

# Zheng Liu

## List of Publications by Year in descending order

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60  
papers

6,177  
citations

117625

34  
h-index

123424

61  
g-index

63  
all docs

63  
docs citations

63  
times ranked

7897  
citing authors

#	ARTICLE	IF	CITATIONS
1	The role of LiO <sub>2</sub> solubility in O <sub>2</sub> reduction in aprotic solvents and its consequences for Li-O <sub>2</sub> batteries. <i>Nature Chemistry</i> , 2014, 6, 1091-1099.	13.6	942
2	A stable cathode for the aprotic Li-O <sub>2</sub> battery. <i>Nature Materials</i> , 2013, 12, 1050-1056.	27.5	677
3	Nanoparticulate TiO <sub>2</sub> (B): An Anode for Lithium-Ion Batteries. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 2164-2167.	13.8	305
4	Facile Synthesis and Characterization of Novel Mesoporous and Mesorelief Oxides with Gyroidal Structures. <i>Journal of the American Chemical Society</i> , 2004, 126, 865-875.	13.7	297
5	Stable Nontrivial $Z^2$ Topology in Ultrathin Bi (111) Films: A First-Principles Study. <i>Physical Review Letters</i> , 2011, 107, 136805.	7.8	292
6	Quantum Anomalous Hall Effect in 2D Organic Topological Insulators. <i>Physical Review Letters</i> , 2013, 110, 196801.	7.8	292
7	Organic topological insulators in organometallic lattices. <i>Nature Communications</i> , 2013, 4, 1471.	12.8	238
8	Epitaxial growth of large-gap quantum spin Hall insulator on semiconductor surface. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 14378-14381.	7.1	205
9	First-Principles Calculations on the Effect of Doping and Biaxial Tensile Strain on Electron-Phonon Coupling in Graphene. <i>Physical Review Letters</i> , 2013, 111, 196802.	7.8	197
10	Flat Chern Band in a Two-Dimensional Organometallic Framework. <i>Physical Review Letters</i> , 2013, 110, 106804.	7.8	191
11	Topological and electronic transitions in a Sb(111) nanofilm: The interplay between quantum confinement and surface effect. <i>Physical Review B</i> , 2012, 85, .	3.2	164
12	Single-ion conducting gel polymer electrolytes: design, preparation and application. <i>Journal of Materials Chemistry A</i> , 2020, 8, 1557-1577.	10.3	154
13	Exotic electronic states in the world of flat bands: From theory to material. <i>Chinese Physics B</i> , 2014, 23, 077308.	1.4	153
14	Tunable spin states in the two-dimensional magnet CrI <sub>3</sub> . <i>Nanoscale</i> , 2018, 10, 14298-14303.	5.6	136
15	Nonflammable organic electrolytes for high-safety lithium-ion batteries. <i>Energy Storage Materials</i> , 2020, 32, 425-447.	18.0	127
16	Progress and Perspective: MXene and MXene-Based Nanomaterials for High-Performance Energy Storage Devices. <i>Advanced Electronic Materials</i> , 2021, 7, 2000967.	5.1	122
17	Electronic Strengthening of Graphene by Charge Doping. <i>Physical Review Letters</i> , 2012, 109, 226802.	7.8	104
18	Prediction of two-dimensional nodal-line semimetals in a carbon nitride covalent network. <i>Journal of Materials Chemistry A</i> , 2018, 6, 11252-11259.	10.3	101



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37	Improving the cycling performance of LiNi <sub>0.8</sub> Co <sub>0.15</sub> Al <sub>0.05</sub> O <sub>2</sub> cathode materials via zirconium and fluorine co-substitution. Journal of Alloys and Compounds, 2019, 806, 136-145.	5.5	32
38	Selectively doping barlowite for quantum spin liquid: A first-principles study. Physical Review B, 2015, 92, .	3.2	31
39	Strain-Engineered Surface Transport in Si(001): Complete Isolation of the Surface State via Tensile Strain. Physical Review Letters, 2013, 111, 246801.	7.8	27
40	Effect of stacking order on the electronic state of $\text{LiTl}_2\text{Te}_4$ . Physical Review B, 2022, 105, .	3.2	27
41	Stabilizing Ni-Rich LiNi <sub>0.92</sub> Co <sub>0.06</sub> Al <sub>0.02</sub> O <sub>2</sub> Cathodes by Boracic Polyanion and Tungsten Cation Doping for High-Energy Lithium-Ion Batteries. ChemElectroChem, 2020, 7, 3811-3817.	3.4	24
42	Magnetic Dirac fermions and Chern insulator supported on pristine silicon surface. Physical Review B, 2016, 94, .	3.2	18
43	Relationships between strain and band structure in Si(001) and Si(110) nanomembranes. Physical Review B, 2009, 80, .	3.2	16
44	Tunable bending modulus and bending limit of oxidized graphene. Nanoscale, 2020, 12, 1623-1628.	5.6	16
45	A first-principle perspective on electronic nematicity in FeSe. Npj Quantum Materials, 2020, 5, .	5.2	15
46	Prediction of intrinsic topological superconductivity in Mn-doped GeTe monolayer from first-principles. Npj Computational Materials, 2021, 7, .	8.7	15
47	Exotic fractional topological states in a two-dimensional organometallic material. Physical Review B, 2014, 89, .	3.2	13
48	In-plane ordering of oxygen vacancies in a high- $T_c$ cuprate superconductor with compressed Cu-O octahedrons: An automated cluster expansion study. Physical Review Materials, 2020, 4, .	2.4	12
49	Role of interstitial hydrogen in $\text{SrCoO}_{2.5}$ antiferromagnetic insulator. Physical Review Materials, 2018, 2, .	2.1	11
50	Reconciling the bulk metallic and surface insulating state in $\text{TaSe}_2$ . Physical Review B, 2022, 105, .	3.2	9
51	Visualizing the evolution from Mott insulator to Anderson insulator in Ti-doped 1T-TaS <sub>2</sub> . Npj Quantum Materials, 2022, 7, .	5.2	9
52	Observation of Rashba splitting on reconstructed surface. Surface Science, 2013, 618, 115-119.	1.9	7
53	First-principles study of the organometallic compound Cu(1,3-bdc). Physical Review B, 2015, 92, .	3.2	7
54	$\pi$ -conjugation in the epitaxial Si(111)- $\sqrt{3}\times\sqrt{3}$ surface: Unconventional $\pi$ -bonding geometry for Si. Physical Review B, 2017, 95, .	3.2	5

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55	An Anode Material for Lithium Storage: Si@N,S-Doped Carbon Synthesized <i>via In Situ</i> Self-Polymerization. ACS Applied Energy Materials, 2021, 4, 3555-3562.	5.1	5
56	Electronic and spin dynamics in the insulating iron pnictide $\text{NaFe}_{1-x}\text{Co}_x\text{P}_2\text{As}_2$ . Physical Review B, 2017, 96, .	3.2	4
57	Renormalization of the Mott gap by lattice entropy: The case of 1T- TaS <sub>2</sub> . Physical Review Research, 2020, 2, .	3.6	4
58	Understanding the flat band in $\text{TaS}_2$ using a rotated basis. Physical Review B, 2021, 104, .	3.2	4
59	Testing density functional theory in a quantum Ising chain. Physical Review B, 2021, 104, .	3.2	2
60	Electron-nuclear hyperfine coupling in quantum kagome antiferromagnets from first-principles calculation and a reflection of the defect effect. Science Bulletin, 2019, 64, 1584-1591.	9.0	0