

Kamalika Ghatak

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/3527946/publications.pdf>

Version: 2024-02-01

12
papers

225
citations

1162367

8
h-index

1372195

10
g-index

12
all docs

12
docs citations

12
times ranked

388
citing authors

#	ARTICLE	IF	CITATIONS
1	Scandium-Decorated MOF-5 as Potential Candidates for Room-Temperature Hydrogen Storage: A Solution for the Clustering Problem in MOFs. <i>Journal of Physical Chemistry C</i> , 2012, 116, 17336-17342.	1.5	50
2	Utilizing van der Waals Slippery Interfaces to Enhance the Electrochemical Stability of Silicon Film Anodes in Lithium-Ion Batteries. <i>ACS Applied Materials & Interfaces</i> , 2018, 10, 13442-13451.	4.0	48
3	Effect of cobalt content on the electrochemical properties and structural stability of NCA type cathode materials. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 22805-22817.	1.3	27
4	Amorphous germanium as a promising anode material for sodium ion batteries: a first principle study. <i>Journal of Materials Science</i> , 2018, 53, 14423-14434.	1.7	23
5	Metal or Nonmetal Cooperation with a Phenyl Group: Route to Catalysis? A Computational Investigation. <i>ACS Catalysis</i> , 2013, 3, 920-927.	5.5	21
6	Controlled edge dependent stacking of WS ₂ -WS ₂ Homo- and WS ₂ -WSe ₂ Hetero-structures: A Computational Study. <i>Scientific Reports</i> , 2020, 10, 1648.	1.6	19
7	A computational investigation of the role of the iridium dihydrogen pincer complex in the formation of the cyclic pentamer (NH ₂ BH ₂) ₅ . <i>Computational and Theoretical Chemistry</i> , 2012, 992, 18-29.	1.1	12
8	Computational investigation on the catalytic activity of Rh ₆ and Rh ₄ Ru ₂ clusters towards methanol activation. <i>Theoretical Chemistry Accounts</i> , 2015, 134, 1.	0.5	10
9	The inherent behavior of graphene flakes in water: A molecular dynamics study. <i>Computational Materials Science</i> , 2019, 162, 140-147.	1.4	7
10	Atomistic modeling by density functional theory of two-dimensional materials. , 2020, , 113-123.		4
11	Exploring the Potential of Doped Zero-Dimensional Cages for Proton Transfer in Fuel Cells: A Computational Study. <i>Journal of Physical Chemistry B</i> , 2012, 116, 9803-9811.	1.2	3
12	Two-dimensional materials and its heterostructures for energy storage. , 2020, , 385-401.		1