

Sai Tang

List of Publications by Year in descending order

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papers

224
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1307594

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1058476

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15
times ranked

258
citing authors

#	ARTICLE	IF	CITATIONS
1	Growth modes of grain boundary precipitate in aluminum alloys under different lattice misfits. <i>Journal of Materials Science</i> , 2022, 57, 2744-2757.	3.7	9
2	Molecular-Level Insights into the Nucleation Mechanism of One-Component Soft Matter Icosahedral Quasicrystal Studied by Phase-Field Crystal Simulations. <i>Crystal Growth and Design</i> , 2022, 22, 2637-2643.	3.0	4
3	Two-dimensional ordering governs the overpotential of Li intercalation and plating on graphene and its variants. <i>Journal of Applied Physics</i> , 2022, 131, .	2.5	1
4	The atomic scale mechanisms of the interaction between pore and grain boundary during sintering. <i>Materials Today Communications</i> , 2022, 32, 103970.	1.9	2
5	Atomic-scale study of compositional and structural evolution of early-stage grain boundary precipitation in Al-Cu alloys through phase-field crystal simulation. <i>Journal of Materials Science</i> , 2021, 56, 12700-12715.	3.7	12
6	Phase-field crystal study on the influence of nanoscale intergranular structures on grain boundary segregation. <i>Modelling and Simulation in Materials Science and Engineering</i> , 2021, 29, 085009.	2.0	1
7	Phase-field simulation of solidification microstructure in Ni and Cu-Ni alloy using the Wheeler, Boettinger and McFadden model coupled with the CALPHAD data. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 2020, 68, 101691.	1.6	4
8	An atomic scale study of two-dimensional quasicrystal nucleation controlled by multiple length scale interactions. <i>Soft Matter</i> , 2020, 16, 5718-5726.	2.7	5
9	Competitive bcc and fcc crystal nucleation from non-equilibrium liquids studied by phase-field crystal simulation. <i>Acta Materialia</i> , 2017, 139, 196-204.	7.9	34
10	Phase-field-crystal investigation of the morphology of a steady-state dendrite tip on the atomic scale. <i>Physical Review E</i> , 2017, 95, 062803.	2.1	7
11	Kinetic Pathways and Mechanisms of Two-Step Nucleation in Crystallization. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 5008-5014.	4.6	50
12	Strain mapping in nanocrystalline grains simulated by phase field crystal model. <i>Philosophical Magazine</i> , 2015, 95, 973-984.	1.6	11
13	Phase-field-crystal simulation of nonequilibrium crystal growth. <i>Physical Review E</i> , 2014, 89, 012405.	2.1	38
14	Orientation selection process during the early stage of cubic dendrite growth: A phase-field crystal study. <i>Acta Materialia</i> , 2012, 60, 5501-5507.	7.9	45
15	A New Multi-Phase Field Model for the Electrochemical Corrosion of Aluminum Alloys. <i>Advanced Theory and Simulations</i> , 0, , 2200299.	2.8	1