

Conference on Modelling, Analysis and Simulation of Molecular Systems

PLENARY SPEAKERS

Eric Cancès (École des Ponts)
Benedict Leimkuhler (Edinburgh)

ORGANIZING COMMITTEE

Gero Friesecke (TU Munich)
Reinhold Schneider (TU Berlin)
Benjamin Stamm (RWTH Aachen)

REGISTRATION

www.moansi.wixsite.com/gamm

CONFIRMED SPEAKERS

Gabriele Ciaramella (University of Konstanz)
Konstantin Fackeldey (TU Berlin)
Heinz-Jürgen Flad (TU Munich)
Gero Friesecke (TU Munich)
Paolo Gatto (RWTH Aachen)
Jan Hamaekers (Fraunhofer SCAI)
Carsten Hartmann (BTU Cottbus-Senftenberg)
Venera Khoromskaia (MPI Leipzig)
Caroline Lasser (TU Munich)
Christian Lubich (University of Tübingen)
Sebastian Matera (FU Berlin)
Max Pfeffer (TU Berlin)
Giulia Rossetti (FZ Jülich)
Anja Schlömerkemper (University of Würzburg)
Bernd Schmidt (University of Augsburg)
Johannes Zimmer (Bath 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.