

Isostable reduction of periodic orbits

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(Received 1 May 2016; revised manuscript received 3 September 2016; published 14 November 2016)

The well-established method of phase reduction neglects information about a limit-cycle oscillator's approach towards its periodic orbit. Consequently, phase reduction suffers in practicality unless the magnitude of the Floquet multipliers of the underlying limit cycle are small in magnitude. By defining isostable coordinates of a periodic orbit, we present an augmentation to classical phase reduction which obviates this restriction on the Floquet multipliers. This framework allows for the study and understanding of periodic dynamics for which standard phase reduction alone is inadequate. Most notably, isostable reduction allows for a convenient and self-contained characterization of the dynamics near unstable periodic orbits.

DOI: [10.1103/PhysRevE.94.052213](https://doi.org/10.1103/PhysRevE.94.052213)

Oscillatory behavior is a ubiquitous natural phenomenon with a wide array of applications, including neurological behavior, circadian rhythms, chemical reactions, mechanical vibrations, and chaotic systems [1–7]. For nearly half a century, phase reduction has been an indispensable tool to aid in the understanding and manipulation of such oscillators [1–6] by allowing periodic solutions of

$$\dot{\mathbf{x}} = F(\mathbf{x}) + G(\mathbf{x}, t), \quad \mathbf{x} \in \mathbb{R}^n \quad (1)$$

to be represented by the single variable system:

$$\dot{\theta} = \omega + Q(\theta)^T G(\mathbf{x}, t). \quad (2)$$

Here $G \in \mathbb{R}^n$ is an external perturbation, $\theta \in [0, 2\pi)$ is the phase of oscillation, $\omega = 2\pi/T$ is the natural frequency with T being the natural period, and $Q(\theta) \in \mathbb{R}^n$ is an infinitesimal phase response curve (PRC), which as its name suggests, is valid for perturbations G with small magnitude. Practically, (2) is valid in a close vicinity of the periodic orbit, γ . Consequently, the amplitude of allowable perturbations is limited by the size of the Floquet multipliers [8]; stable orbits with Floquet multipliers with magnitude close to 1 can only admit relatively small perturbations without the risk of being driven away from the limit cycle over time. In many applications [5, 9–11], however, the efficacy of a given control strategy is directly related to the magnitude of allowable perturbations. Furthermore, for unstable periodic orbits, (2) alone cannot adequately describe the long-term behavior of (1), rendering it unusable. Discrete-time proportional state feedback [7, 12] and delayed feedback control [13] have been used to suppress chaos by means of stabilizing unstable periodic orbits. These classes of chaos control strategies have found practical applications in cardiac, electronic, networked, and optical systems [14–17]. Phase reduction techniques could be of practical interest in the suppression of chaos if appropriate modifications could be made.

This work proposes an augmentation of standard phase reduction which allows for a better understanding of systems with either a stable or unstable periodic orbit. We introduce a set of isostable coordinates and associated isostable response curves (IRC), which represent the distance from the periodic orbit in an appropriate basis. The notion of isostables was introduced in Ref. [18] (cf. Refs. [19] and [20]) to represent sets of initial conditions which converge toward a stable fixed point together in a well-defined sense. The isostable coordinates

used in this manuscript are defined in a manner that is similar in spirit but adapted for use with periodic orbits. Unlike Refs. [21, 22], the strategy proposed here does not require computationally intensive calculations of an *ad hoc* coordinate system with respect to a periodic orbit of dimensionality greater than 2. Furthermore, the strategy presented here does not require the periodic orbit to be exponentially attracting in either forward or backward time. In the examples given in this work, we illustrate the utility of this reduction method in two different oscillatory systems for which (2) alone is insufficient.

Starting with a general system of ordinary differential equations (1), let γ be a T -periodic orbit which exists for $G \equiv 0$. We define a scalar phase variable on γ such that $\theta(\mathbf{x}) : \mathbb{R}^n \rightarrow [0, 2\pi)$ for which $d\theta[\mathbf{x}(t)]/dt = \omega$ and $\theta[\mathbf{x}(t)] = \theta[\mathbf{x}(t + T)]$, and choose an arbitrary point $\mathbf{x}_0 \in \gamma$ for which $\theta(\mathbf{x}_0) = 0$. By solving for $Q(\theta)$ (the gradient of the phase field) using, e.g., adjoint methods [4, 23, 24], this notion of phase can be extended to any \mathbf{x} in a neighborhood of γ by noting that for $\epsilon > 0$, $\theta(\mathbf{x} + \epsilon\mathbf{y}) = \theta(\mathbf{x}) + \mathcal{O}(\epsilon^2)$ for any \mathbf{y} in the null space of $Q[\theta(\mathbf{x})]$. This definition of phase is possible regardless of the stability type of γ . We define isochrons as level sets of the phase field, i.e., $\Gamma_\theta = \{\mathbf{x} | \theta(\mathbf{x}) = \theta\}$. When γ is a stable periodic orbit, isochrons have an intuitive meaning: For any initial conditions $\mathbf{a}(0) \in \gamma$ and $\mathbf{b}(0)$ in the basin of attraction of γ on the same isochron, $\lim_{t \rightarrow \infty} \|\mathbf{a}(t) - \mathbf{b}(t)\| = 0$ [1].

Changing to phase variables using the chain rule, one arrives at the phase reduction (2). Here we are also interested in the transient behavior of (1) near γ which can be understood in terms of Poincaré maps. By construction, any initial condition in Γ_0 first returns to Γ_0 at time T allowing for the definition of a Poincaré map,

$$P : \Gamma_0 \rightarrow \Gamma_0; \quad \mathbf{x} \mapsto \phi(\mathbf{x}), \quad (3)$$

with $P(\mathbf{x}_0) = \mathbf{x}_0$. In a small neighborhood of \mathbf{x}_0 we may approximate ϕ from (3) as

$$\phi(\mathbf{x}) = \mathbf{x}_0 + J_P(\mathbf{x} - \mathbf{x}_0) + \mathcal{O}(\|\mathbf{x} - \mathbf{x}_0\|^2), \quad (4)$$

where $J_P = d\phi/d\mathbf{x}|_{\mathbf{x}_0}$. Suppose J_P is diagonalizable and let $V \in \mathbb{R}^{n \times n}$ be a matrix with columns that form a basis of unit length eigenvectors $\{\mathbf{v}_k, k = 1, \dots, n\}$ of J_P with associated eigenvalues $\{\lambda_k, k = 1, \dots, n\}$. The eigenvalues λ_i are often referred to as Floquet multipliers of the periodic orbit. For any eigenvector v_i with associated $\lambda_i \in \mathbb{R} > 0$ (if any $\lambda_i < 0$, one can define the period to be $2T$ so all eigenvalues are positive),

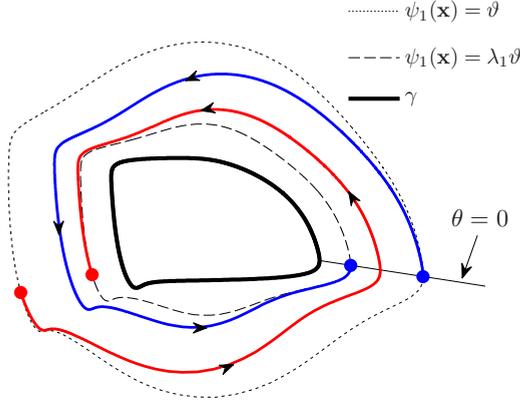


FIG. 1. A sketch of the behavior of a general two-dimensional system near its limit cycle. The red and blue lines represent two trajectories integrated over one period which start on the same isostable level set. Dashed and dotted lines represent two different isostable level sets. When the Poincaré map defined with respect to the $\theta = 0$ isochron is well approximated as a linear map the discontinuity across the $\theta = 0$ isochron is negligible, and trajectories on the same isostable level set cross the successive isostable level sets together. Furthermore, the isostable coordinate decreases at an exponential rate governed by the Floquet multiplier λ_1 . Here the isostable level sets give a sense of the distance from the periodic orbit.

we can define a set of isostable coordinates,

$$\psi_i(\mathbf{x}) = \mathbf{e}_i^T V^{-1}(\mathbf{x}_\Gamma - \mathbf{x}_0) \exp(-\log(\lambda_i)t_\Gamma/T). \quad (5)$$

Here \mathbf{x}_Γ and $t_\Gamma \in [0, T)$ are defined to be the location and time, respectively, at which the trajectory first returns to Γ_0 under the flow $\dot{\mathbf{x}} = F(\mathbf{x})$ and \mathbf{e}_i is a vector with 1 in the i th position and zeros elsewhere. These isostable coordinates $\psi_i(\mathbf{x})$ are defined for all $\mathbf{x} \in \mathbb{R}^n$, not just on the Poincaré section. Intuitively, a trajectory near the periodic orbit will spiral towards or away at a rate determined by each of the Floquet multipliers. This growth (decay) is matched by the decay (growth) of the exponential term from (5), giving a sense of the distance in directions transverse to the periodic orbit. Noting that $dt_\Gamma/dt = -1$ for $\mathbf{x}(t) \notin \Gamma_0$, by direct differentiation of (5) $d\psi_i/dt = \psi_i \log(\lambda_i)/T$, therefore,

$$\nabla \psi_i(\mathbf{x}) \cdot F(\mathbf{x}) = \psi_i \log(\lambda_i)/T. \quad (6)$$

By (4), one can verify that under the flow $\dot{\mathbf{x}} = F(\mathbf{x})$, $\psi_i[\mathbf{x}(t_\Gamma^+)] = \psi_i[\mathbf{x}(t_\Gamma^-)] + \mathcal{O}(\|\mathbf{x} - \mathbf{x}_0\|^2)$. This discontinuity results from the approximation of (3) as a linear mapping. Throughout this analysis, we will assume close-enough proximity to the periodic orbit so this discontinuity is negligible. Figure 1 gives a representation of a general two-dimensional system in terms of its isostable coordinates. For an arbitrary time τ , if we compute the isostable coordinate $\psi_1[\mathbf{x}(\tau)] = \vartheta$, after one revolution, $\psi_1[\mathbf{x}(\tau + T)] = \lambda_1 \vartheta + \mathcal{O}(\|\mathbf{x} - \mathbf{x}_0\|^2)$. Furthermore, close to the periodic orbit, when the Poincaré map (3) is well approximated by a linear mapping, two initial conditions on the same isostable level set will cross successive isostable level sets together on their way to the limit cycle (to leading order)

Letting $\kappa_i \equiv \log(\lambda_i)/T$, and changing variables to isostable coordinates, with the chain rule we find

$$\begin{aligned} \frac{d\psi_i(\mathbf{x})}{dt} &= \nabla \psi_i(\mathbf{x}) \cdot [F(\mathbf{x}) + G(\mathbf{x}, t)] \\ &= \kappa_i \psi_i(\mathbf{x}) + \nabla \psi_i(\mathbf{x}) \cdot G(\mathbf{x}, t). \end{aligned} \quad (7)$$

Evaluating the vector field at $\mathbf{x}^\gamma(\theta, \psi_i)$, which we define as the intersection of the trajectory γ , the $\psi_i(\mathbf{x})$ level set (i.e., isostable), and the $\theta(\mathbf{x})$ level set (i.e. isochron), we have

$$\frac{d\psi_i(\mathbf{x})}{dt} = \kappa_i \psi_i(\mathbf{x}) + \nabla \psi_i[\mathbf{x}^\gamma(\theta, \psi_i)] \cdot G(\mathbf{x}^\gamma(\theta, \psi_i), t). \quad (8)$$

Here, as in Ref. [25] we have ignored an $\mathcal{O}(|G|^2)$ term so (8) is valid for perturbations with small G .

Towards deriving an equation for the numerical computation of $\nabla \psi_i[\mathbf{x}^\gamma(\theta, \psi_i)]$, we will momentarily take $G \equiv \mathbf{0}$ and consider the effect of a small perturbation $\Delta \mathbf{x}$ to a trajectory $\mathbf{x}(t) \in \gamma$. $\Delta \mathbf{x}(t)$ evolves according to

$$\frac{d\Delta \mathbf{x}(t)}{dt} = J[\mathbf{x}(t)]\Delta \mathbf{x}(t) + \mathcal{O}(\|\Delta \mathbf{x}\|^2), \quad (9)$$

where $J[\mathbf{x}(t)]$ is the Jacobian matrix evaluated at $\mathbf{x}(t)$. The corresponding isostable shift, $\Delta \psi_i \equiv \psi_i[\mathbf{x}(t) + \Delta \mathbf{x}(t)] - \psi_i[\mathbf{x}(t)]$, is given by

$$\Delta \psi_i = \nabla_{\mathbf{x}(t)} \psi_i \cdot \Delta \mathbf{x}(t) + \mathcal{O}(\|\Delta \mathbf{x}\|^2), \quad (10)$$

where $\nabla_{\mathbf{x}(t)} \psi_i$ is the gradient of ψ_i evaluated at $\mathbf{x}(t) \in \gamma$. After the initial perturbation at $t = 0$,

$$\begin{aligned} \frac{d\Delta \psi_i}{dt} &= \kappa_i \Delta \psi_i \\ &= \kappa_i \nabla_{\mathbf{x}(t)} \psi_i \cdot \Delta \mathbf{x}. \end{aligned} \quad (11)$$

In the spirit of Refs. [24] and [26], by taking the time derivative of (10) and rearranging, to lowest order in $\|\Delta \mathbf{x}\|$,

$$\begin{aligned} \langle d\nabla_{\mathbf{x}(t)} \psi_i / dt, \Delta \mathbf{x}(t) \rangle & \\ &= -\langle \nabla_{\mathbf{x}(t)} \psi_i, d\Delta \mathbf{x}(t)/dt \rangle + \langle \kappa_i \nabla_{\mathbf{x}(t)} \psi_i, \Delta \mathbf{x} \rangle \\ &= -\langle \nabla_{\mathbf{x}(t)} \psi_i, J[\mathbf{x}(t)]\Delta \mathbf{x} \rangle + \langle \kappa_i \nabla_{\mathbf{x}(t)} \psi_i, \Delta \mathbf{x} \rangle \\ &= -\langle J[\mathbf{x}(t)]^T \nabla_{\mathbf{x}(t)} \psi_i - \kappa_i \nabla_{\mathbf{x}(t)} \psi_i, \Delta \mathbf{x} \rangle. \end{aligned} \quad (12)$$

Here $\langle \cdot, \cdot \rangle$ is the Euclidean inner product (i.e., dot product), and superscript T indicates the transpose (i.e., adjoint) of the real-valued matrix $J[\mathbf{x}(t)]$. Because (12) is valid for any $\Delta \mathbf{x}$,

$$\frac{d\nabla_{\mathbf{x}(t)} \psi_i}{dt} = \{\kappa_i I - J[\mathbf{x}(t)]^T\} \nabla_{\mathbf{x}(t)} \psi_i, \quad (13)$$

where I is the identity matrix. Note the similarity between (13) and the adjoint equation derived in Ref. [21] which was valid only for two-dimensional systems. Recall that for $\mathbf{x} \in \Gamma_0$, $t_\Gamma = 0$ and $\mathbf{x}_\Gamma = \mathbf{x}$, so from (5), $\psi_i(\mathbf{x}) = \mathbf{e}_i^T V^{-1}(\mathbf{x} - \mathbf{x}_0)$. This implies that $\psi_i(\mathbf{x} + \eta \mathbf{v}_i) = \psi_i(\mathbf{x}) + \eta$ for $\eta \in \mathbb{R}$ or, equivalently,

$$\nabla_{\mathbf{x}_0} \psi_i \cdot \mathbf{v}_i = 1. \quad (14)$$

This normalization condition along with T -periodicity defines a unique solution of (13). Equation (1) can then be understood

in reduced form

$$\begin{aligned}\dot{\theta} &= \omega + Q^T(\theta) \cdot G(t) \\ \dot{\psi}_i &= \kappa_i \psi_i + \mathcal{I}_i^T(\theta) \cdot G(t) \text{ for } i = 1, \dots, n-1,\end{aligned}\quad (15)$$

where $Q^T(\theta) \equiv \nabla \theta|_{x^\nu(\theta)}$ is often referred to as the PRC, and $\mathcal{I}_i(\theta) \equiv \nabla \psi_i|_{x^\nu(\theta)}$ will be referred to as an IRC. As shown in Appendix, the magnitude but not the shape of the resulting IRCs depends on the initial choice of $\theta(\mathbf{x}_0) = 0$. Much like in the standard phase reduction (2), Eqs. (15) are valid provided the unreduced state dynamics remain close to γ . In practice, we generally only need to consider a few isostable coordinates: If $|\lambda_k| \approx 0$, any perturbations to ψ_k will be quickly forgotten and this coordinate can simply be ignored. Finally, we note that if J_p is not diagonalizable, a similar reduction analysis can be performed for any eigenvalue for which the geometric and algebraic multiplicity are identical. As we will show in the following examples, isostable reduction is essential for understanding phase reduced systems with eigenvalues close to or greater than 1.

To illustrate the principles derived above, we will first consider a three-dimensional model of gene regulation [27] which has been used to describe the oscillatory behavior of the suprachiasmatic nucleus responsible for the mammalian circadian clock:

$$\begin{aligned}\dot{X} &= v_1 K_1^n / (K_1^n + Z^n) - v_2 X / (K_2 + X) + L(t), \\ \dot{Y} &= k_3 X - v_4 Y / (K_4 + Y), \\ \dot{Z} &= k_5 Y - v_6 Z / (k_6 + Z).\end{aligned}\quad (16)$$

Here X , Y , and Z represent concentrations (expressed in nM) of the messenger ribonucleic acid (mRNA) clock gene *per* or *cry*, the PER or CRY protein, and the nuclear form of the protein, respectively, with all constants taken as the nominal values from Fig. 2 of Ref. [27], and $L(t)$ represents a perturbation from ambient light. Figure 2(a) shows the

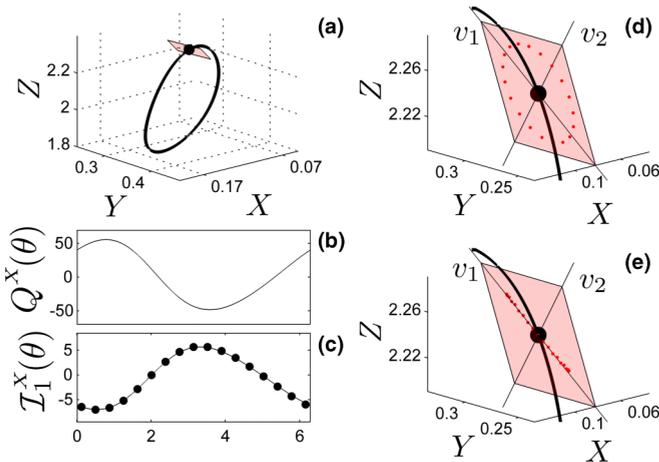


FIG. 2. Panel (a) shows the limit cycle in black with Γ_0 in red. Panels (b) and (c) show the PRC and IRC, respectively. Numerical validation of the IRC through calculation of $\Delta\psi_1/\Delta X$ (represented by black dots) for various phases of perturbation are shown with black dots. Initial conditions (red dots) in panel (d) are mapped to locations in panel (e) after one iteration of the Poincaré map.

limit-cycle solution of (16) when $L(t) = 0$ with a natural period $T_c = 23.54$ h. The value $\theta = 0$ (an arbitrary reference point) is represented with a black dot, and the Γ_0 Poincaré section (i.e., the $\theta = 0$ isochron) is approximated by the red plane near the periodic orbit. Initial conditions represented by red dots in Fig. 2(d) are mapped to the locations in Fig. 2(e). Eigenvalues λ_1 and λ_2 of eigendirections v_1 and v_2 of $P(\mathbf{x})$ are determined numerically to be 0.951 and approximately 0, respectively. Using standard techniques [4,23,24], the PRC, $Q^c(\theta) \equiv [Q^X(\theta) Q^Y(\theta) Q^Z(\theta)]$, is calculated with $Q^X(\theta)$ shown in Fig. 2(b). The IRC, $\mathcal{I}_i^c(\theta) = [\mathcal{I}_1^X(\theta) \mathcal{I}_1^Y(\theta) \mathcal{I}_1^Z(\theta)]$, is calculated using (13) with $\mathcal{I}_1^X(\theta)$ shown in Fig. 2(c).

We can represent the reduced dynamics of this oscillator with two coupled ordinary differential equations:

$$\dot{\theta} = \omega_c + Q^X(\theta)L(t), \quad (17)$$

$$\dot{\psi}_1 = \kappa_c \psi_1 + \mathcal{I}_1^X(\theta)L(t), \quad (18)$$

with $\omega_c = 2\pi/T_c$ and $\kappa_c = \log(\lambda_1)/T_c$. Note that we do not include ψ_2 in the reduction because the stability in this coordinate is very strong. To illustrate the necessity of (18) in the phase reduction, we will test and implement a simple control strategy for entrainment to an external perturbation. Suppose we would like to entrain the oscillation to an external periodic perturbation $L(t) = \mu\delta[\text{mod}(t, T_c + \Delta T)]$, a control objective which is relevant in the treatment in certain types of circadian misalignment (cf. Refs. [28,29]). By simply taking into account the phase reduction (17), one can understand the phase dynamics as a series of maps

$$\bar{\theta}^+ = \bar{\theta} + 2\pi\Delta T/T_c + \mu Q^X(\bar{\theta}), \quad (19)$$

where $\bar{\theta}$ represents the phase immediately prior to the application of a pulsatile stimulus, and $\bar{\theta}^+$ gives the phase at a time $T_c + \Delta T$ later. Equation (19) has a fixed point when $\mu = -2\pi\Delta T/T_c Q^X(\bar{\theta})$. The minimal control effort required occurs when $\theta = \text{argmax}|Q^X(\theta)| \approx 1$. Letting $T_c + \Delta T = 22.2$ h, simple stability analysis reveals that the resulting fixed point is stable. However, as shown in the right panels of Fig. 3, this control strategy does not give stable entrainment and the mean period remains unchanged. From the three-dimensional trajectory in blue, we find that the trajectory does not remain close to the periodic orbit, as would be predicted from the isostable reduction, as $\mathcal{I}_1^X(1)$ is relatively large. Instead, if

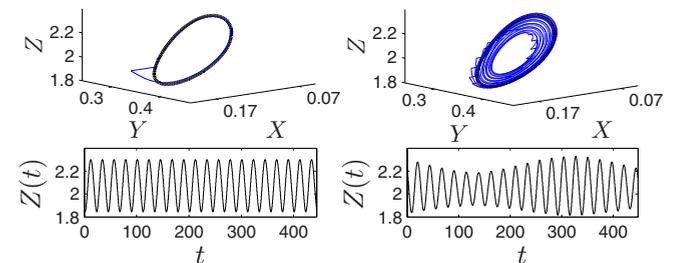


FIG. 3. Left: The circadian oscillator can be entrained (trajectory shown in blue) to periodic δ -function perturbations if they are given when the IRC is approximately zero. Right: Entrainment does not occur even when analysis of (17) alone predicts it should. Equation (18) is essential for using phase reduction in this application.

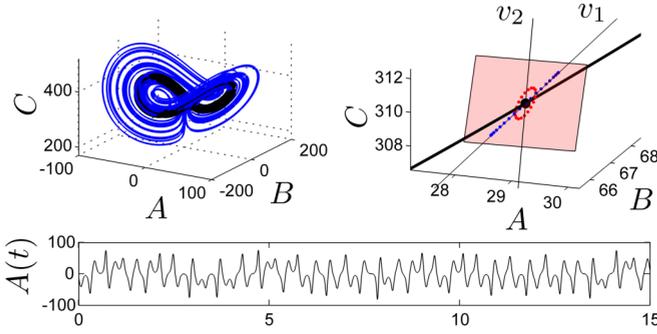


FIG. 4. Top-left: Trajectory along the Lorenz attractor in blue with an unstable periodic orbit outlined in black. Top-right: A small portion of the periodic orbit is shown as a thick black line. Eigendirections of the Poincaré map are shown as thin lines. Initial conditions in red are mapped to the locations in blue after one application of the map. Bottom: an example uncontrolled chaotic trajectory on the Lorenz attractor.

we choose μ as above, numerically we find $\bar{\theta} = 2$ is a stable fixed point of (19), and with $\mathcal{I}_1^X(1) = 0$, the pulsatile stimulus stably entrains the oscillator as shown in the left panels of Fig. 3. Numerically, we find that fixed points $\bar{\theta}_0$ of (17) are unstable unless $\mathcal{I}(\bar{\theta}_0) \approx 0$.

Next, we show that the notion of isostables can be used to understand the effect of small perturbations near an unstable periodic orbit and how they can be used to optimally drive a given trajectory to an unstable periodic orbit. Here we consider the three-dimensional Lorenz equations [30]:

$$\begin{aligned}\dot{A} &= \sigma(B - A) + u(t), \\ \dot{B} &= A(\rho - C) - B, \\ \dot{C} &= AB - \beta C.\end{aligned}\quad (20)$$

Here, A , B , and C are nondimensional system variables; $\sigma = 16$, $\beta = 8/3$, and $\rho = 350$ are constants chosen so the unperturbed dynamics are chaotic; and $u(t)$ is an external perturbation. An unstable periodic orbit, γ_L with period $T_L = 0.372$ and dynamics along the Lorenz attractor are shown in Fig. 4. Along this orbit, $\theta = 0$ is taken to correspond to an arbitrary location along γ_L allowing for the definition of Γ_0 . The resulting Poincaré section is shown in the top-right panel. Eigenvalues λ_1 and λ_2 of the Poincaré map's v_1 and v_2 eigendirections are numerically determined to be 3.33 and approximately 0, respectively. Using (13) the IRC of this unstable eigendirection $\mathcal{I}_1^L(\theta) = [\mathcal{I}_1^A(\theta) \mathcal{I}_1^B(\theta) \mathcal{I}_1^C(\theta)]$ is calculated with $\mathcal{I}_1^A(\theta)$ shown in Fig. 5. We also calculate the PRC $Q^A(\theta)$ for perturbations in the A direction, allowing for the reduction

$$\begin{aligned}\dot{\theta} &= \omega_L + Q^A(\theta)u(t), \\ \dot{\psi}_1 &= \kappa_L \psi_1 + \mathcal{I}_1^A(\theta)u(t),\end{aligned}\quad (21)$$

where $\omega_L = 2\pi/T_L$ and $\kappa_L = \log(\lambda_1)/T_L$. Understanding the Lorenz system in reduced coordinates allows for the formulation of a control problem to drive any initial condition to the unstable periodic orbit provided it is close enough to the orbit so the reduction is valid. Towards this calculus of variations problem formulation [31], we define a cost

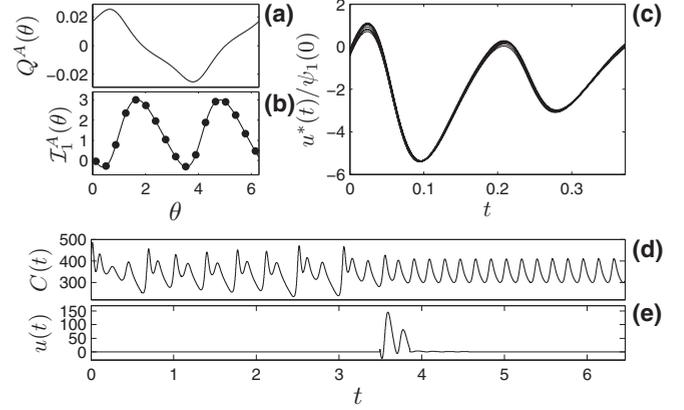


FIG. 5. Panels (a) and (b) give the PRC and IRC for this system. IRC values are numerically validated by calculating $\Delta\psi_1/\Delta A$ at various phases and shown as black dots. Panel (c) shows an optimal stimulus which takes an initial condition in Γ_0 and returns it to Γ_0 with $\psi_1 = 0$ one period later. A control strategy using this information is shown in panels (d) and (e).

functional

$$\begin{aligned}\mathcal{M}[\dot{\Phi}, \Phi, u(t)] &= \int_0^{T_L} \left[u^2(t) + \zeta_1 \{ \dot{\theta} - \omega_L - Q^A(\theta)u(t) \} \right. \\ &\quad \left. + \zeta_2 \{ \dot{\psi}_1 - \kappa_L \psi_1 - \mathcal{I}_1^A(\theta)u(t) \} \right] dt,\end{aligned}\quad (22)$$

with $\Phi(t) = [\theta(t), \psi_1(t), \zeta_1(t), \zeta_2(t)]$. Here, Lagrange multipliers ζ_1 and ζ_2 force the dynamics to satisfy the phase and isostable reduced equations. The associated Euler-Lagrange equations are

$$\frac{\partial \mathcal{M}}{\partial u} = \frac{d}{dt} \left(\frac{\partial \mathcal{M}}{\partial \dot{u}} \right); \quad \frac{\partial \mathcal{M}}{\partial \Phi} = \frac{d}{dt} \left(\frac{\partial \mathcal{M}}{\partial \dot{\Phi}} \right).\quad (23)$$

Optimal solutions to the cost function satisfy (23) with boundary conditions $\theta(0) = 0$, $\theta(T_L) = 2\pi$, $\psi_1(T_L) = 0$, and $\psi_1(0)$ determined from initial data. This two-point boundary problem can be solved, e.g., with a double bisection algorithm and is chosen so after one cycle, the trajectory ends on the stable manifold of the Poincaré section Γ_0 . The numerically determined optimal control $u^*(t)$ is shown in Fig. 5(c) for multiple choices of $\psi_1(0) \in [-20, 20]$. In this range, $u^*(t)$ is approximately proportional to $1/\psi_1(0)$. Using this information, we can devise a control algorithm to drive the Lorenz system to the unstable periodic orbit: For any initial condition, wait until the trajectory crosses Γ_0 close enough to the unstable periodic orbit and calculate $\psi_1(0)$ to determine $u(t)$. Figures 5(d) and 5(e) show the result of this strategy where the control is set to engage when $\psi_1(0) < 20$. After the first control application at $t \approx 3.5$, the system is nearly driven to the periodic orbit, but because (13) provides an approximation to the IRC, a second and third control application (of rapidly decreasing magnitude) are required to bring the system exactly to the periodic orbit.

The chaos control strategy illustrated here differs from discrete time proportional feedback methods (for example, Refs. [7,12]) in that it uses information at all locations near the periodic orbit, not just along the Poincaré surface. This added

information about the system allows for the definition of an optimal control problem which yields a continuous solution. In many control applications, allowing for continuous wave forms can lead to a substantial decrease in energy required to achieve a control objective as compared to discrete time control [5,32]. One drawback of the method developed here is that the periodic orbit must be known before IRCs can be measured and control strategies can be implemented. In this case, in an experimental setting one might envision that the unstable orbit could be found, for instance, with delayed feedback control strategies [13], after which an IRC could be measured with an experimental protocol similar to the “direct method” [33,34] for measuring PRCs. Such a protocol could make a chaos control described here more feasible in an experimental setting.

In summary, we have developed a set of isostable coordinates which allow classical phase reduction to be useful for understanding the dynamics near periodic orbits without Floquet multipliers near 0. The computational complexity of implementing this reduction strategy is comparable to that of standard phase-reduction calculations. The reduced dynamics are particularly useful for the problem of stabilizing unstable periodic orbits as a means of controlling chaotic dynamical systems [7,12–17]. Additional applications of this reduction strategy could include investigating effect noise on a system in the dynamics in directions transverse to the periodic orbit (similar to how the variance of the firing rate of an oscillator was studied in Ref. [35]). Furthermore, it could be of interest to use this isostable coordinate system to investigate how the phase response characteristics of an oscillator change when perturbed from the periodic orbit, which may have applications to memory effects from pacing history [36,37]. Continued development of strategies for understanding isostable coordinates of periodic orbits will assist in the understanding of these systems in problems pertaining to synchronization, entrainment, and stabilization of oscillatory dynamics when (2) alone is insufficient.

ACKNOWLEDGMENTS

Support for this work by National Science Foundation Grants No. NSF-1363243, No. NSF-1264535, and No. NSF-1602841 is gratefully acknowledged.

APPENDIX: RELATIONSHIP BETWEEN THE CHOICE OF POINCARÉ SECTION AND THE RESULTING ISOSTABLE RESPONSE CURVE

In general, the choice of $\theta = 0$ on the periodic orbit used to define the initial Poincaré section will change the isostable coordinates and will affect resulting isostable response curves. However, the shape of the isostable response curve for any isostable field defined as in (5) from the main text is invariant to the choice of θ for which we use to define our Poincaré section provided the algebraic multiplicity of the associated eigenvalue λ_i is unity. This will be shown here by examining the relationship between isostable changes caused by arbitrary perturbations in coordinate systems defined by two different Poincaré sections.

To begin let γ be a T -periodic solution of the vector field

$$\dot{\mathbf{x}} = F(\mathbf{x}) + G(\mathbf{x}, t), \quad \mathbf{x} \in \mathbb{R}^n, \quad (\text{A1})$$

where $F \in \mathbb{R}^n$ represents the unperturbed dynamics and $G \in \mathbb{R}^n$ is an external perturbation. As in the main text, we define a scalar phase variable $\theta(\mathbf{x}) : \mathbb{R}^n \rightarrow [0, 2\pi)$ for which $d\theta[\mathbf{x}(t)]/dt = \omega$ and $\theta[\mathbf{x}(t)] = \theta[\mathbf{x}(t + T)]$. We also define isochrons to be level sets of the phase field, i.e., $\Gamma_\theta = \{\mathbf{x} | \theta(\mathbf{x}) = \theta\}$.

Let \mathbf{x}_1 and \mathbf{x}_2 be any two points on γ such that $\theta(\mathbf{x}_1) = \theta_1$ and $\theta(\mathbf{x}_2) = \theta_2$. The associated isochrons, Γ_{θ_1} and Γ_{θ_2} , can be used to define Poincaré maps

$$P_1 : \Gamma_{\theta_1} \rightarrow \Gamma_{\theta_1} \\ \mathbf{x} \mapsto \phi_1(\mathbf{x}), \quad (\text{A2})$$

$$P_2 : \Gamma_{\theta_2} \rightarrow \Gamma_{\theta_2} \\ \mathbf{x} \mapsto \phi_2(\mathbf{x}). \quad (\text{A3})$$

In the analysis to follow, we will restrict our attention to the dynamics of these maps in a small neighborhood of \mathbf{x}_1 and \mathbf{x}_2 . Taylor expanding, one can show that for any time t_1 (respectively, t_2) for which $\mathbf{x}(t_1) \in \Gamma_{\theta_1}$ (respectively, $\mathbf{x}(t_2) \in \Gamma_{\theta_2}$),

$$\mathbf{x}(T + t_1) - \mathbf{x}_1 = J_1(\mathbf{x}(t_1) - \mathbf{x}_1) + \mathcal{O}(\|\mathbf{x}(t_1) - \mathbf{x}_1\|^2), \quad (\text{A4})$$

$$\mathbf{x}(T + t_2) - \mathbf{x}_2 = J_2(\mathbf{x}(t_2) - \mathbf{x}_2) + \mathcal{O}(\|\mathbf{x}(t_2) - \mathbf{x}_2\|^2),$$

where $J_1 = d\phi_1/d\mathbf{x}|_{\mathbf{x}_1}$ and $J_2 = d\phi_2/d\mathbf{x}|_{\mathbf{x}_2}$. By construction of the phase field, Γ_{θ_2} is the image of Γ_{θ_1} under the flow of the vector field (A1) with $G \equiv 0$. As discussed in Chapter 11 of Ref. [38], this defines a C^r diffeomorphism

$$h : \Gamma_{\theta_1} \rightarrow \Gamma_{\theta_2} \\ \mathbf{x} \mapsto \xi(\mathbf{x}) \quad (\text{A5})$$

and implies that the eigenvalues of J_1 are equal to the eigenvalues of J_2 . Note that $h(\mathbf{x}_1) = \mathbf{x}_2$ so in a small neighborhood of \mathbf{x}_1 , we can approximate (A5) as

$$h(\mathbf{x}) = \mathbf{x}_2 + H(\mathbf{x} - \mathbf{x}_1) + \mathcal{O}(\|\mathbf{x} - \mathbf{x}_1\|^2), \quad (\text{A6})$$

where $H = d\xi/d\mathbf{x}|_{\mathbf{x}_1}$.

Towards the definition of isostables with respect to each map, suppose the J_1 and J_2 are both diagonalizable. Let V_1 (respectively, V_2) $\in \mathbb{R}^{n \times n}$ be matrices with columns that form an orthonormal basis of unit length eigenvectors $\{\mathbf{v}_k^1, k = 1, \dots, n\}$ of J_1 (respectively, $\{\mathbf{v}_k^2, k = 1, \dots, n\}$ of J_2) associated with the eigenvalues $\{\lambda_k, k = 1, \dots, n\}$.

Let \mathbf{e}_i be a vector with 1 in the i th position and zeros elsewhere. Mirroring the definition used in the main text, we define two isostable fields ψ_i^1 and ψ_i^2 with respect to the Poincaré sections Γ_{θ_1} and Γ_{θ_2} :

$$\psi_i^j(\mathbf{x}) = \mathbf{e}_i^T V_j^{-1}(\mathbf{x}_{\Gamma_j} - \mathbf{x}_j) \exp(-\log(\lambda_i)t_{\Gamma_j}/T), \\ j = 1, 2, \quad (\text{A7})$$

where \mathbf{x}_{Γ_1} and t_{Γ_1} (respectively, \mathbf{x}_{Γ_2} and t_{Γ_2}) are the location and time, respectively, at which the trajectory under the flow $\dot{\mathbf{x}} = F(\mathbf{x})$ next returns to Γ_{θ_1} (respectively, Γ_{θ_2}). Here, the superscript -1 denotes the matrix inverse and the superscript

T denotes the matrix transpose. To present the following derivation in a more intuitive way, we will define the scalars $s_i^1(\mathbf{x}) \equiv \mathbf{e}_i^T V_1^{-1} \mathbf{x}$ and $s_i^2(\mathbf{x}) \equiv \mathbf{e}_i^T V_2^{-1} \mathbf{x}$ which give the coordinates of \mathbf{x} in the basis of eigenvectors of J_1 and J_2 , respectively. Using (A4), one can show that

$$\begin{aligned} s_i^1(\mathbf{x}(T + t_1) - \mathbf{x}_1) &= \lambda_i s_i^1(\mathbf{x}(t_1) - \mathbf{x}_1) \\ s_i^2(\mathbf{x}(T + t_2) - \mathbf{x}_2) &= \lambda_i s_i^2(\mathbf{x}(t_2) - \mathbf{x}_2). \end{aligned} \tag{A8}$$

For convenience of notation in (A8), we have dropped the higher-order error terms that would carry through from (A4).

Consider any initial condition $\mathbf{x}(0) \in \Gamma_{\theta_1}$. Written in the basis of eigenvectors of J_1

$$\begin{aligned} \mathbf{x}(0) - \mathbf{x}_1 &= s_1^1(\mathbf{x}(0) - \mathbf{x}_1)v_1^1 \\ &\quad + \cdots + s_n^1(\mathbf{x}(0) - \mathbf{x}_1)v_n^1 \\ \mathbf{x}(T) - \mathbf{x}_1 &= \lambda_1 s_1^1(\mathbf{x}(0) - \mathbf{x}_1)v_1^1 \\ &\quad + \cdots + \lambda_n s_n^1(\mathbf{x}(0) - \mathbf{x}_1)v_n^1, \end{aligned} \tag{A9}$$

where we have used (A8) to obtain the second line. Using (A6), we can write to leading order $\|\mathbf{x}(0) - \mathbf{x}_1\|^2$

$$\begin{aligned} \mathbf{x}(\Delta t) - \mathbf{x}_2 &= H[s_1^1(\mathbf{x}(0) - \mathbf{x}_1)v_1^1 \\ &\quad + \cdots + s_n^1(\mathbf{x}(0) - \mathbf{x}_1)v_n^1] \\ \mathbf{x}(\Delta t + T) - \mathbf{x}_2 &= H[\lambda_1 s_1^1(\mathbf{x}(0) - \mathbf{x}_1)v_1^1 \\ &\quad + \cdots + \lambda_n s_n^1(\mathbf{x}(0) - \mathbf{x}_1)v_n^1]. \end{aligned} \tag{A10}$$

Here $\Delta t = \omega(\theta_2 - \theta_1)$ with $\dot{\theta} = \omega$ under the flow $\dot{\mathbf{x}} = F(\mathbf{x})$. When calculating isostables using (A7) $t_{\Gamma_2} = 0$ for any $\mathbf{x} \in \Gamma_{\theta_2}$. Because $\mathbf{x}(\Delta t) \in \Gamma_{\theta_2}$, $\psi_i^2[\mathbf{x}(\Delta t)] = s_i^2[\mathbf{x}(\Delta t) - \mathbf{x}_2]$. Also, $\mathbf{x}(\Delta t + T) \in \Gamma_{\theta_2}$, and using (A8), $\psi_i^2[\mathbf{x}(\Delta t + T)] = \lambda_i s_i^2[\mathbf{x}(\Delta t) - \mathbf{x}_2]$. With (A10), this implies

$$\begin{aligned} \lambda_i s_i^2(H\{s_1^1[\mathbf{x}(0) - \mathbf{x}_1]v_1^1 + \cdots + s_n^1[\mathbf{x}(0) - \mathbf{x}_1]v_n^1\}) \\ = s_i^2(H\{\lambda_1 s_1^1[\mathbf{x}(0) - \mathbf{x}_1]v_1^1 \\ + \cdots + \lambda_n s_n^1[\mathbf{x}(0) - \mathbf{x}_1]v_n^1\}). \end{aligned} \tag{A11}$$

Equation (A11) holds for any $\mathbf{x}(0)$. We will assume that λ_i has an algebraic multiplicity of 1 so, by linearity of s_i^1 and s_i^2 , (A11) implies

$$s_i^2(Hv_j^1) = 0 \text{ for all } j \neq i \tag{A12}$$

and

$$s_i^2[H(\mathbf{x} - \mathbf{x}_1)] = s_i^1(\mathbf{x} - \mathbf{x}_1)s_i^2(Hv_i^1). \tag{A13}$$

Using this information, we can now show that the isostable response curve is invariant to the value of θ we use for our initial Poincaré section.

Consider any initial condition $\mathbf{z}(0) \in \mathbb{R}^n$ which returns to Γ_{θ_1} at time t_1 . Let $s_i^1(\mathbf{z}(t_1) - \mathbf{x}_1) = \beta$, and then by (A7)

$$\psi_i^1(\mathbf{z}(0) - \mathbf{x}_1) = \beta \exp[-\log(\lambda_i)t_1/T]. \tag{A14}$$

Now consider a perturbed initial condition $\mathbf{y}(0) = \mathbf{z}(0) + \Delta \mathbf{x}$ which returns to Γ_{θ_1} at time $t_1 + p$ with $s_i^1(\mathbf{y}(t_1 + p) - \mathbf{x}_1) = \mu + \beta$. We can write

$$\psi_i^1(\mathbf{y}(0) - \mathbf{x}_1) = (\mu + \beta) \exp[-\log(\lambda_i)(t_1 + p)/T] \tag{A15}$$

so

$$\begin{aligned} \psi_i^1(\mathbf{y}(0) - \mathbf{x}_1) - \psi_i^1(\mathbf{z}(0) - \mathbf{x}_1) \\ = (\mu + \beta) \exp[-\log(\lambda_i)(t_1 + p)/T] \\ - \beta \exp[-\log(\lambda_i)t_1/T] \\ \equiv \Lambda. \end{aligned} \tag{A16}$$

Now consider the isostables with respect to the Γ_{θ_2} Poincaré section. Each trajectory will reach Γ_{θ_2} at a time Δt after it reaches Γ_{θ_1} . Using (A6),

$$\begin{aligned} \mathbf{z}(t_1 + \Delta t) &= H(\mathbf{z}(t_1) - \mathbf{x}_1) + \mathbf{x}_2 \\ &\quad + \mathcal{O}(\|\mathbf{z}(t_1) - \mathbf{x}_1\|^2) \\ \mathbf{y}(t_1 + p + \Delta t) &= H(\mathbf{y}(t_1 + p) - \mathbf{x}_1) + \mathbf{x}_2 \\ &\quad + \mathcal{O}(\|\mathbf{y}(t_1 + p) - \mathbf{x}_1\|^2). \end{aligned} \tag{A17}$$

Using (A13), we find

$$\begin{aligned} \psi_i^2[\mathbf{z}(0) - \mathbf{x}_2] &= \beta s_i^2(Hv_i^1) \\ &\quad \times \exp[-\log(\lambda_i)(t_1 + \Delta t)/T] \\ \psi_i^2(\mathbf{y}(0) - \mathbf{x}_2) &= (\mu + \beta)s_i^2(Hv_i^1) \\ &\quad \times \exp[-\log(\lambda_i)(t_1 + p + \Delta t)/T]. \end{aligned} \tag{A18}$$

Algebraic manipulation yields

$$\begin{aligned} \psi_i^2[\mathbf{y}(0) - \mathbf{x}_2] - \psi_i^2[\mathbf{z}(0) - \mathbf{x}_2] \\ = s_i^2(Hv_i^1) \exp[-\log(\lambda_i)\Delta t/T] \Lambda. \end{aligned} \tag{A19}$$

Note here that $s_i^2(Hv_i^1)$ and $\exp[-\log(\lambda_i)\Delta t/T]$ are both constant terms so Eqs. (A16) and (A19) imply that regardless of which isochron is chosen for the Poincaré section to define isostables, the shape of the resulting isostable response curves will be identical provided λ_i has an algebraic multiplicity of 1. The magnitude of the isostable response curve can vary depending on the geometry of the periodic orbit.

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