Wen-Zhi Xiao

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Electronic structures and magnetic properties in nonmetallic element substituted MoS2 monolayer. Computational Materials Science, 2015, 107, 72-78.	3.0	55
2	Theoretical discovery of novel two-dimensional V ^A -N binary compounds with auxiticity. Physical Chemistry Chemical Physics, 2018, 20, 22027-22037.	2.8	52
3	Two-dimensional H-TiO2/MoS2(WS2) van der Waals heterostructures for visible-light photocatalysis and energy conversion. Applied Surface Science, 2020, 504, 144425.	6.1	48
4	Electronic and magnetic properties of SnS 2 monolayer doped with 4 d transition metals. Journal of Magnetism and Magnetic Materials, 2017, 438, 152-162.	2.3	35
5	Magnetism in undoped ZnS studied from density functional theory. Journal of Applied Physics, 2014, 115, .	2.5	34
6	A first-principles study of the SnO2 monolayer with hexagonal structure. Journal of Chemical Physics, 2016, 145, 174702.	3.0	34
7	New two-dimensional V-V binary compounds with a honeycomb-like structure: a first-principles study. Materials Research Express, 2018, 5, 035903.	1.6	34
8	Electronic structures and magnetic properties in Cu-doped two-dimensional dichalcogenides. Physica E: Low-Dimensional Systems and Nanostructures, 2015, 73, 69-75.	2.7	33
9	Two-dimensional hexagonal chromium chalco-halides with large vertical piezoelectricity, high-temperature ferromagnetism, and high magnetic anisotropy. Physical Chemistry Chemical Physics, 2020, 22, 14503-14513.	2.8	32
10	Ferromagnetic and metallic properties of the semihydrogenated GaN sheet. Physica Status Solidi (B): Basic Research, 2011, 248, 1442-1445.	1.5	27
11	Firstâ€principles study of magnetic properties in Agâ€doped SnO ₂ . Physica Status Solidi (B): Basic Research, 2011, 248, 1961-1966.	1.5	26
12	Unexpected magnetic properties in carbon-doped SnO2 from first-principles calculation. Computational Materials Science, 2014, 83, 5-11.	3.0	26
13	Half-metallic and magnetic properties in nonmagnetic element embedded graphitic carbon nitride sheets. Physical Chemistry Chemical Physics, 2015, 17, 22136-22143.	2.8	25
14	Ab initio study of magnetism in nonmagnetic metal substituted monolayer MoS2. Solid State Communications, 2015, 220, 67-71.	1.9	21
15	Stability and electronic structure of two-dimensional arsenic phosphide monolayer. Materials Science and Engineering B: Solid-State Materials for Advanced Technology, 2018, 228, 206-212.	3.5	21
16	Magnetic properties in CdS monolayer doped with first-row elements: A density functional theory investigation. Physica Status Solidi (B): Basic Research, 2014, 251, 1257-1264.	1.5	19
17	First principles study on magnetic properties in ZnS doped with palladium. European Physical Journal B, 2013, 86, 1.	1.5	18
18	Magnetic properties in BiFeO3 doped with Cu and Zn first-principles investigation. Journal of Alloys and Compounds, 2016, 674, 463-469.	5.5	17

Wen-Zhi Xiao

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19	Half-metallic and magnetic properties of AlN nanosheets doped with nonmagnetic metals: A first-principles study. Computational Materials Science, 2016, 124, 98-105.	3.0	14
20	Magnetism and ferroelectricity in BiFeO3 doped with Ga at Fe sites. Journal of Alloys and Compounds, 2019, 797, 117-121.	5.5	13
21	Elasticity, piezoelectricity, and mobility in two-dimensional BiTeI from a first-principles study. Journal Physics D: Applied Physics, 2020, 53, 245301.	2.8	13
22	First-principles calculations of electronic and magnetic properties in semi-fluorinated CdS sheet. Physics Letters, Section A: General, Atomic and Solid State Physics, 2012, 376, 3402-3406.	2.1	12
23	Magnetic properties in AlN nanosheet doped with alkali metals: A first-principles study. Physica Status Solidi (B): Basic Research, 2016, 253, 1816-1823.	1.5	12
24	Ferromagnetism and controllable half-metallicity of two-dimensional hexagonal CrOX (XÂ=ÂF, Cl, Br) monolayers. Journal of Magnetism and Magnetic Materials, 2020, 515, 167310.	2.3	11
25	Ferromagnetic coupling in Mgâ€doped passivated AlN nanowires: A firstâ€principles study. Physica Status Solidi (B): Basic Research, 2012, 249, 185-189.	1.5	10
26	Magnetic properties in semifluorinated GaN sheet from first principles calculations. Physica Status Solidi (B): Basic Research, 2012, 249, 1465-1469.	1.5	10
27	Magnetic properties in a IIIA-nitride monolayer doped with Cu: a density functional theory investigation. RSC Advances, 2015, 5, 82357-82362.	3.6	10
28	Magnetic properties in BiFeO ₃ doped with nonâ€metallic element: Firstâ€principles investigation. Physica Status Solidi (B): Basic Research, 2016, 253, 279-283.	1.5	10
29	Electronic and magnetic properties of SnS2 monolayer doped with non-magnetic elements. Physica E: Low-Dimensional Systems and Nanostructures, 2018, 99, 182-188.	2.7	9
30	Half-metallicity in carbon-substituted CdS monolayer. Physica E: Low-Dimensional Systems and Nanostructures, 2014, 59, 230-234.	2.7	8
31	First-principles insight into the surface magnetism of Cu-doped SnO ₂ (110) thin film. RSC Advances, 2014, 4, 39860.	3.6	8
32	Electronic and magnetic properties in Mnâ€doped IIIAâ€nitride monolayers. Physica Status Solidi (B): Basic Research, 2016, 253, 2001-2008.	1.5	8
33	Two-dimensional Al2O3 with ultrawide bandgap and large exciton binding energy for solar-blind ultraviolet photodetectors. Computational Materials Science, 2021, 200, 110775.	3.0	8
34	Newtype two-dimensional Cr2O3 monolayer with half-metallicity, high curie temperature, and magnetic anisotropy. Journal of Magnetism and Magnetic Materials, 2022, 543, 168657.	2.3	6
35	Electronic structures and magnetic properties in transition metal adsorbed gt-C3N4 monolayer. Journal of Magnetism and Magnetic Materials, 2020, 493, 165745.	2.3	5
36	Half-metallicity and enhanced Curie temperature of Ti-embedded CrI3 monolayer. Materials Today Communications, 2020, 25, 101438.	1.9	5

WEN-ZHI XIAO

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37	First-principle study on the stability, mechanical, electronic, and optical properties of two-dimensional scandium oxyhalides. Materials Chemistry and Physics, 2022, 287, 126306.	4.0	3
38	Magnetic properties in Nb/Tc adsorbed gt-C3N4 monolayer. Physica E: Low-Dimensional Systems and Nanostructures, 2020, 116, 113753.	2.7	2
39	Nitrogen-induced magnetism in stannates from first-principles calculations. International Journal of Modern Physics B, 2016, 30, 1650236.	2.0	1
40	Oxygenationâ€Induced Twoâ€Dimensional Topological Insulators in Antimony Arsenide. Physica Status Solidi - Rapid Research Letters, 2019, 13, 1900146.	2.4	1
41	Two-dimensional hexagonal LaOF with ultrawide bandgap, large exciton energy, and low lattice thermal conductivity. Physica E: Low-Dimensional Systems and Nanostructures, 2022, 140, 115195.	2.7	1
42	Large exciton binding energy, superior mechanical flexibility, and ultra-low lattice thermal conductivity in Bil ₃ monolayer. Journal of Physics Condensed Matter, 2022, 34, 055302.	1.8	0