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Synchronisation of chaos and its applications

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ABSTRACT

Dynamical networks are important models for the behaviour of complex systems, modelling physical, biological and societal systems, including the brain, food webs, epidemic disease in populations, power grids and many other. Such dynamical networks can exhibit behaviour in which deterministic chaos, exhibiting unpredictability and disorder, coexists with synchronisation, a classical paradigm of order. We survey the main theory behind complete, generalised and phase synchronisation phenomena in simple as well as complex networks and discuss applications to secure communications, parameter estimation and the anticipation of chaos.

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1. Introduction

This survey provides an introduction to the phenomenon of synchronisation in coupled chaotic dynamical systems. Both chaos and synchronisation are important concepts in science, from a philosophical as well as a practical point of view.

Synchronisation expresses a notion of strong correlations between coupled systems. In its most elementary and intuitive form, synchronisation refers to the tendency to have the same dynamical behaviour. Scientists also recognise weaker forms of synchronisation, where some key aspects of dynamical behaviour are the same – like frequencies – or where coupled dynamical behaviours satisfy a specific spatiotemporal relationship – like a constant phase lag.

Synchronisation is fundamental to our understanding of a wide range of natural phenomena, from cosmology and natural rhythms like heart beating [1] and hand clapping [2] to superconductors [3]. While synchronisation is often beneficial, some pathologies of the brain such as Parkinson disease [4,5] and epilepsy [6] are also related to this phenomenon. In ecology, synchronisation of predators can lead to extinction [7,8] while improving the quality of synchronised behaviour of prey can increase the odds to survive [9]. In epidemiology, synchronisation in measles outbreaks can cause social catastrophes [10].

Synchronisation is also relevant to technology. Lasers form an important example. The stability of a laser generally decreases when its power increases. A successful way to create a high-power laser system is by combining many low-power stable lasers. A key challenge is to make

sure that the lasers synchronise [11–13], as without synchronisation destructive interference diminishes power.

Synchronisation can also cause engineering problems. A recent well-publicised example concerns the *London Millennium Bridge*, traversing the River Thames. On the opening day in 2000, the bridge attracted 90,000 visitors, holding up to 2000 visitors on the bridge at the same time. Lateral motion caused by the pedestrians made the bridge lurches to one side, as a result of which the pedestrians would adjust their rhythm to keep from falling over. In turn, this led to increased oscillations of the bridge due to the synchronisation between the bridge's oscillations and pedestrians' gait [14–16]. Eventually, the oscillations became so extensive that the bridge was closed for safety reasons. The bridge was only opened to the public again after a redesign where dampers were installed to increase energy dissipation and thereby impede synchronisation between bridge and pedestrians.

The above examples of synchronisation in coupled systems describe a spontaneous transition to order because of the interaction. Coupled systems are modelled as networks of interacting elements. We often have a detailed understanding about the dynamics of the individual uncoupled elements. For example, we have reasonably good models for individual superconducting Josephson junctions, heart cells, neurons, lasers and even pedestrians. In the systems we consider here, the coupling between elements is assumed to be built up from bilateral interactions between pairs of elements, and a network structure indicating which pairs of elements interact with each other. First, the way the individuals talk to each

other. For example, in the neurons the interaction is mediated by synapses and in heart cells by electrical diffusion. Second, the linking structure describing who is influencing whom. So, it is the network structure that provides the interaction among individual elements. The collective behaviour emerges from the collaboration and competition of many elements mediated by the network structure.

Synchronisation can be effectively used to create secure communication schemes [17–19]. And it can help developing new technologies. Synchronisation is also used for model calibration, that is, the synchronised regime between data and equations can reveal the parameter of the equations [20,21].

Chaos in dynamics is one of the scientific revolutions of the twentieth century that has deepened our understanding of the nature of unpredictability. Initiated by Henri Poincaré in the late-nineteenth century, the chaos revolution took off in the 1980s when computers with which chaotic dynamics can be studied and illustrated, became more widely available. Chaos normally arises when recurrent dynamical behaviour has locally dispersing characteristics, as measured by a positive Lyapunov exponent. In this review, we will not discuss any details of chaotic dynamics in detail. For a comprehensive monograph on this topic, see for instance [22].

At first sight it may appear that the concept of synchronisation, as an expression of order, and the concept of chaos, associated with disorder, could not be more distant from one another. Hence, it was quite a surprise when physicists realised that coupled chaotic systems also could spontaneously synchronise [23,24]. Despite many years of studies into this phenomenon and its applications, many fundamental problems remain open. This survey is meant to provide a concise overview of some of the most important theoretical insights underlying our current understanding of synchronisation of chaos as well as highlighting some of the many remaining challenges.

In this review, we will discuss the basic results for synchronisation of chaotic systems. The interaction can make these systems adapt and display a complicated unpredictable dynamics while behaving in a synchronous manner. Synchronisation in these systems can appear in hierarchy depending on the details of the individual elements and the network structure. We will first discuss this hierarchy in two coupled chaotic oscillators and latter generalise to complex networks. The review is organised as follows. In Section 2 we discuss the synchronisation scenarios between two coupled oscillators. In Section 3 we discuss such applications where the synchronisation phenomenon can be used for prediction and parameter estimation. In Section 4 we discuss synchronisation in complex networks.

2. Synchronisation between two coupled systems

2.1. Synchronisation of linear systems

Before touching upon more topical and interesting settings in which synchronisation is observed in non-linear systems, as an introduction we consider the elementary example of synchronisation between two linearly coupled linear systems. Although simple, this example bears the main ideas of the general case of synchronisation between two or more nonlinearly coupled systems.

We consider two identical linear systems

$$\dot{x}_i := \frac{dx_i}{dt} = ax_i, \quad i = 1, 2$$

with a a non-zero constant. The solution of these linear differential equations with initial condition $x_i(0)$ is $x_i(t) = e^{at}x_i(0)$ so that the ensuing dynamics is simple and all solutions converge exponentially fast to zero if $a < 0$, or diverge to infinity if $a > 0$ unless $x(0) = 0$.

We now consider these linear systems coupled in the following way

$$\begin{aligned} \dot{x}_1 &= ax_1 + \alpha(x_2 - x_1) \\ \dot{x}_2 &= ax_2 + \alpha(x_1 - x_2) \end{aligned} \quad (1)$$

and α is called the *coupling parameter*.

In the context of this model, we speak of *Complete Synchronisation* (CS) if $x_1(t)$ and $x_2(t)$ converge to each other as $t \rightarrow \infty$. In order to study this phenomenon, it is natural to consider the new variable

$$z := x_1 - x_2.$$

In terms of this variable, synchronisation corresponds to the fact that $\lim_{t \rightarrow \infty} z(t) = 0$. As $\dot{z} = \dot{x}_1 - \dot{x}_2$, we find directly by substitution from Equation (1) that

$$\dot{z} = (a - 2\alpha)z,$$

which has the explicit solution $z(t) = z(0)e^{(a-2\alpha)t}$. Hence, we find that $\lim_{t \rightarrow \infty} z(t) = 0$ if and only if $a - 2\alpha < 0$ (unless the initial condition is already synchronised, i.e. $z(0) = 0$). Defining the *critical coupling value* α_c as

$$\alpha_c := \frac{a}{2} \quad (2)$$

we thus obtain synchronisation if the coupling parameter exceeds the critical value: $\alpha > \alpha_c$.

We finally note that as the system synchronises, the coupling term converges to zero and the solution of each of the components behaves in accordance with the underlying uncoupled linear system: to be precise,

$\lim_{t \rightarrow \infty} \left(x_i(t) - \frac{x_1(0) + x_2(0)}{2} e^{at} \right) = 0$ for $i = 1, 2$. It is important to note that the sign of the parameter a here determines a difference between synchronisation to the trivial equilibrium (if $a < 0$) or to an exponentially growing solution (if $a > 0$).

In view of later generalisations, we will go through the above analysis again, exploiting more the linear structure of the problem so that we can appreciate synchronisation in terms of spectral properties of the coupling term.

With $\mathbf{x} := \begin{pmatrix} x_1 \\ x_2 \end{pmatrix}$, (1) can be written as

$$\dot{\mathbf{x}} = [a\mathbf{I} - \alpha\mathbf{L}]\mathbf{x} \quad (3)$$

where

$$\mathbf{I} = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} \text{ and } \mathbf{L} = \begin{pmatrix} 1 & -1 \\ -1 & 1 \end{pmatrix}.$$

L is known as the *Laplacian matrix*. In Section 4, we will generalise it to any network. The solution of (3) with initial condition $\mathbf{x}(0)$ is

$$\mathbf{x}(t) = e^{[a\mathbf{I} - \alpha\mathbf{L}]t} \mathbf{x}(0), \quad \text{where } e^{A^t} := \sum_{n=0}^{\infty} \frac{t^n}{n!} A^n. \quad (4)$$

To solve (4) we note that since \mathbf{I} and L commute,

$$e^{[a\mathbf{I} - \alpha\mathbf{L}]t} = e^{a\mathbf{I}t} e^{-\alpha\mathbf{L}t}$$

and $e^{a\mathbf{I}t} = e^{at} \mathbf{I}$. In order to evaluate $e^{-\alpha\mathbf{L}t}$, it is useful to observe that $\mathbf{v}_1 = (1, 1)^*$ and $\mathbf{v}_2 = (1, -1)^*$ are the eigenvectors of L for its corresponding eigenvalues $\lambda_1 = 0$ and $\lambda_2 = 2$. As $\{\mathbf{v}_1, \mathbf{v}_2\}$ is a basis of \mathbb{R}^2 , we may write any initial condition as $\mathbf{x}(0) = c_1 \mathbf{v}_1 + c_2 \mathbf{v}_2$ with $c_1, c_2 \in \mathbb{R}$, so that

$$e^{-\alpha\mathbf{L}t} \mathbf{x}(0) = c_1 \mathbf{v}_1 + c_2 e^{-\alpha\lambda_2 t} \mathbf{v}_2$$

and

$$\mathbf{x}(t) = e^{[a\mathbf{I} - \alpha\mathbf{L}]t} \mathbf{x}(0) = c_1 e^{at} \mathbf{v}_1 + c_2 e^{(a - \alpha\lambda_2)t} \mathbf{v}_2. \quad (5)$$

Synchronisation corresponds to the phenomenon that $\mathbf{x}(t)$ converges to the *synchronisation subspace* generated by \mathbf{v}_1 . This only happens if $\lim_{t \rightarrow \infty} c_2 e^{(a - \alpha\lambda_2)t} \mathbf{v}_2 = \mathbf{0}$, i.e. if $\alpha > \frac{a}{\lambda_2}$. Thus in view of (2), we define the *critical coupling value*

$$\alpha_c = \frac{a}{\lambda_2} = \frac{a}{2}.$$

We note that the critical coupling value is expressed in terms of the gap between the lowest eigenvalue 0 and smallest nonzero (and positive) eigenvalue of the Laplacian L .

2.2. Complete synchronisation of non-linear systems

We now consider two fully diffusively coupled identical non-linear n -dimensional systems

$$\begin{aligned} \dot{\mathbf{x}}_1 &= \mathbf{f}(\mathbf{x}_1) + \alpha \mathbf{H}(\mathbf{x}_2 - \mathbf{x}_1) \\ \dot{\mathbf{x}}_2 &= \mathbf{f}(\mathbf{x}_2) + \alpha \mathbf{H}(\mathbf{x}_1 - \mathbf{x}_2) \end{aligned} \quad (6)$$

where $\mathbf{f} : \mathbb{R}^n \rightarrow \mathbb{R}^n$ is in general non-linear and $\mathbf{H} : \mathbb{R}^n \rightarrow \mathbb{R}^n$ is a smooth coupling function. We assume that $\mathbf{H}(\mathbf{0}) = \mathbf{0}$ so that the synchronisation subspace $\mathbf{x}_1 = \mathbf{x}_2$ is invariant for all coupling strengths α . Meaning that for any synchronised initial condition the entire solution remains synchronised: as in the synchronised state the diffusive coupling term vanishes, the dynamics is identical to that of the uncoupled system (with $\alpha = 0$). Consequently, the coupling has no influence on the synchronised motion. In particular, it could be the case that the synchronised motion is chaotic, if the uncoupled systems exhibit such behaviour.

We aim to show that if the coupling is sufficiently strong, the system Equation (6) will synchronise $\mathbf{x}_1(t) - \mathbf{x}_2(t) \rightarrow \mathbf{0}$ as $t \rightarrow \infty$. We consider $\mathbf{H} = \mathbf{I}$ (the identity matrix) then the term reads as

$$\alpha \mathbf{H}(\mathbf{x}_2 - \mathbf{x}_1) = \alpha(\mathbf{x}_2 - \mathbf{x}_1).$$

To analyse stability, we consider – as before – the evolution of the difference variable $\mathbf{z} := \mathbf{x}_1 - \mathbf{x}_2$ in terms of which the synchronisation subspace is characterised as $\mathbf{z} = \mathbf{0}$:

$$\dot{\mathbf{z}} = \dot{\mathbf{x}}_1 - \dot{\mathbf{x}}_2 \quad (7)$$

$$= \mathbf{f}(\mathbf{x}_1) - \mathbf{f}(\mathbf{x}_2) - 2\alpha \mathbf{z} \quad (8)$$

The aim is to identify sufficient conditions for the coupling parameter α to guarantee that locally near $\mathbf{z} = \mathbf{0}$ we have $\lim_{t \rightarrow \infty} \mathbf{z}(t) = \mathbf{0}$. To this end, we linearise the equations of motion Equation (8) near $\mathbf{z} = \mathbf{0}$. We note to this extent that near $\mathbf{x}_1 = \mathbf{x}_2$ we obtain by Taylor expansion that

$$\begin{aligned} \mathbf{f}(\mathbf{x}_2(t)) &= \mathbf{f}(\mathbf{x}_1(t)) - D\mathbf{f}(\mathbf{x}_1(t))(\mathbf{x}_2(t) - \mathbf{x}_1(t)) \\ &\quad + O(\|\mathbf{x}_1(t) - \mathbf{x}_2(t)\|^2) \\ &= \mathbf{f}(\mathbf{x}_1(t)) - D\mathbf{f}(\mathbf{x}_1(t))\mathbf{z}(t) + O(\|\mathbf{z}(t)\|^2). \end{aligned}$$

Here $D\mathbf{f}(\mathbf{x}_1(t))$ is the derivative (Jacobian matrix of $\mathbf{f}(\mathbf{x})$) at $\mathbf{x} = \mathbf{x}_1(t)$. We use this to write Equation (8) near $\mathbf{z} = \mathbf{0}$ as

$$\frac{d\mathbf{z}}{dt} = [D\mathbf{f}(\mathbf{x}_1(t)) - 2\alpha\mathbf{I}]\mathbf{z} + O(\|\mathbf{z}\|^2). \quad (9)$$

The linear part of this equation, obtained by ignoring the $O(|z|^2)$ term in Equation (9), is commonly known as the *first variational equation*. It should be noted that this equation is nonautonomous as it depends explicitly on the reference solution $\mathbf{x}_1(t)$. In general it is not easy to analyse nonautonomous differential equations, not even linear ones. Fortunately, we are able to achieve insights without solving this equation because the coupling is rather convenient adding an extra damping term $-\alpha z$.

To simplify the analysis, we introduce a new variable

$$\mathbf{w}(t) = e^{2\alpha t} \mathbf{z}(t) \quad (10)$$

in terms of which the linear part of Equation (9) becomes precisely the variational equation for the solution $\mathbf{x}_1(t)$ of the uncoupled equation of motion $\dot{\mathbf{x}} = \mathbf{f}(\mathbf{x})$:

$$\begin{aligned} \dot{\mathbf{w}}(t) &= 2\alpha e^{2\alpha t} \mathbf{z}(t) + e^{2\alpha t} \dot{\mathbf{z}}(t) \\ &= 2\alpha \mathbf{w} + [D\mathbf{f}(\mathbf{x}_1(t)) - 2\alpha \mathbf{I}]e^{2\alpha t} \mathbf{z} \\ &= [D\mathbf{f}(\mathbf{x}_1(t))] \mathbf{w}. \end{aligned} \quad (11)$$

Let $\Phi(\mathbf{x}_1(t))$ be the fundamental matrix for the variational equation, so that any solution of this nonautonomous equation can be written as $\mathbf{z}(t) = \Phi(\mathbf{x}_1(t))\mathbf{z}(0)$. Let $\{\lambda_j(\mathbf{x}_1(t))\}_{j=1}^n$ be the set of positive square roots of the eigenvalues of the symmetric matrix $\Phi(\mathbf{x}_1(t))^* \Phi(\mathbf{x}_1(t))$ (where $*$ denotes transpose). Then we define

$$\Lambda := \max_j \lim_{t \rightarrow \infty} \frac{1}{t} \lambda_j(\mathbf{x}_1(t)). \quad (12)$$

Λ is known as the *Lyapunov exponent* of the orbit $\mathbf{x}_1(t)$ and it measures the infinitesimal asymptotic divergence rate near this trajectory. We refer to Appendix 7 for more details about the Lyapunov exponent.

The assertion now is that if the orbit $\mathbf{x}_1(t)$ has Lyapunov exponent Λ , then there exists a constant $C > 0$ such that

$$\|\mathbf{w}(t)\| \leq C e^{\Lambda t}. \quad (13)$$

From Equation (13) and using Equation (10) we obtain that

$$\|\mathbf{z}(t)\| \leq C e^{(\Lambda - 2\alpha)t}.$$

Hence,

$$\alpha_c := \frac{\Lambda}{2}$$

is a critical coupling strength for synchronisation, above which observe synchronisation.

A complication with the above analysis, is that the Lyapunov exponent Λ and constant C may depend on the chosen trajectory $\mathbf{x}_1(t)$. The probabilistic (ergodic) theory of dynamical systems, which we will not dwell on here, asserts that often the Lyapunov exponent is constant for almost all trajectories on a given attractor. However,

the constant C may still vary per trajectory, which leads to *non-uniform* synchronisation, implying the potential of large variation of transit times until synchronisation occurs. For similar phenomena, see also [25,26]. We now proceed to apply the above to the examples of coupled Lorenz and Rössler systems.

Lorenz system. The Lorenz system was introduced by Edward Lorenz in 1963 as a simplified model for atmospheric convection:

$$\begin{aligned} \dot{x} &= \sigma(y - x), \\ \dot{y} &= x(\rho - z) - y, \\ \dot{z} &= -\beta z + xy, \end{aligned} \quad (14)$$

where the three coordinates x , y and z represent the state of the system and σ , ρ , β are parameters. When parameter values are chosen as $\sigma = 10$, $\rho = 28$ and $\beta = 8/3$, the equations display unpredictable (chaotic) dynamics. Lorenz used this choice of parameters in his original paper [27]. We use these parameter values as well.

The Lorenz equations are dissipative and all trajectories eventually enter the absorbing domain

$$\Omega = \left\{ \mathbf{x} \in \mathbb{R}^3 : \rho x^2 + \sigma y^2 + \sigma(z - 2\rho)^2 < \frac{\beta^2 \rho^2}{\beta - 1} \right\},$$

see Appendix C or Ref. [28]. For the classical parameters, $\sigma = 10$, $\rho = 28$ and $\beta = 8/3$, inside Ω , trajectory accumulates on the chaotic Lorenz attractor [29], as depicted in Figure 1(b). Close to the attractor nearby trajectories diverge. To see this, we simulate two trajectories with nearly the same initial condition. The initial 10^{-2} difference grows to roughly 10^2 in a matter of only six cycles, see Figure 1(a). Using numerical simulations, we estimate the maximal divergence rate of nearby trajectories $\Lambda \approx 0.906$.

We consider two coupled chaotic Lorenz oscillators, as in Equation (6). We derived above that the critical coupling α_c for synchronisation depends on the Lyapunov exponent Λ . Using the numerical results for Λ we obtain

$$\alpha_c = \frac{\Lambda}{2} \approx 0.453.$$

Simulation confirms that this critical coupling is sharp. Indeed, for $\alpha = 0.4$ there is no synchronisation, and trajectories do not move together, see Figure 2(a). On the other hand, when $\alpha = 0.5$, the trajectories synchronise, see Figure 2(b).

To compare the amount of synchronisation at different parameter values, we may consider the average deviation from synchronisation during a time-interval of length T as

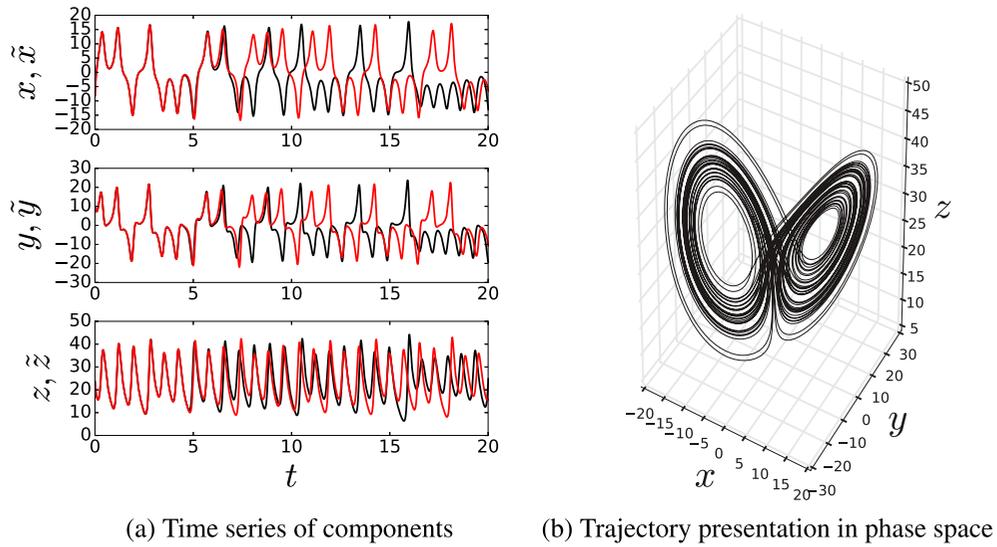


Figure 1. Illustration of the chaotic dynamics of the Lorenz system (14) with parameter values $\sigma = 10$, $\rho = 28$ and $\beta = 8/3$. (a) Two simulations for the Lorenz system starting from two slightly different initial conditions $(x, y, z) = (-10, 10, 25)$ and $(\tilde{x}, \tilde{y}, \tilde{z}) = (-10.01, 10, 25)$. The Lorenz attractor has a positive Lyapunov exponent and the trajectories diverge from each other. (b) Representation of the trajectory of $(x, y, z) = (-10, 10, 25)$ in the phase space. The shape of the attractor resembles a butterfly.

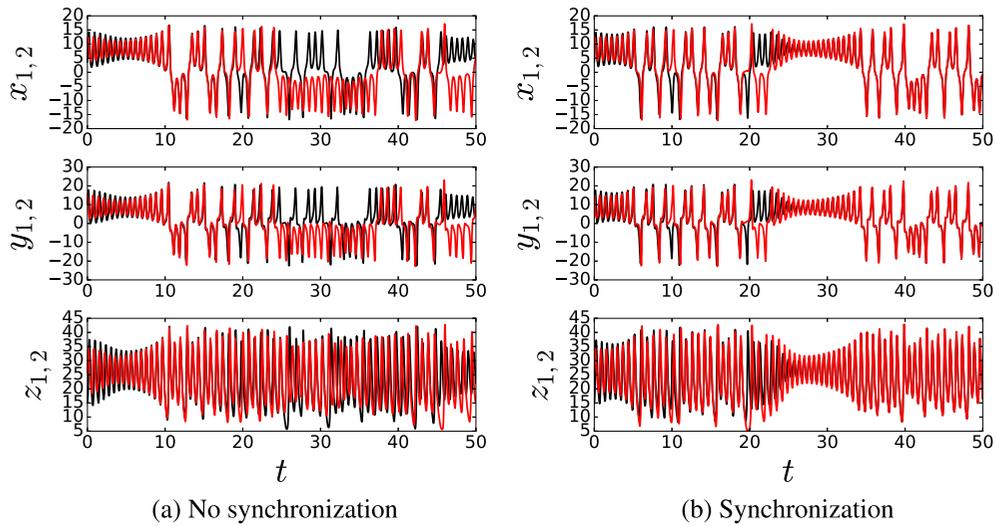


Figure 2. Comparison of trajectories of two initial conditions for the system of two coupled Lorenz systems. The critical transition coupling is $\alpha_c \approx 0.453$ for the classical parameters. The initial conditions are selected as $(x_1, y_1, z_1) = (3, 10, 15)$ and $(x_2, y_2, z_2) = (10, 15, 25)$. (a) When $\alpha = 0.4 < \alpha_c$, there is no synchronisation. (b) If $\alpha = 0.5 > \alpha_c$ one observes synchronisation of trajectories.

$$E = \frac{1}{T} \int_{t=0}^T \|\mathbf{x}_1(t) - \mathbf{x}_2(t)\| dt \quad (15)$$

In Figure 3(a) we present a synchronisation diagram where we plot E against the coupling strength α . We observe a good correspondence with the derived value of α_c . The synchronisation error depends on initial conditions so that we compute the synchronisation diagram via averaging over some realisations.

Examples on different coupling functions. It is worth mentioning that this above analysis works when the coupling

adds a damping term αz in other words when $\mathbf{H} = \mathbf{I}$. Indeed, the damping term in general form is $\alpha \mathbf{H} z$ and in this case the above results can no longer be applied. Therefore the synchronisation depends on the coupling function, we here just illustrate the effect on synchronisation if \mathbf{H} is chosen to be

$$\mathbf{H} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}, \quad (16)$$

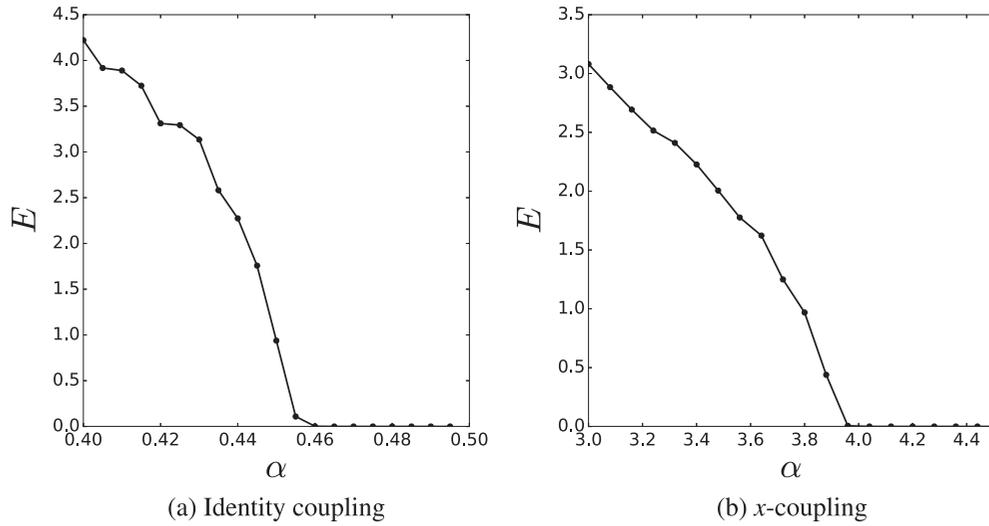


Figure 3. Synchronisation diagram of two coupled Lorenz systems, (a) with coupling matrix $\mathbf{H} = \mathbf{I}$ and (b) with coupling matrix \mathbf{H} as in Equation (16). When $\mathbf{H} = \mathbf{I}$, the observed critical coupling constant corresponds to the theoretically derived value $\alpha_c \sim 0.453$. With coupling matrix (16), synchronisation is observed to set in for coupling strengths larger than ~ 3.75 . The synchronisation error E was averaged over 300 realisations. Each realisation is simulated by a fourth-order Runge–Kutta scheme for 2000 s with 0.01 time step.

implying that the coupling arises only via the first coordinate x . The corresponding synchronisation diagram shows that the critical coupling α_c for x -coupling increases as a result, see Figure 3(b). Importantly, when \mathbf{H} does not commute with the Jacobian matrix along the trajectory, we cannot use the ansatz of Equation (10). In that case we need a different approach to derive the critical coupling, which will be discussed in Section 4 that deals with synchronisation in complex networks.

Rössler System. As a final example, we consider a system introduced by Otto Rössler in 1976:

$$\begin{aligned}\dot{x} &= -y - z, \\ \dot{y} &= x + ay, \\ \dot{z} &= b + z(x - c),\end{aligned}\quad (17)$$

where a, b and c denote parameters.

We consider two coupled Rössler systems with identity coupling function $\mathbf{H} = \mathbf{I}$ and coupling parameter α , as in Equation (6). We consider parameter values $a = 0.2, b = 0.2$ and $c = 5.7$. We numerically find that the corresponding attractor has a Lyapunov exponent $\Lambda \approx 0.071$. Hence, the expected critical coupling for synchronisation is

$$\alpha_c = \frac{\Lambda}{\lambda_2} \approx 0.0355 \quad (18)$$

This is in excellent agreement with the numerical results is shown in the synchronisation diagram Figure 4.

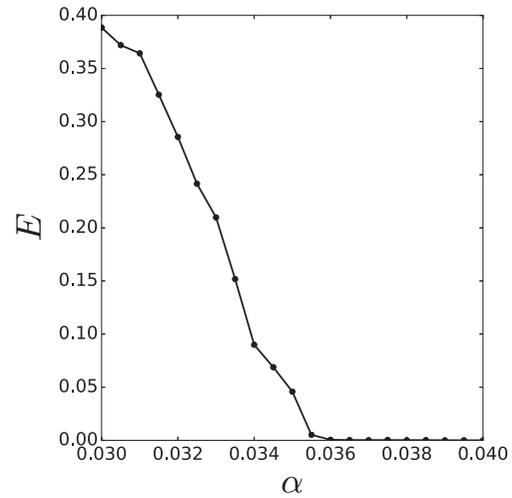


Figure 4. The synchronisation diagram of two coupled Rössler systems with the coupling function $\mathbf{H} = \mathbf{I}$. The theoretical critical coupling constant $\alpha_c \approx 0.0355$ indeed corresponds to the numerically observed one. The synchronisation error E was averaged over 300 realisations. Each realisation is simulated by a fourth-order Runge–Kutta scheme for 2000 s with 0.01 time step.

2.2.1. CS in driven systems

Another possibility is that we use certain sets of variables to drive a subsystem. For appropriate choices, we can observe synchronisation [30]. We illustrate this scheme in the Lorenz system where x -component can be driving signal of another identical system Figure 5.

In this scheme, we consider the variable x for the master the same as in the slave. That is, the x -variable of

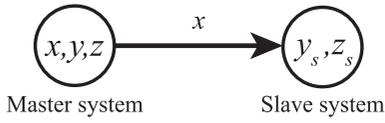


Figure 5. Master–Slave configuration where, x -variable is made identical to the response and thereby it drives the response subsystem.

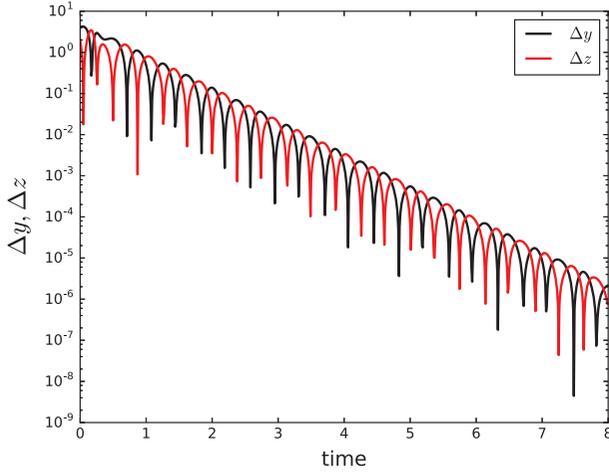


Figure 6. Simulation of master–slave type of coupling.

the master is fully replaced to the x variable in the slave

$$\begin{aligned} \dot{x} &= \sigma(y - x) \\ \dot{y} &= x(\rho - z) - y & \dot{y}_s &= x(\rho - z_s) - y_s \\ \dot{z} &= -\beta z + xy & \dot{z}_s &= -\beta z_s + xy_s \end{aligned} \quad (19)$$

where (x, y, z) are the states of the master system and (y_s, z_s) are the states of the slave system. In order to check the behaviour of the trajectories, we track the simultaneous variation of the trajectories by $\Delta_y(t) = y(t) - y_s(t)$ and $\Delta_z(t) = z(t) - z_s(t)$. For given initial conditions $(x, y, z, y_s, z_s) = (-10.1, 10.1, 10.1, 0.1, 0.1)$, Δ_y and Δ_z goes to zero (Figure 6).

For this particular choice of subsystem it is possible to construct a Lyapunov function for the displacements $\Delta_y = y - y_s$ and $\Delta_z = z - z_s$ for x -driven system (Equation (19)). We obtain

$$\begin{aligned} \dot{\Delta}_y &= -\Delta_y - x\Delta_z \\ \dot{\Delta}_z &= x\Delta_y - \beta\Delta_z. \end{aligned} \quad (20)$$

Next consider the Lyapunov function

$$V = \frac{1}{2} \left(\Delta_y^2 + \Delta_z^2 \right),$$

and along solutions of the subsystem we obtain $\dot{V} = \Delta_y \dot{\Delta}_y + \Delta_z \dot{\Delta}_z$, after some manipulation we obtain

$$\dot{V} = -\Delta_y^2 - \beta\Delta_z^2.$$

Since V is positive and \dot{V} negative Δ_y and Δ_z will converge to zero. So, the slave subsystem will have the same dynamics as the master.

2.3. Phase synchronisation

If there are small mismatches between the systems another type of synchronisation can appear for very small coupling strengths: *Phase Synchronisation* (PS) – which corresponds to a locking of phases of chaotic oscillators

$$|m\phi_1(t) - n\phi_2(t)| < C$$

where ϕ is the phase of the chaotic oscillators, m, n and C are constants. When this holds we have phase synchronisation between the two systems [31,32]. We are considering the phases on the lift, that is, diverging steadily as opposed to consider the phase mod 2π . The phase difference won't be precisely zero because of the chaotic nature of the system. We could consider higher relations of phase locking, however, the higher the relation $m : n$ more difficult is to observe the phase synchronisation. Therefore, our examples will be for 1 : 1 phase synchronisation.

Phase synchronisation is also vast research periodic oscillators [33–36]. In this case, the phases may be perfectly locked. If we are considering periodic oscillators the phase reduction approach will lead to a description of the interaction in terms of the phases alone [37]. The simplest equation in this setting is

$$\dot{\phi}_{1,2} = \omega_{1,2} + \alpha \sin(\phi_{2,1} - \phi_{1,2})$$

where ϕ is the phase along the periodic orbit. Introducing the phase difference $\Phi = \phi_1 - \phi_2$ and $\Delta = \omega_1 - \omega_2$ we obtain

$$\dot{\Phi} = \Delta - 2\alpha \sin \Phi$$

this equation has a stable fixed point $\Phi = \phi_1 - \phi_2 =$ constant if $\alpha > \alpha_c = |\Delta|/2$.

For a chaotic oscillator if coupling strength is small, the amplitudes will remain chaotic but the phase difference will be bounded. Though, it will oscillate as a result of the coupling to the amplitude. In general, it is not straightforward to introduce a phase for a chaotic attractor [38–42]. For a suitable class of attractors it is possible to define a phase in a useful way.

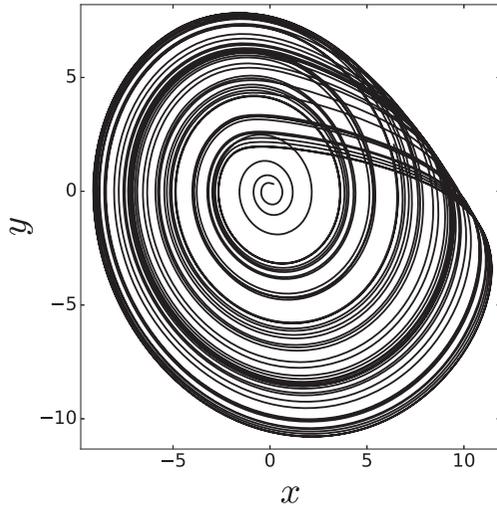


Figure 7. Projection of the Rössler attractor on $x - y$ plane for $a = 0.165$.

We focus on coupled two nonidentical Rössler oscillators, the equation is given by

$$\begin{aligned}\dot{x}_{1,2} &= -\omega_{1,2}y_{1,2} - z_{1,2} + \alpha(x_{2,1} - x_{1,2}) \\ \dot{y}_{1,2} &= \omega_{1,2}x_{1,2} + ay_{1,2} \\ \dot{z}_{1,2} &= b + z_{1,2}(x_{1,2} - c)\end{aligned}\quad (21)$$

where $a = 0.165$, $b = 0.2$ and $c = 10$ are the constants of the Rössler system. ω is the mismatch parameter to make the oscillators *nonidentical* and given as $\omega_{1,2} = \omega_0 \pm \Delta$ where $\omega_0 = 0.97$ and $\Delta = 0.02$. α is the coupling constant, the system is coupled over x components (x -coupling). For certain values of the parameter a , the projection of the attractor on $x - y$ plane resembles a limit cycle and the trajectories rotates around the origin (see Figure 7), and phase and amplitudes are given by

$$\phi_{1,2} = \arctan\left(\frac{y_{1,2}}{x_{1,2}}\right) \quad (22)$$

$$A_{1,2} = \sqrt{x_{1,2}^2 + y_{1,2}^2}. \quad (23)$$

We consider the phase on the lift (growing in time without taking the mod).

To gain insight on the adjustment of rhythm leading to phase synchronisation, we analyse the average frequencies defined as

$$\Omega_{1,2} = \lim_{T \rightarrow \infty} \frac{\phi_{1,2}(T) - \phi_{1,2}(0)}{T}. \quad (24)$$

And the frequency mismatch is

$$\Delta\Omega = \Omega_2 - \Omega_1 \quad (25)$$

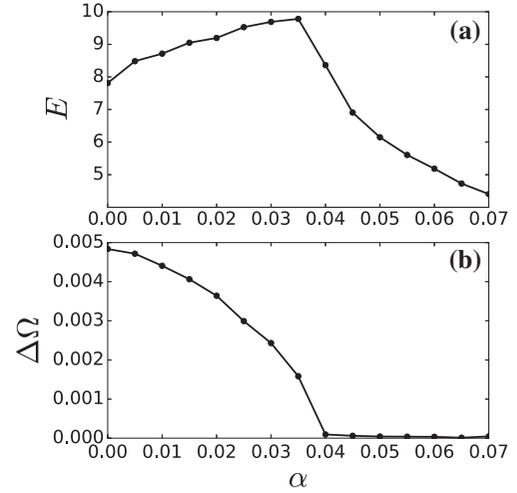


Figure 8. For a weak coupling constant: Although there is no synchrony for the difference of the amplitudes Equation (15) (top panel), there is a tendency towards to phase synchronisation while the coupling constant α increases Equation (24) (bottom panel). The synchronisation error E and frequency mismatch $\Delta\Omega$ were averaged over 300 realisations. Each realisation is simulated by a fourth-order Runge–Kutta scheme for 2000 s with 0.01 time step.

When phase synchronisation occurs $|\phi_1(t) - \phi_2(t)| \leq C$, the average frequency is the same $\Delta\Omega = 0$. The phase difference will not be tend to a constant as the phase nature of the amplitudes acts as a noise in the phases causing mismatches. The comparison of the amplitude difference (Equation (15)) and the phase (Equation (25)) is given in Figure 8. If we increase the coupling constant α .

An approximate theory of phase synchronisation can be obtained by averaging [43]. We write the model Equation (21) in terms of the phase Equations (22) as

$$\dot{\phi}_{1,2} = \frac{x_{1,2}\dot{y}_{1,2} - y_{1,2}\dot{x}_{1,2}}{A_{1,2}^2} \quad (26)$$

In this form, using polar coordinates we have $x_{1,2} = A_{1,2} \cos \phi_{1,2}$ and $y_{1,2} = A_{1,2} \sin \phi_{1,2}$, and using this representation in Equation (26), we obtain

$$\begin{aligned}\dot{\phi}_{1,2} &= \omega_{1,2} + a \sin \phi_{1,2} \cos \phi_{1,2} + \frac{z_{1,2}}{A_{1,2}} \sin \phi_{1,2} \\ &\quad - \alpha \left(\frac{A_{2,1}}{A_{1,2}} \cos \phi_{2,1} \sin \phi_{1,2} - \cos \phi_{1,2} \sin \phi_{1,2} \right)\end{aligned}\quad (27)$$

The idea now is that since the mismatch is small, both phases behave nearly the same. So we can split the dynamics of the phases as an overall increasing trend $\omega_0 t$ and a slow phase dynamics θ . This split is very clear in Figure 14. So, we write

$$\phi_{1,2} = \omega_0 t + \theta_{1,2},$$

To obtain an equation for θ (simpler than the one for ϕ) we use the fact that θ is a slow variable. That is, while $\omega_0 t$ grows a lot θ is nearly constant. Hence, we will average out the contribution of $\omega_0 t$. So we average the phases over $\omega_0 t$ over a period $\frac{2\pi}{\omega_0}$ and keep $\theta_{1,2}$ fixed. After some laborious manipulation we obtain

$$\frac{d}{dt}(\theta_1 - \theta_2) = 2\Delta - \frac{\alpha}{2} \left(\frac{A_2}{A_1} + \frac{A_1}{A_2} \right) \sin(\theta_1 - \theta_2) \quad (28)$$

Both amplitudes $A_{1,2}$ depend on time and display a chaotic behaviour. Lets assume for a moment that they are constant. Then for the phase locking of the Rössler systems, $\frac{d}{dt}(\theta_1 - \theta_2) = 0$, the equation has a stable fixed point,

$$\theta_1 - \theta_2 = \arcsin \frac{4\Delta A_1 A_2}{\alpha(A_1^2 + A_2^2)}. \quad (29)$$

This fixed point only exists when the argument of the arcsin has modulus less than 1. Therefore, we obtain the critical transition coupling

$$\alpha_c \approx 2\Delta.$$

For the given parameters ($\Delta = 0.02$) we find $\alpha_c \approx 0.04$, in agreement with the numerical analysis Figure 8. The chaotic behaviour of the amplitudes leads to fluctuations of the phases around the stable fixed point, and so the phases different will not be identically zero. Close to the critical coupling strength the frequencies exhibit a critical behaviour $\Delta\Omega \propto |\alpha - \alpha_c|^{1/2}$ as observed in Figure 8.

2.4. Generalised synchronisation

When the interacting systems are different, either because of a large parameter mismatch or the systems have distinct dynamics, these two can still exhibit synchronisation in a generalised sense. *Generalised Synchronisation* (GS) can be observed in mutually coupled systems as well as unidirectionally coupled system [30,44–46]. Surprisingly, GS can be a mapped to a complete synchronisation (CS) problem!

Here we will focus on the dynamics of unidirectionally coupled systems. The master x and the slave y systems coupled as

$$\begin{aligned} \dot{x} &= f(x) \\ \dot{y} &= g(y, h(x)) \end{aligned} \quad (30)$$

where $x \in \mathbb{R}^n$, $y \in \mathbb{R}^m$ and $h(x)$ is the coupling. For certain coupling strengths, the dynamics of system y is totally determined by the dynamics of system x . That

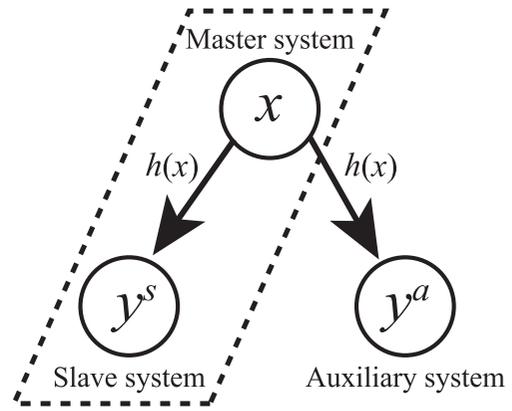


Figure 9. Scheme of the auxiliary system approach for the generalised synchronisation. Originally we only have the system in the dashed line box which is master–slave system as in Section 2.2.1. Then we add an auxiliary (helper) system y^a . If there is CS between y^s and y^a , then the GS occurs between x and $y^{a,s}$.

is, the solutions of, say x can be mapped into solutions of y .

$$y = \psi(x)$$

where ψ is a function from the phase space of the system x to the phase space of system y . When this happens we have generalised synchronisation between these two systems. CS is a particular case of GS when ψ is the identity.

To detect a functional relation between two systems in generalised synchronisation, Rulkov and co-workers proposed a technique called mutual false nearest neighbours [47]. The main idea is the see how nearby points are mapped under the dynamics. By studying the properties of nearby points one can infer the existence of the mapping ψ . Here, we focus on another approach that turns the GS problem into a CS problem. This is the auxiliary system approach [48,49]. The master system drives the slave system and an auxiliary system (copy of the slave). If the two copies of the slave exhibit CS then the master and slave are in GS [48,49]. An illustration of this scheme can be found in Figure 9.

Necessary conditions for the occurrence of GS for the system given by Equation (31) is introduced by Kocarev and Parlitz as following: for all $(x_0, y_0) \in B$, where x_0 and y_0 are states for the master–slave systems at time $t = 0$ and B is the basin where all the trajectories approach to a manifold

$$M_\psi = \{(x, y) : y = \psi(x)\}.$$

If M_ψ is attractive, different trajectories of the slave system will converge to the trajectory lying in M and it is determined only by x . In other words, if the master drives a slave y_0^s and an auxiliary (copy of slave) y_0^a systems simultaneously, the driven ones must be completely synchronised $\forall y_0^s, y_0^a \in B_y$ we have

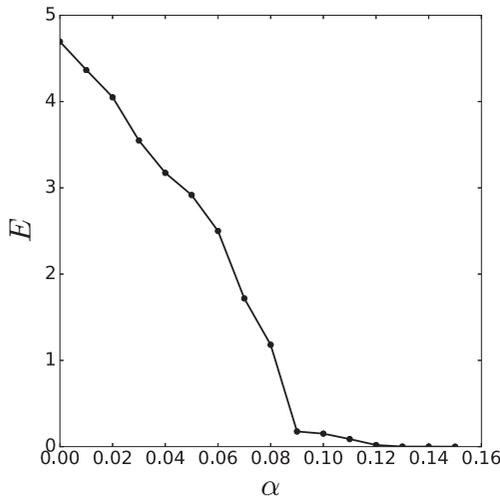


Figure 10. Generalised synchronisation: averaged over 300 realisations, time = 4000 and time step = 0.01.

$$\lim_{t \rightarrow \infty} \|y_t^s - y_t^a\| = 0.$$

Example: Consider two identical Rössler systems (Equation (17)) with the parameters ($a = 0.2$, $b = 0.2$ and $c = 5.7$) are driven by a Lorenz system (Equation (14)) with the classical parameters ($\rho = 28$, $\sigma = 10$ and $\beta = 8/3$) via x -components. We used the auxiliary system approach to detect the critical coupling for GS. Indeed, numerical results showed that $\alpha_c \approx 0.12$ as seen in the synchronisation diagram Figure 10. For given $\alpha = 0.06$ CS is not observed between the slave systems therefore there is no GS between master and slave systems as well Figure 11(a). For a coupling constant larger than the critical one $\alpha = 0.2$ the slave and the auxiliary system display CS. Hence GS can be observed between master–slave system Figure 11(b).

2.4.1. Generalised synchronisation between diffusively coupled oscillators

To gain some insight on this, we will use some ideas put forward by [50,51]. This approach allows us to obtain an analytical understanding of the critical coupling associated with the transition to GS. Consider a master–slave system diffusively coupled.

$$\begin{aligned} \dot{x} &= f(x) \\ \dot{y} &= g(y) + \alpha H(x - y) \end{aligned} \quad (31)$$

where H is a positive definite matrix. We can write the slave equation as

$$\dot{y} = \bar{g}(y) + \alpha Hx$$

where $\bar{g}(y) = g(y) - \alpha Hy$. The equation then splits into contributions solely coming from the slave and the

driver. Now consider two copies of the slaves y_1 and y_2 . Because we know the system will exhibit GS when the copies of the slaves synchronise, we introduce a variable $z = y_1 - y_2$. The system will undergo GS when $z \rightarrow \mathbf{0}$. Differentiating we obtain

$$\dot{z} = U(t)z - \alpha Hz, \quad (32)$$

where we used the mean value theorem [52] to express

$$\begin{aligned} U(t)z &= g(y_1(t)) - g(y_1(t) + z(t)) \\ &= \left(\int_0^1 DG(y_1(t) + s y_2(t)) ds \right) z(t) \end{aligned}$$

Notice that for the difference z the driving term Hx disappears as it is common for both copies of the slave y_1 and y_2 . The only part of the coupling remaining is the term $-\alpha Hz$, which adds an extra damping and provides dissipation. The trivial solution of z is globally stable if the coupling is large enough. Indeed, we can construct a Lyapunov function for z . Indeed, consider

$$V(z) = \frac{1}{2} \langle z, z \rangle,$$

and differentiating the Lyapunov function along the solution $z(t)$ of Equation (32) we obtain

$$\frac{dV(z(t))}{dt} = \langle \dot{z}(t), z(t) \rangle \quad (33)$$

$$\leq (\|Dg\| - \alpha \lambda_{\min}(H)) \|y\|^2 \quad (34)$$

where $\lambda_{\min}(H)$ is the minimum eigenvalue of H . In this last passage, we used the Cauchy Schwartz inequality $|\langle U(t)z, z \rangle| \leq \|U(t)z\| \|z\| \leq \|U\| \|z\|^2$, and noticed that $\|U\| \leq \|Dg\|$.¹ We also used the fact that H is positive $\langle Hz, z \rangle \geq \lambda_{\min}(H) \|z\|^2$. Hence, for

$$\alpha > \alpha_c = \frac{\|Dg\|}{\lambda_{\min}(H)}$$

the derivative of the Lyapunov function is negative and every solution of the system sinks to zero.

What did we learn? When $\alpha > \alpha_c$ we have GS. Any two trajectories of the slave y_1 and y_2 will converge towards the same asymptotic state. This happens because the coupling terms add extra dissipation. The convergence rate is exponential $\|y_1(t) - y_2(t)\| \leq Ke^{-\eta t}$ because $\dot{V} \leq \eta V$ and η is uniform on the trajectories y_1, y_2 and x . As a conclusion, there is a function ψ such that the dynamics of the slave can be as $y = \psi(x)$.

In the literature, a typical way to estimate whether one has GS is to compute the Lyapunov exponents of the slaves. Since we are assuming that the uncoupled systems

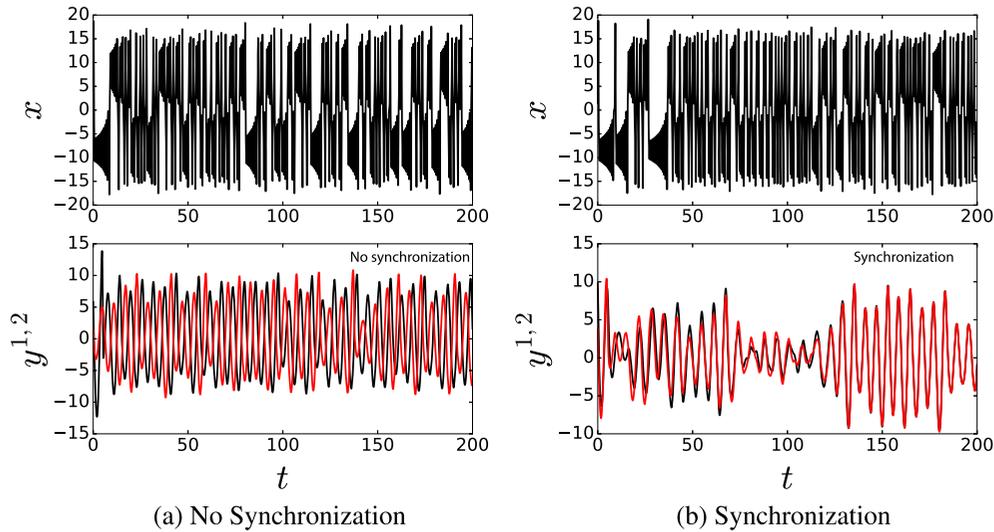


Figure 11. Two simulations for the generalised coupling scheme: a Lorenz system x drives two Rössler systems $y^{s,a}$. The critical transition coupling is $\alpha_c \approx 0.12$. (a) For the coupling constant $\alpha = 0.06$, there is no synchronisation since $\alpha < \alpha_c$ (b) for $\alpha = 0.2$ synchronisation is obtained since $\alpha > \alpha_c$.

are chaotic, for $\alpha = 0$ the slave will have a positive Lyapunov exponents. As we increase α the maximum Lyapunov exponent may become negative for a value α_c . We use this α_c as an estimate for the critical coupling for GS.

2.5. Summary of synchronisation types

We discuss the three cases commonly found in applications. A schematic representation of the cases is found in Figure 12.

Complete synchronisation in identical systems. If the isolated dynamics are identical, $f_1 = f_2$ and diffusively coupled, hence, the subspace $x_1 = x_2$ is invariant under Equation (6). Indeed, the coupling vanishes and both systems will oscillate in unison for all coupling strengths α and all times. Such collective motion is called complete synchronisation (CS). The question is whether CS is attractive, that is, if the oscillators state are nearly the same $x_1(0) \approx x_2(0)$, will they synchronise? Meaning that

$$\lim_{t \rightarrow \infty} \|x_1(t) - x_2(t)\| = 0.$$

See Figure 13 for an illustration. In Section 2.2, the CS was discussed in detail.

Phase synchronisation (PS) when $f_1 \approx f_2$. In this situation the subspace $x_1 = x_2$ is not invariant. And each system will have its own frequency given by their phase dynamics $\phi_{1,2}$. For small coupling strengths the phases can be locked $\phi_1 \approx \phi_2$ while the amplitudes remain uncorrelated Figure 14. This phenomenon is called phase synchronisation. Typically, the critical coupling for PS is

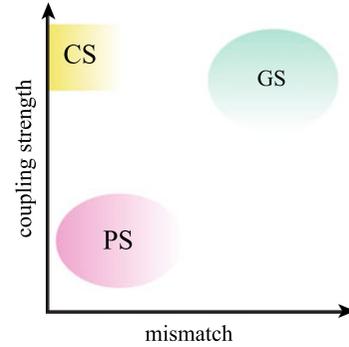


Figure 12. Diagram of synchronisation types for diffusively coupled systems. The horizontal axis depicts the mismatch between the isolated dynamics (f_1 and f_2) and the vertical axis the coupling constant. The diagram shows the typical balance between mismatch and coupling strength and to achieve a certain kind of synchronisation. Complete synchronisation (CS) occurs for identical chaotic systems ($f_1 = f_2$) and large enough coupling strengths. Phase synchronisation (PS) is observed between slightly different systems for small coupling strengths. Generalised synchronisation (GS) is a result of master–slave system and can occur for large mismatch parameters or even between distinct systems when the coupling strength large enough.

proportional to the mismatch $f_1 - f_2$, as illustrated in Figure 12. Further details were given in Section 2.3.

Generalised synchronisation in master–slave configurations. If the vector fields are different $f_1 = f$ and $f_2 = g$, the systems can synchronise, but in a generalised sense. We consider systems coupled in a master–slave configuration. For certain coupling strengths, the dynamics of the master x can determine the dynamics of the slave y , that is $y = \psi(x)$, see Figure 15. This is called *Generalised Synchronisation* (GS). Further details for GS was given in Section 2.4.

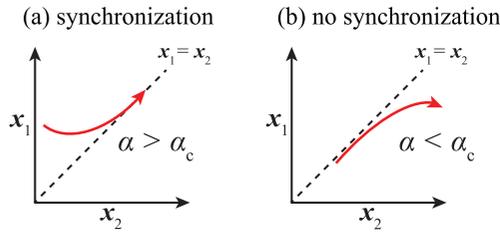


Figure 13. Illustration of complete synchronisation. (a) If the coupling strength is large enough ($\alpha > \alpha_c$), the systems converge to invariant synchronisation manifold ($x_1 = x_2$), (b) otherwise ($\alpha < \alpha_c$), they diverge hence no synchronisation.

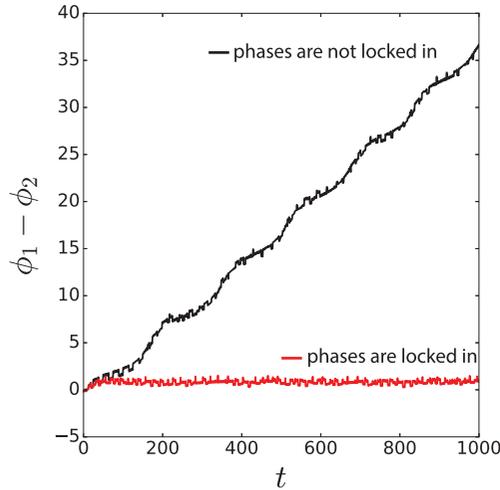


Figure 14. Illustration of phase synchronisation for two coupled slightly different and chaotic systems ($f_1 \neq f_2$). The evolution of the phase differences between the systems for two different coupling constants $\alpha < \alpha_c$ and $\alpha > \alpha_c$.

2.6. Historical notes

Studies on synchronisation dates back to Christiaan Huygens who studied coupled pendulums. In this case, the pendulums are periodic and have distinct frequencies, but due to interaction they adjust their rhythm. In the seventies, thanks to the works of Winfree [53] and Kuramoto [37] the area experienced a boom. In early 2000's many excellent books and reviews were devoted to this subject [1,33,34,54–56].

Chaotic synchronisation on the other hand is younger. To begin with, the establishment and full acceptance of the chaotic nature of dynamics is fairly new [57]. The role of chaos in nature was object of intense debate in the seventies when Ruelle and Takens proposed that turbulence was generated by chaos.

The chaotic dynamics can be fairly complicated. Typically, the evolution never repeats itself, nearby points drift apart exponentially fast, but in the long run the dynamics return arbitrarily close to its initial state. Such dynamics is so unpredictable that modern approach tackles it from a probabilistic perspective.

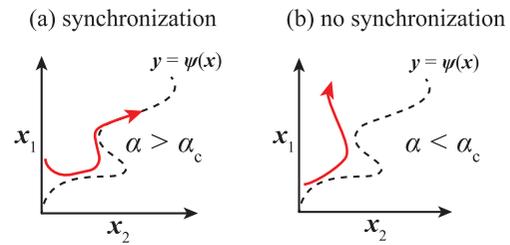


Figure 15. Illustration of generalised synchronisation. If the coupling strength is large enough, a functional relationship ($y = \psi(x)$) exhibits between the dynamical variables x_1 and x_2 . (a) If $\alpha > \alpha_c$, the generalised synchronisation is observed (b) otherwise $\alpha < \alpha_c$, there is no generalised synchronisation.

Given this complexity, many researchers thought it is unlikely one could possibly synchronise two chaotic systems. How could a system with exponential divergence of nearby trajectories have a state were trajectories come together while keeping their chaotic nature? That seemed paradoxal. Chaos and synchronisation should not come together. This view was proven wrong in the late eighties. In fact, we have come to think it as rather natural. Funny enough, before this view was accepted synchronisation of chaos had to be rediscovered a few times.

Back in the eighties, Fujisaka and Yamada had the first results on synchronisation of chaos [23,58,59]. They publish it in Japan, but their results went fairly unnoticed in the west. Just two years later mathematicians and physicists from Nizhny Novgorod exposed many of the concepts necessary for analysing synchronous chaos [24]. This paper is now famous, but back then it also went largely unnoticed.

Only some years later the study of synchronisation of chaos had its boom, largely as a result of the works by Pecora and Carroll [30]. Lou Pecora and co-workers went systematically tackling two coupled systems and then moved on to study chaotic systems coupled on periodic lattices [60,61]. These early results were relying on ideas from Nuclear physics to diagonalise the lattice and stability theory (the Lyapunov methods) to analyse synchronisation.

The nineties proved prolific for synchronisation! Two groups proposed an extension of synchronisation, the so-called generalised synchronisation [47–49,62,63]. Generalised synchronisation *only* asked for a functional relationship between the states, that is, the dynamics of one system is fully determined by the dynamics of the other. Also in the mid nineties, Rosenblum, Pikovsky and Kurths put forward the concept of chaotic phase synchronisation. Here two nearly identical chaotic oscillators can have their phase difference bounded while the amplitudes remain uncorrelated [31,32].

A few years down the road, Pecora and Carroll were able to generalise their approach to undirected networks

of diffusively coupled systems [64]. They also wrote a review about their approach [65]. These results open the door to the understanding of the role of the linking structure on the stability of synchronisation. Barahona and Pecora [66] showed that small-world networks are easier to globally synchronise than regular networks. Motter and coworkers [67–70] showed that heterogeneity in the network structure may hinder global synchronisation. On the other hand, Pereira showed such heterogeneity may enhanced synchronisation of highly connected nodes [71].

3. Applications

In this section, we discuss the role of synchronisation phenomena in various applications including secure communication approaches, parameter estimation of a model from data and prediction.

3.1. Secure communication based on complete synchronisation

The first approach is to send an *analog* message [72]. The key idea is the following: the sender adds the message $m(t)$ on a chaotic signal $x(t)$ and generate a new signal $s(t) = m(t) + x(t)$ (Figure 16). The assumption is that the amplitude of $x(t)$ is much larger than the amplitude of $m(t)$. This method is called the masking information on bearing signals. Because chaotic signals are noise-like and broadband (have many frequencies), it is difficult to read the message. One could then retrieve the message using synchronisation.

Masking of messages on bearing signals does not require encryption. Here is the keystone is selection of the transmitter and receiver systems such that they synchronise. They are also assumed to be identical (this means that the receiver knows the parameters of the transmitter). One can retrieve the message if the parameters are known. So, the parameters play a role of encryption key.

We illustrate this communication scheme using the Lorenz system Equation (14). The Lorenz system has the particularity that it divided into subsystems (x, z) and (y, z) . We can use the variables x or y of a subsystem to drive the other subsystem. In this driving setting, the synchronisation between driver and slave is exponentially stable provided that the parameters σ , ρ and β are identical. Here, we have chosen x component to act as a driver. The message $s(t)$ drives the receiver system as

$$\begin{aligned}\dot{x}_r &= \sigma(y_r - x_r) \\ \dot{y}_r &= s(\rho - z_r) - y_r \\ \dot{z}_r &= -\beta z_r + sy_r.\end{aligned}\quad (35)$$

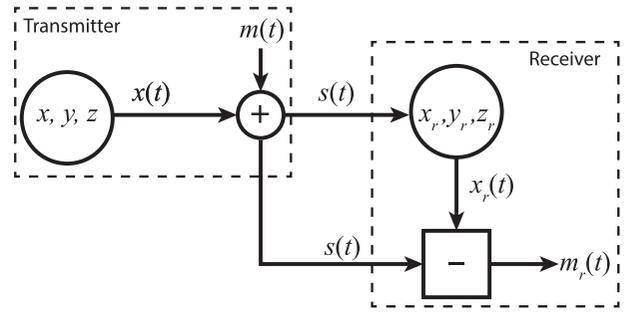


Figure 16. Illustration of the message masking on bearing signal scheme. Transmitter generates a chaotic signal $x(t)$ and add the message $m(t)$ on it. This combined two signals $s(t) = x(t) + m(t)$ is sent to the receiver and both systems synchronise. Discarding the synced signal x_r from the s , the message $m_r \sim m$ is restored.

Since the synchronisation of chaotic systems is exponentially stable for such system [30,73] under low amplitude of noise the synchronisation (coherence) still occurs. Then the chaotic signal x_r can be obtained from Equation (35). Therefore the message can be regenerated by $m(t) = s(t) - x_r(t)$ (Figure 16).

As a message we use the signal

$$m(t) = 0.1 \frac{\sin(1.2\pi \sin^2(t))}{\pi \sin^2(t)} \cos(10\pi \cos(0.9t)) + \xi$$

where ξ is a white noise Figure 17(a). We attach this message on x -component of the Lorenz system Equation (14) with parameters $\sigma = 16.0$, $\rho = 45.2$ and $\beta = 4.0$ then the information is masked on bearing signal s Figure 17(b). By synchronisation we restore the message $m_r \sim m$ (Figure 17(c)). This secure communication application is also experimentally demonstrated using Chua's circuits [74,75] and Lorenz-like circuit [76].

The second approach is the modulation of the parameters for the *digital* communication. In this case, the message $m(t)$ only carries binary-valued signals. The setup is similar to the masking approach but the message is included in the transmitter parameters. The transmitter system has an adjustable parameter $\sigma_a(t) = \sigma + \delta m(t)$ such that we can tune the system into synchronisation when $m(t) = 0$ and out synchronisation when $m(t) = 1$ Figure 18. We retrieve the message $m(t)$ by the synchronisation and desynchronisation pattern.

The dynamics of transmitter and the receiver is given by

$$\begin{aligned}\dot{x}_1 &= \sigma_a(y_1 - x_1) \\ \dot{y}_1 &= x_1(\rho - z_1) - y_1 \\ \dot{z}_1 &= -\beta z_1 + x_1 y_1 \\ \dot{x}_2 &= \sigma(y_2 - x_2) \\ \dot{y}_2 &= x_2(\rho - z_2) - y_2 \\ \dot{z}_2 &= -\beta z_2 + x_2 y_2\end{aligned}$$

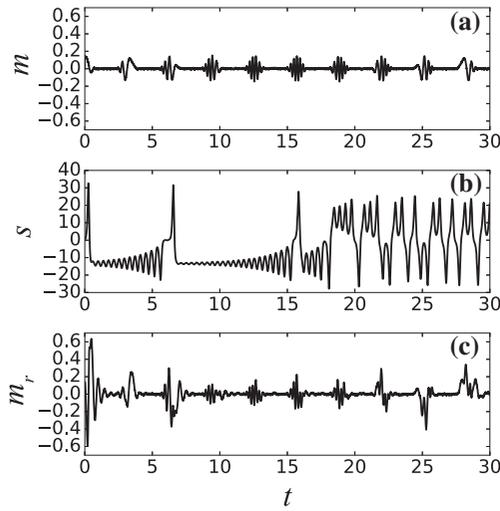


Figure 17. The masking an analog message on bearing chaotic signal. (a) The low-amplitude message, (b) the message embedded into high-amplitude chaotic signal and (c) restored message from the transmitted signal.

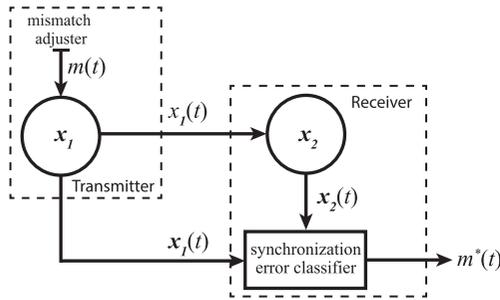


Figure 18. A secure communication scheme: hiding a digital message on a chaotic signal. Changing a parameter of transmitter causes different level of synchronisation errors between the transmitter and the receiver. The amplitude of the error E brings the message out.

where $\sigma_a = \sigma + \delta m(t)$ is the adjustable parameter. Again the key aspect is having mismatch between the transmitter and the receiver does not allow systems to synchronise. Modulating the parameter σ_a by the message $m(t)$ we can produce different levels of synchronisation errors. Choosing the parameters of the transmitter and the receiver identical gives the synchronisation error $E \sim 0$ (CS), this can be assigned binary 0 by $m(t) = 0$. The large mismatch δ causes a certain amount of synchronisation error $E > 0$, this can be assigned binary signal 1 by $m(t) = 1$. Then the digital communication can be set between sender and receiver [76,77].

Using same parameters as in the previous application ($\sigma = 16.0$, $\rho = 45.2$ and $\beta = 4.0$), we illustrate this digital communication. For this example, a digit of the message is set for 10 time units and the message is 0101010101 (Figure 19(a)). For each message time, we change σ_a from σ to $\sigma + \delta$ and other way round. Then the

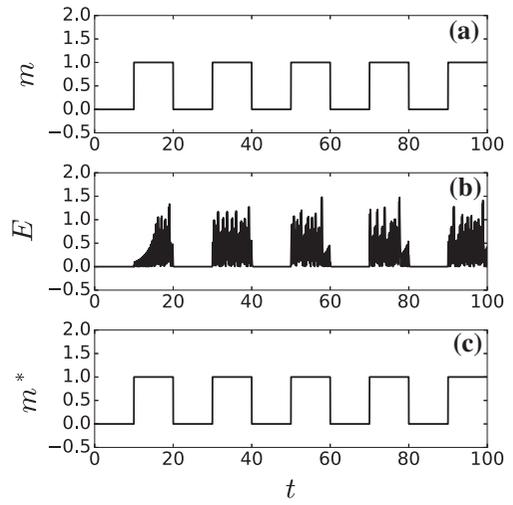


Figure 19. Manipulating the parameter of a transmitter allows digital secure communication. In this application 10 time units are used for a single digit. (a) Digital message (0101010101), (b) the synchronisation error and (c) restored message.

synchronisation error E varies according to this change (Figure 19(b)). Due to change in the E , the message is restored (Figure 19(c)).

There are more communication applications using the synchronisation mechanism e.g. using hyperchaotic systems [61,63,78,79] or volume-preserving maps [80]. The common idea of all these given approaches is the CS phenomena, negative conditional Lyapunov exponent between the systems are needed to exhibit of the synchronisation [30,73].

3.2. Secure communication based on phase synchronisation

Security is an important issue for communication approaches. As might be expected, some methods were improved and reported to break the CS-based communication schemes [81–84]. Then more secure communication scheme demonstrated by means the PS [85].

The scheme based on the PS possesses three chaotic Rössler systems ($x_{1,2,3}$). The transmitter of the scheme consists of two weakly coupled identical systems x_1 and x_2 over their x -components Equation (36) and the receiver x_3 has slightly different dynamics. In this case, we couple the systems with using their phases Equation (36) as presented in Ref. [86]. The phase definition for Rössler system is given by Equation (22). The mean of two systems' phases ϕ_1 and ϕ_2 in transmitter can be used as a spontaneous phase signal ϕ_m to couple the third system as in Equation (36). As distinct from the CS-based schemes, we have three systems and the reason behind these to improve the security. The return maps of the phase ϕ_m is way more complex than ϕ_1 (or ϕ_2), this makes to break dynamics not trivial [85].

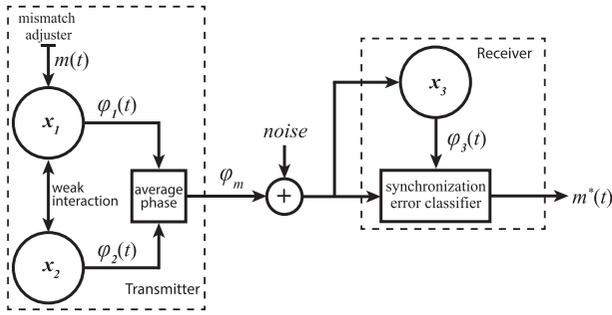


Figure 20. A secure communication scheme by phase synchronisation: hiding a digital message on a chaotic signal.

We illustrate this application by

$$\begin{aligned}
 \dot{x}_{1,2} &= -(\omega + \Delta\omega)y_{1,2} - z_{1,2} + \epsilon(x_{2,1} - x_{1,2}) \\
 \dot{y}_{1,2} &= x_{1,2} + ay_{1,2} \\
 \dot{z}_{1,2} &= b + z_{1,2}(x_{1,2} - c) \\
 \dot{x}_3 &= -y_3 - z_3 + \alpha(r_3 \cos \phi_m - x_3) \\
 \dot{y}_3 &= x_3 + ay_3 \\
 \dot{z}_3 &= b + z_3(x_3 - c)
 \end{aligned} \tag{36}$$

where constant $\omega = 1$ and standard parameters of the Rössler system $a = 0.15$, $b = 0.2$ and $c = 10.0$. Coupling constants $\epsilon = 5 \times 10^{-3}$ is between x_1 and x_2 , and α is between the transmitter and the receiver. r_3 is the amplitude of the response system given by Equation (23). $\Delta\omega$ is the adjustable mismatch parameter, for this illustration we select

$$\Delta\omega = \begin{cases} 0.01 & \text{if bit digit} = 1 \\ -0.01 & \text{if bit digit} = 0. \end{cases}$$

Similar to digital communication by the CS (see Section 3.1), the modulation of parameters in the transmitter would hide a binary message $m(t)$ on ϕ_m . The PS will exhibit between ϕ_m and ϕ_3 . Due to the changes on the adjustable control parameters $\Delta\omega$, the phase difference between ϕ_m and ϕ_3 varies. In other words, the phases are locked on different phase shifts. The message can be retrieved from different phase locking values (Figure 20).

Because of the weak coupling, the CS never occurs Figure 21(a). Every 10 time unit we switch $\Delta\omega$ parameter to create a digital message $m(t)$ (010101...) Figure 21(b). The hidden message on chaotic signal can be restored from the receiver using the phase difference between ϕ_m and ϕ_3 . If the message digit is 0, then the phase difference oscillates most time below 0, otherwise above 0 Figure 21(c). Therefore it is easy to restore associated message m^* Figure 21(d).

In real world examples, it is almost not possible to create identical systems, and the noise is always an issue

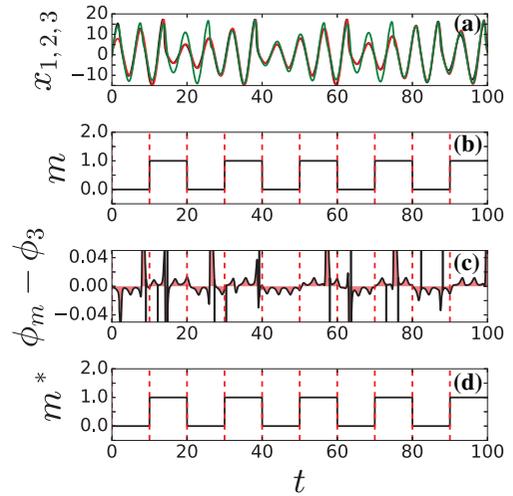


Figure 21. An illustration for the secure communication by phase synchronisation: hiding a digital message on a chaotic signal.

to deal. The phase locking can be still preservable under effect of a certain level of noise (Figure 21).

3.3. Parameter estimation and prediction

Now we have data and we want to learn about the system that generated the data. Thus we will be able to predict the future behaviours and critical transitions. The determining equations of the system are known however the parameters are not. The goal is to find these unknown parameters with using synchronisation phenomena. So, we blend the data with equations. The data are then used to drive the equations. If they are coupled in a proper way (Section 2.2.1), the equations can synchronise with data. The key point is the following: if the parameters of slave system are identical with the master whose produced driving signal, then the CS exhibits (synchronisation error $E = 0$) otherwise no CS (synchronisation error $E > 0$). Therefore, it is possible to estimate the parameters by a strategy to minimise the synchronisation error $E \rightarrow 0$ such as POWELL technique [87].

We assumed that we have a limited data and we want to predict the future of the system. After the parameters are estimated, the state of the synchronised slave matches the data. Because the solution of the equations is then the same as the data, we can use the equations to predict future dynamics.

The second approach is to estimate a slave system's parameters of a master–slave system. In this case, the dynamics of master and slave is distinct. We assume that we have two given data-sets: one of them from master system and the other one is from the slave. The governing dynamics of the master–slave system is given

$$\begin{aligned}
 \dot{x} &= f(x) \\
 \dot{y} &= g(y) + h(x, y)
 \end{aligned}$$

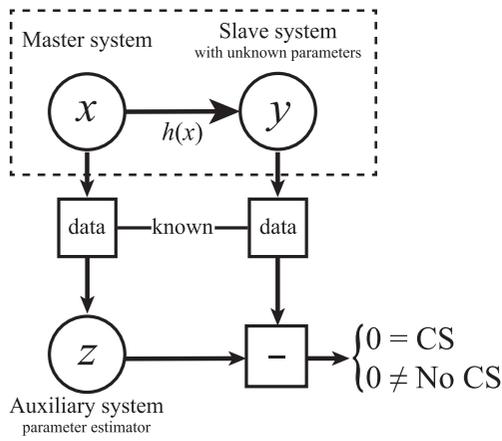


Figure 22. Only two data-sets are known from a master–slave system, into the dashed rectangle, without any info about the parameters. An auxiliary system is driven by the data from the master and measures the amplitude difference between the auxiliary and the data from the slave system. If the difference is 0, then there is a CS that means the parameters of the slave and the auxiliary are identical.

where x and y are the states of master and slave systems respectively. We aim to estimate the parameters of g . Here we cannot use y data to drive another g system directly as in previous approach since y is driven by x . If we know that master–slave system is in the GS and the coupling function h is known, then we can apply the auxiliary system approach which is the master system drives an auxiliary (copy of the slave) z Figure 22. We expect that the CS exhibits between the slave y and auxiliary z systems if the parameters are identical.

Using the GS idea, the problem turned into the CS problem. From now on, minimising the synchronisation error $E \rightarrow 0$ technique can be used to estimate the unknown parameters. Similar to the previous approach, the future of the system can be predicted as well.

Example: Consider a Lorenz system with classical parameters is driven by a Rössler system. We have only two data sets, x_1 -component of the Rössler (Equation (17)) and y_1 -component of the Lorenz (Equation (14)). Then we drive an auxiliary system z by x_1 as

$$\begin{aligned}\dot{z}_1 &= \sigma_e(z_2 - z_1) + \alpha(x_1 - z_1) \\ \dot{z}_2 &= z_1(\rho_e - z_3) - z_2 \\ \dot{z}_3 &= -\beta_e z_3 + z_1 z_2.\end{aligned}$$

The goal is to find the parameters of slave system. The spontaneous synchronisation error is

$$E(t) = \|z_1 - y_1\|. \quad (37)$$

Adjusting the parameters of $z(\sigma_e, \rho_e, \beta_e)$ we minimise the simultaneous error $E(t)$ by Powell's algorithm [87]. This

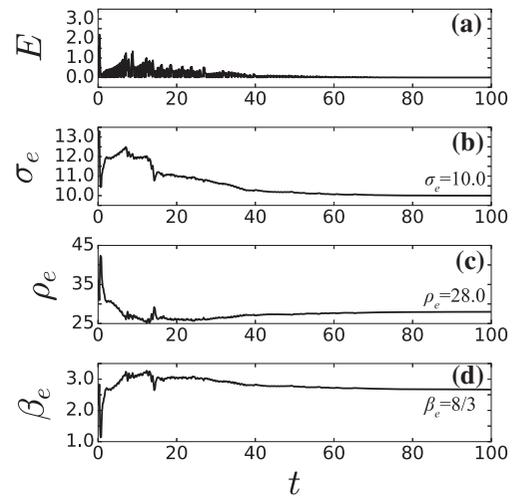


Figure 23. Illustration of parameter estimation. Standard parameters of Lorenz system.

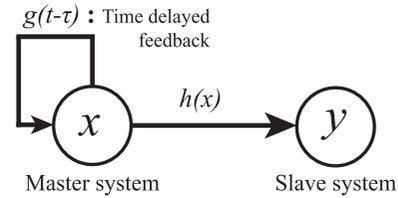


Figure 24. Scheme of anticipating synchronisation with memory in the driver systems.

method allows us to estimate the parameters of the slave system (Figure 23).

3.4. Chaos anticipation

Chaos is unpredictable but synchronisation can help predicting the state of a chaotic system ahead of time. Anticipating synchronisation (AS) is a good approach for the future prediction since the slave system synchronises with the upcoming states of the master system at time $t + \tau$ where τ is a time delay. The occurrence of AS depends on the coupling constant α . Therefore, it is not dependent on isolated dynamics or time delay τ and regarding to type of desired application higher dimensional chaotic systems can be used for an arbitrary time delay. This anticipation of chaos can be used or is used in applications such as semiconductor lasers with optical feedback, secure communications [88].

Consider two chaotic systems in a master–slave interaction and the master has a certain delay τ feedback Figure 24. Because of the internal delay feedbacks, it may well happen that the master x and slave y synchronise but with some time delay

$$x(t) = y(t - \tau)$$

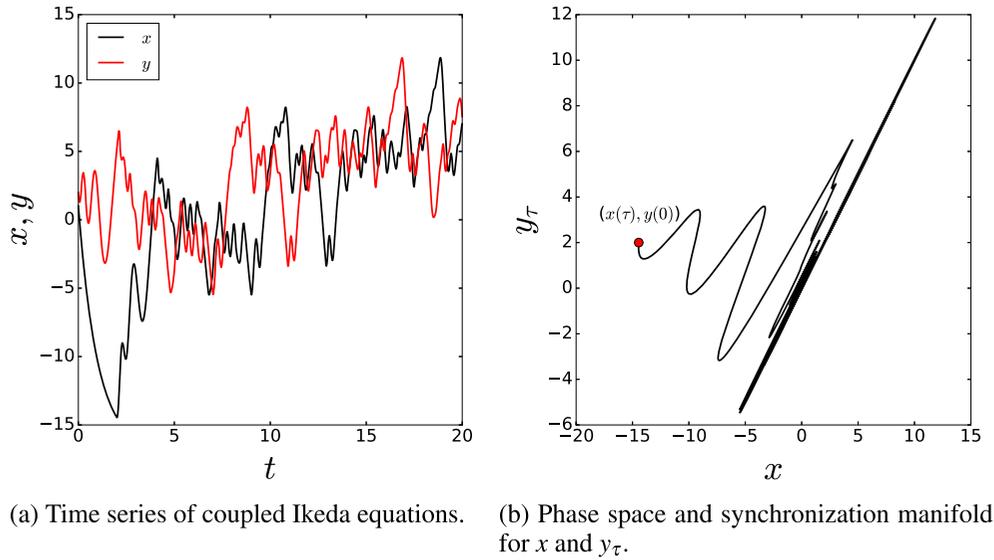


Figure 25. Anticipating chaotic synchronisation. In the beginning, the systems are not in harmony. After a transient time both systems are getting into a time-delay τ synchronisation. For this illustration $\tau = 2$. (a) time series of the systems (b) phase space and synchronisation manifold of the system. The red circle is the initial condition for the trajectory of (x, y_τ) .

When this happens we have

$$y(t) = x(t + \tau).$$

Hence, although the system x is fully chaotic, we can precisely predict its future state from the system y . In other words the slave system anticipates the master system. This kind of synchronisation is called anticipated synchronisation (AS).

Example: We consider two coupled Ikeda equations,

$$\begin{aligned} \dot{x} &= -\alpha x - \mu \sin x_\tau \\ \dot{y} &= -\alpha y - \mu \sin x \end{aligned} \quad (38)$$

We use the notation $y_\tau = y(t - \tau)$. The scheme of the system is given in Figure 24. The synchronisation error for delayed system is given by,

$$z = x - y_\tau$$

and to show that synchronisation is attractive we analyse the first variational equation

$$\begin{aligned} \dot{z} &= \dot{x} - \dot{y}_\tau \\ &= -\alpha x - \mu \sin x_\tau - (-\alpha y_\tau - \mu \sin x_\tau) \\ &= -\alpha z. \end{aligned} \quad (39)$$

The solution is $z(t) = z_0 e^{-\alpha t}$ and for $\alpha > 0$ the synchronisation is globally exponentially stable.

To illustrate AS, we simulate Equation (38) with a fourth-order Runge–Kutta integrator for the

delay-differential equations for the parameters $\alpha = 1$, $\mu = 20$ and $\tau = 2$. The simulation starts from a random initial condition. After enough transient time t , the master x and the slave y systems synchronise with a time delay (τ) Figure 25(a). The transient time can be observed from the phase space of x and y_τ . The initial condition is given by a red circle in Figure 25(b), the trajectory converges to the synchronisation manifold ($x = y_\tau$).

Example: The AS can occur for higher dimensional chaotic system. For such cases the critical coupling constant can be positive ($\alpha_c > 0$). The AS can be obtained without delayed state in the master system, that is, without memory in the master system. The scheme of this model is given in Figure 26. We can demonstrate this case with Rössler system, the governing equations are given by

$$\begin{aligned} \dot{x}_1 &= -x_2 - x_3 \\ \dot{x}_2 &= x_1 + ax_2 \\ \dot{x}_3 &= b + x_3(x_1 - c) \\ \dot{y}_1 &= -y_2 - y_3 + \alpha(x_1 - y_{1,\tau}) \\ \dot{y}_2 &= y_1 + ay_2 \\ \dot{y}_3 &= b + y_3(y_1 - c) \end{aligned} \quad (40)$$

where the parameters are $a = 0.15$, $b = 0.2$ and $c = 10$. We simulate Equation (40) for no AS ($\alpha < \alpha_c$) Figure 27 and AS ($\alpha > \alpha_c$) Figure 27 cases. In this memoryless AS approach, the synchronisation is also dependent on time delay τ .

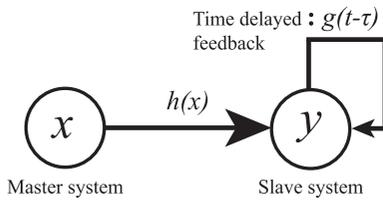


Figure 26. Scheme of anticipating synchronisation without memory in the driver systems.

4. Synchronisation in complex networks

Synchronisation is commonly found in networks of natural and mankind-made systems such as neural dynamics [89–91], cardiovascular systems [92–94], power grids [95], superconducting Joseph junctions [3]. The theory we presented in the previous chapters can be generalised to understand certain aspects of synchronisation in these complex systems.

We will focus on diffusively interacting identical oscillators, so the dynamics of the coupled system reads as

$$\frac{d\mathbf{x}_i}{dt} = \mathbf{f}(\mathbf{x}_i) + \alpha \sum_{j=1}^N A_{ij}[\mathbf{H}(\mathbf{x}_j) - \mathbf{H}(\mathbf{x}_i)], \quad (41)$$

where $\mathbf{f} : \mathbb{R}^n \rightarrow \mathbb{R}^n$ describes the dynamics of the isolated system, α is the overall coupling strength, N is the number of oscillators, $\mathbf{H} : \mathbb{R}^n \rightarrow \mathbb{R}^n$ is the coupling function. Finally, A_{ij} dictates who is interacting with whom. $A_{ij} = 1$ if i receives a connection from j and 0 otherwise. The matrix \mathbf{A} (the dimension is $N \times N$) provides the network linking structure and it is called adjacency matrix. This matrix will play a clear role in the analysis.

The network dynamics of diffusively coupled system in Equation (41) models many physical systems. A few specific examples are:

Electronic Circuits with Resistive interaction. Electric circuits, e.g. Chua, Roessler-like, Lorenz-like, can be coupled over their resistors then Equation (41) models this system [30]. Another case, only one electric circuit can be driven by an external signal as master–slave system (for details, see Section 2.2.1) [96].

Neuron Networks with Electrical Coupling. In brain network, \mathbf{f} can be the isolated neuron dynamics modelled by differential equations with chaotic or periodic behaviour and having different time-scales, that is, the isolated can have burst and single regime [97] and \mathbf{H} the electrical synapses $\mathbf{H}(\mathbf{x}_1 - \mathbf{x}_2) = (x_1 - x_2, 0, 0)$.

Stable Biological System. In biological systems when the isolated dynamics has a stable periodic motion then typically one can rephrase the network dynamics in terms of our model. For instance, the heart consists of millions

of pacemaker cells. Each cell when isolated has its own rhythm, but when put together these cells interact and behave in unison to deliver the strong electrical pulse that make our heart beat [1]. The dynamics of the pacemaker cells are modelled by phase oscillators ϕ_i with distinct frequencies ω_i and the coupling function is a simple sine function $H(\phi_1 - \phi_2) = \sin(\phi_1 - \phi_2)$ [94,98].

Laser Arrays. Lasers can be arranged arrays or complex networks. In this case, one is interested in the electrical field dynamics. Such electrical field is govern by equations with interaction akin to diffusion Ref. [99]. So, the approach presented here can be extrapolated to such lasers under slight changes.

In fact, when we considering periodic oscillators² the above model is a normal form for the networked system. That is, the isolated system has a periodic exponentially attracting orbit, we couple the system, and in the weak coupling regime, the amplitudes will change only slightly but the phases can change by large amounts. So the dynamics can be described only in terms of the phases. The phase description will again fit in our Equation (41).

Our synchronisation results given in the previous sections are exponentially stable. In other words, if once the trajectories are into the synchronisation subset, the solution is robust and persistent to the perturbations. For N coupled nonidentical systems ($\mathbf{f}_1 \neq \mathbf{f}_2 \neq \dots \neq \mathbf{f}_N$), complete synchronisation is not possible. However, because of exponentially stable solutions, highly coherent state around the synchronisation subset can be still observed [100,101].

4.1. Interactions in terms of Laplacian

Because of the diffusive nature of the interaction, it is possible to represent the coupling in terms of the Laplacian matrix \mathbf{L} . Indeed,

$$\begin{aligned} & \sum_{j=1}^N A_{ij}[\mathbf{H}(\mathbf{x}_j) - \mathbf{H}(\mathbf{x}_i)] \\ &= \sum_{j=1}^N A_{ij}\mathbf{H}(\mathbf{x}_j) - \mathbf{H}(\mathbf{x}_i) \sum_{j=1}^N A_{ij} \\ &= \sum_{j=1}^N A_{ij}\mathbf{H}(\mathbf{x}_j) - k_i\mathbf{H}(\mathbf{x}_i) \\ &= \sum_{j=1}^N (A_{ij} - \delta_{ij}k_i)\mathbf{H}(\mathbf{x}_j) \end{aligned}$$

where $k_i = \sum_{j=1}^N A_{ij}$ is the degree of the i th node, δ_{ij} is the Kronecker delta, and recalling that $L_{ij} = \delta_{ij}k_i - A_{ij}$ we obtain

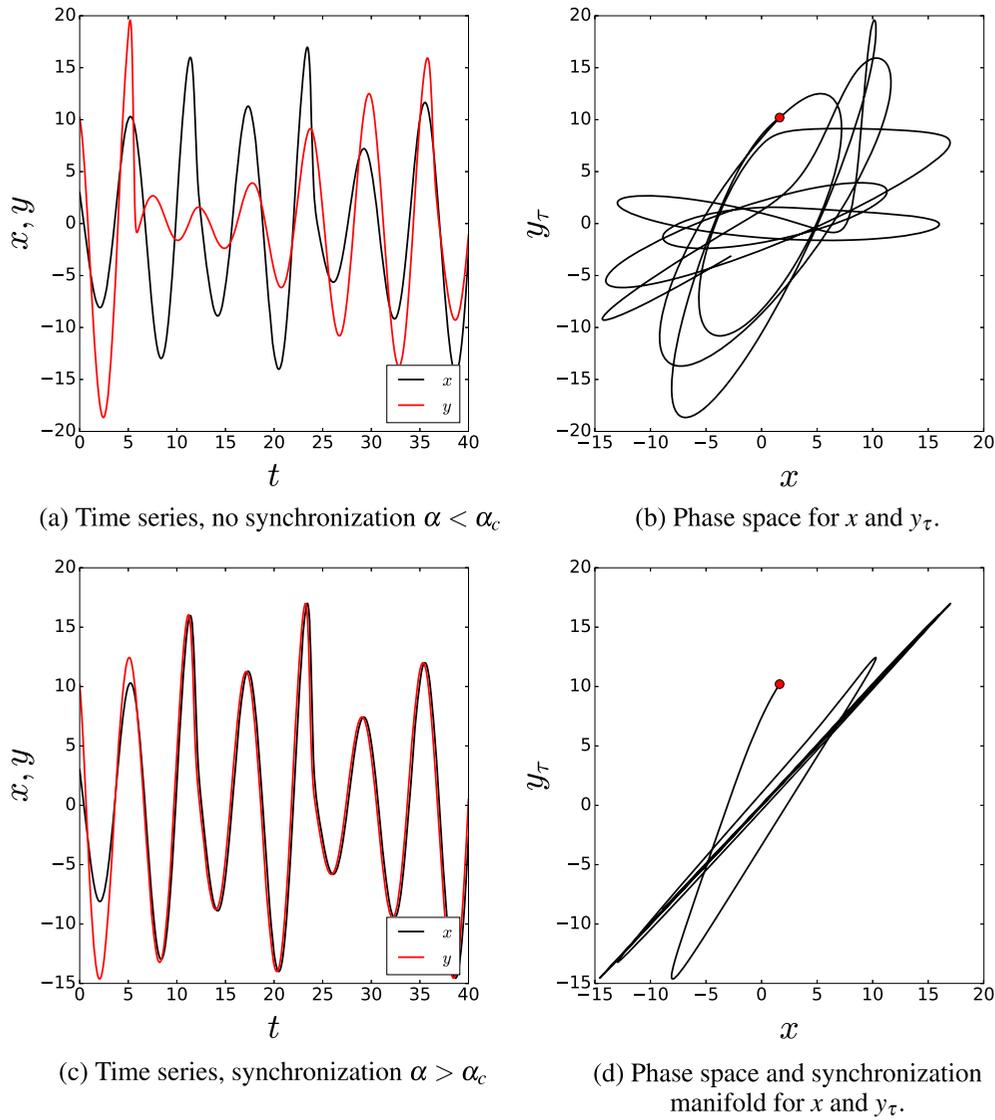


Figure 27. Anticipating synchronisation does not obtain for $(\alpha < \alpha_c)$, otherwise $(\alpha > \alpha_c)$ the AS exhibits. For this illustration the delay is selected as $\tau = 0.2$. (a) time series of the systems where $\alpha < \alpha_c$ and (b) phase space. (c) time series of the systems where $\alpha > \alpha_c$ and (d) phase space and synchronisation manifold of the system. The red circle is the initial condition for the trajectory of (x, y_τ) .

$$\frac{d\mathbf{x}_i}{dt} = \mathbf{f}(\mathbf{x}_i) - \alpha \sum_{j=1}^N L_{ij} \mathbf{H}(\mathbf{x}_j). \quad (42)$$

Our results will depend on this representation and on the spectral properties of \mathbf{L} .

Notice that

$$\mathbf{x}_1(t) = \mathbf{x}_2(t) = \dots = \mathbf{x}_N(t) = s(t),$$

is an invariant state for all coupling strength α , because the Laplacian is zero row sum. When $\alpha = 0$ the oscillators are decoupled, and Equation (52) describes N copies of the same oscillator with distinct initial conditions. Since, the chaotic behaviour leads to a divergence of nearby

trajectories, without coupling, any small perturbation on the globally synchronised motion will grow exponentially fast, and lead to distinct behaviour between the node dynamics. We wish to address the local stability of the globally synchronised state. That is, if all trajectories start close together would become synchronised

$$\lim_{t \rightarrow \infty} \|\mathbf{x}_i(t) - \mathbf{x}_j(t)\| = 0$$

The goal of the remaining exposition is to answer this question. Before, we continue with the analysis, we will briefly review some examples and constructions of graphs and discuss the relevant aspects necessary to tackle for problem.

4.2. Relation to other types of synchronisation

We will focus on the transition to complete synchronisation, which is mainly related to Section 2.2. This is no severe restriction as in certain scenarios other types of synchronisation can often be formulated in terms of our model.

Extension to Phase Synchronisation. As we discussed in the introduction of Section 4, if the dynamics of f is periodic then we can introduce a phase variable which will follow our main system of equations Equation (41) as the phase reduction tells us that generically the interaction between phases are diffusive. Moreover, because our results will yield robust transition to synchronisation, if the oscillators are nearly identical the phase synchronisation will persist.

Extension to Generalised Synchronisation. Roughly speaking a form of generalised synchronisation in networks is the so-called pinning control, where one tries to control the behaviour of synchronised trajectories by driving the network with external nodes. One extends the network to include the driver node. Therefore, the theory necessary to tackle this problem is the same as presented here. The main question is how to connect the driver nodes in such a way that control is effective.

4.3. Complex networks

A *network*, also called graph G in mathematical literature, is a set of N elements, called *nodes* (or *vertices*), connected by a set of M *links* (or *edges*) Figure 28. Networks represent interacting elements and are all around from biological systems, e.g. swarm of fireflies, food webs or brain networks, to mankind made systems, e.g. the world wide web, power grids, transportation networks or social networks.

A *network is called simple* if the nodes do not have self-loops (i.e. nodes have connections to themselves). An illustration of a simple network is found in Figure 29(a). A nonsimple network, therefore, containing connections is depicted in Figure 29(d). We need a bit of technology from graph theory to make sense of our networks. A few basic notions are as follows

A *network is undirected* if there is no distinction between the two nodes associated with each link Figure 29(a).

A *path in a network* is a route (sequence of edges) between connected nodes without repeating i.e. a path can visit a node only once. The length of a path is the number of links in the path. See further details in Refs. [102,103]. In Figure 30 we illustrate the paths of two selected (red) nodes in a network. Between two red nodes there are five different paths and each path is given in subplots of Figure 30. The length of paths are five for Figure 30(a),

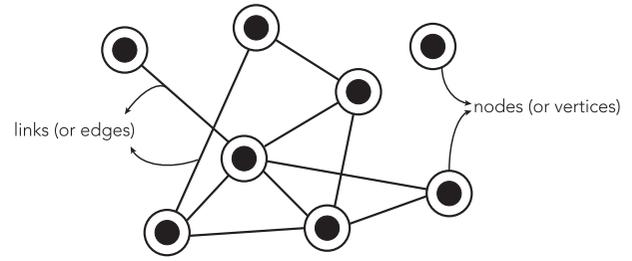


Figure 28. A network visualisation with eight nodes and ten links.

four for Figure 30(b), three for Figure 30(c) and (d), and two for Figure 30(e). Therefore the shortest path length, also called *geodesic path*, between these two red nodes is illustrated in Figure 30(e).

The *network diameter d* is the longest length of the shortest path between all possible pairs of nodes. In order to compute the diameter of a graph, first we find the shortest path between each pair of nodes. The longest length of all these geodesic paths is the diameter of the graph. If there is an isolated node (a node without any connections) or disconnected network components, then the diameter of the network is infinite. A network of finite diameter is called *connected* (Figure 29(a)), otherwise *disconnected* (Figure 29(b)).

A *network is directed* if the links transmit the information towards only associated direction Figure 29(c). If the graph is directed then there are two connected nodes say, u and v , such that u reachable from v , but v is not reachable from u . See Figure 29(c) for an illustration.

A *network is weighted* if links have different importance from each other or the links may carry different amount of information. Such graphs are called *weighted networks* Figure 29(d). Moreover, a network may have self-loops, that is, a node can affect itself as well Figure 29(d).

A graph can be described by its *adjacency matrix A* with $N \times N$ elements A_{ij} . The adjacency matrix A encodes the topological information, and is defined as $A_{ij} = 1$ if i and j are connected, otherwise 0. Therefore, the adjacency matrix of an undirected network is symmetric, $A_{ij} = A_{ji}$. The *degree k_i* of the i th node is the number of edges belongs to the node, defined as

$$k_i = \sum_j A_{ij}.$$

The *Laplacian matrix L* is another way to represent the network, defined as

$$L_{ij} = \begin{cases} k_i & \text{if } i = j \\ -1 & \text{if } i \text{ and } j \text{ are connected} \\ 0 & \text{otherwise.} \end{cases} \quad (43)$$

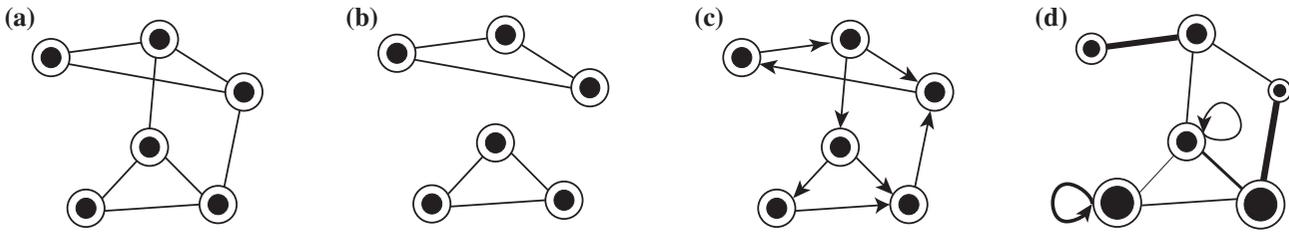


Figure 29. Visualisation of network types: (a) an unweighted simple network, (b) a disconnected network, (c) a directed network, (d) a weighted network with self-loops.

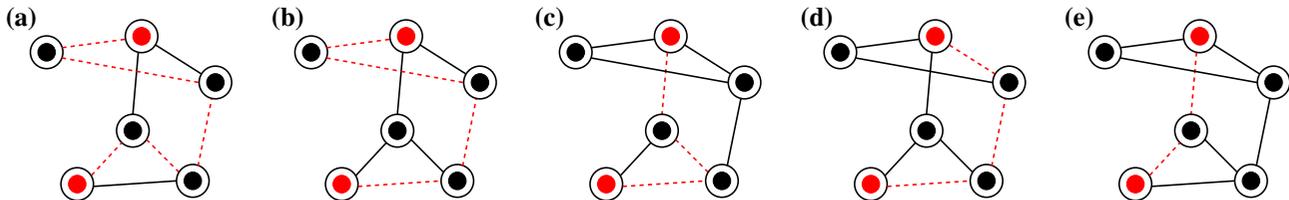


Figure 30. Visualisation of paths (red dashed) between two selected nodes (red) in a network: path length of (a) is five, (b) is four, (c) and (d) are three and (e) is two. Therefore the shortest path length for these two red nodes is two.

There is a direct relationship between the Laplacian L and the adjacency matrix A . In a compact form it reads

$$L_{ij} = \delta_{ij}k_i - A_{ij}$$

where δ_{ij} is the Kronecker delta, which is 1 if $i = j$ and 0 otherwise. We demonstrate some example network sketches with their adjacency A and Laplacian L matrices in Figure 31.

The networks we encounter in real applications have a wilder connection structure. Typical examples are cortical networks, the Internet, power grids and metabolic networks [103,104]. These networks don't have a regular structure of connections such as the ones presented in Figure 31. We say that the network is *complex* if it does not possess a regular connectivity structure.

One of the goals is to understand the relation between the topological organisation of the network and its relation functioning such as its collective motion.

2k Regular Graph is a standard graph model where each node has $2k$ links then the total number of links is $M = kN$ where N is total number of nodes Figure 32(a). This model is rather important one since the connections of spatiotemporal graphs, in general, connected to the nearest neighbours. *2k regular graph* is an alternative representation of such models. It is important to mention that the graph model is fixed with given k and N therefore all properties of the graph is known analytically.

Erdős-Rényi network is generated by setting an edge between each pair of nodes with equal probability p , independently of the other edges Figure 33(a). If $p \gg \ln N/N$, then the network is almost surely connected,

that is, as N tends to infinity, the probability that a graph on n vertices is connected tends to 1. The degree is pretty homogeneous, almost surely every node has the same expected degree [105].

Small World network is a random graph model which possesses the small-world properties; i.e. the average path length is short and clustering is large. The network is generated from a $2k$ regular graph, each link of the graph is rewired with a probability p . Therefore if $p = 0$ then there is no rewiring and the graph is $2k$ regular. For $p = 1$ then each link is rewired i.e. the graph is approaching to Erdős-Rényi network with $p = \frac{kN}{2\binom{N}{2}}$. The small-world properties come true between $0 < p < 1$ Figure 32. In many real world networks, the properties of small-world topology can be obtained.

The Barabasi-Albert network possesses a great deal of heterogeneity in the node's degree, while most nodes have only a few connections, some nodes, termed hubs, have many connections Figure 33(a). These networks do not arise by chance alone. The network is generated by means of the cumulative advantage principle – the rich gets richer. According to this process, a node with many links will have a higher probability to establish new connections than a regular node. The number of nodes of degree k is proportional to $k^{-\beta}$. These networks are called scale-free networks [103,104]. Many graphs arising in various real world networks display similar structure as the Barabasi-Albert network [106–108].

Hypercube graph Q_m is a m -dimensional hypercube formed regular graph (Figure 33(b)). It is a regular graph since each node has m neighbours and total number of nodes is 2^m and edges is 2^{m-1} .

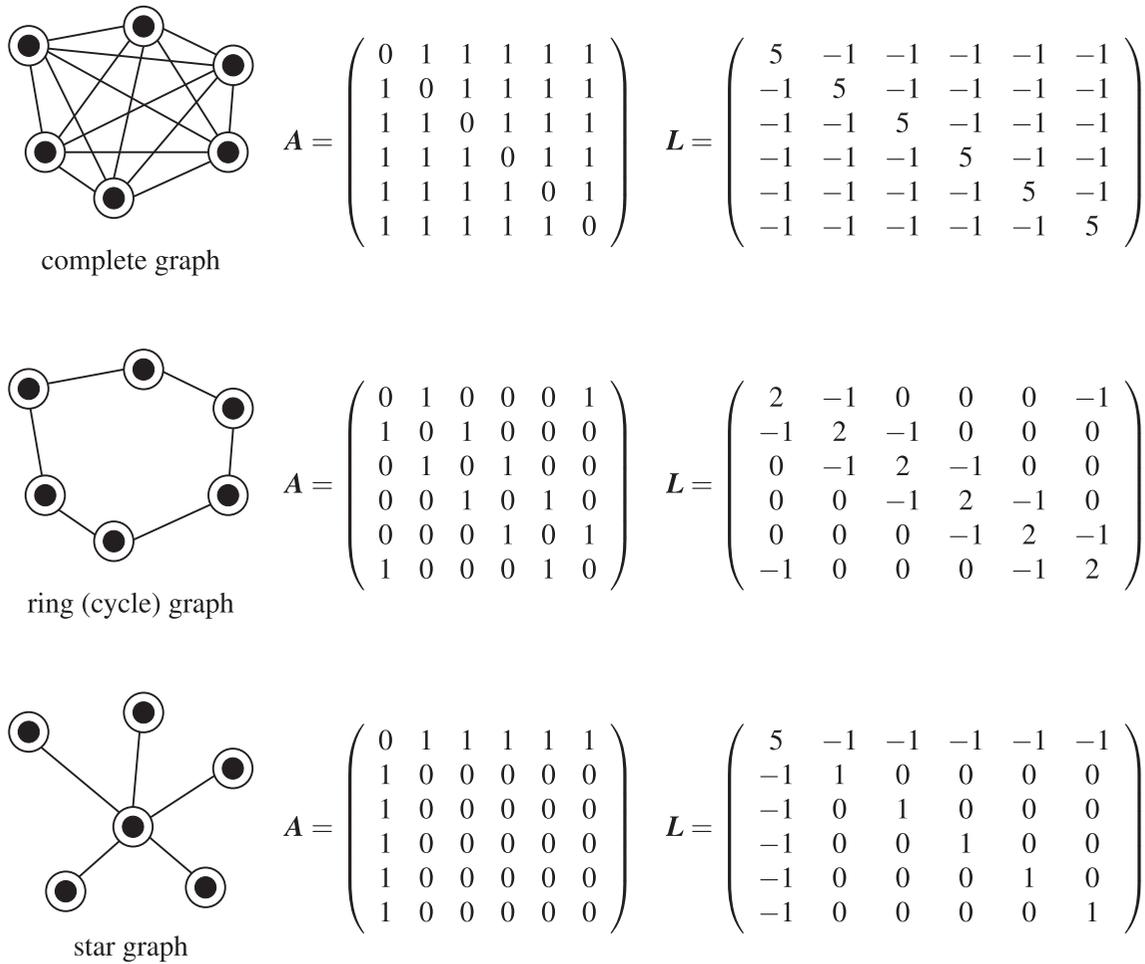


Figure 31. Various network examples with six nodes. Their adjacency A and Laplacian L matrices.

Random networks serve as a proxy to many applications as well as a surrogate. There are many nice ways to construct random network

Configuration Model is a random network model created by the degree distribution $P(k)$. If the degree distribution of a graph is known, then the nodes with associated number of links are known however the connection structure between nodes is unknown. The nodes can be drawn with their stubs (half links) Figure 34(a), then randomly these stubs can be linked and two stubs create a proper link Figure 34(b). This process is a random matching, obviously different network structures can arise from this random process.

Expected degrees Fix a network of N nodes and consider a vector

$$\mathbf{w} = (w_1, w_2, \dots, w_N),$$

In this model, each link A_{ij} between nodes i and j is an independent Bernoulli variable taking value 1 with success probability

$$p_{ij} = w_i w_j \rho,$$

and 0 with probability $1 - p_{ij}$, where

$$\rho = \frac{1}{\sum_{i=1}^N w_i}.$$

To ensure that $p_{ij} \leq 1$ it is assumed that $\mathbf{w} = \mathbf{w}(N)$ is chosen so that

$$\Delta^2 \rho \leq 1. \quad (44)$$

The *degree* $k_i = \sum_j A_{ij}$ of the i th is a random variable. An interesting property of this model is that under this construction w_i is the expected value of k_i , that is,

$$\mathbb{E}_{\mathbf{w}}(k_i) = \sum_j \mathbb{E}_{\mathbf{w}}(A_{ij}) = w_i$$

So, the different to the configuration model is that we do not fix the node degree, but rather obtain the degree probabilistically. This model also has many desirable concentration properties in the large N limit.

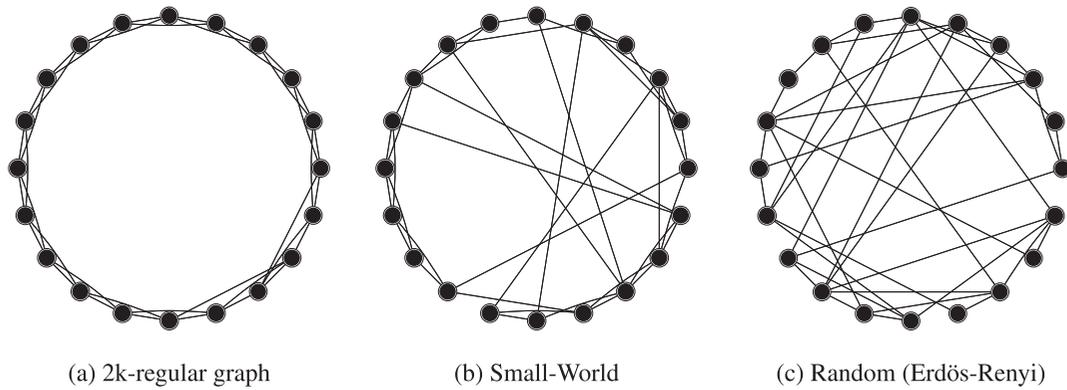


Figure 32. Watts–Strogatz random network approach.

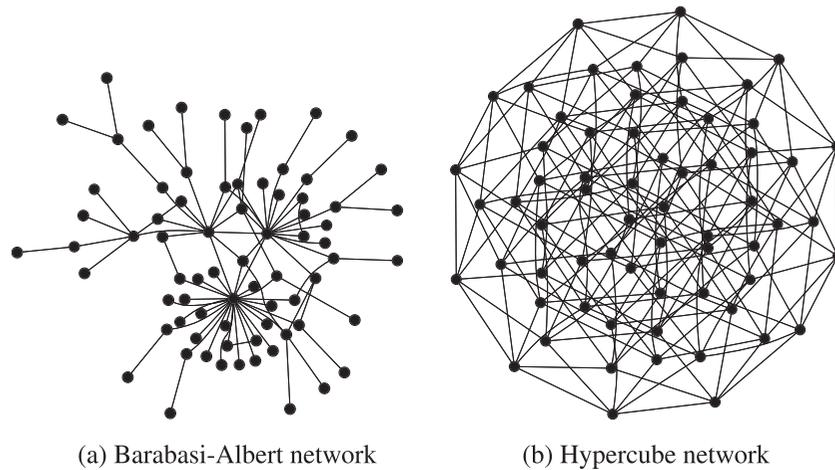


Figure 33. Some examples of complex networks.

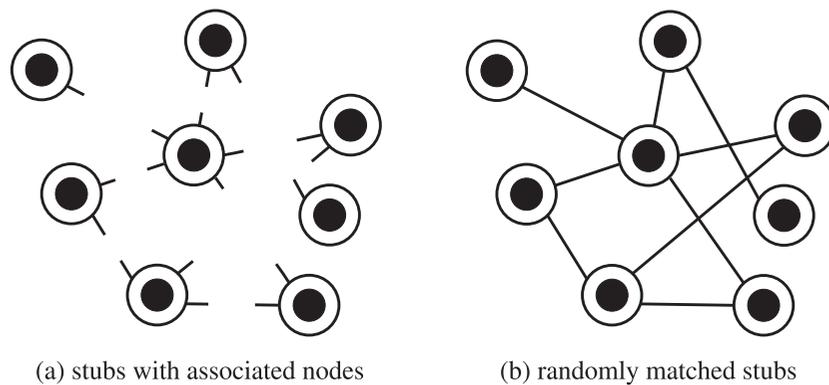


Figure 34. Configuration model.

4.4. Spectral properties of the Laplacian

The eigenvalues and eigenvectors of A and L tell us a lot about the network structure. The eigenvalues of L for instance are related to how well connected is the graph and how fast a random walk on the graph could

spread. In particular, the smallest nonzero eigenvalue of L will determine the synchronisation properties of the network. Since the graph is undirected the matrix L is symmetric its eigenvalues are real, and L has a complete set of orthonormal eigenvectors [109]. The next result characterises important properties of the Laplacian

Theorem 1: Let G be an undirected network and \mathbf{L} its associated Laplacian. Then:

- (a) \mathbf{L} has only real eigenvalues,
- (b) 0 is an eigenvalue and a corresponding eigenvector is $\mathbf{1} = (1, 1, \dots, 1)^*$, where $*$ stands for the transpose.
- (c) \mathbf{L} is positive semidefinite, its eigenvalues enumerated in increasing order and repeated according to their multiplicity satisfy

$$0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_N$$

- (d) The multiplicity of 0 as an eigenvalue of \mathbf{L} equals the number of connect components of G .

Therefore, λ_2 is bounded away from zero whenever the network is connected. The smallest non-zero eigenvalue is known as algebraic connectivity, and it is often called the Fiedler value. The spectrum of the Laplacian is also related to some other topological invariants. One of the most interesting connections is its relation to the diameter, size and degrees.

Theorem 2: Let G be a simple network of size N and \mathbf{L} its associated Laplacian. Then:

- (1) [110] $\lambda_2 \geq \frac{4}{Nd}$
- (2) [111] $\lambda_2 \leq \frac{N}{N-1}k_1$

We will not present the proof of the Theorem here, however, they can be found in references we provide in the theorem. We suggest the reader to see further bounds on the spectrum of the Laplacian in Ref. [112]. Also Ref. [113] presents many applications of the Laplacian eigenvalues to diverse problems. One of the main goals in spectral graph theory is to obtain better bounds by having access to further information on the graphs.

For a fixed network size, the magnitude of λ_2 reflects how well connected is graph.

5. Stability of synchronised solutions

We will state two results on network synchronisation. The first assumes that the coupling function \mathbf{H} is linear and the network is undirected. This assumption facilitates the discussion of the main ideas. Then, we will discuss the case where the coupling is non-linear. Basically, the results are the same with an additional complication as the latter involves the theory of Lyapunov exponents.

Theorem 3: Consider the diffusively coupled network model

$$\dot{\mathbf{x}}_i = \mathbf{f}(\mathbf{x}_i) - \alpha \sum_{j=1}^N L_{ij} \mathbf{H}(\mathbf{x}_j),$$

on an undirected and connected network. Assume that the \mathbf{H} is a positive definite linear operator. The reason of the assumption will appear in Step 5 of the theorem. Then, there is $\Gamma = \Gamma(\mathbf{f}, \mathbf{H})$ such that for any

$$\alpha > \frac{\Gamma}{\lambda_2},$$

the global synchronisation is uniformly asymptotically stable. Moreover, the transient to the globally synchronised behaviour is given the spectral gap λ_2 , that is, for any i and j

$$\|\mathbf{x}_i(t) - \mathbf{x}_j(t)\| \leq C e^{-(\alpha\lambda_2 - \Gamma)t},$$

where C is a constant.

The above result relates the threshold coupling for synchronisation in contributions coming solely from dynamics Γ , and network structure λ_2 .

Definition 1: [Synchronisation Threshold] We call

$$\alpha_c(\mathbf{f}, \mathbf{H}, G) = \frac{\Gamma(\mathbf{f}, \mathbf{H})}{\lambda_2(G)}$$

the critical synchronisation coupling.

Therefore, for a fixed node dynamics \mathbf{f} and coupling \mathbf{H} we can analyse how distinct network facilitates or inhibits global synchronisation. To continue our discussion we need the following

Definition 2: [Better Synchronisable] We say that the network G_1 is more synchronisable than G_2 if for fixed \mathbf{f} and \mathbf{H} we have

$$\alpha_c(G_1) < \alpha_c(G_2)$$

Likewise, we say that G_1 has better synchronisability than G_2 .

5.1. Which networks synchronise best

In this setting, the coupling function \mathbf{H} is positive definite and the network is undirected, the synchronisability depends only on the spectral gap. Using the previous study on the properties of various networks presented in Section 4.4. We consider networks of N nodes then

- A complete graph is the most synchronisable. In fact, $\alpha_c \approx 1/N$. So, the larger the graph the less coupling strength is necessary to synchronise the network.
- A path or ring are poorly synchronisable. For these networks, $\alpha_c \approx N^2$.
- $2k$ -regular graphs share the same properties for also poorly synchronisable when $k \ll N$.
- Erdős–Renyi graphs the synchronisation properties depend only on the mean degree $\alpha_c \approx 1/\langle k \rangle$.

- *Small world networks* are better than regular but worse than random graphs. In the limit of large graphs $\alpha_c \approx 1/s$ where s is the fraction of added random links. In general, adding links to a network favours synchronisation.

5.2. Proof of the stable synchronisation

Now we present the proof of Theorem 3. We omit some details that are not relevant for the understanding of the proof. A full discussion of the proof can be found in [116]. We will show that whenever the nodes start close together they tend to the same future dynamics, that is, $\|\mathbf{x}_i(t) - \mathbf{x}_j(t)\| \rightarrow 0$, for any i and j . For pedagogical purposes we split the proof into six main steps.

Step 1: Kronecker Form. We have N coupled equations each has dimension n . Because of the nice structure of the interaction we can use the Kronecker product to write them as a single block. Given two matrices $\mathbf{A} \in \mathbb{R}^{p \times q}$ and $\mathbf{B} \in \mathbb{R}^{r \times s}$, the Kronecker Product of the matrices \mathbf{A} and \mathbf{B} and defined as the matrix

$$\mathbf{A} \otimes \mathbf{B} = \begin{pmatrix} A_{11}\mathbf{B} & \cdots & A_{1q}\mathbf{B} \\ \vdots & \ddots & \vdots \\ A_{p1}\mathbf{B} & \cdots & A_{pq}\mathbf{B} \end{pmatrix},$$

we introduce the following notation

$$\mathbf{X} = \text{col}(\mathbf{x}_1, \dots, \mathbf{x}_N),$$

where col denotes the vector formed by stacking the columns vectors \mathbf{x}_i into a single vector. Similarly

$$\mathbf{F}(\mathbf{X}) = \text{col}(\mathbf{f}(\mathbf{x}_1), \dots, \mathbf{f}(\mathbf{x}_N)).$$

Then Equation (52) can be arranged into a compact form

$$\frac{d\mathbf{X}}{dt} = \mathbf{F}(\mathbf{X}) - \alpha(\mathbf{L} \otimes \mathbf{H})\mathbf{X}, \quad (45)$$

where \otimes is the Kronecker product. The easiest way to check that this is correct is to compute the i th block of dimension n and compare with the equation for the i th node.

Step 2: Transversal Laplacian Eigenmodes. The Kronecker product has many nice properties such as

$$(\mathbf{A} \otimes \mathbf{B})(\mathbf{C} \otimes \mathbf{D}) = \mathbf{AC} \otimes \mathbf{BD}. \quad (46)$$

And this holds whenever the matrix multiplication make sense. A nice consequence of the multiplication result in Kronecker form is that if

$$\begin{aligned} \mathbf{A}\mathbf{v}_i &= \lambda_i\mathbf{v}_i \quad \text{and} \quad \mathbf{B}\mathbf{u}_j = \mu_j\mathbf{u}_j \quad \text{then} \\ \mathbf{A} \otimes \mathbf{B}(\mathbf{v}_i \otimes \mathbf{u}_j) &= \lambda_i\mu_j\mathbf{v}_i \otimes \mathbf{u}_j \end{aligned}$$

Since we are assuming that \mathbf{L} is undirected and \mathbf{H} is positive definite the eigenvectors $\{\mathbf{v}_i\}_{i=1}^N$ of \mathbf{L} form a basis of \mathbb{R}^N . Likewise, the eigenvectors of \mathbf{H} form a basis of \mathbb{R}^m . This implies that the eigenvectors of $\mathbf{L} \otimes \mathbf{H}$ form a basis of \mathbb{R}^{Nn} . Using this fact we can represent \mathbf{X} as

$$\mathbf{X} = \sum_{i=1}^N \mathbf{v}_i \otimes \mathbf{y}_i$$

where \mathbf{y}_i is the coordinates of \mathbf{X} in the Kronecker basis. For sake of simplicity we call $\mathbf{y}_1 = \mathbf{s}$, and remember that $\mathbf{v}_1 = \mathbf{1}$ is an eigenvector. Hence

$$\mathbf{X} = \mathbf{1} \otimes \mathbf{s} + \mathbf{U},$$

where

$$\mathbf{U} = \sum_{i=2}^N \mathbf{v}_i \otimes \mathbf{y}_i.$$

In this way we split the contribution in the direction of the global synchronisation and \mathbf{U} , which accounts for the contribution of the transversal. Note that if \mathbf{U} converges to zero then the system completely synchronise, that is \mathbf{X} converges to $\mathbf{1} \otimes \mathbf{s}$ which clearly implies that

$$\mathbf{x}_1 = \cdots = \mathbf{x}_N = \mathbf{s}$$

The goal then is to obtain conditions so that $\|\mathbf{U}\| \rightarrow 0$.

Step 3: Variational equations for the Transversal Modes. The equation of motion in terms of the Laplacian modes decomposition reads

$$\begin{aligned} \mathbf{1} \otimes \frac{d\mathbf{s}}{dt} + \frac{d\mathbf{U}}{dt} &= \mathbf{F}(\mathbf{1} \otimes \mathbf{s} + \mathbf{U}) \\ &\quad - \alpha(\mathbf{L} \otimes \mathbf{H})(\mathbf{1} \otimes \mathbf{s} + \mathbf{U}), \end{aligned}$$

We assume that \mathbf{U} is small and perform a Taylor expansion about the synchronisation manifold.

$$\mathbf{F}(\mathbf{1} \otimes \mathbf{s} + \mathbf{U}) = \mathbf{F}(\mathbf{1} \otimes \mathbf{s}) + D\mathbf{F}(\mathbf{1} \otimes \mathbf{s})\mathbf{U} + \mathbf{R}(\mathbf{U}),$$

where $\mathbf{R}(\mathbf{U})$ is the Taylor remainder $\|\mathbf{R}(\mathbf{U})\| = O(\|\mathbf{U}\|^2)$. Using the Kronecker product properties Equation (46) and the fact that $\mathbf{L}\mathbf{1} = \mathbf{0}$, together with

$$\mathbf{1} \otimes \frac{d\mathbf{s}}{dt} = \mathbf{F}(\mathbf{1} \otimes \mathbf{s}) = \mathbf{1} \otimes \mathbf{f}(\mathbf{s})$$

and we have

$$\frac{d\mathbf{U}}{dt} = [D\mathbf{F}(\mathbf{1} \otimes \mathbf{s}) - \alpha(\mathbf{L} \otimes \mathbf{H})]\mathbf{U} + \mathbf{R}(\mathbf{U}) \quad (47)$$

and likewise

$$D\mathbf{f}(\mathbf{1} \otimes s)\mathbf{U} = [\mathbf{I}_N \otimes D\mathbf{f}(s)]\mathbf{U},$$

therefore, the first variational equation for the transversal modes reads

$$\frac{d\mathbf{U}}{dt} = [\mathbf{I}_N \otimes D\mathbf{f}(s) - \alpha\mathbf{L} \otimes \mathbf{I}]\mathbf{U}. \quad (48)$$

Step 4: Decoupling of Transversal Modes. Instead of analysing the full set of equations, we can do much better by projecting the equation into subspace $W_i = \text{span}\{\mathbf{v}_i \otimes \mathbf{I}\}$. Let $P_i : \mathbb{R}^N \otimes \mathbb{R}^n \rightarrow W_i$ be a projection operator given by $P_i = \mathbf{v}_i \mathbf{v}_i^* \otimes \mathbf{I}$, it follows that P_i is an orthogonal projection since \mathbf{v}_i 's are orthonormal. Using Equation (48) and the identity Equation (46) we obtain

$$P_i \frac{d\mathbf{U}}{dt} = [\mathbf{v}_i \mathbf{v}_i^* \otimes D\mathbf{f}(s) - \alpha(\mathbf{v}_i \mathbf{v}_i^* \mathbf{L}) \otimes \mathbf{I}]\mathbf{U}, \quad (49)$$

$$= [\mathbf{v}_i \mathbf{v}_i^* \otimes D\mathbf{f}(s) - \alpha\lambda_i(\mathbf{v}_i \mathbf{v}_i^*) \otimes \mathbf{I}]\mathbf{U} \quad (50)$$

where in the last passage we used that the network is undirected implying $\mathbf{v}_i * \mathbf{L} = \lambda_i \mathbf{v}_i^*$. Using and the fact that $\mathbf{v}_j^* \mathbf{v}_i = \delta_{ij}$, where is δ_{ij} the Kronecker delta, we have that $(\mathbf{v}_i \mathbf{v}_i^* \otimes \mathbf{I})\mathbf{U} = \sum_{j=2}^N \mathbf{v}_i \delta_{ij} \otimes \mathbf{y}_j$. Moreover, since P_i does not depend on time $P_i \dot{\mathbf{U}} = (P_i \mathbf{U})$

$$\sum_{j=2}^N \mathbf{v}_i \delta_{ij} \otimes \frac{d\mathbf{y}_j}{dt} = \sum_{j=2}^N \mathbf{v}_i \delta_{ij} \otimes [D\mathbf{f}(s) - \alpha\lambda_i \mathbf{I}]\mathbf{y}_j$$

the nonzero coefficients give the dynamics in W_i . Hence,

$$\frac{d\mathbf{y}_i}{dt} = [D\mathbf{f}(s) - \alpha\lambda_i \mathbf{I}]\mathbf{y}_i$$

All blocks have the same form which are different only by λ_i , the i th eigenvalue of L . We can write all the blocks in a parametric form

$$\frac{d\mathbf{u}}{dt} = \mathbf{K}(t)\mathbf{u}, \quad (51)$$

where

$$\mathbf{K}(t) = D\mathbf{f}(s(t)) - \kappa \mathbf{I}$$

with $\kappa \in \mathbb{R}$. Hence if $\kappa = \alpha\lambda_i$ we have the equation for the i th block. This is just the same type of equation we encounter before in the example of the two coupled oscillators, see Equation (9).

Step 5. Stability. Because \mathbf{H} is positive definite we can first solve the homogeneous equation $\dot{\mathbf{u}} = -\kappa \mathbf{H} \mathbf{u}$. This equation has an globally attracting trivial solution. Then, we incorporate $D\mathbf{f}$ in terms of the variation of constants

formula. So, first notice that

$$\mathbf{u}(t) = e^{-\kappa \mathbf{H} t} \mathbf{u}_0 \Rightarrow \|\mathbf{u}(t)\| \leq K \|\mathbf{u}_0\| e^{-\kappa \lambda_{\min}(\mathbf{H})}$$

where $\lambda_{\min}(\mathbf{H})$ is the smallest eigenvalue of \mathbf{H} . So, by the variation of constants formulate

$$\mathbf{u}(t) = e^{-\kappa \mathbf{H} t} \mathbf{u}_0 + \int_0^t e^{-\kappa \mathbf{H}(t-\tau)} D\mathbf{f}(s(\tau)) d\tau$$

taking norms

$$\|\mathbf{u}(t)\| = K \|\mathbf{u}_0\| + \int_0^t e^{-\kappa \lambda_{\min}(\mathbf{H})(t-\tau)} \|D\mathbf{f}(s(\tau))\| d\tau$$

where K is a constant and defining $M_f = \sup_t \|D\mathbf{f}(s(t))\|$ by Gronwal inequality

$$\|\mathbf{u}(t)\| = K_1 \|\mathbf{u}_0\| e^{(-\kappa \lambda_{\min}(\mathbf{H}) + M_f)t}.$$

The trivial solution will be exponentially stable when

$$-\kappa \lambda_{\min}(\mathbf{H}) + M_f < 0 \Rightarrow \kappa > \Gamma = \frac{M_f}{\lambda_{\min}(\mathbf{H})}$$

Recall that taking $\kappa = \alpha\lambda_i > \Gamma$ we are stabilising the equation for the i th block. But, once we stabilise the second block all blocks will be stable (because λ_2 is the smallest nonzero eigenvalue)

$$\alpha\lambda_N \geq \dots \geq \alpha\lambda_3 \geq \alpha\lambda_2 \geq \Gamma$$

Hence, the stability condition so that all blocks have exponentially stable trivial solution is

$$\alpha > \frac{\Gamma}{\lambda_2}$$

Step 6: Norm Estimates and Nonlinearities. Using the bounds for the blocks it is easy to obtain a bound for the norm of the evolution operator. Indeed, note that

$$\|\mathbf{U}\|_2 = \left\| \sum_{i=2}^N \mathbf{v}_i \otimes \mathbf{y}_i \right\|_2 \leq \sum_{i=2}^N \|\mathbf{v}_i\| \|\mathbf{y}_i\|_2$$

Therefore,

$$\|\mathbf{U}\|_2 \leq \sum_{i=2}^N \|\mathbf{v}_i\| K_i e^{-\eta_i(t-s)} \|\mathbf{y}_i(s)\|$$

Now using that $e^{-\eta_i(t-s)} \leq e^{-(\alpha\lambda_i - \Gamma)(t-s)}$, so

$$\|\mathbf{U}(t)\|_2 \leq K_2 e^{-\eta(t-s)}$$

with $\eta = \alpha\lambda_2 - \Gamma$ for any $t \geq s$.

Because the trivial solution is exponentially stable (uniformly in $s(t)$) by the principle of linearisation, we conclude that the nonlinearities coming Taylor remainder does not affect the stability of the trivial solution, which correspond to the global synchronisation.

The claim about the transient is straightforward, because all norms are equivalent in finite dimensions we can take

$$\|\mathbf{X}(t) - \mathbf{1} \otimes s(t)\|_\infty \leq K_3 e^{-\eta(t-s)} \|\mathbf{U}(s)\|_\infty$$

implying that $\max_i \|\mathbf{x}_i(t) - s(t)\|_2 \leq K_3 e^{-\eta(t-s)} \|\mathbf{U}(s)\|_\infty$ and in virtue of the triangular inequality

$$\|\mathbf{x}_i(t) - \mathbf{x}_j(t)\|_\infty \leq \|\mathbf{x}_i(t) - s(t)\|_\infty + \|\mathbf{x}_j(t) - s(t)\|_\infty$$

and using the previous bound, we concluding the proof. \square

6. General diffusive coupling and master stability function

Until now we have considered linear coupling functions which are positive definite. This assumption can be relaxed and thereby we are generalise our previous results. The statement will then become rather technical and will be beyond the scope of our review. So, here we will discuss the main ideas but will not give much details on the technical issues. Consider the function $\mathbf{g} : \mathbb{R}^n \times \mathbb{R}^n \rightarrow \mathbb{R}^n$. We say that \mathbf{g} is diffusive if

$$\mathbf{g}(\mathbf{x}, \mathbf{x}) = 0 \quad \text{and} \quad \mathbf{g}(\mathbf{x}, \mathbf{y}) = -\mathbf{g}(\mathbf{y}, \mathbf{x})$$

Hence, we can extend the model to a general diffusive coupling

$$\dot{\mathbf{x}}_i = \mathbf{f}(\mathbf{x}_i) + \alpha \sum_{j=1}^N A_{ij} \mathbf{g}(\mathbf{x}_j, \mathbf{x}_j). \quad (52)$$

We perform the analysis close to synchronisation $\mathbf{x}_i = s + \xi_i$ so

$$\begin{aligned} \mathbf{g}(\mathbf{x}_j, \mathbf{x}_j) &= \mathbf{g}(s + \xi_j, s + \xi_j) \\ &= \mathbf{g}(s, s) + D_1 \mathbf{g}(s, s) \xi_j + D_2 \mathbf{g}(s, s) \xi_j \end{aligned}$$

but because the coupling is diffusive

$$D_2 \mathbf{g}(s, s) = -D_1 \mathbf{g}(s, s)$$

we obtain

$$\mathbf{g}(\mathbf{x}_j, \mathbf{x}_i) = \mathbf{G}(s)(\xi_j - \xi_i) + \mathbf{R}(\xi_i, \xi_j)$$

where $\mathbf{G}(s) = D_1 \mathbf{g}(s, s)$ and $\mathbf{R}(\xi_i, \xi_j)$ contains quadratic terms. So, the first variational equation about the synchronisation manifold

$$\dot{\xi}_i = D \mathbf{f}(s(t)) \xi_i + \alpha \sum_{j=1}^N A_{ij} \mathbf{G}(s)(\xi_j - \xi_i) \quad (53)$$

$$= D \mathbf{f}(s(t)) \xi_i - \alpha \mathbf{G}(s) \sum_{j=1}^N L_{ij} \xi_j. \quad (54)$$

Now we can perform the same steps as before. In fact, Steps 1–4 remain unchanged. The change difference is Step 5, which concerns the stability of the modes. Performing all the steps 1 to 4 we obtain the parametric equation for the modes

$$\dot{\mathbf{u}} = [D \mathbf{f}(s(t)) - \kappa \mathbf{G}(s(t))] \mathbf{u}$$

And we can no longer apply the *trick* of using the coupling function \mathbf{H} to solve the equation. Here, $\mathbf{G}(s(t))$ depends on time and this generality tackling the stability of the trivial solution is challenging.

The main idea is to fix κ compute the maximum Lyapunov exponent $\Lambda(\kappa)$ as

$$\|\mathbf{u}(t)\| \leq C e^{\Lambda(\kappa)t}$$

see Appendix. The map

$$\kappa \mapsto \Lambda(\kappa) \quad (55)$$

is called master stability function. Notice that if $\Lambda(\kappa) < 0$ when $\kappa \in (\alpha_c^1, \alpha_c^2)$ then $\|\mathbf{u}\| \rightarrow 0$.

The stability condition then become

$$\alpha_c^1 \leq \alpha \lambda_2 \leq \dots \leq \alpha \lambda_N \leq \alpha_c^2$$

Or

$$\frac{\alpha_c^2}{\alpha_c^1} \geq \frac{\lambda_N}{\lambda_2} \quad (56)$$

This is a well studied condition. Much energy has been devoted to study the master stability function Equation (55), see e.g. [117].

6.1. Examples of master stability functions

Now let us consider coupled Rössler systems which are coupled through their x -coordinates:

$$\begin{aligned}\dot{x}_i &= -y_i - z_i + \alpha \sum_{j=1}^N A_{ij}(x_j - x_i) \\ \dot{y}_i &= x_i + ay_i \\ \dot{z}_i &= b + z_i(x_i - c).\end{aligned}\quad (57)$$

In order to compute $\Lambda_{\max}(\kappa)$, we find that $D\mathbf{f}$ and $D\mathbf{H}$ are given by

$$\begin{aligned}D\mathbf{f}(\mathbf{s}) &= \begin{pmatrix} 0 & -1 & -1 \\ 1 & a & 0 \\ z^* & 0 & x^* - c \end{pmatrix} \quad \text{and} \\ D\mathbf{H} = \mathbf{H} &= \begin{pmatrix} 1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix},\end{aligned}\quad (58)$$

x and z are the components of \mathbf{s} . The constants are $a = 0.2$, $b = 0.2$ and $c = 5.7$.

To compute $\Lambda(\kappa)$, we first simulate the isolated dynamics $\dot{\mathbf{s}} = \mathbf{f}(\mathbf{s})$ and obtain the trajectory $\mathbf{s}(t)$, then we feed this trajectory to $\dot{\mathbf{u}} = [D\mathbf{f}(\mathbf{s}(t)) - \kappa\mathbf{H}]\mathbf{u}$ and then for each κ estimate the maximal Lyapunov exponent $\Lambda(\kappa)$. The result is depicted in Figure 35.

Stability region where Λ is negative is bounded between $\alpha_c^1 \approx 0.13$ and $\alpha_c^2 \approx 4.4$. So, if the network is a complete graph then $\lambda_2 = \dots = \lambda_N = N$, so the network synchronisation when

$$\frac{\alpha_c^2}{N} > \alpha > \frac{\alpha_c^1}{N}$$

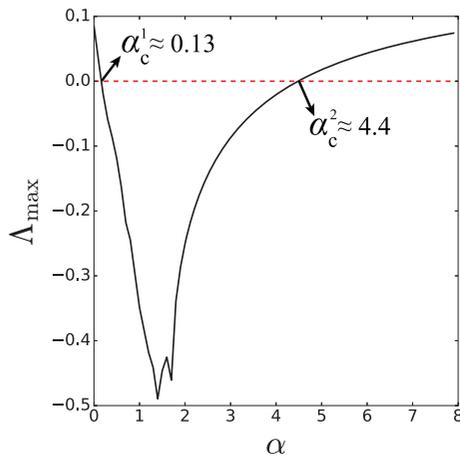


Figure 35. Master stability function for x -coupled Rössler attractors on a network structure.

6.2. Synchronisation conditions and synchronisation loss

In Section 5.1, we discussed the synchronisability of network types when there is only one critical threshold in other words $\alpha_c^2 \rightarrow \infty$. This case is true when \mathbf{H} is a positive matrix. In the general case, there are two finite critical couplings Equation (56) and the master stability function has a finite stability region between these critical points (as discussed in the above section). Now we discuss this synchronisation scenario for various network types. To quantify the synchronisability, we will use the network properties given in Table 1.

2k-regular graphs – Diameter driven synchronisation loss. When $k \ll N$, the mean geodesic distance (shortest path length) between nodes are increasing very fast as N increasing. Hence, while the diameter of the network is increasing $d = 2N/k$, speed of information exchange between the nodes is decreasing drastically. In this case, the complete synchronisation is not stable and it is visible from the Laplacian spectrum of the graph. For $N \rightarrow \infty$ and $k \ll N$, the extremal eigenvalues given in Table 1 can be extended to Taylor expansion and major role playing part can be rewritten as

$$\lambda_2 = \frac{4\pi^2(k + \frac{1}{2})^3}{3N^2} \quad \text{and} \quad \lambda_N = 2k \frac{N}{N-1}$$

The ratio between the largest λ_N and second smallest eigenvalue λ_2 of the Laplacian is not growing in the same scale,

$$\frac{\lambda_N}{\lambda_2} \approx N^2/k^2$$

and the condition in Equation (56) is never satisfied.

ER graphs – Optimal Synchronisation. In this case, the extremal eigenvalues λ_2 and λ_N of the Laplacian matrix increase in the same scale. The diameter of the network increases very slowly $d \approx \log N$ when the network size N increases Table 1. Therefore the synchronisation is stable for any scale of size.

BA networks – Heterogeneity driven synchronisation loss. When the network is too heterogeneous the complete synchronisation is unstable, this is because the extremal eigenvalues λ_2 and λ_N grow in different scales and the condition in Equation (56) is never met. For instance, consider a BA network. Then, the eigenvalues satisfy

$$\lambda_2 \approx m_0 \quad \text{and} \quad \lambda_N \approx m_0 N^{1/2}$$

where m_0 is the mean degree. Hence, the eigenratio becomes

$$\frac{\lambda_N}{\lambda_2} \approx N^{1/2}$$

this should be compared to the stability interval given by the master stability function. Lets consider the example

Table 1. Network of N nodes with m mean degree. Examples of nonrandom networks are depicted in Figure 31 and random networks in Figure 33. For random networks, the mean degree is given by $m = \sum_k kP(k)$ where $P(k)$ is the degree distribution of the network. The randomness of the networks arises via probability p . Consider the degree distribution for the SF network $P(k) = k^{-\gamma}$. C is constant.

| Network | λ_2 | λ_N | k_{\max} | k_{\min} | d | Reference |
|---------------------|--|--|----------------|-----------------------------------|---|-----------|
| Complete | N | N | $N-1$ | $N-1$ | 1 | |
| Ring | $2 - 2 \cos\left(\frac{2\pi}{N}\right)$ | 4 | 2 | 2 | $(N+1)/2$ if N is odd $N/2$ if N is even | [56] |
| Star | 1 | N | $N-1$ | 1 | 2 | |
| Hypercube (Q_m) | 2 | $\frac{2 \ln(N-2)}{m} \left(\frac{m}{N}\right)$ | m | m | m | |
| 2k-regular | $2k + 1 - \frac{\sin\left[(2k+1)\frac{\pi}{N}\right]}{\sin\left(\frac{\pi}{N}\right)}$ | $2k + 1 - \frac{\sin\left[(2k+1)\frac{\pi(N-1)}{N}\right]}{\sin\left(\frac{\pi(N-1)}{N}\right)}$ | $2k$ | $2k$ | $\lceil (N+1)/2k \rceil$ if N is odd $\lfloor N/2k \rfloor$ if N is even | [56] |
| SW network | $\sim 2k + 1 - \frac{\sin\left[(2k+1)\frac{\pi}{N}\right]}{\sin\left(\frac{\pi}{N}\right)} + Cp$ | $\sim 2k + 1 - \frac{\sin\left[(2k+1)\frac{\pi(N-1)}{N}\right]}{\sin\left(\frac{\pi(N-1)}{N}\right)} + Cp$ | $\sim 2k$ | $\sim 2k$ | $O(\log N)$ | [103] |
| SF network | $\sim m_0 \left(1 - \frac{1}{(\gamma-1)}\right)$ | $\sim m_0 N^{\frac{1}{\gamma-1}}$ | $\sim m_0$ | $\sim m_0 N^{\frac{1}{\gamma-1}}$ | $O\left(\frac{\log N}{\log \log k}\right)$ | [114] |
| ER network | $\sim Np$ | $\sim Np$ | $pN(1 + O(1))$ | $(1-p)(1 + O(1))$ | $\frac{\log N}{2 \log Np}$ | [115] |

in the above section with the Rössler Equation (58). The master stability function gives (as seen in Figure 35) an stability interval $\alpha_c^2/\alpha_c^1 \approx 34$. The stability conditions Equation (56) reads as

$$N^{1/2} < 34$$

So, when the BA network is large enough it is not possible to synchronise the system. In particular the critical system size to be able to synchronisation a network of Rössler as in Equation (58) is, therefore,

$$N \approx 10^3.$$

6.2.1. Extensions

There are a few extensions of the model. Here we discuss a few directions.

Directed Networks. The major problem considering directed networks is that they may not be diagonalisable. So, the decoupling of transversal modes by projection is a nontrivial steps. There are a few ways to overcome this. The first, using Jordan decomposition of the Laplacian. The other possibility is to perturb the Laplacian to make the eigenvalues simples. This must be done in such a way that the perturbation does not spoil the stability. In both cases, the stability condition remain unchanged. Only the transients may be longer.

When the network is nondiagonalisable, small perturbations in the network may lead to large perturbations in the eigenvalues (the eigenvalues in this case are not differentiable functions of the perturbations) [118]. Moreover, structural improvements in the network may lead to desynchronisation [119].

Nonidentical Nodes. Here we consider $\mathbf{f}(\mathbf{x}) \mapsto \mathbf{f}(\mathbf{x}) + \mathbf{r}_i(\mathbf{x}, t)$, where \mathbf{r}_i is either a perturbation of the vector field or a signal playing the role of noise. When r_i is very small synchronisation will persist [120]. For general networks, Boltt and co-workers [100] extended the master stability function approach when \mathbf{r}_i is a perturbation of the vector field. Pereira and co-workers [101] study the effect of general perturbations $\|\mathbf{r}_i\| \leq \delta$ and the role that the network structure plays in suppressing the fluctuations. In the case where \mathbf{H} is positive definite and the network is undirected, they showed that the synchronisation error

$$E = \frac{1}{n(n-1)} \sum_{ij} \|\mathbf{x}_i - \mathbf{x}_j\|$$

behaves as

$$E \leq K \frac{\delta}{\alpha \lambda_2 - \alpha_c}$$

For example, if the oscillators where uncoupled and \mathbf{r}_i independent noise then the Central Limit theorem

would yield $E = O(N^{-1/2})$. For complete networks, the interaction and synchronisation yields $E = O(N^{-1})$ which is a large improvement over the naive application of the Central Limit theorem. In certain sense, this shows that interacting maybe better then isolation.

Nonidentical Coupling Functions. In many applications the coupling function are not identical and has fluctuating components [121]. Consider an undirected networks of identical oscillators and coupling function

$$\mathbf{H}_{ij}(\mathbf{x}_i - \mathbf{x}_j, t) = \mathbf{H}(\mathbf{x}_i - \mathbf{x}_j) + \mathbf{P}_{ij}(\mathbf{x}_i - \mathbf{x}_j, t)$$

where $\|\mathbf{P}_{ij}(\mathbf{x}_i - \mathbf{x}_j, t)\| \leq \eta$. In this case, the network structure will play a major on the size of perturbation η . If the network is random and the degree distribution homogeneous, then even for large perturbations δ synchronisation is stable. If the network is heterogeneous degrees such as Barabasi–Albert then typically $\delta_c = O(N^{-\beta})$ is the critical perturbation size. If $\delta > \delta_c$ synchronisation becomes unstable, solely because of the interaction between network structure and perturbations in the coupling function.

Cluster Synchronisation. According to similarities in coupled systems, such as symmetries in network topology or identical dynamics in a diverse population or equally time-delayed nodes in differently distributed feedbacks, the partial or cluster synchronisation can emerge. In order to enlighten the reason of these cluster synchronisation cases, many techniques are developed and experimental observations are analysed [122–127].

The symmetries are easy to detect for some network geometries for instance Bethe lattice is a regular graph which grows from a root (parent) node by p -nodes for ℓ -levels, an example of Bethe lattice given in Figure 36 for $p = 3$ and $\ell = 3$. The nodes in the same level of the Bethe graph, they all symmetric to each other. In Figure 36, the levels of the graph are given in the same colour and the cluster synchronisation occurs for each level.

Recently, Pecora et al. put forward that all the symmetries in network structures are not visible directly. They developed a computational group theory-based method to reveal these hidden symmetries and predicted possible synchronisation patterns [128,129].

Time-Delay Coupling. Simultaneous coupling is not always possible for real world applications in other words some time delays can occur in the interaction process. Therefore it is important to investigate synchronisability and stability of coupled time-delayed systems. The necessary conditions for time-delayed synchronisation is analytically shown by Pyragas [130]. The finding of time-delay synchronisation is used as an application for the anticipating synchronisation (see Section 3.4).

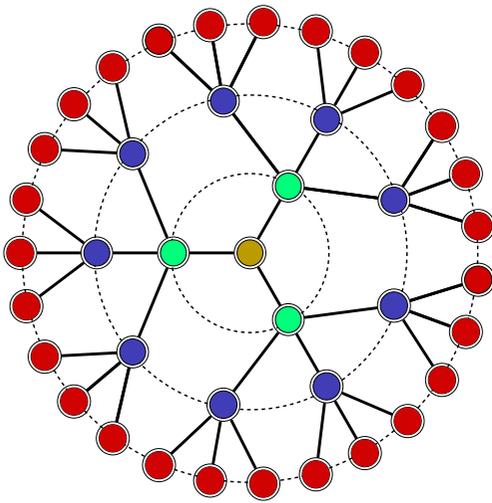


Figure 36. Bethe lattice graph: an example of symmetries in network structure. Nodes in the same level of the Bethe graph are exactly symmetric to each other. Therefore same colour nodes constitute a cluster.

Notes

1. We are using the uniform operator norm $\|U\| = \sup_{t \geq 0} \|U(t)\|$.
2. Or Rössler type oscillator where the phases are well defined and the coupling between chaotic amplitudes and phases are small.

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Appendices

Appendix 1. List of frequently used notions and abbreviations

Table of notions

| | |
|----------------------------|--|
| $ \cdot $ | absolute value |
| $\ \cdot\ $ | norm |
| δ_{ij} | Kroneker delta |
| G | graph |
| L | Laplacian matrix |
| A | adjacency matrix |
| I | identity matrix |
| H | coupling function |
| $\mathbf{1}$ | vector whose every components is 1 |
| α | coupling strength |
| α_c | critical coupling strength for synchronisation |
| f | isolated dynamics (vector field) |
| Λ | Maximum Lyapunov exponent |
| λ_2 | spectral gap the second minimum eigenvalue of Laplacian matrix |
| Df | Jacobian matrix of f |
| ϕ | phase |
| t | time |
| a, b and c | parameters of Rössler system |
| σ, ρ and β | parameters of Lorenz system |
| n | dimension of vector fields |
| i and j | natural numbers |
| N | system size of networks |
| M | total number of links |
| k_i | degree of i th node |
| d | diameter of a network |

Table of abbreviations

| | |
|----|------------------------------|
| CS | complete synchronisation |
| PS | phase synchronisation |
| GS | generalised synchronisation |
| AS | anticipating synchronisation |
| ER | Erdős–Renyi network |
| SF | Scale-free network |
| SW | Small world network |
| BA | Barabasi–Albert network |

Appendix 2. Lyapunov exponent

Sensitive dependence on initial conditions is one of the main characteristics of chaotic systems. The main idea is that nearby orbits diverge at an exponential rate. This rate is called Lyapunov exponents. In this Appendix, we provide the basic notions on the theory of Lyapunov exponents.

If we have a non-linear equation, we can study the properties of a given solution s by linearising the dynamics around the orbit, as we have done in Section 2.2. This procedure leads to a linear nonautonomous equation

$$\mathbf{v}' = A(t) \mathbf{v}$$

where A is continuous and bounded matrix function. The goal is to study the behaviour of solutions. Typically solving the equation explicitly is impossible. So the theory of Lyapunov exponents plays a major role.

Let $\mathbf{v} : \mathbb{R} \rightarrow \mathbb{R}^n$ be a solution $\mathbf{v}' = A(t) \mathbf{v}$ and $T(t, s)$ is the fundamental matrix. The Lyapunov exponent of the solution is defined as

$$\lambda(\mathbf{v}) = \overline{\lim}_{t \rightarrow \infty} \frac{1}{t} \ln \|T(t, s) \mathbf{v}(s)\|$$

We also define $\lambda(\mathbf{0}) = -\infty$. The largest Lyapunov exponents is our main object of study and is given by

$$\Lambda = \overline{\lim}_{t \rightarrow \infty} \frac{1}{t} \ln \|\Pi(t, s)\|$$

it determines the behaviour of solutions asymptotically because for any solution of the equation we have

$$\|\mathbf{v}(t)\| < C_\varepsilon e^{(\Lambda + \varepsilon)t}$$

If $\Lambda < 0$, the trivial solution $\mathbf{v}(t) = 0$ is asymptotically stable. Lyapunov exponents generalises stability criteria for autonomous (given by eigenvalues) and periodic equations (given by Floquet exponents).

Lemma B.1: Let $A \in \text{Mat}(n)$ and v be an eigenvector of $A\mathbf{v} = \beta\mathbf{v}$. Then $\lambda(\mathbf{v}) = \beta$.

If all $\lambda(\mathbf{v}) < 0$, we have $\max_v \{\lambda(\mathbf{v})\} = \Lambda < 0$, and the trivial solution is asymptotically stable. The Lyapunov exponent also generalises the Floquet exponents.

Lemma B.2: Let $A(t)$ be a periodic matrix by the Floquet representation we have $T(t, s) = P(t, s)e^{(t-s)Q(s)}$. Let \mathbf{v} be an eigenvector of $Q(s)$, then $\lambda(\mathbf{v})$ is an eigenvalue of $Q(s)$.

Hence, Lyapunov exponents are the eigenvalues of the monodromy matrix Q . Although, for the synchronisation analysis we care about the maximal Lyapunov exponents, it is important to know that there are at most n distinct Lyapunov exponents because the set $X = \{\mathbf{v}(t) \mid \lambda(\mathbf{v}) \leq \alpha\}$ is a vector space.

Appendix 3. Lyapunov function

One of the main techniques to tackle stability of non-linear system is the Lyapunov function method. The method by Lyapunov allows us to obtain the stability without finding the trajectories by studying properties of the Lyapunov function. We consider a dynamical system is modelled by a differential equation

$$\dot{\mathbf{x}} = \frac{d\mathbf{x}}{dt} = f(\mathbf{x}) \quad (\text{C1})$$

We will study notions relative to connected nonempty subsets Ω of \mathbb{R}^m . A function $V : \mathbb{R}^m \rightarrow \mathbb{R}$ is said to be positive definite with respect to the set B if $V(\mathbf{x}) > 0$ for all $\mathbf{x} \in \mathbb{R}^q \setminus \Omega$. It is radially unbounded if

$$\lim_{\|\mathbf{x}\| \rightarrow \infty} V(\mathbf{x}) = \infty.$$

Note that this condition guarantees that all level sets of V are bounded. This fact plays a central role in the analysis. We also define $V' : \mathbb{R}^m \rightarrow \mathbb{R}$ as

$$V'(\mathbf{x}) = \nabla V(\mathbf{x}) \cdot f(\mathbf{x}).$$

where \cdot denotes the Euclidean inner product. This definition agrees with the time derivative along the trajectories. That is, if $\mathbf{x}(t)$ is a solution of Equation (C1), then by the chain rule we have

$$\frac{dV(\mathbf{x}(t))}{dt} = V'(\mathbf{x}(t)).$$

Notice that applying the chain rule we obtain

$$V'(\mathbf{x}(t)) = \nabla V(\mathbf{x}(t)) \cdot f(\mathbf{x}(t))$$

This has a nice geometric interpretation. Since $\nabla V(\mathbf{x}(t))$ is perpendicular to the level set of V if $V'(\mathbf{x}(t)) < 0$ it means that the vector field is point inwards the level set and trajectories will enter the level set and never leave it. Repeating the argument we obtain stability as the following statement shows

Theorem 4: [Lyapunov] Let $V : \mathbb{R}^n \rightarrow \mathbb{R}$ be radially unbounded and positive definite with respect to the set $\Omega \subset D$. Assume that

$$V'(x) < 0 \quad \text{for all } x \in \mathbb{R}^n \setminus \Omega$$

Then all trajectories of Equation (C1) eventually enter the set Ω , in other words, the system is dissipative.

There are also converse Lyapunov theorems [131]. Typically if the system is dissipative (and have nice properties) then there exists a Lyapunov function. Although the above theorem is very useful, since we don't need knowledge of the trajectories, the drawback is the function V itself. There is no recipe to obtain a function V fulfilling all these properties. One could always try to guess the function, or go for a general form such as choosing a quadratic function V . We assume that the Lyapunov function is given.

Appendix 4. Chaos in Lorenz system

In order to understand the behaviour of a continuous system, we can use the concept of a Poincaré section – a transversal surface to the flow. This method was developed by Henri Poincaré in 1890s. The crossing points are a set of discrete numbers and this number sequence is called Poincaré map. We can study the structure of the crossings of the trajectory to the surface. This reduces the dimension of the system by 1. The structure of crossing points between the plane and the trajectory determines the behaviour of the system. For example, if the trajectory cross the section always at same k -coordinate points and repeat these points in the same order then the system is periodic so-called period- k .

The maxima of z -component of the Lorenz system, which is Poincaré section of velocities, graphically show the chaotic regime clearly. The governing equation of the Poincaré map ($\{z_n\}$) can be plotted as z_n vs z_{n+1} (Figure 1(a)) which resembles the tent map function (Figure 1(b)). The tent map is given by

$$f(x_n) = x_{n+1} = \begin{cases} 2x_n & 0 \leq x_n \leq 1/2 \\ 2 - 2x_n & 1/2 < x_n \leq 1. \end{cases}$$

Lyapunov exponent of the tent map

$$\Lambda = \lim_{t \rightarrow \infty} \frac{1}{t} \ln \|Df^t(x)\| \quad (\text{D1})$$

where Df is the Jacobian of f and $\|Df(x)\| = 2$ for all $x \neq 1/2$ since the function is not differentiable at $x = 1/2$. Therefore the Lyapunov exponent is $\Lambda = \ln 2$ and according to the positive Lyapunov exponent, the behaviour of the system is chaotic.

The definition of chaos given by Devaney is the following let X be a metric space and a continuous map $f : X \rightarrow X$ is chaotic if

- (1) f is transitive (indecomposability); that is any non-empty intervals $U, V \subset X$ there exist a natural number k such that $f^k(U) \cap V$. The transitivity condition ensures the irreducibility of X , means that X is an open invariant set and cannot split into two open invariant sets.
- (2) the periodic trajectories of f are dense in X (an element of regularity); that is every point in the attractor of f will return back to neighbourhood of the initial point. In other words, any trajectory initiating from an interval $I \subset X$, will be back in I in sufficiently long but finite time. Therefore the subset I contains infinitely many periodic points.

- (3) f has sensitive dependence on initial conditions (unpredictability); that is if there is an infinitesimal distance δ_0 between any two point $x, y \in X$ and there exists a non-negative number k such that after n iterations the distance between $f^n(x)$ and $f^n(y)$ is larger than $\delta_n > \delta_0$. These nearly started orbits diverge from each other at a rate Λ (see Appendix 7).

More details can be found in Ref. [132].

Appendix 5. Mathematical structure of generalised synchronisation

For completeness, we include this brief discussion of the mathematical structure of GS and it may be skipped without harm to the remaining sections. Again lets consider the $\psi : \mathbb{R}^m \rightarrow \mathbb{R}^n$ and the manifold

$$M = \{(x, y) \in \mathbb{R}^n \times \mathbb{R}^m : y = \psi(x)\} \subset \mathbb{R}^{n+m}.$$

GS corresponds to the case where M is normally attracting. Lets review the notion of normally attracting invariant manifold (NAIM). M is normally attracting if it is invariant under the flow Φ (of the full system) and the dynamics in the directions normal to M is contracting stronger than in direction tangential to M . That is,

- (1) (Invariance) $\Phi^t(M) = M$ for all t ,
- (2) (Normal Contraction) There is a continuous splitting of the tangent space

$$\forall x \in M : T_x \mathbb{R}^{n+m} = T_x M \oplus E_x^S$$

which is left invariant by $D\Phi^t$. Moreover, there is $\eta > 0$ such that

$$\forall t \geq 0, x \in M, v \in E_x^S : \|D\Phi^t(x)v\| \leq Ce^{-\eta} \|v\|$$

- (3) (Contraction in the normal directions dominate the tangential) There exist $r > 1, \lambda > 0$ and $C > 0$ such that

$$\begin{aligned} \forall t \geq 0, x \in M : \|D\Phi^t(x)\big|_{E_x^S}\| \\ \leq Ce^{-\lambda t} \|D\Phi^{-t}(x)\big|_{T_x M}\|^{-r} \end{aligned}$$

If M is a NAIM for the system F . Then there exist locally invariant stable manifolds $W_{loc}^s(M)$ such that $W_{loc}^s(M)$ is tangent to $TM \oplus E_s$ at M and $W_s(M) \in C^r$. Moreover, $W_{loc}^s(M)$ consists of all points near M whose forward orbit converges to M at rate $e^{-\eta t}$. For each $y \in W_s(M)$ shadows a point $x \in M$ such that $y \in W_{loc}^s(x)$ and

$$\|\Phi^t(y) - \Phi^t(x)\| \leq Ce^{-\eta t} \|y - x\| \quad (\text{E1})$$

Since the orbits of points $x_0, x_1 \in M$ cannot approach each other that fast, we can characterise points $y \in W_{loc}^s(x)$ precisely as those that satisfy Equation (E1). Lets consider the straightening of the manifold. That is, we introduce new coordinates

$$u = y - \psi(x)$$

in this coordinates the manifold corresponds to the x axis and u are the normal directions to M . Lets take two points $u_1, u_0 \in W_{loc}^s(x)$ then

$$u_i = y_i - \psi(x) \Rightarrow u_1 - u_0 = y_1 - y_0$$

Hence, if

$$\|y_1 - y_0\| \leq Ke^{-\eta t}$$

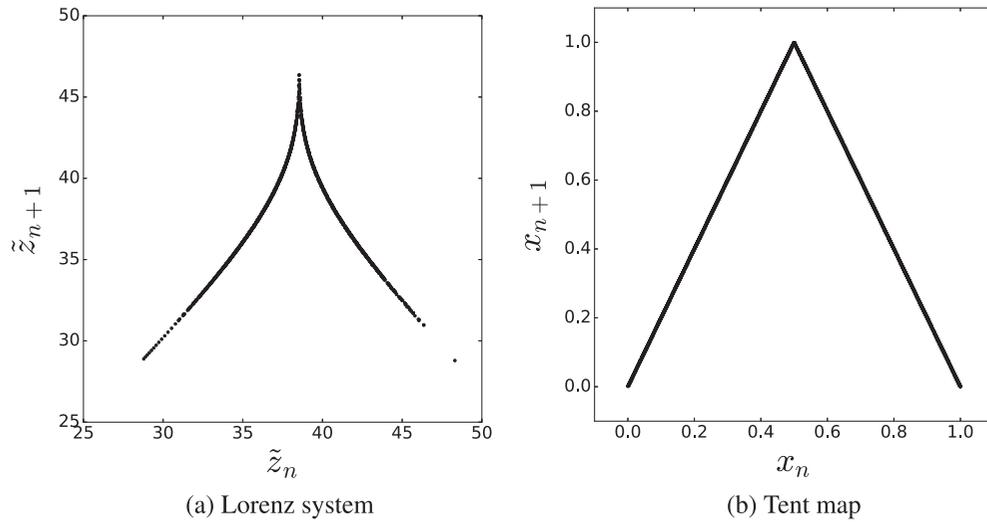


Figure 1. Similarity between Poincaré map of Lorenz system and the tent map.

and η is larger than the smallest Lyapunov exponents of the driver in modulus the manifold M will be normally attracting, according to condition (3). In fact, ψ will be differentiable. Another important fact to the mention is that NAIM persist under small perturbations. For us this means that once we obtain GS small perturbations such as increasing the coupling strength will be destroy GS [133,134]. If the condition is not satisfied then ψ won't be a NAIM. However, it may still happen that when $r = 0$ and M is attracting. In this case, ψ is only continuous. This is called strong and weak generalised synchronisation [135].

Back to our Diffusively driven oscillator. In Section ?? we showed the contraction rate between two nearby trajectories is

$$\eta = \alpha \lambda_{\min} - \|Dg\|$$

On the other hand, the smallest Lyapunov exponent of the driver is at most $-\|Df\|$, hence the condition for normal attraction is

$$\eta > \|Df\| \Rightarrow \alpha > \frac{\|Df\| + \|Dg\|}{\lambda_{\min}}$$

This gives the bound for M to be NAIM.