

Leverhulme Research Centre for Functional Materials Design Materials Discovery Symposium, September 2022

Location: Anfield Stadium

Registration: Main Stand reception, Anfield Stadium

Oral Presentations: The Dugout, 2nd Floor, Main Stand

PROGRAMME: DAY ONE – Wednesday 21st September 2022

09.00 – 10.00	Registration and Coffee
10.00 – 10.15	Welcome & Intro to the LRC Professor Andrew I Cooper <i>Director, Leverhulme Research Centre for Functional Materials Design</i>
Session 1:	Chaired by: TBC
10.15 - 10.45	Translating the enormous promise of complex oxides to working devices Professor Judith Driscoll <i>University of Cambridge, UK</i>
10.45 – 11.15	Data-driven Materials Discovery Professor Jacqui Cole <i>University of Cambridge, UK</i>
11.15 – 11.45	Coffee break
11.45 – 12.15	Coordinating and reconfiguring swarms of objects at extreme dimensions Professor Sándor Fekete <i>Braunschweig University of Technology, Germany</i>
12.15 – 12.45	Designing a Programmable Material Made of Autonomous Robots Professor Julien Bourgeois <i>Université Bourgogne Franche-Comté, FEMTO-ST Institute, France</i>
12.45 – 14.00	Lunch and Poster Session
Session 2:	Chaired by: TBC
14.00 – 14.30	Structure design and band gap engineering from mixed anion building blocks Dr. Houria Kabbour <i>Unité de Catalyse et Chimie du Solide, France</i>
14.30 – 15.00	Employing Chemical Heuristics in Computational Materials Design of Functional Materials Professor David Scanlon <i>University College London, UK</i>
15.00 – 15.30	Scheduling Theory and Its Application to Chemical Production Professor Prudence Wong <i>University of Liverpool, UK</i>

15.30 – 16.00	Coffee Break
16.00 – 16.30	Materials that compute: 21 Molecular algorithms using reprogrammable DNA self-assembly Professor Damien Woods <i>Maynooth University, Ireland</i>
16.30 – 17.00	Self-Optimising Industry 4.0 Chemical Reaction Platforms for Multi-objective and Multistep Process Development Professor Richard Bourne <i>University of Leeds, UK</i>

PROGRAMME: DAY TWO – Thursday 22nd September 2022

Session 3:		Chaired by: TBC
09.00 – 09.30	Coffee Breakfast	
09.30 - 10.00	Exploring Catalyst-Support Interfaces with Faceted Perovskite Nanoparticles Professor Kenneth R. Poeppelmeier <i>Northwestern University, USA</i>	
10.00 – 10.30	Defects and Disorder in Electronic Oxides Professor Elizabeth C. Dickey <i>Carnegie Mellon University, USA</i>	
10.30 – 11.00	Getting into shape – Precision polymer nanoparticles Professor Rachel O'Reilly <i>University of Birmingham, UK</i>	
11.00 – 11.30	Coffee Break	
11.30 – 12.00	Designing polymers for energy in the digital age Dr. Brett Helms <i>Lawrence Berkeley National Laboratory, USA</i>	
12.00 – 12.30	Designing High Performing Semiconducting Polymers Professor Ian McCulloch <i>University of Oxford, UK</i>	
12.30 – 14.00	Lunch & poster session	
Session 4:		Chaired by: TBC
14.00 – 14:30	Accelerating the discovery and evaluation of sorbents for sustainable gas separations Professor Camille Petit <i>Imperial College London, United Kingdom</i>	
14.30 – 15.00	Flexible Automation: A New Approach to Meet Evolving Challenges Dr. Jason Hein <i>University of British Columbia, USA</i>	
15.00 – 15.30	New Ways in Automating Research and Development Laboratories Professor Kerstin Thurow <i>University of Rostock, Germany</i>	

15.30 – 16.00	Coffee Break
16.00 – 16.30	It's a machine's world: How automation and machine learning reshapes polymer chemistry Professor Tanja Junkers <i>Monash University, Australia</i>
16.30 – 17.00	Billions upon billions of molecules Professor Alan Aspru Guzik <i>University of Toronto, Canada</i>
17.00 – 18:45	Anfield Stadium Tours (Pre-booked) & Networking Drinks
19:00 – 22:00	Conference dinner for seating at 19:00 in The Chemistry Suite, 3rd Floor, Main Stand, Anfield Stadium

PROGRAMME: DAY THREE – Friday 23rd September 2018

Session 5:		Chaired by: TBC
09.00 – 09.30	Coffee Breakfast	
09.30 - 10.00	Superconductivity in infinite-layer nickelates Professor Harold Y. Hwang <i>Stanford University, USA</i>	
10.00 – 10.30	Meta-learning adaptive deep kernel Gaussian processes for molecular property prediction Professor José Miguel Hernández-Lobato <i>University of Cambridge, UK</i>	
10.30 – 11.00	Towards molecular design in allosteric processes through unsupervised, multiscale learning on atomistic graphs Professor Sophia Yaliraki <i>Imperial College London, UK</i>	
11.00 – 11.30	Coffee Break	
11.30 – 12.00	Complementary Artificial Intelligence for Sustained Innovation Professor James Evans <i>University of Chicago, USA</i>	
12.00 – 12.30	Advancement of Zintl Phases from Curiosities to Thermoelectric Materials Professor Susan Kauzlarich <i>University of California, Davis Campus, USA</i>	
12.30 – 14.00	Lunch & poster session	
Session 6:		Chaired by: TBC
14.00 – 14:30	Voltage & Current Controlled Nanomagnetism for Memory and Logic Professor Lucas Caretta <i>Brown University, USA</i>	
14.30 – 15.00	Topology and Chirality Professor Claudia Felser <i>Max Planck Institute Chemical Physics of Solids, Germany</i>	

15.00 – 15.30	Excitement in Materials Research: from material design to social implementation Professor Hideo Hosono <i>Tokyo Institute of Technology, Japan</i>
15.30 – 16.00	Poster Prize Presentation & Close