

Aulia Sukma Hutama

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9577331/publications.pdf>

Version: 2024-02-01

8
papers

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citations

1937685

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docs citations

8
times ranked

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citing authors

#	ARTICLE	IF	CITATIONS
1	Photophysical properties of ammonium, pyrrolidinium, piperidinium, imidazolium, and pyridinium as a guide to prepare ionic-organic hybrid materials. <i>Heliyon</i> , 2022, 8, e09121.	3.2	3
2	Theoretical studies on structure and dynamics of anatase TiO ₂ (101)/H ₂ SO ₄ /H ₂ O interface in the early stage of titania sulfation. <i>Structural Chemistry</i> , 2022, 33, 1341-1354.	2.0	1
3	Density-Functional Tight-Binding Parameters for Bulk Zirconium: A Case Study for Repulsive Potentials. <i>Journal of Physical Chemistry A</i> , 2021, 125, 2184-2196.	2.5	2
4	Development of Density-Functional Tight-Binding Parameters for the Molecular Dynamics Simulation of Zirconia, Yttria, and Yttria-Stabilized Zirconia. <i>ACS Omega</i> , 2021, 6, 20530-20548.	3.5	8
5	Investigations of the influence of non-metal dopants on the electronic and photocatalytic properties of ZrTiO ₄ by density functional theory calculations. <i>Computational Condensed Matter</i> , 2021, 29, e00607.	2.1	2
6	Natural resources for dye-sensitized solar cells. <i>Heliyon</i> , 2021, 7, e08436.	3.2	12
7	Investigation of the chemical and optical properties of halogen-substituted N-methyl-4-piperidone curcumin analogs by density functional theory calculations. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2019, 221, 117152.	3.9	11
8	Coupled Cluster and Density Functional Studies of Atomic Fluorine Chemisorption on Coronene as Model Systems for Graphene Fluorination. <i>Journal of Physical Chemistry C</i> , 2017, 121, 14888-14898.	3.1	12