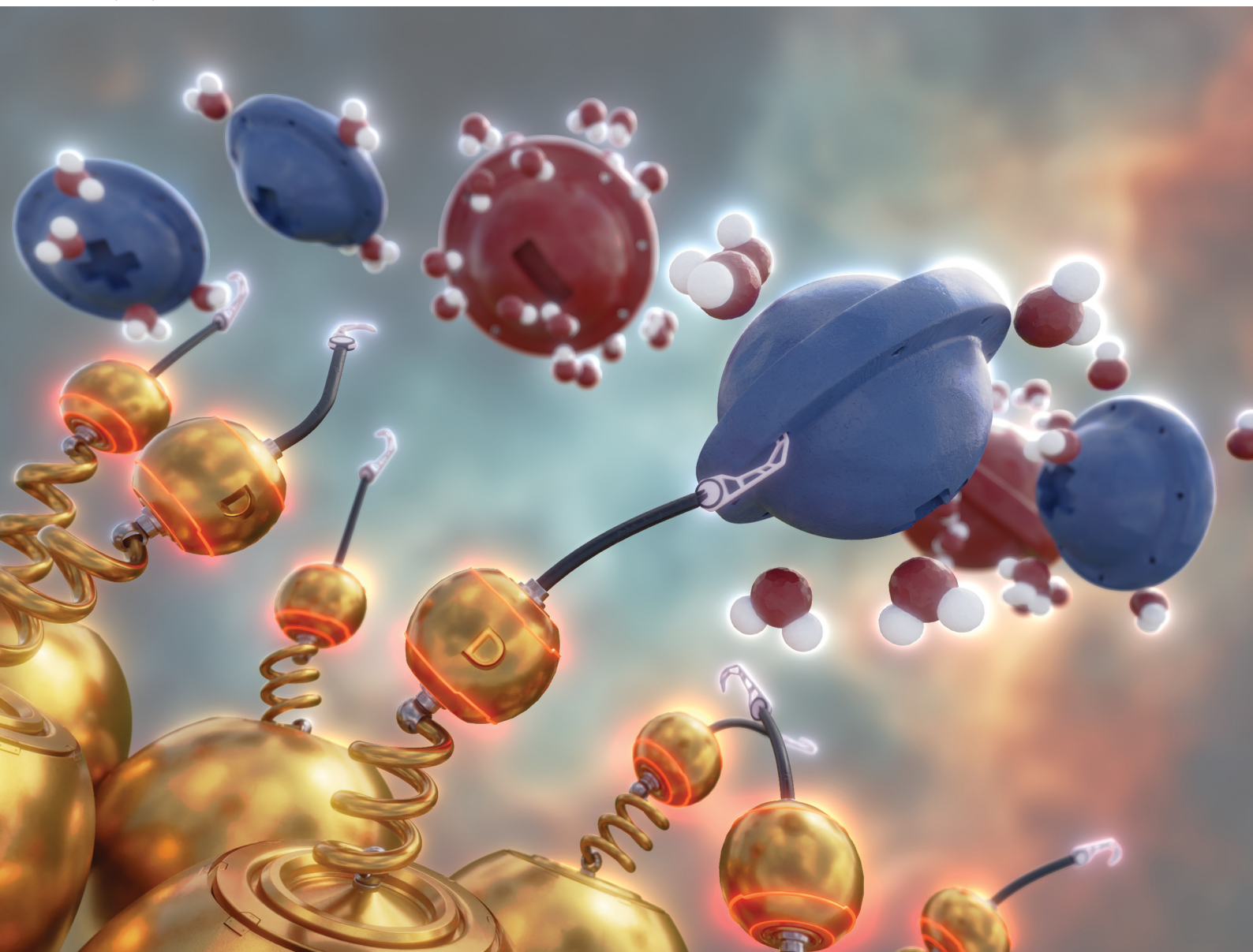


# PCCP

Physical Chemistry Chemical Physics

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ISSN 1463-9076

**PAPER**

Samuel Ntim and Marialore Sulpizi  
Molecular dynamics simulations of electrified interfaces  
including the metal polarisation