

Rehan Ullah

List of Publications by Year in descending order

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14
papers

285
citations

840585

11
h-index

1058333

14
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docs citations

14
times ranked

125
citing authors

#	ARTICLE	IF	CITATIONS
1	Exploring electronic, structural, magnetic and thermoelectric properties of novel Ba ₂ EuMoO ₆ double perovskite. <i>Materials Science in Semiconductor Processing</i> , 2022, 137, 106218.	1.9	10
2	Effect of cation exchange on structural, electronic, magnetic and transport properties of Ba ₂ MReO ₆ (M = Al, Gd). <i>Journal of Magnetism and Magnetic Materials</i> , 2022, 546, 168816.	1.0	14
3	Modeling of bulk modulus of A ₂ BX ₆ cubic crystals (A = K, Cs, Rb, Tl, NH ₄ ; B = tetravalent cation; X = F.) <i>Tj ETQq1 1,0,784314,rgBT /O</i>	1.8	7
4	Influence of the spin-orbit coupling effect on the electronic and thermoelectric properties of Cs ₂ Ml ₆ (M = Zr, Hf) variant perovskites. <i>Materials Research Bulletin</i> , 2021, 134, 111112.	2.7	11
5	An investigation of half-metallic variant perovskites A ₂ NbCl ₆ (A = K, Rb) for spintronic based applications. <i>Journal of Solid State Chemistry</i> , 2021, 293, 121823.	1.4	14
6	An investigation of structural, elastic, mechanical, electronic, magnetic and thermoelectric properties of ferromagnetic half metallic EuCrO ₃ . <i>Materials Science in Semiconductor Processing</i> , 2021, 122, 105487.	1.9	12
7	Pressure dependent elastomechanical stability and thermoelectric properties of A ₃ MYbF ₃ (M = Rb, Cs) materials for renewable energy. <i>International Journal of Energy Research</i> , 2021, 45, 8711-8723.	2.2	48
8	Spin-based transport properties of Cs ₂ WX ₆ (X = Cl, Br) ferromagnets for spin-injected thermoelectric current. <i>European Physical Journal Plus</i> , 2021, 136, 1.	1.2	12
9	Pressure induced structural, electronic, optical and thermal properties of CsYbBr ₃ , a theoretical investigation. <i>Journal of Materials Research and Technology</i> , 2021, 10, 687-696.	2.6	13
10	Modeling of structural, elastic, mechanical, acoustical, electronic and thermodynamic properties of XPdF ₃ (X = Rb, Tl) perovskites through density functional theory. <i>Physica Scripta</i> , 2020, 95, 075705.	1.2	24
11	Ab initio study for the structural, electronic, magnetic, optical, and thermoelectric properties of K ₂ OsX ₆ (X = Cl, Br) compounds. <i>International Journal of Energy Research</i> , 2020, 44, 9035-9049.	2.2	36
12	Insight into pressure tunable structural, electronic and optical properties of "Equation missing" via DFT calculations. <i>European Physical Journal Plus</i> , 2020, 135, 1.	1.2	32
13	The significance of anti-fluorite Cs ₂ NbI ₆ via its structural, electronic, magnetic, optical and thermoelectric properties. <i>International Journal of Energy Research</i> , 2020, 44, 10179-10191.	2.2	25
14	Structural, electronic and optical properties of cubic perovskite RbYbF ₃ under pressure: a first principles study. <i>Materials Research Express</i> , 2019, 6, 125901.	0.8	27