

Zamaan Raza

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/4030550/publications.pdf>

Version: 2024-02-01

15

papers

677

citations

567281

15

h-index

996975

15

g-index

15

all docs

15

docs citations

15

times ranked

1114

citing authors

#	ARTICLE	IF	CITATIONS
1	The microscopic features of heterogeneous ice nucleation may affect the macroscopic morphology of atmospheric ice crystals. <i>Faraday Discussions</i> , 2013, 167, 389.	3.2	80
2	GLYCOLALDEHYDE FORMATION VIA THE DIMERIZATION OF THE FORMYL RADICAL. <i>Astrophysical Journal</i> , 2013, 777, 90.	4.5	62
3	High Energy Density Mixed Polymeric Phase from Carbon Monoxide and Nitrogen. <i>Physical Review Letters</i> , 2013, 111, 235501.	7.8	62
4	Proton ordering in cubic ice and hexagonal ice; a potential new ice phaseâ€”Xlc. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 19788.	2.8	60
5	Lattice Vibrations Change the Solid Solubility of an Alloy at High Temperatures. <i>Physical Review Letters</i> , 2016, 117, 205502.	7.8	60
6	Effect of salt on the H-bond symmetrization in ice. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 8216-8220.	7.1	58
7	Large scale and linear scaling DFT with the CONQUEST code. <i>Journal of Chemical Physics</i> , 2020, 152, 164112.	3.0	55
8	ON THE FORMATION OF GLYCOLALDEHYDE IN DENSE MOLECULAR CORES. <i>Astrophysical Journal</i> , 2012, 750, 19.	4.5	54
9	Temperature-dependent elastic properties of $Ti_{1-x}Al_xN$ alloys. <i>Applied Physics Letters</i> , 2015, 107, .	3.3	46
10	From cellulose to kerogen: molecular simulation of a geological process. <i>Chemical Science</i> , 2017, 8, 8325-8335.	7.4	37
11	Novel superconducting skutterudite-type phosphorus nitride at high pressure from first-principles calculations. <i>Scientific Reports</i> , 2014, 4, 5899.	3.3	29
12	First-principles calculations of properties of orthorhombic iron carbide Fe_{3-x}C at the Earth's core conditions. <i>Physical Review B</i> , 2015, 91, .	3.2	20
13	Impact of anharmonic effects on the phase stability, thermal transport, and electronic properties of AlN. <i>Physical Review B</i> , 2016, 94, .	3.2	20
14	Computer simulations of glasses: the potential energy landscape. <i>Journal of Physics Condensed Matter</i> , 2015, 27, 293201.	1.8	17
15	Highly accurate local basis sets for large-scale DFT calculations in conquest. <i>Japanese Journal of Applied Physics</i> , 2019, 58, 100503.	1.5	17