

# Abdul Hadi

## List of Publications by Year in descending order

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Version: 2024-02-01

51  
papers

1,960  
citations

186265

28  
h-index

254184

43  
g-index

51  
all docs

51  
docs citations

51  
times ranked

724  
citing authors

#	ARTICLE	IF	CITATIONS
1	Optical response, lithiation and charge transfer in Sn-based 211 MAX phases with electron localization function. <i>Journal of Materials Research and Technology</i> , 2022, 18, 2470-2479.	5.8	13
2	Indirect to direct band gap transition through order to disorder transformation of Cs <sub>2</sub> AgBiBr <sub>6</sub> via creating antisite defects for optoelectronic and photovoltaic applications. <i>RSC Advances</i> , 2022, 12, 15461-15469.	3.6	21
3	Structural, mechanical, thermal, and optical properties of inverse-Heusler alloys Cr <sub>2</sub> CoZ (Z = Al, In): A first-principles investigation. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2021, 385, 126967.	2.1	39
4	Insights into the physical properties of a new 211 MAX phase Nb <sub>2</sub> CuC. <i>Journal of Physics and Chemistry of Solids</i> , 2021, 149, 109759.	4.0	26
5	Insights into the physical properties and anisotropic nature of ErPdBi with an appearance of low minimum thermal conductivity. <i>Chinese Physics B</i> , 2021, 30, 016203.	1.4	16
6	Effects of Al substitution by Si in Ti <sub>3</sub> AlC <sub>2</sub> nanolaminate. <i>Scientific Reports</i> , 2021, 11, 3410.	3.3	27
7	Effect of boron incorporation into the carbon-site in Nb <sub>2</sub> SC MAX phase: Insights from DFT. <i>Journal of Materials Research and Technology</i> , 2021, 11, 1969-1981.	5.8	36
8	Changes in the physical properties of Mg <sub>2</sub> Ir <sub>3</sub> Si compared to Li <sub>2</sub> IrSi <sub>3</sub> due to the substitution of Li by Mg and the exchange of positions between Ir and Si. <i>Solid State Communications</i> , 2021, 334-335, 114361.	1.9	0
9	Ab initio exploration of the structural, elastic, electronic and optical properties of a new layered perovskite-type oxyfluoride: CsSrNb <sub>2</sub> O <sub>6</sub> F. <i>Materials Science in Semiconductor Processing</i> , 2021, 131, 105890.	4.0	47
10	Effect of hydrostatic compression on physical properties of Li <sub>2</sub> TmSi <sub>3</sub> (Tm = Ir, Pt, Rh, Os) with ground-state optical features. <i>Journal of Physics and Chemistry of Solids</i> , 2021, 156, 110124.	4.0	5
11	Structural, electronic, mechanical, thermal, and optical properties of UIr <sub>3</sub> under pressure: A comprehensive DFT study. <i>AIP Advances</i> , 2021, 11, .	1.3	14
12	Pressure effects on the structural, elastic, magnetic and thermodynamic properties of Mn <sub>2</sub> AlC and Mn <sub>2</sub> SiC MAX phases. <i>Journal of Alloys and Compounds</i> , 2021, 885, 160998.	5.5	40
13	A first-principles study of a new ternary silicide superconductor Li <sub>2</sub> IrSi <sub>3</sub> . <i>Journal of Physics: Conference Series</i> , 2021, 1718, 012018.	0.4	0
14	A density functional theory approach to the effects of C and N substitution at the B-site of the first boride MAX phase Nb <sub>2</sub> SB. <i>Materials Today Communications</i> , 2021, 29, 102910.	1.9	6
15	DFT investigations into the physical properties of a MAB phase Cr <sub>4</sub> AlB <sub>4</sub> . <i>Journal of Alloys and Compounds</i> , 2020, 821, 153547.	5.5	15
16	Superconducting phases in a remarkable class of metallic ceramics. <i>Journal of Physics and Chemistry of Solids</i> , 2020, 138, 109275.	4.0	84
17	Elastic behaviour and radiation tolerance in Nb-based 211 MAX phases. <i>Materials Today Communications</i> , 2020, 25, 101499.	1.9	10
18	Comprehensive first-principles calculations on physical properties of ScV <sub>2</sub> Ga <sub>4</sub> and ZrV <sub>2</sub> Ga <sub>4</sub> in comparison with superconducting HfV <sub>2</sub> Ga <sub>4</sub> . <i>Materials Today Communications</i> , 2020, 24, 100935.	1.9	23

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19	Insights into the predicted Hf <sub>2</sub> Sn in comparison with the synthesized MAX phase Hf <sub>2</sub> SnC: A comprehensive study. Computational Condensed Matter, 2020, 24, e00485.	2.1	17
20	Chemically stable new MAX phase V <sub>2</sub> SnC: a damage and radiation tolerant TBC material. RSC Advances, 2020, 10, 43783-43798.	3.6	34
21	Physical properties of a novel boron-based ternary compound Ti <sub>2</sub> InB <sub>2</sub> . Materials Today Communications, 2020, 25, 101600.	1.9	14
22	Structural, elastic, thermal and lattice dynamic properties of new 321 MAX phases. Computational Materials Science, 2019, 170, 109144.	3.0	34
23	Mechanical behaviors, lattice thermal conductivity and vibrational properties of a new MAX phase Lu <sub>2</sub> SnC. Journal of Physics and Chemistry of Solids, 2019, 129, 162-171.	4.0	83
24	Electronic structures, bonding natures and defect processes in Sn-based 211 MAX phases. Computational Materials Science, 2019, 168, 203-212.	3.0	17
25	Phase stability and physical properties of (Zr <sub>1</sub> -Nb) <sub>2</sub> AlC MAX phases. Journal of Physics and Chemistry of Solids, 2019, 132, 38-47.	4.0	32
26	Influence of Ni doping in a lead-halide and a lead-free halide perovskites for optoelectronic applications. AIP Advances, 2019, 9, .	1.3	56
27	312 MAX Phases: Elastic Properties and Lithiation. Materials, 2019, 12, 4098.	2.9	20
28	Newly synthesized MgAl <sub>2</sub> Ge <sub>2</sub> : A first-principles comparison with its silicide and carbide counterparts. Journal of Physics and Chemistry of Solids, 2018, 117, 139-147.	4.0	24
29	Intrinsic defect processes and elastic properties of Ti <sub>3</sub> AlC <sub>2</sub> (A = Al, Si, Ga, Ge, In, Sn) MAX phases. Journal of Applied Physics, 2018, 123, .	2.5	31
30	Physical properties and defect processes of M <sub>3</sub> SnC <sub>2</sub> (M = Ti, Zr, Hf) MAX phases: Effect of M-elements. Journal of Alloys and Compounds, 2018, 748, 804-813.	5.5	49
31	Cubic Perovskite Pb(Mg <sub>1/3</sub> Nb <sub>2/3</sub> )O <sub>3</sub> : A Damage Tolerant, Machinable, and Thermal barrier coating material. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 2018, 74, 71-81.	1.5	33
32	Synthesis and physical properties of (Zr <sub>1-x</sub> Ti <sub>x</sub> ) <sub>3</sub> AlC <sub>2</sub> MAX phases. Journal of the American Ceramic Society, 2017, 100, 3393-3401.	3.8	63
33	Elastic and thermodynamic properties of new (Zr <sub>3-x</sub> Ti <sub>x</sub> )AlC <sub>2</sub> MAX-phase solid solutions. Computational Materials Science, 2017, 137, 318-326.	3.0	119
34	Theoretical investigation of structural, elastic, and electronic properties of ternary boride MoAlB. Physica Status Solidi (B): Basic Research, 2017, 254, 1700010.	1.5	113
35	Physical properties of ternary silicide superconductors Li <sub>2</sub> XSi <sub>3</sub> (X = Rh, Os): An ab initio study. International Journal of Modern Physics B, 2017, 31, 1750135.	2.0	5
36	New ternary superconducting compound LaRu <sub>2</sub> As <sub>2</sub> : Physical properties from density functional theory calculations. Chinese Physics B, 2017, 26, 037103.	1.4	39

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37	First-principles Study of Superconducting ScRhP and ScIrP pnictides. Physica Status Solidi (B): Basic Research, 2017, 254, 1700336.	1.5	39
38	Mechanical behavior, bonding nature and defect processes of Mo <sub>2</sub> ScAlC <sub>2</sub> : A new ordered MAX phase. Journal of Alloys and Compounds, 2017, 724, 1167-1175.	5.5	52
39	First hafnium-based MAX phase in the 312 family, Hf <sub>3</sub> AlC <sub>2</sub> : A first-principles study. Journal of Alloys and Compounds, 2017, 727, 616-626.	5.5	95
40	Density functional theory study of a new Bi-based (K <sub>1.00</sub> )(Ba <sub>1.00</sub> ) <sub>3</sub> (Bi <sub>0.89</sub> Na <sub>0.11</sub> ) <sub>4</sub> O <sub>12</sub> double perovskite superconductor. Computational Materials Science, 2017, 138, 160-165.	3.0	45
41	Structural and optical properties of the recently synthesized (Zr <sub>3-<math>\hat{x}</math></sub> Ti <sub><math>\hat{x}</math></sub> )AlC <sub>2</sub> MAX phases. Journal of Materials Science: Materials in Electronics, 2017, 28, 3386-3393.	2.2	27
42	First-principles prediction of mechanical and bonding characteristics of new T <sub>2</sub> superconductor Ta <sub>5</sub> GeB <sub>2</sub> . Physica Status Solidi (B): Basic Research, 2016, 253, 2020-2026.	1.5	61
43	Structural, Elastic, and Electronic Properties of Newly Discovered Li <sub>2</sub> PtSi <sub>3</sub> Superconductor: Effect of Transition Metals. Journal of Superconductivity and Novel Magnetism, 2016, 29, 2503-2508.	1.8	23
44	Physical properties of the recently discovered Zr <sub>2</sub> (Al <sub>1-<math>\hat{x}</math></sub> Bi <sub><math>\hat{x}</math></sub> )C MAX phases. Journal of Materials Science: Materials in Electronics, 2016, 27, 11925-11933.	2.2	71
45	New ordered MAX phase Mo <sub>2</sub> TiAlC <sub>2</sub> : Elastic and electronic properties from first-principles. Chinese Physics B, 2016, 25, 107103.	1.4	36
46	Physical properties of predicted Ti <sub>2</sub> CdN versus existing Ti <sub>2</sub> CdC MAX phase: An ab initio study. Computational Materials Science, 2016, 113, 148-153.	3.0	101
47	New ternary nanolaminated carbide Mo <sub>2</sub> Ga <sub>2</sub> C: A first-principles comparison with the MAX phase counterpart Mo <sub>2</sub> GaC. Computational Materials Science, 2016, 117, 422-427.	3.0	74
48	Structural, elastic, and electronic properties of recently discovered ternary silicide superconductor Li <sub>2</sub> IrSi <sub>3</sub> : An ab-initio study. Chinese Physics B, 2015, 24, 117401.	1.4	33
49	Zirconium metal-based MAX phases Zr <sub>2</sub> AC (A =) Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 International Journal of Modern Physics B, 2014, 28, 1550022.	2.0	41
50	New MAX Phase Superconductor Ti <sub>2</sub> GeC: A First-principles Study. Journal of Scientific Research, 2014, 6, 11-27.	0.3	29
51	BAND STRUCTURE, HARDNESS, THERMODYNAMIC AND OPTICAL PROPERTIES OF SUPERCONDUCTING Nb <sub>2</sub> AsC, Nb <sub>2</sub> InC AND Mo <sub>2</sub> GaC. International Journal of Computational Materials Science and Engineering, 2013, 02, 1350007.	0.7	28