

# Molecular excited states theory and experiment

14-16 September 2026 | Cambridge, UK



## Faraday Discussions

**Monday 14 September 2026**

11:30	Registration and lunch
12:30	<b>Welcome and introductions</b> Michael Bearpark, <i>Imperial College London, UK</i> Javier Segarra Martí, <i>Universitat de València, Spain</i> <i>Co-chairs of Scientific Committee</i>
12:40	<b>Outline of Discussion format</b> <i>Royal Society of Chemistry Publishing Editors</i>
12:45	<b>Introductory lecture – Spiers memorial lecture</b> (Session chair: ) Michael A Robb <i>Imperial College London, UK</i>
	<b>Session 1: The excited state electronic structure problem: new methods and computer architectures</b> (Session chair: )
13:45	<b>Title TBC</b> Katharina Boguslawski <i>Nicolaus Copernicus University, Poland</i>
13:50	<b>Symmetry-projected generalised normal-ordered coupled-cluster theory for excited states</b> Bang Huynh <i>University of Oxford, UK</i>
13:55	<b>Multireference perturbation theory (MRPT) geometry optimizations with scalar-relativistic effects: Effective core potential (ECP) and spin-free X2C Hamiltonian</b> Jae Woo Park <i>Chungbuk National University, South Korea</i>
14:00	<b>Ground- and Excited-State Description via a Hybrid Quantum-Classical Method: Case Study of Three States</b> Benjamin Lasorne <i>Univ Montpellier, France</i>
14:05	Discussion
15:45	Refreshments
16:15	<b>Title TBC</b> Ksenia Bravaya <i>University of Boston, USA</i>
16:20	<b>Benchmarking electronic-structure methods for nonadiabatic dynamics in a propeller-shaped molecular rotor: Insights into conical intersections</b> Rachel Crespo-Otero <i>University College London, UK</i>
16:25	<b>Modular Integration of MRSF-TDDFT, NAMD, and QM/MM for Excited-State Dynamics on the OpenQP Platform</b> Cheol Ho Choi <i>Kyungpook National University, South Korea</i>
16:30	Discussion
17:45	Lightning presentations (by invitation of the Scientific Committee)
18:15	Poster session and wine reception
19:30	Close of sessions

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Tuesday 15 September 2026

	<b>Session 2: Non-adiabatic and ultrafast dynamics: models and measurements</b> (Session chair: )
09:00	<b>Title TBC</b> Nanna List <i>KTH Royal Institute of Technology, Sweden and University of Birmingham, UK</i>
09:05	<b>Time-resolved extreme ultraviolet photoelectron spectroscopy of 4-bromophenolate anions in a liquid jet</b> Daniel Neumark <i>University of California, Berkeley, USA</i>
09:10	<b>Modeling the coupled electron-nuclear dynamics following electronic coherences induced by attosecond and few-femtosecond light pulses</b> Fernando Martin <i>IMDEA Nanoscience and Universidad Autonoma de Madrid, Spain</i>
09:15	<b>Disentangling Ultrafast Reaction Pathways of Furan-derivative with Time-Resolved X-ray Absorption Spectroscopy</b> Vesna Erić <i>Max Planck Institute for Polymer Research, Germany</i>
09:20	Discussion
11:00	Refreshments
11:30	<b>Ultrafast electron diffraction imaging of chemical substitution effects on nonadiabatic nuclear dynamics at conical intersections</b> Kasra Amini <i>Max-Born-Institute, Germany</i>
11:35	<b>The Information Content in Ultrafast Observables: Spectroscopy and Scattering in Excited Norbornadiene</b> Adam Kirrander <i>University of Oxford, UK</i>
11:40	<b>Modeling correlation effects in photoionization spectroscopy and photoelectron circular dichroism of ground- and excited-state molecular systems.</b> Sonia Coriani <i>Technical University of Denmark, Denmark</i>
11:45	Discussion
13:00	Lunch
14:00	<b>Solvent and field effects on photoinduced ring-opening: dynamics of furan as a case study</b> Lea Ibele <i>ICR, Aix Marseille University, CNRS, France</i>
14:05	<b>Direct dynamics simulation of photochemistry without the Born-Oppenheimer approximation</b> Ryan MacDonell <i>Dalhousie University, Canada</i>
14:10	<b>Modelling photoelectron spectra in solution: deoxyadenosine as a test case</b> Roberto Improta <i>Consiglio Nazionale delle Ricerche, Italy</i>
14:15	Discussion
15:30	Refreshments
	<b>Session 3: Developing links with large-scale experiments: computing observables and interpreting measurements in new facilities</b> (Session chair: )
16:00	<b>Real-space imaging of the valence electron density during a photochemical reaction</b> Thomas Wolf

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	<i>LCLS and Stanford PULSE Institute, SLAC National Accelerator Laboratory, USA</i>
16:05	<b>Understanding Resonant Inelastic X-ray Scattering Experiments of Diazines via Quantum Dynamics Simulation</b> Antonia Freibert <i>Technical University of Munich, Germany</i>
16:10	<b>Time-Resolved Photoemission Spectroscopy in the X-ray Domain: Opportunities and Challenges with Mixed-Reference Spin-Flip TDDFT</b> Petr Slavicek <i>University of Chemistry and Technology, Prague, Czech Republic</i>
16:15	<b>Excited state dynamics of haem proteins</b> Majed Chergui <i>Ecole Polytechnique Fédérale de Lausanne, Switzerland</i>
16:20	Discussion
18:00	Close of sessions
18:30	Pre-dinner drinks
19:00	Conference dinner

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## Faraday Discussions

Wednesday 16 September 2026

	<b>Session 3 continued: Developing links with large-scale experiments: computing observables and interpreting measurements in new facilities</b> (Session chair: )
09:00	<b>Title TBC</b> Mary Matthews <i>Imperial College London, UK</i>
09:05	<b>Measurement of Electron Correlations with X-Ray Scattering</b> Peter M. Weber <b>Brown University, USA</b>
09:10	<b>Time-Resolved Core-Level Photoelectron Spectroscopy of Glycine Fragmentation: A Case Study on the Capabilities and Limitations of Site-Selective Probes at XFELs</b> Andre Al haddad <i>Paul Scherrer Institute, Switzerland</i>
09:15	Discussion
10:30	Refreshments
	<b>Session 4: AI and data-driven approaches in molecular excited states</b> (Session chair: )
11:00	<b>Active delta-learning for surface hopping: efficient fitting of potential energy surfaces</b> Pavlo O. Dral <i>Xiamen University, China; NCU in Toruń, Poland; Aitomistic, China</i>
11:05	<b>Beyond Minimum Energy Conical Intersections: A Data-Driven Reconstruction of the Accessible Intersection Seam</b> Elisa Pieri <i>University of North Carolina at Chapel Hill, USA</i>
11:10	<b>Machine Learning and Multireference Configuration Interaction for High-Level Nonadiabatic Dynamics Simulation of Hexatriene</b> Hans Lischka <i>Texas Tech University, USA</i>
11:15	Discussion
12:30	<b>Concluding remarks lecture</b> (Session chair: ) Graham A. Worth <i>University College London, UK</i>
13:00	<b>Acknowledgements</b>
13:15	<b>Close of meeting and lunch</b>

Please note that this is a draft programme and timings may change.