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49
papers

839
citations

19
h-index

27
g-index

49
ext. papers

1,111
ext. citations

2.7
avg, IF

4.99
L-index

#	Paper	IF	Citations
49	Theoretical investigation of structural, elastic, and electronic properties of ternary boride MoAlB. <i>Physica Status Solidi (B): Basic Research</i> , 2017 , 254, 1700010	1.3	79
48	First hafnium-based MAX phase in the 312 family, Hf ₃ AlC ₂ : A first-principles study. <i>Journal of Alloys and Compounds</i> , 2017 , 727, 616-626	5.7	62
47	Recently synthesized (Zr _{1-x} Ti _x) ₂ AlC (0 ≤ x ≤ 1) solid solutions: Theoretical study of the effects of M mixing on physical properties. <i>Journal of Alloys and Compounds</i> , 2018 , 743, 146-154	5.7	45
46	Newly synthesized Zr ₂ AlC, Zr ₂ (Al _{0.58} Bi _{0.42})C, Zr ₂ (Al _{0.2} Sn _{0.8})C, and Zr ₂ (Al _{0.3} Sb _{0.7})C MAX phases: A DFT based first-principles study. <i>Computational Materials Science</i> , 2017 , 131, 139-145	3.2	43
45	Structural, morphological and electrical properties of Sn-substituted Ni-Zn ferrites synthesized by double sintering technique. <i>Journal of Magnetism and Magnetic Materials</i> , 2017 , 424, 148-154	2.8	42
44	Effects of transition metals on physical properties of M ₂ BC (M = V, Nb, Mo and Ta): A DFT calculation. <i>Journal of Alloys and Compounds</i> , 2019 , 770, 523-534	5.7	31
43	First-Principles Study of Superconducting ScRhP and ScIrP pnictides. <i>Physica Status Solidi (B): Basic Research</i> , 2017 , 254, 1700336	1.3	31
42	First principles study of M ₂ InC (M = Zr, Hf and Ta) MAX phases: The effect of M atomic species. <i>Results in Physics</i> , 2018 , 11, 869-876	3.7	30
41	Predicted MAX Phase Sc ₂ InC: Dynamical Stability, Vibrational and Optical Properties. <i>Physica Status Solidi (B): Basic Research</i> , 2018 , 255, 1700235	1.3	28
40	Structural Properties, Impedance Spectroscopy and Dielectric Spin Relaxation of Ni-Zn Ferrite Synthesized by Double Sintering Technique. <i>Journal of Scientific Research</i> , 2015 , 7, 65-75	1.4	28
39	Ni-rich Nitrides ANNi ₃ (A = Pt, Ag, Pd) in Comparison with Superconducting ZnNNi ₃ . <i>Journal of Scientific Research</i> , 2011 , 4, 1	1.4	28
38	An ab initio investigation of vibrational, thermodynamic, and optical properties of Sc ₂ AlC MAX compound. <i>Chinese Physics B</i> , 2016 , 25, 103102	1.2	28
37	Comparative study of Mo ₂ Ga ₂ C with superconducting MAX phase Mo ₂ GaC: First-principles calculations. <i>Chinese Physics B</i> , 2017 , 26, 033102	1.2	24
36	Sulvanite Compounds Cu ₃ TMS ₄ (TM = V, Nb and Ta): Elastic, Electronic, Optical and Thermal Properties using First-principles Method. <i>Journal of Scientific Research</i> , 2014 , 6, 407-419	1.4	24
35	Recently synthesized (Ti Mo) ₂ AlC (0 ≤ x ≤ 0.20) solid solutions: deciphering the structural, electronic, mechanical and thermodynamic properties simulations.. <i>RSC Advances</i> , 2020 , 10, 31535-31546	3.7	22
34	Elastic, thermodynamic and optical behavior of V ₂ AC (A= Al, Ga) MAX phases. <i>Results in Physics</i> , 2017 , 7, 3634-3639	3.7	21
33	Physical properties of new MAX phase borides M ₂ SB (M = Zr, Hf and Nb) in comparison with conventional MAX phase carbides M ₂ SC (M = Zr, Hf and Nb): Comprehensive insights. <i>Journal of Materials Research and Technology</i> , 2021 , 11, 1000-1018	5.5	21

32	New MAX Phase Compound Mo ₂ TiAlC ₂ : First-principles Study. <i>Journal of Scientific Research</i> , 2016 , 8, 109-117	1.4	19
31	Ternary boride Hf ₃ PB ₄ : Insights into the physical properties of the hardest possible boride MAX phase. <i>Journal of Alloys and Compounds</i> , 2021 , 857, 158264	5.7	19
30	Tailoring the properties of Ni-Zn-Co ferrites by Gd ³⁺ substitution. <i>Journal of Magnetism and Magnetic Materials</i> , 2020 , 497, 165978	2.8	18
29	Enhanced dielectric properties of prospective Bi _{0.85} Gd _{0.15} Fe _{1-x} Ir _x O ₃ multiferroics. <i>Results in Physics</i> , 2019 , 12, 1653-1659	3.7	15
28	First-principles Study of Vickers Hardness and Thermodynamic Properties of Ti ₃ SnC ₂ Polymorphs. <i>Journal of Scientific Research</i> , 2015 , 7, 53-64	1.4	15
27	First-principles study of elastic, electronic, optical and thermoelectric properties of newly synthesized K ₂ Cu ₂ GeS ₄ chalcogenide. <i>Journal of Alloys and Compounds</i> , 2019 , 781, 37-46	5.7	15
26	Study of physical properties towards optimizing sintering temperature of Y-substituted Mg-Zn ferrites. <i>Results in Physics</i> , 2019 , 14, 102517	3.7	14
25	Sn _{1-x} BixO ₂ and Sn _{1-x} TaxO ₂ (0 ≤ x ≤ 0.75): A first-principles study. <i>Physica B: Condensed Matter</i> , 2012 , 407, 1020-1026	2.8	14
24	DFT insights into new B-containing 212 MAX phases: Hf ₂ AB ₂ (A = In, Sn). <i>Journal of Alloys and Compounds</i> , 2021 , 860, 158408	5.7	14
23	Structural, elastic, electronic and optical properties of Cu ₃ MTe ₄ (M = Nb, Ta) sulvanites: An ab-initio study. <i>International Journal of Modern Physics B</i> , 2016 , 30, 1650089	1.1	12
22	First-principles study of SnO under high pressure. <i>International Journal of Modern Physics B</i> , 2016 , 30, 1650228	1.1	10
21	Yttrium-substituted Mg ₂ Zn ferrites: correlation of physical properties with Yttrium content. <i>Journal of Materials Science: Materials in Electronics</i> , 2019 , 30, 13258-13270	2.1	9
20	Mechanical behavior, enhanced dc resistivity, energy band gap and high temperature magnetic properties of Y-substituted Mg ₂ Zn ferrites. <i>Materials Research Express</i> , 2020 , 7, 036101	1.7	9
19	Newly synthesized MAX phase ZrSeC: DFT insights into physical properties towards possible applications.. <i>RSC Advances</i> , 2021 , 11, 16892-16905	3.7	9
18	Origin of high hardness and optoelectronic and thermo-physical properties of boron-rich compounds B ₆ X (X = S, Se): A comprehensive study via DFT approach. <i>Journal of Applied Physics</i> , 2021 , 129, 175109	2.5	8
17	Magnetic properties of Sn-substituted Ni ₂ Zn ferrites synthesized from nano-sized powders of NiO, ZnO, Fe ₂ O ₃ , and SnO ₂ . <i>Chinese Physics B</i> , 2017 , 26, 077501	1.2	6
16	Newly Synthesized Ta-Based MAX Phase (Ta _{1-x} Hfx) ₄ AlC ₃ and (Ta _{1-x} Hfx) ₄ Al _{0.5} Sn _{0.5} C ₃ (0 ≤ x ≤ 0.25) Solid Solutions: Unravelling the Mechanical, Electronic, and Thermodynamic Properties. <i>Physica Status Solidi (B): Basic Research</i> , 2021 , 258, 2000307	1.3	6
15	Screen the thermomechanical and optical properties of the new ductile 314 MAX phase boride Zr ₃ CdB ₄ : A DFT insight. <i>Journal of Alloys and Compounds</i> , 2021 , 877, 160248	5.7	6

14	Dielectric and optical properties of Ni-doped LaFeO ₃ nanoparticles. <i>SN Applied Sciences</i> , 2019 , 1, 1	1.8	5
13	Influence of Se doping on recently synthesized NaInS ₂ -xSex solid solutions for potential thermo-mechanical applications studied via first-principles method. <i>Materials Today Communications</i> , 2021 , 26, 101988	2.5	5
12	Dynamical stability, vibrational, and optical properties of anti-perovskite A ₃ BX (Ti ₃ TlN, Ni ₃ SnN, and Co ₃ AlC) phases: A first principles study. <i>AIP Advances</i> , 2020 , 10, 095226	1.5	4
11	NaInX ₂ (X = S, Se) layered materials for energy harvesting applications: first-principles insights into optoelectronic and thermoelectric properties. <i>Journal of Materials Science: Materials in Electronics</i> , 2021 , 32, 3878-3893	2.1	4
10	Impact of Sn ⁴⁺ substitution in Mg ₂ Zn ferrites: Deciphering the structural, morphological, dielectric, electrical and magnetic properties. <i>Materials Chemistry and Physics</i> , 2021 , 263, 124357	4.4	3
9	Physical properties of predicted MAX phase borides Hf ₂ AB (A = Pb, Bi): A DFT insight. <i>Materials Today Communications</i> , 2021 , 27, 102411	2.5	3
8	Newly Synthesized Three-Dimensional Boron-Rich Chalcogenides BX (X = S and Se): Theoretical Characterization of the Physical Properties for Optoelectronic and Mechanical Applications.. <i>ACS Omega</i> , 2021 , 6, 33899-33913	3.9	2
7	Comparative study of predicted MAX phase Hf ₂ AlN with recently synthesized Hf ₂ AlC: a first principle calculations. <i>Indian Journal of Physics</i> , 1	1.4	2
6	Understanding the improvement of thermo-mechanical and optical properties of 212 MAX phase borides Zr ₂ AB ₂ (A = In, Tl). <i>Journal of Materials Research and Technology</i> , 2021 , 15, 2227-2241	5.5	2
5	DFT insights into the new Hf-based chalcogenide MAX phase Hf ₂ SeC. <i>Vacuum</i> , 2022 , 201, 111072	3.7	2
4	Impact of V substitution on the physical properties of Ni ₂ ZnCo ferrites: structural, magnetic, dielectric and electrical properties. <i>Materials Research Express</i> , 2021 , 8, 046102	1.7	1
3	Effect of sintering temperature on structural and magnetic properties of Ni _{0.6} Zn _{0.4} Fe ₂ O ₄ ferrite: synthesized from nanocrystalline powders. <i>Journal of Physics: Conference Series</i> , 2021 , 1718, 012013	0.3	1
2	Optoelectronic, thermodynamic and vibrational properties of intermetallic MgAl ₂ Ge ₂ : a first-principles study. <i>SN Applied Sciences</i> , 2021 , 3, 1	1.8	0
1	Mechanical, optical and high-temperature magnetic properties of Sn-substituted Mg ₂ Zn ferrites. <i>Phase Transitions</i> , 2021 , 94, 23-36	1.3	