

Koussai Lazaar

List of Publications by Year in descending order

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Version: 2024-02-01

11
papers

114
citations

1478280

6
h-index

1372474

10
g-index

11
all docs

11
docs citations

11
times ranked

73
citing authors

#	ARTICLE	IF	CITATIONS
1	Electronic and vibrational properties of TMDs heterogeneous bilayers, nontwisted bilayers silicene/TMDs heterostructures and photovoltaic heterojunctions of fullerenes with TMDs monolayers. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2018, 104, 155-164.	1.3	30
2	DFT study of electronic and optical properties of silicene functionalized with chemical groups. <i>Journal of Molecular Graphics and Modelling</i> , 2019, 91, 72-79.	1.3	23
3	A DFT study of Janus structure of S and Se in HfSSe layered as a promising candidate for electronic devices. <i>Journal of Molecular Graphics and Modelling</i> , 2020, 96, 107511.	1.3	13
4	A first principle study of the pristine InBi honeycomb film functionalized with fluorine atoms. <i>Journal of Fluorine Chemistry</i> , 2018, 212, 171-179.	0.9	12
5	A DFT study of GaSe/AlN(ZnO) two-dimensional vdW heterostructure practiced as an encouraging photocatalyst for water splitting. <i>Computational Materials Science</i> , 2022, 201, 110912.	1.4	12
6	DFT study of the electronic and vibrational properties of silicene/stanene heterobilayer. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2019, 111, 127-129.	1.3	10
7	A theoretical investigation of the effect of fluorination and bromination on the optoelectronic properties of tetrathienophenazine derivatives. <i>Computational Materials Science</i> , 2020, 177, 109578.	1.4	6
8	Electronic and optical properties of $W\text{-Sn-Z}$ and $W\text{-Sn-W}$ monolayers using density functional theory. <i>Solid State Communications</i> , 2020, 321, 114016.	0.9	5
9	DFT study of optoelectronic and magnetic properties of a novel type perovskites. <i>Chemical Physics</i> , 2018, 513, 120-128.	0.9	2
10	Effect of halogenation on the optical and electronic properties of tetrathienoanthracene and tetrathionoacridine derivatives: A DFT study. <i>Computational Condensed Matter</i> , 2021, 26, e00528.	0.9	1
11	Electronic structure of 2D quaternary materials and of their van der Waals heterostructures. <i>Journal of Applied Physics</i> , 2021, 130, 064304.	1.1	0