

Amol B Rahane

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/794343/publications.pdf>

Version: 2024-02-01

12
papers

220
citations

1039406

9
h-index

1199166

12
g-index

12
all docs

12
docs citations

12
times ranked

364
citing authors

#	ARTICLE	IF	CITATIONS
1	Structural and Electronic Properties of $(Al_2O_3)_n$ Clusters with $n = 1-10$ from First Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2011, 115, 18111-18121.	1.5	60
2	B_{84} : a quasi-planar boron cluster stabilized with hexagonal holes. <i>Nanoscale</i> , 2015, 7, 4055-4062.	2.8	44
3	Density Functional Calculations of the Structural and Electronic Properties of $(Y_2O_3)_n$, $n=1-10$ Clusters with $n = 1-10$. <i>Journal of Physical Chemistry A</i> , 2013, 117, 5542-5550.	1.1	22
4	Electronic Origin of the Stability of Transition-Metal-Doped B_{14} Drum-Shaped Boron Clusters and Their Assembly into a Nanotube. <i>Journal of Physical Chemistry C</i> , 2017, 121, 10728-10742.	1.5	19
5	Structural and Electronic Properties of Neutral and Ionic $(Ga_2O_3)_n$ Clusters with $n = 1-10$. <i>Journal of Physical Chemistry C</i> , 2012, 116, 2691-2701.	1.5	17
6	Lithium storage in amorphous TiNi hydride: Electrode for rechargeable lithium-ion batteries. <i>Materials Chemistry and Physics</i> , 2013, 141, 348-354.	2.0	15
7	Analysis of the electron density features of small boron clusters and the effects of doping with C, P, Al, Si, and Zn: Magic B_7P and B_8Si clusters. <i>Physica Scripta</i> , 2016, 91, 053005.	1.2	13
8	First Principles Calculations for Structural, Electronic, and Magnetic Properties of Gadolinium-Doped Alumina Clusters. <i>Journal of Physical Chemistry C</i> , 2012, 116, 6115-6126.	1.5	11
9	Optical Properties of Gallium Oxide Clusters from First-Principles Calculations. <i>Journal of Physical Chemistry A</i> , 2012, 116, 10559-10565.	1.1	11
10	Carbon Doping in Boron Suboxide: Structure, Energetics, and Elastic Properties. <i>Journal of the American Ceramic Society</i> , 2015, 98, 2223-2233.	1.9	5
11	Theoretical study of small clusters of manganese-doped gallium oxide: $Mn(GaO)_n$ and $Mn_2(GaO)_n$ with $n=1-7$. <i>Journal of Nanoparticle Research</i> , 2010, 12, 727-736.	0.8	2
12	Time dependent DFT investigation of the optical response in pristine and Gd doped Al_2O_3 . <i>RSC Advances</i> , 2016, 6, 72537-72543.	1.7	1