

## Poster Session, Monday 04 May 2026, 20:00 CEST

Katrin Bickel	<b>Solvent effects on the reaction entropy of an outer sphere electron transfer process</b>
Tim Bolter	<b>Dynamics of Au electrode surfaces probed by electro catalytic reactions</b>
Marine Bossert	<b>Capillary rise of electrolyte-alkane mixtures between gold plates: wetting transition or electrocapillary effect?</b>
Steffen Braun	<b>Comparing microcalorimetric and temperature-variation approaches for determining the reaction entropy of electrochemical reactions</b>
Yuchen Cai	<b>Confinement effect of cation on proton transport</b>
Alper Tunga Celebi	<b>A thermodynamic origin of ion-specific adsorption at charged interfaces</b>
Hao Chen	<b>Single-molecule detection of small molecules enabled by charge regulation at the solid-liquid interface of silicon nitride solid-state nanopores.</b>
Junxiang Chen	<b>DP<math>\chi</math>: A Constant-Potential Machine-Learning Potential for Large-Scale Electrocatalysis</b>
Lucas de Kam	<b>Poisson-Boltzmann theory in practice</b>
Qiuhan Deng	<b>Electronic and Ionic Charge Redistribution in Electric Double Layers under Curvature and Steric Constraints</b>
Laurin Derr	<b>Entropy Profiling of Hard Carbon/Na Metal Batteries with Stepwise Temperature Changes: Deciphering the Cell Voltage Response</b>
Hannah Dickinson	<b>Impact of Electrolyte pH on the Morphology of Silver Electrodeposition Studied Using in-situ Transmission Electron Microscopy</b>

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- Martin Durner      **Scanning Impedance Microscopy under Potential Control to Probe Local Electrochemical Properties**
- Jordy Eggebeen      **Using electrochemical impedance spectroscopy to deconvolute (pseudo) capacitance from metal oxide interfaces**
- Mikhail Elkhimov      **Influence of the electronics and cell resistances on electrochemical noise measurements**
- Victoria Gamez      **Anion-Water and Hydroxyl Adsorption at the Pt(111) Interface: Influence of pH and Electrolyte Composition**
- Antonia Gerstenberg      **The effect of bias and temperature on interfacial water at pristine Pt(111) electrodes: Insights from atomistic MLIP simulations**
- Greta Grossman      **Surface Electrochemistry of the Au(111)/acetonitrile Interface: Formation of a Solvent Related Adsorbed Layer and its Implications for the Electrochemical Double Layer**
- Siyuan Han      **Machine-Learning-Accelerated First-Principles Simulations Reveal Ion-Correlated Structures at Concentrated Electrochemical Interfaces**
- Zetong Jia      **Long-Range Ordered Pt Facets Govern Methane Electro-oxidation via Interfacial Solvent Dynamics**
- Rick Kort      **Linking Interfacial Water Structure to AFM Forces under Electrochemical Control**
- Justus Leist      **Probing Surface Structure and Vibrational Stark Effects in Ni Anodes with SERS during Alkaline OER**
- Chenkun Li      **Constant-Potential Kohn-Sham-Poisson-Boltzmann Theory for Electrified Metal-Solution Interfaces**
- Antony Litovolis      **Cation Effects On Solvated Electron Generation**

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Julius Lonnes	<b>Disentangling Temperature Effects in Proton Adsorption Barriers at Electrified Pt(111): Roles of Water Structure, Potential, and Vibrational Free Energies</b>
Alessandro Mangiameli	<b>Benchmarking QM/RISM method for the Au-Electrolyte Interface</b>
Ricardo Alonso Martinez Hincapie	<b>Double Layer Capacity at Pt(111) Electrode in Near-Neutral Electrolytes</b>
Daniel Metzger	<b>Concentration dependent entropy determination for sodium deposition</b>
Marcus Frahm Nygaard	<b>Electron Redistribution Between Coadsorbed Species at the Electrochemical Interface</b>
Emmanouil Pervolarakis	<b>The Computational Cation Electrode: A Case Study on CO<sub>2</sub>RR</b>
Jorik Schaap	<b>The influence of the Electric-Double layer and Cathodic Bias on the Ultrafast Dynamics of Silver Nanoparticles</b>
Barbara Sumic	<b>First Steps Matter Most: Free Energy Barriers of the Volmer Step in Hydrogen Evolution at Pt(111)</b>
Yu Chen Sun	<b>Probing Nafion structure at Electrified Pt Interfaces via in-situ Polarization-Modulated Infrared Spectroscopy</b>
Rose Marie Tom Petersen	<b>Machine-Learning Assisted Determination of Reaction Paths for Electrocatalytic Reactions</b>
Harsh Wadhvani	<b>3D porous model systems for electrocatalysis</b>
Haowei Wang	<b>Raman Scattering Measurement of Single Nanoparticle Collision</b>
Xiwei Wang	<b>Density-Potential Functional Theoretic (DPFT) Schemes of Modeling Reactive Solid-Liquid Interfaces</b>

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- Johanna Weidelt      **Towards a DFT-based Workflow to Simulate Competitive Adsorption of Electrolyte Species at a Pt Electrode**
- Lulu Zhang            **Correlating electric double layer structure and electrochemical plasmonics of nanoparticles: a theoretical study**
- Zengming Zhang      **Potential of Zero Stress of Solid-Liquid Interfaces**
- Jia-Xin Zhu            **Reactive electrical double layers under confinement: insights from atomistic to continuum**