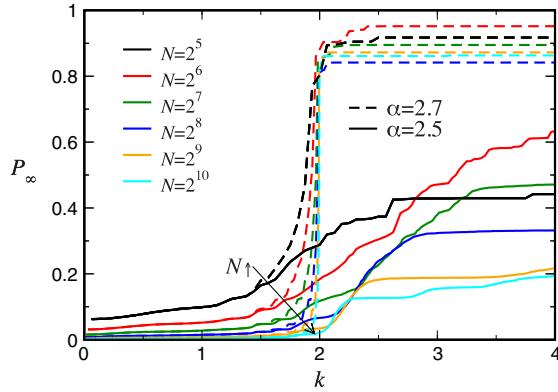


Fig. 10. Transition to EP: the process shows (i) a slow continuous growth for the largest cluster when it favors redundant bonds for $\alpha = 2.5$ (continuous lines), and (ii) an abrupt transition when the merging bonds are favored for $\alpha = 2.7$ (dashed lines). Legend indicates the color code of the curves. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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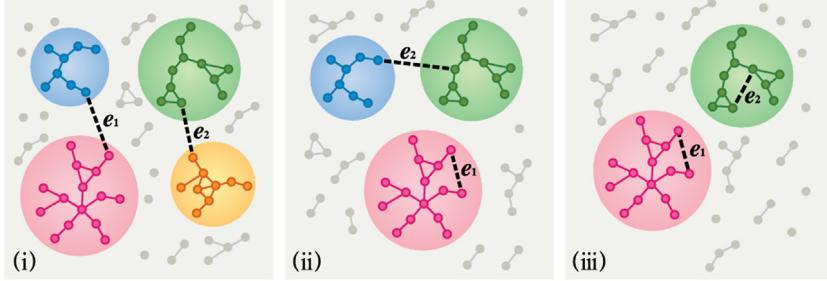


Fig. 11. Classification of types of edge candidate pairs. (i) Both candidates e_1 and e_2 are inter-cluster edges. (ii) e_1 is an intra-cluster edge and e_2 is an inter-cluster edge. (iii) Both e_1 and e_2 are intra-cluster edges.

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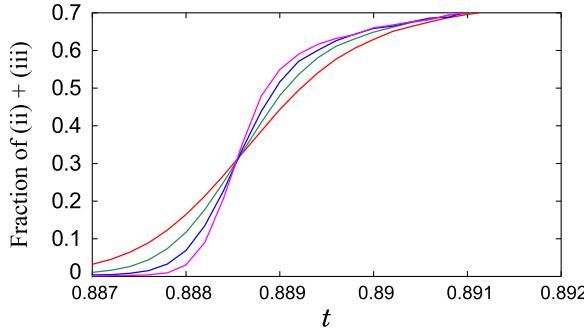


Fig. 12. The fractions of potential edges falling in case (ii) and case (iii) as a function of time t for different system sizes: $N/10^4 = 32$ (red curve), 64 (green curve), 128 (blue curve), and 256 (magenta curve) from the top (bottom) in the small- t (large- t) region. As N increases, the fraction increases dramatically. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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increase in size. As time approaches the percolation threshold, intra-cluster edges can be selected more frequently, which leads to a more frequent failure of suppression principles, as shown in Fig. 12.

In addition, based on the same classification, a fundamental difference between the product and sum rules can be identified. In particular, the PR may not satisfy the suppression principle. A simple example of two inter-cluster connections can be considered, in which $s_{1a}^{(i)} = 2$, $s_{1b}^{(i)} = 7$, $s_{2a}^{(i)} = 4$, and $s_{2b}^{(i)} = 4$. The product of the two inter-clusters is respectively $P_1^{(i)} = 14$ and $P_2^{(i)} = 16$, and therefore, edge e_1 is added to the system. However, the resulting cluster size is 9 in this case, which is larger than the resulting size 8 when e_2 is added. Therefore the SP is not fulfilled.

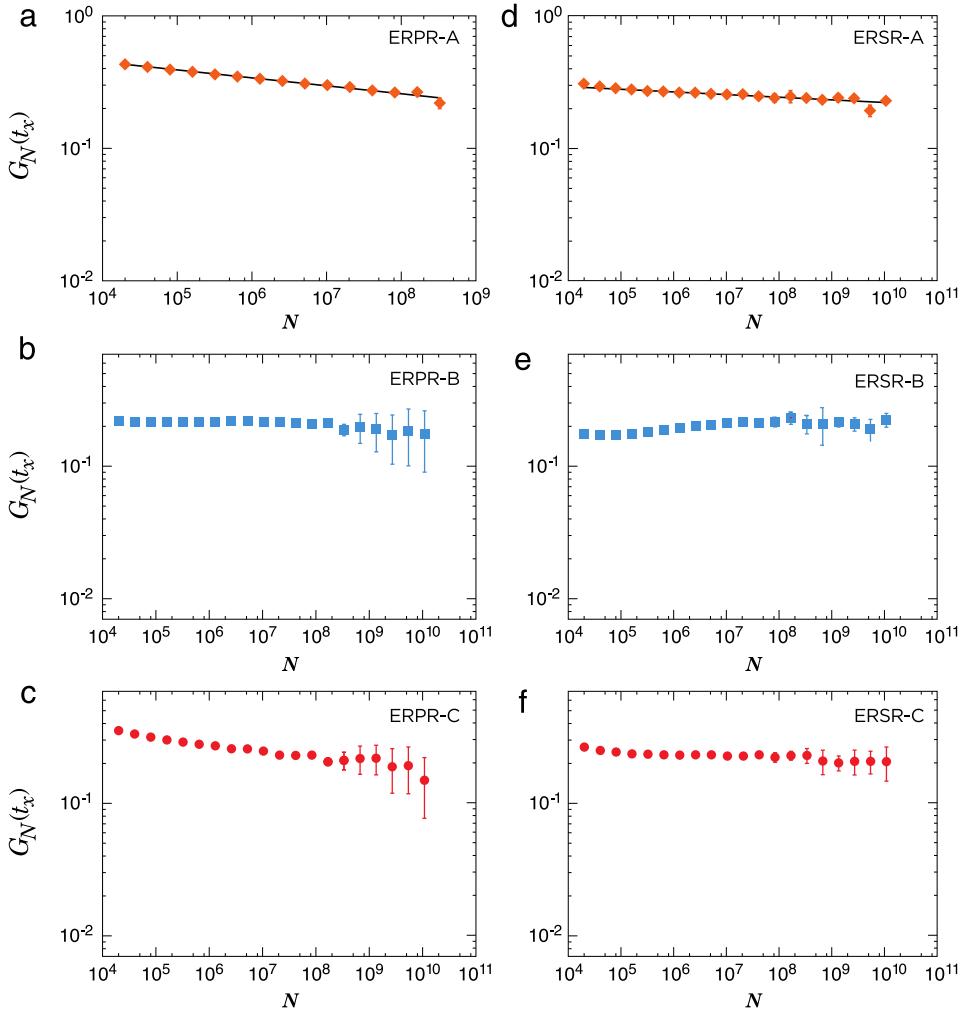


Fig. 13. $G_N(t_x)$ vs. N for models of product rules (a)–(c), and sum rules (d)–(f). (a) and (d) refer to type A models, (b) and (e) refer to type B models, and (c) and (f) refer to type C models. Best linear fits of the data occur for slopes of -0.06 (a), -0.05 (b), -0.05 (c), -0.02 (d), 0 (e), and 0 (f). In the case of sum rules in type B and C models, $G_N(t_x)$ converges to finite values in the thermodynamic limit.

Source: Adapted with permission from Ref. [47]. Courtesy of B. Kahng.

The SR may remove the drawback inherent to the PR, yielding discontinuous percolation transition as observed numerically [47]. To evaluate the discontinuity of the transition, the following two criteria have been proposed: (a) the values ($t_x(N)$, $G_N(t_x)$) remain finite as $N \rightarrow \infty$, where the time $t_x(\infty)$ is regarded as the transition point t_c in the thermodynamic limit, and $G_N(t_x)$ represents the G components at time t_x ; (b) the tangent of the curve of $G_N(t)$ with respect to t at $t_x(N)$ diverges as N increases. More specifically, as shown in Fig. 13(a)–(c), for all models with product rules, $G_N(t_x)$ decreases with increasing N in the whole considered range, suggesting that $G_N(t_c) \rightarrow 0$ in the limit $N \rightarrow \infty$. Furthermore, when the SR is implemented in type A models, a similar convergence of $G_N(t_c) \rightarrow 0$ in the limit $N \rightarrow \infty$ is observed in Fig. 13(d). This indicates that these cases violate the SP, and lead to a continuous percolation transition. However, when the SR is implemented in type B and C models, the asymptotic values of $G_N(t_x)$ look relatively flat, yielding discontinuous percolation transitions within the range of the numerical data (see Fig. 13(e)–(f)).

The original dCDGM model [11] does not follow the suppression principle, because it does not take into account the natural choice of intra-cluster edges. The selection of an intra-cluster edge becomes actually more and more frequent as time approaches the percolation point, and this yields a continuous transition. However, once the dCDGM model is properly modified, discontinuities in the transitions can be observed. The suppression principle plays a fundamental role in achieving discontinuous percolation on random graphs, inter-dependent networks [95,100], adaptive networks [32,116], and fractal networks [103].

For a finite size system, suppressive rules are not a necessary condition for the sharpness of the transition at the percolation point [104]. Rather, a finite size system with aggressive tendency to form a giant cluster may exhibit an instability that is relieved through an abrupt and discontinuous transition to the stable branch. Although the Author of Ref. [104] has tracked analytically such an abrupt transition to percolation, it is worth noticing that discontinuity is especially pronounced

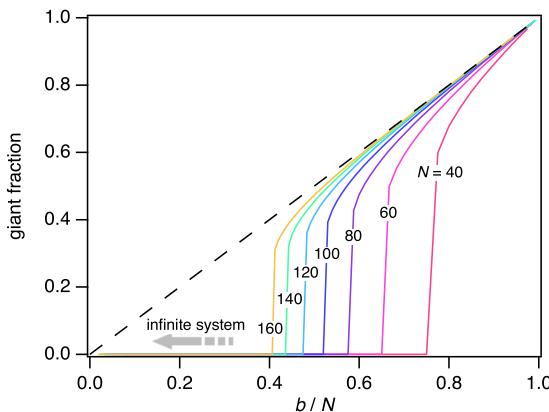


Fig. 14. Effect of system size on the discontinuity of one dimensional models. In the limit $N \rightarrow \infty$, the discontinuity evaporates since the giant cluster forms at $b/N \rightarrow 0$, where b is the number of edges that have been added to the system.

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in small size networks, but it disappears when the size tends to infinite, as shown in Fig. 14. This suggests that in the thermodynamic limit the system actually shows a continuous transition to percolation, a point which will be further discussed in Section 2.7.

Generally speaking, the idea of AP is that of choosing the best among randomly given multiple options, in order to avoid the formation of a certain target pattern [13]. In Ref. [86] the spanning cluster rather than the giant cluster is chosen as the target pattern. Such a latter choice enables one to determine analytically the order of the EP transition in the thermodynamic limit, and in particular for percolation in a Euclidean space. Spanning cluster avoiding models feature either discontinuous or continuous percolation transition, depending on the dimension of the Euclidean space in which the dynamics is implemented. More specifically, the EP is continuous or discontinuous if $d < d_c = 6$, and it is always continuous if $d \geq d_c$ (where d is the spatial dimension), this way explaining some of the confusing and contradictory behavior previously observed [87].

2.6.3. Growth by overtaking

The growth by dominant overtaking has been interpreted as the mechanism responsible for the discontinuous percolation transition of the Bohman–Frieze–Wormald (BFW) model [73]. More specifically, the BFW model has been extended to the case where various constraints exist on the occupation of competitive edges. Such a simple stochastic graph evolution examines only one edge at a time, which leads to a discontinuous transition provided the control parameter β is small enough. The overtaking mechanism means that the size of the largest component does not change by direct growth, but it changes when two smaller components merge and become the new largest component. On the other hand, a continuous transition appears if β is large enough to allow substantial direct growth of the giant component. Notice that the BFW model has been originally proposed to study discontinuous percolation with multiple giant clusters in Ref. [70], and was further studied on finite dimensional square, simple-cubic lattices, and SF networks obtaining similar results [71,72,117,118]. Furthermore, in the generalized BFW model, it has been shown that the emergence of global connectivity is announced by microscopic transitions of the largest component [74].

2.6.4. Recursive and hierarchical structures

Networks with a recursive and hierarchical structure induce explosive percolation dynamics as well, although the formation of an extensive cluster is restricted to ordinary (uncorrelated) and random bond addition [103]. This suggests that a hierarchy of SW bonds grafted onto a one-dimensional lattice results in an EP transition, even if bonds are added sequentially in an uncorrelated manner.

2.7. Debates on explosive percolation

The results of Ref. [9] have triggered a big debate in the community, aiming at definitely clarifying whether the observed transition are discontinuous or not. In equilibrium systems, discontinuous phase transitions are usually identified as first-order transitions, while their continuous counterparts are called second-order transitions. However, the term “explosive percolation” was solely based on the numerical evidence of a jump (a sort of discontinuity from a visualization perspective) in the order parameter at the critical transition point. Therefore, all numerical results in Refs. [33–35,44,48,49,52,56,62,80] initially focused on searching for the existence of such a discrete jump.

2.7.1. Finite-size scaling theory

While the existence of a discontinuous jump in EP was confirmed since the earliest studies on the subject, most other properties of a first-order transition (like cooperation, phase coexistence, and nucleation [12]) had to be proved. From the physical point of view, the competitive choice of a new link suppresses the formation of large clusters, which yields the threshold value significantly delayed in comparison to random percolation. Therefore, the competitive percolation transition can be extremely abrupt with a large (seemingly discontinuous) jump in the connectivity of the graph. Although systems of rather large sizes (i.e., tens of billions of nodes) are used in the original AP model, the heuristic arguments of discontinuity at the percolation transition disappears in the thermodynamic limit. Therefore, the term “explosive” refers here to the unusual feature which distinguishes the critical behavior of the AP from both ordinary and truly discontinuous models [51].

Finite-size scaling theory is useful for characterizing phase transitions [119,120]. Traditionally, the percolation transition has been conceived as a typical continuous phase transition displaying a power law size distribution, and a set of standard scaling properties. Therefore, a percolation transition is continuous if one finds the signature of power law distribution of component sizes. In contrast, it is discontinuous if a power law distribution does not exist.

The first systematic analysis of cluster size distributions for AP by finite-size scaling theory has been reported in Ref. [34], including lattices, random, and scale free networks. For continuous phase transitions, every variable X of interest near the percolation threshold p_c (critical average connectivity of the network in the limit of infinite size) is scale independent, due to the infinite correlation length of the system at p_c , so it has a power-law form

$$X \sim |p - p_c|^\beta, \quad (6)$$

where β is a critical exponent. On a system of a finite size N , the scaling of the variable X is normalized as follows:

$$X = N^{-\beta/\nu} F[(p - p_c)N^{1/\nu}], \quad (7)$$

where ν is another critical exponent, and $F(\cdot)$ is a universal function. Both critical exponents β and ν are associated with the magnetization and the correlation length, respectively. For $p = p_c$, the variable shows a simple scaling $X \sim N^{-\beta/\nu}$, which can be used to deduce β/ν by examining several systems of different sizes. Furthermore, if p_c , β , ν are known, the plot of $XN^{\beta/\nu}$ as a function of $(p - p_c)N^{1/\nu}$ yields the universal function F , which does not depend on N .

In the case of percolation, one chooses X to be the average value $\langle S \rangle$ of the order parameter S , which measures the relative size of the largest connected component with respect to the total system size N (i.e., the number of nodes belonging to the largest connected component divided by N), and which is also called “percolation strength” [11,12,34].

The identification of the percolation threshold p_c is performed in several independent ways [34]. One way is to plot the order parameter $\langle S \rangle$ as a function of the system size N , for a given value of p . A correct value of the percolation threshold can be determined by finding the value of p which yields the best power-law fit. The second method consists in using the scaling of the pseudo-critical points $p_c(N)$, defined as

$$p_c = p_c(N) + bN^{-1/\nu}, \quad (8)$$

where b is a constant to be determined from the fit, together with the other critical parameters ν and p_c . The pseudo-critical value $p_c(N)$ can be simply obtained by the monitoring positions of the peaks of S at different system sizes N .

An alternative quantity that can be adopted as variable X in Eq. (7) is the susceptibility, which again can have two slightly different definitions. The first is to consider the mean cluster size $\chi_1 = \sum_s n_s s^2 / \sum_s n_s$, where n_s stands for the number of clusters of size s per node. The sums run over all possible values of s , except for the one of the largest cluster [34,44]. The second definition is obtained by considering the amplitude of the fluctuations of the order parameter S , i.e. $\chi_2 = N \sqrt{\langle S^2 \rangle - \langle S \rangle^2}$ [44,49]. When these susceptibility measures are used, the critical exponent is indicated by γ , and the scaling ansatz is $N^{\gamma/\nu} F'[(p - p_c)N^{1/\nu}]$, where the universal function F' is different from F .

For ER networks, the order parameter S starts to be defined at the percolation threshold $p_c = 1/2$, and grows with time as $S \sim |p - p_c|^\beta$, with $\beta = 1$. For randomly diluted d -dimensional lattices, β depends on d . At $p = p_c$, the cluster size distribution $n(s)$ (the fraction of finite connected components of s nodes) scales as $n(s) \sim s^{-\tau}$, with $\tau = 5/2$. When the transition is discontinuous, the finite size scaling theory does not apply anymore, and the scaling relation of Eq. (7) trivially holds with $\beta = 0$. The curves of S versus p do not scale as well, and p approaches p_c at increasing N faster than a power law.

Simulations of EP performed for AP-type processes reveal power law distributions of cluster sizes as well, resembling the results of continuous percolation transitions. It turns out that this self-contradicting combination of discontinuity and scaling has made EP one of the most challenging and urgent issues. The first contradicting result has been reported in Ref. [11], which concludes that EP is actually continuous for a modified version of the AP (the dCDGM model), and analytically derives the critical scaling relations based on numerical observations of power-law critical distribution of cluster sizes, with a unique small critical exponent $\beta \approx 0.0555$. The numerical simulation for a very large system (2×10^9 nodes) shows a pronounced power law regime, which indicates a continuous transition (see Fig. 15). The small β exponent makes the transition so sharp that it is almost indistinguishable from a discontinuous one.

Further systematic analysis of finite size scaling was performed for four Achlioptas-type processes [12], showing unusual behaviors with non-analytic scaling functions which demonstrate that AP-like percolation transitions are indeed continuous. More specifically, Ref. [12] characterized the phase transitions in terms of the distribution $P_{p,N}(S)$ of the order parameter S ,

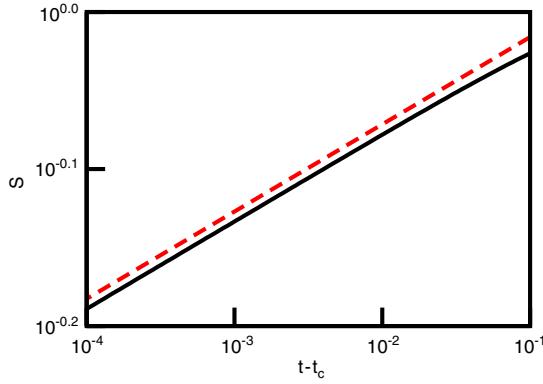


Fig. 15. (Color online). Log–log plot of the relative size S of a percolation cluster vs. $t - t_c$ (black solid curve), showing a power law regime with exponent $\beta = 0.05555$ (red dashed line).

Source: Adapted with permission from Ref. [11]. Courtesy of R. A. da Costa.

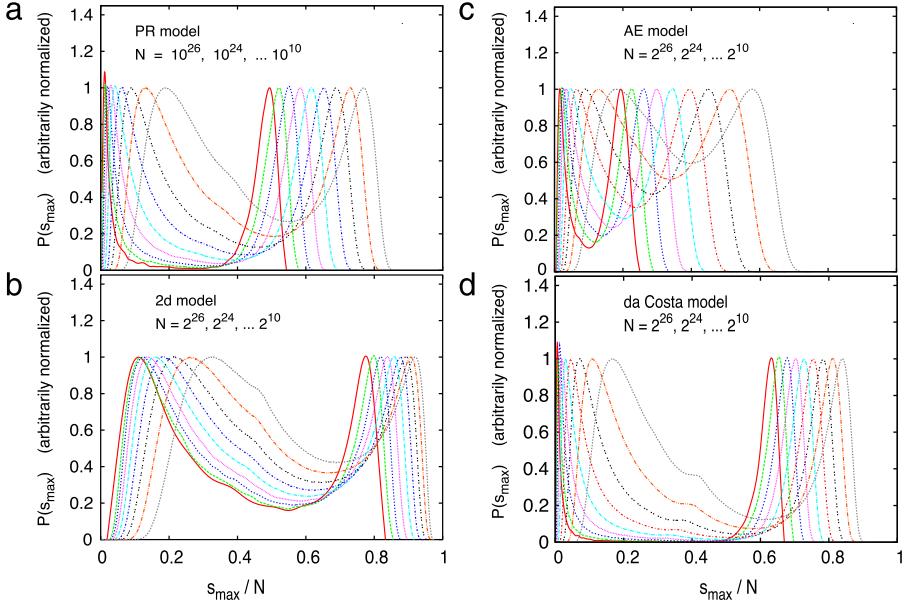


Fig. 16. (Color online). Distributions of the order parameter $S = s_{\max}/N$ for four EP models: (a) the original product rule from Ref. [9]; (b) the product rule on a 2D square lattice; (c) the adjacent edge rule; and (d) the dCDGM rule [11].

Source: Adapted with permission from Ref. [12].

where p is the control parameter and N is the system's size. The distribution $P_{p=p_c,N}(S)$ at criticality scales, for continuous transitions, as

$$P_{p=p_c,N}(S) \sim N^{\beta/\nu} f(SN^{1/\nu}), \quad (9)$$

where f is a universal function. This equation is directly related to the finite size scaling relation of Eq. (7). For a continuous phase transition, the universal function f might be double peaked, but then, as $N \rightarrow \infty$, the dip between two peaks usually does not deepen and the horizontal distance between them shrinks to zero, so that $P_{p=p_c,N}(S)$ becomes single humped. In contrast, in typical first-order transitions, $P_{p=p_c,N}(S)$ is double peaked with a deepening valley between the two peaks. Furthermore, the distance between the peaks tends to be equal to the value of the jump observed in the order parameter S .

The results are shown in Fig. 16. Notice that, while the two peaks seem to indicate an indeed explosive character of the percolation transition, they do not yet prove that EP is discontinuous. For the latter, one must also show that the distance between the peaks does not vanish for $N \rightarrow \infty$. It turns out that the valley between the peaks deepens with increasing N in all cases, but at the same time both peaks shift to the left (see Fig. 16). Ref. [12] also confirms that the critical exponent β is rather small. Furthermore, the distance between the peaks decreases for large N . As a matter of fact, one may extract information on the continuous transition from the depth and position of the valley [121].

In the original AP model with the PR on a complete graph, a careful finite size scaling analysis has shown that EP transition is indeed continuous in the thermodynamic limit [107]. For percolation processes, the main question is whether the gap

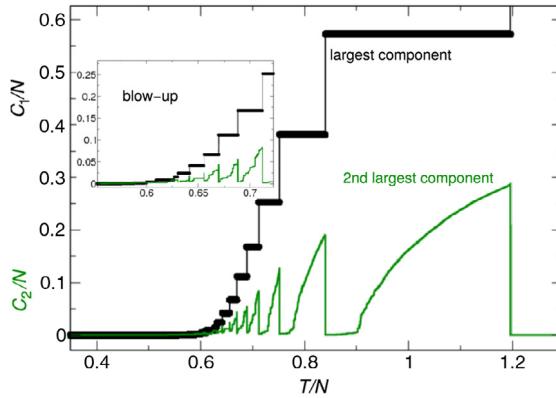


Fig. 17. (Color online). Staircase shape of the multiple discontinuous stochastic process for $N = 655,360$. Source: Figure taken from Ref. [59]. Courtesy of Harold W. Gutch.

$(S|_{p \rightarrow p_c})$ vanishes or not. To this end, the strategy consists in the following steps: (i) the information above the transition point ($p > p_c$) is used to set up lower and upper pseudo-transition points, $p_l(N)$ and $p_u(N)$. This provides an easy way to find p_c , since one would expect that both pseudo-transition points converge to p_c as $N \rightarrow \infty$. (ii) The upper bound for the size increase of the largest cluster between $p_l(N)$ and $p_u(N)$ is found, namely $\Delta G = G_u(N) - G_l(N)$. This is equivalent to consider the gap of the order parameter between the two pseudo-transition points $\Delta S = \Delta G/N$. (iii) Such an upper bound turns out to be sub-linear in N for a continuous percolation transition, and therefore the investigation of the cluster size distribution leads to conclude that the explosive percolation transition is indeed continuous.

Other developments regarding the application of finite size scaling analysis to EP can be found in Refs. [39,44]. For a wide set of representative AP models, an exhaustive investigation has been conducted in Ref. [65], where the full set of scaling relations between critical exponents was found, together with the scaling functions and the upper critical dimension.

2.7.2. Discontinuity of the transition

Finally, the continuity of the EP transition has been rigorously proven in Ref. [13], which indeed mathematically demonstrated that continuity is an essential feature of all mean-field AP models, and of a wide range of their variants. For instance, any competitive rule based on picking a fixed number of edges would lead to a continuous transition. Furthermore, it is shown that the order parameter of the AP percolation models has random fluctuations that do not disappear in the thermodynamic limit [60]. On the other hand, a discontinuous phase transition may appear if the number of nodes sampled grows with the network size [75]. Discontinuous transitions are observed in some models (such as, for instance, the BFW model [70]), and in a hierarchy of SW random bonds model [103]. Notice that the results of Ref. [13] do not imply that the percolation transition at p_c is characterized by a power law divergence of the order parameter, nor they mean that a globally continuous transition cannot be entangled by multiple discontinuous transitions. For instance, for certain l -vertex rules, the continuous percolation transition can have an incomplete devil's staircase with multiple discontinuous steps in the vicinity of p_c [59], as reported in Fig. 17.

2.7.3. Irreversibility of the transition

The last point to be discussed is that EP processes are irreversible, in contrast to traditional continuous percolation. In the latter case indeed, one can reach any state either adding or removing connections. For EP, instead, only adding links makes sense, while the inverse process is impossible [11]. This argument suggests that any efforts in searching hysteresis behavior in EP would fail. Typically, hysteresis is considered to be a necessary condition of a first-order phase transition. For AP processes on 2D square lattices of finite size, hysteresis behavior between the forward (adding edges) and the reverse (removing edges) processes has been numerically observed in Ref. [36]. However, the hysteretic area tends to 0 in the thermodynamic limit, which indirectly shows that the product rule of EP processes generates a continuous transition. In a different model embedded in a Euclidean space (that chooses the spanning cluster instead of the giant cluster as the target pattern) a hysteretic loop has been observed in discontinuous percolation transitions [86].

3. Synchronization as a continuous phase transition

3.1. Introduction and historical overview

Synchronization is possibly one of the most ubiquitous natural phenomena. There, inanimate or living systems [122] tend to adjust their own rhythms due to an existing coupling interaction, or a driving force. It was known for few centuries that two pendulum clocks can synchronize when hanging from the same piece of wood. Since then, synchronization has

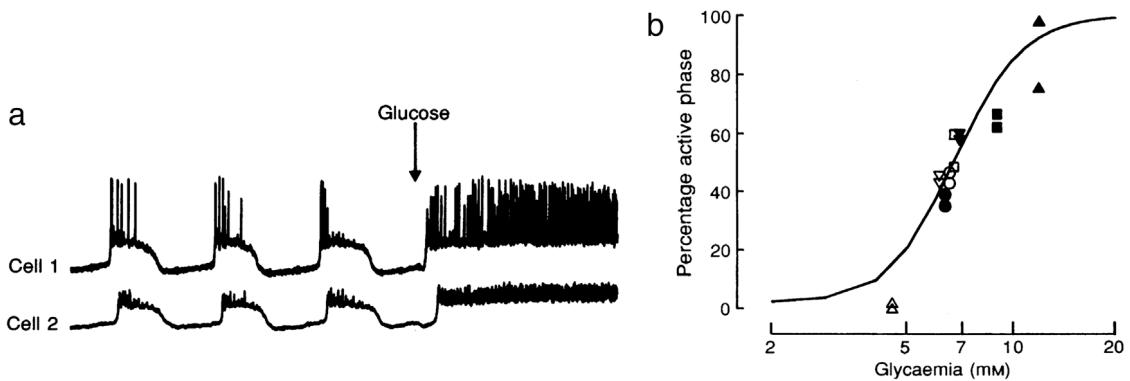


Fig. 18. (a) Simultaneous recordings of the electrical activity of two β -cells within a given islet of Langerhans of an Albino mice's pancreas show synchronous oscillations. (b) Relationship between blood glucose concentration and the percentage of time that pairs of islets spend in the active phase. Source: Reprinted from Ref. [129].

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captivated the interest of scientists and, nowadays, it is investigated in fields as diverse as natural and social sciences [8], or engineering [4]. Popular examples of systems exhibiting synchrony are: birds flocking together, fishes swimming in schools, male fireflies flashing together, people clapping in unison after a performance, pacemaker cells in the heart, β -cells in a pancreatic islet, and neurons in the brain. For comprehensive reviews providing the detailed description of synchronization phenomena we address the reader to Refs. [5–8,123].

When the coupled systems are oscillators, the fate of synchronization always depends on two ingredients: (i) the frequency detuning (the more alike the systems are, the better is synchronization), and (ii) the coupling pattern (i.e. the strength, the network topology, and the output function) through which the network's units communicate. It is not sufficient, indeed, that two systems are alike, it is also necessary that the information exchanged between the systems is of good quality to allow them to synchronize [124–126].

As a coupling strength increases, a transition from an incoherent to a coherent phase takes place generically at a critical value, where the elements follow the same dynamical behavior. The nature of such a macroscopic rhythm can be described by an order parameter as in analogy with phase transitions in statistical mechanics [127]. As first pointed out by Winfree [128], in most of the cases the transition is second-order like, with the order parameter exhibiting power-law divergences at the onset of the synchronization. Fig. 18 shows an example of such a transition in the case of β -cells within a pancreatic islet, which burst all in synchrony, and therefore produce a coordinated insulin release in response to glucose [129].

Abrupt (discontinuous) transitions to synchrony have been recently reported [14,23,130] where an infinitesimal variation of the coupling strength gives rise to a macroscopic synchronous phase exhibiting hysteresis, a feature observed in theoretical models of coupled Josephson junctions [131], as well as in more complex coupled oscillators models [130,132–136]. These abrupt transitions will be the object of the next Section, whereas the present Section has the aim of reviewing the classical works on synchronization, the synchronization types and the main indicators and models used to describe the spontaneous emergence of ordered states.

3.2. Synchronization of two coupled dynamical systems

The case of two coupled dynamical systems constitutes the building block for the later study of networked systems. The problem is properly framed by considering each system as described by a m -dimensional ($m \geq 2$) vector $\mathbf{x}_i(t)$ ($i = 1, 2$). In general, the equations of motion are given by:

$$\frac{d\mathbf{x}_i}{dt} = \mathbf{F}_i(\{\mathbf{x}\}), \quad (10)$$

where $\{\mathbf{x}\} \equiv (\mathbf{x}_1, \mathbf{x}_2)$. Each unit evolves according to the velocity field \mathbf{F}_i , and can reach either a stable fixed point, or a limit cycle, or a chaotic attractor [122,137].

Eq. (10) accounts for the general framework in which the two systems can be of different nature, have different dimensions, and the interaction can be described by any generic function of their dynamical states. However, in order to allow an analytical treatment of the problem, the autonomous dynamics is assumed to be independent from the interaction between the units, and a particular form of interaction has to be assumed. We here focus on the case in which each unit is coupled to a linear superposition of the outputs, and the interaction is symmetrical:

$$\frac{d\mathbf{x}_i}{dt} = \mathbf{f}_i(\mathbf{x}_i) + \lambda \sum_{j=1}^N A_{ij} \mathbf{h}(\mathbf{x}_j), \quad (11)$$

where λ represents the coupling strength, \mathbf{f} and \mathbf{h} are vectorial functions, and A_{ij} is a matrix satisfying the rules of a diffusive coupling ($\sum_j A_{ij} = 0$). For two units, one has therefore $\dot{\mathbf{x}}_{1,2} = \mathbf{f}_{1,2}(\mathbf{x}_i) + \lambda[\mathbf{h}(\mathbf{x}_{2,1}) \mp \mathbf{h}(\mathbf{x}_{1,2})]$.

3.2.1. Complete synchronization

If the systems are identical ($\mathbf{f}_i(\mathbf{x}_i) = \mathbf{f}(\mathbf{x}_i)$), the synchronization is realized by the equality of the state variables as they evolve in time ($\{\mathbf{x}_i(t) = \mathbf{s}(t), i = 1, 2\}$), and is termed as *complete* (or *identical* or *full*, or *conventional*) synchronization [5,138,139].

Complete synchronization (CS), $\mathbf{x}_1 = \mathbf{x}_2$, is a threshold phenomenon and occurs beyond a critical value λ_c of the coupling strength. CS can be observed by monitoring the mean synchronization error, defined as:

$$\langle e \rangle = \lim_{T \rightarrow \infty} \frac{1}{T} \int_0^T \|\mathbf{x}_1(t) - \mathbf{x}_2(t)\| dt \quad (12)$$

i.e., the averaged distance to the synchronization manifold $\mathbf{x}_1(t) = \mathbf{x}_2(t)$.

One of the most popular approaches to assess the stability of CS is the *master stability function* (MSF) method, introduced for arrays of coupled systems in Ref. [140]. The MSF is the largest Lyapunov exponent transverse to the synchronization manifold, and measures the exponential rate at which any infinitesimal perturbation decays or grows. Therefore, a necessary condition for CS to occur is that the MSF is negative.

3.2.2. Phase synchronization

Real systems, however, are not identical, and they are subjected in general to noise or parameter mismatch preventing CS. Weaker forms of synchronization exists, where correlation or equality of the two systems are limited to proper subsets, or functions, of the variables.

When two periodic oscillators are coupled, for instance, synchronization is observed as the entrainment of their phases and frequencies. When the systems are uncoupled, the two stable limit cycles are described by two phases $\theta_{1,2}$ whose dynamics are governed by their respective natural frequencies ω_1 and ω_2 as $\dot{\theta}_{1,2} = \omega_{1,2}$, which in general are incommensurate ($\Delta\omega = n\omega_1 - m\omega_2 \neq 0, \forall n, m \in \mathcal{N}$). Phase locking means here that the phase difference ($|n\theta_1 - m\theta_2|$) is bounded beyond a certain coupling strength, and the corresponding instantaneous frequencies $v_i = \dot{\theta}_i$ are locked, i.e., $n\nu_1 - m\nu_2 = 0$.

The onset of synchronization is therefore a transition from a quasi-periodic motion (with two incommensurate frequencies) to a motion with a single frequency, and can be characterized by means of families of Arnold tongues [141], i.e. synchronization regions where the relationship $|\frac{\Delta\omega}{\lambda}| \leq 1$ holds. Inside those regions, the oscillators mutually adjust their frequencies until attaining the relation $n\nu_1 = m\nu_2$, with the resulting synchronization frequency depending on the initial frequency detuning $\Delta\omega$. Moving away from the border of an Arnold tongue, the phase difference uniformly grows in time.

Phase synchronization has been later extended to chaotic oscillators [142]. When the chaotic oscillators have a well defined center of rotation, as it is the case of phase coherent systems, the instantaneous phase can be reliably calculated by means of several mathematical techniques [5]. The mean square displacement of the phase linearly diffuses in time (with a very small diffusion constant), due to the pronounced variations of the amplitude modulating the rotations. As two phase coherent chaotic oscillators are coupled, a transition from a non-synchronous regime (where the phase difference linearly grows in time) to a synchronous regime (where the phase difference $|\theta_1 - \theta_2|$ is bounded and the mean frequencies are equal) occurs.

When instead the two chaotic oscillators are not phase coherent (and exhibit a broad distribution of time scales), phases do not fully synchronize, and synchronization episodes are interrupted by intermittent phase slips (jumps of 2π in the phase difference). Such a latter phenomenon has been called *imperfect phase synchronization* [143] and it has been observed experimentally in an electronic circuit representing a forced Lorenz system [144].

3.2.3. Lag, anticipated, relay and generalized synchronization

After the setting of a phase synchronization, intermediate values of the coupling strength drive the systems into a stronger degree of synchronization, where also amplitudes start to become correlated. The resulting regime corresponds to the dynamics of one of the oscillators which lags (in *lag synchronization* [145]) or leads (in *anticipated synchronization* [146,147]) the dynamics of the other. Lag synchronization was first shown in symmetrically coupled nonidentical oscillators [145], and the time shifting τ between one and the other state's variable was detected as the minimum of a time dependent similarity function.

The existence of this *achronal synchronization* in mutually delayed-coupled oscillators has been observed in laser systems [148], and studied theoretically in Ref. [149]. It is possible also to have two completely synchronized chaotic units in the presence of a delayed coupling, phenomenon known as *zero-lag* or *isochronal synchronization* [150,151]. Such a regime can be achieved via a self-feedback in both coupled units [152,153], or via an indirect coupling through a relay [154], or through a third driving unit [155–157], or when multiple coupling delay times are used [158].

In the so called *relay systems*, two units achieve CS as a result of an indirect coupling through a third system acting as a relay. Experimental evidence of relay synchronization has been obtained with lasers [150,154,159] and electronic circuits [155,157]. Relay synchronization has been further used as a method for chaos-based communications [151,153,156,160], as well as a synchronizing mechanism among remote neuronal resources [161–163].

Anticipated synchronization, on the other hand, may occur in systems coupled in a drive-response configuration, and it relies on the use of time delay lines in the dynamics of the slave system. It has been studied theoretically and numerically in different contexts including dissipative chaotic flows [146,147,164], chaotic maps [165], excitable spiking neurons [166–169], inertial ratchets [170], and complex Ginzburg–Landau equation [171]. Experimentally, it has been observed in electronic circuits [166,172–174] and in optical systems [175,176]. Furthermore, the fact that behavior of one of the oscillator can be forecasted up to time scales larger than those characteristics of the system's dynamics has been used recently as a method to control and suppress unwanted spikes in excitable systems [177], or to anticipate or suppress extreme desynchronization events [174].

A further form of synchronization, called *generalized synchronization* [178,179], involves the existence of a time-independent functional relationship between the states of the two coupled units (typically nonidentical), such that $\psi(\mathbf{x}_1, \mathbf{x}_2) = 0$. Initially, this type of synchronization was found for unidirectional coupling schemes and proved experimentally by means of the auxiliary system method [180,181]. Recently, generalized synchronization has been reported also for bidirectionally coupled oscillators [182–184]. In particular, Gutiérrez et al. provided evidence that relay synchronization corresponds in fact to the setting of generalized synchronization between the relay and the outer units [184].

Finally, for very large coupling or parameter mismatches a counterintuitive phenomenon may emerge that suppresses the oscillations, termed as *oscillation quenching*. Such a phenomenon can be manifested in two distinct ways, whose origin is structurally different: amplitude death and oscillation death. In the former case, the suppression of oscillations results in the existence of a homogeneous steady state, in the latter a homogeneous unstable steady state splits, giving rise to an inhomogeneous state with two coexisting steady states. Refs. [185,186] are two very recent reviews on the subject.

3.3. Synchronization in ensembles of coupled systems

The scenario described in Section 3.2 can be generalized to ensembles of spatially ordered oscillators, coupled (either locally or globally) in chains and lattices, as well as to spatially extended oscillatory media. Synchronization in these contexts, indeed, is very relevant for the study of cooperative behavior in many physical [187], biological [128,188,189] and chemical systems [23,190–192]. For a review, the reader is addressed to the books in Refs. [8] and [193].

3.3.1. Identical systems

Assemblies of coupled identical systems can exhibit non trivial synchronization patterns. Precisely, two types of synchronous behavior can be produced: *full* or *global* synchronization where all elements evolve identically in time, and *partial* or *cluster* synchronization in which groups of mutually synchronized systems with different synchronization patterns emerge. The specific scenario that is produced (and its features) depends on many factors: the number of units, the way they are mutually coupled, and the initial and boundary conditions.

Formally, the coupling conditions may be classified as local (by setting in Eq. (11) $A_{ij} = 1$ for $j = i \pm 1$, and zero otherwise), global ($A_{ij} = 1/N \forall i \neq j$), and non-local ($A_{ij} = g(|i - j|)/N$, where g is a decreasing function of the distance between oscillators i and j).

From now on, we will focus mostly on the case in which each unit of the ensemble is an oscillator. In such a case, the amplitude of the mean field is taken as an indicator of the onset of coherence. If all the frequencies in the ensemble are different, then the instantaneous phases θ_i are uniformly distributed in the interval $[0, 2\pi]$, and the mean field vanishes. If instead some of the oscillators are locked to the same frequency, the vectors associated to their phases sum up coherently in the unit circle, and the mean field is non-zero. A weaker condition of synchronization is the coincidence of the averaged frequencies $\Omega_i = \langle \dot{\theta}_i \rangle$.

Ensembles of locally *identical* coupled systems have been studied through chaotic maps [194–196] and flows [179,197–200], and limit-cycle oscillators. An example is the complex Ginzburg–Landau equation [201,202] which represents the normal form of an Hopf bifurcation, and describes a wide variety of wave phenomena, including phase turbulence (spatiotemporal chaos), plane waves, targets and spirals in which oscillations are synchronized [201]. As for globally coupled systems, the main studies have concerned chaotic maps [203–206] and flows [197,207,208], rotators with diffusive [136,209–211] or pulsed [212–214] coupling. Experiments instead have been conducted with chaotic electrochemical oscillators [215].

Regardless on the specific framework that is considered, a transition is generically found from global (at large coupling) to cluster (at intermediate coupling) synchronization. Detailed analysis of the mechanisms determining the loss of stability of the global coherent state into a hierarchical cluster dynamics can be found in Refs. [205,206], while Ref. [216] studies the destabilization of the synchronous state for increasing coupling due to the so called short-wavelength bifurcation. Ref. [204] analyzed the cluster structure emerging in a system of randomly coupled identical maps, as a function of the average connectivity.

3.3.2. Chimera states

Together with the emergence of clustered synchronous patterns, an even more surprising phenomenon is the coexistence of clusters of coherent and incoherent phases in ensembles of identical oscillators, which was termed as *chimera state* by Abrams et al. in 2004 [217]. As an example, Fig. 19 shows the coexistence of a synchronized cluster of 64 oscillators

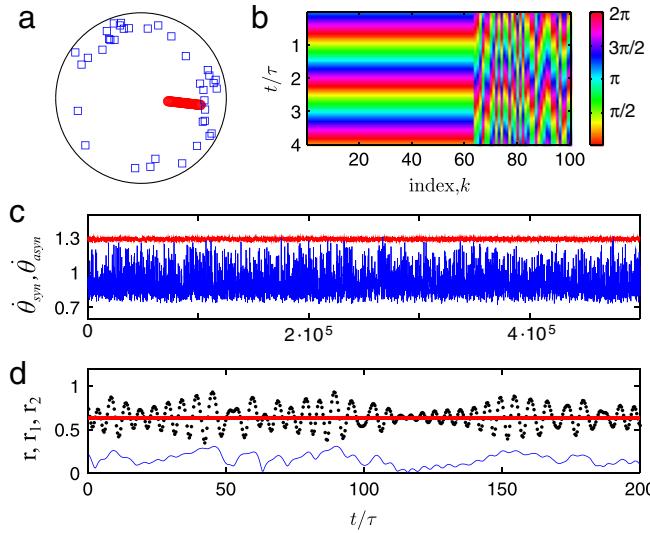


Fig. 19. Chimera states in a homogeneous ensemble of globally coupled oscillators with internal delayed feedback loop [$\dot{\theta}_k = \omega + \lambda \sin(\theta_{\tau,k} - \theta_k) + \epsilon \text{Im}(e^{i\beta} Ze^{-i\theta_k})$]. (a) Snapshot of the phases of the oscillators belonging to the cluster (red circles) and to the cloud (blue squares). The amplitude of the phasor in the circle is set proportional to the oscillator index k to improve visibility. (b) Colormap of the instantaneous phases. (c) Instantaneous frequencies of an oscillator from the cluster (upper red curve) and of an oscillator from the cloud (lower blue curve). (d) Amplitude of the mean fields of the cluster (r_1 , red thick line), of the cloud (r_2 , blue thin line), and of the whole ensemble (r , black dotted line). (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

Source: Adapted with permission from Ref. [211]. Courtesy of M. Rosenblum.

and a cloud of 36 asynchronous ones, which emerge in a homogeneous ensemble of globally coupled oscillators, due to an internal delayed feedback in individual units [211]. First noticed by Battogtokh and Kuramoto [218], the breakdown of spatial coherence in the presence of non-local interaction has been reported theoretically [217,219–222], and numerically [223,224]. Experimental evidences of chimera states have been obtained in chemical [225,226], optical [227] and mechanical [228] systems. For instance, in Fig. 20, a chimera state is observed in a photosensitive Belousov–Zhabotinsky reaction, in which the catalyst is loaded onto N ion-exchange oscillatory particles. The N oscillatory particles are divided into two groups of equal size (A and B), with each member i of group π experiencing the feedback light intensity P_i^π :

$$P_i^\pi = P_0 + k_\pi [\bar{I}_\pi(t - \tau) - I_i(t)] + k_{\pi\pi'} [\bar{I}'_\pi(t - \tau) - I_i(t)],$$

being $\pi = A, B$ and $\pi' = B, A$. \bar{I}_π is the mean intensity in group π , k_π and $k_{\pi\pi'}$ are the intra and inter-group coupling coefficients, P_0 is a normalized background light intensity, and I is the measured intensity scaled from 0 to 1. The delay τ plays the role of the phase frustration term used in the model by Abrams et al. [217]. With such a coupling scheme, group A remains fully synchronized, while B can exhibit three basic states: fully synchronized (Fig. 20(a)), n -cluster ($n = 2$ in Fig. 20(b)) and the unsynchronized chimera (Fig. 20(c)). Panel (d) shows the behaviors found at different intra- and inter-group coupling strengths, under the setting of $k_A = k_B$ and $k_{AB} = k_{BA}$.

Coexistence of coherence and incoherence has also been observed in networks of coupled nonidentical phase oscillators [229,230], and in globally coupled populations of identical oscillators [231] with internal delayed feedback [211,219]. A recent review on the subject [232] gives a comprehensive history and overview on chimera states.

3.3.3. Non-identical systems

Real-world systems often feature noise and various degrees of heterogeneity. Therefore, synchronization phenomena have been explored in large ensembles of interacting non-identical oscillators [23].

The same routes of two coupled oscillators can also be explored for ensembles of locally or globally coupled (periodic or chaotic) systems. As long as oscillators are phase coherent, there is no qualitative difference between periodic and chaotic case [233–235], with the exception that chaotic assemblies can display phase synchronization while their amplitudes remain uncorrelated.

The transition to synchronization in chains of periodic [234,236] and phase coherent chaotic [233] oscillators with different frequencies can occur in a soft or a hard way, depending on the frequency mismatch of the oscillators. Namely, for a linear distribution with a small mismatch, the transition is smooth, while for large mismatches the transition is characterized by an abrupt formation of clusters with phase slips between them and whose size is increasing with the coupling. In the case of randomly distributed frequencies, only the hard transition is possible, and the onset of synchronization appears for lower values of the coupling.

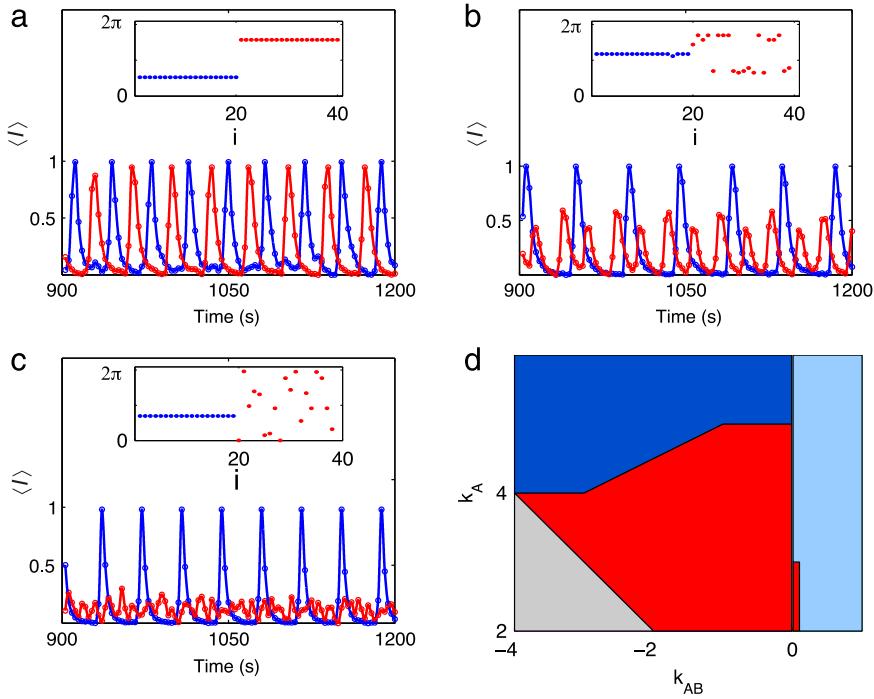


Fig. 20. Chemical chimeras in the photosensitive Belousov–Zhabotinsky (BZ) reaction. (a)–(c) Time evolution of the normalized mean intensity $\langle I \rangle$ measured in the BZ particles of groups A (blue) and group B (red), and snapshot of the phases of the oscillators in each group (insets). (a) Synchronized out-of-phase state, $k_A = 3.0$, $k_{AB} = -1.8$; (b) synchronized 1–2 phase cluster state, $k_A = 4.0$, $k_{AB} = -1.0$; (c) chimera state, $k_A = 3.0$, $k_{AB} = -0.2$. (d) Regions of chimera (red), in-phase (light blue), and out-of-phase (dark blue) synchronized states as a function of k_A and k_{AB} . (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)
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3.3.4. The Kuramoto model

Despite the peculiarity of the system under study, a large population of weakly interacting oscillators can be described by considering their phases, and assuming a homogeneous interaction among them. With such an assumption, Kuramoto formulated his celebrated equation [23,187] (based on the previous Winfree's works with phase models [128]) which transforms Eq. (11) into a mathematically tractable one.

The Kuramoto model [23,187] considers a system of N oscillators having natural frequencies ω_i distributed according to a probability density $g(\omega)$ which (for the time being) is uni-modal and symmetric around its mean frequency $\Omega(t=0) = \sum \omega_i/N$. The instantaneous phases θ_i obey the following equations:

$$\dot{\theta}_i = \omega_i + \frac{\lambda}{N} \sum_{j=1}^N \sin(\theta_j - \theta_i) \quad (13)$$

where the normalization constant $1/N$ ensures the boundedness of the coupling term in the thermodynamic limit $N \rightarrow \infty$. A transition to synchronization occurs for values of the coupling λ above a critical threshold λ_c , which depends on the distribution $g(\omega)$. A motivating survey of the Kuramoto model is offered in Ref. [237]. Tutorials and reviews of the main achievements on the Kuramoto model can be found in Refs. [238–240].

The monitoring of the synchronization onset is performed by introducing a complex mean field of the population, known as the *order parameter* [187]

$$r(t)e^{i\psi(t)} = \frac{1}{N} \sum_{j=1}^N e^{i\theta_j(t)}. \quad (14)$$

Eq. (14) is the vector centroid of all oscillators represented as points in the unit circle. ψ is the average phase, and r its modulus, which implies $r \sin(\psi) = \sum_j \sin(\theta_j)$ and $r \cos(\psi) = \sum_j \cos(\theta_j)$. r measures the level of coherence in the population: if $r \sim 1$ the population is synchronized, and if $r \sim 1/\sqrt{N}$ the points are uniformly distributed in the unit circle.

Using the parameters r and ψ , Eq. (13) can be rewritten as

$$\dot{\theta}_i = \omega_i + \lambda r \sin(\psi - \theta_i). \quad (15)$$

Now, the dynamics of each oscillator is implicitly coupled to all the others through the mean field in a positive feedback way: the more coherent is the population, the larger is r and the more effective is the coupling strength λr , which drives therefore more and more oscillators into the synchronization cluster. The threshold value λ_c can be derived by self consistency conditions applied to Eq. (15). If one assumes a priori a solution where $r(t)$ is constant and ψ is rotating at a constant frequency Ω ($\psi = \Omega t$), then Eq. (15) can be expressed as a function of the angle $\phi_i = \theta_i - \psi$ using a rotating reference frame at frequency Ω :

$$\dot{\phi}_i = \omega_i - \Omega - \lambda r \sin \phi_i. \quad (16)$$

The scalar vector field $f(\phi) = \omega - \Omega - \lambda r \sin \phi$ for $\phi \in [0, \pi]$ exhibits a saddle-node bifurcation at $\lambda r = |\omega_i - \Omega|$ or $\phi_i = \pi/2$. If $|\omega_i - \Omega| \leq \lambda r$, the solution $\phi_i = \sin^{-1} \frac{\omega_i - \Omega}{\lambda r}$ is stable, and oscillators are locked to the mean frequency Ω . The oscillators instead with $|\omega_i - \Omega| > \lambda r$ are uniformly spread around the unit circle.

A rigorous analysis of Eq. (16) has been performed in Refs. [237,238]. Here we just highlight the main steps in such an analysis. By introducing the probability density function $\rho(\phi, t, \omega)$, the fraction of oscillators lying within the arc $[\phi_a, \phi_b]$ with natural frequencies $\omega \in [\omega_a, \omega_b]$ is

$$\int_{\phi_a}^{\phi_b} \int_{\omega_a}^{\omega_b} \rho(\phi, t, \omega) g(\omega) d\omega d\phi,$$

and the order parameter can be expressed as:

$$r(t) e^{i\psi(t)} = \int_0^{2\pi} \int_{-\infty}^{\infty} e^{i\phi} \rho(\phi, t, \omega) g(\omega) d\omega d\phi. \quad (17)$$

$\rho(\phi, t, \omega)$ must obey the continuity equation

$$\frac{\partial}{\partial t} \rho + \frac{\partial}{\partial \phi} (\rho v) = 0, \quad (18)$$

where $v = \omega - \Omega - \lambda r \sin(\phi)$ is the angular velocity of a given oscillator with phase ϕ (in the co-rotating frame) and natural frequency ω . One obtains therefore the following self-consistency condition

$$r = \lambda r \int_{-\pi/2}^{\pi/2} \cos^2 \phi \cdot g(\Omega + \lambda r \sin \phi) d\phi. \quad (19)$$

The saddle-node bifurcation manifests as a phase transition from the incoherent state (at $r \sim 0$ with density $\rho = 1/2\pi$) to the partially synchronized state (with $0 < r < 1$) characterized by a group of phase oscillators rotating in unison at Ω (those whose frequencies obey $|\omega - \Omega| \leq \lambda r$), while the remaining ones are rotating incoherently. These solutions satisfy

$$1 = \lambda \int_{-\pi/2}^{\pi/2} \cos^2 \phi \cdot g(\Omega + \lambda r \sin \phi) d\phi. \quad (20)$$

By expanding the integrand $g(\Omega + \lambda r \sin \phi)$ in a Taylor series, one finds that the amplitude of the mean field r (at the onset of synchronization) is given by

$$r^2 \sim \frac{-16}{\pi \lambda_c^4 g''(\Omega)} (\lambda - \lambda_c), \quad (21)$$

with

$$\lambda_c = \frac{2}{\pi g(\Omega)}. \quad (22)$$

Therefore, the transition is continuous and the order parameter behaves (in the limits $N, t \rightarrow \infty$) as

$$r = \begin{cases} 0 & \lambda < \lambda_c \\ (\lambda - \lambda_c)^\beta & \lambda \geq \lambda_c \end{cases} \quad (23)$$

which resembles a continuous second-order phase transition characterized by the critical exponent $\beta = 1/2$.

Finally, exploiting the extensive symmetries of the continuum-limit model (17)–(18), Ott and Antonsen derived an exact reduction to a finite set of nonlinear differential equations for the evolution of the order parameter [241,242]. This latter approach has been used in several recent studies [221,229,243,244].

3.4. Synchronization in complex networks

Units in the ensemble may, however, interact through specific coupling patterns (or connection topologies) which can be conveniently encoded in an interaction network. The seminal paper by Watts and Strogatz [245] proved that the connection topology of networks can be organized in between a regular and a random architecture, and that such a topological arrangement (called SW property) endows the system with enhanced synchronizability. Since then, a flurry of studies has been focused on synchronization in complex networks, and the reader is here addressed to a series of classic [246–248] and recent [239,249] reviews on the subject.

We here follow the same path of the previous two Sections, and provide a schematic overview of synchronization on networks, for both cases of identical and non-identical oscillators.

3.4.1. Identical systems

Complete, cluster, and chimera states of synchronization have been reported so far in networks of identical flows or maps. As far as complete synchronization is concerned, one of the main research lines has been the investigation on the stability of this state for different network topologies and dynamical functions.

For a generic network of N diffusively coupled identical units there are two main approaches to assess the stability of the synchronous state. The first one is the extension of the MSF [140,250,251] to complex networks [252], which related the local stability of the synchronization manifold $\mathbf{s}(t) = \mathbf{x}_1(t) = \mathbf{x}_2(t) = \dots = \mathbf{x}_N(t)$ to the spectral properties of the Laplacian matrix associated to the underlying network structure. A comprehensive derivation of this formalism can be found in Ref. [246]. Precisely, the analysis reduces to plot the maximum Lyapunov exponent Λ_{\max} transverse to the synchronization manifold as a function of a generic coupling parameter μ which encodes both the local function $\mathbf{f}(\mathbf{x})$ and the coupling function $\mathbf{h}(\mathbf{x})$ in Eq. (11). Different choices of local and coupling functions give rise to three possible classes of MSF. For instance, for a bounded MSF (the so called case of class-III MSF), where $\Lambda_{\max}(\mu) < 0$ within a finite interval (μ_1, μ_2) , the necessary condition for stability can be attained only if $\frac{\lambda_N}{\lambda_2} < \frac{\mu_2}{\mu_1}$, being λ_2 and λ_N the minimal and maximal non-zero eigenvalues of the Laplacian matrix. The interested reader is here referred to Ref. [246] for the complete discussion on the three possible classes of MSF.

The second method is known as *the connection graph stability* method, and was derived in Ref. [253]. It combines the Lyapunov function approach with graph theoretical reasoning, and provides a bound for the minimum coupling strength needed for synchronization which explicitly depends on the average path length of the coupling graph.

MSF arguments led to the study of the propensity towards synchronization of different topologies regardless on the nature of the coupled dynamical systems. Ref. [254] introduced a family of graphs, called entangled networks, having an extremely homogeneous structure which optimizes synchronizability: degree, node distance, betweenness, and loop distributions are all very narrow. Therefore, synchronization is shown to be affected by the heterogeneity of the degree distribution, clustering, characteristic path length or betweenness.

Ref. [252] compared the synchronizability of a SW network against an interpolated ordered ring structure, and a completely random ER graph, as implemented in Ref. [245]. It was shown that SW graphs exhibit better synchronizability than regular and completely random ones, when the system is far from the complete graph limit. In contrast, Ref. [255] showed that the onset and stability of complete synchronization in networks of pulse coupled oscillators (neurons) only depends on the number of inputs a neuron receives regardless on the details of the network topology, as long as the degree distribution is a δ function. The effect of link directionality in SW networks has been studied in Ref. [256].

Ref. [257] further elaborated on the role of heterogeneity, and realized that heterogeneous structures may display very large ratios λ_N/λ_2 (λ_N scales with the hub's degree, i.e. the largest degree in the graph [258]). The fact that heterogeneous networks may be more difficult to synchronize is known as the *paradox of heterogeneity* [257]. The effect is due to the overloading of the hubs, which may receive independent inputs and may hamper effective communication between oscillators. The paradox stimulated many scientists to explore mechanisms to improve synchronizability, by the implementation of asymmetric weighting procedures. In Refs. [259,260], enhanced synchronization is obtained by normalizing the coupling to the power of the node degree k_i (i.e. $\frac{\lambda}{k_i^2} \sum_j A_{ij} \mathbf{h}(\mathbf{x}_j)$), while in Ref. [261] the coupling strength between two nodes is re-scaled to the load of the edge connecting them through the edge's betweenness centrality b_{ij} (i.e. $\frac{\lambda}{\sum_j b_{ij}^2} \sum_j b_{ij}^2 A_{ij} \mathbf{h}(\mathbf{x}_j)$). Both methods implement an asymmetric arrangement of the coupling, and promote a dominant interaction from the hubs to the non-hubs nodes [262]. This consideration can be straightforwardly generalized to weighted networks, by showing that synchronizability decreases when the node's strengths become more heterogeneous [263].

The existence of degree correlations may also affect the propensity for synchronization [264]. In particular, synchronizability is enhanced in disassortative SF networks when smaller nodes drive larger degree ones, while the opposite holds for assortative networks if the hubs are driving the small degree nodes [265,266]. Synchronizability is therefore enhanced if hubs are strongly inter-connected and drive the nodes with smaller degrees, that is, if the network displays a rich club of large degree nodes.

At a more microscopic scale, Ref. [268] showed that the stability of the synchronous states supported by 3-node and 4-node motifs is correlated with their relative abundance in undirected networks, while in directed graphs the correlation only holds for some specific motifs. Recently, stability of synchronization has been investigated also in relationship with topological criteria which apply to network's meso-scales [269].

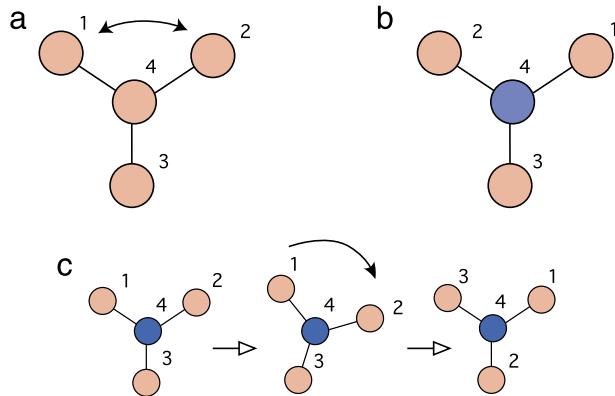


Fig. 21. (Color online). Relationship between group symmetry and dynamics in networks of oscillators. (a) A star network of four identical oscillators. (b) Nodes 1 and 2 are reflected. (c) The same network in (a) after a rotation operation.

Source: Adapted from Ref. [267].

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Another extension is that covered by adaptive networks [270], whose structure is shaped by the microscopic organization patterns [263,271,272]. To achieve global synchronization, one plausible assumption is to consider that each node tries to synchronize to its neighbors by increasing the coupling strength from them. Inspired by the anti-Hebbian learning rule in neural systems, Ref. [272] studies how the feedback from dynamical synchronization shapes network structure by adding, at every time step, a link between two randomly chosen nodes i and j with a probability proportional to the rate of their dynamical differences with the mean activity of the network. The evolution of the network spontaneously forms a SF structure with negative degree-degree correlation, a typical feature of technological and biological networks.

On the other hand, Refs. [263,271] introduced a dynamical feedback principle by means of a nonlinear differential equation for the evolution of the weights which is coupled to the dynamics of the nodes while the connectivity matrix is fixed a priori. In this latter case, when complete synchronization is achieved, the coupling strength becomes weighted and correlated with the topology due to a hierarchical transition to synchronization in heterogeneous networks.

MSF approach has also inspired some pinning control strategy for synchronization [273,274] where nodes are sequentially perturbed to improve as much as possible the eigenratio of λ_N/λ_2 .

More recently, several studies revealed that the symmetry groups in the architecture of complex networks can determine constraints for the appearance (or stability) of a given cluster solution [275]. Cluster synchronization has attracted considerable attention in SW networks of coupled maps [250], chaotic flows [276], synaptically coupled networks of bursting neurons [277] (for diffusively coupled neurons only complete synchronization is possible [255]), or in networks of Stuart-Landau oscillators [278].

In particular, Refs. [267,279] have studied the emergence and stability of cluster synchronization in networks of coupled oscillators for arbitrary topologies and individual dynamics. Using the tools of computational group theory, Ref. [279] addressed the general case of networks where the intrinsic symmetries are neither *ad hoc* produced nor easily observed. Fig. 21 schematically describes a case where a strong relationship exists between symmetry and dynamics: a symmetry (a permutation matrix like a reflection in Fig. 21(b) or rotation in Fig. 21(c)) applied to an adjacency matrix leaves it unchanged (i.e. the equations of motion are the same when mapped into each other by the symmetry transformation). In the example, nodes 1, 2, and 3 constitute one cluster and node 4 another one. The main result is a technique, based on the cluster variational equations, to evaluate the stability of all the dynamically valid cluster synchronization patterns. The conclusion is that the range of stability typically becomes smaller for those synchronization clusters that are characterized by higher symmetry. Moreover, the validation of such patterns is performed by an optoelectronic experiment on a five-node graph. The same approach was independently developed in Ref. [280], and recently an application for the control of specific clusters along the synchronization path has been reported in Ref. [281] (see Fig. 22).

Ref. [282] also relates network's symmetries with the emergence of cluster synchronization states in a Kuramoto model of identical oscillators, when the interactions include a phase frustration preventing full synchronization. In these conditions, the system organizes into a regime of remote synchronization where pairs of nodes with the same network symmetry are fully synchronized, despite their distance on the graph. An application to brain networks suggests that anatomical symmetry may play a role in neural synchronization by determining correlated functional modules across distant locations.

In the framework of networks of communities [283] and networks of networks [249], there are several works seeking for the synchronization performance as a function of the competition between individual communities and the whole network [284–286]. In particular, Ref. [286] discusses how the degree of the nodes through which two networks are connected influences the ability of the whole system to synchronize. It is shown that connecting the high-degree (low-degree) nodes of each network turns out to be the most (least) effective strategy to achieve synchronization, and the existence of the optimal connector link weights for the different interconnection strategies is reported. On the other hand, Ref. [285] reaches similar

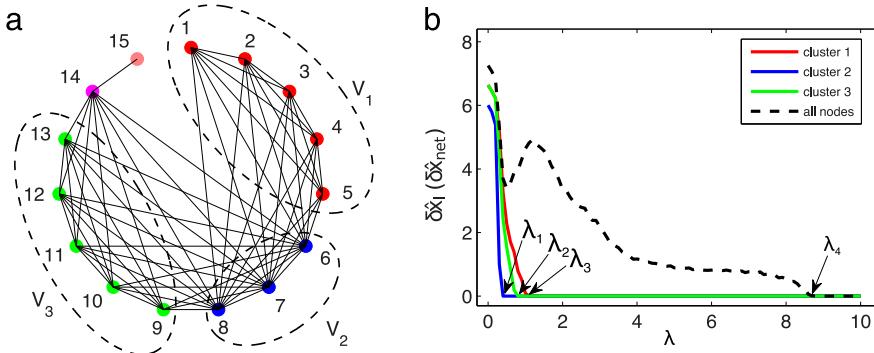


Fig. 22. Cluster synchronization in the Nepal power-grid network. (a) The structure of the Nepal power-grid consisting of 15 nodes (power stations) and 62 links (power lines). Computational group theory partitions the nodes into 5 clusters [$V_1 = \{1, 2, 3, 4, 5\}$ (red), $V_2 = \{6, 7, 8\}$ (blue), $V_3 = \{9, 10, 11, 12, 13\}$ (green) and two trivial clusters of just one node $V_4 = \{14\}$ (pink) and $V_5 = \{15\}$ (yellow)]. (b) Using Lorenz chaotic oscillators as nodal dynamics and coupling function $\mathbf{h} = (0, x, 0)$, the cluster synchronization error $\delta\hat{x}_l$ ($l = 1, 2, 3$) and the network synchronization error $\delta\hat{x}_{\text{net}}$ (black dashed line) are reported as functions of the coupling strength λ . Clusters 2, 3, and 1 are synchronized at $\lambda_1 \approx 0.4$, $\lambda_2 \approx 0.8$, and $\lambda_3 \approx 1.1$, respectively. For $\lambda > \lambda_4 \approx 8.9$, the network is globally synchronized. $\delta\hat{x}_l = \sum_i (|x_i - \bar{x}_l|)/n_l$, with $\bar{x}_l = \sum_i x_i/n_l$ being the averaged state of the oscillators in cluster l ; $\delta\hat{x}_{\text{net}} = \sum_i (|x_i - \bar{x}|)/N$, with $\bar{x} = \sum_i x_i/N$ being the network averaged state. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

Source: Adapted with permission from Ref. [281]. Courtesy of X. Wang.

conclusions, and observes also intermediate states of inter-community connectivity for which there is a balance between segregation (clustered) and integration (global synchronization) dynamics.

To conclude, chimera states have also been reported in complex networks (for an updated compendium of references the reader is addressed to Ref. [232]). In Ref. [287], the Authors considered two globally coupled populations of identical Kuramoto oscillators with inter-population links randomly switched on or off, resulting in the appearance of stable, breathing, and alternating chimera states. Ref. [288] investigated non-locally coupled identical phase oscillators in SF and ER networks, where a coupling function dependence on the shortest length d_{ij} between oscillators is introduced ($A_{ij} = \lambda e^{-\kappa d_{ij}}$, with λ being the global coupling strength and κ the strength of the non-local coupling). With these settings, chimera states spontaneously emerge out of arbitrary initial conditions for both types of complex networks and oscillators. These states tend to include nodes with high degrees on SF networks, while there is no preference for degree for oscillators on ER ones (see Fig. 23).

3.4.2. Non-identical systems

In real systems, populations of interacting oscillators are neither homogeneous nor regular in frequencies, and we already saw that a widely studied paradigm is the Kuramoto model, described in Section 3.3.4. The first attempt to investigate the role of non-regular topologies was performed in Ref. [289]. It was shown there that introducing some variability in a lattice (by randomly rewiring a few local interactions to long-range connections) leads to a more rapid and robust phase locking than nearest-neighbor couplings, or locally dense connection patterns. When the Kuramoto model is considered on top of a complex network of interactions, Eq. (13) translates to

$$\dot{\theta}_i = \omega_i + \lambda \sum_{j=1}^N A_{ij} \sin(\theta_j - \theta_i), \quad (24)$$

where λ is again a global coupling parameter, and A_{ij} is a matrix defining the oscillators' interactions (a complete graph is recovered by setting $A_{ij} = 1/N$ for all pairs).

There are two important issues to be discussed about the extension of Kuramoto model to complex networks: the re-scaling factor in the coupling term of the dynamics, and the issue of properly defining an order parameter. In Refs. [247,290] the implications of different normalizations used in the literature are investigated. One of the most used prescriptions consists in taking $A_{ij} = A_{ij}/N$, which however has the problem that (in the thermodynamic limit) the coupling term vanishes for all nodes except for those whose degree scales as N . Another prescription, used to solve the *paradox of heterogeneity* (discussed in the previous Subsection), consists in taking $A_{ij} = A_{ij}/k_i$. In particular, Ref. [290] considers $A_{ij} = A_{ij}/k_{\max}$, being k_{\max} the largest degree in the network, this way decoupling the heterogeneity of the network and the interaction dynamics and, at the same time, having a non diverging limit for $N \rightarrow \infty$. As for the definition of the order parameter, it is worth noticing that the asymptotic magnitude r defined in Eq. (14) allows for a complete graph the measurement of the average performance of synchronization. Tailoring r for non-complete graphs is the object of many studies, as e.g. those realized by Ichinomiya [291], Restrepo et al. [292], or by Sonnenschein and Schimansky-Geier [293] in the context of complex networks of noisy oscillators.

Early works investigating synchronization in Eq. (24) are those by Watts and Strogatz [245] and Hong et al. [294] for SW networks, and by Ichinomiya [291] and Moreno and Pacheco [295] for SF networks. In particular, Ref. [294] showed (by means of extensive calculations) that any finite probability p of rewiring in the Watts–Strogatz model [245] gives rise to a

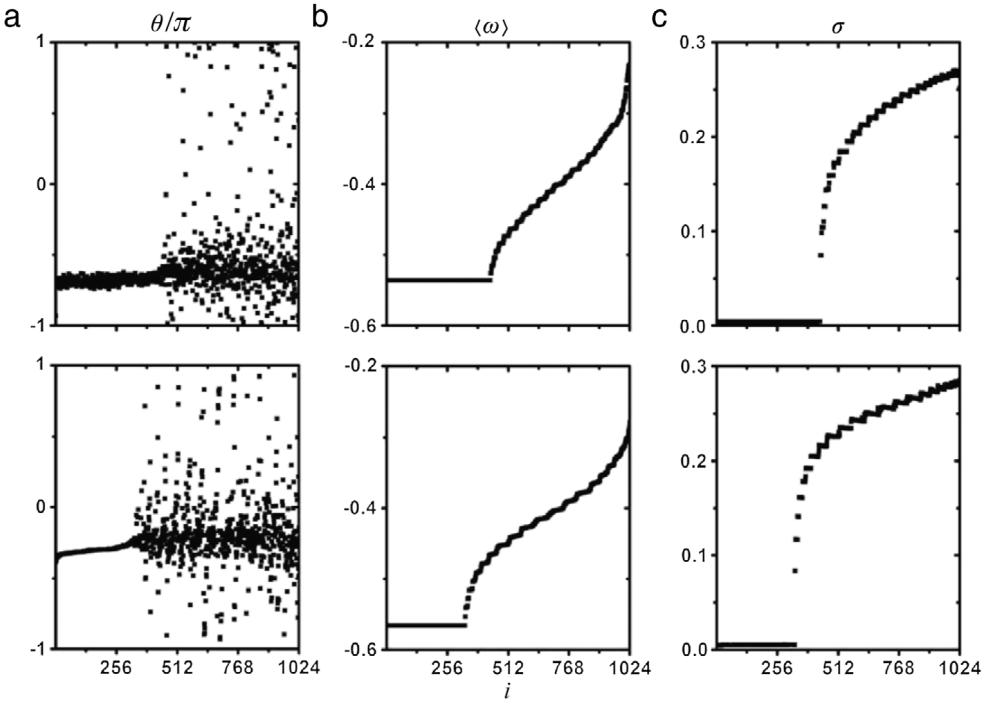


Fig. 23. Chimera states in complex networks. (Column a) Snapshots of the oscillator phases θ_i , (column b) Corresponding time average frequency, and (column c) frequency variance. Top (bottom) panels correspond to ER (SF) networks. The size is $N = 1024$, and the mean degree is $\langle k \rangle = 4$. Other parameters are $\lambda = 1$ and $\kappa = 0.1$.

Source: Reprinted with permission from Ref. [288].

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finite critical coupling λ_c , and that the scaling is indeed compatible with the mean-field of the globally connected graph. Moreover, λ_c diverges as $p \rightarrow 0$, and the time needed to achieve the coherent state decreases with p . In the presence of large degree fluctuations, Ichinomiya proposed to rescale r as follows

$$r(t) = \left| \frac{1}{\sum_l k_l} \sum_j k_j e^{i\theta_j(t)} \right|, \quad (25)$$

and found that the critical coupling for uncorrelated random networks with arbitrary degree distribution is

$$\lambda_c = \lambda_c^0 \frac{\langle k \rangle}{\langle k^2 \rangle}, \quad (26)$$

where λ_c^0 is the critical point for Eq. (22). When the degree fluctuations are bounded, both $\langle k^2 \rangle$ and λ_c are finite in the thermodynamic limit. However, for power-law degree distributions $k^{-\gamma}$, the critical coupling vanishes when $2 < \gamma < 3$. A refinement of the latter result has been provided by Restrepo et al. [292], who generalized the mean-field approach, and showed that the coupling strength at which the transition takes place is determined by the largest eigenvalue of the adjacency matrix ($\lambda_c = \lambda_c^0/\lambda_N$). As for the role of the hubs in the synchronization process on top of a SF network, Ref. [295] showed that the relaxation time τ for the hubs is shorter than that of the less connected nodes, scaling as $\tau \sim k^{-1}$. Hence, the more connected a node is, the more stable it is.

The mean-field approach allows also to compute the scaling of r [296]. For finite λ_c , one has

$$r \sim \Delta^\beta, \quad (27)$$

where $\Delta = (\lambda - \lambda_c)/\lambda_c$. For SF networks with $P(k) \sim k^{-\gamma}$, the exponent takes the value $\beta = 1/2$ if $\gamma > 5$, and $\beta = 1/(\gamma - 3)$ for $3 < \gamma < 5$.

Finally, we briefly review clustering and modular synchronization in networks of oscillators [284,292,297–301]. Since the route to complete synchronization is made up of groups of synchronized oscillators that grow and coalesce, structural properties at the meso-scale of the network can be detected by a fine tuning of the coupling strength. In particular, Ref. [302] compared the synchronization patterns in ER and SF networks and showed that even in the incoherent solution, the system self-organizes following different paths: while in SF networks the giant component of the synchronized pairs is gradually increasing with the coupling strength around the largest degree nodes, in ER networks the coalescence of many small clusters leads to a giant component whose size is of order of the system size once the incoherent state destabilizes. Oh et al. [297]

studied the Kuramoto model on two different types of modular networks, finding that the synchronization transition crucially depends on the type of inter-modular connections. McGraw and Menzinger [298] found that the promotion of dynamical clusters oscillating at different frequencies hinders in general synchronization.

In Li et al. [300], the interfaces between synchronized clusters emerging due to frustration are explored. An algorithm is there provided able to detect the frustrated nodes located at the overlapping structures of real-world networks. The algorithm was then used to unveil protein functions in a protein–protein interaction network in Ref. [303]. Further, an easily computable measure was introduced in Ref. [301] which locates the effective crossover between segregation and integration in modular networks.

Relay (or remote) synchronization of pairs of nodes that are not directly connected via a physical link (nor via any sequence of synchronized nodes) has been observed in star like motifs [304], and in arbitrary networks through the action of mismatched units [305–307].

New forms of synchronous solutions have also been found in multi-layer networks [249,308]. In this latter context, several scenarios have been described: intra-layer synchronization, i.e. a state in which the nodes of a layer are synchronized between them without being necessarily synchronized with those of the other layer [309]; inter-layer synchronization, which, on the contrary, refers to a state where all nodes of a layer are synchronous with their replicas in the other layer regardless of whether or not nodes in each layer are synchronized with the other members of the layer [310,311]; breathing synchronization (a regime resulting from the competition between an instantaneous intra-layer and a delayed inter-layer coupling [312]), where, depending on the couplings and delay, two frequency groups of oscillators emerge within the same layer, each one synchronized with its mirror in a breathing mode.

Most of the works dedicated to the study of the interplay between network topology and coupled dynamics have been performed considering dynamical models of pure phase oscillators and diffusive couplings. However, many biological systems interact in a rather episodic, or pulse-like, way and the amplitude dynamics of the oscillators are usually not constant, as it occurs in coupled neuronal networks. Lago-Fernández et al. [313] compared the synchronization patterns exhibited by networks of nonidentical Hodgkin–Huxley neurons coupled through pulses and different connectivity topologies. It is found that regular topologies produce coherent oscillations with long transients and slow response times, while completely random topologies give rise to a fast but not coherent response to external stimuli, and SW configurations sustain both coherent global oscillations with a fast response.

In a posterior work, it was shown that different types of neural excitability (different amplitude dynamics) endow neural assemblies with very different dynamical properties [314]. Recently, Ref. [307] compared the behavior of networks with dynamical units having both amplitude and phase dynamics and that of networks of pure phase oscillators. The comparison highlights that synchronization is enhanced when the dynamics of the units is close to a Hopf bifurcation, and that the emerging behavior crucially depends on the topology and on the node frequency distribution [307,314].

Synchronization phenomena have been used as a useful (and easy to compute) topology probing tool. Several works focused on how synchronization can unveil structural details of the underlying network where the dynamics is developing [315–317]. It turns out that, in the transient of a synchronization process, topological hierarchies are related to different time scales [315]. On the other hand, based on cluster de-synchronization, Ref. [316] was able to fully detect the modular structure of a network by dynamically changing the weights of the graph's links. Furthermore, when considering negative interactions (a fraction of which has been shown to be beneficial in Ref. [318]) synchronization in signed networks can be used to unveil the organization of an ecological system, and to analyze the role of each species [317].

Finally, several works dealt with synchronization in networks in which structural and dynamical features coevolve [319–325]. By introducing different mechanisms for governing the co-evolution (as, for instance, the competition between the reinforcement of those interactions between pairwise synchronized units [319] and the conservation of the total input strength received by each unit [322,323]), the adaptive network evolves to a state characterized by local synchronization patterns together with an underlying network structure featuring SF and/or modular structures.

4. Explosive synchronization

4.1. Earlier studies on discontinuous transitions in the Kuramoto model

We start this Section with a journey on several early studies aimed at showing the possibility of discontinuous transitions in the Kuramoto model of Eq. (13). The main ideas and conclusions of these studies, indeed, have stimulated the recent advancements on explosive synchronization in networked systems, which will then be reviewed in Sections 4.2, 4.3 and 4.5.

4.1.1. The importance of the frequency distribution

In 1984, Kuramoto proved that there are always two branches of solutions in the thermodynamic limit of Eq. (13) [23]. The first one is the trivial solution $r = 0$ (corresponding to incoherence), and the second one bifurcates continuously at $\lambda = \lambda_c$, with $\lambda_c := \frac{2}{\pi g(\omega_0)}$, being $g(\omega)$ the distribution of natural frequencies, characterized by $g'(\omega_0) = 0$. Kuramoto also found the first evidence of an abrupt synchronization [23]: close to λ_c , there is a small-bifurcating solution [238]

$$r = \sqrt{-\frac{16(\lambda - \lambda_c)}{\pi \lambda_c^4 g''(\omega_0)}},$$

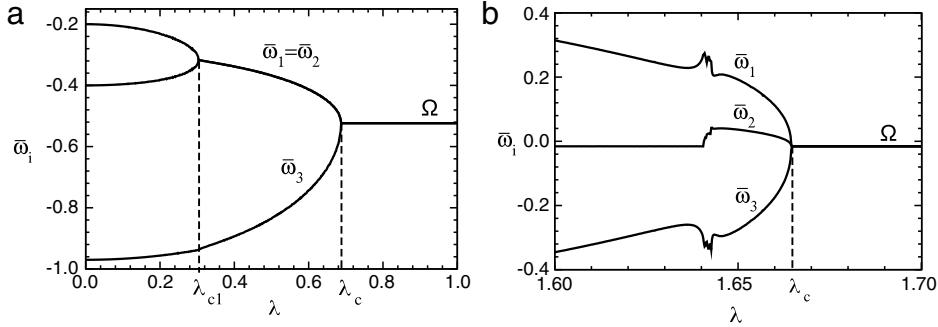


Fig. 24. Frequency-splitting bifurcation diagrams for the Kuramoto model with 3 oscillators. (a) When the natural frequencies are $\{-0.97, -0.4, -0.2\}$, and (b) when they are almost perfectly evenly distributed ($\{-1.0, 0.0, 0.953\}$).
Source: Adapted with permission from Ref. [326]. Courtesy of Y. Maistrenko.

which is supercritical for $g''(\omega_0) < 0$ (i.e., when there is maximum at ω_0) and subcritical if $g''(\omega_0) > 0$ (i.e., when ω_0 is a local minimum).

The behavior of Eq. (13) with a discrete frequency distribution was then studied using a Lyapunov function formalism [327]. Precisely, the Lyapunov function $\mathcal{H} := -\frac{\lambda}{2N} \sum_{i,j} \cos(\theta_i - \theta_j) - \sum_i (\omega_i - \Omega) \theta_i$ was proposed (with Ω being the average frequency) and interpreted as a Hamiltonian of a XY ferromagnet of strength λ in a random field. The approach allowed to prove that the order parameter is always $r \geq 0.5$, and therefore a continuous transition from the phase locked to the incoherent state is impossible. Later on, Ref. [328] proved that the appearance of a macroscopic cluster of mutually phase-locked oscillators takes place continuously when the bifurcation is supercritical, and discontinuously (with hysteresis and bistability) otherwise.

Ref. [326] studied how the Kuramoto model de-synchronizes. A transition called frequency-splitting bifurcation is observed: when the coupling strength decreases below λ_c only one of the phases splits off from the others, but if the natural frequencies are evenly distributed in an interval, the synchronized state splits into several clusters (see Fig. 24, where the effect is shown for three oscillators). The conclusion is that, although a discrete frequency distribution produces a discontinuous transition in the thermodynamic limit, the transition at any finite size depends on how the natural frequencies are chosen.

Following the same concepts and ideas, Pazó focused on understanding the finite-size behavior of the Kuramoto model [14]. Ref. [14] proved (analytically and numerically) that several frequency distributions (with support on a finite interval) lead to a discontinuous phase transition. First, the Kuramoto model is analyzed in the limit $N \rightarrow \infty$ for a uniform frequency distribution $U(-\gamma, \gamma)$, and the critical coupling $\lambda_c := 4\gamma/\pi$ is deduced for which the order parameter undergoes an abrupt transition from zero to $r_c := \pi/4$ (associated with the emergence of meta-stable states [329]). Then, the finite size effects are studied with natural frequencies that are evenly distributed in the interval $(-\gamma, \gamma)$. With such a choice of frequencies, Ref. [14] derives a self-consistent equation for a finite population of oscillators that accurately reproduces the numerical results.

Ref. [330] further studied the case of a unimodal frequency distribution $g(\omega)$ (depicted in Fig. 25(a)) with a plateau in the interval $(-\omega_0, \omega_0)$ and symmetric tails behaving as $g(\omega) := g(\omega_0) - C(\omega - \omega_0)^m H(\omega - \omega_0)$, where C is a constant, $m > 0$ is a parameter and H is the unit step Heaviside function. Fig. 25(b) shows that a critical coupling value λ_c exists where the system undergoes a discontinuous transition. The order parameter r jumps from 0 to $r_c := \frac{\pi}{2}\omega_0 g(0)$. The transition is discontinuous for any value of m , as long as the plateau is a finite flat region (i.e., $\omega_0 > 0$), and only when $\omega_0 = 0$ the transition becomes continuous.

The latter result was extended to asymmetric frequency distributions in 2008 [331]. Going beyond earlier studies on the impact of asymmetries in the frequency distributions [332,333], Ref. [331] found that a necessary condition for a transition to be discontinuous is to have a flat interval in the region of the frequency distribution where the synchronization seed is generated.

Originally, when Kuramoto considered the subcritical case, $g(\omega)$ was assumed to be a symmetric function around the local minimum at ω_0 , featuring two local maxima at ω_{\pm} . The existence of the two local maxima implies that the incoherence solution, $r = 0$, cannot persist up to λ_c even as a meta-stable state. Therefore, the lower critical value for the onset of nucleation around ω_{\pm} is $\lambda'_c := \frac{2}{\pi g(\omega_{\pm})} < \lambda_c$. As λ is increased, the two clusters grow independently of each other, like two giant oscillators, until for some large coupling value they eventually synchronize. An exact description of the thermodynamic limit of Eq. (13) with a bimodal frequency distribution was found in 2009 [243] for $g(\omega) := \frac{\Delta}{2\pi} \left(\frac{1}{(\omega - \omega_0)^2 + \Delta^2} + \frac{1}{(\omega + \omega_0)^2 + \Delta^2} \right)$, being $\pm\omega_0$ the center frequencies of two Lorentzians, and Δ the half width at half maximum, which must verify $\omega_0/\Delta > 1/\sqrt{3}$. Together with the well-known incoherent and partially synchronized states, a third state (consisting of two symmetric clusters of synchronized oscillators near the distribution's maxima) is found, which later on was termed as the standing wave state [334]. However, the bifurcation sequence suggested by Kuramoto (incoherence \rightarrow

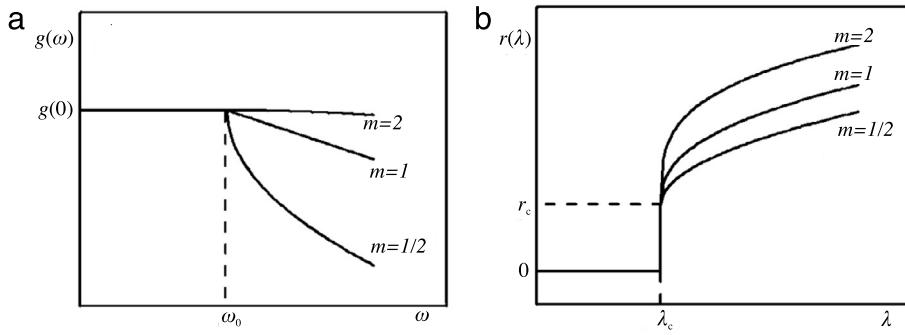


Fig. 25. (a) Frequency distributions considered in Ref. [330], for three values of the parameter m , and (b) the order parameter $r(\lambda)$ for each distribution in (a).

Source: Adapted with permission from Ref. [330]. Courtesy of L. Basnarkov.

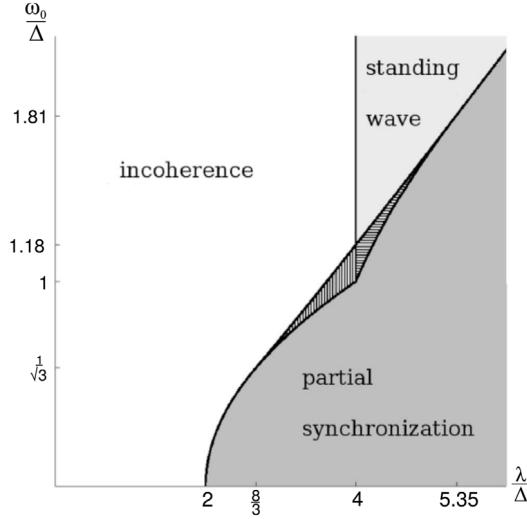


Fig. 26. Stability diagram of the Kuramoto model with a bimodal frequency distribution showing the long-term behavior in each region of parameter space. White, incoherence; dark gray, partial synchronization; light gray, standing wave; vertical lines, coexistence of incoherent and partially synchronized states; horizontal lines, coexistence of partial synchronization and standing waves.

Source: Adapted with permission from Ref. [243]. Courtesy of E.A. Martens.

standing wave → partial synchronization, as sketched in the upper part of Fig. 26) is not the unique possibility. Actually, such a sequence occurs when the two maxima in the frequency distribution are sufficiently far apart. Furthermore, the conjectured formula for λ'_c is incorrect in general, being only valid asymptotically in the limit of widely separated peaks. When the frequency distribution is barely bimodal, instead, the bifurcation sequence does not include the standing wave state, and there is hysteresis in the transition. In the intermediate regime, where the peaks of $g(\omega)$ are neither too far apart nor too close, the forward bifurcation occurs via the sequence predicted by Kuramoto, while the backward bifurcation skips the standing wave state and results directly from partial synchronization to incoherence.

These results were complemented by Pazó & Montbrió [335] with a study in which the minimum and the distance between maxima can be independently modified. In particular, a normalized distribution is considered of the form $g(\omega) := \frac{1}{\pi(1-\xi)} \left(\frac{1}{\omega^2+1} - \xi \frac{\gamma}{\omega^2+\gamma^2} \right)$, and similar bifurcation sequences are obtained. The case $\xi = \gamma$ is of special interest (as there the value at the minimum is always $g(0) = 0$, independently of how close the two maxima are), and results in a sequence incoherence → standing wave → partial synchronization without hysteresis (even for γ extremely close to zero).

4.1.2. Effects of noise

The study on how noise affects the Kuramoto model started with Ref. [336], where an independent white noise stochastic source ξ_i was added to Eq. (13), with $\langle \xi_i(t) \rangle = 0$ and $\langle \xi_i(t)\xi_j(t') \rangle = 2D \delta(t-t') \delta_{ij}$ (D being the noise intensity and angular brackets denoting ensemble average over independent noise realizations). Assuming a Gaussian frequency distribution, Ref. [336] found that noise makes the critical coupling strength larger but does not change the continuous nature of the transition.

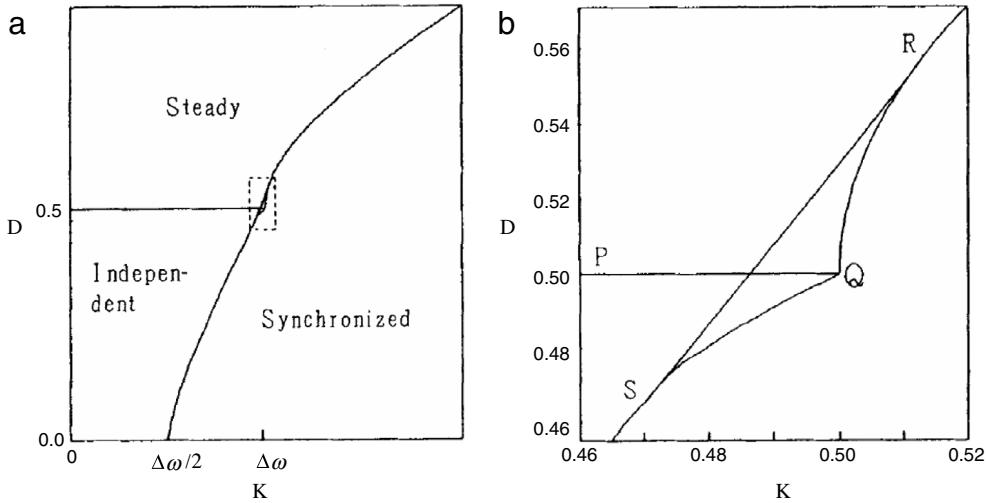


Fig. 27. (a) Bifurcation diagram of the Kuramoto model with Gaussian white noise of intensity D and a discrete frequency distribution consisting of two equiprobable frequencies such that $\Delta\omega = \omega_2 - \omega_1 = 0.5$. (b) is a magnification of the square enclosed with broken lines in (a). The boundaries of steady (i.e., asynchronous), synchronized and independent regimes are represented by the curves linking PQR, RS and PQS, respectively. Therefore, the system exhibits bistability in the region enclosed by QRS. Notice that here K stands for the coupling strength λ .
Source: Reprinted from Ref. [339] with permission of PTP.

Despite seeming an obvious fact, it is actually hard to prove that the solution $r = 0$ is stable for $\lambda < \lambda_c$ and unstable for $\lambda > \lambda_c$. Ref. [337] revealed that the incoherent solution is unstable for $\lambda > \lambda_c$, but neutrally stable for $\lambda < \lambda_c$. Surprisingly, when the problem is studied for small noise, and then the limit is considered of noise intensities going to zero, the results are completely different: the relaxation of $r(t)$ to its steady state follows an exponential law for $\lambda > \lambda_c$, whereas for $\lambda < \lambda_c$ the decay rate depends on the frequency distribution [338]. If $g(\omega)$ has a compact support, $r(t)$ tends to zero slower than an exponential, but if $g(\omega)$ is supported on the reals, $r(t)$ is known only in some cases.

Back in 1991, Okuda & Kuramoto [339] considered the case of Gaussian white noise and a discrete frequency distribution consisting of two equiprobable frequencies $\omega_2 > \omega_1$. A critical value D_c of the noise intensity is found, below which the system synchronizes either partially (only oscillators with the same frequency synchronize) or globally. Consequently, the bifurcation diagram, shown in Fig. 27(a), has three regimes: asynchronous (referred to as steady), partially synchronized (referred to as independent), and globally synchronized, which overlap in the region QRS depicted in Fig. 27(b), indicating that the transition to synchrony occurs through a region with bistability.

Therefore, with or without noise, the Kuramoto model presents a rather complex behavior as soon as the frequency distribution has two peaks. The case of two peaks equidistant to zero $\omega_2 - \omega_1 := 2\omega_0$ was studied in Ref. [340]. If the maxima are close, $\omega_0 < D/\sqrt{2}$, the transition to the synchronous state is continuous. When $\omega_0 \in (D/\sqrt{2}, D)$, the transition is discontinuous with hysteresis. Finally, when the peaks are separated enough, $\omega_0 > D$, the transition is also discontinuous with hysteresis, but now the stable synchronized state is characterized by having a time-periodic order-parameter. The case of a more general frequency distribution $g_\varepsilon(\omega) := \frac{\varepsilon}{2\pi} \left(\frac{1}{(\omega+\omega_0)^2+\varepsilon^2} + \frac{1}{(\omega-\omega_0)^2+\varepsilon^2} \right)$ (which consists in two Lorentzians centered at $\pm\omega_0$ with the same width parameter ε) was studied in Ref. [334]. A critical value for the separation between the maxima, $\omega_d := (\varepsilon + D)\sqrt{\frac{\varepsilon+2D}{3\varepsilon+4D}}$ is found, that determines the behavior of the system. For $\omega_0 < \omega_d$, the bifurcation is supercritical to a stable synchronized state. For $\omega_d < \omega_0 < \varepsilon + D$, there is a subcritical bifurcation to a synchronized state with no prior onset of synchronization in the peaks of the distribution. And for $\omega_0 > \varepsilon + D$, there is a transition to a state consisting of two symmetric clusters of synchronized oscillators, called standing waves. As the peaks move far apart (i.e., $\omega_0 \rightarrow \infty$), the oscillators in the peaks indeed synchronize, but as $\omega_0 \rightarrow \varepsilon + D$, they synchronize at frequencies that are substantially shifted away from the peaks of the native distribution. Actually, Fig. 28, where these results are schematically shown, is remarkably similar to Fig. 26, but the effect of the noise added is to increase the width of each Lorentzian in D , having an effective width $\varepsilon + D$.

The asymmetric case was studied in Ref. [341] which considered the bimodal distribution $g(\omega) := \alpha \delta(\omega - \omega_0) + (1 - \alpha) \delta(\omega + \omega_0)$. Here, $\alpha \in (0, 1)$ entails essentially different features with respect to the symmetric case ($\alpha = 1/2$). By comparing Fig. 29 (which shows the stability boundaries for $\alpha = 0.49$ and $\alpha = 0.3$) with Fig. 28, one can appreciate how the asymmetry of the frequency distribution changes the stability boundaries: both the region of incoherence and standing waves become smaller, whereas the partially synchronized area grows. Finally, Ref. [342] considered a discrete version of the Kuramoto model to explore the transition to synchronization in a population of weakly coupled oscillators. When the coupling strength exceeds some noise-dependent threshold, the population develops spike-like solutions with noisy boundary layers, and its behavior is similar to that of the continuum limit. However, a much more complicated

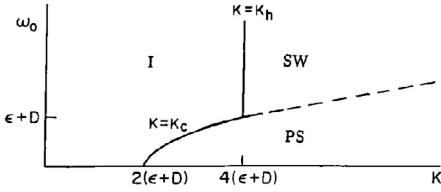


Fig. 28. States of the Kuramoto model with Gaussian white noise of intensity D and a bimodal frequency distribution consisting of two Lorentzians centered at $\pm\omega_0$ with the same width parameter ε : the incoherent state (I), the partially synchronized state (PS), and the standing waves (SW). The boundary between the standing waves and the partially synchronized state is shown schematically as a dashed line because the precise nature and location of this boundary have not been determined. Notice that here K stands for the coupling strength λ .
Source: Reprinted from Ref. [334] with permission of Springer.

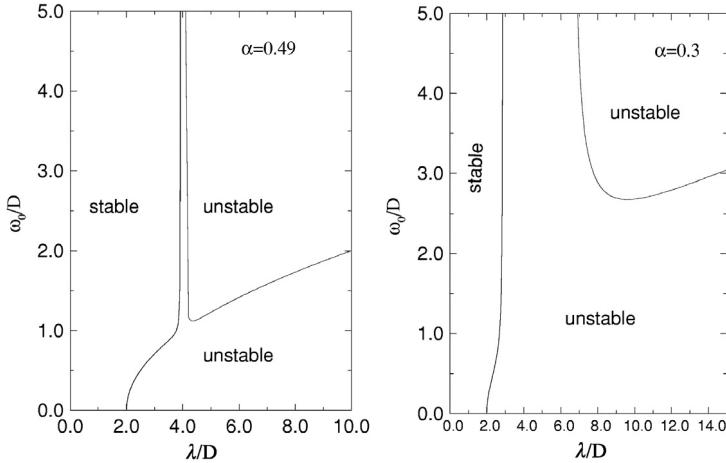


Fig. 29. Stability boundaries for the incoherent solution for the asymmetric bimodal frequency distribution when $\alpha = 0.49$ and $\alpha = 0.3$.
Source: Adapted with permission from Ref. [341]. Courtesy of R. Spigler.

behavior occurs for sub-threshold coupling strengths, where the finite population displays complicated temporal dynamics (ephemeral coherent structures, switching patterns).

4.1.3. Generalizations of the coupling pattern

Another streamline of research analyzed the effects of changing the original (uniform, all-to-all) connectivity of Eq. (13). Ref. [343], for instance, generalized the Kuramoto model to account for intersegmental coordination of neural networks, responsible for generating locomotion in the isolated spinal cord of lamprey. Ref. [343] considered that each oscillator only has two nearest neighbors, and assumed that all upward and downward coupling strengths are equal. Not only most of the experimental observations were reproduced, but also stable phase locked motions (corresponding to traveling waves) as well as drifting motions were found.

A more general approach was followed in Ref. [344], which assumed a Gaussian $g(\omega)$ and an underlying topology of a d -dimensional hypercubic lattice. Namely, $\lambda_{ij} = \lambda$ is taken only if i and j are nearest neighbors in the lattice. The d -lattices self-organize in clusters of entrained oscillators, for which the lattice dimension determines if some of them develop into a macroscopic size.

Ref. [345] revisited the problem with the aim to give more accurate predictions by means of a real-space renormalization-group analysis. Ref. [345] considered the case of a generic coupling pattern $h_{ij}(\theta, \lambda)$ satisfying $h_{ij}(-\theta, \lambda) = -h_{ij}(\theta, \lambda)$ (therefore the Kuramoto model is encompassed as the case $h_{ij}(\theta, \lambda) = \lambda \sin \theta$). Ref. [345] first showed that the global behavior of the lattice strongly depends on the frequency distribution. If the variance of $g(\omega)$ is not defined, its asymptotic behavior can be written as $g(\omega) = \Theta(|\omega|^{-a-1})$, with $a \in (0, 2]$, while if a frequency distribution has finite variance, it can be studied by taking $a = 2$. Assuming that the whole d -dimensional lattice is divided into a set of hypercubes with an equal linear scale L , Ref. [345] demonstrated that synchronization only arises in a d -lattice when $d > a/(a-1)$. In particular, this implies that synchronization is impossible in any d -lattice if $a \in (0, 1]$ (e.g., for a Lorentzian), while it is possible for any frequency distribution with finite variance (e.g., for a Gaussian) if the lattice has $d > 2$.

Later on, however, Ref. [346] found that global synchronization is actually impossible for all d -lattice. Precisely, given a Gaussian frequency distribution, and N oscillators in a d -lattice coupled with strength λ , the probability $P(N, \lambda, d)$ of phase-locking is found to verify $\lim_{N \rightarrow \infty} P(N, \lambda, d) = 0$, for all finite λ and d . Furthermore, dropping the assumption of a Gaussian distribution and assuming only that $g(\omega)$ has finite mean and finite non-zero variance, the former result remains valid. Therefore, there is no critical dimension for phase-locking, because the probability tends to zero exponentially fast for any d -lattice.

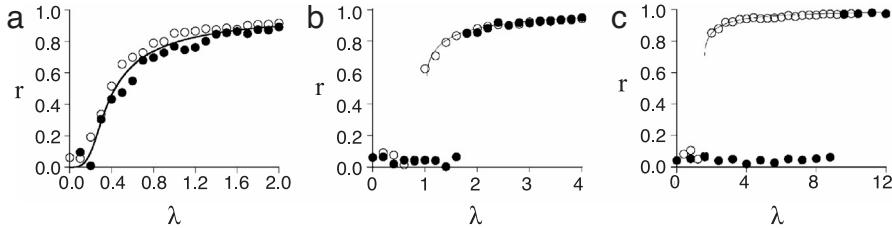


Fig. 30. Numerical and analytical results of the Kuramoto model proposed in Ref. [131] for various feedback strengths: (a) $z = 0.7$, (b) $z = 2$, (c) $z = 3$. The solid lines represent the analytic prediction, the filled (open) circles show the forward (backward) transition numerically computed. The disorder is set to $\gamma = 0.05$.

Source: Adapted with permission from Ref. [131]. Courtesy of G. Filatrella.

Ref. [347] proposed one of the first models in which the interactions among the oscillators are random. In order to have an analytically tractable model, it is considered that the interaction λ_{ij} between oscillators i and j (which are placed on a lattice) is $\lambda_{ij} = \frac{\lambda}{z} s_i s_j$, where λ is a control parameter, z is the number of interacting neighbors of each oscillator, and s_i are random parameters taken from a distribution $P(s)$. For the particular case $P_a(s) = a \delta(s - 1) + (1 - a) \delta(s + 1)$ (being $a \in [0, 1]$ a parameter and δ the Dirac function), the order parameter r_d comes out to be proportional to the original Kuramoto parameter r , $r_d = |2a - 1| r$, thus randomness causes a decrease in the order parameter. The symmetric case (i.e., $a = 1/2$) results in a disordered system even for $\lambda > \lambda_c$, which was called a spurious glass-like phase. Ref. [348] proposed another model in which the interaction λ_{ij} is assumed to be an independent random variable with normal probability distribution, $P(\lambda_{ij}) = \mathcal{N}(0, \lambda^2/N)$ (where λ is the control parameter). Numerical evidence is obtained of a new type of ordered phase that is analogous to those observed in a variety of glassy systems.

Ref. [349] studied the case in which the underlying network topology is a tree with branching ratio b , being the oscillators at the leaves, under the hypotheses that the coupling strengths vary with the distance and are distributed over a hierarchy of values. Specifically, it was assumed that intra-cluster interactions are weaker than inter-cluster ones. It is shown that distinct regimes are emerging depending on whether the branching ratio b is smaller or larger than the critical value b_c .

Ref. [350] provided a stability analysis of the Kuramoto model in terms of appropriate Lyapunov functions, recovering all the previous results for fully connected graphs, and obtaining new ones for generic topologies. Finally, Filatrella et al. [131] proposed a modification of the Kuramoto model to account for the effective change in the coupling constant suggested by some experiments on Josephson junctions [351,352], laser arrays, and mechanical systems. Namely, Ref. [131] proposed a variation of the Kuramoto model that accounts for change of the coupling with the number of active oscillators: the natural frequencies are taken from a Lorentzian with zero average and γ width, and the coupling λ_{ij} is taken to be $\lambda_{ij} = \lambda(r) = \lambda r^{z-1}$, where the parameter z is a heuristic measure of the strength of the feedback mechanism. The resulting model is analytically tractable, and predicts that both first and second order phase transitions are possible, depending upon the value of the parameter z that tunes the coupling among the oscillators. For $z \leq 1$ the evolution from the incoherent value $r = 0$ to the partially coherent state is continuous (see Fig. 30(a)). When $z > 1$, the critical value where the transition to synchrony occurs changes for the forward and backward transitions, yielding hysteresis, whose width is larger for the larger z (as it is evident by comparing the case $z = 2$ in Fig. 30(b) to the case $z = 3$ in Fig. 30(c)).

4.2. Explosive synchronization in complex networks

Unfortunately, the scientific community was initially somehow apathetic to the many results summarized in the previous Section, and such a situation (of almost indifference) persisted lamentably for longtime, up to when the scientific community started to be interested in the study of the transition to synchrony in complex networks.

4.2.1. Explosive synchronization in networks of frequency-degree correlated oscillators

In Ref. [15], an unweighted and undirected network of N Kuramoto oscillators [187] is considered, whose equations of motion are

$$\dot{\theta}_i = \omega_i + \lambda \sum_{j=1}^N A_{ij} \sin(\theta_j - \theta_i), \quad (28)$$

where ω_i stands for the natural frequency of oscillator i , λ is the coupling strength, and A_{ij} are the elements of the network's adjacency matrix.

Ref. [15] actually introduced a specific (ad-hoc) form of correlation between the natural frequencies and the degrees of the oscillators. Namely, the natural frequency ω_i of node i was imposed to be proportional to its degree k_i ($\omega_i = k_i$). Once such a correlation is artificially imposed, different networks (generated by the algorithm of Ref. [353]) are studied, having the same average connectivity but variable heterogeneity properties (i.e. ranging from ER to Barabási-Albert (BA) graphs, depending on the value of a single parameter $\alpha = [0, 1]$).

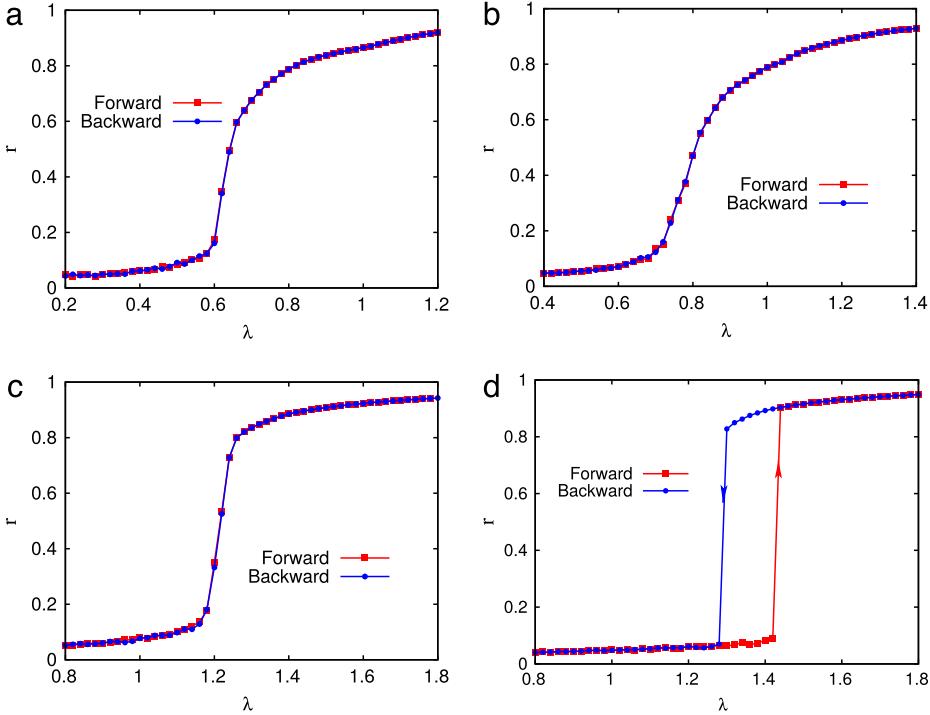


Fig. 31. (Color online). $r(\lambda)$ for various networks constructed with the algorithm in Ref. [353], $N = 10^3$ and $\langle k \rangle = 6$. The α values are: 1 (a, ER), 0.6 (b), 0.2 (c), and 0 (d, BA). The panels report both the forward and backward transitions using steps $\delta\lambda = 0.2$.

Source: Reprinted with permission from Ref. [15].

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The results are summarized in Fig. 31. The synchronization diagrams for several values of α are plotted as a function of the coupling λ , for both forward and backward transitions. The transition is soft and reversible for most values of α , but when $\alpha = 0$ (corresponding to a BA network) an abrupt behavior appears, with the parameter r remaining close to zero up to a point at which it suddenly jumps to $r \sim 1$. As the critical coupling is different for the forward and the backward diagrams, the transition is irreversible and displays a hysteresis loop.

Ref. [15] offers also an analytical study of a specific star configuration, composed of a central node (the hub h) and K peripheral nodes (the leaf nodes), each one labeled by the index i ($i = 1, \dots, K$). By setting the frequencies of the central node ($\omega_h = K\omega$), and that of the leaves ($\omega_i = \omega \forall i$), the equations of motion can be written as

$$\begin{aligned} \dot{\theta}_i &= \omega + \lambda \sin(\theta_h - \theta_i), \\ \dot{\theta}_h &= K\omega + \lambda \sum_{i=1}^K \sin(\theta_i - \theta_h). \end{aligned} \quad (29)$$

Then, using the average phase of the system ($\Psi(t) = \Psi(0) + \Omega t$, where Ω is the average frequency) as a rotating frame, the equation for the hub evolution can be rewritten as

$$\dot{\theta}_h = (K - \Omega) + \lambda(K + 1)r \sin(\theta_h). \quad (30)$$

The condition for stability of the hub locking solution ($\dot{\theta}_h = 0$) allows one to obtain the *backward* critical coupling value as $\lambda_c^b = (K - 1)/(K + 1)$, and the order parameter at the critical point is found to be $r_c = K/(K + 1)$, that is $r_c > 0$ at the critical point, confirming the existence of a discontinuity in the transition.

The same degree-frequency correlated star of Eqs. (29) was studied in Ref. [20], where the origin of the hysteresis loop was rooted in the properties of the basin of attraction of the synchronization manifold. Ref. [20] rewrote Eq. (29) as $\dot{\Phi} = \Omega_K + \lambda H(\Phi)$, where $\Phi = (\theta_1, \dots, \theta_h)$, $\Omega_k = (\omega, \dots, \omega, K\omega)$ and $H(\Phi) = (\sin(\theta_h - \theta_1), \sin(\theta_h - \theta_2), \dots, \sum_{j=1}^K \sin(\theta_j - \theta_h))$. Then, the locking manifold can be defined as $M_a = \{\Phi \in T^{K+1} : \theta_1 = \theta_2 = \dots = \theta_K; \theta_h - \theta_1 = a\}$, where a is a constant. The condition $\dot{\Phi} = 0$ yields the minimal coupling that enables the existence of M_a : $\lambda_c^b = \frac{(K-1)\omega}{K+1}$, which corresponds to the *backward* transition.

To obtain insight on the *forward* transition, Ref. [20] provided numerical evidences that (for $\lambda > \lambda_c^f$) the local attractiveness of M_a becomes global, i.e. any state (even incoherent) is attracted to M_a for coupling strengths above λ_c^f . This fact (together with the theory developed in Ref. [354]) can be used to reformulate the phase locking problem between

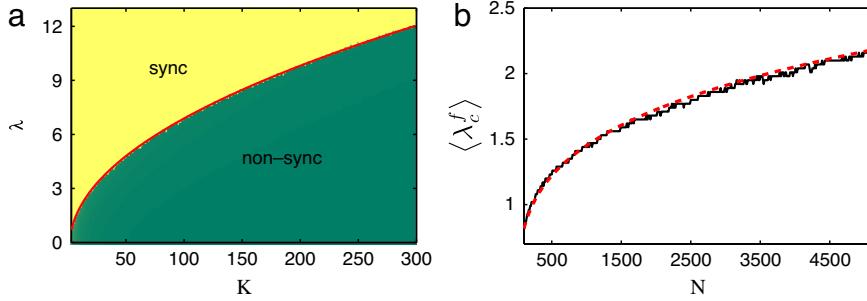


Fig. 32. (a) Order parameter r vs. the space of parameters (K, λ) for the forward transition. (b) The red thick line indicates the solution of Eq. (31), for $1/B = 0.6989$. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)
Source: Adapted with permission from Ref. [20]. Courtesy of J. Kurths.

the hub and leaves as a perturbation of an identical synchronization problem. In such a latter framework, the condition for coherence gives

$$\lambda_c^f \simeq \left(\frac{K-1}{\sqrt{K}} \frac{1}{B} \right) \omega, \quad (31)$$

where B is a constant which depends on the chosen initial conditions. In Fig. 32(a) it is shown how numerical simulations are in excellent agreement with the predictions of Eq. (31).

The results can be extended to the case of SF networks with small mean degrees, that can be considered, indeed, as a collection of stars of different sizes. For $K \gg 1$, $\lambda_c^b \rightarrow \omega$, but $\lambda_c^f \rightarrow \sqrt{K}\omega$. In a SF network with degree distribution $P(k) = k^{-\gamma}$ the expected degree of the largest hub scales as $N^{\frac{1}{\gamma-1}}$. Therefore, the expected λ_c^f scales as

$$\langle \lambda_c^f \rangle \approx N^{\frac{1}{2(\gamma-1)}}. \quad (32)$$

As it can be seen in Fig. 32(b) the approximation gives a very good fit for SF networks with $\gamma = 3$.

The problem of calculating the forward critical point in a frequency-degree correlated star has been very recently reconsidered by several groups, which applied dimensional reduction techniques [355] and the Ott–Antonsen method [241]. The first approach was applied independently by Vlasov et al. [356] and Xu et al. [357]. Taking $\phi_j = \theta_h - \theta_j$, the Eq. (29) can be reduced to

$$\dot{\phi}_j = \delta\omega - \lambda \sum_{i=1}^K \sin \phi_i - \lambda \sin \phi_j \quad j = 1, \dots, K, \quad (33)$$

where $\delta\omega$ is the frequency difference between the hub and leaf node. By defining $z = re^{i\phi} = \frac{1}{K} \sum_{j=1}^K e^{i\phi_j}$, Eq. (33) is rewritten as

$$\dot{\phi}_j = fe^{i\phi_j} + g + \bar{f}e^{-i\phi_j}, \quad (34)$$

where $f = i\frac{\lambda}{2}$, and $g = \delta\omega - \lambda Kr \sin(\Phi)$. The Watanabe–Strogatz approach can be used in systems of identical oscillators driven by a common force (having the general form $\psi_j = f(t) + \text{Im}(F(t)e^{-i\psi_j})$), as it is the present case. The phase dynamics of the K nodes can be constructed by K constants as

$$e^{i\phi_j} = \frac{\beta + e^{i(\psi + \epsilon_j)}}{1 + \bar{\beta}e^{i(\psi + \epsilon_j)}}, \quad (35)$$

where $\beta = \beta(t)$ and $\psi = \psi(t)$ are global variables, and ϵ_j are constants depending on the initial condition of the systems. In the thermodynamic limit ($K \rightarrow \infty$) such a transformation allows to separate the evolution of the $\beta(t)$ and $\psi(t)$ variables, and gives $\beta(t) = z(t)$. Therefore, the equation of the order parameter can be written as [356,357]

$$\dot{z} = -\frac{\lambda}{2}z^2 + i(\delta\omega - \lambda Kr \sin \Phi)z + \frac{\lambda}{2}, \quad (36)$$

which describes the collective dynamics of system (29) in terms of the ensemble order parameter, and can be used even for small values of K in case of random initial distributions. In the phase space of the ensemble order parameter, the synchronous state corresponds to a fixed point (with $r = 1$) and a fixed phase Φ . All the other solutions of Eq. (36) represent various incoherent states. If Eq. (36) is rewritten in its Cartesian coordinates $z = x + iy$, the null-clines $\dot{x} = \dot{y} = 0$ of the system can be studied as a function of the coupling λ , allowing the calculation of the forward critical coupling [356,357]

$$\lambda_c^f = \frac{\delta\omega}{\sqrt{K}} \frac{1}{\sqrt{2 + K^{-1}}}. \quad (37)$$

It can be seen that the approximation (31) is a good estimation of the latter result for large K , where $\lambda_c^f = \frac{\delta\omega}{\sqrt{2K}}$.

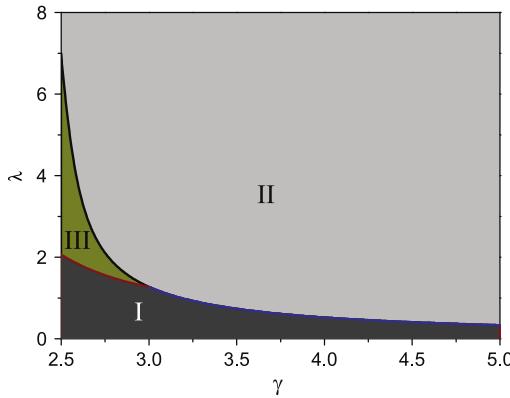


Fig. 33. Phase diagram $\gamma - \lambda$ of the Kuramoto model for frequency-degree correlated SF networks. In region I there is no spontaneous synchronization, and the order parameter vanishes. Synchronization appears in region II, in which the order parameter is larger than zero. Region III is the region of hysteresis with a meta-stable state and one stable state.

Source: Adapted with permission from Ref. [359]. Courtesy of S.N. Dorogovtsev.

A similar dimensional reduction has been provided in Ref. [358], in which an extension of the model is presented where random perturbations affect the frequencies of leaves. The results show that for small perturbations the synchronization phase transition is still discontinuous, while for large perturbations the transition becomes continuous.

In Ref. [359] the case of a star graph with arbitrary natural frequencies is studied. Then, the systems' equations become

$$\begin{aligned} \dot{\theta}_i &= \omega_i + \lambda \sin(\theta_h - \theta_i), \\ \dot{\theta}_h &= \omega_h + \lambda \sum_{i=1}^K \sin(\theta_i - \theta_h). \end{aligned} \quad (38)$$

The transition to synchrony of this system is discontinuous when $\omega_h - \langle \omega_i \rangle > \omega_c$, where ω_c is a certain critical frequency difference.

Beyond a star graph, analytic results for general networks (under the assumption that $g(\omega) = P(k)$) were given by Peron et al. [360]. In the continuous limit, the equation of the system is

$$\dot{\theta} = k + \lambda k \int dk' \int d\theta \frac{k' P(k')}{\langle k \rangle} \rho(\theta'|k) \sin(\theta' - \theta). \quad (39)$$

Ref. [360] shows that (in such a limit) the forward critical coupling is $\lambda_c^f = \frac{2}{\pi \langle k \rangle P(\langle k \rangle)}$.

Notice that this critical coupling differs from the usual value $\lambda_c^0 = \frac{2}{\pi g(0)} \frac{\langle k \rangle}{\langle k^2 \rangle}$ for symmetric $g(\omega)$.

Even though the calculations in Ref. [360] identify the frequency and degree distributions, they do not account for degree-frequency correlations, and therefore fail in predicting if the transition is explosive or continuous. The issue was addressed in Ref. [359], where the annealed network approximation was used to solve Eq. (28) under the assumption $\omega_i = k_i$. An implicit solution for the order parameter

$$\langle k \rangle - \omega = \int_{|k_i - \Omega| < \alpha rk} dk P(k) (k - \Omega) \sqrt{1 - \left(\frac{\lambda rk}{k - \Omega} \right)^2}, \quad (40)$$

$$r = \frac{1}{\langle k \rangle} \int_{|k_i - \Omega| > \alpha rk} dk P(k) k \sqrt{1 - \left(\frac{k - \Omega}{\lambda rk} \right)^2}, \quad (41)$$

is found. Following such a solution, ER networks always undergo second order transitions, whereas the nature of the transition in SF networks depends on the heterogeneity parameter γ , as shown in Fig. 33: a first-order like transition is found for $2 < \gamma < 3$, but a second-order transition occurs for $\gamma > 3$. A hybrid phase transition is found for $\gamma = 3$, combining a jump of the order parameter and critical phenomena at the critical point. This is actually a first indication that not all SF networks are able to sustain explosive synchronization (ES), even when they are fully frequency-degree correlated, a subject that will be reconsidered afterwards [22].

In a number of works, various generalizations of the frequency-degree correlation have been analyzed. In Ref. [361], a more general joint distribution $P(\omega, k) = P(k)[\delta(\omega - \alpha k^\beta) + \delta(\omega + \alpha k^\beta)]$ is considered. The phase-locking behavior of the system is examined, and it is found that the criterion for locking is $\alpha k^{\beta-1} < \lambda r$, being r the order parameter. From this expression one can deduce that the nodes with low degree are the last ones to get phase-locked for sub-linear correlation ($\beta < 1$), whereas super-linear correlation ($\beta > 1$) induces the last drifting nodes to be the hubs. These two qualitatively

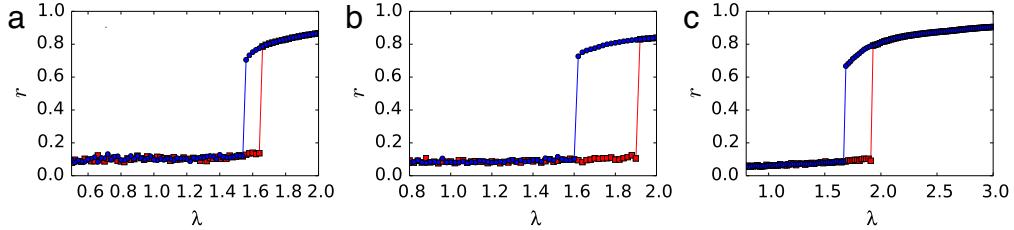


Fig. 34. (Color online). Synchronization diagrams $r(\lambda)$ for networks built with the mechanism proposed in Ref. [353] with $\alpha = 0.2$ (a), 0.1 (b) and 0 (c), where only 10% of the vertices with largest degree show degree–frequency correlations.
Source: Adapted with permission from Ref. [362]. Courtesy of R.S. Pinto.

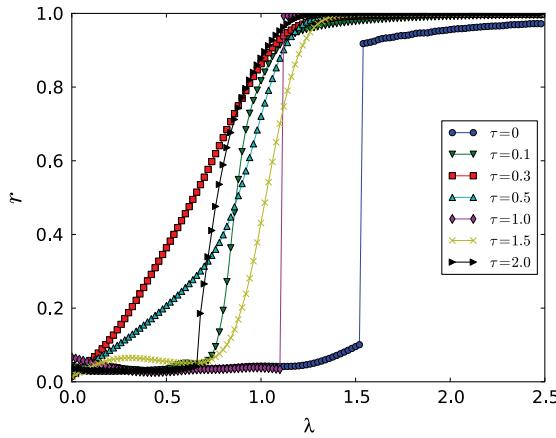


Fig. 35. $r(\lambda)$ for a BA model with $N = 1000$ and $\langle k \rangle = 6$, for several values of the delay τ (see legend for color and symbol code). (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)
Source: Adapted with permission from Ref. [364]. Courtesy of F.A. Rodrigues.

different behaviors are separated by the critical case of linear correlations ($\beta = 1$) for which k disappears, and the oscillators lock simultaneously.

In Ref. [363], instead, the possibility of a partial frequency–degree correlation is considered, and it is shown that a small fraction of correlated nodes is enough to obtain ES in a SF network. The issue of partial correlation has been dealt with also in Ref. [362], where only vertices with degree k larger than a certain threshold k_* are supposed to exhibit the degree–frequency correlation, whereas the other vertices have random natural frequencies with distribution $g(\omega)$. This latter feature can be formally enunciated through a joint probability distribution (for a node of degree k and natural frequency ω) given by $G(\omega, k) = [\delta(\omega - k)P(k) - g(\omega)P(k)]H(k - k_*) + g(\omega)P(k)$, where $\delta(x)$ and $H(x)$ are, respectively, the Dirac delta and the Heaviside step functions, and $P(k)$ is the degree distribution. A mean-field calculation shows that the correlation of the 10% highest degree nodes is enough to promote ES in BA networks. The results are shown in Fig. 34: by imposing a partial degree–frequency correlation not only it is possible to keep ES in the cases where it already happens with full correlation (Fig. 34(b)–(c)) but, unexpectedly, ES emerges also in cases in which full correlation would have prevented it (Fig. 34(a)).

Other works have tested how ES is affected by modifications to the original Kuramoto model. In Ref. [101] is reported that ES persists (and is even enhanced) in modular BA, k – ω correlated networks. In Ref. [364], the model of Ref. [15] is modified to include delay times phase shifts [357], and it is found that an appropriate choice of parameters can enhance or suppress ES. An example of the effect of time delay is shown in Fig. 35.

Ref. [365] generalized the k – ω relationship as $\omega_i = k_i/\beta$, where $\beta = \sum_i^N k_i^\alpha / \sum_i^N k_i$ is a normalization to avoid homogenization of the frequencies when $|\alpha|$ is small, and α is in the range $[-1, 1]$, thus including the possibility that k_i and ω_i are anti-correlated. For positively k – ω correlated networks ($\alpha > 0$) ES remains unmodified for up to $\alpha = 0.1$, but it becomes a second-order transition for any $\alpha < 0$ value. The effect of degree mixing in the model is also considered, and it is pointed out that assortativity has a strong destructive effect on ES. In Ref. [366] a frequency–degree correlated BA network is rewired to introduce assortativity, and it was found that a moderate disassortativity favors the ES, which instead is destroyed for assortative mixing.

A detailed study of the role of the assortativity in ES is provided in Ref. [22]. With the aim of inspecting whether or not ES depends on the chosen network model, Ref. [22] comparatively considers ensembles of networks obtained with the BA algorithm and networks displaying the same SF distributions, but constructed by the so called configuration model (CM, [367]). For both networks, parameters were assumed to be $N = 10^3$ and $\langle k \rangle = 6$, and the oscillators' frequencies were distributed so as to determine a direct correlation with the node degree ($\omega_i = k_i$).

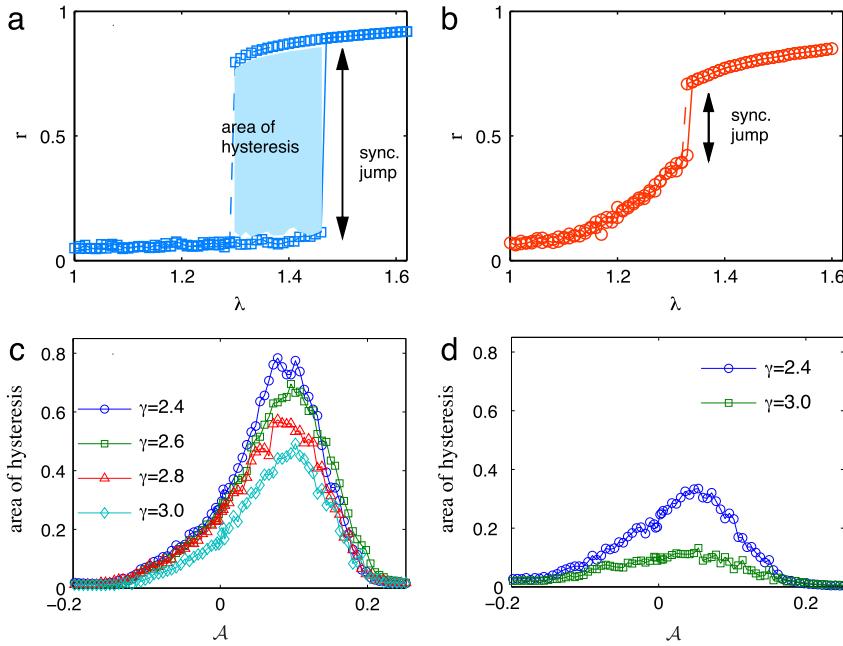


Fig. 36. (Color online). Comparison of ES between two SF networks belonging to two different configuration ensembles: (a), (c) preferential attachment (PA) and (b), (d) configuration (CM) models. (a)–(b) Forward (solid lines) and backward (dashed lines) synchronization curves for a PA (a) and CM (b) SF network, with exactly the same degree distribution. The area of the hysteresis is depicted as a blue shaded area in (a). The synchronization jumps are marked in both cases. (c)–(d) Area of the hysteretic region delimited by the forward and backward synchronization curves vs. the Pearson correlation coefficient \mathcal{A} for PA (c) and CM (d) SF networks with different values (reported in the legend) of the exponent γ for the degree distribution $P(k) \sim k^{-\gamma}$. Each point is an average over 10 different simulations.

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The results reported in Fig. 36 show a dramatic dependence of the ES behavior on the underlying SF network model used, despite displaying the same $P(k)$. The forward and backward continuations of the order parameter r are totally different for BA (Fig. 36(a)) and CM (Fig. 36(b)) networks, and indicate that a crucial condition to obtain irreversibility is having an underlying growing process through which the SF topology is shaped.

One customary way to quantify the amount of degree correlation with a single parameter is using the Pearson correlation coefficient \mathcal{A} , which can be calculated as [368]:

$$\mathcal{A} = \frac{L^{-1} \sum_i j_i k_i - \left[L^{-1} \sum_i \frac{1}{2} (j_i + k_i) \right]^2}{L^{-1} \sum_i \frac{1}{2} (j_i^2 + k_i^2) - \left[L^{-1} \sum_i \frac{1}{2} (j_i + k_i) \right]^2}, \quad (42)$$

where j_i and k_i are the degrees of the nodes at the ends of the i th link, with $i = 1, \dots, L$. Actually, one has that $-1 \leq \mathcal{A} \leq 1$, with positive (negative) values of \mathcal{A} quantifying the level of assortative (disassortative) mixing. The BA model does not exhibit any form of mixing in the thermodynamic limit ($\mathcal{A} \rightarrow 0$ as $(\log^2 N)/N$ as $N \rightarrow \infty$), while a random CM produces highly disassortative networks.

Finally, Ref. [22] studies the impact of increasing/decreasing the assortativity mixing on a network with a given degree sequence $\{k_i\}$ taken from a power law distribution $k^{-\gamma}$. SF networks with given and tunable levels of degree mixing can be generated by the rewiring algorithm of Xulvi-Brunet and Sokolov [369]. Two ensembles of BA and CM networks are then constructed, all of them having the same degree distribution, but different values of the assortativity coefficient \mathcal{A} . Fig. 36(c)–(d) illustrate the effect of degree mixing on ES. Extensive numerical simulations were performed at various values of \mathcal{A} , for slopes γ ranging from 2.4 to 3.0, and for the same mean degree $\langle k \rangle = 6$. The most relevant result is that, regardless of the specific SF network model, the hysteresis of the phase transition is highly enhanced (weakened) for positive (negative) values of \mathcal{A} , and that there is an optimal (positive) value of \mathcal{A} where the irreversibility of the phase transition is maximum. The enhancement is far more pronounced in BA (Fig. 36(c)) than in CM (Fig. 36(d)) networks.

4.2.2. Explosive synchronization in frequency-disassortative networks

While earlier works concentrated on ad-hoc imposed correlations between the node degree and the corresponding oscillator's natural frequency, posterior studies highlighted that a sharp and discontinuous phase transition is by no

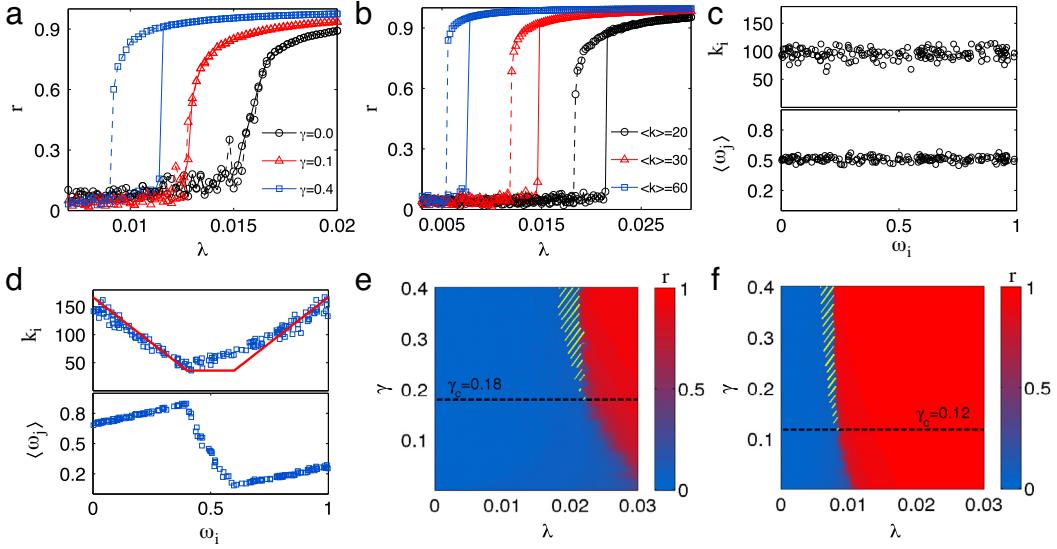


Fig. 37. (a) Phase synchronization level r vs the coupling strength λ , for different values of the frequency gap γ (see, Eq. (43)) at $\langle k \rangle = 40$. (b) Same as in (a), but for different values of the average degree $\langle k \rangle$ at $\gamma = 0.4$. In both panels, the legends report the color and symbol codes for the different plotted curves. In (c) and (d), the degree k_i that each node achieves after the network growth is completed (upper plots) and the average of the natural frequencies $\langle \omega_j \rangle$ of the neighboring nodes ($j \in \mathcal{N}(i)$, bottom plots) are reported vs. the node's natural frequency ω_i , for $\langle k \rangle = 100$ and frequency gaps $\gamma = 0.0$ (c), and $\gamma = 0.4$ (d). The red solid line in (d) is a sketch of the theoretical prediction $f(\omega)$. Panels (e) and (f) show r (color coded according to the color bar) in the parameter space (λ, γ) for (e) $\langle k \rangle = 20$ and (f) $\langle k \rangle = 60$. The horizontal dashed lines mark the separation between the region of the parameter space where a second-order transition occurs (below the line) and that in which the transition is instead of the first order type (above the line). The yellow striped area delimits the hysteresis region. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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means restricted to such rather limited cases, but it constitutes, instead, a generic feature in synchronization of networked oscillators. Precisely, Ref. [17] gave a condition for the transition from unsynchronized to synchronized states to be abrupt, and demonstrates how such a condition is easy to attain in many circumstances, and for a wide class of frequency distributions.

The basic idea, which evokes the Achlioptas process [9] described in Section 2.2, is avoiding that oscillators behave as cores of a clustering process, where neighboring units begin to aggregate smoothly and progressively. As the synchronization seeds are the aggregations of nodes with close frequencies, the procedure introduces then a *frequency disassortativity* in the networks. The practical realization of such a condition is made by imposing explicit constraints in the frequency difference between each node i and the whole set $\mathcal{N}(i)$ of oscillators belonging to its neighborhood [17]:

$$|\omega_i - \omega_j| > \gamma_c, j \in \mathcal{N}(i). \quad (43)$$

In practice, a threshold γ_c for the frequency gap is set, and the network is then grown by means of a *conditional CM* approach: after having initially distributed the oscillator's frequency from a given frequency distribution, pairs of unconnected nodes are randomly selected, and a connection between the nodes is established only if they verify the condition (43). The process is repeated until the network acquires a given, desired, mean degree $\langle k \rangle$, and the resulting adjacency matrix is used to simulate the Kuramoto model (28).

Fig. 37 reports the results obtained by setting $g(\omega)$ as a uniform frequency distribution in the interval $[0, 1]$. Panels (a) and (b) show the phase synchronization index as a function of the coupling strength. In particular, Fig. 37(a) (resp. (b)) illustrates the case of a fixed mean degree $\langle k \rangle = 40$ (resp. of a fixed frequency gap $\gamma = 0.4$), and reports the results for the forward and backward simulations at different values of γ ($\langle k \rangle$). A first important result is the evident first-order character acquired, in all cases, by the transitions for sufficiently high values of γ .

A second relevant result of Ref. [17] is the spontaneous emergence of degree–frequency correlation features associated to the passage from a second- to a first-order like phase transition. While such a correlation was imposed *ad hoc* in Refs. [15,16], here the condition (43) creates for each oscillator i a frequency barrier around ω_i , where links are forbidden. The final degree k_i will be then proportional to the total probability for that oscillator to receive connections from other oscillators in the network, and therefore to $1 - \int_{\omega_i - \gamma}^{\omega_i + \gamma} g(\omega') d\omega'$. This is shown in (c) and (d) of Fig. 37, where the degree k_i that each node achieves after the network construction is completed is reported as a function of its natural frequency ω_i , for $\langle k \rangle = 100$. Precisely, the upper plot of Fig. 37(c) refers to the case $\gamma = 0$ in which no degree–frequency correlation is present. In the upper plot of Fig. 37(d), instead, it is reported the case $\gamma = 0.4$ (a value for which a first-order phase transition

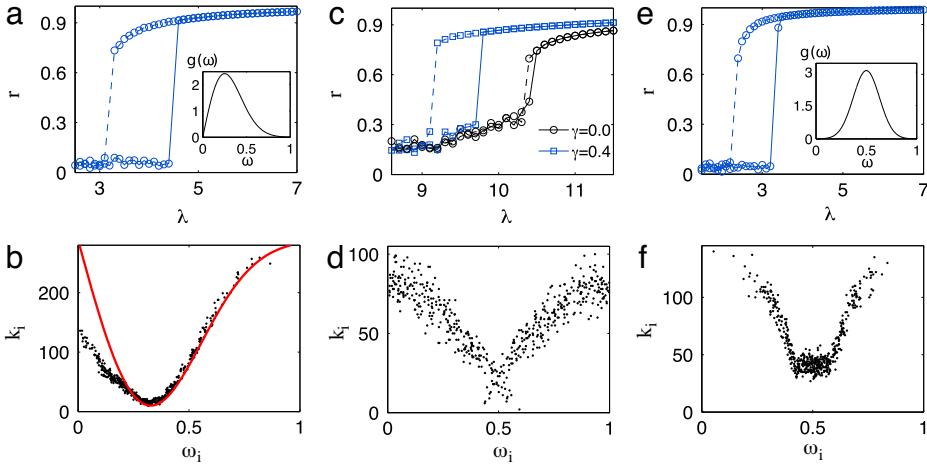


Fig. 38. (Top row) r vs. λ . Synchronization schemes for different frequency distributions or network construction rules. (Bottom row) The corresponding distribution of the final node degree k_i vs. the corresponding oscillator's natural frequency. (a)–(b) Rayleigh distribution for $\gamma = 0.3$. In (b), the red solid line depicts the theoretical prediction $f(\omega)$, obtained with the same method of the red solid line in panel (d) of Fig. 37; (c)–(d) uniform frequency distribution, but network constructed accordingly to a local mean field condition (see text) for $\gamma = 0$, and $\gamma = 0.4$ (see legend for color code). In panel (d) $\gamma = 0.4$; (e)–(f) Gaussian distribution with $Z = 0.7$. The insets in panels A and E report the corresponding distributions $g(\omega)$. In all cases, $\langle k \rangle = 60$. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)
Source: Adapted from Ref. [17].

occurs) and the (conveniently normalized) function $f(\omega) = 1 - \int_{\omega-\gamma}^{\omega+\gamma} g(\omega') d\omega'$, with $g(\omega) = 1$ for $\omega \in [0, 1]$, and $g(\omega) = 0$ elsewhere, which gives evidence of the emergence of a very pronounced V-shape relationship between the frequency and the degree of the network's nodes. The latter fact marks a noticeable difference between the approach of Ref. [17] and those described in the previous Section. Here, indeed, a pattern of frequency-degree correlation emerges spontaneously, and is not artificially imposed. Moreover, the emerging pattern (i) is clearly non-linear; (ii) it depends explicitly on the frequency distribution $g(\omega)$; (iii) it is in general V-shaped.

Finally, Fig. 37(e) and (f) report r in the $\lambda-\gamma$ space, for $\langle k \rangle = 20$ and $\langle k \rangle = 60$, and show that the rise of a first-order like phase transition is, indeed, a generic feature in the parameter space. The horizontal dashed lines in panels (e) and (f) mark the values of γ_c , separating the two regions where a second-order transition (below the line) and a first-order transition (above the line) occur. The fulfillment of Eq. (43) leads to an explosive transition for a very wide class of distributions of the oscillators' natural frequencies, as shown in Fig. 38(a)–(b) for a Rayleigh distribution.

Furthermore, Ref. [17] discusses also several ways of even softening the condition of Eq. (43). For instance, a frequency gap in the network growth can be introduced as $|\omega_i - \langle \omega_j \rangle| > \gamma_c$, where $\langle \cdot \rangle$ indicates the average value over the ensemble $\mathcal{N}(i)$. The results of this (local mean field) gap in a homogeneous frequency distribution are shown in Fig. 38(c)–(d). It is worth noticing that a strict application of the gap condition for non uniform frequency distributions implies that oscillators at different frequencies would in general have a different number of available neighbors in the network. That is why a natural extension is to consider a frequency-dependent gap $\gamma(\omega)$ defined by $\int_{\omega-\gamma}^{\omega+\gamma} g(\omega') d\omega' = Z$. The gap condition for the construction of the network is now to fix the value of Z , and accept the pairing of nodes when $|\omega_i - \omega_j| > \frac{1}{2}[\gamma(\omega_i) + \gamma(\omega_j)]$. The generic case is again that of an explosive transition, with pronounced frequency-degree correlation features, as long as $g(\omega)$ is symmetrical. Panels (e) and (f) of Fig. 38 report the case of a Gaussian distribution limited to the frequency range $[0, 1]$, centered at $\omega = 0.5$, and given by $g(\omega) = \frac{1}{\sigma\sqrt{2\pi}} e^{-\frac{(\omega-0.5)^2}{2\sigma^2}}$, with $\sigma = 0.13$.

Ref. [370] also focused on the idea that ES phenomenon is rooted in preventing oscillators from behaving as the cores of clustering processes. By the use of a constructive gap scheme similar to that of Ref. [17], it is found that frequency disassortativity is accompanied by a strong increase of degree disassortativity. The frequency-equivalent of the Pearson coefficient \mathcal{A} in Eq. (42) is

$$\mathcal{A}_\omega = \frac{L^{-1} \sum_i p_i q_i - \left[L^{-1} \sum_i \frac{1}{2} (p_i + q_i) \right]^2}{L^{-1} \sum_i \frac{1}{2} (p_i^2 + q_i^2) - \left[L^{-1} \sum_i \frac{1}{2} (p_i + q_i) \right]^2}, \quad (44)$$

where q_i, p_i are now the frequencies of the nodes at the ends of link i . From Fig. 39, it can be easily seen that, as the gap parameter γ increases, both degree- and frequency-assortativity coefficients \mathcal{A} and \mathcal{A}_ω stay close to zero up to the critical value γ_c , and then become increasingly negative.

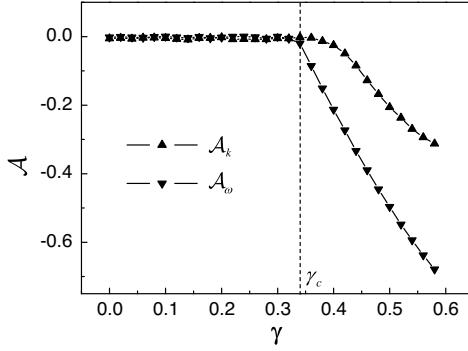


Fig. 39. The degree- and frequency-assortativity coefficients A and A_ω vs. the continuous frequency gap γ .
Source: Adapted with permission from Ref. [370]. Courtesy of D. Shi.

The same idea of a gap condition is used in Ref. [371] in a completely different context, i.e. an ensemble of coupled maps:

$$X_i^{n+1} = f(X_i^n) + \frac{d}{k_i} \sum_{j=1}^N a_{ij} |f(X_i^n) - f(X_j^n)| \quad (45)$$

where $f(X_i^n) = \mu_i X_i^n (1 - X_i^n)$ is the logistic map, and the parameter $\mu_i \in [3, 4]$ is the dimensionless grow factor of element i . In Ref. [371], the connectivity on the ensemble is generated by imposing the gap restriction $|\mu_i - \mu_j| < \gamma$, and ES and hysteresis are observed as a consequence. This latter fact gives a clue on the vast generality of the gap approach, which provides, indeed, a rather good comprehension of the inner mechanisms of ES, and can be extended to a broad variety of systems.

4.2.3. Explosive synchronization in frequency-weighted networks of oscillators

Motivated by the results reviewed in the previous Subsection, Ref. [18] further extended the study to the case of a network with given frequency distribution and architecture, for which the only action that an external operator can perform is a weighting procedure on the already existing links. Namely, Ref. [18] combined the information on the frequency mismatch of the two end oscillators of a link with that of the link betweenness (in the more general case), and demonstrated that such a weighting mechanism has the effect of inducing (or enhancing) ES for both homogeneous and heterogeneous graph topologies, as well as for any (symmetric or asymmetric) frequency distribution.

We follow the path of reasoning of Ref. [18], and start by modifying Eq. (28) as

$$\dot{\theta}_i = \omega_i + \frac{\lambda}{\langle k \rangle} \sum_{i=1}^N \Omega_{ij}^\alpha \sin(\theta_j - \theta_i), \quad (46)$$

where

$$\Omega_{ij}^\alpha = a_{ij} |\omega_i - \omega_j|^\alpha, \quad (47)$$

is the weight factor to be applied to the (existing) link between nodes i, j , being a_{ij} the elements of the adjacency matrix that uniquely defines the network, and α a constant parameter which possibly modulates the weights. The strength of the i th node (the sum of all its links weights) is then $s_i = \sum_j \Omega_{ij}^\alpha$.

Fig. 40(a) reports the results for a ER network of size $N = 500$, and several frequency distributions $g(\omega)$ within the range $[0, 1]$. For the simplest case of a uniform frequency distribution, the un-weighted network ($\alpha = 0$ in Eq. (46)) displays a smooth, second-order like transition [dark blue curve in Fig. 40(a)], whereas the effect of a linear weighting factor ($\alpha = 1$) is that of inducing ES in the system, with an associated hysteresis in the forward (solid line) and backward (dashed line) transitions.

More importantly, such a drastic change in the nature of the transition seems to be independent of the specific frequency distribution $g(\omega)$, as long as they are defined in the same frequency range $[0, 1]$. The results are identical for symmetric (homogeneous, Gaussian, bimodal derived from a Gaussian) and for asymmetric (Rayleigh, a Gaussian centered at 0 but with just having the positive half) frequency distributions, indicating the existence of a sort of *universal* behavior for homogeneous graph's topologies.

Fig. 40(b) shows the existence of a parabolic relationship between the strengths and the natural frequencies of the oscillators (associated with the passage from a smooth to an explosive transition). This relationship has been obtained analytically (see Eq. (54)), and perfectly fits the numerical results. The emergent correlation features shape a bipartite-like network, where low and high frequency oscillators are the ones with maximal overall strength.

Furthermore, the weighting procedure inducing ES appears to be really general, because it applies for a large family of detuning dependent functions. As an example, Fig. 41 describes the case of nonlinear weighting procedures [$\alpha \neq 1$ in

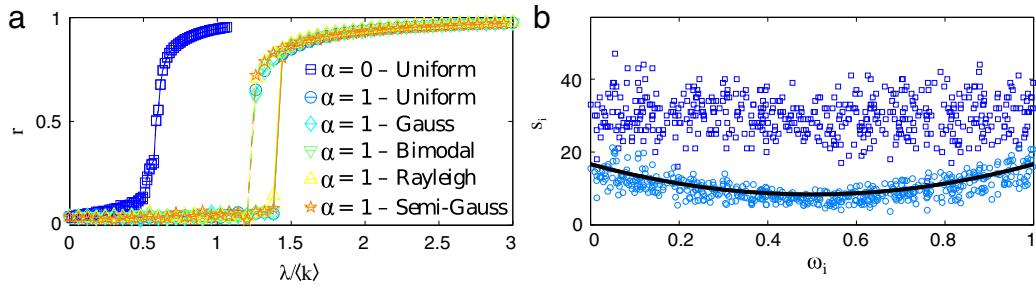


Fig. 40. (a) Synchronization transitions (r vs. $\lambda/\langle k \rangle$) for ER networks of size $N = 500$, $\langle k \rangle = 30$. The Figure reports the un-weighted case ($\alpha = 0$) (blue squares), and linearly weighted cases ($\alpha = 1$) for several frequency distributions within the range $[0, 1]$ (detailed in the legend). (b) Node strengths s_i (see text for definition) vs. natural frequencies ω_i , for the un-weighted (dark blue squares) and weighted (light blue circles) networks reported in (a). The solid line corresponds to the analytical prediction. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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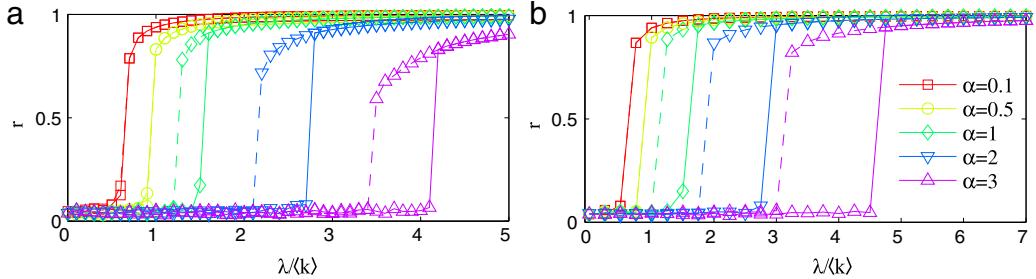


Fig. 41. (Color online). Synchronization transitions for ER networks ($N = 500$), with uniformly distributed frequencies in the $[0, 1]$ range and nonlinear weighting functions Ω_{ij}^α . Reported curves refer to several α values, from sub-linear to super-linear weighting (see legend in panel b). (a) ER networks, $\langle k \rangle = 30$, (b) regular random networks, $k = 30$. In all cases, forward and backward simulations correspond respectively to solid and dashed lines.

Source: Adapted from Ref. [18].

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Eq. (46)]. There, $N = 500$ and $\langle k \rangle = 30$ are set, and ER graphs (Fig. 41(a)), together with random networks (where each node has exactly the same number of connections, $k_i = \langle k \rangle = 30$) are considered (Fig. 41(b)) (the latter case is obtained by a simple configuration model, imposing a δ -Dirac degree distribution).

A generic non-linear function of the frequency mismatch is able to induce ES in all topologies, and the effect of a super-linear ($\alpha > 1$) weighting (a sub-linear ($\alpha < 1$) weighting) is that of enhancing (reducing) the width of the hysteretic region.

When networks are heterogeneous, Ref. [18] suggests to use, as a new weighting function,

$$\tilde{\Omega}_{ij} = a_{ij} |\omega_i - \omega_j| \frac{\ell_{ij}^\beta}{\sum_{j \in \mathcal{N}_i} \ell_{ij}^\beta}, \quad (48)$$

with β being an extra tuning parameter, and ℓ_{ij} being the edge betweenness associated to the link a_{ij} , defined as the total number of shortest paths (between pairs of nodes in the network) that makes use of that edge. The results are reported in Fig. 42. While the case $\beta = 0$ corresponds to a smooth transition, moderate (positive or negative) values of β establish abrupt transitions to synchronization.

In order to rigorously predict the onset (and nature) of the explosive transition, Ref. [18] analytically examines the thermodynamic limit in which N oscillators form a fully connected graph: $\dot{\theta}_i = \omega_i + \frac{\lambda}{N} \sum_{j=1}^N |\omega_i - \omega_j| \sin(\theta_j - \theta_i)$. By assuming the following definitions,

$$\frac{1}{N} \sum_{j=1}^N \Omega_{ij} \sin \theta_j := A_i \sin \phi_i, \quad (49)$$

$$\frac{1}{N} \sum_{j=1}^N \Omega_{ij} \cos \theta_j := A_i \cos \phi_i, \quad (50)$$

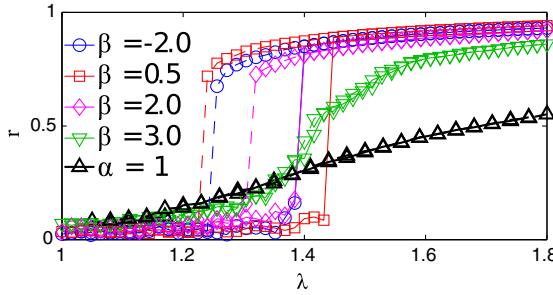


Fig. 42. (Color online). Synchronization transition for SF networks using different values of the β parameter [in Eq. (48)]. A clear passage from a second-order to a first-order like transition is seen.

Source: Adapted from Ref. [18].

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the evolution equations can be expressed in terms of trigonometric functions as $\dot{\theta}_i = \omega_i + \lambda A_i \sin(\phi_i - \theta_i)$. In the thermodynamic limit, the solution (in the co-rotating frame) is

$$\omega = \lambda A_\omega \sin(\theta_\omega - \phi_\omega). \quad (51)$$

The definition of A_ω and ϕ_ω implies that

$$\begin{aligned} F(\omega) &:= A_\omega \sin \phi_\omega = \int g(x)|\omega - x| \sin \theta(x) dx, \\ G(\omega) &:= A_\omega \cos \phi_\omega = \int g(x)|\omega - x| \cos \theta(x) dx. \end{aligned} \quad (52)$$

When all oscillators are close to synchronization, one can assume that $\cos \theta(x) \approx r$, and therefore $G(\omega) \simeq rs(\omega)$, where $s(\omega)$ is the strength of a node with frequency ω . As a consequence, Eq. (51) takes the form

$$\frac{2}{r\lambda} g(\omega)\omega = F''(\omega)s(\omega) - F(\omega)s''(\omega), \quad (53)$$

which is a second order ordinary differential equation, whose integration yields $F(\omega)$. For instance, given a uniform distribution $g(\omega)$ in the interval $[-a/2, +a/2]$, the resulting strength is the second order polynomial

$$s(\omega) = a \left[\left(\frac{\omega}{a} \right)^2 + \frac{1}{4} \right], \quad (54)$$

which fits the numerical results of Fig. 40(b). The analysis allows determining the dependence of r on λ as $r = \int g(x) \cos \theta(x) dx = \int g(x) \sqrt{1 - \sin^2 \theta(x)} dx$, where $H(z) := \frac{4}{4+\pi} \left[\frac{z}{1+z^2} + \arctan(z) \right]$. In Ref. [18] it is shown that r is indeed a multi-valued function (for values of λ above a given threshold). Similar results are also reported in Ref. [372], which used a slightly modified weighting function

$$\Omega_{ij} = \frac{a_{ij} N \langle k \rangle |\omega_i - \omega_j|^\beta}{\sum_{p=1}^N \sum_{q=1}^N a_{pq} |\omega_p - \omega_q|^\beta}. \quad (55)$$

In Ref. [61], Zhang et al. proposed a simpler alternative weighting method:

$$\dot{\theta}_i = \omega_i + \frac{\lambda |\omega_i|}{\langle k_i \rangle} \sum_{i=1}^N a_{ij} \sin(\theta_j - \theta_i). \quad (56)$$

For symmetric frequency distributions $g(\omega)$ (such as a random, a Lorentzian, or a Gaussian distribution), the model of Eq. (56) induces ES.

The model is, however, less effective in generating ES for asymmetric frequency distributions. In the mean field approximation, Eq. (56) yields $r \sin(\Psi - \theta_i) = \frac{1}{N} \sum_j \sin(\theta_j - \theta_i)$. Therefore, one can write

$$\dot{\theta}_i = \omega_i + \lambda |\omega_i| r \sin(\Psi - \theta_i). \quad (57)$$

Then, setting as reference frame the one rotating with the average phase of the system, we have $\Psi(t) = \Psi(0) + \langle \omega \rangle t$, where $\langle \omega \rangle$ is the average frequency of the oscillators. For a symmetric $g(\omega)$, we have $\langle \omega \rangle = 0$. Letting $\Delta\theta_i = \theta_i - \Phi$, Eq. (57) becomes

$$\Delta\dot{\theta}_i = \omega_i - \lambda |\omega_i| r \sin(\Delta\theta_i). \quad (58)$$

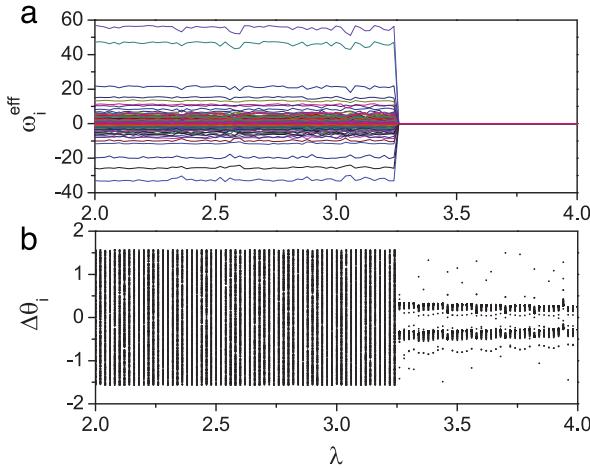


Fig. 43. (Color online). The evolutionary dynamics along the forward continuation in the network model of Eq. (56). (a) ω_i^{eff} vs. λ , (b) $\Delta\theta_i = \theta_i - \psi$ vs. λ .
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In the phase locking state, $\Delta\dot{\theta}_i = 0$, and therefore one has

$$\Delta\theta_i = \begin{cases} \Delta\theta_+ = \arcsin\left(\frac{1}{\lambda r}\right) & \omega_i > 0, \\ \Delta\theta_- = \arcsin\left(-\frac{1}{\lambda r}\right) & \omega_i < 0, \end{cases} \quad (59)$$

which implies that two equal clusters form: one including the clockwise, and the other the counterclockwise oscillators. Both values of $\Delta\theta_+$ and $\Delta\theta_-$ gradually approach zero with increasing λ . This fully confirms the observation in Fig. 43(b). Then, the order parameter takes the form $r = \frac{1}{2}(e^{i\Delta\theta_+} + e^{i\Delta\theta_-})$. Substituting Eq. (59) in such a latter expression, one obtains $r^2 = \frac{\lambda + \sqrt{\lambda^2 - 4}}{2\lambda}$, which is independent on $g(\omega)$, provided it is symmetric. This result implies that ES exists only for $\lambda \geq \lambda_c^b = 2$, where r has a strong discontinuity from $r > \sqrt{5} \approx 0.707$ to $r \approx 0$.

In the case of asymmetric frequency distributions, a similar calculation gives instead

$$r(x) = \int \sqrt{1 - \left(\frac{\omega - \langle\omega\rangle}{x\omega}\right)^2} g(\omega) H\left(1 - \left|\frac{\omega - \langle\omega\rangle}{x\omega}\right|\right) d\omega,$$

where $x = r\lambda$, and $H(x)$ is the Heaviside step function. In this case, there is a strong dependence on $g(\omega)$, and Ref. [61] shows that continuous or discontinuous transitions emerge for different frequency distributions.

The analytic study of Eq. (56) was tackled by Hu et al. in Ref. [373], where the forward critical coupling was calculated with the use of the method introduced in Ref. [237]. In the limit $N \rightarrow \infty$ a density function $\rho(\theta, \omega, t)$ can be defined, which denotes the fraction of oscillators with frequency ω whose phases have values between θ and $\theta + d\theta$ at time t . $\rho(\theta, \omega, t)$ satisfies the normalization condition $\int_0^{2\pi} \rho(\theta, \omega, t) d\theta = 1$ for all θ and t . The evolution of $\rho(\theta, \omega, t)$ is governed by the continuity equation:

$$\frac{\delta\rho}{\delta t} + \frac{\delta(\rho v)}{\delta\theta} = 0. \quad (60)$$

Using the expression obtained in Eq. (57), the velocity in the thermodynamical limit can be written as

$$v = \omega + \lambda|\omega|r \sin(\phi - \theta). \quad (61)$$

When the coupling strength is relatively small, all oscillators are rotating in the unit circle according (almost exclusively) to their natural frequencies. This corresponds to an incoherent state $\rho(\theta, \omega, t) = 1/(2\pi)$. Therefore, the point at which such a state loses its stability can be extracted, providing a prediction for the forward critical point. Specifically, let one consider a small perturbation from the incoherent state, i.e. $\rho(\theta, \omega, t) = \frac{1}{2\pi} + \epsilon\eta(\theta, \omega, t)$. Then, the linearized form of the continuity equation (60) is

$$\frac{\delta\eta}{\delta t} = -\omega \frac{\delta\eta}{\delta\theta} + \frac{\lambda r' |\omega| \sin(\phi - \theta)}{2\pi}, \quad (62)$$

Table 4

Critical equations and critical points λ_f^c (calculated in Ref. [373]) for a triangular, a Lorentzian, a Gaussian and a bimodal Lorentzian frequency distribution $g(\omega)$.

Frequency distribution $g(\omega)$	Critical equation	λ_f^c
$g(\omega) = (D - \omega)/D^2, \omega < D, 0 \text{ otherwise}$	$1 = \lambda \left[\frac{z}{2} \ln \left(1 + \frac{1}{z^2} \right) - z + z^2 \arctan \left(\frac{1}{z^2} \right) \right], z = \mu/D$	~ 2.65
$g(\omega) = \frac{\Delta}{\pi(\omega^2 + \Delta)}$	$\frac{\pi}{\lambda} = \frac{z \ln z}{z^2 - 1}, z = \mu/\Delta$	4
$g(\omega) = \frac{1}{\pi\sigma} \exp \left(-\frac{\omega^2}{\pi\sigma^2} \right)$	$1 = \frac{\lambda z}{2\sqrt{2\pi}} \exp(z^2/2) [-E_i(-z^2/2)], z = \mu/\sigma, E_i(z) = \int_{-\infty}^z \frac{e^u}{u} du$	~ 2.68
$g(\omega) = \frac{1}{2\pi} \left[\frac{\Delta}{(\omega - \omega_0)^2 + \Delta^2} + \frac{\Delta}{(\omega + \omega_0)^2 + \Delta^2} \right]$	$\frac{2\pi}{\lambda} = \ln \left(\frac{\sqrt{1+\delta^2}}{\mu} \right) \left[\frac{(1-\mu)\mu}{(1-\mu)^2 + \delta^2} + \frac{(1+\mu)\mu}{(1+\mu)^2 + \delta^2} \right], \delta = \omega_0/\Delta$	$4/\sqrt{1 + \delta^2}$

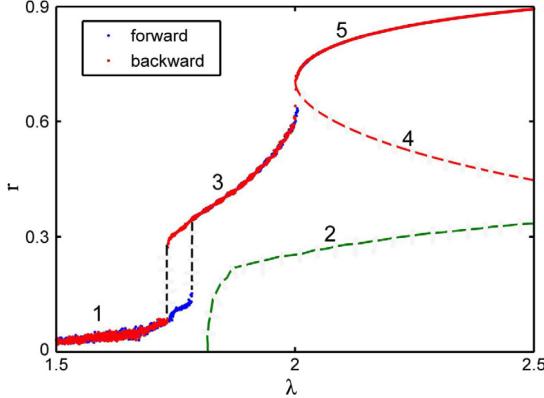


Fig. 44. Characterization of various coherent states in the phase diagram r vs. λ , for a uniform frequency distribution $g(\omega) = 1/2, \omega \in (-1, 1)$, and for a fully connected $N = 50,000$ Kuramoto ensemble. The distinct branches correspond to: the incoherent state (1), the (unstable) traveling wave (TW) state predicted by the mean-field theory (2), the standing wave state (3), the unstable (4) and the stable (5) two-cluster synchronous states. The blue and red lines denote the forward and the backward transitions, respectively. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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where

$$re^{i\phi} = \epsilon r'e^{i\phi} = \epsilon \int_0^{2\pi} \int_{-\infty}^{\infty} e^{i\theta} \eta(\theta, \omega, t) d\omega d\theta. \quad (63)$$

Now $\eta(\theta, \omega, t)$ can be expanded into a Fourier series as $\eta(\theta, \omega, t) = b(\omega)e^{i\mu t} e^{i\theta} + b^*(\omega)e^{i\mu t} e^{-i\theta} + O(\theta, \omega, t)$, where $O(\theta, \omega, t)$ stands for higher harmonics. The equation becomes

$$\mu b(\omega) = -i\omega b(\omega) + \frac{\lambda|\omega|}{2} \int_{-\infty}^{\infty} b(v)g(v)dv, \quad (64)$$

and can be solved in a self-consistent way. In particular, when $g(\omega)$ is an even function, Eq. (64) becomes

$$1 = \frac{\lambda}{2} \int_{-\infty}^{\infty} \frac{\mu|\omega|}{\mu^2 + \omega^2} g(\omega)d\omega, \quad (65)$$

i.e., when $\text{Re}(\mu)$ changes from negative to positive, the incoherent state loses its stability. For any specific $g(\omega)$ one can then use this condition to determine the critical coupling strength λ_f^c for the forward phase transition. Several examples of this calculation are reported in Table 4.

Xu et al. [374] completed the analysis initiated in Ref. [373], and calculated all possible steady states in the model of Eq. (56), which include the incoherent state and the two-cluster synchronous state already deduced in Ref. [373], but also a traveling and a standing wave state. Ref. [374] demonstrated that the incoherent state is only neutrally stable below the synchronization threshold. The amplitude equations near the bifurcation point are derived according to the center-manifold reduction, and non-stationary standing wave states are predicted to exist in the model. Fig. 44 reports all possible solutions for a fully connected Kuramoto ensemble with uniform frequency distribution $g(\omega) = 1/2, \omega \in (-1, 1)$.

More detailed information on the microscopic mechanisms underlying ES in the model of Eq. (56) can be found in Ref. [19], where a local order parameter, r_{ij} , is introduced

$$r_{ij} = \left| \lim_{T \rightarrow \infty} \frac{1}{T} \int_t^{t+T} e^{(\theta_i(t) - \theta_j(t))} dt \right|, \quad (66)$$

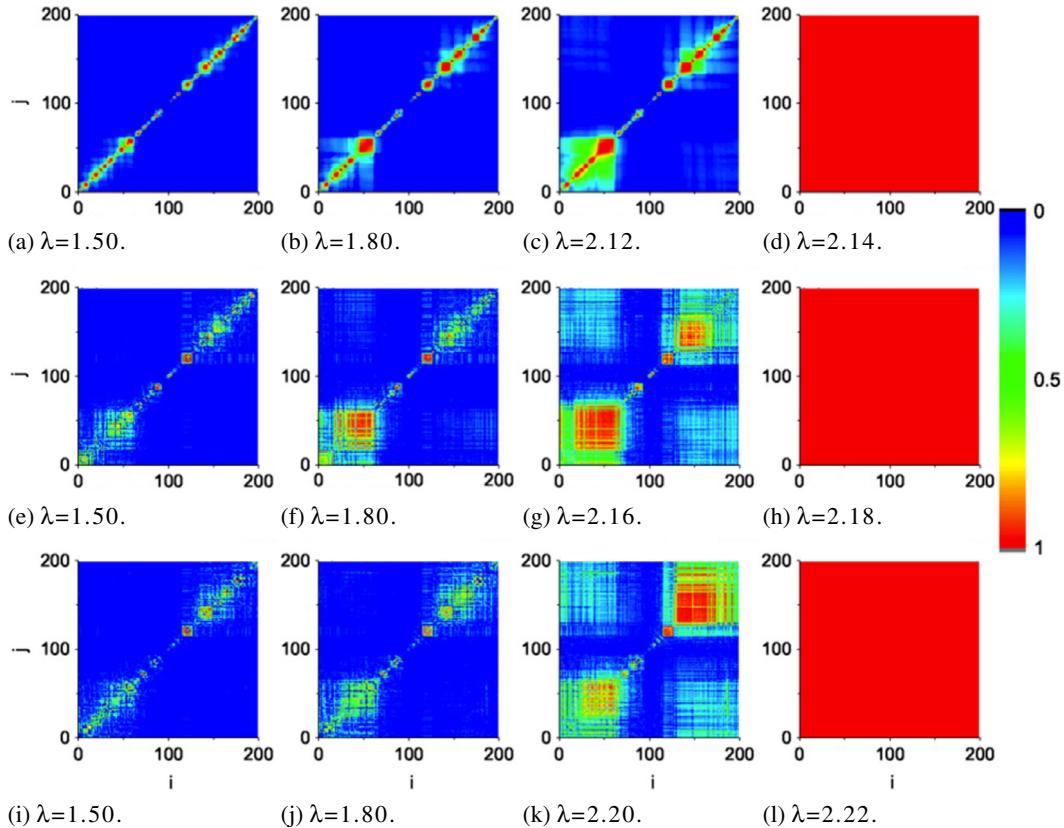


Fig. 45. Plots of the matrix r_{ij} (Eq. (66)) for fully connected (first line), ER (second line) and uncorrelated CM (third line) networks. Oscillators are labeled in the ascending order of the frequency ω_i . The coupling strengths are $\lambda = 1.5, 1.8, 2.12$, and 2.14 in (a)–(d) ($\lambda_c = 2.13$); $\lambda = 1.5, 1.8, 2.16$, and 2.18 in (e)–(h) ($\lambda_c = 2.17$); and $\lambda = 1.5, 1.8, 2.20$, and 2.22 in (i)–(l) ($\lambda_c = 2.21$).

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where T is a sufficiently long time window. r_{ij} is 1 for any two phase-locked oscillators, zero for all pairs of fully uncorrelated oscillators, and will take a value between 0 and 1 for any two partially correlated oscillators. Fig. 45 reports the values of r_{ij} for: (a)–(d) fully connected, (e)–(h) ER and (i)–(l) uncorrelated CM networks, for four representative λ values. Only small synchronized clusters of oscillators exist for $\lambda < \lambda_c$, while a giant synchronized cluster shows up suddenly right after λ_c , indicating that the small synchronized clusters suddenly merge together right at λ_c .

The similarity with the Achlioptas process of EP is even clearer from looking at Fig. 45, where it is seen that, below λ_c , links are always generated among those nodes with close frequencies. As a consequence, separated clusters form for λ up to λ_c , where instead (and suddenly) all clusters merge together to mold a giant one. To quantify the analogy with EP, an equivalent suppressing rule is introduced. From Eq. (57), the evolution of the phase difference $\Delta\theta_{ij} = \theta_i - \theta_j$ is given by

$$\dot{\Delta\theta}_{ij} = \omega_i + |\lambda_i|r \sin(\Psi - \theta_i) - \omega_j + |\lambda_j|r \sin(\Psi - \theta_j). \quad (67)$$

When the two oscillators i and j are phase-locked, one has $\dot{\Delta\theta}_{ij} = 0$ and therefore

$$\omega_i - \omega_j = |\lambda_i|r \{\sin(\Psi - \theta_i) - \sin(\Psi - \theta_j)\}, \quad (68)$$

whose maximum value gives a necessary condition for the phase-locking between two oscillators:

$$\frac{|\omega_i - \omega_j|}{|\omega_i| + |\omega_j|} \leq \lambda r. \quad (69)$$

Eq. (69) can be considered as a *suppressive rule* for pair wise synchronization: as r takes a rather small value when $\lambda < \lambda_c$, only those pairs of oscillators with smaller frequency differences can satisfy the condition (69) and thus form synchronized clusters. When all the free oscillators have been attracted to the synchronized clusters, the further increase of λ cannot make the synchronized clusters become larger, but makes the clusters to merge each other suddenly, this way producing the significant jump on r observed in the explosive transition.

Ref. [19] shows also that the suppressive rule (69) can actually guide manipulative processes to enhance or quash ES. For instance, one can randomly pick a pair of nodes i, j and then artificially re-adjust the network by exchanging their frequencies

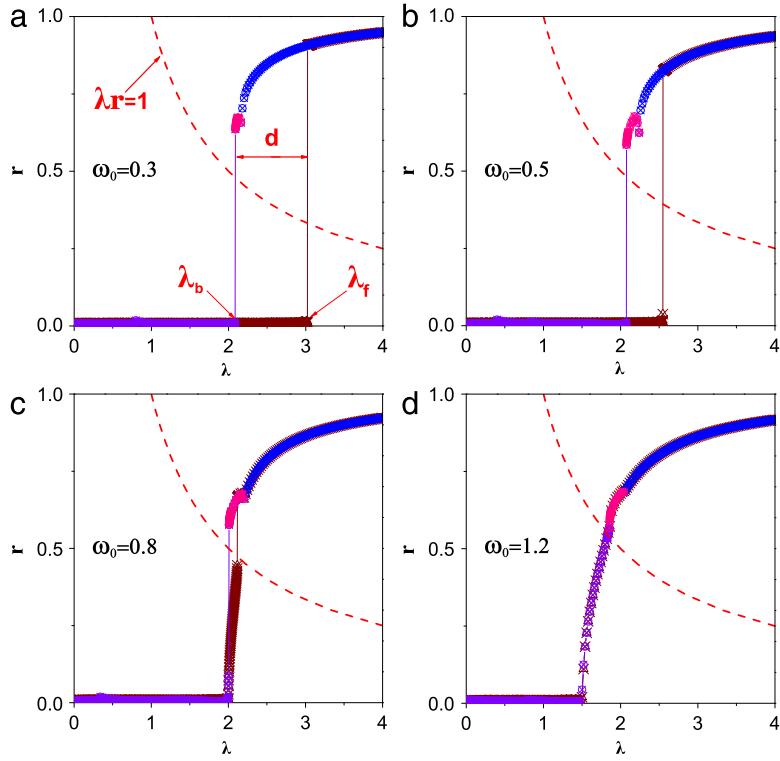


Fig. 46. (Color online). r vs. λ . With the increase of ω_0 , the synchronization transition converts from a first-order type into a second-order one. From (a) to (d), $\omega_0 = 0.3, 0.5, 0.8, 1.2$. The dashed lines correspond to $\lambda r = 1$. A Lorentzian frequency distribution is used, and $N = 10,000$.

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if the quantity

$$S_{ij} = \sum_{l=1}^{\Gamma_i} \frac{|\omega_i - \omega_l|}{|\omega_i| + |\omega_l|} + \sum_{l=1}^{\Gamma_j} \frac{|\omega_j - \omega_l|}{|\omega_j| + |\omega_l|} \quad (70)$$

decreases (or increases), where Γ_i, Γ_j stand for the neighborhoods of nodes i, j , resulting in a enhanced (suppressed) ES, as it occurs for the Achlioptas algorithm in EP.

Ref. [375] considers the effect of using frequency distributions $g(\omega)$ not centered in 0 (see Eq. (56)), and specifically studies the case of uni-modal frequency distributions (Lorentzian, Gaussian, triangle), displacing their central frequency from 0 to $\omega_0 \neq 0$, which is then used as a parameter.

Fig. 46 plots the phase diagrams when increasing ω_0 , and shows that the hysteresis area significantly shrinks, mainly due to the decrease of the forward transition point λ_c^f , while the backward transition point λ_c^b remains unchanged.

An interesting view on microscopic processes in both the first-order and second-order cases can be obtained from Fig. 47. Here, the incoherent steady state is explored for two values of the central frequency: $\omega_0 = 0.5$ (Fig. 47(a)–(b), which causes ES), and $\omega_0 = 1.2$ (Fig. 47(c)–(d), for which the transition is reversible). In all the cases of Fig. 47 the coupling is $\lambda = 1.5$, i.e. below the synchronization threshold, and therefore $r \simeq 0$. However, essential microscopic differences exist between these incoherent states. While the phases are randomly distributed just below the forward threshold in the ES case (Fig. 47(a)), the incoherent state of the corresponding backward continuation (Fig. 47(b)) is non-trivial, showing the remaining of small clusters with fine structures. In the second order case with $\omega_0 = 1.2$, these microscopic structures appear in the incoherent state of both the forward and the backward (Fig. 47(c)–(d)) continuations. Such microscopic differences reveal that, as the frequencies distribution is not symmetric around 0, formation of micro-clusters occurs, and prevents ES.

This is actually due to the suppression rule of Eq. (69). All the pairs of nodes i, j whose frequencies ω_i, ω_j , have different sign fulfill $|\omega_i - \omega_j|/(|\omega_i| + |\omega_j|) = 1$, and therefore, being $r \simeq 0$ in the incoherent state, they do not synchronize. An asymmetrical frequency distribution will reduce the number of possible pairs of nodes with frequencies of opposite sign, and therefore an increasing number of pairs will be able to form synchronization seeds, so that eventually ES is destroyed for large enough ω_0 .

Following the same line of investigation of microscopic mechanisms underlying ES, Ref. [376] proposed the use of an effective topological network whose structure explicitly reflects the interplay between the topology and dynamics of the

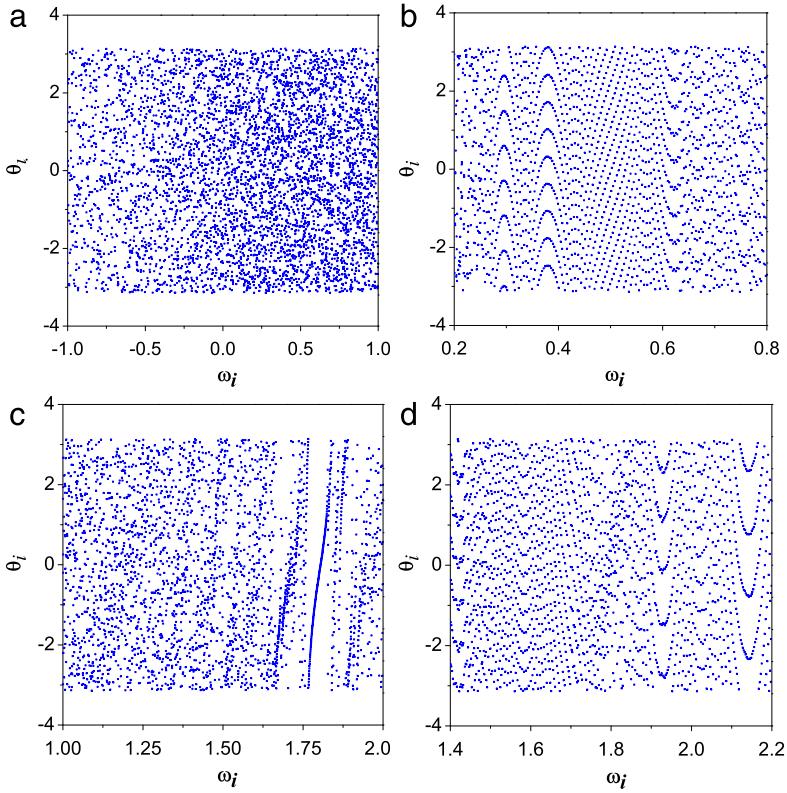


Fig. 47. Instantaneous phase distribution θ_i vs. natural frequency ω_i for all the nodes of a ER of $N = 10,000$ nodes, coupling strength $\lambda = 1.5$, for Lorentzian frequency distribution centered in $\omega_0 = 0.5$ ((a) and (b) corresponding to a system able to achieve ES), and $\omega_0 = 1.2$ ((c) and (d) corresponding to a system with second order transition). Left (right) panels correspond to forward (backwards) continuations.

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original system in an individual basis, and therefore, is able to easily identify the nodes which actually act as synchronization seeds. The effective adjacency matrix is given by

$$C_{ij} \equiv A_{ij} \left(1 - \frac{\Delta\omega_{ij}}{\Delta\omega_{max}} \right), \quad (71)$$

where $\Delta\omega_{ij} = |\omega_i - \omega_j|$ is the frequency detuning, and $\Delta\omega_{max}$ the maximum possible detuning present in the system in order to guarantee $C_{ij} \geq 0$.

In order to quantify the role of each node in the synchronization process, Ref. [376] extracts the most important nodes in the network defined by \mathbf{C} , i.e. the standard eigenvector centrality measure of \mathbf{C} , obtaining the *effective centrality* vector \mathbf{A}^C , whose i th component quantifies the potential of node i to behave as a seed of synchronization.

Fig. 48 reports the comparison between \mathbf{A}^C and its topological counterpart \mathbf{A}^A , the eigenvector centrality extracted from the original adjacency matrix \mathbf{A} . For ER networks (Fig. 48(a)), the distribution of the components of the vector \mathbf{A}^C as a function of the corresponding node's natural frequencies shows the existence of many seeds of synchronization with natural frequencies close to $\Omega_s = 0$. This allows characterizing the connection between the micro-scale and the macro-scale of the network in a much better way than \mathbf{A}^A , whose components are instead uniformly distributed. For SF networks (Fig. 48(b)), the synchronization seeds are the hubs, and therefore \mathbf{A}^C and \mathbf{A}^A provide essentially the same information.

4.2.4. Noise induced explosive synchronization

As discussed in Section 4.2.1, a simple (yet artificially engineered) degree–frequency correlation gives explosive synchronization on sufficiently heterogeneous networks [15,16]. An extension of such artificial process is proposed in Ref. [377], where the frequencies of oscillators are correlated with (but not fully determined by) the local network's properties, such as the degree. More precisely, the evolution of phase oscillator θ_i is taken to be

$$\dot{\theta}_i = \omega'_i + \lambda \sum_{i=1}^N A_{ij} \sin(\theta_j - \theta_i), \quad (72)$$

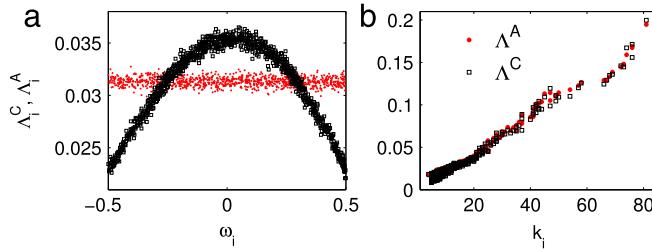


Fig. 48. (Color online). Comparison between synchronization centrality Λ_i^C (black squares) and topological centrality Λ_i^A (red dots). (a) ER networks, $\langle k \rangle = 50$, Λ_i^C and Λ_i^A are reported vs. the nodes' natural frequencies ω_i ; (b) SF networks, $\langle k \rangle = 12$, Λ_i^C and Λ_i^A are plotted vs. the node degrees k_i . All data refer to ensemble averages over 100 different network realizations.

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where $\omega'_i = k_i + \xi_i$, where ξ_i is a noise term taken from a random homogeneous distribution in the range $[-\varepsilon, \varepsilon]$. $\varepsilon = 0$ returns the case of Ref. [15]. Ref. [377] considers realistic *C. elegans* neural networks, stretched exponential networks and SF networks with mild heterogeneity (i.e. $\gamma > 3$ in the degree distribution $P(k) \sim k^{-\gamma}$). Given that the synchronized solution is composed of a cluster with uniform angular velocity Ω , the rotating reference frame involving the change of variable $\phi_i = \theta_i - \Omega t$ is defined as follows

$$\dot{\phi}_i = (\omega_i - \Omega) + \lambda \sum_{j=1}^N A_{ij} \sin(\phi_j - \phi_i), \quad (73)$$

and the local order parameter is defined as

$$r_i e^{i\psi_i} = \sum_{j=1}^N A_{ij} e^{i\phi_j}, \quad (74)$$

having a magnitude $r_i \in [0, k_i]$. Substituting Eq. (74) into Eq. (73), one gets

$$\dot{\phi}_i = (\omega_i - \Omega) + \lambda r_i \sin(\psi_i - \phi_i). \quad (75)$$

From Eq. (75) it is clear that, if $|\omega_i - \Omega| \leq \lambda r_i$, ϕ_i can reach a steady point expressed by $\sin(\phi_i - \psi_i) = (\omega_i - \Omega)/\lambda r_i$ and becomes phase-locked; otherwise the oscillator i drifts forever.

The self-consistency condition for the steady-state order parameter r is [377]

$$r = \langle k \rangle^{-1} \int \int_{|\omega - \Omega| \leq \lambda r k} P(k, \omega) k \sqrt{1 - \left(\frac{\omega - \Omega}{\lambda r k} \right)^2} d\omega dk, \quad (76)$$

which can be used to predict the evolution of synchronization, and where $P(k, \omega)$ is the so-called joint probability distribution of degree and frequency.

In the absence of noise, the transition from incoherent to synchronized states seems to be second-order like for both real and synthetic networks. The trend can be accurately predicted by the self-consistency condition (76), and is shown in Fig. 49(b). For sufficient large noise, however, ES takes place with a clear bistable region on the less heterogeneous networks (see Fig. 49(d)). If λ_c^f and λ_c^b denote the forward and backward critical coupling strengths, the width of hysteresis loop is $\lambda_c^f - \lambda_c^b$. From Fig. 49(e) and (f), one sees that the transition from incoherent to synchronized phases is second-order like until a critical value of the noise ε_c (denoted by the dashed lines). Therefore, ES can be promoted by noise mechanisms.

Ref. [378] suggested to focus on the angular velocity v_i of oscillator i . With Gaussian noise $\eta_i(t)$, the dynamics is ruled by

$$\begin{aligned} \dot{\theta}_i &= v_i, \\ m\dot{v}_i &= -\gamma v_i + \lambda r \sin(\psi - \theta_i) + \gamma \omega_i + \sqrt{\gamma} \eta_i(t), \end{aligned} \quad (77)$$

where m accounts for an inertia term, γ denotes the friction constant, and r is the order parameter. In particular, Ref. [378] assumed $\langle \eta_i(t) \rangle = 0$, $\langle \eta_i(t) \eta_j(t') \rangle = 2T\delta_{ij}\delta(t-t')$, where T is a stochastic noise (like a temperature in units of the Boltzmann constant). Furthermore, a uni-modal $g(\omega)$ is considered, with width σ . At $\sigma = 0$, the dynamics resembles that of a Brownian mean-field model [379]. In the absence of inertia, $T = 0$ reproduces the Kuramoto model [23,237,238], while $T \neq 0$ gives the model of Ref. [336].

Fig. 50(a) reports the phase diagram in the parameter space. The shaded blue surface is a first-order transition surface. Inside (outside) this surface, the synchronization transition is abrupt and explosive (continuous). To furnish a better intuitive illustration, Fig. 50(b) shows the behavior of r as a function of noise T . For large T , the change of r is continuous and

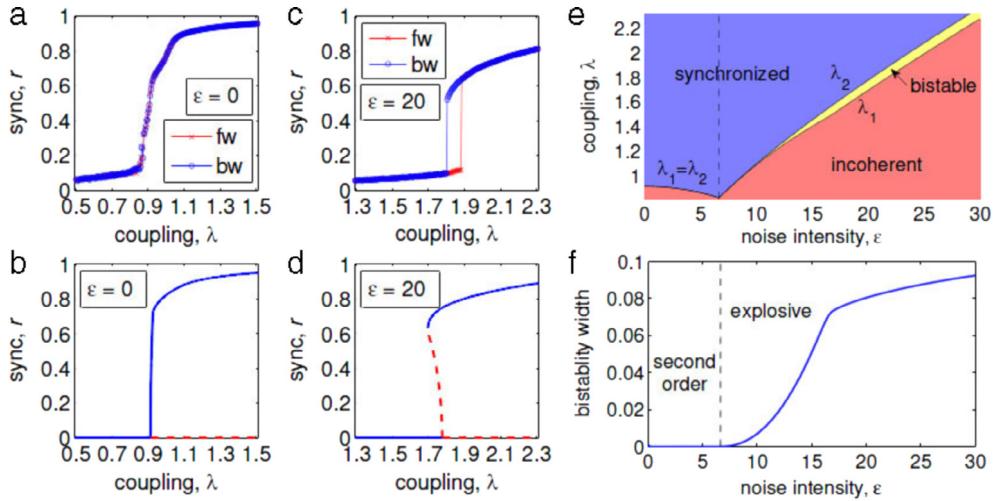


Fig. 49. (Color online). (a)–(d) r vs. λ for different values of ε from simulations (top) and theoretical analysis (bottom). (e) Phase space with critical coupling strength λ_c^f and λ_c^b separating incoherent, synchronized and bistable regions. (f) Bistability width $\lambda_c^f - \lambda_c^b$ as a function of noise λ . The used networks are SF networks with size $N = 10^3$.

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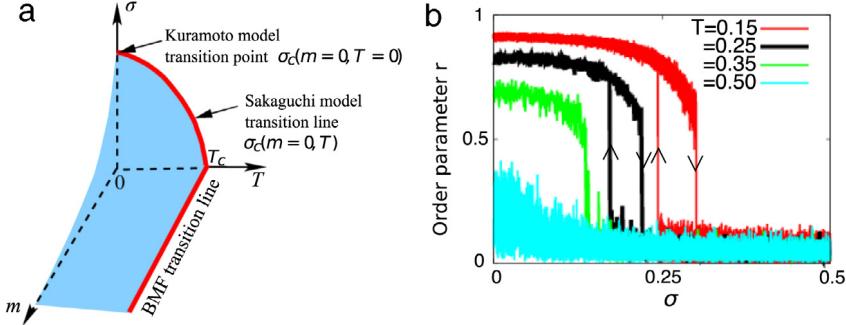


Fig. 50. (a) Phase diagram of synchronization in the parameter space of inertia m , noise T and width σ of the frequency distribution. The shaded blue surface (red line) indicates the first-order (second-order) transition. (b) Synchronization level r vs. σ for different T with fixed inertia $m = 10$. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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overlapped between forward and backward transitions. With decreasing T , however, ES takes place and the width of hysteresis loop becomes larger.

The effect of noise on synchronization was actually observed in other types of synchronization models [380–386]. For instance, for a positive correlation and sufficient wide frequency distribution, an intermediate noise can lead to the maximal hysteresis loop in the synchronization diagram of FitzHugh–Nagumo oscillators operating on top of SF networks [63].

4.3. Explosive synchronization in generalized Kuramoto models

In the original Kuramoto model (Eq. (13)), the sign of the coupling strength is assumed to be positive, implying that the interactions among oscillators are always attractive. However, in many practical circumstances, the interaction among the constituents of an ensemble can be repulsive or suppressive, which would correspond to a negative coupling in the model [348,387–390].

For this purpose, Ref. [391] generalized Eq. (13), by allowing the coupling strength λ to assume both positive and negative values. The model is described by

$$\dot{\theta}_i = \omega_i + \frac{\lambda_i}{N} \sum_{j=1}^N \sin(\theta_j - \theta_i), \quad i = 1, \dots, N, \quad (78)$$

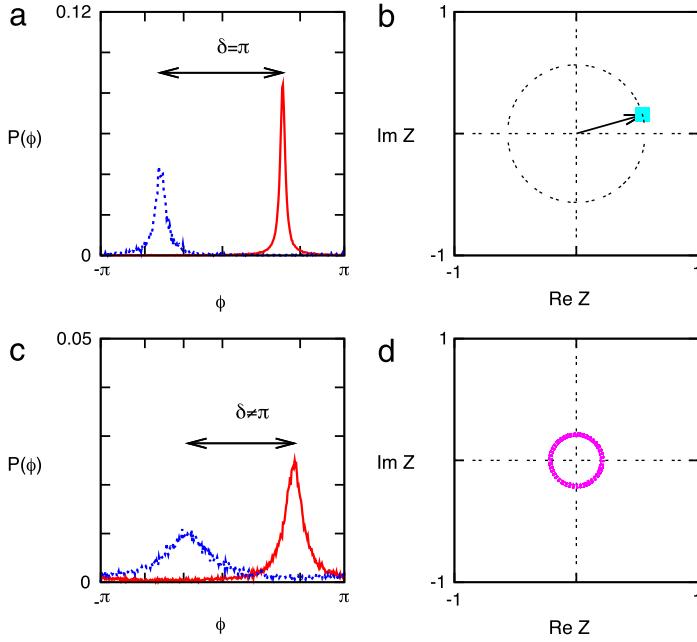


Fig. 51. The π and TW states in model (78). $N = 25, 600$, $\gamma = 0.05$, $\lambda_1 = -0.5$, $\lambda_2 = 1.0$. (a) Phase distribution for the π state; the mean phase difference δ between the conformist (red) and contrarian (blue) oscillators is given by $\delta = \pi$. (b) The order parameter for the π state corresponds to a fixed point. (c) Phase distribution for the TW state; $\delta \neq \pi$. (d) The order parameter for the TW state traces a circle, implying a non-zero mean phase velocity. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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with a Lorentzian frequency distribution $g(\omega) = \gamma/[\pi(\omega^2 + \gamma^2)]$. Here, λ_i (the coupling strength of the i th oscillator to the mean field) is a binary variable, i.e., either $\lambda_i = \lambda_1 < 0$ or $\lambda_i = \lambda_2 > 0$. Accordingly, oscillators in the ensemble are divided into two populations: those with positive λ_i will behave like *conformists*, whereas those with negative λ_i will tend to act as *contrarians*. Initially, the system is set in the incoherent state where only contrarian oscillators exist. Then contrarians are gradually flipped into conformists according to certain strategies. When the proportion of conformists, denoted by p , exceeds a certain threshold, the system undergoes a transition to the coherent state.

Specifically, there are three typical strategies for changing contrarians into conformists: (1) a fraction p of contrarians is randomly chosen and then flipped into conformists; (2) contrarians are ranked according to the absolute value of their natural frequencies ($|\omega_i|$). Then, they are flipped into conformists from the largest $|\omega_i|$ to a threshold ω_0 , i.e. the coupling strength of the i th oscillator will be $\lambda_i = \lambda_2$ if $|\omega_i| > \omega_0$ and $\lambda_i = \lambda_1$ otherwise. Therefore $1 - p = \int_{-\omega_0}^{\omega_0} g(\omega)d\omega$; (3) the opposite of the process described in (2), i.e. $\lambda_i = \lambda_2$ if $|\omega_i| < \omega_0$ and $\lambda_i = \lambda_1$ otherwise, which leads to $p = \int_{-\omega_0}^{\omega_0} g(\omega)d\omega$. Mathematically, the three strategies correspond to three different correlations between the coupling strength and the natural frequencies as follows:

$$\text{Case 1: } \Gamma_1(\lambda) = (1 - p) \delta(\lambda - \lambda_1) + p \delta(\lambda - \lambda_2), \quad (79)$$

$$\text{Case 2: } \Gamma_2(\omega, \lambda) = H(\omega_0 - |\omega|) \delta(\lambda - \lambda_1) + H(|\omega| - \omega_0) \delta(\lambda - \lambda_2), \quad (80)$$

$$\text{Case 3: } \Gamma_3(\omega, \lambda) = H(|\omega| - \omega_0) \delta(\lambda - \lambda_1) + H(\omega_0 - |\omega|) \delta(\lambda - \lambda_2), \quad (81)$$

where $H()$ is the Heaviside function.

Ref. [391] considers Case 1, and shows that the system has two types of coherent states: the π state and the traveling wave (TW) state (see Fig. 51). In the π state, two clusters of conformists and contrarians keep a constant phase difference π . The π state is stationary, and the order parameter is a fixed point in phase space. For the TW state, the phase distribution spontaneously travels at a constant speed along the phase axis, always maintaining a constant separation $\delta \neq \pi$. As a result, the order parameter traces a circle around the origin at a constant angular velocity.

As the control parameter p varies, transitions are observed among the incoherent and the coherent states. The results can be split into two scenarios, depending on whether the conformists or the contrarians are more strongly affected by the mean field. Fig. 52 reports the phase diagram when the contrarians are more strongly affected by the mean field, i.e. $Q = -\lambda_1/\lambda_2 > 1$. For small p , the system is dominated by contrarians, and therefore is in the incoherent state. Once p exceeds a certain threshold, there are enough conformists for a consensus to emerge. At that point, the system jumps discontinuously to the π state, where it is polarized into two groups (see Fig. 51(a)). Inversely, starting from the coherent

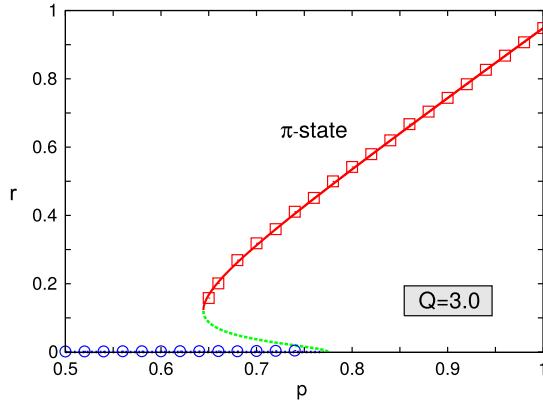


Fig. 52. (Color online). Order parameter r vs. p for $Q = 3$, computed by numerical integration (symbols) and compared to the theoretical prediction (lines). The solid lines denote the stable states; the dotted line individuates the unstable π state. $N = 25,600$, $\gamma = 0.05$, $\lambda_1 = -3.0$, $\lambda_2 = 1.0$.
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state and adiabatically decreasing p yields a discontinuous and hysteretic return to the incoherent state. On the other hand, if $Q < 1$, the transitions become continuous. Typically, as p increases from 0, the system goes from the incoherent to the π state, then to the TW state, and finally settles on the π state (see Fig. 3 of Ref. [391]).

Based on the Ott–Antonsen ansatz [241], the dynamics of Eq. (78) can be effectively reduced into a low-dimensional system describing the complex order parameters for the contrarians and conformists. By means of the analysis of the reduced system, the three steady states and their bifurcation points can be analytically obtained [391].

In Ref. [392], numerical simulations of Case 2 and Case 3 are presented, and both the π state and the TW state are found. In a certain parameter regime of Case 3, furthermore, a special explosive behavior is taking place: when the control parameter ω_0 is decreased to a critical threshold, the system bifurcates from the incoherent state into the TW state via a first-order like transition, while starting from the TW state, if ω_0 increases, the order parameter exhibits a continuous and hysteretic return to the incoherent state (see Fig. 7(b) in Ref. [392]).

Recently, Qiu et al. provided a detailed and complete analytical treatment for model (78), based on linear stability analysis and mean-field analysis [393]. It has been proved that in this model the incoherent state is neutrally stable below the synchronization threshold, and the critical points for synchronization are derived for all the three cases. Moreover, all possible stationary coherent states in the model are predicted, including the π state and the two types of TW states.

In Fig. 53, a detailed phase diagram characterizing the synchronization transitions in model (78) is reported. Depending on the correlations between the coupling strength and the natural frequencies (as well as on whether the conformists or contrarians are more strongly affected by the mean field), the system exhibits either first-order or second-order like transitions among different steady states. Besides the two examples of ES studied in Refs. [391] and [392] (which corresponds to Fig. 53(a) and (i), respectively) it is further revealed that a first-order transition can also occur in Case 2. As illustrated in Fig. 53(d), when p increases, the system experiences a first-order transition from the incoherent state into the coherent state, i.e., the π state. Inversely, when the system starts from the coherent π state, as p decreases, the system does not directly go back to the incoherent state. The π state first bifurcates into a new coherent state named the Bellerophon state (see next Section for the full characterization of such a novel state) via a second-order transition. Then, as p further decreases, the Bellerophon state loses its stability and the system jumps into the incoherent state via a first-order transition.

4.4. The Bellerophon state

In the previous Sections, a variety of coherent states emerging in the classical Kuramoto model have been discussed and characterized, such as the partially coherent state [237], the standing-wave [243,335] and TW [243,394] states, and the chimera state [217,219,395,396]. In such coherent states, oscillators inside each one of them display typically frequency-locked patterns, and therefore behave like a giant oscillator as a whole. Here, instead, we describe the main properties of a recently discovered novel state: the Bellerophon state.

4.4.1. The Bellerophon state in frequency-weighted Kuramoto models

In Ref. [397], ES was studied in a frequency-weighted Kuramoto model

$$\dot{\theta}_i = \omega_i + \frac{\lambda|\omega_i|}{N} \sum_{j=1}^N \sin(\theta_j - \theta_i), \quad i = 1, \dots, N, \quad (82)$$

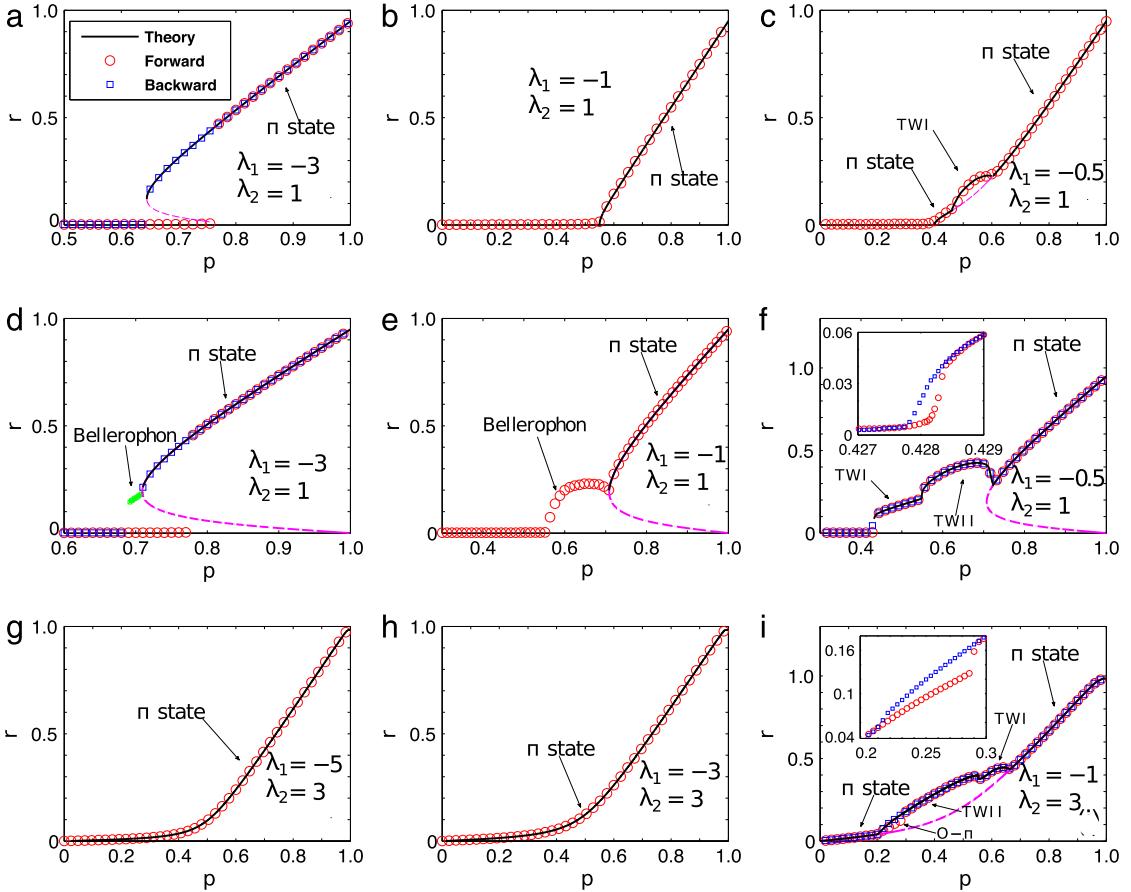


Fig. 53. Scenarios of synchronization in Eq. (78) as the proportion of conformists increases. $N = 50,000$ and $\gamma = 0.05$. From top to bottom, three rows correspond to Case 1, Case 2 and Case 3, respectively. From left to right, the three columns correspond to the case of $|\lambda_1| > \lambda_2$, $|\lambda_1| = \lambda_2$, and $|\lambda_1| < \lambda_2$, respectively. Both the forward (red circles) and the backward (blue squares) transitions are reported as obtained in an adiabatic way, while the black (pinkish red) curves correspond to the theoretical predictions of the stable (unstable) stationary coherent states, including the π state, the TW-I state, and the TW-II states. The O- π state denotes the oscillating π state. The numerical results are perfectly consistent with the theoretical predictions. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

Source: Adapted from Ref. [393].

with a bimodal Lorentzian distribution

$$g(\omega) = \frac{\Delta}{2\pi} \left[\frac{1}{(\omega - \omega_0)^2 + \Delta^2} + \frac{1}{(\omega + \omega_0)^2 + \Delta^2} \right]. \quad (83)$$

Here, θ_i and ω_i are the instantaneous phase and the natural frequency of the i th oscillator, respectively, and λ is the coupling strength. Δ is the width parameter (half width at half maximum) of each Lorentzian distribution, and $\pm\omega_0$ are their central frequencies. As shown in Fig. 54, system (82) sustains both a first- and a second-order transition to synchronization. According to the theoretical analysis [373,397], the critical point for the backward transition λ_c^b is always 2, while the critical point for the forward transition is

$$\lambda_c^f = \frac{4}{\sqrt{1 + (\omega_0/\Delta)^2}}. \quad (84)$$

As the parameter ω_0/Δ increases, the characteristic hysteresis width $\lambda_c^f - \lambda_c^b$ shrinks, which induces eventually the conversion of a continuous transition to synchronization into an ES.

A new coherent state is unveiled in system (82), which is actually non-stationary and essentially different from all the coherent states studied before. In such a state, oscillators form quantized, time-dependent clusters, where neither their phases nor their instantaneous frequencies are locked. Their instantaneous speeds are different within the clusters, but they form a characteristic cusped pattern and, more importantly, they behave periodically in time so that their average values are the same. Given its intrinsic specular nature with respect to chimera states, the new state has been named the Bellerophon state in Ref. [397].

In Fig. 55 four typical steady states are illustrated, corresponding to the λ values denoted by letters A, B, C, and D in Fig. 54(f). Among them, Fig. 55(a)–(c) are Bellerophon states while Fig. 55(d) is the fully coherent state. In Bellerophon

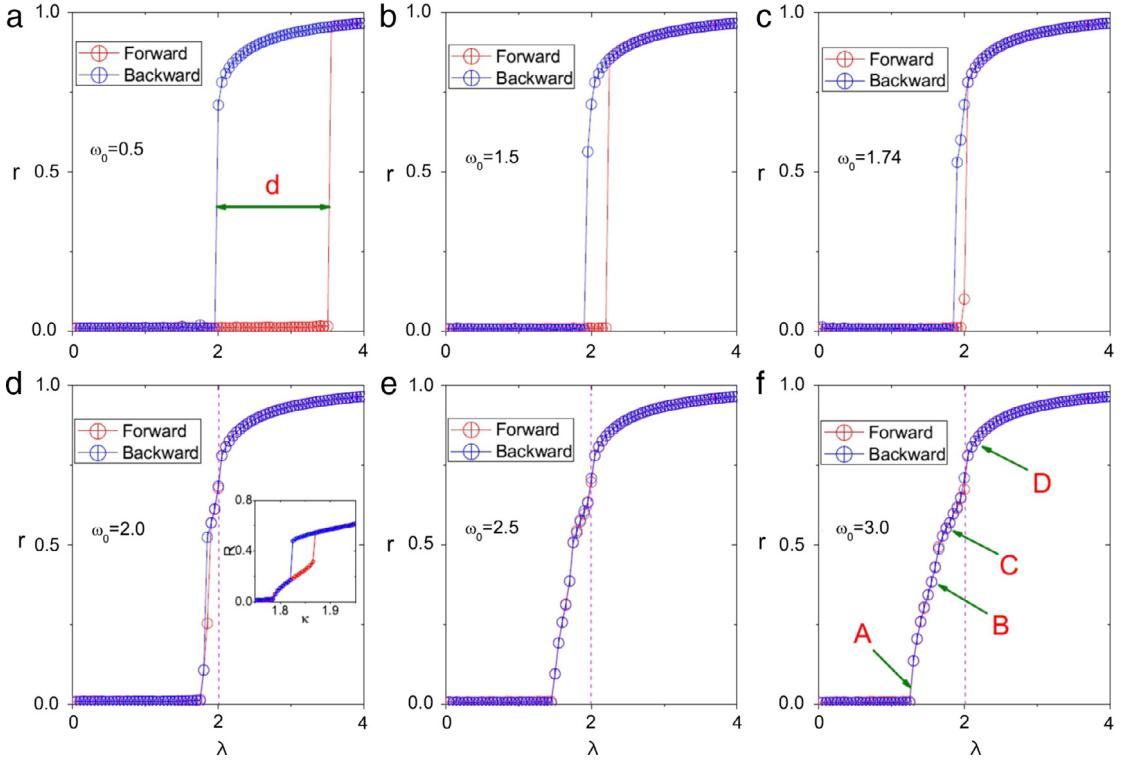


Fig. 54. (Color online). From explosive to continuous transition. r vs. λ for model (82). $N = 10,000$ and $\Delta = 1$. $\omega_0 = 0.5$ (a), 1.5 (b), 1.74 (c), 2.0 (d), 2.5 (e), and 3.0 (f), respectively.
Source: Adapted from Ref. [397].

states, oscillators form quantized clusters in terms of the average frequencies. The crucial point here is that although the average speeds of oscillators inside each cluster are equal to each other, their instantaneous speeds are generally different and quite heterogeneous. Furthermore, the instantaneous speeds of oscillators in each cluster are correlated and form the characteristic cusped pattern [Fig. 55(a3), (b3), and (c3)] analogous to that featured by the average frequencies of the oscillators within the chimera state. Therefore, inside each cluster, oscillators are neither phase- nor frequency-locked in a common sense. Rather, they seem to correlate in a higher order collective way: in terms of average speed oscillators inside each cluster are locked, but in terms of instantaneous speed each oscillator evolves uniquely. In fact, very interesting collective motion of oscillators can be observed in the unit circle, as illustrated in Ref. [397] (and, more specifically, in the animated movies which form part of the Supplementary Material of that Reference).

Bellerophon states can occur in system (82) also for different frequency distributions. For instance, in Ref. [398], model (82) was investigated with an asymmetric Lorentzian distribution:

$$g(\omega) = \frac{\Delta}{\pi[(\omega - \omega_0)^2 + \Delta^2]}, \quad (85)$$

where $\Delta = 1$ is kept as a constant, ω_0 changes in order to shift the frequency distribution along the positive axis. It is found that in the intermediate regime between incoherence and full synchronization [398], the Bellerophon states emerge, as illustrated in Fig. 56. Furthermore, the evolution of oscillators exhibits periodic intermittency following a synchronous pattern of bursting in short periods and resting in long periods [398].

In Ref. [399], model (82) was investigated with a uniform frequency distribution:

$$g(\omega) = \begin{cases} \frac{1}{2\gamma} & \text{for } |\omega| \leq \gamma, \\ 0 & \text{for } |\omega| > \gamma. \end{cases} \quad (86)$$

It is revealed that, as the coupling strength increases, the system undergoes two transitions. First, a first-order phase transition occurs from the incoherent state to a Bellerophon state, and then the Bellerophon state bifurcates into the two-cluster synchronous state via a continuous transition, when the coupling strength further increases.

4.4.2. The Bellerophon state in coupled conformist and contrarian oscillators

The Bellerophon state is also observed in the model of Eq. (78), which is essentially different with respect to the frequency-weighted model (82). Let us focus on a typical example, namely that of Fig. 53(d). As the control parameter p

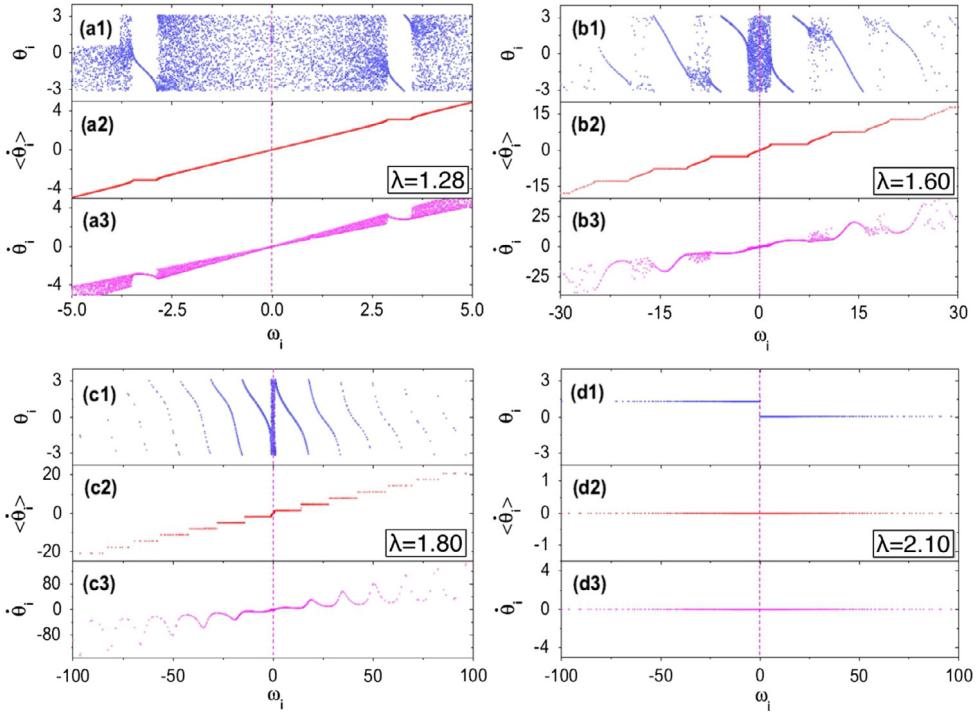


Fig. 55. (Color online). The Bellerophon states observed in the model (82) with bimodal frequency distribution. Snapshots of the instantaneous phase θ_i (upper plots), the average speed $\langle \theta_i \rangle$ (middle plots), and the instantaneous speed $\dot{\theta}_i$ (lower plots) vs. natural frequencies $\{\omega_i\}$ of the oscillators. $\lambda = 1.28$ (a), 1.60 (b), 1.80 (c), and 2.10 (d, the fully synchronized state). Panels (a)–(c) refer to Bellerophon states.
Source: Adapted from Ref. [397].

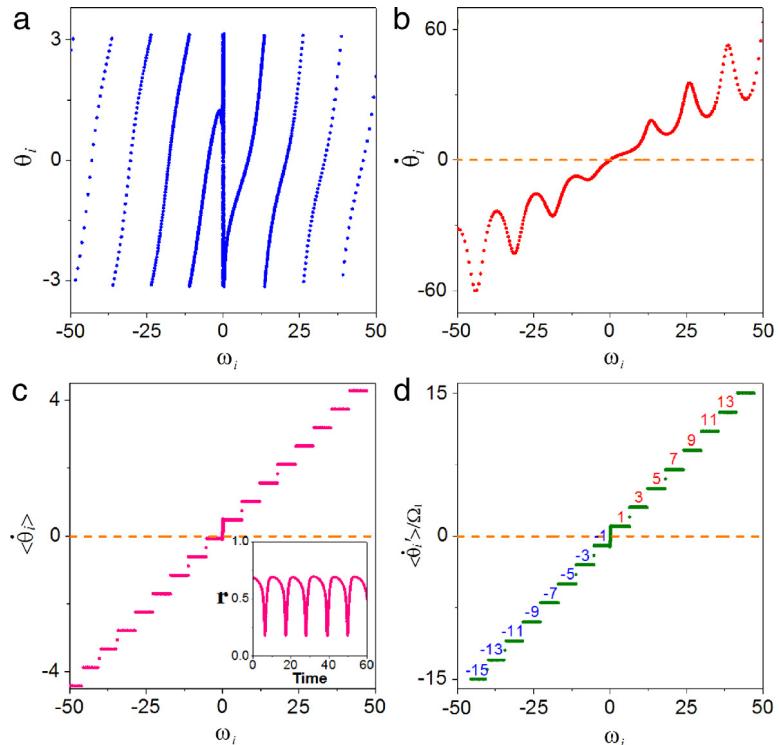


Fig. 56. (Color online). The Bellerophon states observed in the model (82) with asymmetric frequency distributions and $N = 10,000$, $\omega_0 = 0.5$, and $\lambda = 2.08$. (a)–(c) The phases (a), the instantaneous frequencies (speeds) (b), and the average frequencies (average speeds) (c) vs. the natural frequencies of oscillators. (d) The counterpart of (c) in a suitable rotating frame. Notice that the average frequencies are normalized by the fundamental frequency Ω_1 .
Source: Reprinted from Ref. [398].

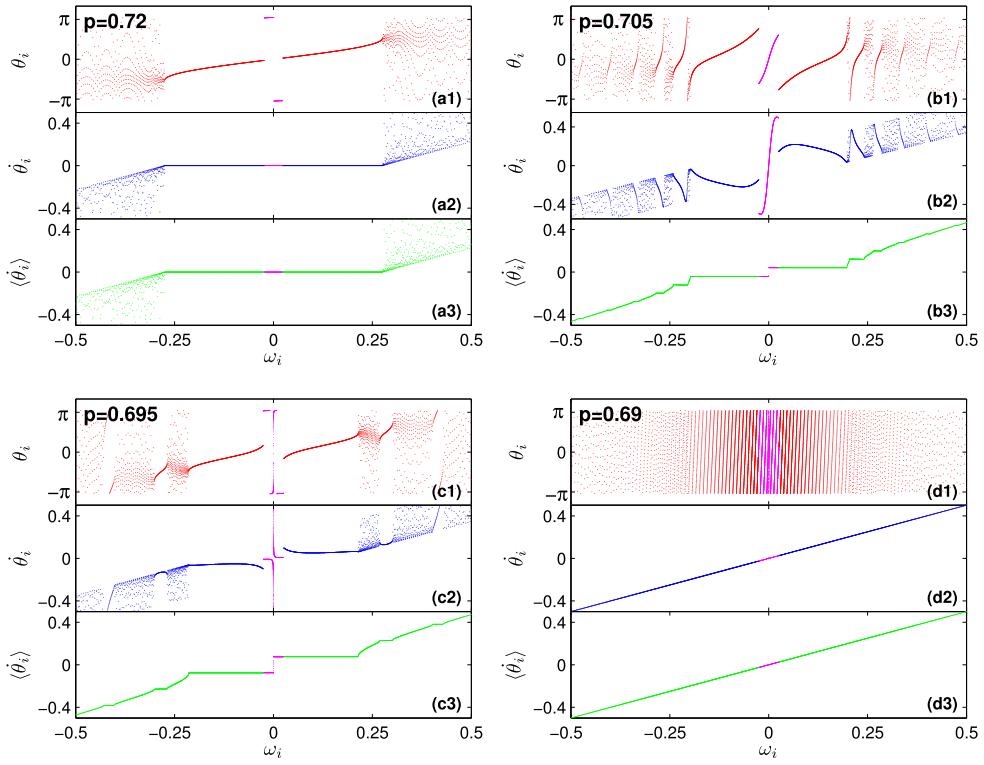


Fig. 57. (Color online). Typical coherent states and the incoherent state in Case 2, corresponding to the backward transition in Fig. 53(d). $N = 50,000$ and $\gamma = 0.05$. Snapshots of the instantaneous phase θ_i (upper plots), the instantaneous frequency (speed) $\dot{\theta}_i$ (middle plots), and the average frequency (average speed) $\langle \dot{\theta}_i \rangle$ (lower plots) vs. natural frequencies $\{\omega_i\}$ of the oscillators. (a) The π state with $p = 0.72$. (b)–(c) The Bellerophon states with $p = 0.705$ and $p = 0.695$, respectively. (d) The incoherent state with $p = 0.69$.

Source: Reprinted from Ref. [393].

increases, the system experiences a first-order like transition from the incoherent state to the coherent state, i.e. the π state. Inversely, when the system starts from the coherent π state, as p decreases, the system does not directly go back to the incoherent state. As shown in Fig. 53(d), the π state first bifurcates into the Bellerophon state via a second-order transition. Then, as p further decreases, the Bellerophon state loses its stability and the system finally jumps into the incoherent state via a first-order like transition. In Fig. 57, the typical Bellerophon states corresponding to Fig. 53(d) are characterized. Except for the fact that there are clusters of both conformists and contrarians, the observed Bellerophon states share all the other features of the frequency-weighted model described above.

4.5. Explosive synchronization in multi-layer networks

So far, our journey on explosive synchronization has been limited to the case of single-layer, or mono-layer, networks. Real complex systems, however, have far more complicated forms of interactions, which points to the fact that the hypothesis of a single-layer network may actually result in an overestimation (or an underestimation) of the problem under study. A more accurate analysis resorts on multi-layer networks, and dynamical processes on top of them are attracting more and more attention (see, for instance, Ref. [249] for a recent and rather comprehensive review). Therefore, examining ES on multi-layer networks became a topic of the utmost significance.

4.5.1. Multi-layer networks with the correlation property $\omega_i = k_i$

Jiang et al. recently proposed a Kuramoto model on multiplex networks [400], which are composed of one SF network α (as the base layer) and one community structured network β (as the external layer), and where each node constructs local links with probability $1 - \mu$ and long-range connections with probability μ (which is then the mixing parameter). Ref. [400] mainly focused on how the coupling between layers affects the phase transition of oscillators in the based layer α , for which the evolution equation reads as

$$\dot{\theta}_i^\alpha = \omega_i + \lambda_\alpha \sum_{j=1}^N A_{ij}^\alpha \sin(\theta_j^\alpha - \theta_i^\alpha) + \lambda_{\alpha\beta} \sin(\theta_i^\beta - \theta_i^\alpha), \quad (87)$$

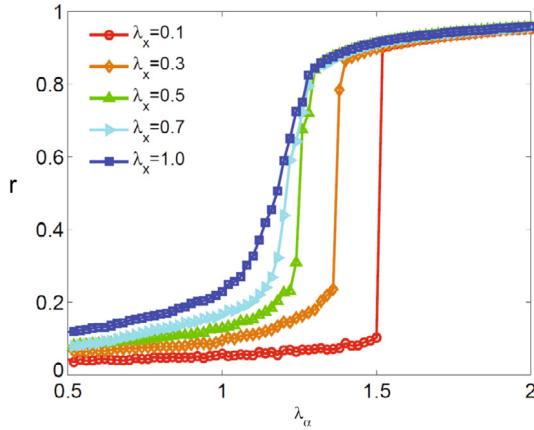


Fig. 58. (Color online). Synchronization level r of the base layer vs. intra-layer coupling strength λ_α for different values of inter-layer coupling strength λ_x . $N = 10^3$, and $\lambda_\beta = 1$.
Source: Reprinted from Ref. [400] with permission of Springer Ed.

where $\omega_i^\alpha = k_i^\alpha$. The first and second terms are the same of those referring to a single-layer network, and only involve intra-layer interactions in the base layer α . The third term accounts for the additional inter-layer interactions controlled by the coupling strength parameter $\lambda_{\alpha\beta}$. For simplicity, the hypothesis of symmetric coupling between both layers ($\lambda_{\alpha\beta} = \lambda_{\beta\alpha}$) is assumed, and $\lambda_{\alpha\beta}$ is denoted by λ_x .

The inter-layer coupling λ_x plays a significant role in determining the phase transition of the base layer. As shown in Fig. 58, at weak coupling strengths ($\lambda_x = 0.1$) ES emerges. With increasing λ_x , however, the transition becomes more and more continuous: there is an interim period at $\lambda_x = 0.5$, and a second-order-like transition occurs at $\lambda_x = 1.0$, which indicates that strong inter-layer couplings hinder the emergence of ES.

Moreover, it was unveiled that other externalities (mainly the parameters μ and λ_β) have great effect on ES. In details, increasing the mixing part μ puts off ES of the base layer, and requires larger critical coupling strengths λ_α^c , irrespectively of other parameters. This point is contrary to the findings in traditional modular networks [101,284,297], where long-range, inter-modular connections speed up the coherent state. As for coupling strength λ_β , its increase can accelerate ES of the base layer, though this effect seems to be very restricted.

4.5.2. Multi-layer networks with partial and weak correlation

A generative model of inter-dependent networks (including subnetworks G_1 and G_2) was proposed in Ref. [401]. At each time step, a node i enters into both subnetworks, and constructs m different links on each subnetwork. That is to say, nodes on both subnetworks have one-to-one interdependency with the same index. Then, like the preferential attachment rule in SF networks [402], the probability of producing one new link reads as $P_i^1 \sim \alpha_1 k_i^1 + (1 - \alpha_1) k_i^2$ and $P_i^2 \sim \alpha_2 k_i^1 + (1 - \alpha_2) k_i^2$, where α_1 and α_2 control the mutual correlation between networks. It is clear that the connectivity of each layer depends on itself and its counterpart, and both subnetworks have the same power-law degree distribution.

The Kuramoto model is then introduced. Instead of a strong correlation, the natural frequency of oscillator i in layer G_1 (G_2) is taken to be $\omega_i^1 = k_i^2$ ($\omega_i^2 = k_i^1$), which means that the dynamics of oscillator is controlled by the local topology and connectivity of the other layer. Such a weak hypothesis of correlation can induce ES when α_1 and α_2 are sufficiently large. As α_1 and α_2 decrease, the abrupt transition is replaced by a continuous transition.

4.5.3. Multi-layer networks in the absence of correlation

Zhang et al. [21] proposed to consider the following system

$$\dot{\theta}_i = \omega_i + \lambda \alpha_i \sum_{j=1}^N A_{ij} \sin(\theta_j - \theta_i), \quad (88)$$

where the new parameter α_i accounts for an adaptive control of nodes.

Initially, a fraction f of the nodes is randomly chosen, and $\alpha_i = r_i$ is assumed for each of the selected nodes, where r_i is the instantaneous local order parameter of the i th oscillator, defined as $r_i(t)e^{i\phi} = (1/k_i) \sum_{j=1}^{k_i} e^{i\theta_j}$, and ϕ denotes the phase averaged over the ensemble of neighbors. For the remaining fraction $1-f$ of the nodes, $\alpha_i = 1$. Obviously, $f = 0$ returns the traditional Kuramoto model [23,237,238,295], while $f > 0$ indicates that a fraction of nodes are adaptively controlled by the local order parameter. Besides, another significant point is the choice of the natural frequencies ω_i of oscillators, which are taken from a random homogeneous distribution $g(\omega_i)$ in the range $[-1, 1]$. Based on these assumptions, it was reported that an abrupt (with a hysteresis loop) transition takes place iff f is over a critical value f_c . Namely, the absence of correlation features can also promote the occurrence of ES in single-layer networks.

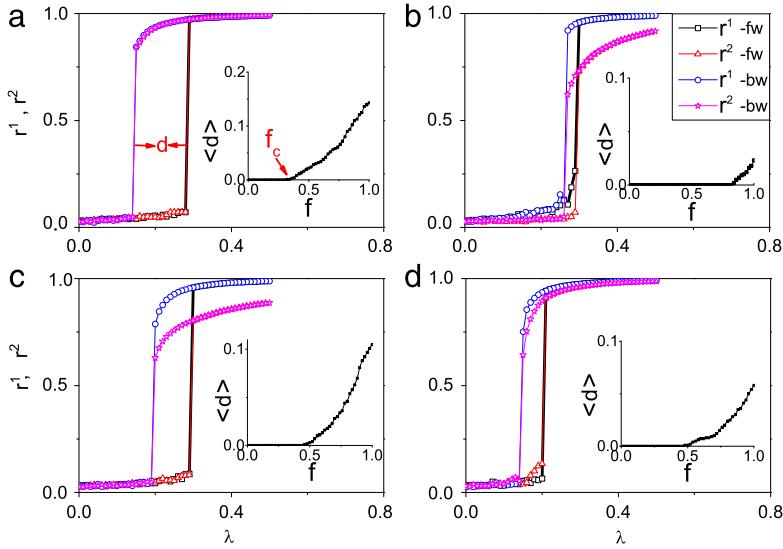


Fig. 59. (Color online). r^1 and r^2 vs. λ for two inter-dependent networks with $N = 10^3$ and $f = 1$. Squares and circles (triangles and stars) denote the forward (backward) transitions, and the insets report the average width (d) of the hysteresis loop as a function of f . Here layer 1 is a ER network with $\langle k \rangle = 12$. From (a) to (b), layer 2 is ER network with $\langle k \rangle = 12$ and $\langle k \rangle = 6$, and $g(w_i^2)$ is a random homogeneous distribution in the range $[-1, 1]$. From (c) to (d), layer 2 is ER network and SF network with $\langle k \rangle = 12$, and $g(w_i^2)$ is Lorentzian distribution and random homogeneous fashion.

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In order to explore the case of inter-dependent networks, Ref. [21] extended the analysis to the situation where each node has a one-to-one partner with the same index i . The equations become

$$\begin{aligned} \dot{\theta}_i^1 &= \omega_i^1 + \lambda \alpha_i^1 \sum_{j=1}^{k_i^1} A_{ij} \sin(\theta_j^1 - \theta_i^1), \\ \dot{\theta}_i^2 &= \omega_i^2 + \lambda \alpha_i^2 \sum_{j=1}^{k_i^2} A_{ij} \sin(\theta_j^2 - \theta_i^2), \end{aligned} \quad (89)$$

where the superscripts 1 and 2 denote the network layers 1 and 2, α_i^1 and α_i^2 involve the coupling of two layers via dependency links. In details, if node i falls into the fraction f , α_i^1 and α_i^2 refer to the local order parameters r_i^2 and r_i^1 , i.e. $\alpha_i^1 = r_i^2$ and $\alpha_i^2 = r_i^1$ ($\alpha_i^1 = \alpha_i^2 = 1$ otherwise), where $r_i^1(t)e^{i\phi^1} = (1/k_i^1) \sum_{j=1}^{k_i^1} e^{i\theta_j^1}$ and $r_i^2(t)e^{i\phi^2} = (1/k_i^2) \sum_{j=1}^{k_i^2} e^{i\theta_j^2}$. The frequency function $g(\omega_i^1)$ is still random homogeneous and in the range $[-1, 1]$, while $g(\omega_i^2)$ can change from a random to a Lorentzian distribution.

Fig. 59 shows the dependency of synchronization levels r^1 and r^2 on λ for different choices of the layer 2 and $g(\omega_i^2)$. ES occurs and seems very robust against the difference in topology and frequency distribution, which indicates that the correlation between oscillator's frequency and node's degree is not the essential condition for the emergence of ES. If f is below the critical value f_c , synchronization returns to a second-order transition. More importantly, these observations can be quantitatively verified via the mean-field theory. In the traditional second-order transition, oscillators with close frequency firstly collapse into small synchronization clusters, which gradually converge towards a giant cluster at the critical coupling strength. However, under ES, small clusters are prevented and all free oscillators are abruptly attracted to the giant cluster. Recently, furthermore, this kind of ES has been shown to coexist with the standard phase of the Kuramoto model in the thermodynamic limit [403].

5. Applications

5.1. Explosive percolation in real physical systems

Most of the studies on Explosive Percolation reviewed in Section 2 are explicitly conducted in the thermodynamic limit, i.e. under the hypothesis of an infinite system size (or a size of the order of the Avogadro's number, $N \sim 10^{23}$). However, real world systems are always finite-sized. For instance, the Internet, world-wide airline network and online social networks (just to mention three examples of very large sized real networks) have all values of N considerably smaller than 10^{18} [27]. In

this sense, it is interesting to reveal how the AP may induce significant discrete jumps in the realm of real-world networks, where EP is generally considered to be a phenomenon of cascading failures [94,106].

One of the major events of electrical blackout was affecting much of Italy on 28 September 2003: the shutdown of power stations directly led to the failure of nodes in the Internet communication network, which in turn caused further breakdown of power stations. This iterative process of cascading failures can also occur in a single power-grid network such as the blackout in Northwestern America in August 1996, and the blackout in Northeastern America and Canada in August 2003, where one overloaded node caused other nodes to also become overloaded and thus disabled the entire network. Other examples include the financial crises (in the network of global financial market) and the spreading of information and rumors through online social networks, such as Facebook or Twitter, where a failing node can cause its neighbors to fail as well. A common feature of these cascading failures is that they represent an inverse process of EP, which has a devastating effect on the network stability, and may cause huge economic losses.

Although a general answer to the question of how to suppress or control a cascading failure in real systems is still missing, some progress in this direction has been achieved [10,45,50,58,85,113,404], illustrating how the AP rule can actually be used in real systems.

For example, Rozenfeld et al. considered EP in the human protein homology network (H-PHN) and showed that the emergence of a spanning cluster exhibits similar features to a AP [50]. Their results indicate that the evolutionary-based processes shaping the topology of the H-PHN through duplication–divergence events may occur in sudden steps, similarly to what is seen in first-order like phase transitions.

Cho et al. considered the diffusion-limited cluster aggregation (DLCA) model of the sol–gel transition, and found that a discontinuous percolation transition can be observed when particles are Brownian, in which cluster velocity depends on cluster size as $v_s \sim s^\eta$ with $\eta = -0.5$ [10,85].

Pan et al. considered social networks community structures and assortativity (such as a mobile phone call network and a large arXiv coauthorship network), and found that the percolation transition depends on the structural properties of the network [58].

Oliveira et al. discovered that the electric breakdown due to pollution with metallic powder can become explosive if the inhibition of adsorption due to a local electric field becomes too strong [113].

Later on, Radicchi considered the case of coupled networks and found that the dependency links have an important role in the phase transition [404]. For a high density of dependency links, the network disintegrates in a form of a first-order phase transition, whereas for a low density of dependency links, the network disintegrates in a second-order transition. In the following, we briefly linger in each one of these cases.

5.1.1. The human protein homology network

First, we discuss the case of H-PHN. Based on real data,¹ Rozenfeld et al. found that H-PHN is composed of highly connected clusters of homologous nodes, while links between nodes of low homology generate inter-cluster connections, which is similar to the presence of strong links within communities and weak links between communities [50]. Further, they found that H-PHN is a weighted modular network, with the weights denoting the degree of similarity (homology) between two proteins. Ref. [50] claimed that such a weighted modular network can be explained on the basis of a product rule (PR) model. The idea is the following: suppose that H-PHN is grown up from a spanning skeleton of the network which connects all the different network areas. A new protein has a much larger weight to generate links within a module rather than with proteins that are further away. These dense modules are then connected with each other through weaker links. In the terminology of the PR model, this corresponds to an increase probability of connections between small clusters, if compared to the growth of already large clusters.

A quantity $0 \leq \zeta \leq 1$ is introduced to detect the degree of similarity between two optimally aligned proteins [50]. $\zeta = 1$ indicates a perfect alignment between the two proteins, or in other words, a short genetic distance between the proteins. The value of ζ for a protein pair in the network is taken to be the weight of the corresponding link. Then, the concrete process to construct the H-PHN can be implemented by considering an empty network of all proteins, and adding one link at a time in decreasing order of the weight, quantified by ζ . The process initially leads to well-connected families of highly homologous proteins, that are inter-connected at later stages by links with smaller values of ζ , resembling the requirements for EP in Ref. [9].

The results are reported (black line with “circles”) in Fig. 60. S/N represents the fraction of nodes in the largest cluster of the H-PHN, and p denotes the percentage of links. One clearly sees that the largest cluster in the network remains small until p reaches 0.8, and then quickly increases for $p > 0.8$. The behavior is very similar to an AP, i.e. links are added locally and without a significantly large spanning cluster before p_c , while the subsequent addition of links after p_c causes the small clusters to merge into a spanning entity. For comparison, the red line with “diamonds” in the same figure shows the results of a pure AP, for the same number of nodes. One sees that the two curves have a high-degree similarity, and the PR model has a sharper transition than the H-PHN.

For a deep understanding of the nature of the transition, Rozenfeld et al. reported the results for the cases of adding links in increasing order of ζ (green curve with “squares” in Fig. 60) and adding links randomly (the blue curve with “triangles” in

¹ The H-PHN was obtained from the Similarity Matrix of Proteins (SIMAP) project, <http://boinc.bio.wzw.tum.de/boincsimap>.

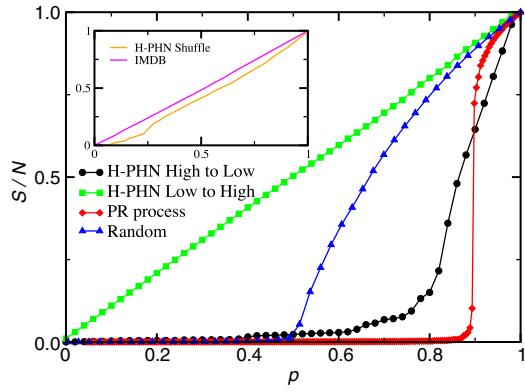


Fig. 60. (Color online). Fraction of nodes in the largest cluster of the H-PHN vs. the percentage p added links according to different rules: (a) links are added in decreasing weight order (black circles), (b) links are added in increasing weight order (green squares), (c) links are chosen according to the PR model (red diamonds), and (d) links are added randomly (blue triangles). Inset: same quantity as in the main figure but (a) the structure of the H-PHN remains the same and link weights are randomly redistributed (orange line), and (b) a weighted IMDB co-acting network is considered (violet line). (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)
Source: Reprinted from Ref. [50] with permission of Springer Ed.

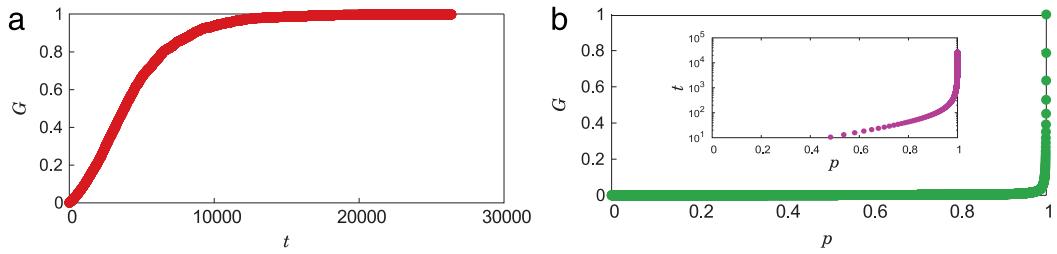


Fig. 61. (Color online). (a) The giant cluster size G vs. t for Brownian particles. Notice that here the giant cluster grows continuously from $t = 0$. (b) G vs. p . Here G grows drastically near $p_f = 1 - 1/N$. Inset: plot of the relationship between t and p . t increases drastically, but in a power-law manner as p approaches p_f . Simulations are carried out with $N = 8000$ mono-particles at $t = 0$ on a 400×400 square lattice.
Source: Reprinted with permission from Ref. [10].

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Fig. 60. The former strategy induces a percolation transition at a very early stage, and the largest cluster S increases linearly with p , i.e. almost every new link attaches to the largest cluster. In the latter strategy, the largest component S remains small before its critical point p_c , and then increases approximately linearly after p_c , i.e. following a second-order transition. By comparison of all the four curves in Fig. 60 one concludes that the case of decreasing weight in the H-PHN is much steeper than the case of random addition, and is much closer to the case of PR processes, indicating that the evolution of a modular network can be explained through the idea of EP.

The sharp transition of the H-PHN is not a universal property ascribed to weighted networks. The inset of Fig. 60 shows that, when the link weights of the H-PHN are shuffled without modifying the structure of the network, a smooth percolation transition is observed at an early stage. Similar growth patterns can be observed in other modular networks, such as the network of movie actors from IMDB in which two actors are connected if they co-acted in a movie (see the inset of Fig. 60).

5.1.2. The DLCA model

As a second case, we briefly discuss the case of the DLCA model. Initially, N single particles are placed randomly in a $L \times L$ square lattice. Assume that the particles are Brownian and the cluster velocity depends on the cluster size as $v_s \sim s^\eta$. For convenience of numerical simulations, one first lets the velocity of a cluster be inversely proportional to the square root of its size. The simulations are performed as follows: at each time step, (i) a s -sized cluster is selected with the probability $q \equiv s^{0.5}/(\sum_s N_s s^{0.5})$, and is moved to the nearest neighbor, where N_s is the number of s -sized clusters. When two distinct clusters are placed at the nearest-neighbor positions, these clusters are regarded as being merged, forming a larger one. (ii) The time is advanced by $\delta t = 1/(\sum_s N_s s^{0.5})$. Steps (i) and (ii) are repeated until the network percolates.

Let p be the number of cluster aggregation events per total particle number. Whenever two clusters merge, p is increased by $\delta p = 1/N$. Since $N - 1$ aggregation events occur during all aggregation processes, the aggregation event stops at

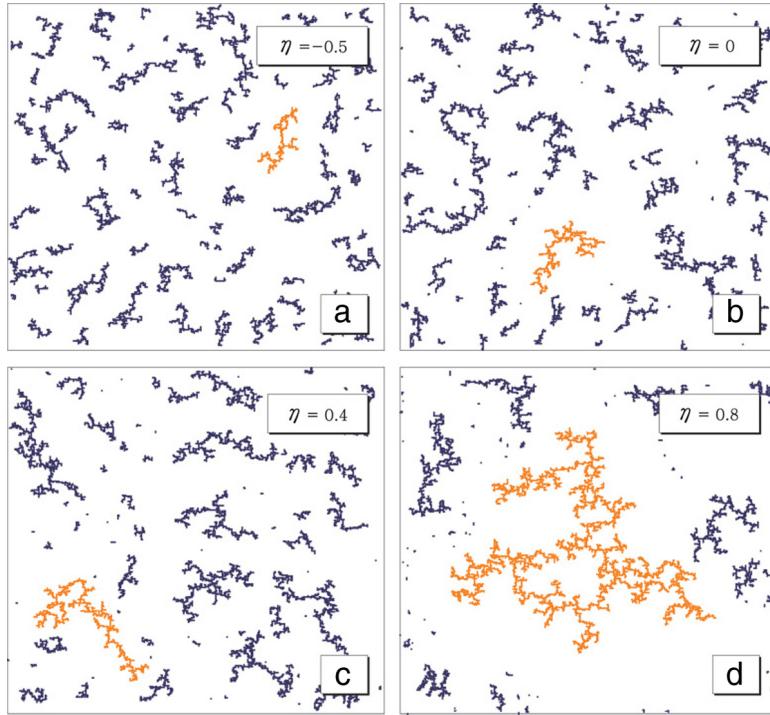


Fig. 62. (Color online). Snapshots of the system for various values of η at $p = 0.99$. The velocity of each cluster is given as $v_s \sim s^\eta$, where s is the cluster size. Numerical simulations are carried out for $N = 8000$ particles on a $L \times L = 400 \times 400$ square lattice. Since p is fixed, the number of clusters for each case is equal. The giant cluster is represented in a different color (gray/orange). The cluster-size distribution becomes more heterogeneous as η increases. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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$p_f = 1 - \delta p$ [10]. Fig. 61(b) shows the dependence of the giant cluster size G on p . One sees that it increases drastically, exhibiting a discontinuous phase transition. For comparison, the dependence of G on t is shown for the original DLCA model in Fig. 61(a), where one sees that G increases monotonically with t . The different behavior between Fig. 61(a) and (b) is caused by the nonlinear relationship between t and p , as shown in the inset of Fig. 61(b). When p is small, t increases almost linearly with p . At variance, as p approaches p_f , t increases drastically in a power-law manner, resulting in a discontinuous phase transition.

Let us now move to the general case of $v_s \sim s^\eta$. Here, a cluster is picked up with a probability proportional to s^η while the other rules in the numerical simulations are held. When clusters merge, p is increased by $\delta p = 1/N$. The time is given by $\delta t = 1/(\sum_s N_s s^\eta)$. Intuitively, when η is small or negative, fewer large-sized clusters are selected, so that their growth is suppressed. Medium-sized clusters are abundant even close to p_f , and they merge suddenly, resulting in a discontinuous transition. In contrast, when η is positively large, larger-sized clusters are selected and they have larger chances of colliding with other ones, merging into a bigger cluster. The result is that these latter clusters can grow faster than smaller ones, so the giant cluster grows continuously and the transition becomes continuous. The presence of a tri-critical point η_c is expected across which the transition type is changed. Fig. 62 shows snapshots of the system for different values of η at $p = 0.99$, supporting the above arguments.

The DLCA model was later extended to a reaction-limited cluster aggregation model in which clusters diffuse following Brownian motion, and when two clusters come into contact with each other, they merge with a certain probability r and remain separated with the remaining probability $1 - r$ [85]. It is found that a discontinuous transition can be observed in two and three dimensions, while the transition remains continuous in four dimensions.

5.1.3. Social networks

Next, we linger on the case of social networks [58], such as the empirical mobile phone call (MPC) network, and a large arXiv coauthorship (CA) network, where the nodes represent people and links denote their interactions. These social networks share common features such as community structures and assortativity.

The MPC data consist of 325×106 voice calls over a period of 120 days. An aggregated undirected weighted network can be constructed, where edges represent bidirectional calls between users and weights represent the total number of calls. The largest connected component (LCC) is then extracted, with 4.6×10^6 nodes and 9.1×10^6 edges.

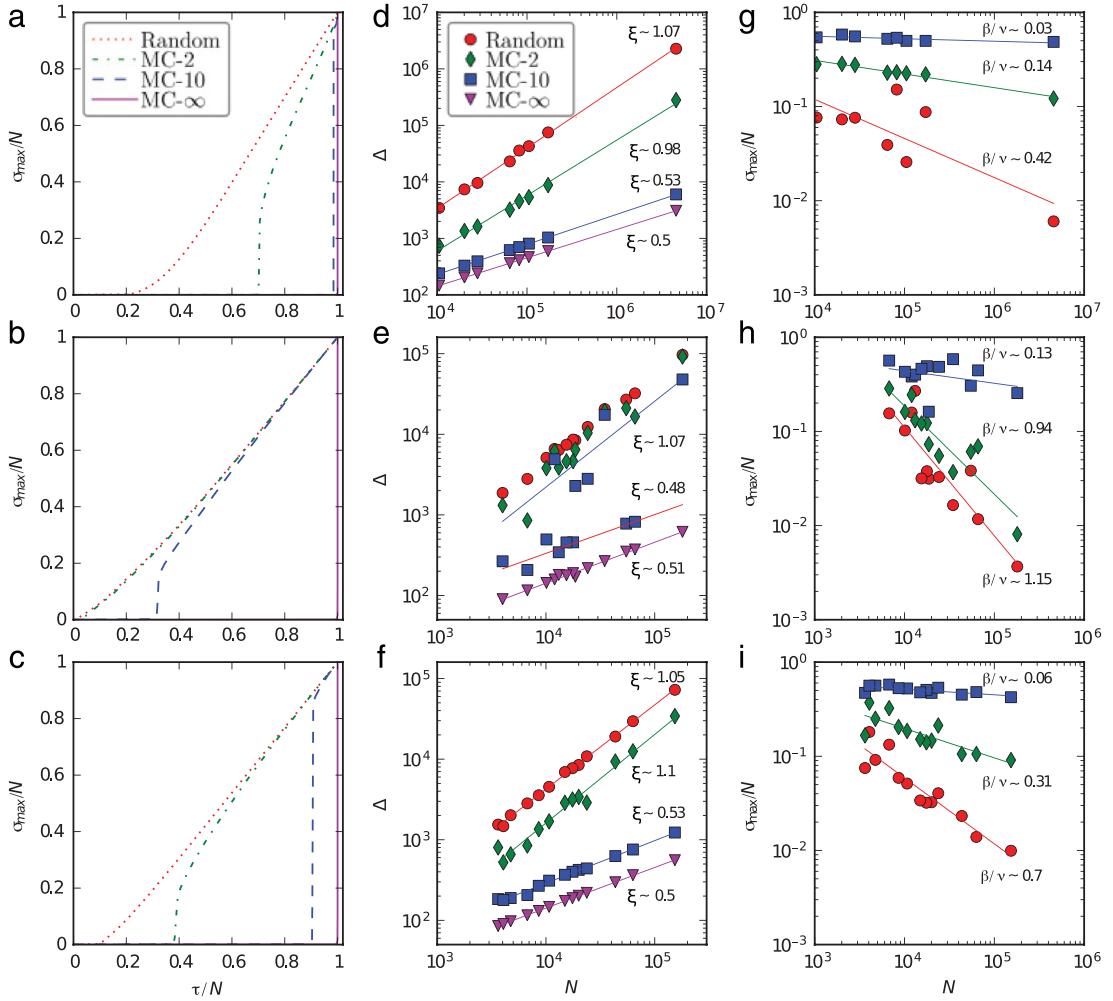


Fig. 63. (Color online). Variation of the relative size of the giant component (s_{max}/N) vs. the scaled number of inter-cluster edges τ/N for the (a) MPC, (b) CA, and (c) small CA networks. The corresponding variations in the gap, $\Delta \equiv \tau(N/2) - \tau(\sqrt{N})$, as a function of system sizes are shown in (d)–(f), for the Random, MC-2, MC-10, and MC- ∞ rules (see legend for color codes). Solid lines indicate fitted scaling exponents ξ . The variation of the order parameter (s_{max}/N) as a function of the system size N is shown for (g) MPC, (h) CA, and (i) SCA networks. For each system the order parameter is calculated at the critical point. The solid line indicates the best fit obtained with a power-law scale with exponent β/v . All curves are averaged over 10^3 runs. Source: Adapted with permission from Ref. [58]. Courtesy of M. Kivelä.

The collaboration data is from the arXiv² and contains all e-prints in “physics” until March 2010. There are 4.8×10^5 article headers, from which one extracts the authors. In the CA network two authors are connected if they signed together articles whose number determines the link weight. The LCC is extracted, with 1.8×10^5 nodes and 9.1×10^6 edges. In addition, a filtered version of the CA can be constructed, where articles with more than 10 authors ($\sim 2\%$ of the total number) are ignored. This is to remove the very large cliques from papers with $\sim 10^3$ authors in fields such as hep-ex or astro-ph, where the principles behind collaboration network formation appear different. The LCC of the resulting smaller CA network has 1.5×10^5 nodes and 9.1×10^5 edges. Note that, although the number of nodes is not much smaller than for that of CA, the number of edges is an order of magnitude smaller.

For describing the percolation process, Pan et al. used the min-cluster (MC- m) SR, with different values of m [58]. Initially, all the edges of the empirical network are considered unoccupied. Then, at each time step, m unoccupied edges are drawn at random. Out of these edges, the one minimizing the size of the component formed if it were occupied is chosen. Intra-component edges are always favored against inter-component edges, as they do not increase the size of any cluster. When comparing two inter-component edges, the one for which the sum of the connected cluster sizes is minimized is selected. As intra-cluster edges do not affect cluster growth, one considers only the number of inter-cluster edges τ . In particular, three variants of the MC rule (MC-2, MC-10, and MC- ∞) as well as random link percolation are compared [58]. Fig. 63(a)–(c) show the variation of the fraction $s_{max}(\tau)/N$ against the scaled number of inter-component edges, τ/N . For all the three

² <http://arxiv.org>.

networks, the transition of the order parameter is smooth for the random case, while for the extreme case, MC- ∞ , the transition appears abrupt. For MC-2 and MC-10, the situation is more complicated, and in the following will be described in more detail.

One can study the dependence of the width of the transition window $\Delta \equiv \tau(N/2) - \tau(\sqrt{N})$ on the system size, where $\tau(N/2)$ and $\tau(\sqrt{N})$ are the lowest values of τ for which $s_{max} > N/2$ and $s_{max} > \sqrt{N}$, respectively [9]. In general, the width scales as a power law with the system size ($\Delta \sim N^\xi$). For classical percolation, $\xi = 1$. It is argued that $\xi < 1$ for EP, and the rescaled width of the transition region ($\Delta/N \sim N^{\xi-1}$) vanishes in the limit of large N . In order to apply finite-size scaling to empirical networks, samples of different sizes are needed. As call networks are geographically embedded, one extracts sub-networks of users in chosen cities, based on postal codes of their subscriptions. For the CA networks, one extracts sub-networks of authors with articles in the same subject class. It turns out that, for all networks $\Delta\tau \sim N^\xi$, with $\xi \sim 1$ for random and $\xi \sim 0.5$ for the MC- ∞ case (see Fig. 63(d)–(f)). Thus, the exponent ξ clearly differentiates the explosive transition from random-link percolation. Furthermore, for all three networks, $\xi \sim 1$ for the MC-2, resembling an ordinary percolation transition. However, for MC-10, the scaling exponent behaves differently for the three networks. For the MPC and smaller CA networks, $\xi \sim 0.5$, indicating EP. For the CA collected data, points seem to indicate that there is no scaling. However, a closer inspection shows that they cluster around two straight lines with $\xi \sim 1$ and $\xi \sim 0.5$.

According to Ref. [34], the order parameter s_{max}/N has a finite-size scaling

$$\frac{s_{max}}{N} = N^{-\beta/v} F[(\tau - \tau_c)N^{1/v}], \quad (90)$$

where F is some universal function, τ is the control parameter, τ_c is the critical point of the transition, β is the critical exponent of the order parameter, and v is another critical exponent. One chooses the critical value τ_c of the control parameter as the value of τ where the susceptibility (that is, the average cluster size) has its maximum. For the MPC network [Fig. 63(g)], the scaling at τ_c of s_{max}/N yields a very small exponent $\beta/v \sim 0.03$ for the MC-10 case, while for MC-2 and random percolation, the exponents are larger ($\beta/v \sim 0.14$ and $\beta/v \sim 0.42$, respectively). The exponents for the SCA network behave similarly [Fig. 63(h)] with a low value $\beta/v \sim 0.06$ for the MC-10 case and relatively high values $\beta/v \sim 0.31$ and $\beta/v \sim 0.70$ for MC-2 and random percolation, respectively. In contrast, for the CA network, the exponents have higher values for all cases [Fig. 63(i)] ($\beta/v \sim 0.13$, $\beta/v \sim 0.94$, and $\beta/v \sim 1.15$) for MC-10, MC-2, and random percolation, respectively.

5.1.4. Inter-dependent networks

Recently, percolation in real systems was extended to the case of inter-dependent networks [404]. The functioning of a real network depends not only on the reliability of its own components, but also on the simultaneous operation of other real networks coupled with it. Percolation transitions in inter-dependent networks can be understood by decomposing these systems into uncoupled graphs: the intersection among the layers, and the remainders of the layers. When the intersection dominates the remainders, an inter-connected network undergoes a smooth percolation transition. Conversely, if the intersection is dominated by the contribution of the remainders, the transition becomes abrupt even in small networks.

As an example of real systems, Radicchi showed the transition for two inter-connected systems of interest in the biology: the *H. sapiens* protein-protein interaction network and the *C. elegans* connectome, finding that the inter-connected systems undergo smooth percolation transitions [404]. The reason is that these organisms have developed inter-connected networks sharing a core of “high quality” edges to prevent catastrophic failures. The same properties seem to characterize also the multi-layer air transportation network within the USA major airports, where the set of “high” quality’ edges that avoid truly catastrophic changes in the connectedness of the entire inter-dependent system are constituted.

5.2. Explosive synchronization in real physical systems

5.2.1. Power grids

Among the many real-world systems where ES can be observed, power grids are actually the ones in which cascading of failures have occurred numerous times.

One way to produce a cascade of failures is by the overload of lines: an outage of a line may lead to overload on other lines, thereby eventually leading to their outage. For instance, in August 2003, an undetected initial power line failure in Ohio caused a massive power blackout rolled across Northeastern USA and Canada, spanning eight states and two provinces, and affecting some 55 million people [405].

Another way is overloading the power stations: an outage of a node may lead to overload on other nodes, also leading to their outage. For example, a small initial power shutdown in El Paso (Texas) caused the August 1996 blackout in Northwestern America in which the power outage spread through six states between Oregon and California, leaving 7.5 million customers without electricity.

Around the world, the number of outages affecting large populations has steadily risen in the past decades, causing enormous economic losses.

The cascade of failures in power grids can be also explained in terms of their underneath inter-dependent network structure. For example, in September 2003, Italy was affected by a country-wide blackout: the shutdown of power stations

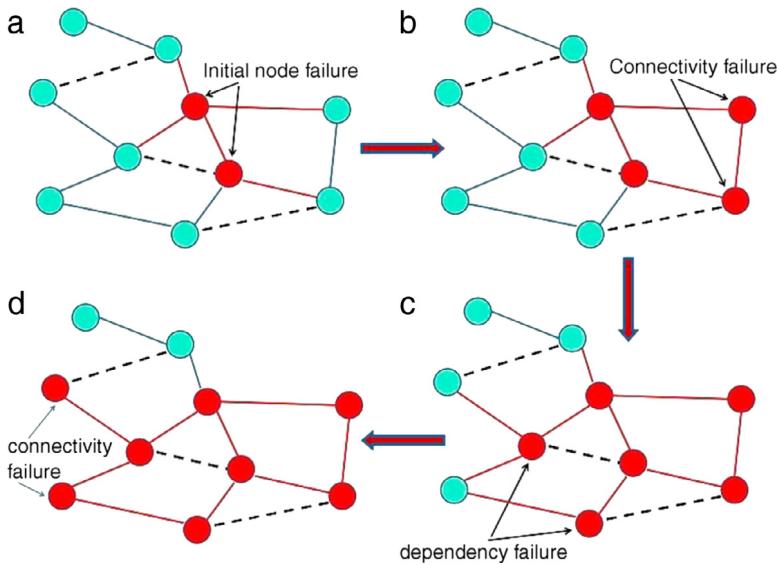


Fig. 64. (Color online). Synergy between the percolation process and the failures caused by dependency links. The network contains two types of links: connectivity links (solid lines) and dependency links (dashed lines). (a) The process starts with the initial failure of two nodes (marked in red). The connectivity links attached to them also fail (marked in red). (b) A percolation process takes place where all the nodes and the connectivity links attached to the failed elements and that are not connected to giant cluster (largest cluster) by connectivity links, also fail (marked in red). (c) The nodes connected by dependency links to the failed nodes also fail (marked in red). (d) Another step of connectivity failure in which two more nodes fail because they are not connected to the largest cluster (currently containing only two nodes). (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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directly led to the failure of nodes in the Internet communication network, which in turn caused further breakdown of power stations and left 57 million Italians in the dark [106].

Large-scale and/or long term failures have devastating effects on almost every aspect in modern life, and especially on inter-dependent systems (such as telecommunication, gas and water supply, and transportation systems). The iterative process of cascading failures can be described by the scheme of Fig. 64 where the outbreak in panel (d) is initiated by the node failure in panel (a) through a few intermediate steps [94].

A power grid is a network where links are power lines and nodes may be either generators (power plants) or consumers. The generators produce the electric energy which needs to travel long distances in high-voltage transmission lines (e.g. 400 kV) in order to arrive the consumers or loads [406–409]. The nodes may therefore consume, produce or distribute power, whereas the links transport power and may include passive elements with resistance, capacitance and inductance. A power grid is embedded in a 2D real space. Fig. 65 is a representation of the Northern European power grid with $N = 236$ nodes and $L = 320$ connections [410].

A fundamental requisite for a power grid to operate in a stable way is the maintenance of a phase-locked state within the entire system, thus synchronization is understood as the ability of a power grid to keep a global state where generators and loads (with intrinsic different frequencies) run with the same effective frequency (e.g. 50/60 Hz). If loads are too strong (or unevenly distributed) or if some major fault (or a lightning) occurs, a node may lose synchronization. In that situation, the synchronization landscape may change drastically and a blackout may occur.

Therefore, power grids are generally planned and operated to withstand the occurrence of certain disturbances. For this purpose, the following conditions of dynamic security are being guaranteed [407]: (1) when any of a specified set of disturbances occurs, the system survives the ensuing transient and moves into a steady-state condition; (2) no bus voltage magnitudes during transients are outside their permissible ranges; (3) in this new steady-state condition, no control devices, equipment or transmission lines are overloaded and no bus voltage magnitudes are outside their permissible ranges (say 5% of nominal).

Now, it happens that a power system is continuously experiencing disturbances, such as outages, short-circuits, sudden large load changes, or a combination of such events. Because of economic and environmental pressures, power grids tend to run at their full utilization, which increases the effects of these disturbances on their security. For example, when several voltages collapse, a power system must operate even closer to the limits of stability and then the risk of a major breakdown is increased [409].

Several models have been proposed to analyze the stability of power grids [239,406,408,411–414]. In these models, each node- i has the same equation of motion, and is characterized by a parameter P_i quantifying the generated ($P_i > 0$) or consumed ($P_i < 0$) power. The state of each node is determined by its phase angle $\phi_i(t)$ and velocity $\dot{\phi}_i(t)$. During regular operations, generators and consumers run with the same frequency $\Omega = 2\pi \times 50$ Hz or $\Omega = 2\pi \times 60$ Hz. The phase of

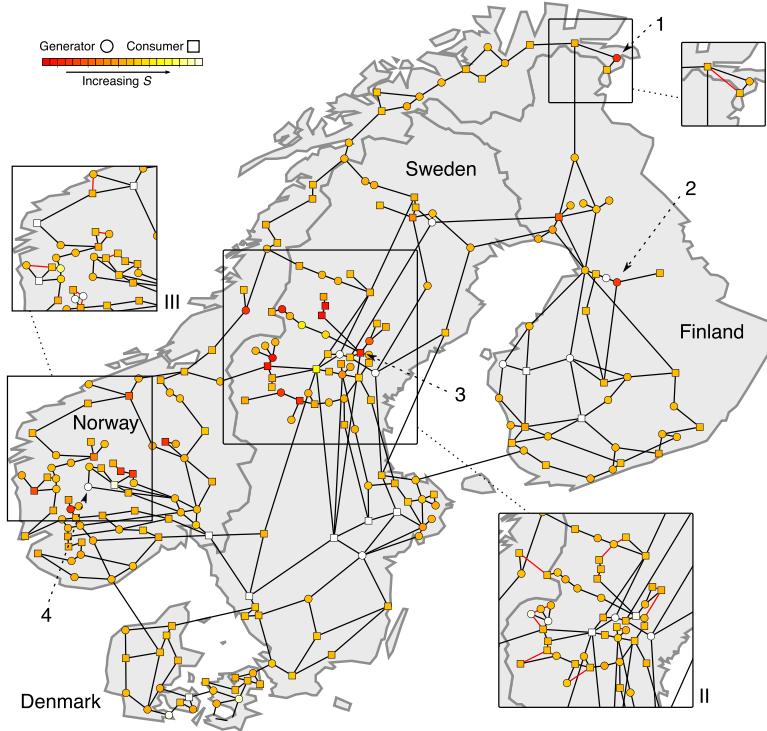


Fig. 65. (Color online) Schematic representation of the Northern European power grid. The grid has $N = 236$ nodes and $L = 320$ transmission lines. For simplicity, the load scenario was chosen randomly, with squares (circles) depicting $N/2$ net consumers (net generators). Source: Reprinted with permission from Macmillan Publishers Ltd: Ref. [410]. © 2014.

each node- i is then written as

$$\phi_i(t) = \Omega t + \theta_i(t), \quad (91)$$

where θ_i is a phase fluctuation.

For simplicity, one may assume that the nodes can be modeled as having the same dissipative coefficient α , the same moment of inertia, and that power is not lost during transmission. By energy conservation, each generated or consumed power P_i must be equal to the sum of the power taken from the grid (P_i^{trans}) plus the accumulated (P_i^{acc}) and dissipated (P_i^{diss}) power, i.e.

$$P_i = P_i^{\text{trans}} + P_i^{\text{acc}} + P_i^{\text{diss}}. \quad (92)$$

Now, during its rotation a turbine dissipates energy at a rate proportional to the square of its angular velocity [408]

$$P_i^{\text{diss}} = k(\dot{\phi}_i)^2, \quad (93)$$

and it accumulates kinetic energy $1/2I(d\phi_i/dt)^2$ at a rate

$$P_i^{\text{acc}} = \frac{1}{2}I \frac{d}{dt}(\dot{\phi}_i)^2, \quad (94)$$

where I is the moment of inertia. The condition for power transmission is that the two devices do not operate in phase, i.e. there is a phase difference between the two nodes connected by a line $\Delta\theta_{ij} = \theta_j - \theta_i$. The transmitted power is proportional to the sinus of the phase difference between the voltages of the nodes at the two end-points of the line [413], which gives

$$P_i^{\text{trans}} = \sum_j -P_{ij}^{\text{max}} \sin(\phi_j - \phi_i), \quad (95)$$

where P_{ij}^{max} is an upper bound for the transmission capacity of a line. Substituting Eqs. (93), (94) and (95) into Eq. (92) one has

$$I\Omega\ddot{\theta}_i = P_i - k\Omega^2 - 2k\Omega\dot{\theta}_i + \sum_j P_{ij}^{\text{max}} \sin(\theta_j - \theta_i), \quad (96)$$

which can be written into normalized units as

$$\ddot{\theta}_i = P'_i - \alpha \dot{\theta}_i + \sum_{j=1}^n W_{ij} \sin(\theta_j - \theta_i), \quad (97)$$

where P'_i , α and $W_{ij} = \lambda A_{ij}$ are the normalized parameters, and A_{ij} are the elements of the adjacency matrix which accounts for the heterogeneity of the network. P'_i is related to both generated and dissipated power at the node, while λ (the strength of the coupling) is related to the maximum transmitted power. Eq. (97) is the basic equation considered in power grids dynamics. Notice that, when the inertia I is negligible, one may ignore the term $\ddot{\theta}_i$ in Eq. (97), which then returns to the classical 1970 Kuramoto model.

In addition to load balance, synchronization between all elements of the grid is a crucial aspect to the system. Recently, ES in Eq. (97) has been extensively studied [415–418]. The stable operation of a power grid requires that all machines run at the same frequency Ω . The phases of the machines will generally be different, but the phase differences are constant in time. To study the influence of dynamics and structure on global synchronization, P'_i is generally assumed to satisfy a distribution with zero average. For example, P'_i can be explained as the natural frequency of node i , and assumed to be of the form [415,417,418]

$$P'_i = D(k_i - \langle k \rangle), \quad (98)$$

where k_i is the degree of node i , $\langle k \rangle$ the network average degree, and D a proportionality constant. That is, the natural frequency P'_i of a node i is assumed to be proportional to its degree.

The choice of Eq. (98) assumes that, in SF topologies, many nodes play the role of consumers (nodes with $k_i < \langle k \rangle$), while nodes with high degrees play the role of power producers (nodes with $k_i > \langle k \rangle$). In this framework, the relation $\sum_i P'_i = 0$ is satisfied, which means that the total consumed power is equivalent to the total generated power. In this latter case, all oscillators try to rotate independently at their own natural frequencies, while the coupling λ tends to synchronize them to a common phase. Besides the degree distribution $P(k)$, it is known that real world network can display either assortativity or disassortativity degree mixing behavior. When degree-degree correlations are accounted for, it is found that the synchronization diagrams have a strong dependence on the network assortativity, indicating that one is able to control the hysteretic behavior of the second-order Kuramoto model by tuning the network properties [415].

In numerical simulations, all the initial networks are constructed through the BA model with $\langle k \rangle = 6$ and $N = 10^3$. Eq. (97) is integrated and the Kuramoto order parameter $r(t)$ (Eq. (14)) is calculated after a transient process. Fig. 66 shows the forward and backward synchronization diagrams $r(\lambda)$ for different values of α within the interval $[0.2, 1]$. Panels (a)–(d) represent the cases of $\alpha = 0.2, 0.4, 0.6$ and 0.8 , respectively. It is easy to see that there is a hysteresis loop in all the four cases, indicating that ES is taking place in the network. From Fig. 66 we also see that the area of hysteresis and the critical coupling for the onset of synchronization in the forward branch (solid lines) tends to decrease as α is increased, which also contributes to increase the maximal value of the order parameter.

The network's degree-degree correlations can be quantified by the assortativity coefficient \mathcal{A} defined in Eq. (42). In order to inspect the influence of network's degree-degree correlation on ES in Eq. (97), one needs to change the assortativity of the network, which can be implemented as follows [415]: at each step, two edges are selected at random and the four nodes associated to the selected pair of edges are ordered from the lowest to the highest degree. Given a target assortative mixing ($\mathcal{A} > 0$), with probability p a new edge is formed between the first and second nodes and another one between the third and fourth nodes. If one of the two new edges already exists, the step is discarded and a new pair of edges is chosen. A similar heuristic mechanism can also generate disassortative networks ($\mathcal{A} < 0$). Fig. 67 shows the synchronization diagram $r(\lambda)$ for networks with different values of assortativity, where panels (a)–(d) represent the cases of $\mathcal{A} = -0.3, -0.1, 0.1$ and 0.3 , respectively. As \mathcal{A} increases, the hysteresis becomes smaller and smaller, and the critical coupling of the increasing branch is weakly affected.

The natural frequencies can also be chosen from different distributions. For instance, Olmi et al. considered the case of a Gaussian distribution $g(P') = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-P'^2/2\sigma^2}$ with zero average and unitary standard deviation σ [416]. It is observed that, in the hysteretic region, clusters of locked oscillators of various sizes and different levels of synchronization coexist.

It is important to remark that the process of ES in Eq. (97) is quite different from that of the Kuramoto model. It is revealed that in the case of uncorrelated SF networks, nodes join the synchronous component progressively, grouped into cluster of nodes with the same degree, starting from small degrees. The phenomenon was therefore called *cluster explosive synchronization* [417,418], and the difference with the case of the Kuramoto model is that in the latter the average frequency $\langle \omega \rangle_k$ jumps to $\langle k \rangle$ for all k at the same time. Fig. 68 reports the synchronization process and highlights the progressive presence of clustering.

5.2.2. Smart grids

An imperative condition for the reliable functioning of power-grids is that the generators remain synchronized. This condition is relatively easy to be attained in traditional power grids, where the power plants are mainly based on coal, gas, oil or nuclear power. However, in recent years, more and more distributed energy sources and self-sufficient micro-grids are included in the network, such as wind turbines, photovoltaic arrays, bio-gas power generators, and other renewable

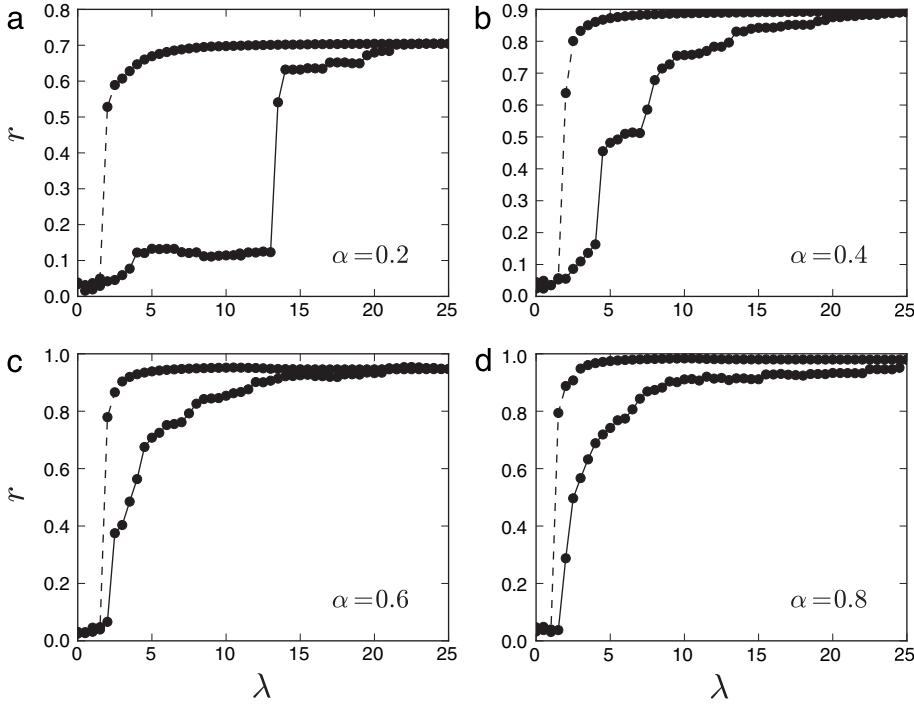


Fig. 66. Synchronization diagram $r(\lambda)$ for (a) $\alpha = 0.2$, (b) $\alpha = 0.4$, (c) $\alpha = 0.6$, and (d) $\alpha = 0.8$. The network has assortativity coefficient $A = -0.3$. With increasing α , the onset of synchronization and the hysteresis decrease. The natural frequency of each oscillator is $P'_i = k_i - \langle k \rangle$ and the networks have $N = 10^3$ and $\langle k \rangle = 6$. The degree distribution follows a power law $P(k) \sim k^{-\gamma}$, where $\gamma = 3$. Curves in which points are connected by solid (dashed) line correspond to the forward (backward) continuations of the coupling strength λ .

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energy sources. Such modern grids often consist of thousands of power substations and generators linked across thousands of kilometers, and are therefore much larger (and more complicated) than the traditional ones.

They are called *smart grids*, because they make use of information and communication technologies (ICTs) to collect information (such as the behaviors of suppliers, consumers, and prosumers) and to take actions in an automated fashion to improve the efficiency, reliability, economics, and sustainability of the production and distribution of electricity [406,412–414,419].

A key problem of smart grids is their stability. For instance, any person who has a solar panel at home, may become an energy trader, i.e. buy energy when the price is low, store it and then resell when the price is high, resulting in a new market within smart grids. As markets are known to be unstable, their influence to power grids is serious. Another element of instability is the fluctuation of renewable energy sources, as for instance wind turbines which depend significantly on the environment and season. All these factors introduce disturbances and fluctuations in production and demand, which may trigger desynchronization of power generators. Once it happens, the connected generators cannot be in pace and thus induce a cascading failure, which may be the reason behind the increasing number of power blackouts reported in recent years.

A characteristic feature of smart grids is their large fluctuations in energy. Notice that in the case of traditional power grids, we have a balance between the generators and the consumers, i.e. from Eq. (98) we have $\sum P'_i = 0$. However, in the case of smart grids, this balance may not hold, and thus one has

$$\sum P'_i > 0 \quad \text{or} \quad \sum P'_i < 0. \quad (99)$$

In the following, without loss of generality, we describe the case of $\sum P'_i > 0$.

The simplest non trivial grid is a two-element system consisting of one generator and one consumer [412]. In the phase-locked state, both derivatives $\dot{\theta}_i$ and $\ddot{\theta}_i$ are zero, and from Eq. (97) one has

$$0 = P'_i + \sum_{j=1}^2 W_{ij} \sin(\theta_j - \theta_i). \quad (100)$$

For the sum over all the equations gives,

$$0 = \sum_i P'_i + \sum_i W_{ij} \sin(\theta_j - \theta_i) + \sum_i W_{ij} \sin(\theta_i - \theta_j) = \sum_i P'_i. \quad (101)$$

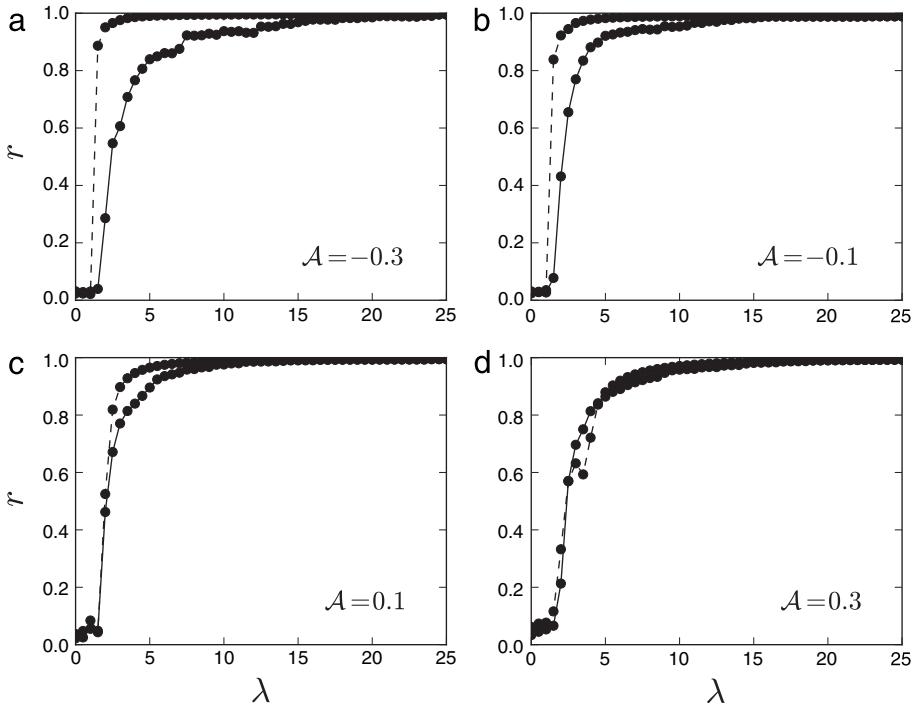


Fig. 67. Synchronization diagram $r(\lambda)$ for (a) $\mathcal{A} = -0.3$, (b) $\mathcal{A} = -0.1$, (c) $\mathcal{A} = 0.1$, and (d) $\mathcal{A} = 0.3$. The dissipation coefficient is fixed at $\alpha = 1$. All other network parameters are in the caption of Fig. 66.

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Therefore, the system can only reach equilibrium when Eq. (101) is satisfied, i.e. when $P'_1 + P'_2 = 0$. When instead $P'_1 + P'_2 > 0$, one defines $\Delta P = P'_1 + P'_2$ and $W_{ij} = \lambda a_{i,j}$, with λ being the coupling strength. Eq. (97) can be then rewritten as

$$\begin{aligned} \Delta \dot{\chi} &= \Delta P - \alpha \Delta \chi - 2\lambda \sin(\Delta\theta), \\ \Delta \dot{\theta} &= \Delta \chi, \end{aligned} \quad (102)$$

where $\Delta\theta = \theta_2 - \theta_1$ and $\Delta\chi = \Delta\dot{\theta}$. Fig. 69(a) shows the case of $2\lambda > \Delta P$, where there are two fixed points. One fixed point is stable and the other is unstable, therefore all trajectories converge to the stable fixed point. Fig. 69(b) shows the case of $2\lambda < \Delta P$, where there are no fixed points, and all trajectories converge to a limit cycle. Fig. 69(c) shows the case of $2\lambda \approx \Delta P$, where the fixed point and the limit cycle coexist such that the dynamics depends crucially on the initial conditions. Fig. 69(d) shows the parameter space of the systems. The majority of power grids are operating close to the edge of stability, i.e. in the region of coexistence, at least during periods of high loads [412]. Therefore, the dynamics depends crucially on the initial conditions, and static power grid models are yet insufficient to fully capture the complexity of the system.

5.2.3. Brain dynamics

Synchronization plays an important role in sustaining basic brain functions such as emotions, complex thoughts, memory, language comprehension, consciousness, etc. [420]. For instance, the importance of synchronization of oscillatory phases between different brain regions in memory processes has been demonstrated [421]. On the other side, clinical evidence points out that abnormal synchronization of a small group of neurons may result in ruining some of the brain functions, leading to pathological behaviors such as Parkinson disease, tremor, and epilepsy.

Let us take epilepsy as an example. The disease manifests itself with recurrent unprovoked seizures resulting from a wide variety of causes, and is also the world's most prominent serious brain disorder. Unless other neurological problems exist, patients with epilepsy typically have normal neurological function in between the seizures. Understanding the transition between the normal function and the seizure is useful for both improving the treatment of epilepsy and providing early warnings to patients.

Epilepsy has four phases: the interictal period, the onset of the seizure, and the propagation and termination phases. It is found that the coupling between brain areas during seizures changes with time, increasing or decreasing at the seizure onset in different cases [422]. At variance, when seizure termination is approached, the coupling of brain activity always increases. Both the onset and termination processes are very fast, indicating a jumping phase transition at critical coupling. The sudden emergence and termination of seizures suggested therefore that ES may play a role [423].

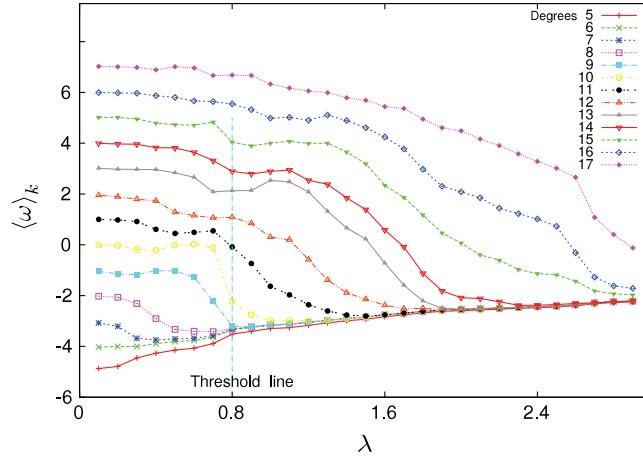


Fig. 68. Cluster explosive synchronization in uncorrelated SF networks. Vertical axis reports the average frequency of nodes of various degrees vs. the coupling strength λ (see the legend for the color codes of the different curves). (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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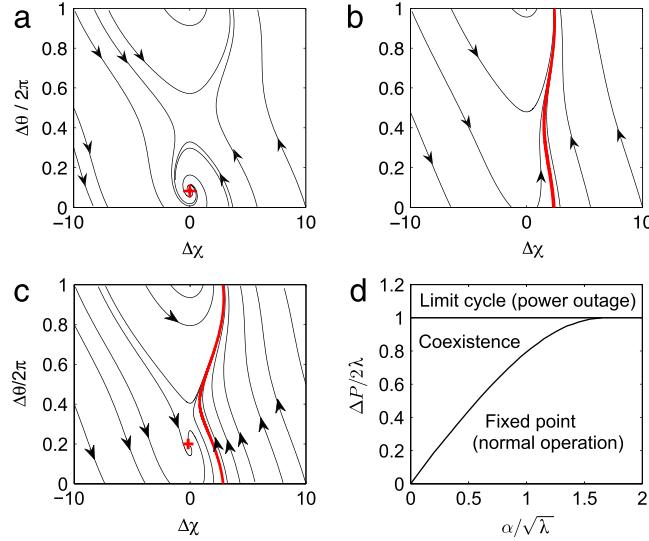


Fig. 69. Scenarios of the dynamics for an elementary network with one generator and one consumer for $\alpha = 1$. (a) Globally stable phase locking for $\Delta P = \lambda = 2$. (b) Globally unstable phase locking (limit cycle) for $\Delta P = 2$ and $\lambda = 0.5$. (c) Coexistence of phase locking (normal operation) and limit cycle (power outage) for $\Delta P = 2$ and $\lambda = 1.1$. (d) Stability phase diagram in the parameter space. Both stationary solutions (fixed point and limit cycle) are reported in red. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

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A seizure originates in a focal region, and then spreads to other brain regions. The individuation of the focal region from which an epileptic seizure is originated was performed in Ref. [424] in terms of phase synchronization indicators. During seizures, there is an overall increase of synchrony between the thalamus and temporal lobe structures. The phenomenon of seizure can be studied by means of a network approach, where a node can be either an individual neuron, or a population of neurons within a given structure, or an entire brain structure or region [423].

Guye et al. selected thirteen patients undergoing pre-surgical evaluation of drug-resistant temporal lobe epilepsy, and checked their recordings of intracerebral electrodes [425]. Fig. 2 of Ref. [425] reveals that, during the seizure, synchronization values are significantly higher than those of the background period, indicating a jumping transition. The abruptness of the transition in the seizure has been also confirmed by many other studies [422,423,426,427].

Another example is the anesthetic-induced unresponsiveness, which may result from specific interactions of anesthetics with the neural circuits regulating sleep and wakefulness. One common belief is that emergence from anesthesia is

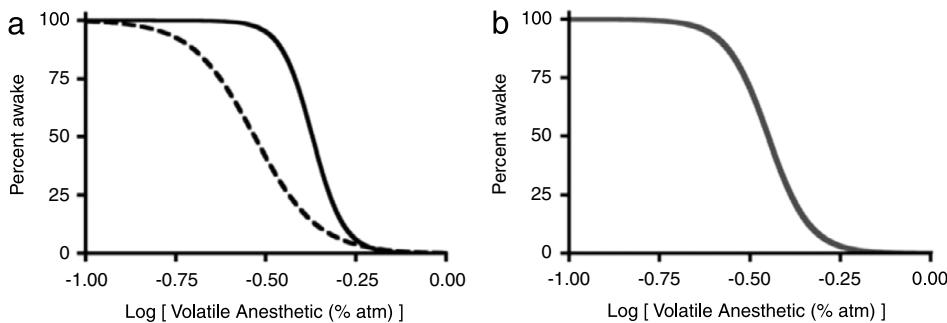


Fig. 70. Path-dependent and path-independent state transitions. (a) The solid black curve represents a population of individuals entering the state of unconsciousness as a function of the anesthetic dose. The dashed black curve depicts the reverse transition. The situation is here clearly characterized by the presence of a hysteresis loop. (b) In the absence of hysteresis, the forward and reverse paths are superimposed (thick gray curve). Source: Reprinted with permission from Ref. [428].

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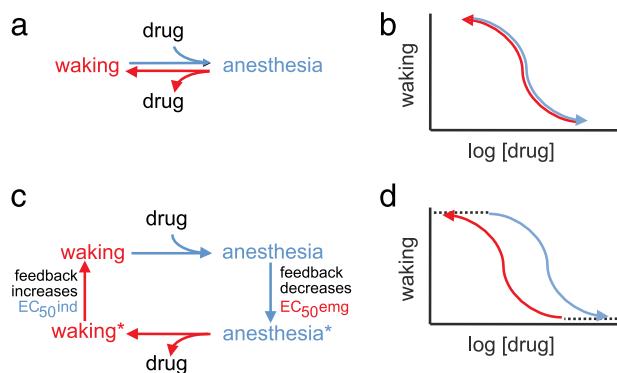


Fig. 71. (Color online). (a) Simple kinetic model describing drug-dependent behavioral state changes in the absence of bistability. (b) In the absence of feedback and bistability, dose-response curves for anesthesia induction and emergence are independent of the history of the prior behavioral state, and thus coincide. (c) Addition of feedback upon binding or unbinding of drug leads to additional, more stable, anesthesia and waking states. (d) Feedback and bistability lead to hysteresis in dose-response curves for anesthesia induction and emergence. Source: Reprinted with permission from Ref. [429].

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the inverse process of induction. Friedman et al. made experiments on transitions between conscious and unconscious states [428]. By generating anesthetic dose-response data in both insects and mammals, Ref. [428] demonstrated that the forward and backward paths through which anesthetic-induced unconsciousness arises and dissipates are not identical. Instead, a hysteresis is exhibited that is not fully explained by pharmacokinetics, as previously thought. Fig. 70 shows the results for both the cases of path-dependent (a) and path-independent (b) state transitions. In particular, Fig. 70(a) displays a hysteresis loop, in contrast to the common belief.

Joiner et al. provided further experimental evidence of the existence of hysteresis in the anesthetic-induced unresponsiveness [429]. The study proposes that processes selectively contributing to neural inertia may be impaired in pathophysiological conditions. Certain features exist in a minimal neural circuit that underly neural inertia, and can be used to model the dependence of anesthesia on feedback bistability. Fig. 71 represents a sketch of the models from Ref. [429]. The sketch (a) is a simple kinetic model describing the transitions between two states, one unbound and the other bound to drug; (b) is the resulting dose-response curves for the forward and reverse reactions; (c) is the bistable situation where distinct feedback mechanisms are activated to shift drug sensitivity towards stabilization of the state; and (d) is the hysteresis in dose-response curves for anesthesia induction and emergence.

Kim et al. recently hypothesized that the conditions for ES in human brain networks would be present in the anesthetized brain just over the threshold of unconsciousness [430]. To test this hypothesis, Ref. [430] constructed functional brain networks from multi-channel electroencephalogram recordings in seven healthy subjects across conscious, unconscious, and recovery states, and demonstrated for the first time that the network conditions for ES (especially the suppressive rule of Ref. [19]) are present in empirically-derived functional brain networks.

Together with the experimental studies for the hysteresis loop in brain functioning, some theoretical works have also tried to understand the underlying mechanism of ES in neural systems [63]. For instance, Chen et al. investigated the property of phase synchronization transition of coupled FitzHugh–Nagumo oscillators in BA SF networks, and reported the existence

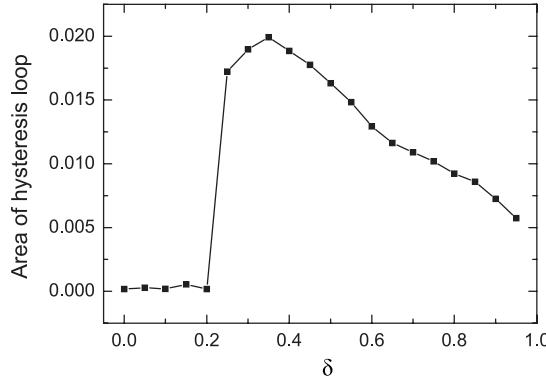


Fig. 72. Area of hysteresis loop (as a function of δ) in the synchronization diagram of Eqs. (104) for BA SF networks. Parameters are $N = 200$, $\epsilon = 0.01$, $D_x = 0$, $D_y = 0.005$, and the average degree $\langle k \rangle = 6$.
Source: Reprinted with permission from Ref. [63].

of ES. The model reads as

$$\begin{aligned} \epsilon \dot{x}_i &= x_i - x_i^3 - y_i + \lambda \sum_{j=1}^N A_{ij}(x_j - x_i) + \xi_i^{(x)}(t), \\ \dot{y}_i &= x_i + a_i + \xi_i^{(y)}(t), \end{aligned} \quad (103)$$

where $i = 1, 2, \dots, N$, $\epsilon = 0.01$, and x and y are the fast and the slow variables, respectively. The parameter a_i describes the excitability of the i th unit. If $|a_i| > 1$, the system is excitable, while $|a_i| < 1$ implies that the system is oscillatory. A_{ij} are the elements of the adjacency matrix of the network, λ is the coupling constant, and $\xi_i^{(\alpha)}(t)$ is a Gaussian noise that is independent for different units and satisfies $\langle \xi_i^{(\alpha)}(t) \rangle = 0$ and $\langle \xi_i^{(\alpha)}(t) \xi_j^{(\alpha)}(t') \rangle = 2D_\alpha \delta_{ij} \delta(t - t')$ with noise intensity D_α , and $\alpha \in \{x, y\}$. The natural frequency ω_i is assumed to be an increasing function of the degree k_i , by taking a_i as follows

$$a_i = 0.99 - \delta \frac{k_i - k_{min}}{k_{max} - k_{min}}, \quad (104)$$

where k_{max} and k_{min} are the maximum and minimum degrees in the network, respectively. The factor δ determines the slope of the linear expression. The larger is δ , the wider the distribution of a_i .

The synchronization diagram can be obtained by performing both the forward and the backward simulations. Chen et al. showed that, for BA networks, the nature of synchronization transition drastically changes with δ [63]. When δ is relatively small, the synchronization transition is continuous. Increasing δ to a middle value, one can see that as λ increases the order parameter r abruptly jumps from $r \approx 0$ to $r \approx 1$ at $\lambda_c^f = 0.044$, and therefore a sharp transition takes place at the onset of synchronization. On the other hand, the curve corresponding to the backward simulations also shows a sharp transition from the synchronized state to the incoherent one at $\lambda_c^b = 0.021$. The two sharp transitions occur at different values of λ , determining a strong hysteresis loop. To measure the influence of δ on ES, Chen et al. calculated the area of hysteresis loop, and found that, increasing further δ to $\delta = 0.9$, the first-order nature of the phase transition is still present, but the area of the hysteresis loop becomes smaller [63]. Fig. 72 shows how the area changes with δ : one can see that the area vanishes when $\delta = 0.2$, implying that the synchronization transition is there of second-order type. When δ is increased to $\delta = 0.25$, this area drastically changes to a non-zero value, indicating that the transition changes from a second-order type to a first-order one at around $\delta \approx 0.25$. With further increasing δ , this area shows a non-monotonic dependence on δ , and a maximum area occurs at $\delta = 0.35$.

5.2.4. Biological systems

Together with the example of the FitzHugh–Nagumo model of Eq. (103), other models of relevance for biological systems have been shown to display ES. Recently Chen et al. [76] reported that ES (as well as frequency–degree correlation properties) can take place in a very wide range of oscillatory networks, without any particular structure constraint. In Ref. [76], diffusively coupled SF networks of complex Ginzburg–Landau oscillators (CGLE) [431] are used,

$$\dot{z}_j = z_j - (1 + i\omega_0)|z_j|^2 z_j + \lambda(1 + i\beta) \sum_{l=1}^N a_{lj}(z_l - z_j), \quad (105)$$

where ω_0 is the oscillator's natural frequency, β is a dispersion parameter. Both parameters are taken to be identical for all nodes. CGLE can generate rather complicated dynamics, including chaos, and different ways exist for the definition of

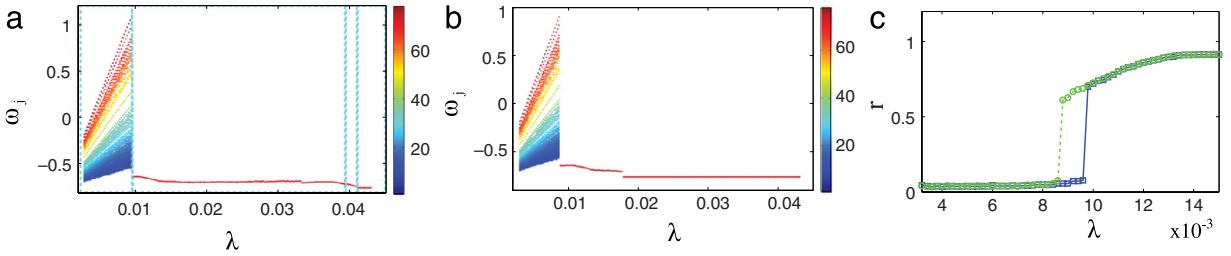


Fig. 73. Principle frequencies ω_j vs. the coupling λ for all nodes corresponding to (a) the forward transition, and (b) the backward transition for SF networks. The colors represent the degree of the nodes. Panel (c) reports the corresponding phase order parameter of panels (a) and (b) for the forward (blue squares) and backward (green circles) transitions. Parameters used in Eq. (105): $\omega_0 = 0.75$, $\beta = -1.7$, $N = 200$, and $\langle k \rangle = 20$. (For interpretation of the references to colour in this figure legend, the reader is referred to the web version of this article.)

Source: Adapted with permission from Ref. [76]. Courtesy of S. Wang.

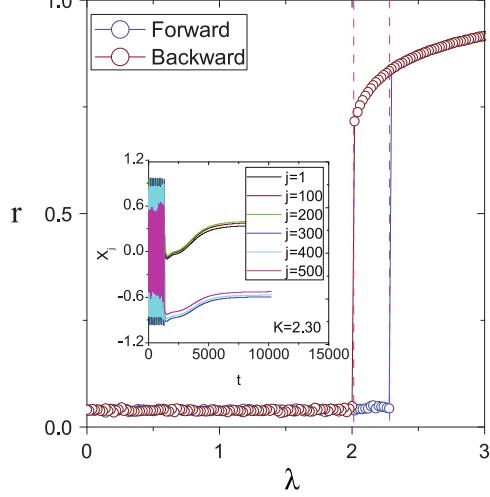


Fig. 74. (Color online). Characterization of explosive oscillation death in a Stuart–Landau model (Eq. (106)). Order parameter r vs. coupling strength λ for a triangular $g(\omega)$, $N = 500$. Inset: evolution of the variable $x_j(t)$ for all the nodes at the transition point.

Source: Adapted from Ref. [432] with permission of IOP.

frequency. In Ref. [76] the *principal frequency* of node j (ω_j) is defined as the value where the Fourier spectrum of a long time series of the dynamics of node j has the highest peak.

Fig. 73(a), (b) report the main results of the integration of Eq. (105). The principal frequencies ω_j of all nodes are there reported vs. the coupling strength λ . Panel (a) corresponds to the forward transition, panel (b) to the backward continuation. One observes a ES transition in the frequency domain. Before the transition, increasing of the coupling strength spontaneously induces a strongly inhomogeneous and well-ordered frequency distribution, with a positive correlation with the node degree. Defining the phase of node j as $\theta_j = \arctan[\text{Im}(z_j)/\text{Re}(z_j)]$, the authors compute also the phase order parameter r . The result is shown in Fig. 73(c), where one notices the presence of a well pronounced hysteresis.

In Ref. [432] Bi et al. investigated explosive oscillation death in another well known model of amplitude oscillators. The oscillation death refers to the complete suppression of oscillations in coupled systems. This phenomenon has been extensively studied, both theoretically and experimentally [433,434]. Ref. [432] considered a model of globally coupled Stuart–Landau oscillators with the frequency-weighted coupling term inspired by Ref. [61]:

$$\dot{z}_j = (a + i\omega_j - |z_j|^2)z_j + \lambda \frac{|\omega_j|}{N} \sum_{l=1}^N (z_l - z_j), \quad (106)$$

where $z_j(t) = x_j(t) + iy_j(t)$ is a complex variable, and a acts as the control parameter for the individual Stuart–Landau oscillator: the dynamics settles on a limit cycle if $a > 0$, and on a fixed point if $a < 0$. At variance with Ref. [76], oscillators here are not identical, and their natural frequencies are randomly taken from a frequency distribution $g(\omega)$.

Explosive oscillation dead was reported to occur in the solutions of Eq. (106). Fig. 74 shows a clear discontinuous jump in the forward and backward transitions for the order parameter $r e^{i\psi} = \sum_{j=1}^N z_j(t)/N$. The inset reports the evolution of the variable $x_j(t)$ for all the nodes at the transition point. It can be observed that not all the nodes collapse to the same fixed point.

5.2.5. Controlled laboratory experiments

Some few experimental evidences of ES have been given in controlled laboratory systems, which actually are of importance as they ultimately demonstrate the robustness and ubiquitousness of the phenomenon.

The first of these experiments was published by Leyva et al. in Ref. [16], and involved the search for explosive phase synchronization in frequency-degree correlated heterogeneous networks of electronic circuits. Initially, Ref. [16] performed numerical simulations of an ensemble of $N = 1000$ piece-wise linear Rössler units, interacting in a SF network via a bidirectional diffusive-like coupling [435,436]:

$$\begin{aligned}\dot{x}_i &= -\alpha_i \left[\Gamma \left(x_i - \lambda \sum_{j=1}^N a_{ij}(x_j - x_i) \right) + \beta y_i + \rho z_i \right], \\ \dot{y}_i &= -\alpha_i(-x_i + \nu y_i), \\ \dot{z}_i &= -\alpha_i(-g(x_i) + z_i),\end{aligned}\tag{107}$$

where the piece-wise linear part is given by

$$g(x_i) = \begin{cases} 0 & \text{if } x_i \leq 3 \\ \mu(x_i - 3) & \text{if } x_i > 3. \end{cases}\tag{108}$$

Every node (indexed by $i = 1, \dots, N$) is here represented by an associated three-dimensional vector $\mathbf{x}_i(t) \equiv (x_i(t), y_i(t), z_i(t))$. The parameters are: $\Gamma = 0.05$, $\beta = 0.5$, $\rho = 1$, $\mu = 15$, and $\nu = 0.02 - \frac{10}{R}$, where R is a tunable quantity that regulates the dynamical state of the system. In particular, R induces a chaotic dynamics in the range $R = [55, 110]$ [435].

Finally, the natural oscillation frequency of node i depends linearly on the parameter α_i . To impose a degree-frequency positive correlation in this system, the α_i values (and therefore, the oscillators' frequencies) are distributed following the relation

$$\alpha_i = \alpha \left(1 + \Delta\alpha \frac{k_i - 1}{N} \right),\tag{109}$$

where $\alpha = 10^4$, and $\Delta\alpha$ is a factor that determines the slope of the linear distribution. The range of frequencies in the ensemble becomes wider and wider as $\Delta\alpha$ and k_{\max} (the maximum degree in the network) are increased, i.e. as more heterogeneous degree distributions (and/or steeper slopes) are considered.

Ref. [16] shows that the system (107) gives rise to a ES transition, for an appropriate selection of the parameters. Precisely, the instantaneous phase for each oscillator i is defined as $\theta_i(t) = \arctan(y_i(t)/x_i(t))$, and the usual order parameter r is monitored as a function of λ . The results are shown in Fig. 75(a), for $N = 1000$, $\langle k \rangle = 6$ SF network obtained by the BA growing algorithm for different values of the power-law exponent γ of the degree distribution. The inset reports the amplitude average synchronization error $\langle e \rangle$ in the proximity of the transition for $\gamma = 3.0$, $R = 100$ and $\Delta\alpha = 8.0$, to confirm that the system does not reach full synchronization when phase ES occurs.

The specific chaotic state of the nodes is a relevant variable, as it can be seen from the comparison (see Fig. 75(a)) of the case $\Delta\alpha = 6.0$, $R = 70$ (black circles, second order phase transition) with the case $\Delta\alpha = 10.0$, $R = 100$ (blue diamonds, first order phase transition) for $\gamma = 3.0$. A more exhaustive description of ES emerges from the exploration of the full parameter space $\lambda-R$. The results are shown in Fig. 75(b), where the values of r , for both forward and backward simulations are reported. As it can be seen, the plane $\lambda-R$ can be clearly divided in two areas (denoted as I and II in the Figure) where the transition is of the first and second order, respectively. The striped portion of the area where the transition is of the first order marks the region where the hysteresis phenomenon is observed.

Motivated by the numerical study of Eqs. (107), Ref. [16] provided the first experimental evidence of ES, with the electronic network schematically shown in Fig. 76(a). The experiment consists of six piecewise Rössler circuits operating in the chaotic regime, which are labeled as $N1, N2, \dots, N6$. The details of the circuit construction, as well as the qualitative equivalence with the model of Eqs. (107) are available in Refs. [435,436].

The circuits are arranged in a star-like configuration, which, on its turn, represents the *maximally heterogeneous structure* available for small ensembles. All the chaotic oscillators have the same internal R_{\exp} , to ensure that they work in an almost identical dynamical regime. By a fine tuning of the values of the capacitors, circuits are configured such that the central node $N1$ oscillates with a mean frequency of 3333 Hz, and the leaves nodes $N2, \dots, N6$ are set with frequencies in the range of 2240 ± 200 Hz. Notice that, due to the experimental variability, the frequencies of the oscillators suffer also from unavoidable dispersion.

Once the data are stored, processing of them is made to obtain the equivalent instantaneous phases. These latter quantities are calculated by defining the instantaneous phase $\theta_i(t)$ of each oscillator i as $\theta_i(t) = 2\pi l_i + 2\pi \frac{t-t_l}{t_{l+1}-t_l}$ in each interval $t_l \leq t < t_{l+1}$, where t_l is the instant at which the l th crossing of the i th oscillator with its Poincaré section occurs. Fig. 76(b) reports the experimental synchronization diagram for three values of the parameter R_{\exp} , both for forward and backward variations of the coupling strength λ . As in the numerical case, the dynamical regime tuned by R_{\exp} determines whether the transition is of the first or second order.

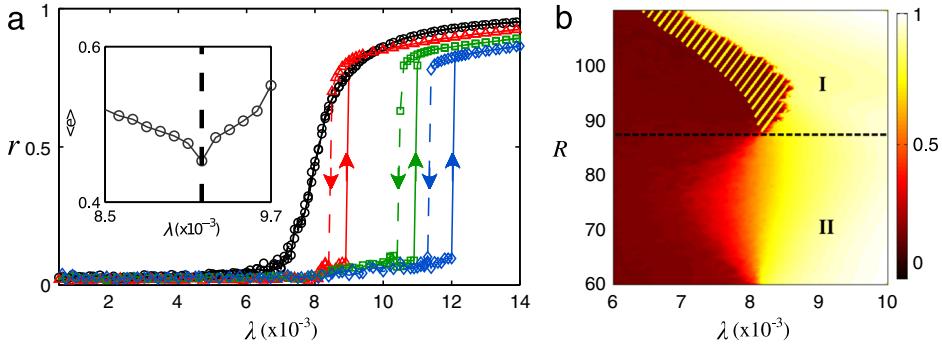


Fig. 75. (Color online). (a) Phase synchronization degree r as a function of the coupling strength λ for different SF networks of size $N = 1000$, and average degree $\langle k \rangle = 6$, $\gamma = 2.2$ (red triangles), $\gamma = 2.5$ (green squares), $\gamma = 3.0$ (blue diamonds and black circles). The correlation between node degree and natural frequency is set via Eq. (109). $\Delta\alpha = 6.0$ and $R = 70$ for the black circles case, while $\Delta\alpha = 10.0$ and $R = 100$ for the other networks. Continuous (dashed) lines mark the forward (backward) simulations, as d is increased (decreased) in steps of $\Delta\lambda = 3 \times 10^{-4}$. The inset reports the average synchronization error $\langle e \rangle$ vs. λ in the proximity of the first-order transition occurring for $\gamma = 3.0$, $R = 100$ and $\Delta\alpha = 8.0$. (b) Mean synchronization degree r in the parameter space λ - R . A second-order transition occurs for values of R below the horizontal dashed line. Above the same dashed line, the transition is instead of the first-order type, exhibiting the typical hysteresis (striped area) in the forward and backward simulations with $\Delta\lambda = 5 \times 10^{-5}$. The entire phase diagram refers to the case of a scale free network with $N = 1000$, $\langle k \rangle = 6$, $\gamma = 3$ and $\Delta\alpha = 6.0$.

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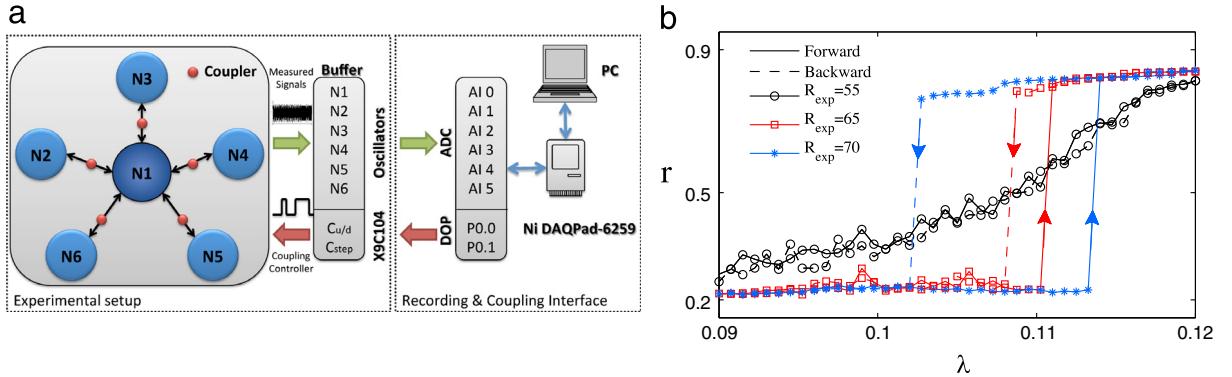


Fig. 76. (Color online). (a) Sketch of the experimental setup, where six Rössler circuits (blue nodes) are bidirectionally connected in a star configuration. (b) r vs. λ for the configuration shown in panel (a). The values of the internal resistance are: $R_{\text{exp}} = 55$ (black circles, second-order phase transition), $R_{\text{exp}} = 65$ (red squares, first order phase transition with narrow hysteresis), and $R_{\text{exp}} = 70$ (blue stars, first order phase transition with wide hysteresis).

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Experimental evidence of ES has been recently given also in coupled chemo-mechanical systems, and more specifically in mercury beating-heart (MBH) oscillators [437]. Ref. [437] considered a star network configuration, where a central MBH oscillator is connected with other three MBH leaves oscillators. Furthermore, the experimental parameters are tuned properly, in a way that the natural frequencies of each oscillator is proportional to the number of its links. Not only Ref. [437] gives evidence that a gradual increase of the coupling strength results in an abrupt and irreversible (first-order like) transition, but it also proved how to engineer magnetic-like states of synchronization, by the use of an external signal.

5.2.6. Control of explosive synchronization

A characteristic feature of continuous phase transitions is that after the transition point, the order parameter increases gradually with the coupling strength, which makes the dynamics of the system rather predictable. In contrast, the order parameter of a discontinuous transition has a jump at a critical point, which renders predictability of the dynamics more difficult to achieve. Control of ES means to change the transition from first-order to second-order like, avoiding big variations at the transition point.

In the first studies of ES, it was shown that the phase transition is first-order-like if the degree distribution exponent γ of the SF network satisfies $2 < \gamma < 3$, is of the second order if $\gamma > 3$, while for $\gamma = 3$ a hybrid phase transition is found [15, 359]. Later on, it was revealed that the conversion can be also induced by other factors such as the degree-degree correlation, degree-frequency correlation, disassortativity, etc. [22, 101, 362, 373, 375, 380]. For instance, Su et al. extended the condition of ES from strong correlation ($\omega_i = k_i$) to weak correlation ($\omega_i = f(k_i)$), and revealed that there is a conversion from continuous to discontinuous that takes place depending on the choice of the correlation function f [373]. Li et al. showed

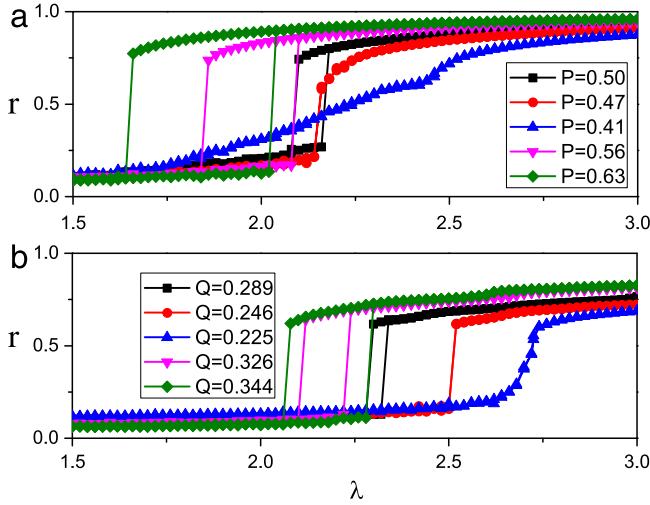


Fig. 77. (Color online). Synchronization diagrams obtained by exchanging the natural frequencies of connected nodes (see the text for the used exchanging strategy). (a) $g(\omega)$ is taken as a Gaussian distribution with both positive and negative ω . “Squares” denote the case where no exchanges of frequencies are performed, “circles” and “up triangles” represent the cases of decreasing P to $P = 0.47$, and 0.41 , respectively, and “down triangles” and “diamonds” represent the cases of increasing P to $P = 0.56$, and 0.63 , respectively. (b) $g(\omega)$ is taken as a Lorentzian distribution with only $\omega > 0$. “Squares” denote the case where no exchanges of frequencies are performed with $Q = 0.289$, “circles” and “up triangles” represent the cases of decreasing Q to $Q = 0.246$, and 0.225 , respectively, and “down triangles” and “diamonds” represent the cases of increasing Q to $Q = 0.326$, and 0.344 , respectively.

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that local degree-degree correlations (and precisely the degree of disassortative mixing) contribute primarily to ES [101]. Sendiña-Nadal et al. discussed the effects of degree correlations, and showed that high levels of positive and negative mixing consistently induce a second-order phase transition, while moderate values of assortative mixing enhance the irreversible nature of ES in SF networks [22]. Zhou et al. discussed explosive synchronization with asymmetric frequency distributions and found that the synchronization transition converts from the first order to the second order as the central frequency shifts towards the positive direction [375]. Pinto and Saa found a conversion when only a small part of the vertices of the network is subjected to a degree-frequency correlation [362].

However, in real situations the network properties (such as the degree distribution and assortativity) are often unchangeable, and therefore, other control techniques have to be implemented to attain the conversion. A first work in this line was done by Zhang et al., where the control is implemented by breaking the suppressive rule of Ref. [19] already introduced in Section 4.2.3 (Eq. (69)).

According to the suppressive rule, local synchronization can only occur for those neighboring pairs of nodes satisfying Eq. (69), while those nodes violating Eq. (69) will actually suppress the formation of local synchronized clusters. Ref. [19] showed that an effective way to break the suppressive rule is by randomly exchanging the frequencies of two nodes i and j .

The idea of Ref. [19] is as follows: if exchanging ω_i and ω_j would result in a value of S_{ij} [Eq. (70)] smaller than the original one, the exchange procedure is accepted. Otherwise, no operation is made. Ref. [19] further introduced the quantity

$$P = \frac{1}{2} - \frac{1}{2N\langle k \rangle} \sum_{i,j} A_{ij} \operatorname{sgn}(\omega_i \omega_j), \quad (110)$$

which represents the fraction of the pairs of connected oscillators with opposite sign in ω and which measures the effect of the exchanging. Fig. 77(a) reports the transitions to synchronization observed at different values of P . It is easy to see that decreasing P initially destroys the hysteretic loop (see the red curve with “circles”), and eventually produces a transition with no jump (see the blue curve with “up triangles”), i.e. it converts ES into a second-order phase transition.

Together with P , Zhang et al. also introduced the quantity

$$Q = \frac{1}{N\langle k \rangle} \sum_{i,j} A_{ij} Y_{ij}, \quad (111)$$

where $Y_{ij} \equiv \frac{|\omega_i - \omega_j|}{|\omega_i| + |\omega_j|}$. The quantity Q measures the average connection between nodes with large and small ω , and repeating the exchanging process driven by Eq. (70), it is possible to decrease or increase Q . Fig. 77(b) shows the resulting synchronization transitions at different values of Q . Similarly to Fig. 77(a), the progressive decreasing of Q has the effect of eventually leading to a second-order transition.

In some circumstances, however, the natural frequencies of the oscillators are also not changeable. To overcome this limitation, Zhang et al. recently presented another efficient method to control ES [438]. The idea is to introduce a small

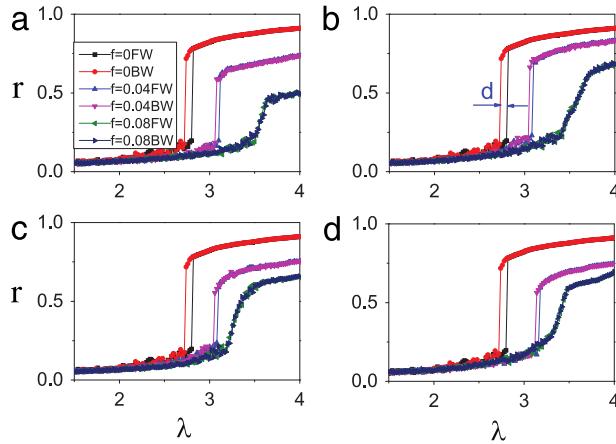


Fig. 78. (Color online). Forward and backward synchronization transitions for an Erdős-Rényi network with $N = 500$ and $\langle k \rangle = 6$, where “squares” and “circles” represent the forward and backward curves at $f = 0$, the “up triangles” and “down triangles” represent the forward and backward curves at $f = 0.04$, and the “left triangles” and “right triangles” represent the forward and backward curves at $f = 0.08$, respectively. The panels report the following cases: (a) the random-case1, (b) the random-case2, (c) the hetero-case1, and (d) the hetero-case2. See text for definition.
Source: Reprinted from Ref. [438] with permission of IOP.

fraction of contrarians to suppress the growing of larger clusters, in contrast to the conformists. There are two approaches to define a conformist or a contrarian. In the first method, a contrarian oscillator will receive interactions from its neighbors via a negative coupling strength, while a conformist oscillator will receive interactions from its neighbors via a positive coupling strength [391,439]. The model is as follows

$$\dot{\theta}_i = \omega_i + \frac{\lambda_i |\omega_i|}{k_i} \sum_{j=1}^N A_{ij} \sin(\theta_j - \theta_i), \quad i = 1, \dots, N \quad (112)$$

where the conformists have positive λ_i while the contrarians have negative λ_i . For simplicity, the coupling strength λ_i is set at the same amplitude $\lambda > 0$ with $\lambda_i = \lambda$ for all conformists and $\lambda_i = -\lambda$ for all contrarians.

In the second method, a contrarian oscillator gives negative coupling to each of its neighbors, while a conformist oscillator gives a positive coupling to each of its neighbors [387,440,441]. The model can be written as follows

$$\dot{\theta}_i = \omega_i + \frac{|\omega_i|}{k_i} \sum_{j=1}^N \lambda_j A_{ij} \sin(\theta_j - \theta_i), \quad i = 1, \dots, N \quad (113)$$

where the conformists contribute positive λ_j , while the contrarians contribute negative λ_j . Once again, all strengths λ_j are set at the same amplitude $\lambda > 0$, with $\lambda_j = \lambda$ for all conformists and $\lambda_j = -\lambda$ for all contrarians.

The first case considered is that of random uniformly distributed contrarians, with the definition of contrarian in Eq. (112) (this setting is called *the random-case1*). Let f represents the fraction of contrarians, i.e. the ratio between the number of contrarians and the total number of oscillators. Fig. 78(a) shows how the order parameter r changes with λ for $f = 0$, 0.04 and 0.08, respectively. It is easy to see that there are hysteretic loops for the cases of $f = 0$ and $f = 0.04$, but no loop is present for the case of $f = 0.08$, indicating that the first-order transition to synchronization has been changed into a second-order transition. A similar scenario, reported in Fig. 78(b), characterizes the case of random uniformly distributed contrarians with the definition of Eq. (113), called *the random-case2*. In order to inspect how the distribution of contrarians influences the controlling effect, contrarians are heterogeneously distributed in the network. Fig. 78(c) (Fig. 78(d)) shows the case of heterogeneously distributed contrarians with the definition given in Eq. (112) (Eq. (113)) called *the hetero-case1* (*the hetero-case2*). Comparing the four panels of Fig. 78, it is clear that in all cases a second-order phase transition is obtained when $f = 0.08$.

From Fig. 78, it can be also noticed that the transition point λ_c increases with the increase of f . In this way, the induced second-order transition point λ_{c2} will be much larger than the original first-order transition point λ_{c1} . This means that not only the first-order transition is suppressed but also the transition point will be postponed, which is a relevant result in real systems with potential cascading risk.

This finding may be also useful in explaining the normal functioning of brain. It is well known that there are both excitatory and inhibitory neurons in cortical neural networks. The excitatory neurons comprise the majority (80%–90%) of the neuronal population, and are largely homogeneous. The inhibitory neurons comprise only 10% of the neuron population, but are extremely heterogeneous, and take the role of controlling and coordinating the activity of large populations of local neurons [442,443]. One can then imagine, that the excitatory and inhibitory neurons are in fact the conformists and contrarians, respectively, and that the existence of inhibitory neurons in brain network is necessary not only for sustaining its normal behavior, but also for preventing its abnormal functioning.

6. Conclusions and future perspectives

After having revisited the main theory (and some applications) of explosive percolation and synchronization in complex networks, we end this review with a few concluding remarks, that have the aim of spotting problems that still remain open to future progresses, as well as of offering a few thoughtful considerations about what are the questions of relevance that should (in our opinion, hopefully soon) attract the attention of scientists in this area.

Explosive percolation was largely examined in Section 2, and its main applications in physical systems discussed in Section 5.1. The process of percolation refers to the emergence of large-scale connectivity on an underlying network, or lattice. The extensive studies of EP only started on 2009, following a seminal work by Achlioptas, where it was pointed out that introducing competitive mechanisms in the selection of bonds to be added to a graph, the threshold of the percolation process can be significantly postponed, and moreover, the giant cluster emerges after a number of steps that is much smaller than the system size. As a consequence, the order parameter exhibits an extremely abrupt jump at the percolation point, which makes this situation drastically different from classical, random, percolation.

As reported in full details in Section 2, many other algorithms and models have been introduced and studied since 2009, and so far EP has been intensively investigated on many networks topologies, including 2D lattices, 3D or high-D lattices, Bethe lattices, ER networks, SF networks, modular networks, and real-world networks. All these works have triggered a big debate in the community, aiming at clarifying whether the observed transitions are really discontinuous or not. While the existence of a seemingly discontinuous jump in the order parameter of the original Achlioptas model is universally recognized, careful numerical investigations of the cooperation, phase coexistence, and nucleation (as well as strict theoretical proofs) have demonstrated that EP in that model is actually continuous in the thermodynamic limit, but it belongs to a universality class different from any other previously considered.

Furthermore (and notably), consensus has been found on the fact that some other models of percolation (such as the k -core percolation model, the spanning cluster-avoiding model, the two-species cluster aggregation model, or the models of jamming on low-dimensional lattices) lead to truly discontinuous transitions. Therefore, it is now fully accepted that the term “explosive” has to refer to the unusual feature which distinguishes the critical behavior of Achlioptas processes from both ordinary and truly discontinuous models.

There are issues that we believe still need efforts from the scientific community. One is that, undoubtedly, more theoretical and numerical studies on the scaling properties and critical exponents in EP are in order, to clarify which new features characterize the critical behavior associated to explosive percolation. Another is the need of exploring EP in other realistic situations (real-world systems or experiments), such as, for instance, the cascading failure processes in power grids. And finally, the hope is that the gathered knowledge will help devising methods to control EP, i.e. to enhance or suppress it, when necessary or when desired.

On the other side, ES has been largely discussed in Section 4, and its main applications described in Section 5.2. From a historical point of view, we saw that the discovery of ES went through a somehow tortuous path: the community was initially apathetic to the early evidences of first order transitions to synchronization in the thermodynamic limit of a model of globally coupled phase oscillators. Such a situation of guilty indifference persisted for long time, up to when it was clarified that a positive correlation between the natural frequencies of oscillators and the degrees of nodes could produce abrupt cooperative behaviors. For the first time, it was realized that a complete understanding of the nature of phase transitions intimately depended on an explicit interaction between the network topology and the characteristic dynamics of oscillators, and therefore much attention has been recently paid to ES, which begot today great progresses, including theoretical understandings and experimental confirmations.

Let us concentrate on a few issues that possibly will become hot topics on ES.

The first is extending the study of explosive transitions to the case of partial and cluster synchronization, a large area of research, indeed, having to do with cases (like brain functioning, or stock market dynamics, to quote just two examples) for which synchronization on the whole system scale is not possible nor desirable, but rather collective arrangements emerge at the system's micro- and meso-scales. How to realize explosive transitions to these latter states from the incoherent behavior is still an open question.

The second is bridging ES with specific mixed states of coupled oscillators. Recently, indeed, Zhang et al. proposed a bridge between ES and chimera state [444], and found that they can both coexist within the same coupled system. Further studies on the relationship between ES and chimera state are certainly to be expected soon. Furthermore, since it is now understood that ES can be considered as a dynamical percolation process in phase space, an issue of the utmost relevance is clarifying how one can use the results of explosive percolation for a better control of ES. Therefore, further studies on the relationship between ES and explosive percolation are certainly necessary.

It is convenient to spend a few words on the novel, non-stationary, state that has been described in Section 4.4, the *Bellerophon* state. In globally coupled nonidentical oscillators, this state can indeed emerge, made of quantized clusters where oscillators are neither phase- nor frequency-locked. Oscillators' instantaneous speeds are different within the clusters, but they are correlated and, more importantly, they behave periodically in time so that their average values are the same. So far, it has been found that the Bellerophon states occur in fact in a pretty wide context of globally coupled oscillators, and for various typical frequency distributions. Physically, such a state can be regarded as a weaker form of coherence achieved by the coupled oscillators when the control parameter is at an intermediate value, or (in other words) as a transitional state between the incoherent and the phase-locked state: on the one hand the control parameter is not

strong enough to completely entrain the system into the phase-locked state, on the other hand it is adequately large to maintain certain correlations among the instantaneous frequencies of oscillators.

Certainly, a lot of future work is in order to unveil Bellerophon states in other oscillatory models, as well as to check if (and under which conditions) the hypothesis of a global coupling can be relaxed to that of a complex network architecture. On the other side, we are totally confident that the discovery of Bellerophon states will stimulate the search for a better theoretical description and characterization, as well as a special care in seeking for such novel states in experimental and natural systems.

Finally, another interesting issue is deepening and enhancing the knowledge of ES in physical systems, where the space embedding (and therefore the lengths of physical links, such as synapses in real neural systems or power lines in power grids), or the multiple level character of interactions between the units (which would require a multi-layer network representation) may play a key role. The study of ES in physically embedded and multi-layer networks is still in its infancy, and we firmly believe that it should instead attract soon a lot of attention, due to its evident relevance for practical applications.

Although beyond the scope of this review, it is worth mentioning that other abrupt transitions have been recently described in networked systems, having to do with spreading processes (like epidemics with co-infection [445,446] or re-infection [447] mechanisms), spontaneous recovery in dynamical networks [448], and social games using the inertial majority-vote model [449].

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