

Yarub Al-Douri

List of Publications by Year in descending order

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287
papers

9,884
citations

34105

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all docs

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docs citations

295
times ranked

4713
citing authors

#	ARTICLE	IF	CITATIONS
1	XPS and optical studies of different morphologies of ZnO nanostructures prepared by microwave methods. <i>Ceramics International</i> , 2013, 39, 2283-2292.	4.8	819
2	Structural, magnetic, electronic and mechanical properties of full-Heusler alloys Co ₂ YAl (Y = Fe, Ti): First principles calculations with different exchange-correlation potentials. <i>Journal of Magnetism and Magnetic Materials</i> , 2018, 448, 208-220.	2.3	245
3	First-principles calculations to investigate magnetic and thermodynamic properties of new multifunctional full-Heusler alloy Co ₂ TaGa. <i>Indian Journal of Physics</i> , 2020, 94, 767-777.	1.8	169
4	FP-APW+lo calculations of the elastic properties in zinc-blende III-P compounds under pressure effects. <i>Computational Materials Science</i> , 2009, 45, 474-479.	3.0	137
5	Electronic and thermoelectric properties of the layered BaF _{Ag} Ch (Ch = S, Se and Te): First-principles study. <i>Journal of Alloys and Compounds</i> , 2018, 759, 32-43.	5.5	136
6	First-Principles Calculations to Investigate Structural, Electronic, Elastic, Magnetic, and Thermodynamic Properties of Full-Heusler Rh ₂ MnZ (Z = Zr, Hf). <i>Journal of Superconductivity and Novel Magnetism</i> , 2021, 34, 269-283.	1.8	132
7	Structural, electronic, optical and thermodynamic investigations of NaXF ₃ (X = Ca and Sr): First-principles calculations. <i>Chinese Journal of Physics</i> , 2018, 56, 131-144.	3.9	125
8	First-principles calculations to investigate structural and thermodynamic properties of Ni ₂ LaZ (Z = As, Sb, Bi). <i>Journal of Superconductivity and Novel Magnetism</i> , 2021, 34, 121-127.	1.8	121
9	First-principles calculations of pressure and temperature dependence of thermodynamic properties of anti-perovskite BiNBa ₃ compound. <i>Chinese Journal of Physics</i> , 2017, 55, 2144-2155.	3.9	118
10	First-principles predictions of the structural, electronic, optical and elastic properties of the zintl-phases AE ₃ GaAs ₃ (AE = Sr, Ba). <i>Solid State Sciences</i> , 2021, 114, 106563.	3.2	118
11	Structural, elastic, thermodynamic and electronic properties of LuX (X = N, Bi and Sb) compounds: first principles calculations. <i>Phase Transitions</i> , 2016, 89, 1236-1252.	1.3	115
12	Electronic, optical and thermoelectric investigations of Zintl phase AE ₃ AlAs ₃ (AE = Sr, Ba): First-principles calculations. <i>Chinese Journal of Physics</i> , 2018, 56, 870-879.	3.9	112
13	First-Principles Calculations to Investigate the Refractive Index and Optical Dielectric Constant of Na ₃ Sb ₄ X ₄ (X = S, Se) Ternary Chalcogenides. <i>Physica Status Solidi (B): Basic Research</i> , 2019, 256, 1900131.	1.5	106
14	First-principles computations of Y _{1-x} Ga _{1-x} As-ternary alloys: a study on structural, electronic, optical and elastic properties. <i>Bulletin of Materials Science</i> , 2020, 43, 1.	1.7	105
15	Electronic and thermoelectric properties of the layered Zintl phase CaIn ₂ P ₂ : first-principles calculations. <i>Philosophical Magazine</i> , 2020, 100, 3023-3039.	1.6	105
16	Half-Metallic Ferrimagnetic Characteristics of Co ₂ YZ (Z = P, As, Sb, and Bi) New Full-Heusler Alloys: a DFT Study. <i>Journal of Superconductivity and Novel Magnetism</i> , 2018, 31, 241-250.	1.8	100
17	Elastic, electronic, optical and thermodynamic properties of Ba ₃ Ca ₂ Si ₂ N ₆ semiconductor: First-principles predictions. <i>Physica B: Condensed Matter</i> , 2020, 589, 412213.	2.7	100
18	Structural, Elastic, Electronic and Optical Properties of LaOAgS-Type Silver Fluoride Chalcogenides: First-Principles Study. <i>Journal of Electronic Materials</i> , 2017, 46, 4539-4556.	2.2	98

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19	First-principles Calculations of Structural, Magnetic Electronic and Optical Properties of Rare-earth Metals TbX (X=N, O, S, Se). Journal of Superconductivity and Novel Magnetism, 2017, 30, 3471-3479.	1.8	96
20	First-Principle Calculations of Structural, Elastic, and Electronic Properties of Intermetallic Rare Earth R ₂ Ni ₂ Pb (R = Ho, Lu, and Sm) Compounds. Journal of Superconductivity and Novel Magnetism, 2018, 31, 395-403.	1.8	90
21	Challenges in improving the performance of eddy current testing: Review. Measurement and Control, 2019, 52, 46-64.	1.8	90
22	Structural, morphological and electrical properties of Cd ²⁺ -doped MgFe _{2-x} O ₄ ferrite nanoparticles. Journal of Alloys and Compounds, 2017, 726, 179-186.	5.5	88
23	Structural, electronic and thermodynamic investigation of Ag ₂ GdSi, Ag ₂ GdSn and Ag ₂ GdPb Heusler alloys: First-principles calculations. Materialpruefung/Materials Testing, 2021, 63, 537-542.	2.2	85
24	Electronic, optical, elastic, thermoelectric and thermodynamic properties of the spinel oxides ZnRh ₂ O ₄ and CdRh ₂ O ₄ . Journal of Alloys and Compounds, 2019, 774, 299-314.	5.5	84
25	Calculation of structural, optical and electronic properties of ZnS, ZnSe, MgS, MgSe and their quaternary alloy Mg _{1-x} Zn _x SySe _{1-y} . Materials Science and Engineering B: Solid-State Materials for Advanced Technology, 2003, 100, 163-171.	3.5	83
26	Structural, elastic and lattice dynamical properties of the alkali metal tellurides: First-principles study. Physica B: Condensed Matter, 2017, 521, 204-214.	2.7	83
27	First-principle study of structural, electronic and elastic properties of beryllium chalcogenides BeS, BeSe and BeTe. Computational Materials Science, 2006, 37, 292-299.	3.0	80
28	First-Principle Investigation of Structural, Electronic and Magnetic Properties in Mn ₂ RhZ (Z = Si, Ge). Tj ETQq0 0 0 rgBT /Overlock 10 Tf 5	1.8	80
29	Investigation of the optical properties of Mg(OH) ₂ and MgO nanostructures obtained by microwave-assisted methods. Journal of Alloys and Compounds, 2012, 521, 71-76.	5.5	79
30	Structural, mechanical and electronic properties of sodium based fluoroperovskites NaXF ₃ (X=Mg,). Tj ETQq0 0 0 rgBT /Overlock 10 Tf 5	4.0	79
31	A needle-like Cu ₂ CdSnS ₄ alloy nanostructure-based integrated electrochemical biosensor for detecting the DNA of Dengue serotype 2. Mikrochimica Acta, 2017, 184, 2211-2218.	5.0	75
32	Elastic, electronic, optical and thermoelectric properties of the novel Zintl-phase Ba ₂ ZnP ₂ . Solid State Sciences, 2022, 128, 106893.	3.2	75
33	Gd impurities effect on $\text{Co}_{1-x}\text{CrSi}_x$ alloy: first-principle calculations. Bulletin of Materials Science, 2018, 41, 1.	1.7	74
34	Full-potential calculations of structural, elastic and electronic properties of MgAl ₂ O ₄ and ZnAl ₂ O ₄ compounds. Physics Letters, Section A: General, Atomic and Solid State Physics, 2005, 344, 271-279.	2.1	73
35	New optical features to enhance solar cell performance based on porous silicon surfaces. Applied Surface Science, 2011, 257, 6112-6117.	6.1	73
36	First-principle calculations of structural, electronic and magnetic investigations of Mn ₂ RuGe _{1-x} Sn _x quaternary Heusler alloys. Chinese Journal of Physics, 2018, 56, 567-573.	3.9	72

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37	Structural and optical insights to enhance solar cell performance of CdS nanostructures. Energy Conversion and Management, 2014, 82, 238-243.	9.2	71
38	Copper substitution effect on the structural properties of nickel ferrites. Ceramics International, 2014, 40, 14413-14419.	4.8	71
39	Theoretical investigations of structural, mechanical, electronic and optical properties of NaScSi alloy. Emergent Materials, 2021, 4, 1465-1477.	5.7	71
40	Structural and optoelectronic properties of NiTiX and CoVX (X = Sb and Sn) half-Heusler compounds: An ab initio study. Optik, 2013, 124, 570-574.	2.9	70
41	Improved room temperature dielectric properties of Gd ³⁺ and Nb ⁵⁺ co-doped Barium Titanate ceramics. Journal of Alloys and Compounds, 2021, 883, 160836.	5.5	68
42	First-principles calculations of the elastic, electronic, and optical properties of the filled skutterudites CeFe ₄ P ₁₂ and ThFe ₄ P ₁₂ . Physical Review B, 2007, 75, .	3.2	66
43	Investigation of the Structural, Elastic, Electronic, and Optical Properties of Half-Heusler CaMgZ (Z = Tl, Bi, Sb, As, P, N, As, Sb, Bi, Tl, Pb, Sn, Ge, Si, C). Journal of Applied Physics, 2014, 115, 093701.	2.2	65
44	An ab initio study of the electronic structure and optical properties of CdS _{1-x} Te alloys. Solar Energy, 2010, 84, 1979-1984.	6.1	63
45	Doping-Induced Half-Metallic Ferromagnetism in Vanadium and Chromium-Doped Alkali Oxides K ₂ O and Rb ₂ O: Ab Initio Method. Journal of Superconductivity and Novel Magnetism, 2017, 30, 2197-2210.	1.8	61
46	Structural, mechanical, magnetic, electronic, and thermal investigations of Ag ₂ YB (Y = Nd, Sm, Gd) full-Heusler alloys. Emergent Materials, 2021, 4, 1769-1783.	5.7	61
47	Ab initio study of the structural and optoelectronic properties of the half-Heusler CoCrZ (Z = Al, Ga). Canadian Journal of Physics, 2014, 92, 1105-1112.	1.1	60
48	Optical investigations using ultra-soft pseudopotential calculations of Si _{0.5} Ge _{0.5} alloy. Solid State Communications, 2008, 148, 521-524.	1.9	59
49	Chalcogenides-based quantum dots: Optical investigation using first-principles calculations. Materials Science in Semiconductor Processing, 2015, 39, 276-282.	4.0	59
50	Structural, dielectric and low temperature magnetic response of Zn doped cobalt ferrite nanoparticles. AIP Advances, 2019, 9, .	1.3	58
51	First-principle calculations to investigate the elastic and thermodynamic properties of A ₃ B ₂ Rh ₃ (A = Sc, Y and La) perovskite compounds. Molecular Physics, 2012, 110, 121-128.	1.07	55
52	Detecting the DNA of dengue serotype 2 using aluminium nanoparticle doped zinc oxide nanostructure: Synthesis, analysis and characterization. Journal of Materials Research and Technology, 2020, 9, 5515-5523.	5.8	55
53	Structural, elastic, electronic and optical properties of the newly synthesized monoclinic Zintl phase BaIn ₂ P ₂ . Solid State Sciences, 2014, 29, 12-23.	3.2	54
54	Ab initio study of the pressure dependence of mechanical and thermodynamic properties of GeB ₂ O ₄ (B) Tj ETQq0 0 0 rgBT /Overlock 10	2.1	54

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55	Structural, elastic, electronic, chemical bonding and optical properties of Cu-based oxides ACuO (A=Li, Na, K and Rb): An ab initio study. Computational Materials Science, 2014, 81, 561-574.	3.0	53
56	Cadmium effect on optical properties of Cu ₂ Zn _{1-x} Cd _x SnS ₄ quaternary alloys nanostructures. Solar Energy, 2015, 114, 39-50.	6.1	53
57	Structural, Elastic, Thermodynamic, Electronic, and Magnetic Investigations of Full-Heusler Compound Ag ₂ CeAl: FP-LAPW Method. Journal of Superconductivity and Novel Magnetism, 2018, 31, 3183-3192.	1.8	53
58	Optical investigation of nanophotonic lithium niobate-based optical waveguide. Applied Physics B: Lasers and Optics, 2015, 121, 107-116.	2.2	52
59	Investigation of structural, elastic, electronic, and magnetic proprieties for X ₂ LuSb (X= Mn and Ir) full-Heusler alloys. Emergent Materials, 2022, 5, 537-551.	5.7	52
60	Electronic and optical properties of Zn Cd _{1-x} Se. Materials Chemistry and Physics, 2003, 82, 49-54.	4.0	51
61	Optical properties of Cauliflower-like Bi ₂ O ₃ nanostructures by reactive pulsed laser deposition (PLD) technique. Solar Energy, 2014, 107, 523-529.	6.1	51
62	Structural, elastic, electronic and thermodynamic properties of uranium filled skutterudites UFe ₄ P ₁₂ : First principle method. Materials Science in Semiconductor Processing, 2014, 27, 368-379.	4.0	50
63	Ab initio method of optical investigations of CdS _{1-x} Te alloys under quantum dots diameter effect. Solar Energy, 2015, 115, 33-39.	6.1	50
64	Predictive study of structural, electronic, magnetic and thermodynamic properties of XFeO ₃ (X = Ag, Zr and Ru) multiferroic materials in cubic perovskite structure: first-principles calculations. Materials Science-Poland, 2015, 33, 402-413.	1.0	50
65	Two symmetric n-type interfaces SrTiO ₃ /LaAlO ₃ in perovskite: Electronic properties from density functional theory. Journal of Applied Physics, 2016, 119, .	2.5	50
66	Synthesis and evaluation of the structural, optical, and antibacterial properties of copper oxide nanoparticles. Applied Physics A: Materials Science and Processing, 2019, 125, 1.	2.3	49
67	Optical investigations of photonics lithium niobate. Solar Energy, 2015, 120, 381-388.	6.1	48
68	Structural and optical investigations of cadmium sulfide nanostructures for optoelectronic applications. Solar Energy, 2012, 86, 3234-3240.	6.1	47
69	Improved efficiency of Cu(In,Ga)Se ₂ thinfilm solar cells using a buffer layer alternative to CdS. Solar Energy, 2019, 178, 150-156.	6.1	47
70	Ab initio exploration of the structural, elastic, electronic and optical properties of a new layered perovskite-type oxyfluoride: CsSrNb ₂ O ₆ F. Materials Science in Semiconductor Processing, 2021, 131, 105890.	4.0	47
71	Electronic and Magnetic Properties of Co ₂ CrGa _{1-x} Si _x Heusler Alloys. Journal of Superconductivity and Novel Magnetism, 2017, 30, 421-424.	1.8	46
72	Fabrication and characterizations of Al nanoparticles doped ZnO nanostructures-based integrated electrochemical biosensor. Journal of Materials Research and Technology, 2020, 9, 857-867.	5.8	46

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73	Correlation between the ionicity character and the heteropolar band gap in semiconductors. <i>Physica B: Condensed Matter</i> , 2001, 301, 295-298.	2.7	45
74	Density functional study of optical properties of beryllium chalcogenides compounds in nickel arsenide B8 structure. <i>Physica B: Condensed Matter</i> , 2012, 407, 286-296.	2.7	45
75	Synthesis of SnO ₂ nanostructures employing Nd:YAG laser. <i>Applied Physics A: Materials Science and Processing</i> , 2015, 120, 725-730.	2.3	44
76	Stirrer time effect on optical properties of nanophotonic LiNbO ₃ . <i>Materials Chemistry and Physics</i> , 2018, 203, 243-248.	4.0	43
77	GaNO colloidal nanoparticles synthesis by nanosecond pulsed laser ablation: Laser fluence dependent optical absorption and structural properties. <i>Powder Technology</i> , 2017, 320, 457-461.	4.2	42
78	First-principles calculations to investigate the structural, electronic and optical properties of Zn _{1-x} Mg _x Te ternary alloys. <i>Chinese Journal of Physics</i> , 2017, 55, 1018-1031.	3.9	41
79	Structural, Mechanical and Thermodynamic Properties under Pressure Effect of Rubidium Telluride: First Principle Calculations. <i>Archives of Metallurgy and Materials</i> , 2017, 62, 865-871.	0.6	41
80	First-principles calculations of a half-metallic ferromagnet zinc blende Zn _{1-x} V _x Te. <i>Journal of Magnetism and Magnetic Materials</i> , 2015, 378, 41-49.	2.3	40
81	Pressure and temperature dependence of the structural, elastic and thermodynamic properties of potassium telluride: First-principles calculations. <i>Chinese Journal of Physics</i> , 2017, 55, 769-779.	3.9	40
82	Effect of stirring time on the structural parameters of nanophotonic LiNbO ₃ deposited by spin-coating technique. <i>Optik</i> , 2018, 156, 886-890.	2.9	40
83	The Elastic, Electronic and Thermodynamic Properties of a New Cd Based Full Heusler Compounds - A Theoretical Investigation Using DFT Based FP-LMTO Approach. <i>Acta Physica Polonica A</i> , 2019, 136, 127-134.	0.5	40
84	Correlation between the bulk modulus and the charge density in semiconductors. <i>Physica B: Condensed Matter</i> , 2001, 305, 186-190.	2.7	39
85	First principle study of mechanical stability and thermodynamic properties of anti-fluorite Li ₂ O and Rb ₂ O under pressure and temperature effect. <i>Chinese Journal of Physics</i> , 2016, 54, 678-694.	3.9	39
86	Ultrasonic effect on optical, structural, topographical and morphological studies of Cu ₂ CdSnS ₄ quaternary alloy nanostructures. <i>Journal of Alloys and Compounds</i> , 2016, 686, 883-895.	5.5	37
87	Structural, elastic, electronic, magnetic, optical, and thermoelectric properties of the diamond-like quaternary semiconductor CuMn ₂ InSe ₄ . <i>Journal of Superconductivity and Novel Magnetism</i> , 2020, 33, 1091-1102.	1.8	37
88	Empirical formula relating the bulk modulus to the lattice constant in tetrahedral semiconductors. <i>Materials Chemistry and Physics</i> , 2004, 87, 14-17.	4.0	36
89	Optical properties of (Pb _{1-x} Mn _x S) _{1-y} Fe _y materials from first-principles calculations. <i>Chinese Journal of Physics</i> , 2017, 55, 1032-1043.	3.9	36
90	Comparative study of Fe doped ZnO based diluted and condensed magnetic semiconductors in wurtzite and zinc-blende structures by first-principles calculations. <i>Materials Science in Semiconductor Processing</i> , 2016, 43, 123-128.	4.0	35

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91	Fabricated Optical Strip Waveguide of Nanophotonics Lithium Niobate. IEEE Photonics Journal, 2016, 8, 1-10.	2.0	35
92	Aluminium nanoparticles size effect on the optical and structural properties of ZnO nanostructures synthesized by spin-coating technique. Results in Physics, 2017, 7, 1190-1197.	4.1	34
93	Effective Synthesis of Silicon Carbide Nanotubes by Microwave Heating of Blended Silicon Dioxide and Multi-Walled Carbon Nanotube. Materials Research, 2017, 20, 1658-1668.	1.3	34
94	External temperature and pressure effects on thermodynamic properties and mechanical stability of yttrium chalcogenides YX (X=S, Se and Te). Physica B: Condensed Matter, 2013, 428, 78-88.	2.7	33
95	Application of Gold Nanoparticles for Electrochemical DNA Biosensor. Journal of Nanomaterials, 2014, 2014, 1-7.	2.7	33
96	Calculation of bulk moduli of semiconductor compounds. Physica B: Condensed Matter, 2002, 322, 179-182.	2.7	32
97	Correlation between the bulk modulus and the transition pressure in semiconductors. Materials Letters, 2005, 59, 2032-2034.	2.6	32
98	Calculated optical properties of GaX (X=P, As, Sb) under hydrostatic pressure. Applied Physics A: Materials Science and Processing, 2011, 104, 1159-1167.	2.3	32
99	Characterisation, analysis and optical properties of nanostructure ZnO using the sol-gel method. Micro and Nano Letters, 2012, 7, 163.	1.3	32
100	Review on the energy and renewable energy status in Iraq: The outlooks. Renewable and Sustainable Energy Reviews, 2014, 39, 816-827.	16.4	32
101	Structural and Optical Properties of Nanophotonic LiNbO3 under Stirrer Time Effect. Journal of Optical Communications, 2018, 39, 297-306.	4.7	32
102	Nanosecond pulsed laser ablation to synthesize GaO colloidal nanoparticles: Optical and structural properties. Optik, 2019, 178, 337-342.	2.9	32
103	Structural and electronic properties of GaN x As ^{1-x} alloys. Applied Physics A: Materials Science and Processing, 2012, 106, 687-696.	2.3	31
104	Morphology and optical investigations of ZnO pyramids and nanoflakes for optoelectronic applications. Optik, 2014, 125, 2560-2564.	2.9	31
105	Analytical investigations of CdS nanostructures for optoelectronic applications. Optik, 2015, 126, 5109-5114.	2.9	31
106	Synthesis of carbon-based quantum dots from starch extracts: Optical investigations. Luminescence, 2018, 33, 260-266.	2.9	31
107	Effect of Temperature on the Physical, Electro-Chemical and Adsorption Properties of Carbon Micro-Spheres Using Hydrothermal Carbonization Process. Nanomaterials, 2018, 8, 597.	4.1	31
108	Theoretical investigation of structural, electronic, elastic, magnetic, thermodynamic, and thermoelectric properties of Ru ₂ MnNb Heusler alloy: FP-LMTO method. Emergent Materials, 2022, 5, 1065-1073.	5.7	31

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109	Structural, elastic, electronic and thermodynamic investigations of neptunium chalcogenides: First-principles calculations. Chinese Journal of Physics, 2016, 54, 33-41.	3.9	30
110	Recent Progress in the Development of Biosensors for Chemicals and Pesticides Detection. IEEE Access, 2020, 8, 82514-82527.	4.2	30
111	Further optical properties of CdX (X=S, Te) compounds under quantum dot diameter effect: Ab initio method. Renewable Energy, 2012, 45, 232-236.	8.9	29
112	Theoretical investigation of the structural, elastic, electronic, and optical properties of the ternary tetragonal tellurides KBT ₂ (B = Al, In). Materials Science in Semiconductor Processing, 2020, 114, 105085.	4.0	29
113	Electronic and Magnetic Investigations of Rare-Earth Tm-doped AlGa _N Ternary Alloy. Journal of Superconductivity and Novel Magnetism, 2018, 31, 1767-1771.	1.8	28
114	Computational Prediction of Structural, Electronic, Elastic, and Thermoelectric Properties of FeVX (X=As, P) Half-Heusler Compounds. Journal of Electronic Materials, 2020, 49, 4916-4922.	2.2	28
115	First-principles calculations of structural, elastic, electronic, and optical properties of CaY ₂ P ₂ O ₁₄ (Y = Cu, Ag). Journal of Applied Physics, 2014, 116, 083701.	5.7	28
116	Electronic properties of orthorhombic LiGaS ₂ and LiGaSe ₂ . Applied Physics A: Materials Science and Processing, 2009, 94, 315-320.	2.3	27
117	Structural, elastic, electronic and thermodynamic properties of the filled skutterudite CeOs ₄ Sb ₁₂ determined by density functional theory. Materials Science in Semiconductor Processing, 2013, 16, 1508-1516.	4.0	27
118	First-principle investigations of structural, electronic and thermodynamic properties of CdS _{1-x} Te _x ternary alloys: (0.0 x 1.0). Materials Express, 2014, 4, 521-532.	0.5	27
119	Optical investigations of blue shift in ZnS quantum dots. Superlattices and Microstructures, 2015, 88, 662-667.	3.1	27
120	Synthesis and Characterization of Cu ₂ CdSnS ₄ Quaternary Alloy Nanostructures. International Journal of Electrochemical Science, 2018, 13, 6693-6707.	1.3	27
121	Exploiting of geothermal energy reserve and potential in Saudi Arabia: A case study at Ain Al Harrah. Energy Reports, 2019, 5, 632-638.	5.1	27
122	The pressure effect of the bulk modulus seen by the charge density in CdX compounds. Materials Chemistry and Physics, 2003, 78, 625-629.	4.0	26
123	Synthesis and Characterization of Natural Extracted Precursor Date Palm Fibre-Based Activated Carbon for Aluminum Removal by RSM Optimization. Processes, 2019, 7, 249.	2.8	26
124	Structural phase transition of boron nitride compound. Solid State Communications, 2004, 132, 465-470.	1.9	25
125	Elastic and thermodynamic properties of the SiB ₂ O ₄ (B=Mg, Zn and Cd) cubic spinels: An ab initio FP-LAPW study. Materials Science in Semiconductor Processing, 2015, 38, 192-202.	4.0	25
126	Analysis and characterization of Cu ₂ CdSnS ₄ quaternary alloy nanostructures deposited on GaN. Indian Journal of Physics, 2018, 92, 695-703.	1.8	25

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127	Half-metallic ferromagnetism in $\text{Be}_{1-x}\text{V}_x\text{Te}$ alloys: an Ab-initio study. Indian Journal of Physics, 2015, 89, 1251-1263.	1.8	24
128	Optoelectronic properties of naphtho[2, 1-b:6, 5-b ϵ^2]difuran derivatives for photovoltaic application: a computational study. Journal of Molecular Modeling, 2016, 22, 248.	1.8	24
129	Structural investigations through cobalt effect on ZnO nanostructures. Optik, 2016, 127, 10102-10107.	2.9	24
130	Structural, elastic, electronic and optical properties of the novel quaternary diamond-like semiconductors $\text{Cu}_2\text{MgSiS}_4$ and $\text{Cu}_2\text{MgGeS}_4$. Solid State Sciences, 2017, 70, 21-35.	3.2	24
131	Effect of nitrogen doping on structural and optical properties of $\text{Mg}_x\text{Zn}_{1-x}\text{O}$ ternary alloys. Optical Materials, 2019, 89, 554-558.	3.6	24
132	Electronic and positron properties of zinc-blende structure of GaN, AlN, and their alloy $\text{Ga}_{1-x}\text{Al}_x\text{N}$. Journal of Applied Physics, 2003, 93, 9730-9736.	2.5	23
133	Phase transition of Nowotny ϵ - Zn_2X ($\text{X}=\text{P}, \text{As}$ and Sb) compounds at high pressure: Theoretical investigation of structural, electronic and vibrational properties. Computational Materials Science, 2014, 87, 187-197.	3.0	23
134	Characterization and analysis of wheat-like CdS nanostructures under temperature effect for solar cells applications. Optik, 2016, 127, 8907-8915.	2.9	23
135	First-principles calculations to investigate optical properties of $\text{ByAl}_x\text{In}_{1-x}\text{N}$ alloys for optoelectronic devices. Superlattices and Microstructures, 2012, 51, 404-411.	3.1	22
136	Structural, Analysis and Optical Studies of Cadmium Sulfide Nanostructured. Procedia Engineering, 2013, 53, 217-224.	1.2	22
137	First-principles prediction of the structural, elastic, thermodynamic, electronic and optical properties of $\text{Li}_4\text{Sr}_3\text{Ge}_2\text{N}_6$ quaternary nitride. Journal of Alloys and Compounds, 2015, 618, 84-94.	5.5	22
138	Fabrication, analysis and characterization of $\text{Cu}_2\text{Zn}_{1-x}\text{Cd}_x\text{SnS}_4$ quaternary alloy nanostructures deposited on GaN. Journal of Materials Science, 2016, 51, 6876-6885.	3.7	22
139	A novel quaternary alloy ($\text{Cu}_2\text{Zn}_{1-x}\text{Cd}_x\text{SnS}_4$) nanostructured sensor for biomedical diagnosis. Materials Research Express, 2016, 3, 085022.	1.6	22
140	Optical analysis of lens-like $\text{Cu}_2\text{CdSnS}_4$ quaternary alloy nanostructures. Applied Physics A: Materials Science and Processing, 2016, 122, 1.	2.3	22
141	Structural, optical and electrical investigations of $\text{Cu}_2\text{Zn}_{1-x}\text{Cd}_x\text{SnS}_4/\text{Si}$ quaternary alloy nanostructures synthesized by spin coating technique. Microsystem Technologies, 2017, 23, 2223-2232.	2.0	22
142	Etching time effect on optical properties of porous silicon for solar cells fabrication. Optik, 2017, 147, 343-349.	2.9	22
143	Investigated optical studies of Si quantum dot. Solar Energy, 2011, 85, 2283-2287.	6.1	21
144	Solar energy status in Iraq: Abundant or not ϵ ”Steps forward. Journal of Renewable and Sustainable Energy, 2016, 8, 025905.	2.0	21

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145	Synthesis, purification and microstructural characterization of nickel doped carbon nanotubes for spintronic applications. <i>Ceramics International</i> , 2016, 42, 5600-5606.	4.8	21
146	Ab initio prediction of the elastic, electronic and optical properties of a new family of diamond-like semiconductors, Li ₂ HgMS ₄ (M = Si, Ge and Sn). <i>Journal of Alloys and Compounds</i> , 2020, 843, 155991.	5.5	21
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