

Chemistry and Material Science on the Computer: Wavefunctions, Orbitals, and Electron Densities in Spectroscopy, Catalysis and Synthesis

Görling Group

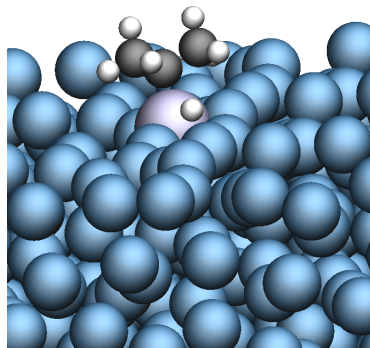
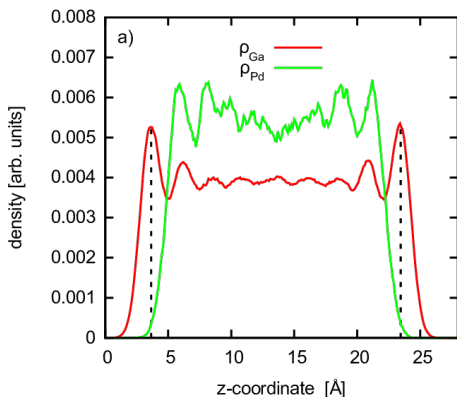
Chair of Theoretical Chemistry

Development and application of quantum chemical methods
for investigation of
molecules, clusters, surfaces, and solids
with respect to

- energetics and structure
- reactivity (catalysis)
- electronic structure (orbitals, band structures, STM)
- spectroscopy (UV/Vis, IR, NMR, non-linear optical properties)

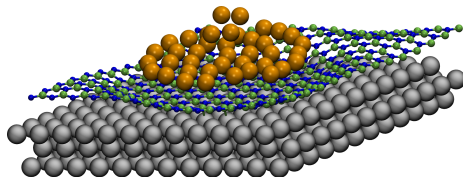
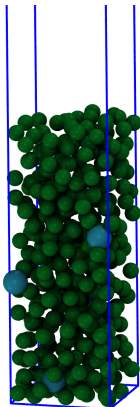
Liquid Pd/Ga or Rh/Ga mixtures as catalyst for hydrocarbon de-hydrogenation

Methodology: Machine-learned force field trained on-the-fly by DFT data



Within DFG collaborative research center 1452 "Catalysis at liquid interfaces"

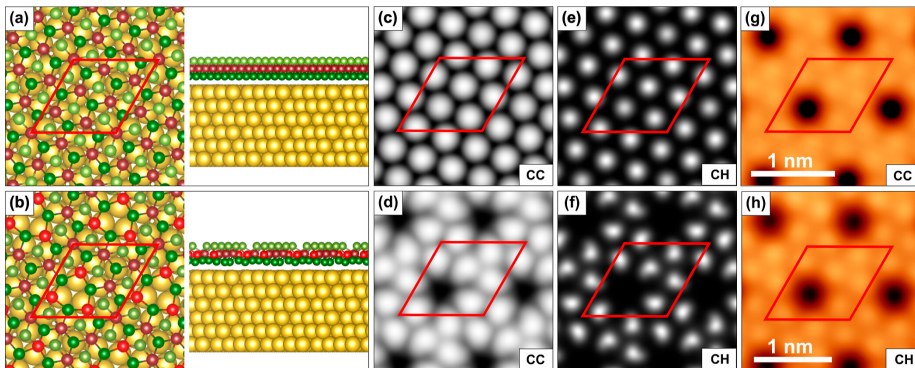
Influence of dynamic processes



Structure formation in komplex systems

Calculation of diffusion coefficient to
interpret results from neutron diffraction

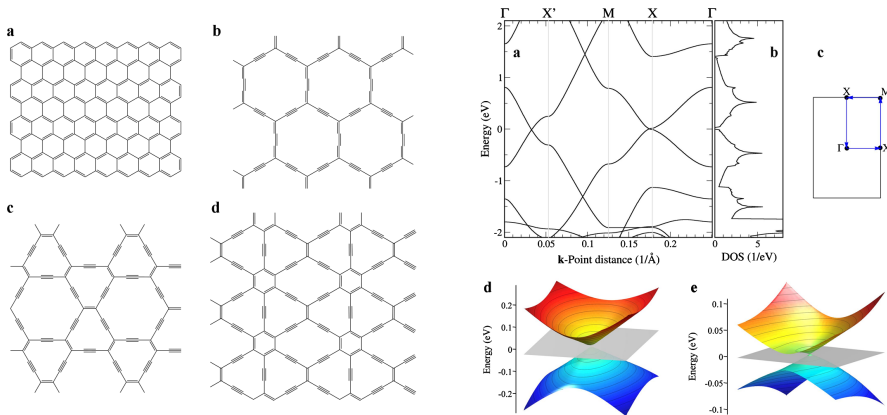
Intrinsically patterned two-dimensional transition halides



Structure of pristine and patterned FeBr₂ films on Au(111)

Collaboration with S. Maier

Alternatives to graphene with equally amazing electronic properties



Within DFG collaborative research center 953 "Synthetic carbon allotropes"

- 1  Organic molecules and ionic liquids on metal surfaces Julien Steffen

- 2  Liquid metal catalysis (ab-initio dynamics simulations) Andreas Mölkner

- 3  Novel 2D materials Neeta Bisht, Christian Reiß

- 4  Test of novel electronic structure methods Steffen Fauser