

Najim Tahiri

List of Publications by Year in descending order

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55
papers

930
citations

471061

17
h-index

500791

28
g-index

56
all docs

56
docs citations

56
times ranked

348
citing authors

#	ARTICLE	IF	CITATIONS
1	Performance analysis of MAPbI ₃ based perovskite solar cells employing diverse charge selective contacts: Simulation study. <i>Solar Energy</i> , 2019, 193, 948-955.	2.9	218
2	Complexity of vesicle microcirculation. <i>Physical Review E</i> , 2011, 84, 041906.	0.8	58
3	On the problem of slipper shapes of red blood cells in the microvasculature. <i>Microvascular Research</i> , 2013, 85, 40-45.	1.1	42
4	A DFT study of the electronic structure, optical, and thermoelectric properties of halide perovskite KGeI ₃ -xBr _x materials: photovoltaic applications. <i>Applied Physics A: Materials Science and Processing</i> , 2021, 127, 1.	1.1	39
5	Electronic, optical, and thermoelectric properties of perovskite BaTiO ₃ compound under the effect of compressive strain. <i>Chemical Physics</i> , 2021, 544, 111105.	0.9	38
6	Importance of spin-orbit coupling on photovoltaic properties of Pb-free vacancy ordered double		

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19	A Monte Carlo study of the spin-1 Blume-Emery-Griffiths phase diagrams within biquadratic exchange anisotropy. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2014, 407, 295-302.	1.2	17
20	Magnetic, magnetocaloric and transport properties in AlMn ₃ antiperovskite compound. <i>Journal of Alloys and Compounds</i> , 2018, 741, 1196-1202.	2.8	16
21	How the strain effects decreases the band gap energy in the CsPbX ₃ perovskite compounds?. <i>Phase Transitions</i> , 2020, 93, 455-469.	0.6	15
22	Magnetocaloric and thermoelectric properties of the perovskite LaMnO ₃ material: A DFT study and Monte Carlo technique. <i>Phase Transitions</i> , 2021, 94, 826-834.	0.6	15
23	Multilayer transition in a spin-1 Blume-Capel model with RKKY interaction and quantum transverse anisotropy. <i>Chinese Physics B</i> , 2011, 20, 017501.	0.7	14
24	Magnetic Properties of NiFe ₂ O ₄ Compound: Ab Initio Calculation and Monte Carlo Simulation. <i>Journal of Superconductivity and Novel Magnetism</i> , 2020, 33, 1369-1375.	0.8	14
25	Chalcogens TM impurities and a single F-center in perovskite SrHfO ₃ compound: Ab initio calculations. <i>Materials Science in Semiconductor Processing</i> , 2022, 138, 106271.	1.9	14
26	A non-magnetic spacer layer effect on spin layers (7/2,3) in a bi-layer ferromagnetic dendrimer structure: Monte Carlo study. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2016, 462, 1067-1074.	1.2	12
27	Theoretical investigation of electronic, magnetic and magnetocaloric properties of Bi ₂₅ FeO ₄₀ compound. <i>Phase Transitions</i> , 2021, 94, 147-158.	0.6	11
28	Calcium hafnate perovskite from an insulator to a semiconductor for photovoltaic and photocatalytic hydrogen production from water splitting applications. <i>Superlattices and Microstructures</i> , 2021, 160, 107058.	1.4	11
29	Structural, electronic, magnetic, and magnetocaloric properties in intermetallic compound TbCu ₂ Si ₂ . <i>Journal of Magnetism and Magnetic Materials</i> , 2019, 481, 72-76.	1.0	10
30	Magnetocaloric effect in metallic antiperovskite Mn ₃ InC compound: Ab-initio study and Monte Carlo simulations. <i>Solid State Communications</i> , 2020, 309, 113841.	0.9	10
31	Magnetic properties of a Dendrimer structure with RKKY interactions. <i>Chinese Journal of Physics</i> , 2016, 54, 115-120.	2.0	9
32	Physical properties of perovskite SrHfO ₃ compound doped with S for photovoltaic applications: the ab initio study. <i>Applied Physics A: Materials Science and Processing</i> , 2020, 126, 1.	1.1	8
33	The effect of chalcogens-doped with dilation strain on the electronic, optic, and thermoelectric properties of perovskite BaSnO ₃ compound. <i>Journal of the Korean Ceramic Society</i> , 2022, 59, 715-728.	1.1	7
34	Theoretical investigations of electronic structure and optical properties of S, Se or Te doped perovskite ATiO ₃ (A=Ca, Ba, and Sr) materials for eco-friendly solar cells. <i>Superlattices and Microstructures</i> , 2022, 163, 107124.	1.4	6
35	Earth-abundant nontoxic ternary calcium nitrides inverse perovskites for single-junction solar cells: Ab-initio simulations. <i>Materials Science in Semiconductor Processing</i> , 2022, 150, 106959.	1.9	6
36	Ab Initio Study of Electronic and Magnetic Properties of Ga _{1-x} Co _x N (Doped) and Ga _{1-x-y} Co _x Cr _y N (Co-doped). <i>Journal of Superconductivity and Novel Magnetism</i> , 2017, 30, 165-170.	0.8	5

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37	Phase diagrams of spin $\hat{A}^{1/2}$ Ashkinâ€“Teller model with Dzyaloshinskiiâ€“Moriya interaction. Journal of Magnetism and Magnetic Materials, 2015, 394, 27-31.	1.0	4
38	RKKY Interactions in a Bilayer Olympicene Structure: A Monte Carlo Study. Journal of Superconductivity and Novel Magnetism, 2016, 29, 2793-2798.	0.8	4
39	Theoretical investigation of electronic and optical properties of the $\text{CuIn}_{1-x}\text{Ga}_x\text{Se}_2$: Ab initio calculation. Optik, 2020, 207, 163881.	1.4	4
40	Electronic and Magnetic Properties of ZnO Doped and Co-doped with (Co, Cr). Journal of Superconductivity and Novel Magnetism, 2016, 29, 3167-3173.	0.8	3
41	Superlattice Film with Ferromagnetic and Antiferromagnetic Layers Under the Effect of RKKY Interactions: a Monte Carlo Study. Journal of Superconductivity and Novel Magnetism, 2016, 29, 1887-1892.	0.8	3
42	Phase diagrams of 2D Ashkinâ€“Teller model within the effect of crystal field and quantum transverse field. Physica A: Statistical Mechanics and Its Applications, 2018, 492, 2310-2315.	1.2	3
43	Electronic, transport and optical properties in perovskite compound LaGaO_3 . Materials Research Express, 2020, 7, 035501.	0.8	3
44	Doping effect of iodine on electronic and optical properties of perovskite CsPbBr_3 compound for photovoltaic applications: Ab initio calculations. Journal of Electron Spectroscopy and Related Phenomena, 2021, 247, 147043.	0.8	3
45	Magnetic properties of a tri-decorated graphene structure: Monte Carlo study. International Journal of Modern Physics B, 2016, 30, 1650233.	1.0	2
46	Theoretical investigation of physical properties of the spinel ZnFe_2O_4 compound: Ab-initio calculation. Phase Transitions, 2021, 94, 134-146.	0.6	2
47	Magnetic properties and large magnetocaloric effect in the perovskite Mn_3GeC compound: Ab initio and Monte Carlo calculations. Phase Transitions, 2022, 95, 10-18.	0.6	2
48	Magnetic, magnetocaloric and thermoelectric properties of the intermetallic LaMn_2Si_2 compound: a theoretical study. Phase Transitions, 2022, 95, 387-397.	0.6	2
49	Ferromagnetism and Anti-ferromagnetism in Nano-films with Alternate Crystal Fields: Monte Carlo Study. Journal of Superconductivity and Novel Magnetism, 2016, 29, 2829-2833.	0.8	1
50	Effect of anisotropic Dzyaloshinskiiâ€“Moriya interactions on phase diagrams of the Ashkinâ€“Teller model. Physica A: Statistical Mechanics and Its Applications, 2016, 455, 92-97.	1.2	1
51	Ferrimagnetism in a Dendrimer Structure. Journal of Superconductivity and Novel Magnetism, 2016, 29, 375-381.	0.8	1
52	Ab Initio Study of Electronic and Magnetic Properties in ZnO-Doped and Co-doped by Vanadium and Silver. Journal of Superconductivity and Novel Magnetism, 2018, 31, 2201-2206.	0.8	1
53	Strain effect on physical properties of the multiferroic Mn_3Sn material: a first-principles calculations. Philosophical Magazine, 2022, 102, 1305-1319.	0.7	1
54	Rheology of particulate suspensions in a Poiseuille flow. Physical Review E, 2010, 82, 026306.	0.8	0

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55	Magnetic properties of a Lie symmetry double square nanostructure: Monte Carlo study. Phase Transitions, 2016, 89, 1006-1018.	0.6	0