fethallah Dahmane

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Insight into the Structural, Magneto-electronic, and Mechanical Characteristic of Y2MnZ (Z = Al, Ga,) Tj	ет <u>о</u> я110).7 <u>8</u> 4314 rg
2	First-principles study of half-metallic properties in X2CrAl (XÂ=ÂCo and Mn) FullHeusler and their quaternary MnCoCrAl and CoMnCrAl compounds. Inorganic Chemistry Communication, 2022, 139, 109408.	3.9	3
3	Structural properties, electronic band structure, magnetic and mechanical characteristic of XFeGe (XÂ=ÂCo, Cr, Ni) half Heusler compounds: Insights from DFT calculation. Inorganic Chemistry Communication, 2022, 142, 109675.	3.9	1
4	Insight view of Hf2CrZ (ZÂ= B, Ga, In, Si, Ge, Sn) Heusler materials via DFT calculations: A study on structural, electronic and magnetic properties. Computational Condensed Matter, 2021, 26, e00518.	2.1	4
5	First-principles investigation of half-metallic ferromagnetism of Fe2YSn (Y = Mn, Ti and V) Heusler alloys. Condensed Matter Physics, 2021, 24, 23703.	0.7	2
6	Prediction of electronic and half metallic properties of Mn2YSn (Y = Mo, Nb, Zr) Heusler alloys. Condensed Matter Physics, 2021, 24, 13703.	0.7	0
7	Structural, electronic, magnetic and mechanical properties of the full-Heusler compounds Ni ₂ Mn(Ge,Sn) and Mn ₂ NiGe. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 2021, 76, 693-702.	1.5	5
8	Theoretical Characterization of Thermodynamic, Magnetic and Electronic Proprieties of Full-Heusler Co ₂ YGa (Y = V, Cr and Mn) Alloys. Spin, 2020, 10, .	1.3	4
9	Phase stability, mechanical, electronic and thermodynamic properties of the Ga3Sc compound: An ab-initio study. Inorganic Chemistry Communication, 2020, 122, 108304.	3.9	17
10	Structural stability, mechanical, electronic and thermal behaviour of Ru2CrZ (Z=Sb, Si, Pb, Ge) Heusler alloys. Chinese Journal of Physics, 2020, 66, 124-134.	3.9	6
11	Ab initio study of structural, electronic and magnetic properties of XSn3 (X = Gd, Cm) and GdxCm1-xSn3 compounds. Condensed Matter Physics, 2020, 23, 33705.	0.7	0
12	Theoretical study of phase stability, electronic and magnetic properties of Rh2CrGe1-xAlx (x = 0, 0.25,) Tj ETQq0	0 0.rgBT /0	Overlock 10
13	First-principles study of half-metallic properties in XVSi (XÂ= Ti, Co) and their quaternary TiCoVSi and CoTiVSi compounds. Computational Condensed Matter, 2019, 19, e00369.	2.1	8
14	A comparative study between Hg2CuTi and Cu2MnAl type structures for Zr2CoZ (ZÂ=ÂAl, Ga, In) Heusler alloys. Chinese Journal of Physics, 2019, 60, 450-461.	3.9	16
15	Structural, electronic and magnetic properties of new full Heusler alloys Rh2CrZ (Z = Al, Ga, In): First-principles calculations. Chinese Journal of Physics, 2019, 59, 281-290.	3.9	26
16	First-principle calculations of structural, electronic and magnetic investigations of Mn2RuGe1-xSnx quaternary Heusler alloys. Chinese Journal of Physics, 2018, 56, 567-573.	3.9	72
17	Ab Initio Investigation of Structural Stability and Electronic and Magnetic Properties of the Half-Heusler Alloys: MTiSb (M = Fe, Co, and Ni). Journal of Superconductivity and Novel Magnetism, 2018, 31, 2991-2998.	1.8	8
18	The effect of Sn substitution on the Al sites in full Heusler compound Fe2VAl. Indian Journal of Physics, 2018, 92, 1403-1411.	1.8	1

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19	First-principle study of the electronic, magnetic and structural characteristics of the Mn2CoAs1â^'xAlx (x =â€0,0.25,0.50,0.75) Heusler alloys. Chinese Journal of Physics, 2018, 56, 1764-1771.	3.9	3
20	Structural, elastic and optoelectronic properties of Sr-based perovskite-type oxides SrXO3 (Mâ€=†Th, Zr) via first-principles calculations. Chinese Journal of Physics, 2018, 56, 1515-1524.	3.9	10
21	Theoretical study of the structural stability, electronic and magnetic properties of XVSb (X = Fe, Ni,) Tj ETQq1 1	0.784314 0.7	rgBT /Overloc
22	Electronic and Ferromagnetic Properties of 3d(V)-Doped (BaS) Barium Sulfide. Journal of Superconductivity and Novel Magnetism, 2017, 30, 917-923.	1.8	25
23	Ab initio study of structural and electronic properties of (GaN)n/(AlN)n superlattices. Journal of Physics: Conference Series, 2016, 758, 012025.	0.4	2
24	First-principle calculations of electronic and ferromagnetic properties of \$\$hbox {Al}_{1-x}hbox {V}_{x}hbox {Sb}\$\$ Al 1 - x V x Sb. Journal of Computational Electronics, 2016, 15, 1255-1262.	2.5	11
25	Band Structure Simulations of the Structural, Electronic, Magnetic, and Half-Metallic Features of the Ti 2 CoAl1â [~] x Sn x (x = 0, 0.25, 0.50, 0.75, 1) Heusler Alloys. Journal of Superconductivity and Novel Magnetism, 2016, 29, 3193-3199.	1.8	3
26	Theoretical investigation of the structural, magnetic and band structure characteristics of Co2FeGe1â^'x Si x (x = 0, 0.5, 1) full-Heusler alloys. Journal of the Korean Physical Society, 2016, 69, 1462-1468.	0.7	22
27	Elastic and electronic properties calculations of the filled skutterudite CeOs ₄ P ₁₂ . Journal of Physics: Conference Series, 2016, 758, 012010.	0.4	3
28	Structural, electronic and magnetic properties of Fe2-based full Heusler alloys: A first principle study. Journal of Magnetism and Magnetic Materials, 2016, 407, 167-174.	2.3	67
29	Investigations of the Structural, Electronic, Magnetic, and Half-Metallic Behavior of Co2MnZ (Z = Al,) Tj ETQq1 1 809-817.	0.78431 1.8	4 rgBT /Overlo 41
30	A Novel Theoretical Investigation of Electronic Structure and Half-Metallic Ferromagnetism in 3d (V)-Doped InP for Spintronic Applications. Journal of Superconductivity and Novel Magnetism, 2016, 29, 1813-1819.	1.8	8
31	Electronic structure, magnetism and stability of Co2CrX (X =Al, Ga, In) ab initio study. Modern Physics Letters B, 2016, 30, 1550265.	1.9	20
32	Half-Metallic Ferromagnetic Property Related to Spintronic Applications in 3d (V, Cr, and Mn)-Doped GaP DMSs. Journal of Superconductivity and Novel Magnetism, 2015, 28, 3163-3172.	1.8	18
33	First-Principles Study of Structural, Electronic, Magnetic and Half-Metallic Properties of the Heusler Alloys Ti2ZAl (Z = Co, Fe, Mn). Journal of Superconductivity and Novel Magnetism, 2015, 28, 3099-3104.	1.8	22
34	Ab Initio Investigation of Half-Metallic Behaviour in the Full-Heusler X2MnGe (X = Sc, Fe, Ni). Journal of Superconductivity and Novel Magnetism, 2015, 28, 2063-2069.	1.8	6
35	First-principle investigation of half-metallic ferromagnetism in octahedrally bonded Cr-doped rock-salt SrS, SrSe, and SrTe. European Physical Journal B, 2015, 88, 1.	1.5	41
36	A novel theoretical design of electronic structure and half-metallic ferromagnetism in the 3d (V)-doped rock-salts SrS, SrSe, and SrTe for spintronics. RSC Advances, 2015, 5, 92328-92334.	3.6	32

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37	First-principles investigations on ferromagnetic behaviour of Be1â^'xVxZ (ZÂ=ÂS, Se and Te) (xÂ=Â0.25). Superlattices and Microstructures, 2015, 88, 139-149.	3.1	24
38	First-Principles Investigation of Half-metallic Ferromagnetism in V-doped BeS, BeSe, and BeTe. Journal of Superconductivity and Novel Magnetism, 2014, 27, 293-300.	1.8	42
39	First principles study of the electronic structures and magnetic properties of transition metal-doped cubic indium nitride. Materials Science in Semiconductor Processing, 2014, 21, 66-73.	4.0	21
40	First-Principle Calculations of Structural, Electronic, and Magnetic Properties of Cubic Al1â^'x TM x N (TM = V, Cr, Mn, Fe). Journal of Superconductivity and Novel Magnetism, 2014, 27, 2647-2654.	1.8	11
41	First-Principle Investigations of Structural, Electronic, and Half-Metallic Ferromagnetic Properties in In1â~'xTM x P (TM = Cr, Mn). Journal of Superconductivity and Novel Magnetism, 2014, 27, 1603-1614.	1.8	20
42	Structural, Electronic and Magnetic Properties of Zinc-Blende Ga1â^'x TM x N (TM = Cr, Mn, Fe, V). Journal of Superconductivity and Novel Magnetism, 2013, 26, 3339-3348.	1.8	15
43	Investigations of Structural, Electronic, and Half-metallic Ferromagnetic Properties in (Al, Ga, In)1â^'x M x N (M = Fe, Mn) Diluted Magnetic Semiconductors. Journal of Superconductivity and Novel Magnetism, 2013, 26, 515-525.	1.8	41