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MOLECULAR SIMULATIONS OF COMPLEX FLUIDS AND INTERFACES



Recent Advances and Challenges

FEBRUARY 21 - 23, 2020
IIT KANPUR



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About



Welcome

Welcome to the international symposium "Molecular Simulations of Complex Fluids and Interfaces", hosted at IIT Kanpur.

The behaviour of interfaces plays an important role in several industrial and natural processes. Molecular simulations can reveal microscopic insights into the structure and properties of solid-liquid interfaces. This meeting aims to provide a forum for exchanging ideas and sharing recent scientific advances from several perspectives. It is hoped that the current state of simulation methodologies will be established, paving the way for the future development of computational tools and research.

Topics

- Molecular Simulations: New Methodologies and Applications
- Advances in Coarse-Graining and Challenges
- Soft Matter Simulations
- Confined Fluids
- Wetting and Interfacial Phenomena
- Biomolecular Simulations
- Specific Systems and Models

Timetable

KL: Keynote Lecture, IS: Invited Speaker, ST: Sponsored Talk.

February 21, Friday

15:00–16:00		Registration	
16:00–16:15		Inauguration Ceremony	
16:15–17:15	KL	Florian Müller-Plathe Technische Universität Darmstadt	Wetting, Drying, Adhesion
17:15–17:40	IS	Neelanjana Sengupta IISER Kolkata	Modulating Self-Assembled Amyloidogenic States via Solvent and Temperature: Insights from Computer Simulations
17:40–18:05	IS	Harshwardhan Katkar IIT Kanpur	Multiscale Modeling of Actin Filaments
18:05–18:30	IS	Manjesh K. Singh IIT Kanpur	Rheology of Nonequilibrium Polymer Melts
19:00–21:30		Reception Dinner	

February 22, Saturday

9:30–10:30	KL	Edward Maginn University of Notre Dame	Computational Design of New Materials for Separations and Energy Storage
10:30–11:00	Tea Break		
11:00–11:25	IS	Beena Rai Tata Research Development and Design Center	Skin Lipids and their Interfaces: A Computational Approach towards Mimicking Skin Barrier Function
11:25–11:50	IS	Sudip Roy National Chemical Lab, India	Bridging Scales for Simulation of Lipids
11:50–12:15	IS	Swati Bhattacharya IIT Bombay	Molecular Dynamics Investigations of anti HIV-1 protein SAMHD1
12:15–12:40	IS	R. Sankararamakrishnan IIT Kanpur	TBA
12:40–14:00	Lunch Break		
14:00–15:00	KL	Balasubramanian Sundaram JNCASR	Modelling Supramolecular Polymers
15:00–15:25	IS	Divya Nayar IIT Kharagpur	Microscopic View of the Crowding Effects on Hydrophobic Collapse
15:25–15:50	IS	Nisanth N. Nair IIT Kanpur	Exploration of High Dimensional Free Energy Landscapes of Chemical Reactions
15:50–16:20	Tea Break		
16:20–16:45	IS	Rajat Srivastava Politecnico di Torino	Thermo-Physical Properties of Graphene Reinforced Thermoplastics: A Coarse-Grained Modelling Approach
16:45–17:10	IS	Tamal Banerjee IIT Guwahati	Reactive Force Field Simulations on the Degradation of Quinoline
17:10–17:30	ST	Ashwini Kumar NEC Corp.	Vector Computing- Simulation and A.I.
17:30–18:55	Poster Presentation Session		
18:55–19:00	Announcement of Springer Best Poster Awards		

February 23, Sunday

9:30 –10:30	KL	David A. Kofke SUNY, Buffalo	'Mapped Averaging' Method for deriving Ensemble Averages: Application to Crystals
10:30–11:00	Tea Break		
11:00–11:25	IS	Shantanu Maheshwari Shell Technology	Nucleation and Growth of a Nanobubble on Rough Surfaces
11:25–11:50	IS	Kaustubh Rane IIT Gandhinagar	The Role of Solid-Liquid Interfacial Fluctuations in the Spontaneous Motion of Droplets
11:50–12:15	IS	Sandip Khan IIT Patna	The Wetting Behavior of Imidazolium Based Ionic Liquids using Molecular Dynamics Simulation
12:15–12:40	IS	Jhumpa Adhikari IIT Bombay	Phase Equilibria of Binary Mixtures of Triangle-Well Fluids : Bulk vs Confined Systems
12:40–14:00	Lunch Break		
14:00–14:25	IS	Sudeep Punnathanam IISC Bangalore	Computing Solid-Liquid Interfacial Free Energy via Thermodynamic Integration
14:25–14:50	IS	Subir K. Das JNCASR	Kinetics of Clustering in an Assembly of Vicsek-Like Active Particles
14:50–15:15	IS	Tarak Patra IIT Madras	Correlation between Glass Formation and Ion Conductivity in Polymeric Ionic Liquids
15:15–15:40	IS	Prateek Kumar Jha IIT Roorkee	Multiscale Modeling Approaches in Excipient Design for Oral Drug Delivery
15:40–16:10	Tea Break		
16:10–16:35	IS	Sk. Musharaf Ali BARC, Mumbai	Microscopic Assessment of Liquid-Liquid Extraction System in Bulk and at the Interface
16:35–17:00	IS	Vishwanath Dalvi ICT, Mumbai	Study of Water Extraction by Phosphate Ligands by Experiments and Molecular Simulations
17:00–17:25	IS	Rajat Desikan Invictus Oncology Pvt. Ltd.	Accurate Computational Calorimetry of Lipid Membranes by void-induced Melting
17:25–17:35	Springer Poster Awardee Talk-1		

17:35–17:45	Springer Poster Awardee Talk-2
17:45–18:00	Vote of Thanks
19:00–21:30	Symposium Dinner

List of Abstracts – Talks

February 21, Friday

Wetting, Drying, Adhesion

F. Müller-Plathe

KL

Technische Universität Darmstadt

We will review recent advances on the simulation of small-molecule fluids as well as polymers at solid interfaces. We will cover wetting of surfaces by liquids as well as the evaporation of liquid droplets from them. We will also look at the structure formation of polymers near solid surfaces and the resulting interphase properties, such as encountered in coatings and nanocomposites.

February 22, Saturday

Skin Lipids Their Interfaces: A Computational Approach Towards Mimicking Skin Barrier Function

B. Rai

IS

Tata Research Development and Design Center, India

Human skin is a vital organ acting as an interface between us and our surroundings. It is one of the largest organ of our body comprising of three layers – Epidermis, Dermis and Hypodermis. While epidermis is responsible for overall appearance and texture of skin, its outermost layer called stratum corneum (SC) controls its barrier function. Dermis, primarily made up of collagen and elastin, provides structural support and elasticity. The deepest layer – hypodermis, is composed of adipose tissues and provides the heat resistance. SC is composed of 15~20 layers of flattened cells called corneocytes (Bricks) which are embedded in a lipid matrix (Mortar) composed of ceramides, cholesterol, and fatty acids. The “bricks and mortar” structure of SC makes it selectively permeable thus to protecting underlying tissue from infection, dehydration, chemicals and mechanical stress. While corneocytes of SC remain almost impermeable, \approx 95% of the transport across skin happens through skin lipid matrix. Hence, the main task in design of transdermal drug delivery formulations or personal care products remains as effective manipulation of skin lipids interface. Addition of specific chemical additives (permeation enhancers) in the formulation is the most common method followed. However, current industry practice to arrive at the most suitable additive is largely experimental involving trial and error in-vitro and in-vivo tests which obviously becomes time consuming and expensive. An in-silico model, which could mimic skin barrier function at molecular scale and help screen/design permeation enhancers, will be of immense value for both pharma and cosmetics industries. We have developed a computational model of SC which is able to mimic its barrier function. A multiscale modelling framework linking molecular (micro) to continuum (macro) scale is employed to study molecular transport across the skin. A bilayer mixture of ceramides, fatty acids and cholesterol molecule is simulated using molecular dynamics (MD) simulations. The transport properties obtained from MD simulations are further incorporated in computational fluid dynamics to compute the macroscopic transport of molecules across the skin. In this talk, I shall briefly describe this model as well illustrate its utility in product design and testing by taking a few examples from pharma and cosmetics industry.

Molecular Dynamics Investigations of Anti HIV-1 Protein SAMHD1

S. Bhattacharya

IS

Indian Institute of Technology Bombay, India

HIV-1 is restricted in macrophages and certain quiescent myeloid cells due to a “Scorched Earth” dNTP starvation strategy attributed to the Sterile Alpha Motif and HD domain protein – SAMHD1. Active SAMHD1 tetramers are assembled by GTP-Mg+2-dNTP cross bridges and cleave the triphosphate groups of dNTPs at a Km of 10 uM, which is consistent with dNTP concentrations in cycling cells, but far higher than the equivalent concentration in quiescent cells. HIV-2 and SIV have evolved defences against SAMHD1, underscoring its role in restriction. The regulation of SAMHD1 by phosphorylation (at T592) and Glutathione driven redox reactions has been experimentally studied and yet not properly understood. We have used correlation network analysis along with molecular dynamics techniques to study the flow of allosteric information across the active complex. I will discuss how our MD analysis has yielded insights that complement and expand X-Ray and biochemistry based models of SAMHD1 activity.

Microscopic View of the Crowding Effects on Hydrophobic Collapse

D. Nayar

IS

Centre for Computational and Data Sciences, IIT Kharagpur, India

A living cell is a crowded milieu comprising of large-sized macromolecules, small co-solutes and ions with less free water [1,2]. Crucial biological processes involving hydrophobic collapse such as protein folding and other self-assemblies occur in this environment. However, a molecular-level understanding of these effects remains elusive. It has been widely accepted that these effects are induced due to size (steric) effects of crowders and the solvent excluded volume effects that are entirely entropic in nature [1]. This excludes the role of any direct solute-crowder or crowder-crowder attractive interactions, which however has been shown to play a crucial role recently [3]. Therefore, molecular mechanisms associated with these effects need to be further explored. We investigate the crowding effects of small (tri)peptides on collapse equilibria of a generic hydrophobic polymer. Advanced molecular dynamics simulations and statistical mechanics solvation theories are used to examine solvation thermodynamics of polymer collapse. The unresolved role of crowder intermolecular interactions is examined. Our results show that weak polymer-crowder attractions lead to strong polymer collapse only at high crowder volume fractions, involving entropic depletion of crowders from polymer surface, in accordance with the widely known depletion mechanism. Interestingly, on making the polymer-crowder attractions stronger, polymer collapses at low volume fractions and that too via preferential adsorption of the crowders on the polymer surface. Strongly interacting crowders weaken the polymer collapse at high crowder volume fractions. A transition from enthalpy-dominated to entropy-dominated polymer collapse is observed with increasing crowder concentrations. Our results provide new insights into the existing theories of crowding effects on macromolecular collapse and the talk will discuss its implications on macromolecular collapse/aggregation processes.

References

- [1] Zhou, H.-X.; Rivas, G. ; Minton, A.P. *Annu. Rev. Biophys.* 2008, 37, 375-397.
- [2] Wang, K.-H.; Chang, C.-W. *Phys. Chem. Chem. Phys.* 2015, 17, 23140–23146
- [3] Sukenik, S.; Politi, R.; Ziserman, L.; Danino, D.; Friedler, A.; Harries, D. *PloS ONE*, 2011, 6, e15608.

February 23, Sunday

The Role of Solid-Liquid Interfacial Fluctuations in the Spontaneous Motion of Droplets

U. Saxena, S. Chouksey, K. Rane

IS

Indian Institute of Technology, Gandhinagar, India

The density fluctuations at the solid-liquid interfaces have received great attention in recent decades in the context of protein folding. In this presentation, I will discuss them in the context of droplet motion on heterogeneous surfaces. I will first discuss our efforts to understand how the above fluctuations affect the variation of solid-liquid interfacial free energy with the nature of crystalline surface. We used the grand canonical Monte Carlo (GCMC) simulations and the cumulant expansions of the interfacial free energy to relate the above fluctuations to the interfacial entropy [1]. We observed that interfacial entropy is important for the motion of droplets when the temperature varies spatially, or temporally. I will also discuss a model system where the fluctuations are expected to strongly affect the motion of droplet [2]. Here, we used the molecular dynamics simulations to study the motion of droplet, and GCMC simulations to compute the interfacial free energies. I will end the talk by discussing our efforts to rationally design the solid surfaces having the desired solid-liquid interfacial fluctuations by using the principle of Maximum Entropy.

References

- [1] Chouksey, S., Rane, K. Transverse correlations near solid-liquid interface: Influence of the crystal structure of solid. *Chemical Physics* 517, 188–197 (2019).
- [2] Saxena, U., Chouksey, S., Rane, K. Spontaneous translation of nanodroplet over a heterogeneous surface due to thermal cycles: role of solid–liquid interfacial fluctuations. *Molecular Physics* 1–14 (2019). doi:10.1080/00268976.2019.1657191

Microscopic Assessment of Liquid-Liquid Extraction System in Bulk and at the Interface

A. Das^{1,3}, P. Sahu^{2,3}, Sk. Musharaf Ali^{2,3}

IS

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The recent trend toward renewal of interest in nuclear power as a futuristic energy alternative demands a close cycle spent fuel reprocessing methods with a long term vision for safe management of spent fuel. Nuclear fuel cycle is important not only to reduce the high-active solid waste but also to produce the fresh fuel for 2nd generation nuclear reactor. The extent of reprocessing depends on the efficiency of the liquid-liquid extraction processes in which the radionuclides are separated from acidic aqueous solution. The most extensively used solvent in spent fuel reprocessing is tri-n-butyl phosphate (TBP) with n-dodecane as the diluent in the PUREX process. In spite of great success, there is a demand and continuing search for an alternative of TBP, specifically for spent fuels from fast breeder reactor. Among many, tri-isoamyl phosphate (TIAP) has been considered as competitor which has similar chemical and radiological stability like TBP in presence of nitric acid. In this context, thermo-physical and dynamical properties of ligand-solvent systems are required to predict the extraction efficiency for a representative system which can be obtained either by performing experiments or by molecular dynamics (MD) simulations. In that perspective, MD simulations assist to arrive at a reasonable conclusion in minimum no of trials of experiments and thus reduce the cost as well as time. Furthermore, there is a continuing endeavour at the molecular level understanding of liquid-liquid bi-phasic extraction of metal ions due to its wide level of application from pharmaceuticals to nuclear industry. Microscopic understanding of the interface between two immiscible or partially miscible liquids of any biphasic system not only has a great interest in view of mass transfer processes but also has considerable technological values in the field of chemistry, physics and biology. Due to inherent difficulty, the understanding of molecular details at liquid–liquid interface using only experimental technique is inadequate to establish the interfacial behaviour. This is mostly due to the fluidity of the interface and its concealed environment, which restricts the experimental facts. The contribution of intrinsic thickness and broadening induced by capillary waves are responsible for total thickness but the determination of these two values are perhaps not encountered for three component system. MD simulations provide a microscopic analysis of the interfacial properties of water–organic interface. The present talk will focus on evaluation of structural, thermo-physical and dynamical properties of the liquid-liquid extraction system in bulk and at the interface using molecular dynamics simulations.

References

- [1] Sk. Musharaf Ali, A Boda, AKS Deb, P Sahu, KT Shenoy, Computational Chemistry Assisted Design and Screening of Ligand-Solvent Systems for Metal Ion Separation, *Frontiers in Computational Chemistry*: 2017, Volume: 3 3, 75, Bentham Science.
- [2] P Sahu, Sk. Musharaf Ali, KT Shenoy, *Physical Chemistry Chemical Physics*, 2016, 18, 23769-23784.
- [3] A. Das, P. Sahu, Sk. Musharaf Ali, *J. Chem. Eng. Data* 2017, 62, 22802295.
- [4] Arya Das and Sk. Musharaf Ali, *J. Chem. Phys.* 2018, 148, 074502.
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List of Posters

Presenter	Poster Title
Shubham Tiwari	Insight to the Mechanism of Nanoparticle Induced Suppression of Detergency: Theory and Simulations
Hrushikesh M. Gade	Water-mediated curvature change in graphene by single-walled carbon nanotube: A molecular dynamics study
Avula Venkata Siva Nikhil	Atomistic Modeling of Binary Ionic Liquid Mixtures
Prosun Halder	High Throughput Screening of Hypothetical Metal Organic Frameworks for Ethane-Ethylene Separation
Sonali Gore	Tuning the Adsorption Behaviour of the Material at the Molecular Scale to Get the Desired Macroscopic Behaviour by Using Statistical Mechanics and Molecular Simulations
Shakkira Erimban	Cold Adaptation of Cell Membrane of a Psychrotolerant Bacteria: Investigation using Molecular Dynamics Simulation
Vikas Dubey	Mechanism of Hydroxide Ion Transfer through Anion Exchange Membrane in Anion Exchange Membrane Fuel Cell: Investigation using Molecular Dynamics Simulation
Nirali Dhiren Desai	New Age Antimicrobial Peptides: Revealing Mode of Actions of Multifunctional AMPs Using Molecular Dynamics Study
Shivam Dubey	Role of Translational Jump-diffusion in the Breakdown of the Stokes-Einstein relation in Supercooled Water and its Binary Mixture with Glycerol
Omkar Singh	Characterization of Biological Water at Interface of Antimicrobial Peptide in Presence of Salts Solution
Arya Das	Molecular Dynamics Simulations on Interfacial Structure in Presence of Third Component
Projesh Kumar Roy	Microscopic structure and CO ₂ adsorption properties of 6FDA/BPDA-DAM polymeric membrane
Gauri Tekbahadur Thapa	Molecular Dynamics Simulation of Anti-HIV Protein SAMHD1

Presenter	Poster Title
Bharti	Melting in Two-Dimensional Gay-Berne Liquid Crystals
Manjinder Singh	A Comparative Study of Tackifying Monomers to Develop Bio-Based Pressure-Sensitive Adhesive: A Computational Approach
Ravi Kumar Reddy A	Uncovering the Molecular Mechanism of Solvent Induced Polymorphism in Crystal Nucleation from Solution
Jyoti Kuntail	Understanding the Adsorption Mechanism of Arsenous acid on Magnetite (311) Surface through Molecular Dynamics simulations
Rajneesh Kashyap	Oil Detachment from Rock Surface using Nanoparticles, Surfactant and low salinity brine: A Molecular Dynamic Study
Shivanand Kumar Veesam	Molecular Modelling of Phase Equilibrium of Gas Hydrates
Jagroop Kaur	Temperature Dependent Interaction of Soft Repulsive Wall with the Thermotropic Liquid Crystals
Krishna Jaiswal	A Functional Force Field Model for Water based on Gaussian Charges
Sanchari Bhattacharjee	Effect of Alkyl chain on The Wetting Behaviour of Aqueous Ionic Liquids: A Molecular Dynamics Study
Shubhangra Tripathi	A Temperature Accelerated Sliced Sampling study of Drug Binding/Unbinding
Sagar Dinkar Kamble	Investigation of Cholesterol Influence in Fluid Phase and Gel Phase Lipid Bilayers by Coarse-grained Molecular Dynamic Simulation
Showkat Mir	Electronic Properties of High CO_2 Capture Ability of Two-Dimensional Metal Nitrides (XN ; $X=Al, Ga, In$): A Computational Study
Amrita Goswami	Formulation and Implementation of General Topological Network Criteria for Exploring the Structures of Confined Ice

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