

INVITATION TO SYMPOSIUM

The DFG Research Unit 618 is pleased to invite you to the international symposium

Aggregation of Small Molecules – from Dimers to Crystals

The symposium will discuss recent work of internationally leading scientists in the field of non-covalent interactions, and present the scientific achievements of the Research Unit 618 over the past six years.

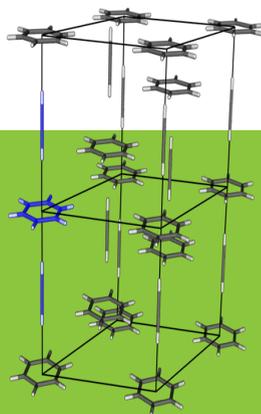
The Symposium will take place at the Ruhr-University of Bochum from **May 29th to 31st 2012**.

Guests are highly welcome! There are no conference fees, but please register at www.rub.de/for618/ before April 15th 2012.

DFG RESEARCH UNIT 618

The aggregation of molecules via non-covalent interactions to form larger structures is a process of fundamental importance to many aspects of chemistry. Although this has been recognized since many years, only now the experimental and theoretical tools are available to gain the detailed insight necessary for an understanding of aggregation processes at the molecular level.

Researchers at the Universities of Bochum, Duisburg-Essen, and Düsseldorf combined their expertise in the fields of molecular beam and low temperature spectroscopy, synthesis, crystal engineering, electronic structure and ab initio simulation techniques, and founded the Research Unit 618 in January 2006. This project is funded by the Deutsche Forschungsgemeinschaft from 2006 to 2012.



SCIENTIFIC PROGRAM

INVITED SPEAKERS

- Roland Boese - University of Duisburg-Essen
Crystalline Organic Hydrates - Frozen Stages of the Dissolution Process?
- Gautam Desiraju - Indian Institute of Science, Bangalore
The Structural Landscape in Crystal Engineering
- Nikos Doltsinis - University of Münster
Simulating aggregation from first principles
- Wolfgang Domcke - Technical University of Munich
Ultrafast Nonadiabatic Photochemistry of Hydrogen Bonds in Organic and Biological Chromophors
- Gary E. Douberly - University of Georgia, USA
Radical Containing Clusters in Helium Nanodroplets
- Stefan Grimme - University of Bonn
Dispersion Corrected Density Functional Theory
- Martina Havenith-Newen - Ruhr-Universität Bochum
Rock and Roll at 0.37 K
- Pavel Hobza - Academy of Sciences of the Czech Republic
Noncovalent Interactions: QM and MM approaches
- Christopher Hunter - University of Sheffield, UK
The Anatomy of Complex Recognition Interfaces
- Georg Jansen - University of Duisburg-Essen
Properties of Small Molecular Aggregates from Analytical Model Potentials obtained through Quantum Chemistry
- Mark Johnson - Yale University, USA
Capturing Reaction Intermediates with Cryogenic Ion Spectroscopy
- Manfred Kappes - Karlsruhe Institute of Technology
The periodic table at 55 (and some structures of other atomic cluster sizes)
- Karl Kleinermanns - University of Duesseldorf
Isomer Selective Vibronic Spectroscopy of Benzene-Acetylene Clusters – Towards a Better Understanding of Seed Crystal Formation

- Dominik Marx - Ruhr-Universität Bochum
Aggregation-Induced Chemical Reaction: HCl-Water Aggregates in the Gas Phase and in Superfluid Helium
- Klaus Merz - Ruhr-Universität Bochum
The Importance of Deuterium and Fluorine-Substituents on Molecular Aggregation Processes
- Karina Morgenstern - Ruhr-Universität Bochum
Aggregation of molecules on surface: From coverage to chirality dependence
- Wolfram Sander - Ruhr-Universität Bochum
Aggregation and Solvation of Radicals and other Reactive Intermediates
- Friedrich Temps - University of Kiel
Noncovalent Interactions in the Ultrafast Dynamics of Electronic Excited DNA Building Blocks

GENERAL INFORMATION

LOCATION

The Symposium will take place at **RUB Conference Center**, floor 04, room 1. The location is marked on the enclosed map as „Veranstaltungszentrum“. For parking please use Car park P9.

PRELIMINARY PROGRAM

May 29	18:00 - 21:00	Arrival, Reception, Evening Lectures
May 30	9:00-12:00	Lectures
	12:00-13:00	Lunch
	13:00 - 18:30	Lectures
	18:30	Dinner/ Poster Session
May 31	9:00 - 12:00	Lectures
	12:00	Closing Session/ Lunch

