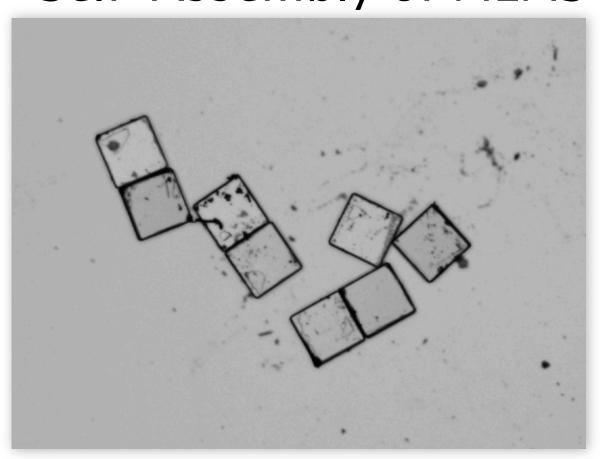


Controlling and Exploiting Self-Assembly of MEMS





Speaker: Grégory Mermoud^{‡ *}

Team: Vahid Fakhfouri*, Prof. Alcherio Martinoli *, Prof. Juergen Brugger*

* Swarm Intelligent Systems Group (I&C)

*Laboratory of Microsystems and Nanoengineering (STI)





A lot of bricks, but where is the plan?



Today

An avalanche of beautiful micro/ nanobricks: nanowires, nanorods, nanotubes, DNA strands, nanoparticles, microspheres, nanocrystals, cells, etc.

A lot of bricks, but where is the plan?



Today

An avalanche of beautiful micro/ nanobricks: nanowires, nanorods, nanotubes, DNA strands, nanoparticles, microspheres, nanocrystals, cells, etc.



The challenge of tomorrow

Finding a way to (self-)assemble these bricks into functional micro/nano-systems

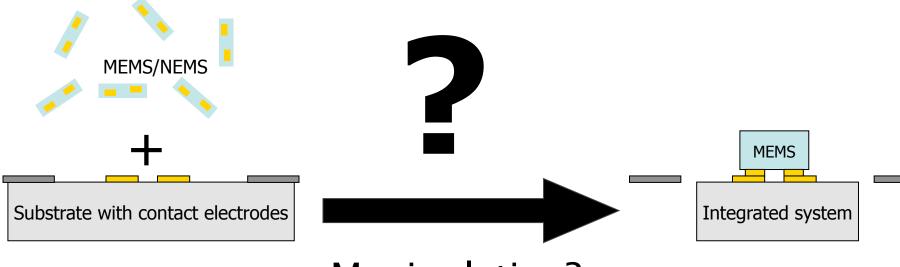
2 fields + 2 motivations + 2 frameworks = 1 challenge

Technology (LMIS1 expertise)

Methodology (SWIS expertise)

Motivation: technology

- Lithography: great advances in engineering of devices in the size range of 0.1–100 μm, but...
- How to build non-monolithic, hybrid systems assembled from a set of components which are coming from different fabrication processes?

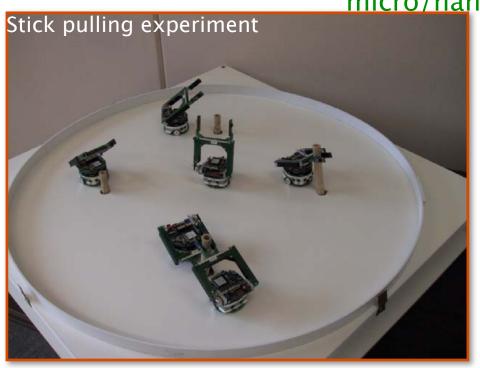


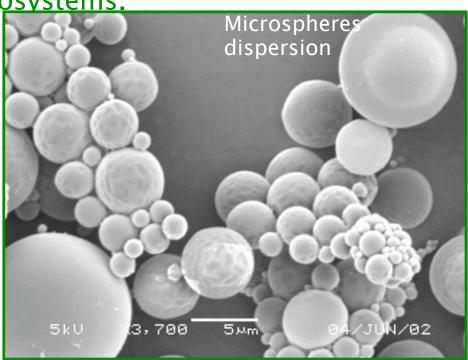
Manipulation? Alignment? Assembly?

Motivation: methodology

We aim at applying our expertise in model-based design and optimization of swarm robotics system to the design of self-assembling

micro/nanosystems.



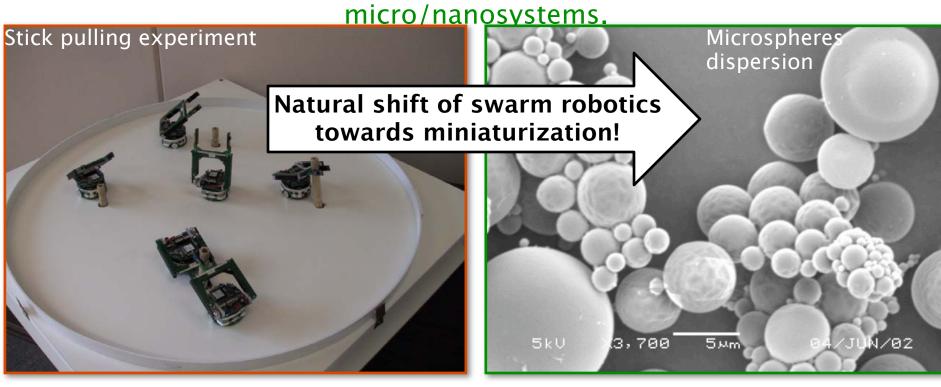


- Typical length scale: a few centimeters
- Typical swarm size: tenth to hundreds
- Sensing, computation, communication

- Typical length scale: 1 nm to 500 um
- Typical swarm size: 10³ to 10⁶ units
- No sensing, no computation, no communication, but local interactions

Motivation: methodology

We aim at applying our expertise in model-based design and optimization of swarm robotics system to the design of self-assembling

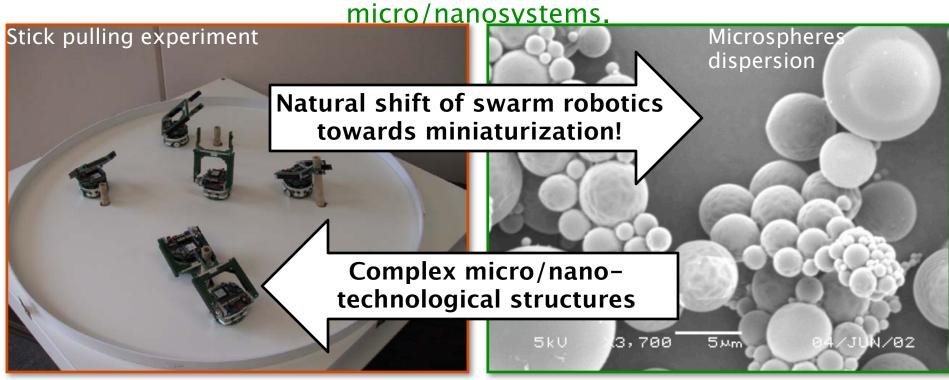


- Typical length scale: a few centimeters
- Typical swarm size: tenth to hundreds
- Sensing, computation, communication

- Typical length scale: 1 nm to 500 µm
- Typical swarm size: 10³ to 10⁶ units
- No sensing, no computation, no communication, but local interactions

Motivation: methodology

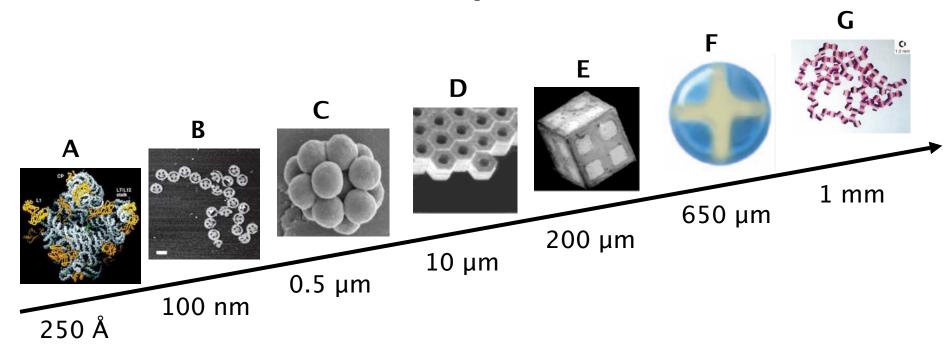
We aim at applying our expertise in model-based design and optimization of swarm robotics system to the design of self-assembling



- Typical length scale: a few centimeters
- Typical swarm size: tenth to hundreds
- Sensing, computation, communication

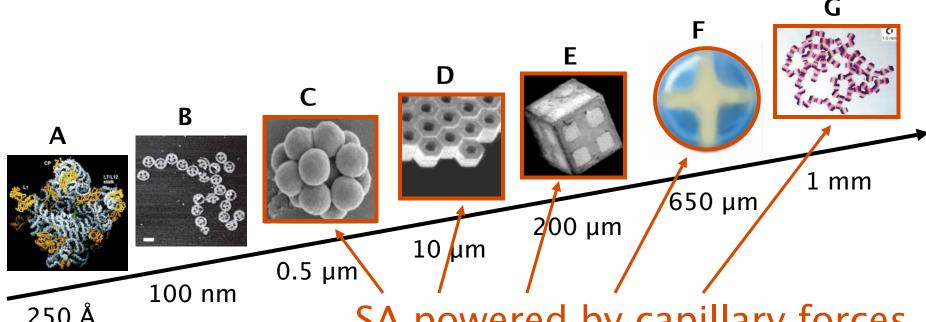
- Typical length scale: 1 nm to 500 µm
- Typical swarm size: 10³ to 10⁶ units
- No sensing, no computation, no communication, but local interactions

Self-assembly at all scales



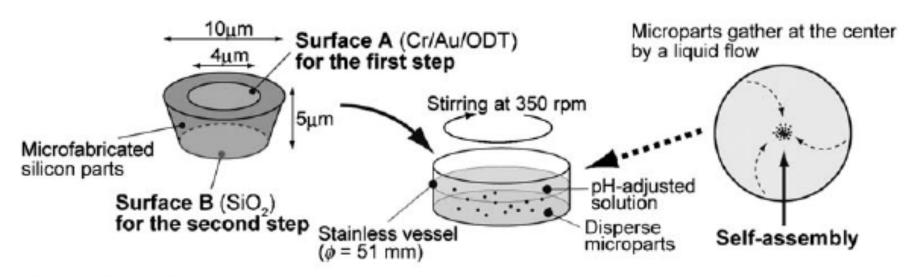
- A. Crystal structure of a ribosome [Yusupov et al., 2001]
- B. Scaffolded DNA origami of 100 nm in diameter [Rothemund et al., 2006]
- C. Clusters of 0.5-µm-sized spheres packed by surface tension [Manoharan, 2003]
- D. Capillary SA of 10-µm-sized hexagonal plates into crystal structures [Clark, 2001]
- E. Capillary SA of 200-µm-sized polyhedra [Gracias et al., 2002]
- F. Capillary SA of polymeric spheres of complex internal structures from pipetted droplets [Fialkowsi et al., 2005]
- G. SA of 1-mm-sized hexagonal plates using "capillary bond" [Bowden et al., 2001]

Self-assembly at all scales

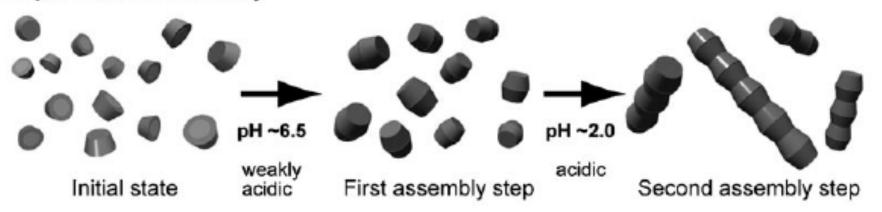


- 250 Å
 A. Crystal structure of a ribosome [Yusupov et al., 2001]
- B. Scaffolded DNA origami of 100 nm in diameter [Rothemund et al., 2006]
- C. Clusters of 0.5-µm-sized spheres packed by surface tension [Manoharan, 2003]
- D. Capillary SA of 10-µm-sized hexagonal plates into crystal structures [Clark, 2001]
- E. Capillary SA of 200-µm-sized polyhedra [Gracias et al., 2002]
- F. Capillary SA of polymeric spheres of complex internal structures from pipetted droplets [Fialkowsi et al., 2005]
- G. SA of 1-mm-sized hexagonal plates using "capillary bond" [Bowden et al., 2001]

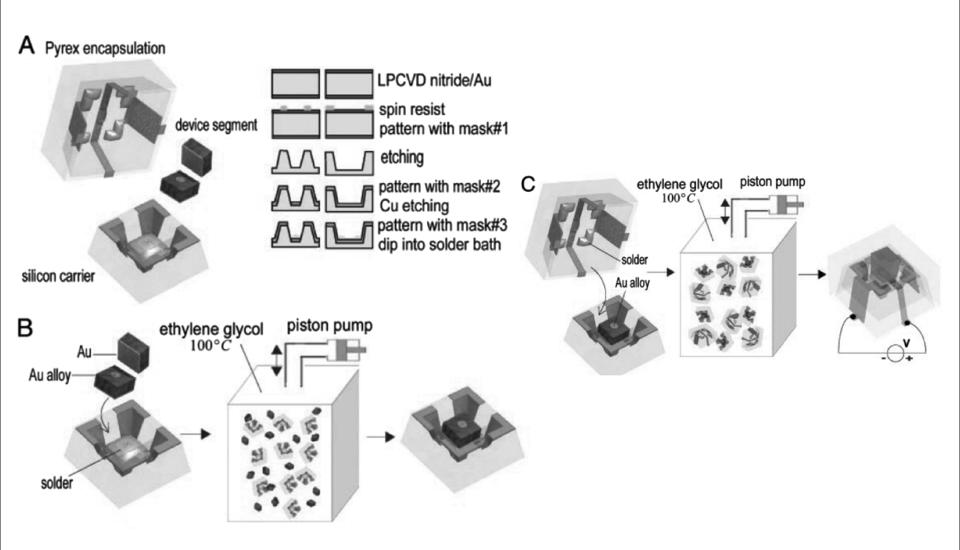
Hydrophobic interaction (part-topart)



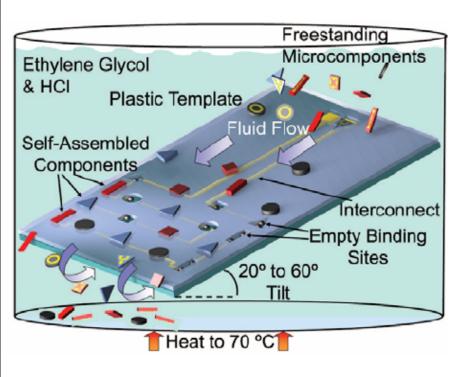
Sequential self-assembly

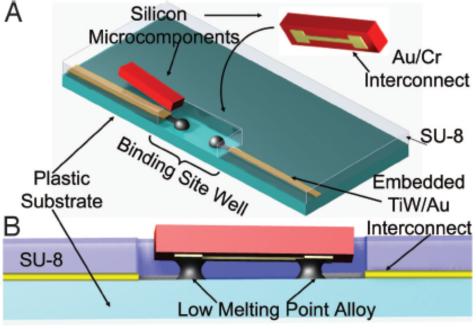


Capillary interaction (part-to-part)



Capillary interaction (part-to-



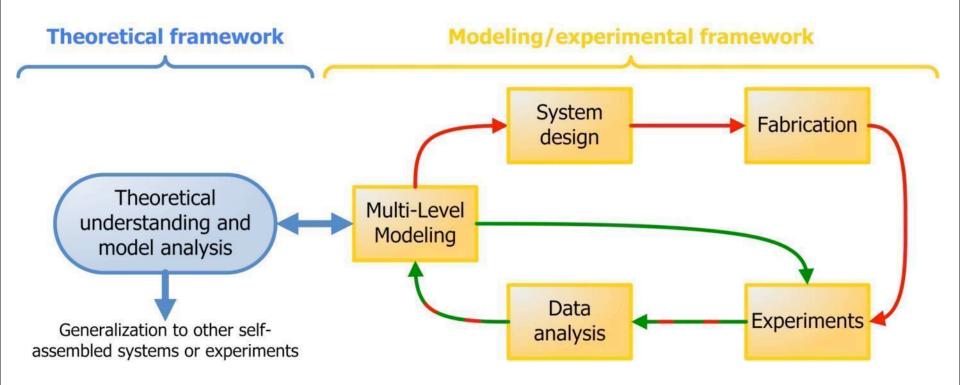


Putting several mechanisms together:

- A dramatic increase in interaction variety (in magnitude, length- and time-scales)
- The apparition of competing interactions (attractive vs repulsive)
- Dynamical steady states resulting from a balance between competing interactions
- Non-linear positive or negative feedbacks resulting from the intricacy of the interaction scheme

These features motivate an efficient and appropriate methodological framework

Methodological framework



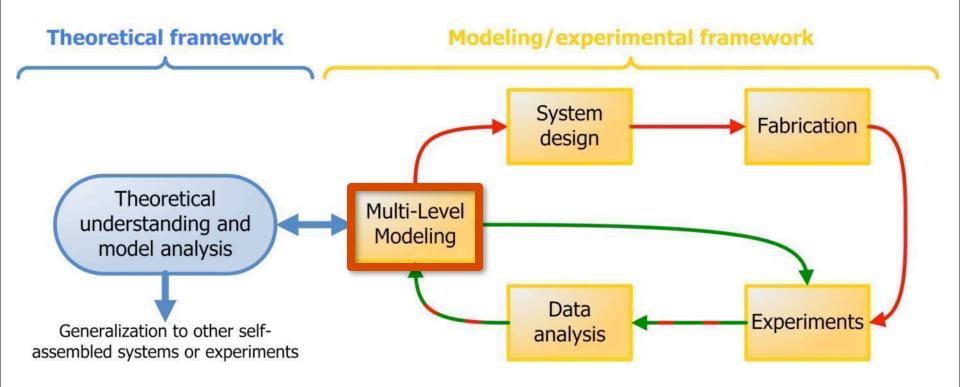
Two ingredients:

- Multi-level modeling
- Experimental work

Two experimental loops:

- Fabrication (slow)
- Modeling (fast)

Methodological framework



Two ingredients:

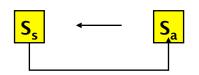
- Multi-level modeling
- Experimental work

Two experimental loops:

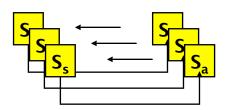
- Fabrication (slow)
- Modeling (fast)

Multi-Level Modeling Methodology

$$\frac{dN_n(t)}{dt} = \sum_{n} W(n \mid n', t) N_n'(t) - \sum_{n} W(n' \mid n, t) N_n(t)$$



Macroscopic: rate equations, mean field approach, whole swarm



Microscopic: multi-agent models,only relevant robot feature captured,1 agent = 1 robot



Realistic: intra-robot (e.g., S&A) and environment (e.g., physics) details reproduced faithfully



Physical reality: Info on controller, S&A, morphology and environmental features

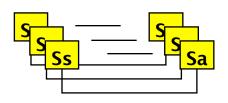
$$AB + A \Longrightarrow AAB$$

$$K_5 = 3 K_{AA} K_{AB}^2 = \frac{[AAB]}{[A][A][B]}$$

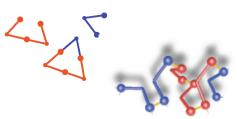
Macroscopic 1: Chemical equilibrium is completely defined by equilibrium constants K of each reaction (law of mass action)

$$\frac{d[\mathsf{AAA}]}{dt} = [\mathsf{AA}][\mathsf{A}] \, k_{f\mathsf{AA}}^2 - [\mathsf{AAA}] \, k_{r\mathsf{AA}}^2$$

= $[AA][A] k_{fAA}^2 - [AAA] k_{rAA}^2$ \longrightarrow Macroscopic 2: Reactions kinetics describes how a reaction occurs and at which speed (differential equations)



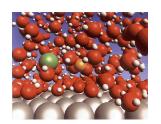
Microscopic 1: Agent-Based model, molecules geometry abstracted, 1 agent = 1 aggregate



Microscopic 2: Agent-Based model,

molecules

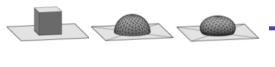
2D- and 3D geometry captured, 1 agent = aggregate



Physical reality: microscopic

(e.g., crystallography) and macroscopic measurements (chemical reaction)

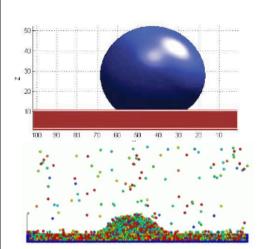
MLMM applied to Droplet Motion



Macroscopic 1: Surface evolver, surface energy minimization by gradient descent, no dynamics

$$o\left(rac{\partial \mathbf{v}}{\partial t} + \mathbf{v}\cdot
abla \mathbf{v}
ight) = -
abla p +
abla \cdot \mathbb{T} + \mathbf{f}$$

 $\rho\left(\frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v}\right) = -\nabla p + \nabla \cdot \mathbf{T} + \mathbf{f} \longrightarrow \mathsf{Macroscopic 2: Navier-Stokes equations}$



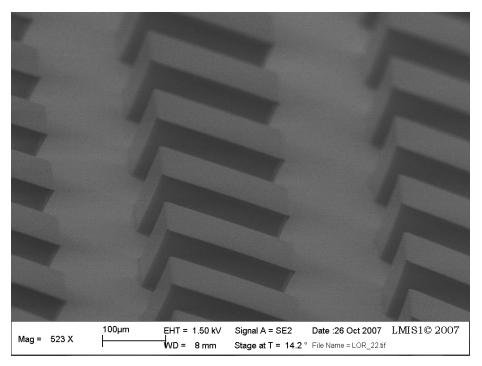
Mesoscopic: Lattice Boltzmann Method, particle distribution, discrete velocities and directions

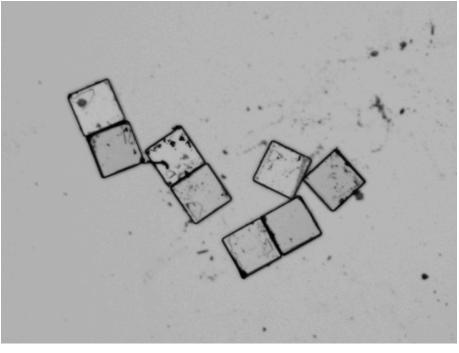
Microscopic: Molecular dynamics, 1 agent = 1 molecule



Physical reality: contact angles measurements, dynamics imaging with high speed camera

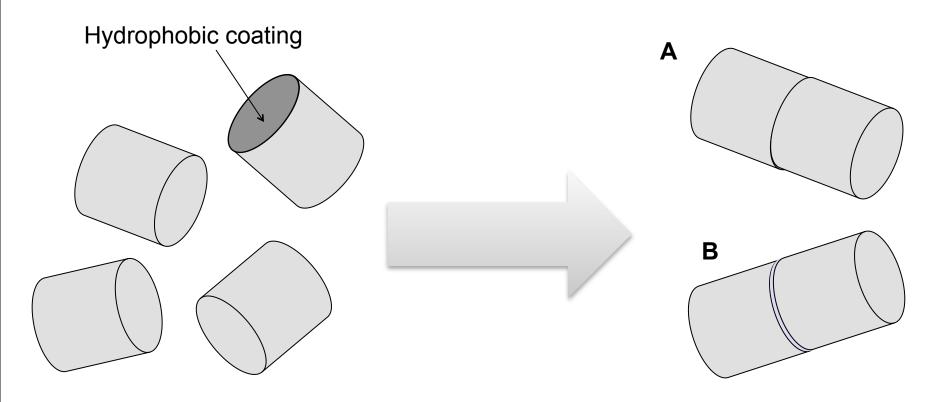
Our preliminary case study





- 200µm-sized SU-8 building blocks
- Fabrication and release is under control
- Next batch: smaller, barrel-shaped building blocks to be assembled into pairs

Our preliminary case study



- Hydrophobic interaction (A) or capillary interaction mediated by an hydrophobic solder (B)
- Same methodology as in robotics: multi-level modeling, imaging (including tracking)