

A Prakash, Aruna Prakash

List of Publications by Year in descending order

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31
papers

853
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times ranked

906
citing authors

#	ARTICLE	IF	CITATIONS
1	Simulation of micromechanical behavior of polycrystals: finite elements versus fast Fourier transforms. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2009, 17, 064010.	2.0	122
2	Organic Memory Device Fabricated Through Solution Processing. <i>Proceedings of the IEEE</i> , 2005, 93, 1287-1296.	21.3	98
3	Assessment and optimization of the fast inertial relaxation engine (fire) for energy minimization in atomistic simulations and its implementation in lammers. <i>Computational Materials Science</i> , 2020, 175, 109584.	3.0	88
4	Atom probe informed simulations of dislocationâ€“precipitate interactions reveal the importance of local interface curvature. <i>Acta Materialia</i> , 2015, 92, 33-45.	7.9	79
5	Modeling the evolution of texture and grain shape in Mg alloy AZ31 using the crystal plasticity finite element method. <i>Computational Materials Science</i> , 2009, 45, 744-750.	3.0	66
6	A multiscale simulation framework of the accumulative roll bonding process accounting for texture evolution. <i>Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing</i> , 2015, 631, 104-119.	5.6	42
7	Quantifying eigenstrain distributions induced by focused ion beam damage in silicon. <i>Materials Letters</i> , 2016, 185, 47-49.	2.6	36
8	Influence of grain boundary structure and topology on the plastic deformation of nanocrystalline aluminum as studied by atomistic simulations. <i>International Journal of Plasticity</i> , 2017, 97, 107-125.	8.8	36
9	Twinning Models in Self-Consistent Texture Simulations of TWIP Steels. <i>Steel Research International</i> , 2008, 79, 645-652.	1.8	34
10	Atomistic simulations of basal dislocations in Mg interacting with Mg ₁₇ Al ₁₂ precipitates. <i>Materialia</i> , 2019, 7, 100355.	2.7	31
11	Nano: A methodology for generating complex realistic configurations for atomistic simulations. <i>MethodsX</i> , 2016, 3, 219-230.	1.6	24
12	Atomistic simulations of focused ion beam machining of strained silicon. <i>Applied Surface Science</i> , 2017, 416, 86-95.	6.1	24
13	In-situ observation of the initiation of plasticity by nucleation of prismatic dislocation loops. <i>Nature Communications</i> , 2020, 11, 2367.	12.8	23
14	High Lightweight Potential of Ultrafineâ€“Grained Aluminum/Steel Laminated Metal Composites Produced by Accumulative Roll Bonding. <i>Advanced Engineering Materials</i> , 2019, 21, 1800286.	3.5	21
15	FE2ATâ€“finite element informed atomistic simulations. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2013, 21, 055011.	2.0	16
16	Idealized vs. Realistic Microstructures: An Atomistic Simulation Case Study on $\hat{\Gamma}_3/\hat{\Gamma}_3$ â€“Microstructures. <i>Materials</i> , 2017, 10, 88.	2.9	16
17	Origins of strengthening and failure in twinned Au nanowires: Insights from inâ€“situ experiments and atomistic simulations. <i>Acta Materialia</i> , 2020, 187, 166-175.	7.9	15
18	A hierarchical multi-scale model for hexagonal materials taking into account texture evolution during forming simulation. <i>Computational Materials Science</i> , 2014, 82, 464-475.	3.0	13

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19	Influence of intrinsic strain on irradiation induced damage: the role of threshold displacement and surface binding energies. <i>Materials and Design</i> , 2016, 111, 405-413.	7.0	12
20	Atomistic Simulations of Compression Tests on $\hat{\beta}$ -Precipitate Containing Ni ₃ Al Nanocubes. <i>Metallurgical and Materials Transactions A: Physical Metallurgy and Materials Science</i> , 2018, 49, 4158-4166.	2.2	12
21	Grain segmentation in atomistic simulations using orientation-based iterative self-organizing data analysis. <i>Materialia</i> , 2022, 21, 101314.	2.7	9
22	Chances and Challenges in Fusing Data Science with Materials Science. <i>Praktische Metallographie/Practical Metallography</i> , 2018, 55, 493-514.	0.3	8
23	Enhanced monotonic and cyclic mechanical properties of ultrafine-grained laminated metal composites with strong and stiff interlayers. <i>International Journal of Fatigue</i> , 2018, 116, 379-387.	5.7	8
24	Experimental and Numerical Investigation of Texture Development during Hot Rolling of Magnesium Alloy AZ31. <i>Materials Science Forum</i> , 2007, 539-543, 3448-3453.	0.3	5
25	On solution mapping and remeshing in crystal plasticity finite element simulations: application to equal channel angular pressing. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2019, 27, 075001.	2.0	4
26	OptiMic: A tool to generate optimized polycrystalline microstructures for materials simulations. <i>SoftwareX</i> , 2021, 15, 100708.	2.6	4
27	Sintering of Alumina Nanoparticles: Comparison of Interatomic Potentials, Molecular Dynamics Simulations, and Data Analysis. <i>Modelling and Simulation in Materials Science and Engineering</i> , 0, , .	2.0	3
28	Determination of thermal accommodation coefficients on C/a and S/a using molecular d. <i>International Journal of Heat and Mass Transfer</i> , 2022, 183, 122219.	4.8	2
29	Fatigue Life Optimized Layer Architecture of Ultrafine-Grained Al-Ti Laminates Under Bending Stresses. <i>Advanced Engineering Materials</i> , 2022, 24, .	3.5	2
30	8-inch wafer-scale HfO ₂ -based RRAM for 1S-1R cross-point memory applications. , 2014, , .		0
31	9. FE2AT “ finite element informed atomistic simulations. , 2016, , 167-190.		0