

Sree Ganesh Balasubramani

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2595106/publications.pdf>

Version: 2024-02-01

9

papers

905

citations

1307594

7

h-index

1474206

9

g-index

9

all docs

9

docs citations

9

times ranked

1161

citing authors

#	ARTICLE	IF	CITATIONS
1	Transition Path Sampling Based Calculations of Free Energies for Enzymatic Reactions: The Case of Human Methionine Adenosyl Transferase and <i>< i>Plasmodium vivax</i></i> Adenosine Deaminase. <i>Journal of Physical Chemistry B</i> , 2022, 126, 5413-5420.	2.6	4
2	Exploring the Solvation of Acetic Acid in Water Using Liquid Jet X-ray Photoelectron Spectroscopy and Core Level Electron Binding Energy Calculations. <i>Journal of Physical Chemistry B</i> , 2021, 125, 8862-8868.	2.6	6
3	High-Resolution X-ray Photoelectron Spectroscopy of Organometallic $(C_{5}H_{4}SiMe_3)_3Ln^{III}$ and $[(C_{5}H_{4}SiMe_3)_3Ln^{II}]^{+}$ Complexes ($Ln = Sm, Eu, Cd, Tb$). <i>Journal of the American Chemical Society</i> , 2021, 143, 16610-16620.	13.7	17
4	TURBOMOLE: Modular program suite for <i>< i>ab initio</i></i> quantum-chemical and condensed-matter simulations. <i>Journal of Chemical Physics</i> , 2020, 152, 184107.	3.0	616
5	Formation of the End-on Bound Lanthanide Dinitrogen Complexes $[(R_{2}N)_{3}Ln^{+}N_{2}^{+}]^{2+}$ from Divalent $[(R_{2}N)_{3}Ln]^{+}$ Salts ($R = SiMe_3$). <i>Journal of the American Chemical Society</i> , 2020, 142, 9302-9313.	13.7	15
6	Variational generalized Kohn-Sham approach combining the random-phase-approximation and Green's-function methods. <i>Physical Review A</i> , 2019, 99, .	2.5	39
7	Synthesis, Structure, and Magnetism of Tris(amide) $[Ln\{N(SiMe_3)_2\}_2]^{1+}$ Complexes of the Non-traditional +2 Lanthanide Ions. <i>Chemistry - A European Journal</i> , 2018, 24, 7702-7709.	3.3	64
8	Random-Phase Approximation Methods. <i>Annual Review of Physical Chemistry</i> , 2017, 68, 421-445.	10.8	127
9	Noble gas encapsulation into carbon nanotubes: Predictions from analytical model and DFT studies. <i>Journal of Chemical Physics</i> , 2014, 141, 184304.	3.0	17