

# Till Junge

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1527676/publications.pdf>

Version: 2024-02-01

10  
papers

319  
citations

1162367

8  
h-index

1372195

10  
g-index

10  
all docs

10  
docs citations

10  
times ranked

411  
citing authors

#	ARTICLE	IF	CITATIONS
1	Efficient implementation of atom-density representations. <i>Journal of Chemical Physics</i> , 2021, 154, 114109.	1.2	32
2	The emergence of small-scale self-affine surface roughness from deformation. <i>Science Advances</i> , 2020, 6, eaax0847.	4.7	48
3	Atomistic-continuum coupling of random alloys. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2019, 27, 075004.	0.8	5
4	Molecular probes reveal deviations from Amontons's law in multi-asperity frictional contacts. <i>Nature Communications</i> , 2018, 9, 888.	5.8	88
5	The Coupled Atomistic/Discrete-Dislocation method in 3d part I: Concept and algorithms. <i>Journal of the Mechanics and Physics of Solids</i> , 2018, 118, 152-171.	2.3	31
6	Ab initio modelling of the early stages of precipitation in Al-6000 alloys. <i>Acta Materialia</i> , 2017, 140, 240-249.	3.8	14
7	Dynamic stability of displacement-based atomistic/continuum coupling methods. <i>Journal of the Mechanics and Physics of Solids</i> , 2015, 80, 103-120.	2.3	8
8	Toward a 3D coupled atomistic and discrete dislocation dynamics simulation: dislocation core structures and Peierls stresses with several character angles in FCC aluminum. <i>Advanced Modeling and Simulation in Engineering Sciences</i> , 2015, 2, .	0.7	27
9	Plastic activity in nanoscratch molecular dynamics simulations of pure aluminium. <i>International Journal of Plasticity</i> , 2014, 53, 90-106.	4.1	57
10	Molecular dynamics nano-scratching of aluminium: a novel quantitative energy-based analysis method. <i>Procedia IUTAM</i> , 2012, 3, 192-204.	1.2	9