

# Zhifeng Liu

## List of Publications by Year in descending order

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33  
papers

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citations

758635

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525886

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34  
docs citations

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times ranked

752  
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#	ARTICLE	IF	CITATIONS
1	Robust Dirac spin gapless semiconductors in a two-dimensional oxalate based organic honeycomb-kagome lattice. <i>Nanoscale</i> , 2022, 14, 2023-2029.	2.8	5
2	Cluster- and energy-separated extreme states in a synthesized superatomic solid. <i>Physical Review B</i> , 2022, 105, .	1.1	1
3	Intrinsic spin-valley-coupled Dirac state in Janus functionalized $\hat{I}^2$ -BiAs monolayer. <i>Nanoscale Horizons</i> , 2021, 6, 283-289.	4.1	9
4	MnNBr Monolayer: A High-Temperature Ferromagnetic Half-Metal with Type-II Weyl Fermions. <i>Physica Status Solidi - Rapid Research Letters</i> , 2021, 15, 2100115.	1.2	7
5	MnNBr Monolayer: A High-Temperature Ferromagnetic Half-Metal with Type-II Weyl Fermions. <i>Physica Status Solidi - Rapid Research Letters</i> , 2021, 15, 2170027.	1.2	0
6	Recent progress on 2D magnets: Fundamental mechanism, structural design and modification. <i>Applied Physics Reviews</i> , 2021, 8, .	5.5	202
7	Strain-Induced Ideal Topological Semimetal in $Ort_{32}$ Holding Parallel Arc-Like Nodal Lines and Anisotropic Multiple Weyl Fermions. <i>Physica Status Solidi - Rapid Research Letters</i> , 2021, 15, 2100324.	1.2	0
8	Metallic subnanometer porous silicon: A theoretical prediction. <i>Physical Review B</i> , 2021, 103, .	1.1	13
9	T-C56: a low-density transparent superhard carbon allotrope assembled from C16 cage-like cluster. <i>Journal of Physics Condensed Matter</i> , 2020, 32, 165701.	0.7	6
10	Room temperature ferromagnetic half metal in Mn doped cluster-assembled sodalite phase of III-N compounds. <i>Journal of Magnetism and Magnetic Materials</i> , 2020, 499, 166295.	1.0	4
11	Palgraphyne: A Promising 2D Carbon Dirac Semimetal with Strong Mechanical and Electronic Anisotropy. <i>Physica Status Solidi - Rapid Research Letters</i> , 2020, 14, 1900670.	1.2	14
12	Three-dimensional borophene: A light-element topological nodal-line semimetal with direction-dependent type-II Weyl fermions. <i>Physical Review B</i> , 2020, 102, .	1.1	9
13	Palgraphyne: A Promising 2D Carbon Dirac Semimetal with Strong Mechanical and Electronic Anisotropy. <i>Physica Status Solidi - Rapid Research Letters</i> , 2020, 14, 2070020.	1.2	4
14	Phase transition, magnetic and electronic properties of iron mononitride: First-principles calculations. <i>Journal of Alloys and Compounds</i> , 2019, 771, 322-326.	2.8	10
15	Two-dimensional spin-valley-coupled Dirac semimetals in functionalized SbAs monolayers. <i>Materials Horizons</i> , 2019, 6, 781-787.	6.4	38
16	All-Silicon Topological Semimetals with Closed Nodal Line. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 244-250.	2.1	24
17	Hexagonal $M_{2}C_{3}$ (M = As, Sb, and Bi) monolayers: new functional materials with desirable band gaps and ultrahigh carrier mobility. <i>Journal of Materials Chemistry C</i> , 2018, 6, 12689-12697.	2.7	42
18	Cluster-Assembled Semiconductor CdO Polymorph with Good Ductility, High Carrier Mobility, and Promising Optical Properties. <i>Journal of Physical Chemistry C</i> , 2018, 122, 24287-24294.	1.5	6

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19	YN2 monolayer: Novel p-state Dirac half metal for high-speed spintronics. Nano Research, 2017, 10, 1972-1979.	5.8	120
20	Computational search for two-dimensional intrinsic half-metals in transition-metal dinitrides. Journal of Materials Chemistry C, 2017, 5, 727-732.	2.7	61
21	Design of superhalogens using a core-shell structure model. Nanoscale, 2017, 9, 18781-18787.	2.8	12
22	Room-Temperature Ordered Spin Structures in Cluster-Assembled Single V@Si <sub>12</sub> Sheets. Journal of Physical Chemistry C, 2015, 119, 1517-1523.	1.5	50
23	A new diluted magnetic semiconductor based on the expanded phase of ZnS: surmounting the random distribution of magnetic impurities. Physical Chemistry Chemical Physics, 2015, 17, 13117-13122.	1.3	7
24	First-principle study on thermodynamic property of superhard BC <sub>2</sub> N under extreme conditions. Journal of Materials Research, 2014, 29, 1326-1333.	1.2	0
25	First-principle study of phase stability, electronic structure and thermodynamic properties of cadmium sulfide under high pressure. Journal of Physics and Chemistry of Solids, 2014, 75, 662-669.	1.9	12
26	Enhancement of hydrogen binding affinity with low ionization energy Li <sub>2</sub> F coating on C <sub>60</sub> to improve hydrogen storage capacity. International Journal of Hydrogen Energy, 2014, 39, 15639-15645.	3.8	21
27	From the ZnO Hollow Cage Clusters to ZnO Nanoporous Phases: A First-Principles Bottom-Up Prediction. Journal of Physical Chemistry C, 2013, 117, 17633-17643.	1.5	45
28	Low-density nanoporous phases of group-III nitrides built from sodalite cage clusters. Physical Chemistry Chemical Physics, 2013, 15, 8186.	1.3	19
29	New nanomaterials based on In <sub>12</sub> As <sub>12</sub> cages: an ab initio bottom-up study. RSC Advances, 2013, 3, 1450-1459.	1.7	11
30	Structures, stabilities, and electronic properties of GaAs tubelike clusters and single-walled GaAs nanotubes. Chinese Physics B, 2012, 21, 123601.	0.7	3
31	Structures, stabilities and electronic properties of InAs double-walled tubelike clusters and nanotubes. Wuli Xuebao/Acta Physica Sinica, 2012, 61, 243101.	0.2	0
32	Structures, Stabilities and Electronic Properties of InAs Tubelike Clusters and Single-Walled InAs Nanotubes. Wuli Huaxue Xuebao/ Acta Physico - Chimica Sinica, 2011, 27, 2079-2087.	2.2	3
33	Structures and electronic properties of Mo <sub>2</sub> N <sub>2</sub> (Tj ETQq1 1.0784314 rgBT /Ov	0.7	5