

# Amol B Rahane

## List of Publications by Year in descending order

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12  
papers

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citations

1039406

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1199166

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g-index

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docs citations

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times ranked

364  
citing authors

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