## Master 2: International Centre for Fundamental Physics

## INTERNSHIP PROPOSAL

Laboratory name: PHENIX (Physico-chimie des Électrolytes et Nanosystèmes Interfaciaux)

CNRS identification code: UMR 8234

Internship director's urname: Pierre ILLIEN (co-supervisors: V. DAHIREL, M. JARDAT)

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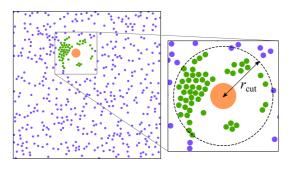
Internship location: Sorbonne Université, campus Jussieu, Paris 5ème

Thesis possibility after internship: YES

Funding: École doctorale (to be discussed with the candidate)

## Spontaneous propulsion of a particle in a phase-separating environment

Self-propelled particles, like **active colloids**, have been the subject of numerous theoretical and experimental studies [1, 2]. Among the possible routes to locomotion, the design of anisotropic colloids, which interact with self-generated gradients of solute concentration, temperature, or electric fields, has been particularly fruitful. However, there was a more recent interest in the propulsion of **isotropic colloids**. Several examples revealed that **built-in asymmetry** of the colloid is actually not necessary to achieve directed motion over long timescales and that a **spontaneous polarization** of its environment can be sufficient.



In a recent work, we have proposed and studied numerically a model for an isotropic colloid in a bath of solute particles that interact with each other [3]. Within our model, a finite-size mesoscale phase transition of a Lennard-Jones fluid can be locally maintained close to the colloid by a non-equilibrium chemical reaction. As a result of phase coexistence, strong density fluctuations can

emerge and considerably enhance the diffusion coefficient of the colloid. The propulsion mode that we have characterized through Brownian dynamics simulations does not require a structural anisotropy of the colloidal particle, as it arises from persistent density fluctuations in its environment. Strikingly, activity is therefore solely driven by **attractive interactions** within the solvent bath.

The goal of this internship is to make progress towards an **analytical model** for the propulsion phenomenon observed in the numerical simulations reported in [3]. A first step would consist in calculating the diffusion coefficient of a tracer particle coupled to a binary suspension that displays **phase separation**. Depending on the taste of the student, the analytical approach will be completed by **numerical simulations** (Brownian and/or molecular dynamics).

## **References:**

- [1] P. Illien, R. Golestanian, A. Sen, Chem. Soc. Rev. 46, 5508 (2017). arXiv:2104.03878.
- [2] C. Bechinger et al., Rev. Mod. Phys. 88, 045006 (2016). arXiv:1602.00081.
- [3] J. Decayeux, V. Dahirel, M. Jardat, P. Illien, Phys. Rev. E. 104, 034602 (2021). arXiv:2103.13244.

Condensed Matter Physics: YES	Soft Matter and Biological Physics:	YES
Quantum Physics: NO	Theoretical Physics:	YES