Jan Janssen

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PROFESSIONAL APPOINTMENTS

2021 – present Postdoctoral Research Associate

Theoretical Division (T-1)

Los Alamos National Laboratory

Los Alamos, NM, USA

2015 – 2021 PhD Candidate

Computational Materials Design

Max-Planck-Institut für Eisenforschung

Düsseldorf, Germany

2017 Invited Fellow

Institute for Pure and Applied Mathematics

University of California Los Angeles, CA, USA

EDUCATION

2015 – 2021 PhD in Theoretical Physics, Paderborn University, Germany

2010 – 2011 Advanced degree in Theoretical Physics (Master's equivalent), Technical University of Kaiserslautern, Germany

AWARDS & HONORS

2019 Runner-up for the Heinz Billing Award of 2019

PUBLICATIONS

L.F. Zhu, **J. Janssen**, S. Ishibashi, F. Körmann, B. Grabowski and J. Neugebauer. A fully automated approach to calculate the melting temperature of elemental crystals. *Computational Materials Science*. doi:10.1016/j.commatsci.2020.110065.

• Code: • pyiron/pyiron_meltingpoint

T.D. Swinburne, **J. Janssen**, M. Todorova, G. Simpson, P. Plechac, M. Luskin and J. Neugebauer. Anharmonic free energy of lattice vibrations in fcc crystals from a mean field bond. *Physical Review B.* doi:10.1103/PhysRevB.102.100101.

• Code: O tomswinburne/BLaSA

2019

- **J. Janssen**, S. Surendralal, Y. Lysogorskiy, M. Todorova, T. Hickel, R. Drautz and J. Neugebauer. pyiron: an integrated development environment for computational materials science. *Computational Materials Science*. doi:10.1016/j.commatsci.2018.07.043.
- Code: pyiron
- Y. Lysogorskiy, T. Hammerschmidt, **J. Janssen**, J. Neugebauer and R. Drautz. Transferability of interatomic potentials for molybdenum and silicon. *Modelling and Simulation in Materials Science and Engineering*. doi:10.1088/1361-651X/aafd13

2016

J. Janssen, N. Gunkelmann and H. M. Urbassek. Influence of C concentration on elastic moduli of $\alpha' - Fe_{1-x}C_x$ alloys. *Philosophical Magazine*. doi:10.1080/14786435.2016.1170224

OPEN-SOURCE SOFTWARE

2015 – present **pyiron**

An integrated development environment for computational materials science

• Role: Project Lead Developer

Code: pyironWebsite: pyiron.org

2018 – present Conda-Forge

A community-led collection of recipes, build infrastructure and distributions for the conda package manager.

• Role: Maintainer for 400+ Materials Science Software Packages

Code: conda-forgeWebsite: conda-forge.org

WORKSHOPS

2021 Workflows for Atomistic Simulation

Max-Planck-Institut für Eisenforschung, Düsseldorf, Germany (online).

• Code: pyiron/potentials-workshop-2021

2020 Software Tools from Atomistics to Phase Diagrams

Max-Planck-Institut für Eisenforschung, Düsseldorf, Germany (online).

• Code: pyiron/phasediagram-workshop-2020

PRESENTATIONS

INVITED TALKS

2021

J. Janssen, T. Hickel, J. Neugebauer. pyiron - an integrated development environment for materials science, *CECAM Workshop: Simulation Workflows in Materials Modelling (SWiMM)*, Lausanne, Switzerland (online).

2020

- **J. Janssen**, T. Hickel, J. Neugebauer. Uncertainty quantification for ab initio thermodynamics, *Group Seminar Professor R. Drautz*, Bochum, Germany (online).
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated ab-initio determination of materials properties at finite temperatures with pyiron, *NIST Workshop: Atomistic simulations for industrial needs*, Gaitgersburg, MD, USA (online).

2019

- **J. Janssen**, T. Hickel, J. Neugebauer. Automated ab-initio determination of materials properties at finite temperatures with pyiron, *CLNS Seminar Los Alamos National Laboratory*, Los Alamos, NM, USA.
- **J. Janssen**, T. Hickel, J. Neugebauer. pyiron an integrated development environment for computational materials science, *Group Seminar Professor G. Kresse*, Vienna, Austria.

TALKS AT INTERNATIONAL CONFERENCES

- 2019
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated uncertainty quantification for ab initio thermodynamics, *MRS Fall Meeting*, Boston, MA, USA.
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated sensitivity analysis for high-throughput ab initio calculations, *IPAM Workshop*, Los Angeles, CA, USA.
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated error analysis and control for ab initio calculations, *DPG Spring Meeting*, Regensburg, Germany.
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated sensitivity analysis for high-throughput ab initio calculations, *TMS Spring Meeting*, San Antonio, TX, USA.
- 2018
- **J. Janssen**, T. Hickel, J. Neugebauer. Generation of ab initio datasets with predefined precision using uncertainty quantification, *DPG Spring Meeting*, Berlin, Germany.
- 2017
- **J. Janssen**, T. Hickel, J. Neugebauer. Towards an uncertainty quantification for ab initio thermodynamics, *MRS Fall Meeting*, Boston, MA, USA.
- **J. Janssen**, T. Hickel, J. Neugebauer. Sensitivity analysis for large sets of density functional theory calculations, *DPG Spring Meeting*, Dresden, Germany.
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated convergence and error analyses for high-precision DFT calculations, *TMS Spring Meeting*, San Diego, CA, USA.
- 2016
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated convergence checks with the python based library pyiron, *DPG Spring Meeting*, Regensburg, Germany.
- **J. Janssen**, T. Hickel, J. Neugebauer. Automated convergence checks with the python based workbench pyiron, *TMS Spring Meeting*, Nashville, TN, USA.