

# NIC Symposium 2010

**24 – 25 February 2010**

**Forschungszentrum Jülich, Germany**

## **Programme**

**Venue: Auditorium (Großer Hörsaal) at Forschungszentrum Jülich, Jülich, Germany**

### **Wednesday, 24 February 2010**

- 8.30        Transfer from Jülich
- 9.00        Registration
- 9.30        Welcome Address by S. Schmidt, Board of Directors, Forschungszentrum Jülich
- 9.45        Th. Lippert, Forschungszentrum Jülich  
Recent Developments in Supercomputing
- 10.30       Coffee
- 11.00       Th. Kaczmarek, Universität Köln  
Protoplanetary Discs in Young Dense Clusters
- 11.45       G. Herdoiza, NIC, DESY Zeuthen  
QCD Simulations with Light, Strange and Charm Dynamical Quark Flavours
- 12.30       Lunch
- 14.00       H. Elbern, Universität Köln  
IMACCO: Virtual Institute for Inverse Modelling of the Atmospheric Chemical  
Composition
- 14.45       N. Peters, RWTH Aachen  
Geometrical Properties of Small Scale Turbulence
- 15.30       Coffee
- 16.00       P. Gibbon, Forschungszentrum Jülich  
Plasma Simulation with Parallel Kinetic Particle Codes
- 16.45       B. Trieu, Forschungszentrum Jülich  
Threshold Determination for Reliable Quantum Error Correction
- 17.30       Poster Session and Reception
- 19.00       Transfer to Jülich

## Thursday, 25 February 2010

- 8.30 Transfer from Jülich
- 9.00 S. Schmitt, Technische Universität Dortmund  
Quantum Transport Through Strongly Interacting Nanodevices
- 9.45 St. Blügel, Forschungszentrum Jülich  
Van der Waals Interaction of Molecules on Surfaces from First Principles
- 10.30 Coffee
- 11.00 M. Magiera, Universität Duisburg-Essen  
Simulation of Magnetic Friction
- 11.45 K. Daoulas, Georg-August-Universität Göttingen  
Studying Directed Assembly of Diblock Copolymers on Patterned Substrates by  
Soft, Coarse-Grained Models
- 12.30 Lunch
- 14.00 M. Brehm, Universität Leipzig  
Ionic Liquids from *ab initio* Molecular Dynamics on Supercomputers
- 14.45 G. La Penna, Institute for Chemistry of Organo-Metallic Compounds (ICCOM)  
The Inhibition of Hydrogenase by Dioxygen
- 15.30 Coffee
- 16.00 M. Walter, Universität Freiburg  
Determination of Structure and Electronic Properties of Free, Supported and Ligand  
Protected Metal Clusters by Density Functional Theory
- 16.45 M. Fleck, Forschungszentrum Jülich  
Pattern Formation in Complex Systems
- 17.30 Transfer to Jülich