

Peyman Aghdasi Mazandarani

List of Publications by Year in descending order

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14
papers

275
citations

840776

11
h-index

1058476

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all docs

14
docs citations

14
times ranked

79
citing authors

#	ARTICLE	IF	CITATIONS
1	A DFT investigation on the mechanical and structural properties of halogen- and metal-adsorbed silicene nanosheets. <i>Materials Chemistry and Physics</i> , 2022, 283, 126029.	4.0	11
2	Investigating elastic and plastic characteristics of monolayer phosphorene under atomic adsorption by the density functional theory. <i>Physica B: Condensed Matter</i> , 2021, 600, 412603.	2.7	18
3	Analysis of quantum effects of fine scaling on the axial buckling of MWCNTs based on the density functional theory and molecular mechanics method. <i>Applied Physics A: Materials Science and Processing</i> , 2021, 127, 1.	2.3	11
4	A DFT study on the mechanical properties of hydrogenated and fluorinated germanene sheets. <i>Superlattices and Microstructures</i> , 2021, 152, 106854.	3.1	13
5	Torsional buckling analysis of MWCNTs considering quantum effects of fine scaling based on DFT and molecular mechanics method. <i>Journal of Molecular Graphics and Modelling</i> , 2021, 104, 107843.	2.4	17
6	On the derivation of coefficient of Morse potential function for the silicene: a DFT investigation. <i>Journal of Molecular Modeling</i> , 2021, 27, 190.	1.8	14
7	Investigation of elastic properties, buckling and vibration of antimonene nanosheets through DFT-based finite element modeling. <i>Materials Science and Engineering B: Solid-State Materials for Advanced Technology</i> , 2021, 271, 115219.	3.5	11
8	Structural and mechanical properties characterization of arsenene nanosheets under doping effect of transition metals: A DFT study. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2020, 124, 114349.	2.7	28
9	A DFT-based finite element approach for studying elastic properties, buckling and vibration of the arsenene. <i>Journal of Molecular Graphics and Modelling</i> , 2020, 101, 107725.	2.4	21
10	Structural and mechanical properties of Sb and SbX (X=H, F, Cl and Br) monolayers. <i>Solid State Communications</i> , 2020, 311, 113849.	1.9	25
11	Structural and mechanical properties of pristine and adsorbed puckered arsenene nanostructures: A DFT study. <i>Superlattices and Microstructures</i> , 2020, 139, 106414.	3.1	26
12	Influence of F and H adsorption on the elasto-plastic properties of silicene: A DFT investigation. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2020, 119, 113984.	2.7	31
13	Investigating the effects of H and F adsorption on the elastic and plastic properties of arsenene nanosheets. <i>Physica B: Condensed Matter</i> , 2019, 574, 411672.	2.7	23
14	On the elastic and plastic properties of the bismuthene adsorbed by H, F, Cl and Br atoms. <i>Superlattices and Microstructures</i> , 2019, 135, 106242.	3.1	26