

Roberto Gomes de Aguiar Veiga

List of Publications by Year in descending order

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Version: 2024-02-01

27
papers

550
citations

687363

13
h-index

610901

24
g-index

28
all docs

28
docs citations

28
times ranked

631
citing authors

#	ARTICLE	IF	CITATIONS
1	Atomistic study of metallurgical bonding upon the high velocity impact of fcc core-shell particles. Computational Materials Science, 2021, 186, 110045.	3.0	14
2	Crystal orientation and grain boundary effects on plastic deformation of FCC particles under high velocity impacts. Materialia, 2021, 15, 101004.	2.7	11
3	Effect of Sodium Chloride on Internal Quasi-Liquid Layers in Ice I _h . Journal of Physical Chemistry C, 2021, 125, 18526-18535.	3.1	5
4	Investigating the kinetics of the formation of a Cottrell atmosphere around a screw dislocation in bcc iron: a mixed-lattice atomistic kinetic Monte-Carlo analysis. Journal of Physics Condensed Matter, 2021, 33, 065704.	1.8	1
5	Deformation of copper particles upon impact: A molecular dynamics study of cold spray. Computational Materials Science, 2020, 171, 109219.	3.0	52
6	Structural changes and kinetics of shear banding in metallic glass composites. Journal of Alloys and Compounds, 2020, 819, 153046.	5.5	7
7	Atomistic Study of the Role of Defects on ϵ Phase Transformations in Iron under Hydrostatic Compression. Metals, 2019, 9, 1040.	2.3	15
8	Elastic constants of ice I _h as described by semi-empirical water models. Journal of Chemical Physics, 2019, 150, 044503.	3.0	10
9	Anomalous diffusion of water molecules at grain boundaries in ice I _h . Physical Chemistry Chemical Physics, 2018, 20, 13944-13951.	2.8	15
10	Local atomic environment and shear banding in metallic glasses. Computational Materials Science, 2018, 155, 129-135.	3.0	8
11	Interaction between interstitial carbon atoms and a $\frac{1}{2}$ self-interstitial atoms loop in an iron matrix: a combined DFT, off lattice KMC and MD study. Journal of Physics Condensed Matter, 2018, 30, 335901.	1.8	2
12	Lattice Instabilities and Phase Transformations in Fe from Atomistic Simulations. Journal of Phase Equilibria and Diffusion, 2017, 38, 185-194.	1.4	1
13	Stability of nanocrystalline Ni-based alloys: coupling Monte Carlo and molecular dynamics simulations. Modelling and Simulation in Materials Science and Engineering, 2017, 25, 075005.	2.0	4
14	Formation of carbon Cottrell atmospheres and their effect on the stress field around an edge dislocation. Scripta Materialia, 2017, 129, 16-19.	5.2	35
15	Adsorption of metal-phthalocyanine molecules onto the Si(111) surface passivated by δ -doping: Ab initio calculations. Physical Review B, 2016, 93, .	3.2	17
16	Self-assembly of NiTPP on Cu(111): a transition from disordered 1D wires to 2D chiral domains. Physical Chemistry Chemical Physics, 2015, 17, 18344-18352.	2.8	12
17	Monte Carlo and molecular dynamics simulations of screw dislocation locking by Cottrell atmospheres in low carbon Fe-C alloys. Scripta Materialia, 2015, 108, 19-22.	5.2	50
18	Comments on Atomistic modeling of an Fe system with a small concentration of C. Computational Materials Science, 2014, 82, 118-121.	3.0	35

#	ARTICLE	IF	CITATIONS
19	Atomistic modeling of carbon Cottrell atmospheres in bcc iron. Journal of Physics Condensed Matter, 2013, 25, 025401.	1.8	35
20	Thermal conductivity of nanocrystalline SiGe alloys using molecular dynamics simulations. Journal of Applied Physics, 2013, 114, 164310.	2.5	10
21	Modified embedded-atom method interatomic potential and interfacial thermal conductance of Si-Cu systems: A molecular dynamics study. Journal of Applied Physics, 2013, 113, .	2.5	6
22	Comparison of atomistic and elasticity approaches for carbon diffusion near line defects in $\hat{\pm}$ -iron. Acta Materialia, 2011, 59, 6963-6974.	7.9	49
23	Structural, electronic, and magnetic properties of pristine and oxygen-adsorbed graphene nanoribbons. Applied Surface Science, 2010, 256, 5776-5782.	6.1	12
24	Molecular dynamics study of the ordering of carbon in highly supersaturated $\hat{\pm}$ -Fe. Physical Review B, 2010, 81, .	3.2	50
25	Effect of the stress field of an edge dislocation on carbon diffusion in $\hat{\pm}$ -iron: Coupling molecular statics and atomistic kinetic Monte Carlo. Physical Review B, 2010, 82, .	3.2	42
26	Quenching of local magnetic moment in oxygen adsorbed graphene nanoribbons. Journal of Chemical Physics, 2008, 128, 201101.	3.0	29
27	Ab initio study of TCNQ-doped carbon nanotubes. Physical Review B, 2006, 73, .	3.2	23