

Meeting Day 1: Thursday, July, 14th*Speaker setup: 8:30 - 9:00 AM***Session 1: Colloids**

09:00 - 09:10				Welcome
09:10 - 09:40	Gerd	Schroeder-Turk	Murdoch University	TBA
09:40 - 10:10	Asaph	Widmer-Cooper	University of Sydney	TBA
10:10 - 10:40	Amelia	Liu	Monash University	Local stability and local structure of colloidal glasses

*Morning Tea: 10:40 - 11:10 AM**Speaker setup***Session 2: Scattering**

11:10 - 11:40	David	Paganin	Monash University	Fokker-Planck equation for optical beams
11:40 - 12:10	Jing	Fu	Monash University	Recent advances in nanoscale tomography of soft matters
12:10 - 12:40	Andrew	Martin	RMIT University	Introducing the Pair-Angle Distribution Function: measuring multi-atom statistics of disordered materials

*Lunch: 12:40 - 14:00**Speaker setup***Session 3 A: Simulations - Principles**

14:00 - 14:30	Ellie	Hajizadeh	University of Melbourne	ML-based optimisation for accelerated and targeted multiscale soft matter design
14:30 - 15:00	David	Huang	University of Adelaide	Systematic coarse-graining of molecular simulation models with anisotropic particles
15:00 - 15:30	Luca	Maffioli	Swinburne University of Technology	Measuring the response of highly confined fluids in a Couette flow: the TTCF formalism

*Afternoon Tea: 15:30 - 16:00**Speaker setup***Session 3 B: Simulations - Principles**

16:00 - 16:30	Stephen	Sanderson	University of Queensland	Machine learning a Time-Local Fluctuation Theorem for non-equilibrium steady-states
16:30 - 17:00	Tim	Duignan	University of Queensland	Accurate first principles simulation of salt water using deep learning.

*Informal Dinner***Meeting Day 2: Friday, July, 15th***Speaker setup: 8:30 - 9:00 AM***Session 5: Active Matter**

09:00 - 09:30	Prabhakar	Ranganathan	Monash University	Cluster and conquer: invasion of a soft substrate by colonies of rod-shaped cells
09:30 - 10:00	Rahil	Valani	University of Adelaide	Generating active matter from strange attractors
10:00 - 10:30	Shibu	Saw	Other (add more details on next page)	Role of the configurational temperature in active-matter models

*Morning Tea: 10:30 - 11:00**Speaker setup***Session 6 A: Simulations - Applications**

11:00 - 11:30	Shern	Tee		When Is A Supercapacitor Like A Protein?
11:30 - 12:00	Mikhail	Suyetin	Other (add more details on next page)	Proof of concept: Molecular Dynamics study of memory nanodevice based on Zn-MOF-74.
12:00 - 12:30	Dominic	Robe	Monash University	Linear viscoelasticity of associating star polymers

*Lunch: 12:30 - 14:00**Speaker setup***Session 6 B: Simulations - Applications**

14:00 - 14:20	Patrick	Adams	RMIT University	Using the pair-angle distribution function to analyse protein structure.
14:20 - 14:40	Isaac	Pincus	Monash University	Modelling dilute solutions of semiflexible polymers with excluded-volume and hydrodynamic interactions
14:40 - 15:00	Silpa	Mariya	Monash University	Are soft dendrimers dynamically equivalent to hard spheres?
15:00 - 15:20	Michael	Rinaudo	University of Sydney	Simulating nanoparticle superlattices with Molecular Dynamics

*Afternoon Tea: 15:20 - 15:40**Speaker setup***Session 6 C: Simulations - Applications**

15:40 - 16:00	Sobin	Alosious	Swinburne University of Technology	The effect of electrostatic interactions on Kapitza resistance in nano-confined water.
16:00 - 16:20	Christian	Zuluaga-Bedoya	University of Queensland	Entry resistance in flexible zeolitic nanosheets
16:20 - 16:40	Kyle	Stevens	University of Newcastle	Lennard-Jones potential and continuum modelling for heterogeneous molecules
16:40 - 17:00	Marltan	Wilson	University of Adelaide	General high-dimensional neural networks for anisotropic coarse-grained molecular simulation models
17:00 - 17:20				Closing