

# 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

51  
citations

1937685

4  
h-index

1720034

7  
g-index

8  
all docs

8  
docs citations

8  
times ranked

40  
citing authors

#	ARTICLE	IF	CITATIONS
1	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
2	Natural resources for dye-sensitized solar cells. <i>Heliyon</i> , 2021, 7, e08436.	3.2	12
3	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
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	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
6	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
7	Investigations of the influence of non-metal dopants on the electronic and photocatalytic properties of ZrTiO <sub>4</sub> by density functional theory calculations. <i>Computational Condensed Matter</i> , 2021, 29, e00607.	2.1	2
8	Theoretical studies on structure and dynamics of anatase TiO <sub>2</sub> (101)/H <sub>2</sub> SO <sub>4</sub> /H <sub>2</sub> O interface in the early stage of titania sulfation. <i>Structural Chemistry</i> , 2022, 33, 1341-1354.	2.0	1