

Conference on Modelling, Analysis and Simulation of Molecular Systems

PLENARY SPEAKERS

Eric Cancès (École des Ponts)

Benedict Leimkuhler (Edinburgh)

LOCAL ORGANIZER

Benjamin Stamm (RWTH Aachen)

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REGISTRATION

www.moansi.wixsite.com/gamm

CONFIRMED SPEAKERS

Mercedes Alfonso-Prieto (FZ Jlich)

Markus Bachmayr (University of Bonn)

Peter Benner (MPI Magdeburg)

Eric Cancs (Ecole des Ponts ParisTech)

Gabriele Ciaramella (University of Konstanz)

Sandra Dpking (FU Berlin)

Heinz-Jrgen Flad (TU Munich)

Gero Friesecke (TU Munich)

Paolo Gatto (RWTH Aachen University)

Jan Hamaekers (Fraunhofer SCAI)

Carsten Hartmann (BTU Cottbus-Senftenberg)

Felix Henneke (FU Berlin)

Caroline Lasser, TU Munich

Laura Lauerbach (University of Wrzburg)

Ben Leimkuhler (University of Edinburgh)

Christian Lubich (University of Tbingen)

Wenping Lyu (RWTH Aachen University)

GENERAL INFORMATION

This conference is the first annual meeting of the GAMM (Gesellschaft für Angewandte Mathematik und Mechanik) activity group on *Modelling, Analysis and Simulation of Molecular Systems* (MOANSI). The aim of the MOANSI activity group is to promote mathematical aspects of modelling and numerical computations in theoretical chemistry, physics, materials science and computational biology. We hope that this conference will provide an opportunity for researchers in these fields to share ideas and discuss the challenges and problems that they are facing.

DATES

October 12-13, 2017.

VENUE

Seminar Room 008

Pontdriesch 14-16

52062 Aachen.