

# Multiscale Mechanics of Polymers, Soft Matter and Network Materials

- **Organizers**

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- **Symposium Description**

Polymers and soft matter have been of critical importance in a broad variety of engineering and biological contexts at a wide range of scales. Recently, there have also been tremendous progresses in advanced functional network materials (e.g. vitrimers, hydrogels), that achieve excellent performance (e.g. toughness, self-healing) through the designs of atomistic, molecular and supramolecular structures within reversible or double networks. The multi-scale and multi-physical processes in these materials are strongly coupled, often resulting in highly nonlinear and emergent phenomena. To elucidate the coupled behaviors in these materials across the scales, multi-scale mechanical modeling approaches are essential.

The main goal of this Symposium is to bring together scientists and engineers from a broad variety of engineering, physics and chemistry disciplines to discuss state-of-the-art in the emerging fields of multi-scale mechanics of polymers, soft matters and network materials.

We welcome combined approaches of theoretical modeling, experimentation and computation that address (1) structure-property-function relationships, (2) macroscopic mechanical properties and their connections to micromechanical features (3) multi-scale interplays between geometry, topology and mechanics, (4) multi-physically coupled phenomena, e.g. electromagnetic fields, biological processes, chemical reactions and deformations in polymers, soft matter and network materials, and (5) advanced/additive manufacturing processes for these materials (e.g. electrospinning, 3D printing).

- **Topics (but are not to limited to)**

1. Mechanics and rheology of polymers and network materials
2. Stimuli-responsive polymers, polymeric composites, hydrogels, vitrimers, cellular materials, foams and mechanical meta-materials for new applications in engineering and biological contexts
3. Multiscale computational methods: molecular simulations (classical molecular dynamics simulations, Monte-Carlo methods), coarse-graining strategies, continuum simulations (non-conventional finite elements), *ab-initio* or density functional for polymers and soft materials
4. Mechanochemistry and multi-physically coupled deformations in polymers and soft materials

- **Invited Speakers**

We propose ~ 40 oral presentations for this symposium. Further, we have 7 confirmed invited lecturers (we may have 1-2 more invited lecturers):

**Kaushik Bhattacharya**, California Institute of Technology

**Masao Doi**, Beihang University and the University of Tokyo

**Maurizio Fermeglia**, University of Trieste

**Jian Ping Gong**, Hokkaido University

**Ludwik Leibler**, ESPCI Paris

**Thao (Vicky) Nguyen**, Johns Hopkins University

**Jerry H. Qi**, Georgia Institute of Technology