

# Zahid Ali

## List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Structural and optoelectronic properties of CsLnZnTe <sub>3</sub> (Ln= La, Pr, Nd and Sm). Journal of Rare Earths, 2023, 41, 388-396.	4.8	4
2	Hybrid DFT study of structural, electronic, magnetic and elastic properties of laves phase binary intermetallics RFe <sub>2</sub> (R= La, Ce, Pr and Nd). Journal of Rare Earths, 2023, 41, 1367-1375.	4.8	7
3	Electronic structure and magnetic properties of the perovskites SrTMO <sub>3</sub> (TM = Mn, Fe, Co, Tc, Ru, Rh.) Tj ETQq1 1 0.784314 rgBT /Over	2.7	21
4	First-principles study of the structural and optoelectronic properties of ANbO <sub>3</sub> (A = Na, K and Rb) in four crystal phases. Materials Science in Semiconductor Processing, 2022, 139, 106364.	4.0	17
5	Structural, Electronic, Elastic and Magnetic Properties of Ln <sub>3</sub> QIn (Ln = Ce, Pr and Nd; Q = C and N) anti-perovskites. Journal of Electronic Materials, 2022, 51, 2819-2827.	2.2	3
6	New anti-ferromagnetic tri-transition quaternary perovskites for magnetic cloaking and spintronic applications. Materials Chemistry and Physics, 2022, 282, 125915.	4.0	9
7	Electronic structure, elastic and magnetic properties of the binary intermetallics RFe <sub>2</sub> (R=Eu, Gd and) Tj ETQq1 1 0.784314 rgBT /Over	2.5	3
8	First principle study of opto-electronic and thermoelectric properties of Zintl Phase XIn <sub>2</sub> Z <sub>2</sub> (X=Ca, Sr) Tj ETQq0 0 0 rgBT /Over	2.3	8
9	Optoelectronic, elastic and thermoelectric properties of the perovskites (Sr <sub>3</sub> N)Sb and (Sr <sub>3</sub> N)Bi. Materials Science in Semiconductor Processing, 2022, 147, 106734.	4.0	2
10	<sub>n</sub>-Type narrow band gap A<sub>3</sub>InAs<sub>3</sub> (A= Sr and Eu) Zintl phase semiconductors for optoelectronic and thermoelectric applications. Journal of Taibah University for Science, 2022, 16, 660-669.	2.5	10
11	Structural, electronic, optical and thermoelectric properties in the phases of AgTaO <sub>3</sub> . Materials Science in Semiconductor Processing, 2021, 122, 105467.	4.0	7
12	Elastic and Optoelectronic Properties of Cs <sub>2</sub> NaMCl <sub>6</sub> (M = In, Tl, Sb, Bi). Journal of Electronic Materials, 2021, 50, 456-466.	2.2	33
13	The effect of potassium insertion on optoelectronic properties of cadmium chalcogenides. Materials Science in Semiconductor Processing, 2021, 122, 105466.	4.0	7
14	Spin-orbit coupling effect on the optoelectronic and thermoelectric properties of the perovskites A <sub>3</sub> SnO (A = Ca, Sr and Ba). Materials Science in Semiconductor Processing, 2021, 132, 105905.	4.0	11
15	Electronic structure, optical and magnetic properties of double Perovskites La <sub>2</sub> MTiO <sub>6</sub> (M = Co, Ni, Cu) Tj ETQq1 1 0.784314 rgBT /Over	4.0	15
16	Electronic Structure, Mechanical and Magnetic Properties of the Quaternary Perovskites CaA <sub>3</sub> V <sub>4</sub> O <sub>12</sub> (A=Mn, Fe, Co, Ni and Cu). Journal of Electronic Materials, 2020, 49, 1230-1242.	2.2	12
17	Electronic structure and magnetic properties of the Mg-rich intermetallic NdNiMg <sub>5</sub> by hybrid density functional theory. Intermetallics, 2020, 127, 106969.	3.9	1
18	Electronic structure and magnetic properties of the quaternary perovskites LnMn<sub>3</sub>V<sub>4</sub>O<sub>12</sub> (Ln=La, Nd and Gd). Philosophical Magazine, 2020, 100, 2386-2401.	5	16

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19	Effects of A-Site cation on the Physical Properties of Quaternary Perovskites AMn <sub>3</sub> V <sub>4</sub> O <sub>12</sub> (A= Ca, Ce) Tj ETQq1 1 0.784314 rgBT /Overall	4.0	14
20	Theoretical Investigations of Quaternary Semiconductors CsInCdTe <sub>3</sub> (Ln <sub>4</sub> =La, Pr, Nd and Sm). Journal of Electronic Materials, 2020, 49, 3357-3366.	2.2	16
21	Effects of Ni Substitution on the Electronic Structure and Magnetic Properties of Perovskite SrFeO <sub>3</sub> . Journal of Electronic Materials, 2020, 49, 3780-3790.	2.2	19
22	Optoelectronic properties of the double perovskites Ba <sub>2</sub> MM <sup>2+</sup> O <sub>6</sub> (M= Sc, Y, La; M <sup>2+</sup> = Nb, Ta) by modified Becke-Johnson potential. , 2020, , .	0	
23	Structural, optoelectronic and elastic properties of quaternary perovskites CaPd <sub>3</sub> B <sub>4</sub> O <sub>12</sub> (B = Ti, V). International Journal of Modern Physics B, 2019, 33, 1950212.	2.0	6
24	First-principles study of BiFeO <sub>3</sub> and BaTiO <sub>3</sub> in tetragonal structure. International Journal of Modern Physics B, 2019, 33, 1950231.	2.0	40
25	First principles studies of CsLnCdTe <sub>3</sub> (Ln <sup>3+</sup> = Gd <sup>3+</sup> Tm) for green energy resources. Computational Condensed Matter, 2019, 21, e00427.	2.1	7
26	First-Principles Study of Perovskite Molybdates AMoO <sub>3</sub> (A <sub>2</sub> =Ca, Sr, Ba). Journal of Electronic Materials, 2019, 48, 1730-1739.	2.2	23
27	First principle studies of structural, magnetic and elastic properties of orthorhombic rare-earth diaurides intermetallics RAu <sub>2</sub> (R=La, Ce, Pr and Eu). Materials Chemistry and Physics, 2018, 212, 44-50.	4.0	11
28	First-Principles Study of Electronic Structure, Mechanical, and Thermoelectric Properties of Ternary Palladates CdPd <sub>3</sub> O <sub>4</sub> and TiPd <sub>3</sub> O <sub>4</sub> . Journal of Electronic Materials, 2018, 47, 1871-1880.	2.2	4
29	HKUST-1 Supported on Zirconium Phosphate as an Efficient Catalyst for Solvent Free Oxidation of Cyclohexene: DFT Study. Catalysts, 2018, 8, 546.	3.5	3
30	Theoretical studies of the electronic structure and magnetic properties of aluminum-rich intermetallic alloy Al <sub>13</sub> Fe <sub>4</sub> . International Journal of Modern Physics B, 2018, 32, 1850201.	2.0	4
31	Effects of cobalt substitution on the physical properties of the perovskite strontium ferrite. Materials Chemistry and Physics, 2017, 196, 222-228.	4.0	26
32	First-principle studies of the optoelectronic properties of ASnF <sub>3</sub> (A = Na, K, Rb and Cs). International Journal of Modern Physics B, 2017, 31, 1750148.	2.0	18
33	Structural, Mechanical and Optoelectronic Properties of Y <sub>2</sub> M <sub>2</sub> O <sub>7</sub> (M = Ti, V and Nb) Pyrochlores: A First Principles Study. Journal of Electronic Materials, 2017, 46, 4640-4648.	2.2	5
34	Electron correlation and spin-orbit coupling effects in scandium intermetallic compounds ScTM (TM) Tj ETQq0 0 0 rgBT /Overall	2.0	5 Tf 5
35	First principle studies of electronic and magnetic properties of Lanthanide-Gold (RAu) binary intermetallics. Journal of Magnetism and Magnetic Materials, 2017, 422, 458-463.	2.3	13
36	The Influence of Oxygen Substitution on the Optoelectronic Properties of ZnTe. Journal of Chemistry, 2016, 2016, 1-8.	1.9	1

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37	Electronic structure of the LiAA <sub>x</sub> O <sub>6</sub> (A=Nb, Ta, and W, Mo) ceramics by modified Becke-Johnson potential. <i>Optical Materials</i> , 2016, 58, 466-475.	3.6	10
38	First-principle studies of the ternary palladates CaPd <sub>3</sub> O <sub>4</sub> and SrPd <sub>3</sub> O <sub>4</sub> . <i>Bulletin of Materials Science</i> , 2016, 39, 1861-1870.	1.7	6
39	First-principles studies of pure and fluorine substituted alanines. <i>International Journal of Modern Physics B</i> , 2016, 30, 1650079.	2.0	6
40	Thermoelectric studies of IV-VI semiconductors for renewable energy resources. <i>Materials Science in Semiconductor Processing</i> , 2016, 48, 85-94.	4.0	58
41	Theoretical studies of the osmium based perovskites AOsO <sub>3</sub> (A=Ca, Sr and Ba). <i>Journal of Physics and Chemistry of Solids</i> , 2015, 86, 114-121.	4.0	22
42	Theoretical studies of the paramagnetic perovskites MTaO <sub>3</sub> (M=Ca, Sr and Ba). <i>Materials Chemistry and Physics</i> , 2015, 162, 308-315.	4.0	38
43	Structural and optoelectronic properties of Mg substituted ZTe (Z=Zn, Cd and Hg). <i>Journal of Physics and Chemistry of Solids</i> , 2015, 83, 75-84.	4.0	32
44	Structural and magnetic properties of TiTF <sub>3</sub> (T=Fe, Co and Ni) by hybrid functional theory. <i>Journal of Magnetism and Magnetic Materials</i> , 2015, 388, 143-149.	2.3	16
45	Structural, Mechanical and Magneto-Electronic Properties of the Ternary Sodium Palladium and Platinum Oxides. <i>Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences</i> , 2015, 70, 815-822.	1.5	1
46	Magneto-electronic studies of the inverse-perovskite (Eu <sub>3</sub> O)In. <i>Journal of Magnetism and Magnetic Materials</i> , 2015, 381, 34-40.	2.3	5
47	Magneto-electronic studies of anti-perovskites NiNMn <sub>3</sub> and ZnNMn <sub>3</sub> . <i>Computational Materials Science</i> , 2014, 81, 141-145.	3.0	14
48	Density functional studies of magneto-optic properties of CdCoS. <i>Journal of Magnetism and Magnetic Materials</i> , 2014, 351, 60-64.	2.3	10
49	Structural and optoelectronic properties of the zinc titanate perovskite and spinel by modified Becke-Johnson potential. <i>Physica B: Condensed Matter</i> , 2013, 420, 54-57.	2.7	44
50	Band Profile Comparison of the Cubic Perovskites CaCoO <sub>3</sub> and SrCoO <sub>3</sub> . <i>Journal of Electronic Materials</i> , 2013, 42, 438-444.	2.2	27
51	Robust Half-Metallicity and Magnetic Properties of Cubic Perovskite CaFeO <sub>3</sub> . <i>Chinese Physics Letters</i> , 2013, 30, 047504.	3.3	15
52	First principle optoelectronic studies of visible light sensitive CZT. <i>Superlattices and Microstructures</i> , 2013, 63, 91-99.	3.1	15
53	Comparison of the electronic band profiles and magneto-optic properties of cubic and orthorhombic SrTbO <sub>3</sub> . <i>Physica B: Condensed Matter</i> , 2013, 423, 16-20.	2.7	15
54	Conversion of optically isotropic to anisotropic CdS <sub>x</sub> Se <sub>1-x</sub> (0<x<1) alloy with S concentration. <i>Computational Materials Science</i> , 2013, 77, 145-152.	3.0	48

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55	GGA+U studies of the cubic perovskites BaMO <sub>3</sub> (M=Pr, Th and U). <i>Physica B: Condensed Matter</i> , 2013, 410, 217-221.	2.7	41
56	Comparison of band profiles and magnetic properties of the different phases of BaTbO <sub>3</sub> . <i>Computational Materials Science</i> , 2013, 67, 151-155.	3.0	9
57	Electronic structure of cubic perovskite SnTaO <sub>3</sub> . <i>Intermetallics</i> , 2012, 31, 287-291.	3.9	55
58	Theoretical studies of structural and magnetic properties of cubic perovskites PrCoO <sub>3</sub> and NdCoO <sub>3</sub> . <i>Physica B: Condensed Matter</i> , 2011, 406, 3800-3804.	2.7	48
59	Bandgap engineering of Cd <sub>1-x</sub> SrxO. <i>Physica B: Condensed Matter</i> , 2011, 406, 2509-2514.	2.7	33