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I. INTRODUCTION

The color of the background represents the order parameter r of the system. The color of the snapshots represents the phase of the oscillators. The color of the arrows represents the direction of the velocity of the oscillators. The size of the arrows represents the speed of the oscillators.

II. MODEL

Oscillators have a spatial position $\mathbf{r}_i = (x_i, y_i)$ and an internal phase θ_i which evolve according to equations:

$$\dot{x}_i = v \cos \theta_i, \quad (1)$$

$$\dot{y}_i = v \sin \theta_i, \quad (2)$$

$$\dot{\theta}_i = \omega_i + \lambda \sum_{j=1}^N A_{ij} \sin(\theta_j - \theta_i) \quad (3)$$

for $i = 1, 2, \dots, N$, where N is the number of oscillators. As per Eq. (1) and (2), each oscillator moves with a constant speed v in the direction of its current phase θ_i . The phase θ_i evolves according to Eq. (3), where ω_i is the natural frequency of the i th oscillator, λ is the coupling strength, and A is the adjacency matrix of the network, with $A_{ij} = 1$ if there is a connection from i th to j th oscillator, and $A_{ij} = 0$ otherwise. We can consider Eq. (1)-(3) as a generalization of the Kuramoto model and the Vicsek model in the sense that it includes both the phase and the spatial position of the oscillators.

Each oscillator i is connected to all the oscillators within a action radius d_0 of its position. The adjacency matrix A is defined as:

$$A_{ij} = \begin{cases} 1, & |\mathbf{r}_i - \mathbf{r}_j| \leq d_0 \\ 0, & |\mathbf{r}_i - \mathbf{r}_j| > d_0 \end{cases} \quad (4)$$

where $|\mathbf{r}_i - \mathbf{r}_j|$ is the Euclidean distance between the i th and j th oscillators.

For simplicity, we consider oscillators are initially distributed uniformly in a two-dimensional square with side length L and periodic boundary conditions. Their positions $\mathbf{r}_i(t) = (x_i(t), y_i(t))$ at given time t are given by:

$$\begin{aligned} x_i(t + \Delta t) &= x_i(t) + v \cos \theta_i(t) \Delta t \bmod L, \\ y_i(t + \Delta t) &= y_i(t) + v \sin \theta_i(t) \Delta t \bmod L, \end{aligned} \quad (5)$$

where Δt is the discrete time step. When two oscillators are on opposite sides of the square, the absolute value of the difference between one of their coordinates is greater than $L/2$. In this case, we take the minimum distance between them, which is the distance between the two points in the periodic boundary conditions. For a given pair of points \mathbf{r}_i and \mathbf{r}_j , the distance between them is $|\mathbf{r}_i - \bar{\mathbf{r}}_j|$, where $\bar{\mathbf{r}}_j = (\bar{x}_j, \bar{y}_j)$ is the adjusted position of the j th oscillator, given by:

$$\bar{x}_j = \begin{cases} x_j, & |x_i - x_j| \leq L/2 \\ x_j + L, & x_i - x_j > L/2 \\ x_j - L, & x_i - x_j < -L/2 \end{cases}, \quad (6)$$

$$\bar{y}_j = \begin{cases} y_j, & |y_i - y_j| \leq L/2 \\ y_j + L, & y_i - y_j > L/2 \\ y_j - L, & y_i - y_j < -L/2 \end{cases}. \quad (7)$$

$|\mathbf{r}_i - \bar{\mathbf{r}}_j|$ can be proved to be the minimum distance between \mathbf{r}_i and \mathbf{r}_j in the periodic boundary conditions (see the proof in Appendix A).

Finally, we consider that the natural frequencies ω_i are distributed in two symmetric uniform distributions. Exactly half of the oscillators have natural frequencies in the range $[\omega_{\min}, \omega_{\max}]$ ($\omega_i \sim U(\omega_{\min}, \omega_{\max}), i = 1, 2, \dots, N/2$) and the other half in the range $[-\omega_{\max}, -\omega_{\min}]$ ($\omega_i \sim U(-\omega_{\max}, -\omega_{\min}), i = N/2 + 1, N/2 + 2, \dots, N$).

III. BEHAVIOR

We performed numerical simulations of the model to probe the behavior of its solutions (see Appendix B for details on the

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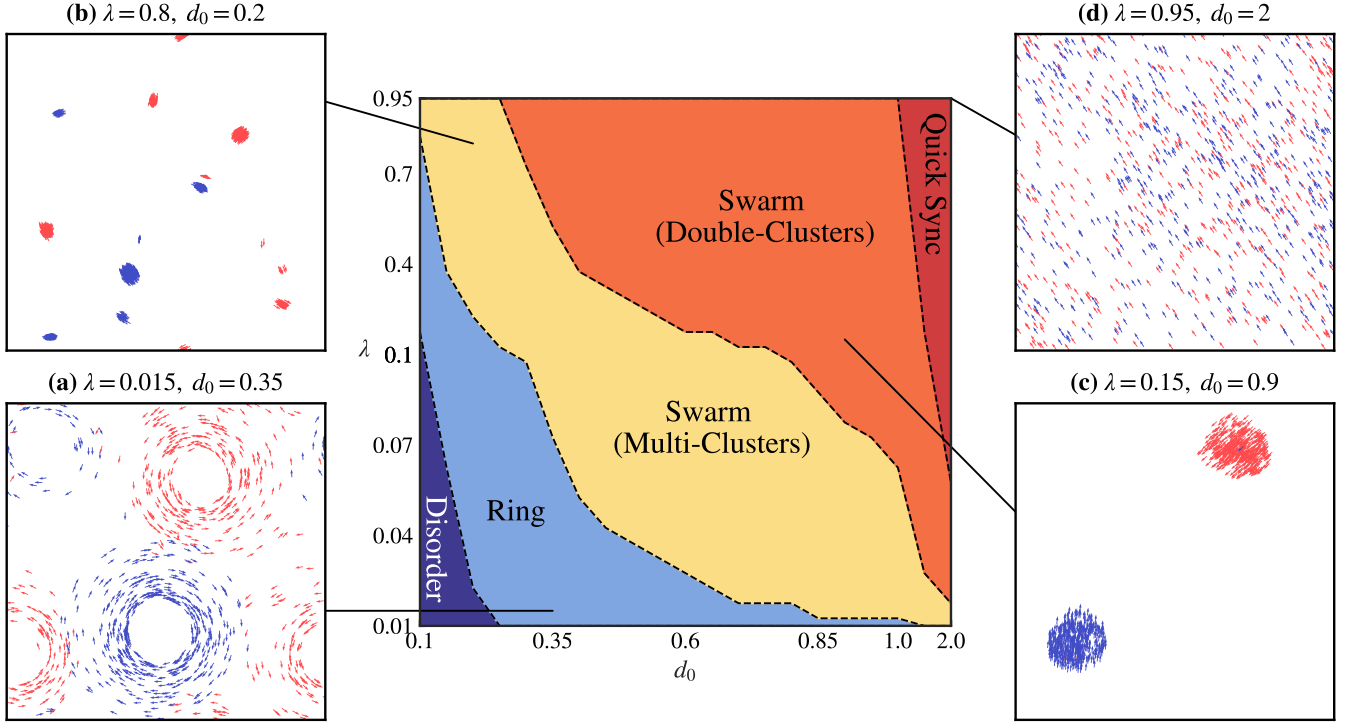


FIG. 1. Phase diagram of model Eq. (1)-(3) in the $(\lambda-d_0)$ plane. The boundaries between states is analytical approximations given by Section V. For the sake of clarity, the scale of λ and d_0 is non-linear (For λ in $[0.01, 0.1]$ and $[0.1, 1]$, step length is 0.1 and 0.05, respectively. For d_0 in $[0.1, 1]$ and $[1, 2]$, step length is 0.05 and 0.5, respectively). (a), The snapshots of Ring ($\lambda = 0.015$, $d_0 = 0.35$). (b), Swarm (Multi-Clusters) ($\lambda = 0.8$, $d_0 = 0.2$). (c), Swarm (Double-Clusters) ($\lambda = 0.15$, $d_0 = 0.9$). (d), Quick Sync ($\lambda = 0.95$, $d_0 = 2$). Two types of chiral oscillators are represented by red ($\omega_i > 0$) and blue ($\omega_i < 0$) arrows, respectively.

numerical methods). $N = 1000$ oscillators were distributed uniformly in the square of length $L = 10$ and their natural frequencies were distributed in the range $[\omega_{\min}, \omega_{\max}] = [1, 3]$ and $[-\omega_{\max}, -\omega_{\min}] = [-3, -1]$. Two-parameter of coupling strength λ and action radius d_0 are presented in the phase diagram in Fig. 1. We found the system settles into five states: **Disorder**, **Ring**, **Swarm** (which can be further divided into **Multi-Clusters** and **Double-Clusters**), and **Quick Sync**. In Fig. 1 we show the snapshots of the last three states and where

these states are located in the phase diagram. The Disorder state is shown in Fig. 2a. We next discuss these five states.

A. Disorder State

Disorder state occurs when both λ and d_0 are small. In this state, the oscillators are not asynchronous (phase histogram is uniform, like Fig. 3a) and move in a way which similar to uncoupled oscillators ($\lambda = 0$), as shown in Fig. 2a. According to Eq. (1)-(3), when $\lambda = 0$, the equations of oscillators' motion can be written as:

$$\begin{aligned} x_i(t) &= x_i(0) + \frac{v}{\omega_i} \sin \omega_i t, \\ y_i(t) &= y_i(0) + \frac{v}{\omega_i} \cos \omega_i t. \end{aligned} \quad (8)$$

In such a setup, oscillators move in a circular trajectory with radius v/ω_i and the phases θ_i increase linearly with time, as show in Fig. 2b. To calculate the real-time rotational radius, we first estimate real-time centers $\mathbf{c}(t)$ of the circular trajectory with method in Fig. 4 and then solve the following linear equations:

$$\begin{aligned} \mathbf{c}_i(t_1) \cdot \mathbf{v}_i(t_1) &= \mathbf{x}_i(t_1) \cdot \mathbf{v}_i(t_1), \\ \mathbf{c}_i(t_2) \cdot \mathbf{v}_i(t_2) &= \mathbf{x}_i(t_2) \cdot \mathbf{v}_i(t_2). \end{aligned} \quad (9)$$

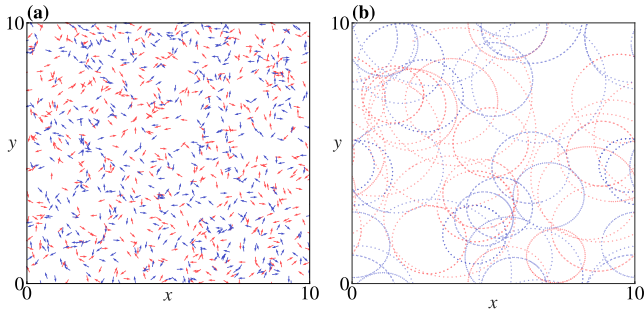


FIG. 2. Key properties of the Disorder state. (a), The snapshot of the Disorder state ($\lambda = 0.01$, $d_0 = 0.1$, $T = 60000$). (b), The scatter plot of last 100 time steps of 20 positive chirality oscillators and 20 negative chirality oscillators.

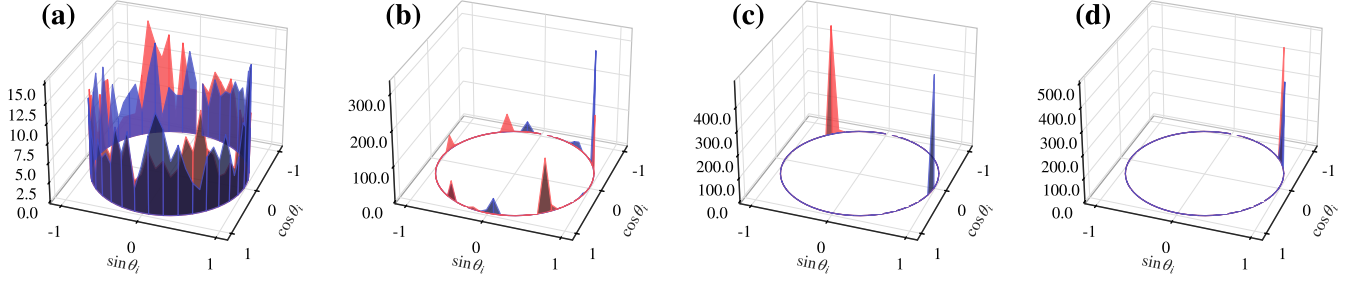


FIG. 3. Histogram of the oscillators' phases. (a), Ring state ($\lambda = 0.015$, $d_0 = 0.35$). (b), Swarm state (Multi-Clusters, $\lambda = 0.8$, $d_0 = 0.2$). (c), Swarm state (Double-Clusters, $\lambda = 0.15$, $d_0 = 0.9$). (d), Quick Sync state ($\lambda = 0.95$, $d_0 = 2$). The histograms are calculated with 70 bins.

The real-time rotational radius is $r_i(t) = |\mathbf{c}_i(t) - \mathbf{r}_i(t)|$. We found that the real-time rotational radius is almost constant and close to v/ω_i for each oscillator in the Disorder state, as shown in Fig. 5a.

B. Ring State

The Ring state is characterized by the oscillators forming several rings with thickness, each of which is composed of oscillators with the same chirality, as is shown in Fig. 1a. Similar to Disorder state, the oscillators in the same ring cluster move in a circular trajectory with a constant rotational radius calculated in Fig. 5a. The oscillators' phase is uniformly distributed in the range $[-\pi, \pi]$ (cf. Fig. 3a), which leads to oscillators uniformly located on the circular trajectory. Fig. 6a shows there is a long transient time before this state is reached, and in this transient time, the trajectories of oscillators with the same chirality aggregate slowly. Conversely, the oscillators with different chirality repel each other. For parameter plane in Fig. 1, the number of ring decreases and local swarms appear at high-frequency ($|\omega_i| \rightarrow \omega_{\max}$) oscillators with λ and d_0 increasing, as is shown in Fig. 7a.

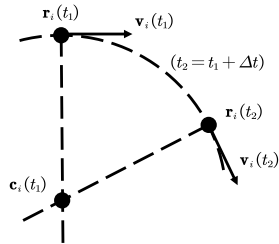


FIG. 4. Estimation for real-time centers. $\mathbf{v}_i(t_1)$ and $\mathbf{r}_i(t_1)$ are the velocity and position of i th oscillator at t_1 , respectively. According to Eq. (1)-(3), we can calculate $\mathbf{v}_i(t_2)$ and $\mathbf{r}_i(t_2)$, ($t_2 = t_1 + \Delta t$). The line $\mathbf{c}_i(t)\mathbf{v}_i(t)$ is perpendicular to $\mathbf{v}_i(t)$.

C. Swarm State

Swarm State is a state where the oscillators form spatial clusters and align into several clusters [Fig. 1b, c and Fig. 6b]. When λ and d_0 increases, the number of clusters decreases by 2, which is named by as Double-Clusters state, and other states are named by Multi-Clusters state. The clusters are composed of oscillators with the same chirality, and the phase θ_i of the oscillators in the same cluster is synchronized as seen in Fig. 3b and 3c, which means that the oscillators in the cluster move with the same velocity $\mathbf{v}_i = (v/\theta_s, \sin \theta_s)$ and rotational radius $r_i = v/\theta_s$, where θ_s is the oscillators' phase in the cluster. Based on this property, we can calculate θ_s and r_i with Eq. (3):

$$\begin{aligned} N_s \omega_s &= \sum_{i=1}^{N_s} \left(\omega_i + \lambda \sum_{j=1}^{N_s} A_{ij} \sin(\theta_j - \theta_i) \right) \\ \omega_s &= \frac{1}{N_s} \sum_{i=1}^{N_s} \omega_i + \frac{\lambda}{N_s} \sum_{i=1}^{N_s} \sum_{j=1}^{N_s} A_{ij} \sin(\theta_j - \theta_i) \\ &= \frac{1}{N_s} \sum_{i=1}^{N_s} \omega_i, \end{aligned} \quad (10)$$

where N_s is the number of oscillators in the cluster. As $\omega_i \sim U(\omega_{\min}, \omega_{\max})$ and $\omega_i \sim U(-\omega_{\max}, -\omega_{\min})$ for two types of chirality, we can calculate θ_i , ω_s and r_i with ω_i for Double-Clusters state:

$$\begin{aligned} \theta_i = \omega_s &= \begin{cases} (\omega_{\max} + \omega_{\min})/2, & i = 1, 2, \dots, N/2 \\ -(\omega_{\max} + \omega_{\min})/2, & i = N/2 + 1, \dots, N \end{cases}, \\ r_i &= \frac{v}{\omega_s}, \end{aligned} \quad (11)$$

as shown in Fig. 5b. But for Multi-Clusters, due to which oscillators are synchronized within each cluster is not accurately known, we can only calculate the real-time rotational radius of the them. As seen in Fig. 5b, similar to Double-Clusters, some local platforms appear in the real-time rotational radius due to synchronization.

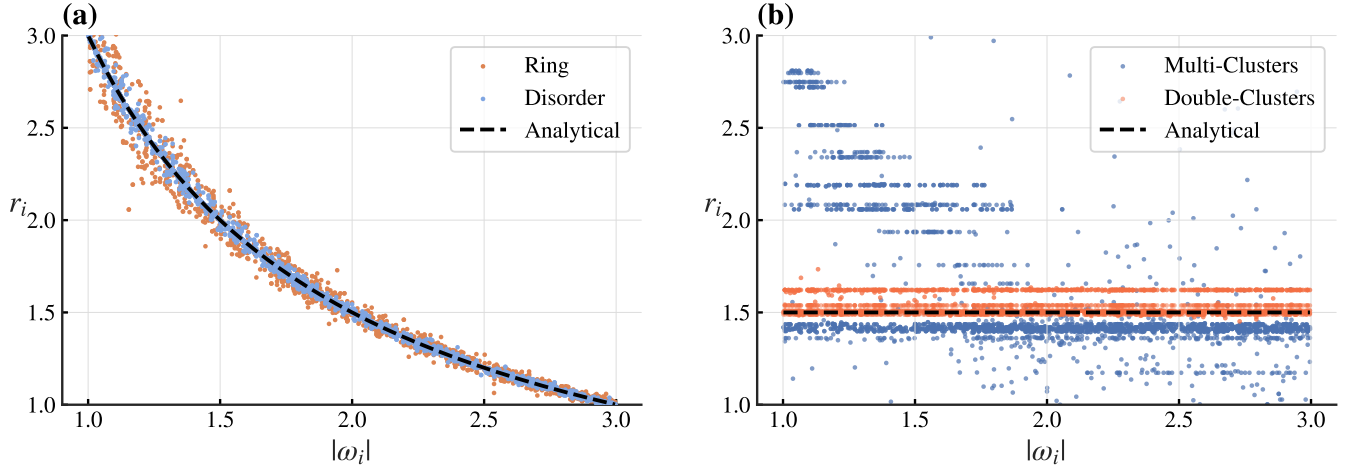


FIG. 5. The real-time and analytical rotational radius. **(a)**, Radius for the Disorder ($d_0 = 0.1$, $\lambda = 0.01 : 0.06$) and Ring ($d_0 = 0.1$, $\lambda = 0.06 : 0.1$). The real-time rotational radius is almost constant and close to v/ω_i for each oscillator. **(b)**, Radius for Swarm (Multi-Clusters, $d_0 = 0.15 : 0.25$, $\lambda = 0.95$) and (Double-Clusters, $d_0 = 2$, $\lambda = 0.02 : 0.05$). Line of Analytical is only for Double-Clusters. All the above simulations are calculated at $t = 60000$.

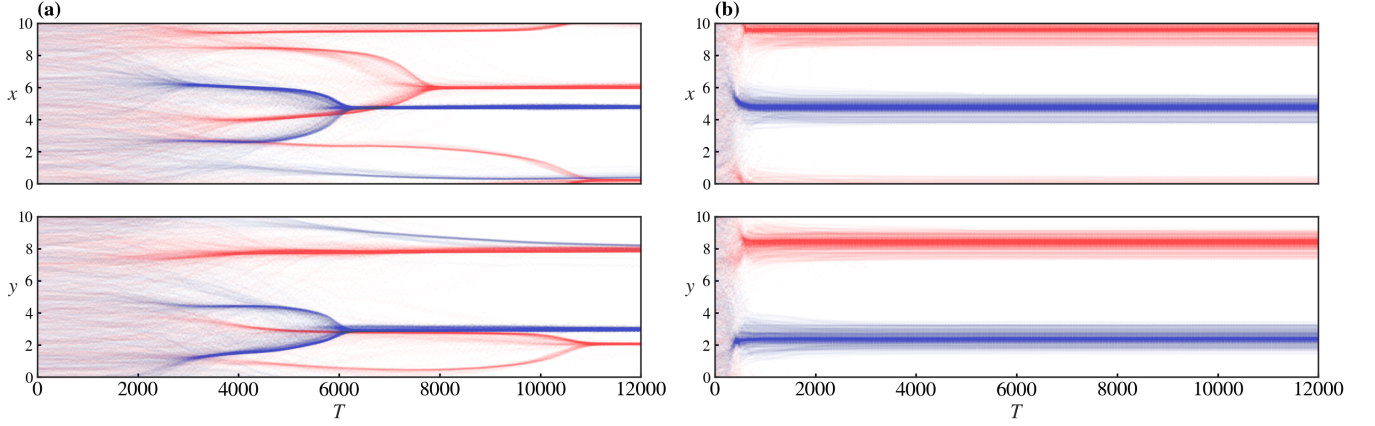


FIG. 6. Scatter plot of the real-time centers position. **(a)**, centers position of Ring ($\lambda = 0.02$, $d_0 = 0.4$). As time goes on, the centers of oscillators with the same chirality converge. **(b)**, centers position of Swarm ($\lambda = 0.01$, $d_0 = 2$). Unlike Ring, the centers converge quickly. The centers position are estimated with method in Fig. 4 and Eq. (9).

D. Quick Sync State

Quick Sync state is a simple state where total oscillators are synchronized quickly, as shown in Fig. 1d. and 3d. The oscillators are synchronized in an extremely short time, which leads them have no time to form clusters (can also be considered as a special case of Swarm state). Due to the two types of chirality oscillators are synchronized and the distributions of them is symmetric, the phase velocities of total oscillators are close to zero according to Eq. (10).

IV. ORDER PARAMETER

Having described the four states of our system, we next discuss how to distinguish them. We use the order parameter R

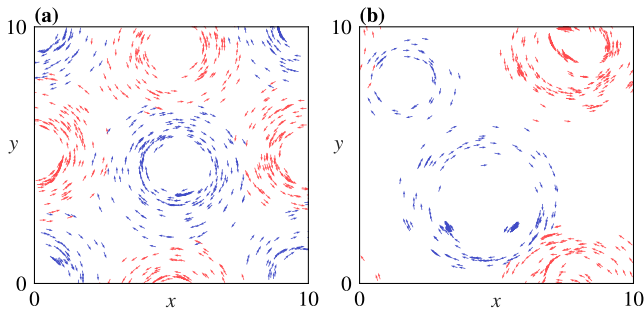


FIG. 7. Snapshots of the Ring state. **(a)**, Four rings ($\lambda = 0.02$, $d_0 = 0.4$). **(b)**, Three rings ($\lambda = 0.025$, $d_0 = 0.45$).

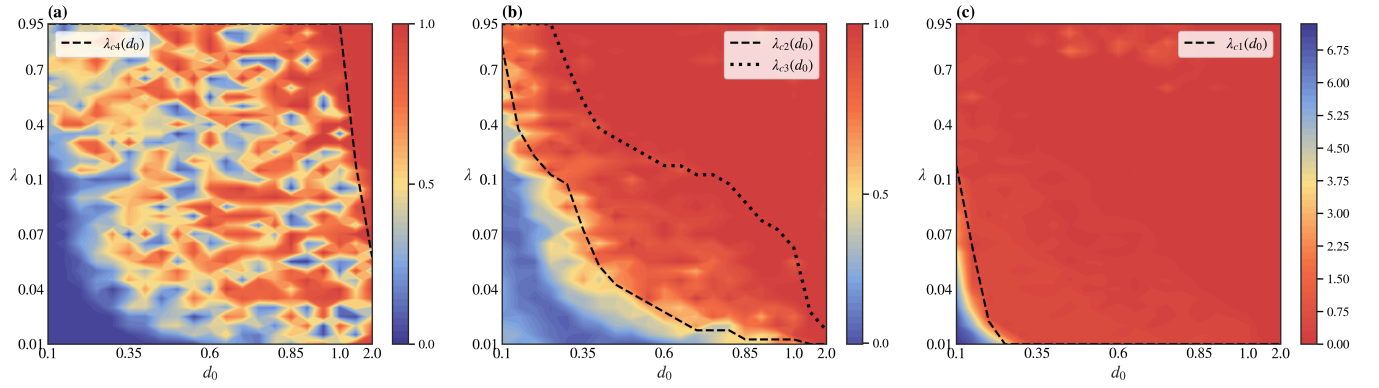


FIG. 8. Order parameter heatmaps of (λ, d_0) plane. (a), order parameter R and critical line of λ_{c4} . (b), order parameter R_s and critical lines of $\lambda_{c2}, \lambda_{c3}$. All order parameters are calculated at $t = 60000$

to metric global synchronization. The order parameter R is defined as:

$$R = \left| \frac{1}{N} \sum_{j=1}^N e^{i\theta_j} \right|. \quad (12)$$

The order parameter R is absolute value of the mean of the complex numbers $e^{i\theta_i}$, which can be interpreted as the mean direction of the oscillators. When $R = 1$, the oscillators are completely synchronized, and when $R = 0$, the oscillators are completely desynchronized. Fig 8a shows the order parameter R in the parameter plane. The order parameter R is close to 1 in the Quick Sync state, close to 0 in the Disorder state and most of the Ring state, and between 0 and 1 in other states. In these states, We see the order parameter R changes non-monotonically in the sense that phases in these states are not globally synchronized and each cluster's phase velocity $\omega_s \neq 0$. When the phases of different clusters are exactly equal, the order parameter R is close to 1, and when they are exactly opposite, R is close to 0.

Having realized that the order parameter R is not enough to distinguish the states with clusters, we next define the following order parameter R_s to metric the local synchronization,

$$R_s = \frac{1}{N_s} \sum_{k=1}^{N_s} \left| \frac{1}{|C_k|} \sum_{j \in C_k} e^{i\theta_j} \right|, \quad (13)$$

where N_s is the number of clusters, C_k is the k th cluster, and $|C_k|$ is the number of oscillators in the k th cluster. We can consider R_s as an order parameter that studies the spatial position and internal phase simultaneously. To estimate the number of clusters, we use the following method: we first calculate the relative center distance matrix $D_{ij} = |\mathbf{c}_i - \bar{\mathbf{c}}_j|$, where $\bar{\mathbf{c}}_j = (\bar{x}_j, \bar{y}_j)$ is the adjusted position of the j th oscillator's rotational center calculated by Eq. (6), (7) and (9). Then we use the DBSCAN algorithm to cluster the oscillators. The DBSCAN algorithm is a density-based clustering algorithm, which can find clusters of arbitrary shapes and sizes. We set the minimum number of oscillators in a cluster to be 10 and

the maximum distance between two oscillators in the same cluster to be 0.3 (see Appendix C for details on the determination of these parameters). We then calculate the order parameter R_s for each cluster. The order parameter R_s is close to 1 in the Swarm state and Quick Sync state, and close to 0 in Disorder state and most of the Ring state, between 0 and 1 in other Ring states with local clusters, as shown in Fig. 8b.

Combining the order parameter R and R_s , we can find only the distinction between Ring and Disorder states has not been resolved.

V. ANALYTICAL APPROXIMATIONS

VI. CONCLUSIONS

Appendix A: PROOF OF THE ADJUSTED POSITION

Appendix B: NUMERICAL METHODS

All the simulations of the model Eq. (1)-(3) were run on Python using Euler integration, with a time step $\Delta t = 0.01$, and a total time of $T = 60000$.

Appendix C: Determination of DBSCAN's parameters

DBSCAN (Density-Based Spatial Clustering of Applications with Noise) is a density-based clustering algorithm, which can find clusters of arbitrary shapes and sizes. The algorithm has two parameters: ϵ and m . ϵ is the maximum distance between two samples for one to be considered as in the neighborhood of the other, and m is the minimum number of samples in a neighborhood for a point to be considered as a core point.

We set the minimum number of oscillators in a cluster to be 10 and the maximum distance between two oscillators in the same cluster to be 0.3. We determined these parameters as follows. We traverse all values between 0.1 and 0.5 with

a step length of 0.1, and for each value of ε , we calculate the number of clusters with $m = 0$.

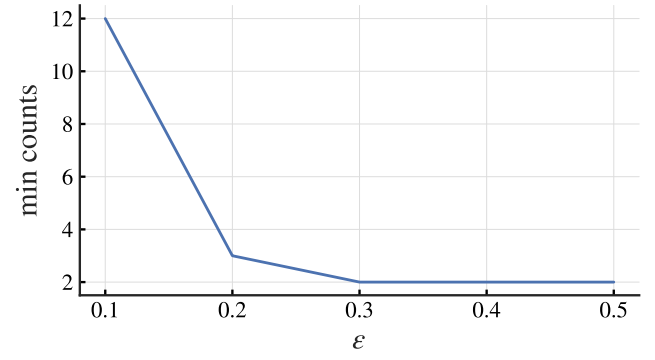


FIG. 9. The number of clusters with different ε , $m = 0$. The number of clusters is calculated with the DBSCAN algorithm.

As shown in Fig. 9, the minimum counts of clusters 2 (corresponding to Double-Clusters state of Swarm) first appear at $\varepsilon = 0.3$. Then we set ε to be 0.3.

Then we set m to be 10 which is 1% of the population ($N = 1000$) of the system.