Curriculum Vitae

Personal Information

DATE OF BIRTH: Wiesbaden, Germany | April 1st, 1992

EMAIL: jung.gerhard@umontpellier.fr

HOMEPAGE: gerhardjung.github.io

Main Areas of Research

Computer simulations in soft matter physics, non-Markovian dynamics, dynamic coarse-graining, glass transition, confinement, rheology, nonequilibrium dynamics

Academic Experience

| 10/2021 - Present | CNRS, MONTPELLIER (FRANCE) Postdoctoral researcher Advisor: Prof. Ludovic Berthier, Prof. Giulio Biroli Area of study: Glass transition, machine learning, amorphous defects |
|-------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| 06/2021 | University of Kyoto (Japan) JSPS Fellow |
| 10/2021 | Advisor: Prof. Ryoichi Yamamoto Area of study: Active particles in viscoelastic media |
| 03/2019 | University of Innsbruck (Austria) Postdoctoral researcher |
| 05/2021 | Advisor: Prof. Thomas Franosch Area of study: Glass transition, crystallization, confined geometry |
| 05/2018 | Durham University (UK) Visiting researcher |
| 10/2018 | Advisor: Prof. Suzanne Fielding Area of study: Soft glassy materials, yielding transition, rheology |
| 10/2014 | UNIVERSITY OF MAINZ (GERMANY) Doctoral and postdoctoral researcher |
| 02/2019 | Advisor: Prof. Friederike Schmid Area of study: Non-Markovian dynamics, systematic coarse-graining, rheology, nonequilibrium dynamics |

Education

01/10/2014 | UNIVERSITY OF MAINZ

Doctor rerum naturalium (fast-track program)

13/12/2018 Advisor: Prof. Friederike Schmid

Thesis title: 'Frequency-dependent phenomena and memory in soft matter systems.' Grade: 0.7

(summa cum laude)

21/12/2017 UNIVERSITY OF MAINZ

Master of Science

Advisor: Prof. Friederike Schmid

Thesis title: 'Frequency-dependent hydrodynamic interaction between two solid spheres.' Grade: 1.0

24/10/2011 | UNIVERSITY OF MAINZ

Bachelor of Science

15/08/2014 Advisor: Prof. Friederike Schmid

Thesis title: 'Phase diagrams of model lipid bilayers.' Grade: 1.0 (with distinction)

12/08/2002 | GYMNASIUM ELTVILLE

High School

07/06/2011

Awards

2021: JSPS 'short-term postdoctoral fellowship'

2019: Dr. rer. nat. with 'summa cum laude'

2016 - 2019: Member of the 'Graduate School of Excellence Materials Science in Mainz'

2013 - 2015: Admission to the 'Studienstiftung des deutschen Volkes' (scholarship)

2013 - 2014: Recipient of the 'Deutschlandstipendium' (scholarship)

Teaching and Mentoring

LECTURER | Advanced statistical physics

TUTORIALS | Mathematical methods of physics, modelling (computer science), statistical

physics, electrodynamics, computer simulations in statistical physics

SUPERVISOR | Co-supervision of one PhD student

Co-supervision of one Master student

Supervision of two Bachelor students and one Master student

Organization of Academic Events

2018: CECAM workshop on 'Dynamic coarse-graining and memory effects in soft matter systems' (2 days)

2017: SFB TRR146 students retreat (5 days)

Languages

ENGLISH: fluent

GERMAN: mother tongue FRENCH: intermediate

Publications

Publications in peer-reviewed journals

- 1. Fluctuation-dissipation relations far from equilibrium: A case study (Open Access) **G. Jung**, F. Schmid, Soft Matter **17**, 6413 (2021)
- 2. Introducing memory in coarse-grained molecular simulations (Review Article, Open Access) V. Klippenstein, M. Tripathy, **G. Jung**, F. Schmid, N. van der Vegt, JCPB **125**, 4931 (2021)
- 3. Tagged-particle motion in quasi-confined colloidal hard-sphere liquids
 L. Schrack, C. F. Petersen, **G. Jung**, M. Caraglio, T. Franosch, J. Stat. Mech **043301** (2021)
- 4. Model reduction techniques for the computation of extended Markov parameterizations for generalized Langevin equations (Open Access)
 - N. Bockius, J. Shea, G. Jung, F. Schmid, M. Hanke, JCMP 33, 214003 (2021)
- 5. Wall slip and bulk yielding in soft particle suspensions **G. Jung**, S. Fielding, Journal of Rheology **65**, 199 (2021)
- 6. An improved integration scheme for mode-coupling-theory equations M. Caraglio, L. Schrack, **G. Jung**, T. Franosch, *Comm. Comp. Phys.* **29**, 628 (2021)
- 7. Tagged-particle dynamics in confined colloidal liquids **G. Jung**, L. Schrack, T. Franosch, *Phys. Rev. E* **102**, 032611 (2020)
- 8. Confinement-induced demixing and crystallization **G. Jung**, C. F. Petersen, *Phys. Rev. Res.* **2**, 033207 (2020)
- 9. Dynamic properties of quasi-confined colloidal hard-sphere liquids near the glass transition L. Schrack, C. F. Petersen, **G. Jung**, M. Caraglio, T. Franosch, *J. Stat. Mech.* 093301 (2020)
- Dynamical properties of densely packed confined hard-sphere fluids
 G. Jung, M. Caraglio, L. Schrack, T. Franosch, *Phys. Rev. E* 102, 012612 (2020)
- 11. Scaling equations for mode-coupling theories with multiple decay channels **G. Jung**, T. Voigtmann, T. Franosch, *J. Stat. Mech.* **7**, 073301 (2020)
- 12. Frequency-dependent dielectric polarizability of flexible polyelectrolytes in electrolyte solution: A Dissipative Particle Dynamics simulation
 - G. Jung, S. Kasper, F. Schmid, J. of the Electrochemical Soc. 166, B3194 (2019)
- 13. Generalized Langevin dynamics: Construction and numerical integration of non-Markovian particlebased models
 - **G. Jung**, M. Hanke, F. Schmid, *Soft Matter* **14**, 9368 (2018).
- 14. Frequency-dependent hydrodynamic interactions between two solid spheres
 - **G. Jung**, F. Schmid, *Phys. of Fluids* **29**, 126101 (2017)
- Iterative reconstruction of memory kernels
 G. Jung, M. Hanke, F. Schmid, J. Chem. Theory and Comp. 13, 2481 (2017).
- Computing bulk and shear viscosities from simulations of fluids with dissipative and stochastic interactions
 - **G. Jung**, F. Schmid, *J. Chem. Phys.* **144**, 204104 (2016)

Preprints

1. Non-Markovian systems out of equilibrium: Exact results for two routes of coarse graining **G. Jung**, submitted. https://arxiv.org/abs/2111.13153

Review activities

Reviewer for international journals: Physical Review Letters, Journal of Chemical Physics, Europhysics Letters, Physical Review E, Molecular Simulation, Macromolecular Theory and Simulations