

Jiajie Zhu

List of Publications by Year in descending order

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46
papers

1,875
citations

393982

19
h-index

264894

42
g-index

46
all docs

46
docs citations

46
times ranked

3409
citing authors

#	ARTICLE	IF	CITATIONS
1	SnSe ₂ 2D Anodes for Advanced Sodium Ion Batteries. <i>Advanced Energy Materials</i> , 2016, 6, 1601188.	10.2	243
2	High-Rate and Ultralong Cycle-Life Potassium Ion Batteries Enabled by In Situ Engineering of Yolk-Shell FeS ₂ @C Structure on Graphene Matrix. <i>Advanced Energy Materials</i> , 2018, 8, 1802565.	10.2	207
3	Boosting the Yield of MXene 2D Sheets via a Facile Hydrothermal-Assisted Intercalation. <i>ACS Applied Materials & Interfaces</i> , 2019, 11, 8443-8452.	4.0	178
4	Active Edge Sites Engineering in Nickel Cobalt Selenide Solid Solutions for Highly Efficient Hydrogen Evolution. <i>Advanced Energy Materials</i> , 2017, 7, 1602089.	10.2	171
5	Two-Dimensional SnO Anodes with a Tunable Number of Atomic Layers for Sodium Ion Batteries. <i>Nano Letters</i> , 2017, 17, 1302-1311.	4.5	118
6	Silicene: Recent theoretical advances. <i>Applied Physics Reviews</i> , 2016, 3, .	5.5	94
7	S-functionalized MXenes as electrode materials for Li-ion batteries. <i>Applied Materials Today</i> , 2016, 5, 19-24.	2.3	89
8	Silicene for Na-ion battery applications. <i>2D Materials</i> , 2016, 3, 035012.	2.0	82
9	Multistimuli-Responsive Display Materials to Encrypt Differentiated Information in Bright and Dark Fields. <i>Advanced Functional Materials</i> , 2019, 29, 1906068.	7.8	79
10	Intrinsic Defects and H Doping in WO ₃ . <i>Scientific Reports</i> , 2017, 7, 40882.	1.6	65
11	P and Si functionalized MXenes for metal-ion battery applications. <i>2D Materials</i> , 2017, 4, 025073.	2.0	62
12	Nb-based MXenes for Li-ion battery applications. <i>Physica Status Solidi - Rapid Research Letters</i> , 2015, 9, 726-729.	1.2	61
13	Silicene/germanene on MgX ₂ (X = Cl, Br, and I) for Li-ion battery applications. <i>Nanoscale</i> , 2016, 8, 7272-7277.	2.8	61
14	Structural and Electronic Properties of Silicene on MgX ₂ (X = Cl, Br, and I). <i>ACS Applied Materials & Interfaces</i> , 2014, 6, 11675-11681.	4.0	55
15	Stability and electronic properties of silicene on WSe ₂ . <i>Journal of Materials Chemistry C</i> , 2015, 3, 3946-3953.	2.7	37
16	Perovskite Quantum Wells Formation Mechanism for Stable Efficient Perovskite Photovoltaics—A Real-Time Phase-Transition Study. <i>Advanced Materials</i> , 2021, 33, e2006238.	11.1	30
17	Stability and electronic properties of carbon in $\hat{\Gamma}$ -Al ₂ O ₃ . <i>Journal of Physics and Chemistry of Solids</i> , 2014, 75, 379-383.	1.9	27
18	Silicene on MoS ₂ : role of the van der Waals interaction. <i>2D Materials</i> , 2015, 2, 045004.	2.0	22

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19	Structural and electronic properties of CuI doped with Zn, Ga and Al. Journal of Physics and Chemistry of Solids, 2013, 74, 1122-1126.	1.9	19
20	Functionalized NbS ₂ as cathode for Li- and Na-ion batteries. Applied Physics Letters, 2017, 111, .	1.5	19
21	Band Gap Opening in Silicene on MgBr ₂ (0001) Induced by Li and Na. ACS Applied Materials & Interfaces, 2014, 6, 19242-19246.	4.0	13
22	Effect of Li doping on the O vacancies in Lu ₂ SiO ₅ :Ce phosphors. Materials Letters, 2018, 228, 372-374.	1.3	13
23	Stress-enhanced lithiation in MAX compounds for battery applications. Applied Materials Today, 2017, 9, 192-195.	2.3	12
24	Potential of B/Al-doped Silicene Electrodes in Na/K-ion Batteries. Advanced Theory and Simulations, 2018, 1, 1800017.	1.3	12
25	The phase transition and elastic and optical properties of polymorphs of CuI. Journal of Physics Condensed Matter, 2012, 24, 475503.	0.7	11
26	Stability and electronic properties of polar and non-polar surfaces of CuI. Applied Surface Science, 2013, 268, 87-91.	3.1	11
27	First-principles study on stability of Li, Na and Ca in Lu ₂ SiO ₅ . Journal of Luminescence, 2013, 139, 1-5.	1.5	11
28	Studies on phase stability, mechanical, optical and electronic properties of a new Gd ₂ CaZnO ₅ phosphor system for LEDs. CrystEngComm, 2014, 16, 1652.	1.3	10
29	B-doping-enhanced Stability of Phosphorene/Graphene Heterostructures. Advanced Theory and Simulations, 2019, 2, 1800176.	1.3	9
30	Structure and role of carbon-related defects in yttrium aluminum garnet. Optical Materials, 2021, 111, 110561.	1.7	8
31	First-principles calculations of oxygen vacancies and cerium substitution in lutetium pyrosilicate. Journal of Luminescence, 2012, 132, 164-170.	1.5	5
32	Phase transition and elastic and optical properties of Lu ₂ SiO ₅ . Optical Materials, 2013, 35, 1659-1663.	1.7	5
33	Phosphorene as cathode for metal-ion batteries: Importance of F decoration. Materials Today Energy, 2018, 10, 141-145.	2.5	5
34	Stability and electronic properties of O vacancies and Ce ⁴⁺ in Lu ₂ SiO ₅ tuned by C doping. Optical Materials, 2019, 93, 15-18.	1.7	5
35	Effect of Carbon Doping on F-type Defects in YAG and YAG:Ce Crystals. Physica Status Solidi (B): Basic Research, 2021, 258, 2100325.	0.7	5
36	Structural properties of Lu ₂ SiO ₅ doped with rare-earth elements. Materials Letters, 2019, 256, 126410.	1.3	4

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37	Reconstructive Phase Transformations in Body-Centered Cubic Titanium. Physica Status Solidi (B): Basic Research, 2020, 257, 2000193.	0.7	4
38	CO ₂ capture by Li-functionalized silicene. Physica Status Solidi - Rapid Research Letters, 2016, 10, 458-461.	1.2	3
39	Oxygen Doping Enhanced Lithiation in MgCl ₂ for Battery Applications. Physica Status Solidi (B): Basic Research, 2019, 256, 1900166.	0.7	3
40	Martensitic transformations of γ -phase in zirconium. Journal of Applied Physics, 2021, 129, .	1.1	3
41	Effect of cation doping on tuning intrinsic defects in Lu ₃ . Journal of Luminescence, 2019, 212, 238-241.	1.5	2
42	Dynamic instability of lithiated phosphorene. RSC Advances, 2020, 10, 32259-32264.	1.7	2
43	Elemental Two-Dimensional Materials Beyond Graphene. ChemistrySelect, 2017, 2, .	0.7	0
44	11. Elemental Two-Dimensional Materials Beyond Graphene. , 2017, , 219-228.		0
45	Condensed Matter in Energy, Environment, and Beyond. Advances in Condensed Matter Physics, 2019, 2019, 1-2.	0.4	0
46	First-Principles Calculations on the Diffusion and Electronic Properties of CuI Doped by Cation and Anion. Results in Physics, 2022, , 105595.	2.0	0