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Microemulsions: a new model for organic phases involved in ion separation methods

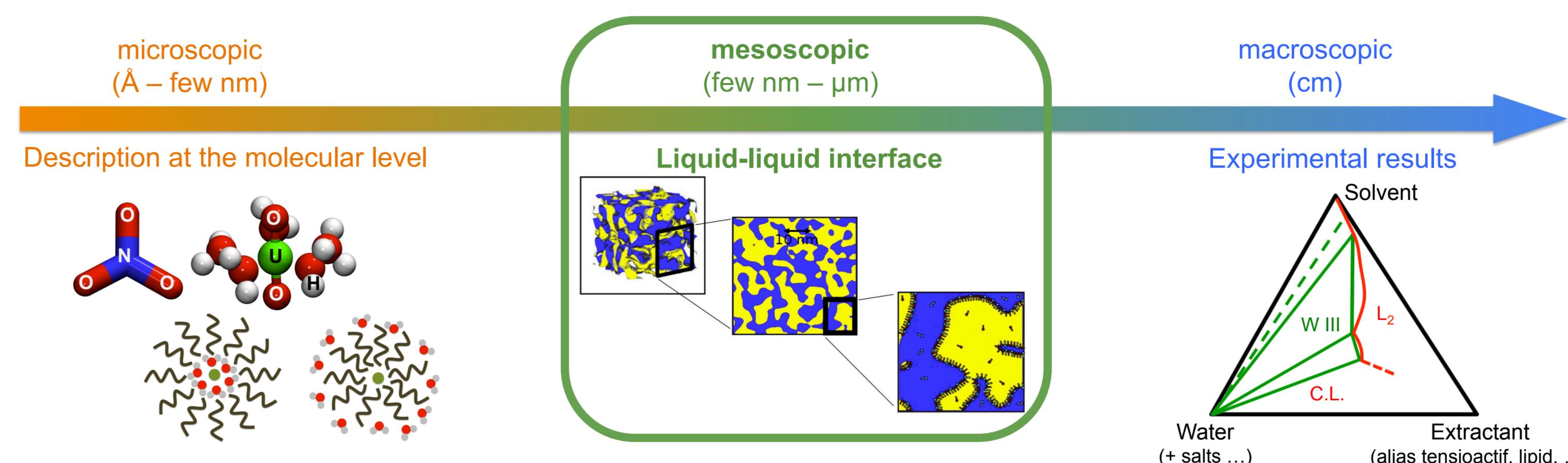
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Introduction

Context

- ✓ Improvement of the reprocessing of spent nuclear fuel
- ✓ Liquid-liquid extraction equilibria
- ✓ Microemulsion model to describe the organic phase
- ✓ Multi-scale approach



Method

- ✓ The physics of a system composed of **soft interfaces** (low bending energy compared to the thermal one, negligible long-range electrostatic and steric interactions) is **dominated** by the **Helfrich Hamiltonian** of the surfactant interface (between oil and water) [1].
- ✓ In our model, the **surfactant interface** is described by appropriate **two level-cuts** of a **Gaussian random field** based on wavelets proposed by Arleth *et al.* [2].
- ✓ The **free energy density** of the Gaussian random field with level-cuts can be approximated by:

$$f_{\text{free}} = \frac{S}{V} [2\kappa \langle (H - H_0)^2 \rangle + \bar{\kappa} \langle K \rangle] - \frac{1}{2\pi^2} \int_k k^2 \ln \nu(k)$$

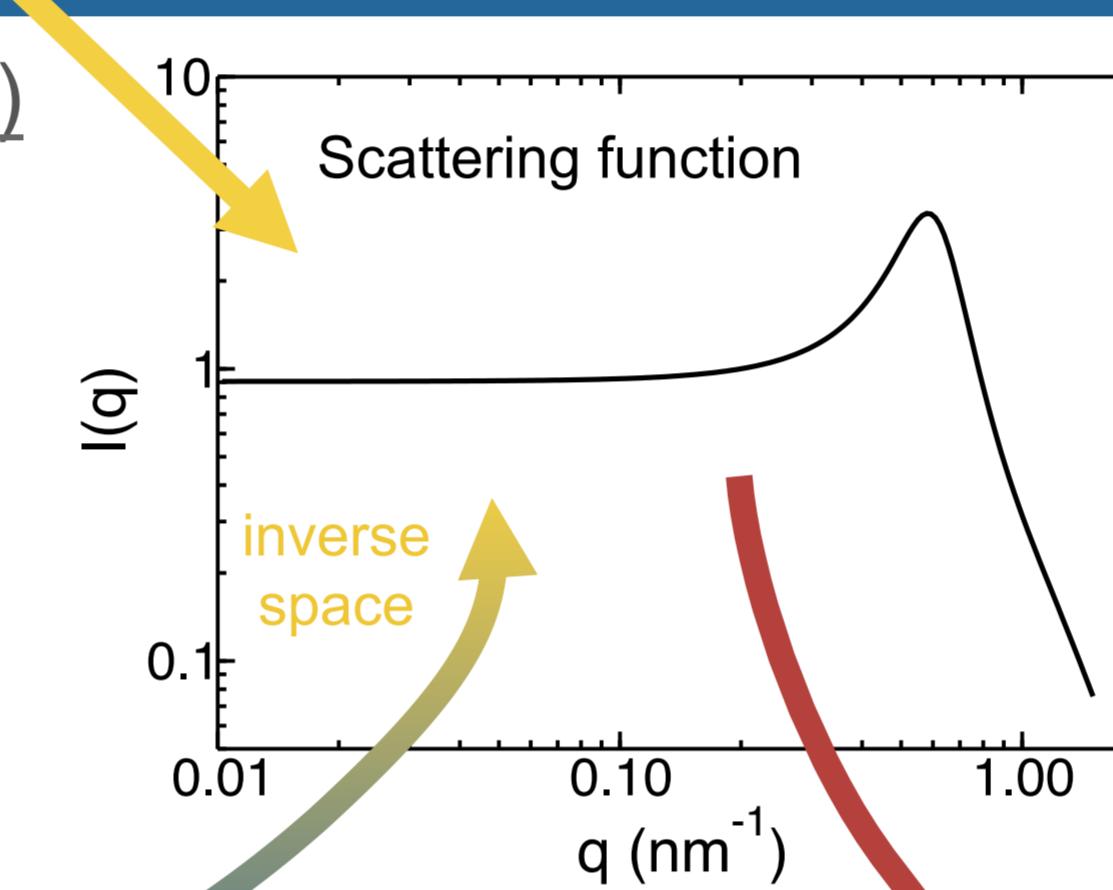
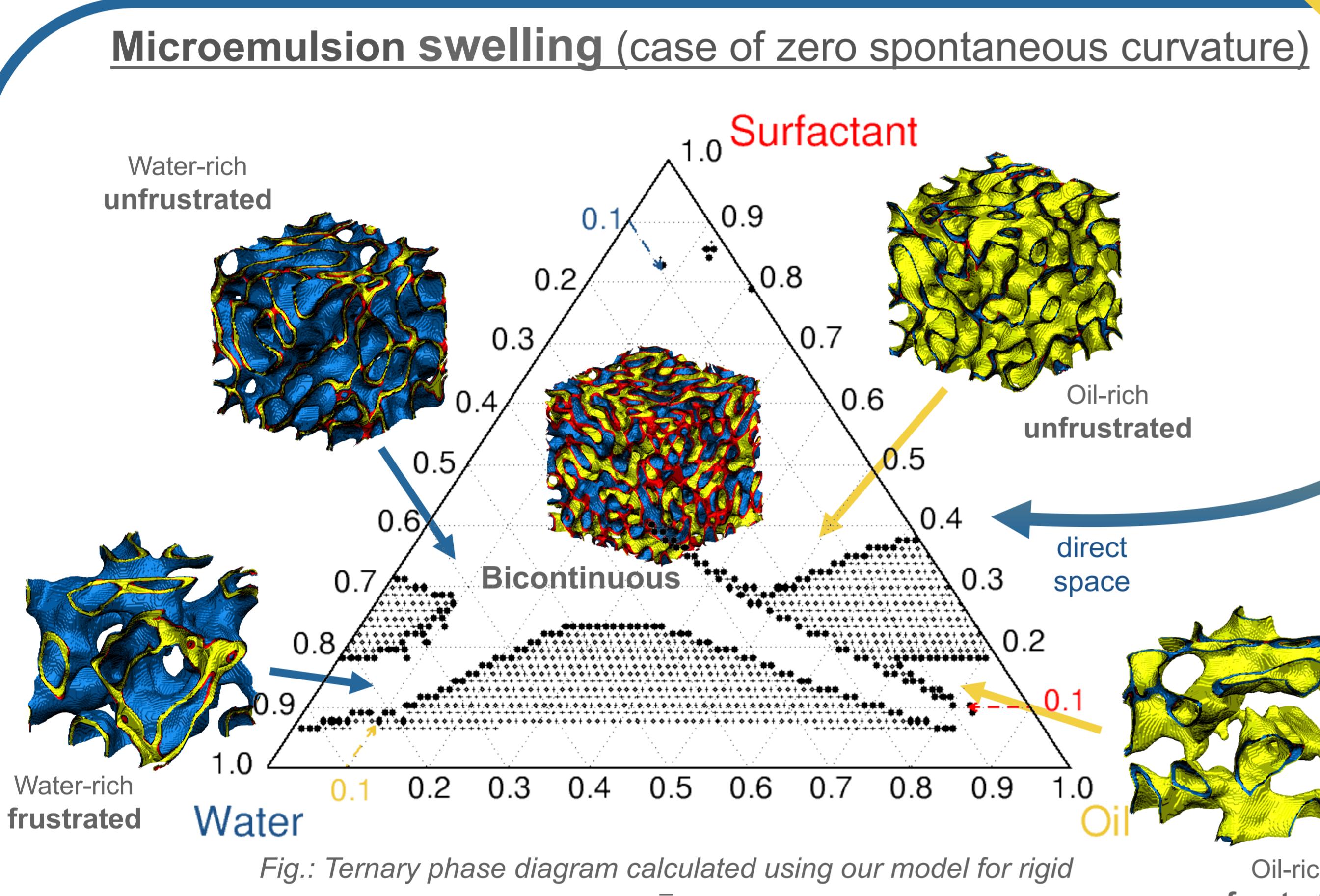
Rigidity controlled by κ and $\bar{\kappa}$

- ✓ Two constants:
 - the bending elastic constant κ
 - the Gaussian elastic constant $\bar{\kappa}$
 - ✓ One unique constant:
 - $\kappa^* = 2\kappa + \bar{\kappa}$
 - Analogy with the packing parameter p_0
- $$f_{\text{free}} = \frac{1}{2} \kappa^* (p - p_0)^2$$

Curvatures:

- ✓ Spontaneous curvature H_0 imposed
 - ✓ Mean average curvature H
$$H = \frac{1}{2} (c_1 + c_2)$$
 - ✓ Gaussian curvature K
$$K = (c_1 \times c_2)$$
-

Results



$$D^* \Sigma = \frac{2\pi}{q_{\max}} \frac{\phi_s}{l_s}$$

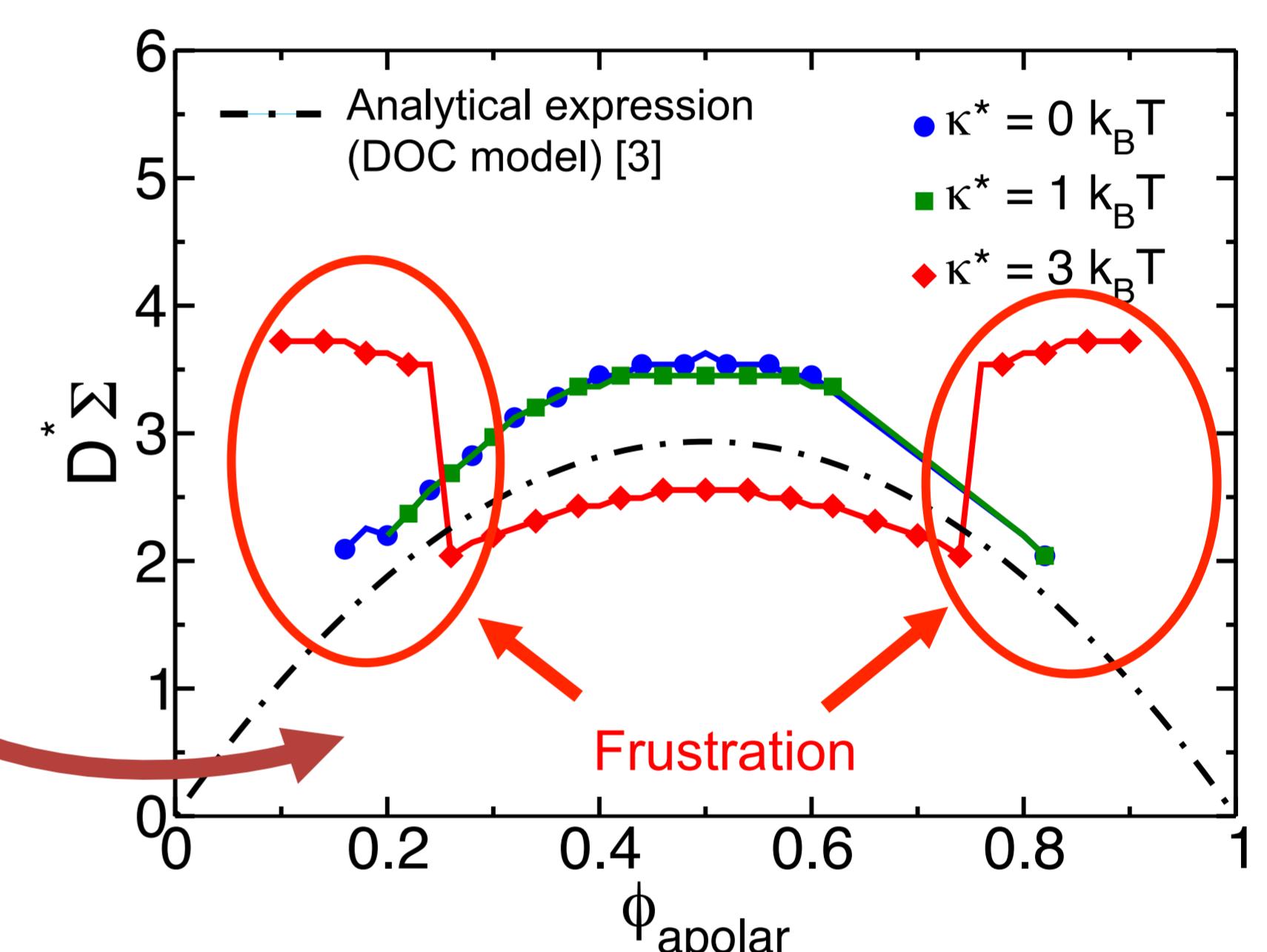


Fig.: Scaled peak position for flexible (blue and green) and rigid (red) microemulsion plotted as a function of the volume fraction of the apolar phase for a volume fraction of surfactant $\Phi_s = 0.15$.

- ✓ Spinodal instabilities calculations (case of rigid surfactant film)
- ✓ Thermalized unfrustrated, frustrated and bicontinuous microemulsions

- ✓ Frustration appears only in the case of **rigid microemulsions**
- ✓ Unfrustrated bicontinuous microemulsions for $0.2 < \Phi_{\text{apolar}} < 0.8$

Influence of the spontaneous curvature (case of rigid microemulsions)

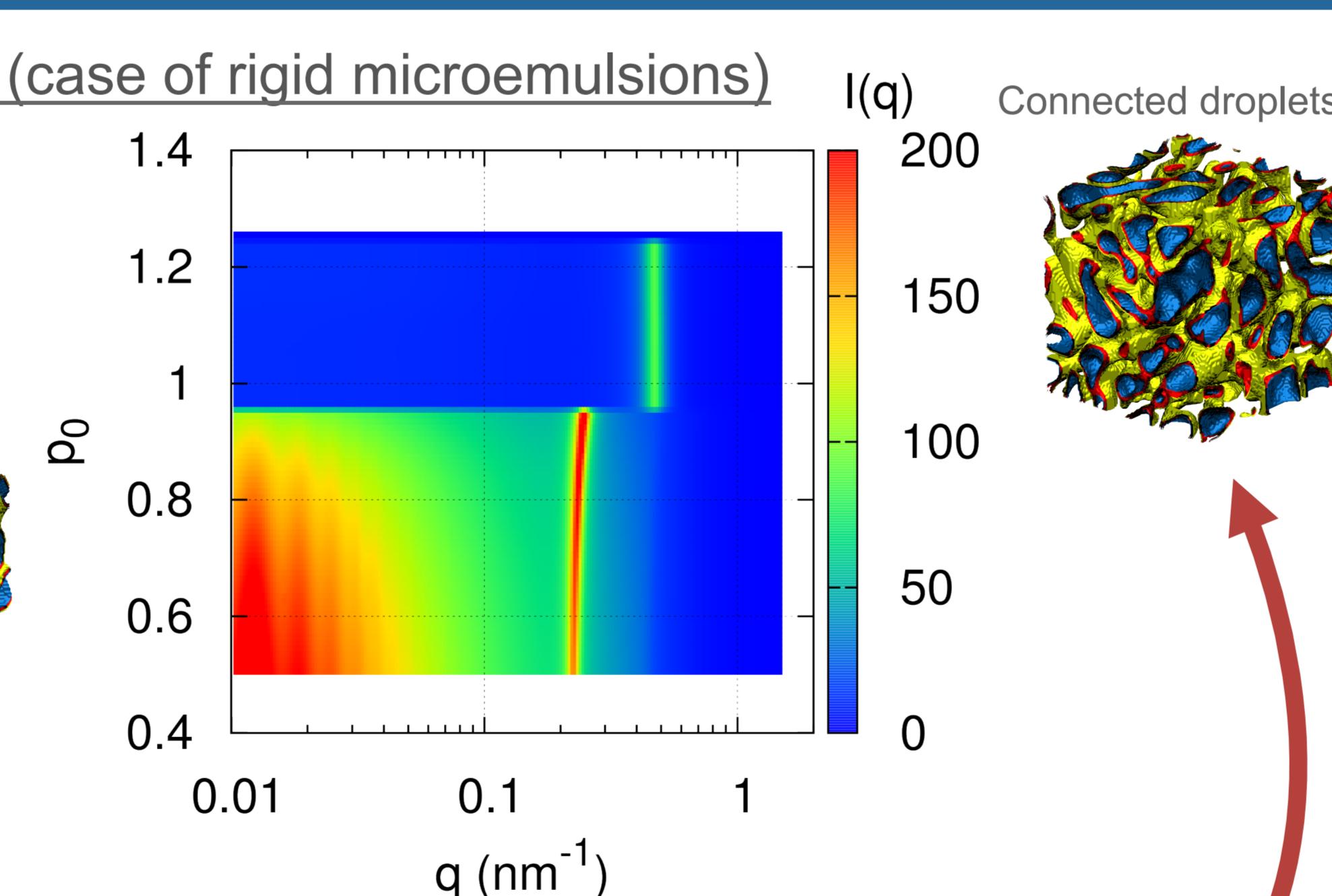
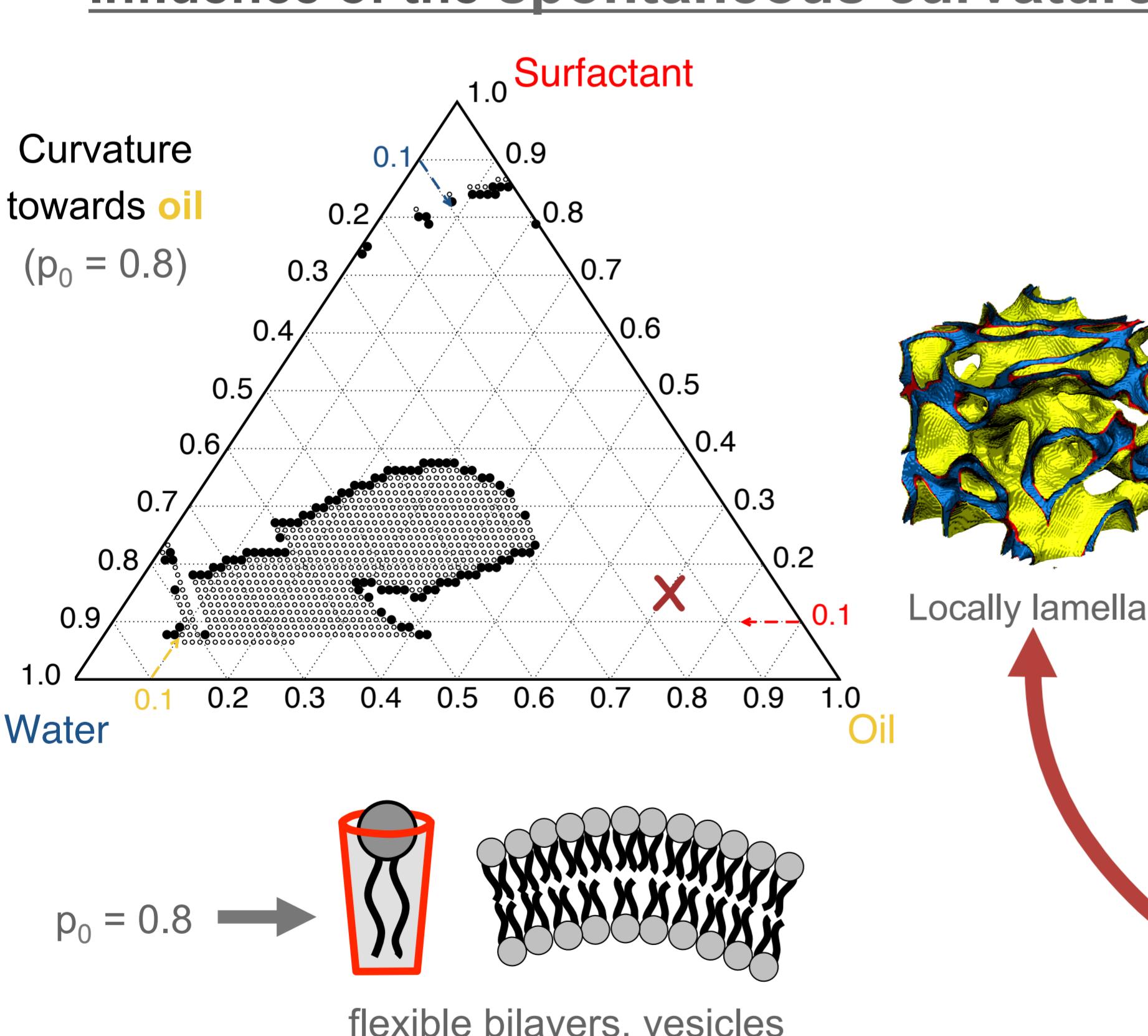
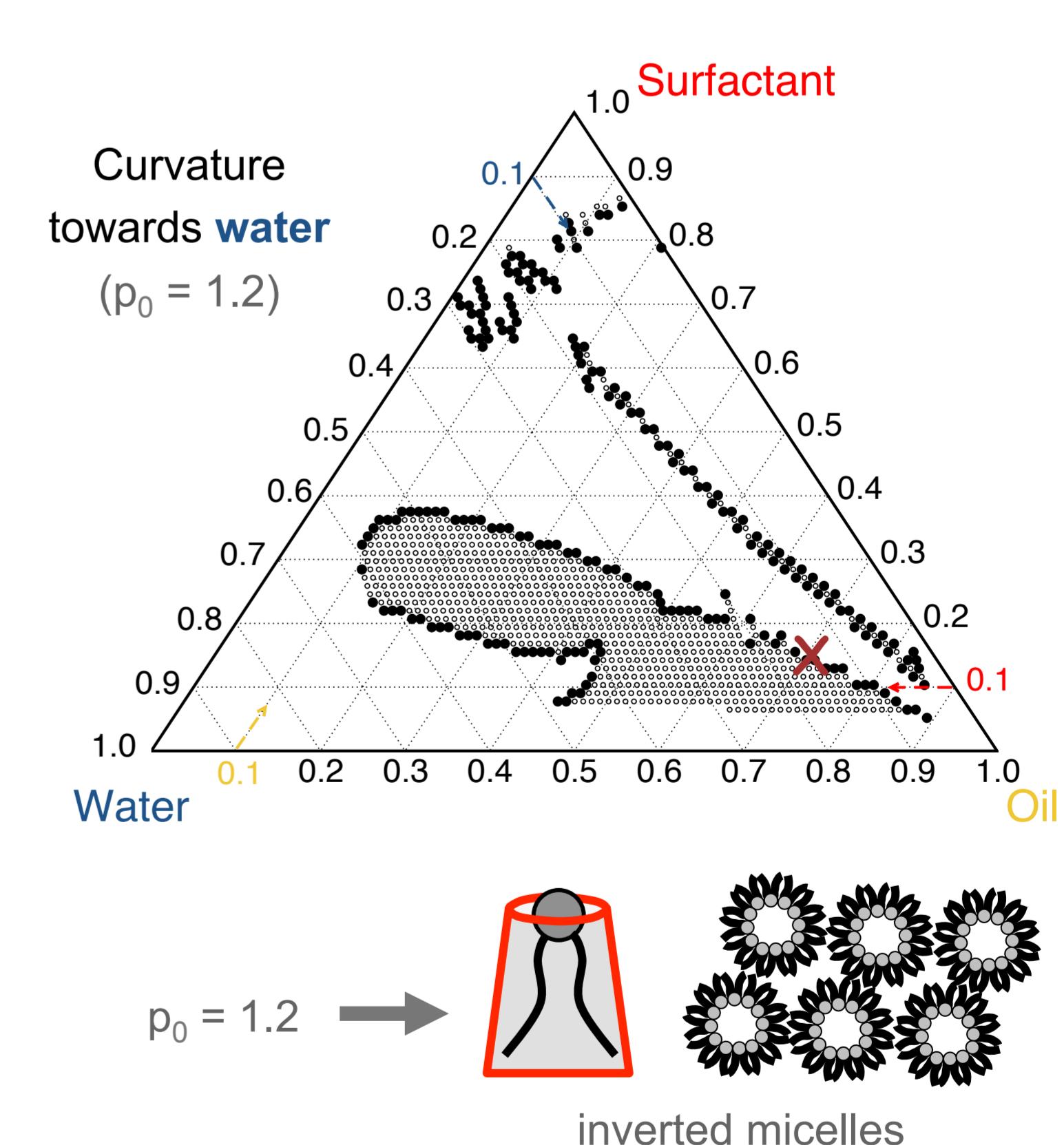


Fig.: 2D map of the scattering functions as a function of the spontaneous packing parameter p_0 calculated for $x_{\text{water}} \approx 0.15$, $x_{\text{oil}} \approx 0.70$ and $x_{\text{surfactant}} \approx 0.15$.

Non-continuous transition



Acknowledgments

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References

- [1] W. Helfrich, *Z. Naturforsch. C* **28**, 693 (1973)
- [2] L. Arleth *et al.*, *J. Chem. Phys.* **115**, 3923 (2001)
- [3] J. Jouffroy *et al.*, *J. Phys.* **43**, 1241 (1982)

Conclusions and outlooks

Our model predicts:

| Microstructures | Scattering | Phase diagrams |
|-----------------|------------|----------------|
| | | |

Comparison with experiments

Improvement of our model to take into account:

- ✓ the presence of **charged species** (ions and ionic surfactant),
- ✓ the **aggregation** of surfactant molecules.

