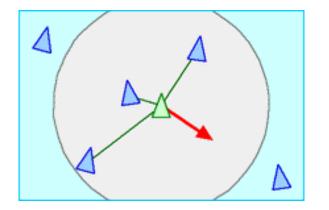
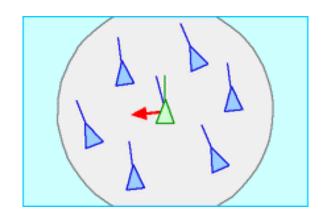


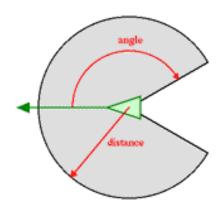
Boids Model

Simulated Flocks

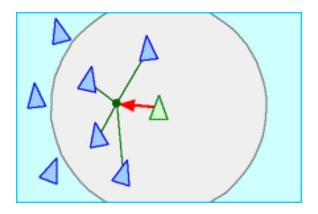
- Collision Avoidance: avoid collisions with nearby flockmates
- Velocity Matching: attempt to match velocity with nearby flockmates
- Flock Centering: attempt to stay close to nearby flockmates





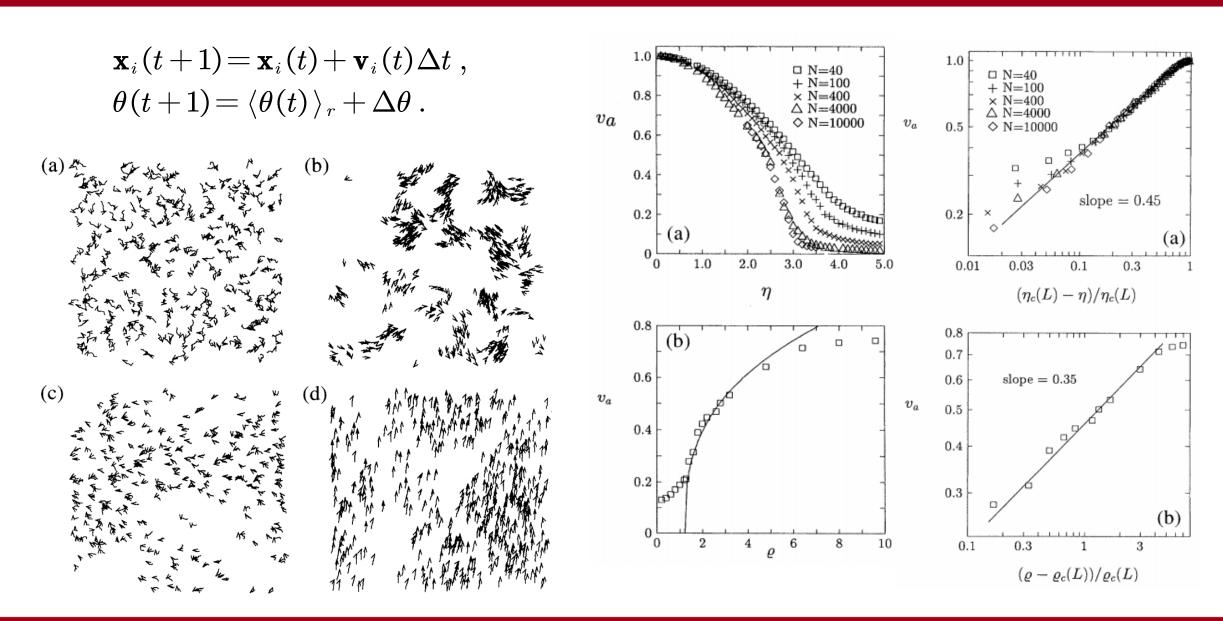


a boid's neighborhood



J. Toner and Y. Tu, Flocks, Herds, and Schools: A Quantitative Theory of Flocking, Phys. Rev. E 58, 4828 (1998).

Vicsek Model



T. Vicsek, A. Czirók, E. Ben-Jacob, I. Cohen, and O. Shochet, Phys. Rev. Lett. 75, 1226 (1995).

Self-Propelled Particles

 $\dot{r}_i \!=\! c\, \mathbf{e}_{ heta_i}$

$$\begin{split} \dot{\theta}_{i} &= \gamma \sum_{j=1}^{N} F(\theta_{j} - \theta_{i}, r_{j} - r_{i}) + \sqrt{2\epsilon} \, \tilde{\boldsymbol{\eta}}_{i}(t), \\ & = \begin{cases} \frac{\sin(\theta)}{\pi R^{2}}, & |r| < R \\ 0, & \text{otherwise} \end{cases} \end{split}$$

$$f(\mathbf{r}, \theta) = \sum_{j=-1}^{N} \delta(\mathbf{r} - r_{j}) \delta(\theta)$$

$$\partial_{t} f(\mathbf{r}, \theta) + \mathbf{e}_{\theta} \cdot \nabla[vf]$$

$$= \epsilon \frac{\partial^{2} f}{\partial \theta^{2}} - \frac{\partial}{\partial \theta} \sqrt{2\epsilon f} \, \boldsymbol{\eta} - \gamma \frac{\partial}{\partial \theta} \int d\theta' d \, \mathbf{r}' f(\mathbf{r}', \theta') \\ \times f(\mathbf{r}, \theta) F(\theta' - \theta, \mathbf{r} - \mathbf{r}'). \end{split}$$

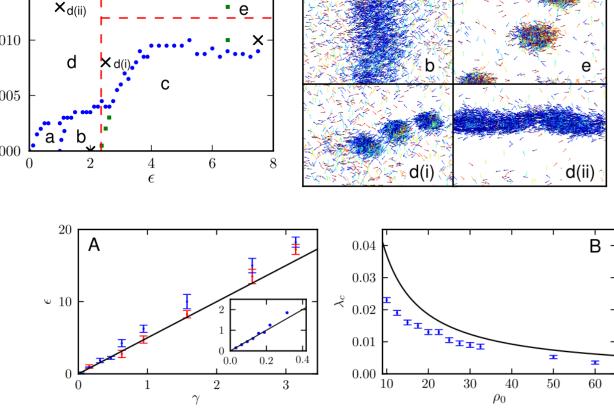


FIG. 2 (color online). (a) Phase boundary for the flying XY model when $\lambda = 0$, showing the critical value of ϵ as a function of γ . Blue (darker gray) points for $\nu = 2.0$, red (lighter gray) for $\nu = 0.5$. Inset: data for $\nu = 2.0$ for smaller values of γ . (b) Phase boundary for $\epsilon = 5$, $\gamma = 0.16$. In all cases L = 10 and N = 1000.

(j)

Uniform disordered phase

3

0

 $-\frac{g \operatorname{Pe}_{\mathbf{r}} \Omega}{\operatorname{QL}} [3 \nabla_{\perp} \mathbf{w}^2 - 6 \mathbf{w} (\nabla_{\perp} \cdot \mathbf{w}) - 10 (\mathbf{w} \cdot \nabla_{\perp}) \mathbf{w}]. \quad (3)$

More CAPs Models

[1] D. Levis, I. Pagonabarraga, and B. Liebchen, *Activity Induced Synchronization: Mutual Flocking and Chiral Self-Sorting*, Phys. Rev. Research 1, 023026 (2019).

$$\dot{\boldsymbol{r}}_{\alpha} = v\boldsymbol{n}_{\alpha},\tag{1}$$

$$\dot{\theta}_{\alpha} = \omega_{\alpha} + \frac{K}{\pi R^2} \sum_{\nu \in \partial_{\alpha}} \sin(\theta_{\nu} - \theta_{\alpha}) + \sqrt{2D_r} \eta_{\alpha}. \tag{2}$$

[3] B. Ventejou, H. Chaté, R. Montagne, and X. Shi, *Susceptibility of Orientationally Ordered Active Matter to Chirality Disorder*, Phys. Rev. Lett. **127**, 238001 (2021).

$$\dot{\mathbf{r}}_i = v_0 \, \mathbf{e}(\theta_i), \tag{1a}$$

$$\dot{\theta}_i = \omega_i + \kappa \langle \sin \alpha(\theta_j - \theta_i) \rangle_{j \sim i} + \sqrt{2D_r} \eta_i,$$
 (1b)

[2] D. Levis and B. Liebchen, *Simultaneous Phase Separation and Pattern Formation in Chiral Active Mixtures*, Phys. Rev. E **100**, 012406 (2019).

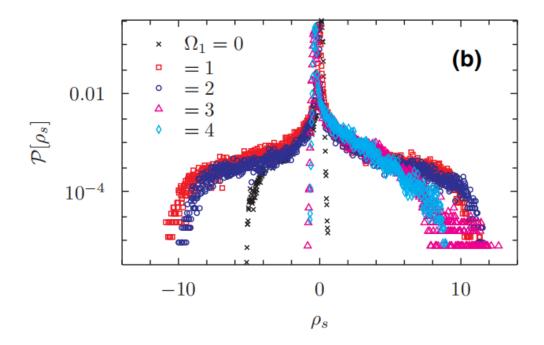
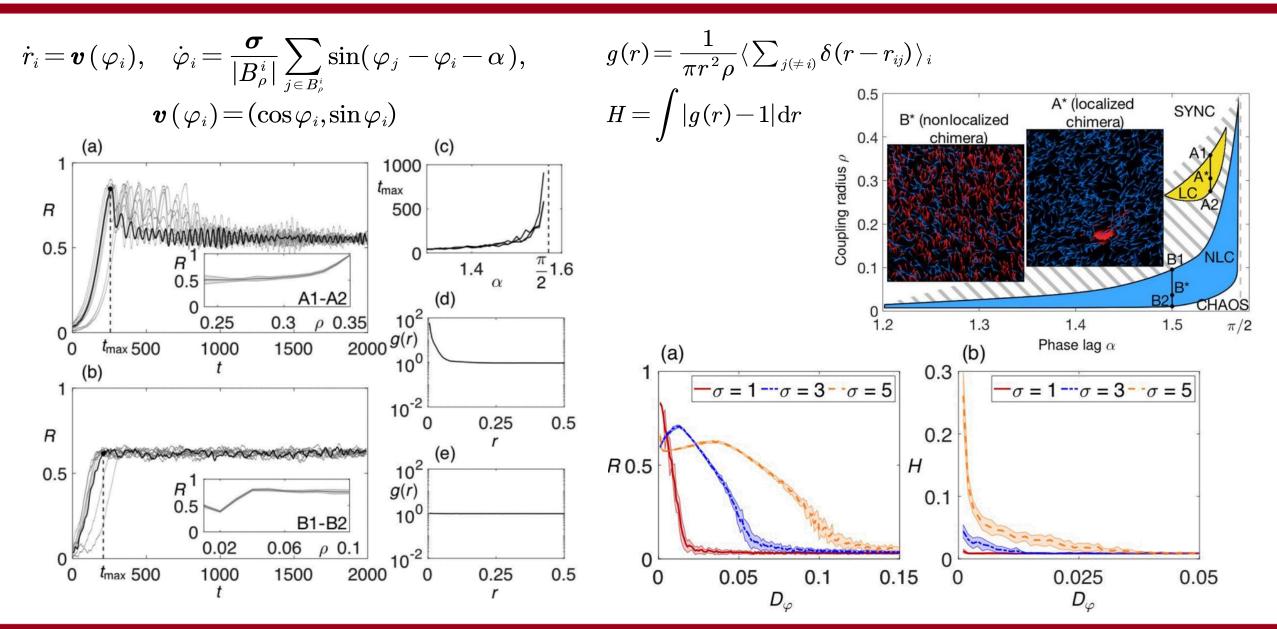


FIG. 3. Distribution of the local segregation factor $\mathcal{P}(\rho_s)$ with $\rho_s = (\rho_1 - \rho_2)/\rho_0$ for several values of Ω_1 (see key) for monochiral mixtures (fixed $\Omega_2 = 1$) (a) and for bichiral mixtures ($\Omega_2 = -1$) (b). The two dotted lines in the top panel correspond to two Gaussian distributions centered in zero with different variance.

Self-propelled chimeras



N. Kruk, Y. Maistrenko, and H. Koeppl, Self-Propelled Chimeras, Phys. Rev. E 98, 032219 (2018).

General Chemotactic Model of Oscillators

$$\dot{X}_i(t) = f(X_i) + kg(S(r_i, t)), \tag{1}$$

$$m\ddot{\mathbf{r}}_i(t) = -\gamma \dot{\mathbf{r}}_i - \sigma(\mathbf{X}_i) \nabla S|_{\mathbf{r}=\mathbf{r}_i}, \tag{2}$$

$$\tau \partial_t S(\mathbf{r}, t) = -S + d\nabla^2 S + \sum_i h(\mathbf{X}_i) \delta(\mathbf{r} - \mathbf{r}_i).$$
 (3)

$$\dot{\psi}_{i}(t) = \sum_{j \neq i} e^{-|\mathbf{R}_{ji}|} \sin(\Psi_{ji} + \alpha |\mathbf{R}_{ji}| - c_1), \tag{15}$$

$$\dot{\mathbf{r}}_{i}(t) = c_{3} \sum_{j \neq i} \hat{\mathbf{R}}_{ji} e^{-|\mathbf{R}_{ji}|} \sin(\Psi_{ji} + \alpha |\mathbf{R}_{ji}| - c_{2}), \quad (16)$$

where $\mathbf{R}_{ji} \equiv \mathbf{r}_j - \mathbf{r}_i$, $\hat{\mathbf{R}}_{ji} \equiv \mathbf{R}_{ji}/|\mathbf{R}_{ji}|$, and $\Psi_{ji} \equiv \psi_j - \psi_i$. These equations contain the four real parameters $c_1 \equiv \arg(c\kappa b/\rho) - \pi/2$, $c_2 \equiv \arg(-b) - \pi/2$, $c_3 \equiv \mathrm{Re}\rho|\rho/c\kappa|(>0)$, and $\alpha \equiv \mathrm{Im}\rho/\mathrm{Re}\rho(>0)$. Note that c_3 is equal to the ratio of the time scales of ψ_i and \mathbf{r}_i . Also, note that although the density of the elements does not appear explicitly in these equations, it is an important parameter.

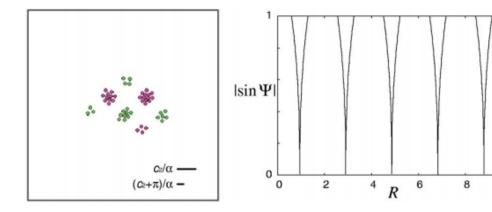


FIG. 1 (color). (left) Clustered clusters (or modular networks): snapshot of the element distribution in two-dimensional space exhibited by Eqs. (15) and (16), with $c_1 = c_2 = c_3 = 1.5$ and $\alpha = 1.6$. The colors represent the internal states ψ of the elements. The spatial size of the system is 30×30 , and it is shown in its entirety. The two scale bars represent c_2/α and $(c_2 + \pi)/\alpha$ (see main text). Although we have adopted a point element in this Letter, we plot its position with a finite size to facilitate visualization. (right) Invariant curve Eq. (17) with E = 1.3. The parameter values are the same as above. When the oscillators synchronize, i.e. $\sin \Psi = 0$, the distance R must be approximately 1, 3, 5, Because $|\sin \Psi| \le 1$, values of R in the neighborhood of 0, 2, 4, ... are avoided. In this way, an effective excluded volume (or zone) appears even though the elements are point objects.

Chemotactic Active Matter

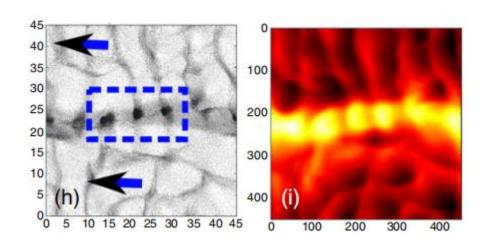
[1] B. Liebchen, D. Marenduzzo, I. Pagonabarraga, and M. E. Cates, *Clustering and Pattern Formation in Chemorepulsive Active Colloids*, Phys. Rev. Lett. **115**, 258301 (2015).

[2] B. Liebchen, M. E. Cates, and D. Marenduzzo, *Pattern Formation in Chemically Interacting Active Rotors with Self-Propulsion*, Soft Matter **12**, 7259 (2016).

$$\dot{\rho} = -\nu_0 \nabla \cdot (\rho \mathbf{p}) + D_\rho \nabla^2 \rho + K \nabla^2 \rho^3 \tag{1}$$

$$\dot{\phi} = \omega + \beta \hat{\mathbf{p}} \times \nabla c; \quad \mathbf{p} = (\cos \phi, \sin \phi)^T$$
 (2)

$$\dot{c} = k_0 \rho - k_d c + D_c \nabla^2 c + \varepsilon (c_0 - c)^3. \tag{3}$$



[3] B. Liebchen, D. Marenduzzo, and M. E. Cates, *Phoretic Interactions Generically Induce Dynamic Clusters and Wave Patterns in Active Colloids*, Phys. Rev. Lett. **118**, 268001 (2017).

$$\dot{\mathbf{r}}_i(t) = v\mathbf{p}_i,\tag{1}$$

$$\dot{\theta}_i(t) = \beta \mathbf{p}_i \times \nabla c + \sqrt{2D_r} \eta_i(t). \tag{2}$$

$$\dot{c}(\mathbf{r},t) = D_c \nabla^2 c - k_d c + \sum_{i=1}^N \oint d\mathbf{x}_i \delta(\mathbf{r} - \mathbf{r}_i(t) - R_0 \mathbf{x}_i) \sigma(\mathbf{x}_i).$$

$$\dot{\rho} = -\text{Pe}\nabla \cdot \mathbf{w},$$

$$\dot{\mathbf{w}} = -\mathbf{w} + \frac{B\rho}{2}\nabla c - \frac{\text{Pe}}{2}\nabla\rho + \frac{\text{Pe}^2}{16}\nabla^2\mathbf{w} - \frac{B^2|\nabla c|^2}{8}\mathbf{w}$$

$$+ \frac{\text{Pe}B}{16}[3(\nabla\mathbf{w})^T \cdot \nabla c - (\nabla c \cdot \nabla)\mathbf{w} - 3(\nabla \cdot \mathbf{w})\nabla c],$$

$$\dot{c} = \mathcal{D}\nabla^2c + K_0\rho + \nu\frac{K_0}{2}\nabla \cdot \mathbf{w} - K_dc.$$
(4)

Overview

