

Yifan Nie

List of Publications by Year in descending order

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31
papers

1,923
citations

331670

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434195

31
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32
all docs

32
docs citations

32
times ranked

3798
citing authors

#	ARTICLE	IF	CITATIONS
1	Polarity governs atomic interaction through two-dimensional materials. <i>Nature Materials</i> , 2018, 17, 999-1004.	27.5	182
2	Realizing Large-Scale, Electronic-Grade Two-Dimensional Semiconductors. <i>ACS Nano</i> , 2018, 12, 965-975.	14.6	172
3	Systematic study of electronic structure and band alignment of monolayer transition metal dichalcogenides in Van der Waals heterostructures. <i>2D Materials</i> , 2017, 4, 015026.	4.4	160
4	Alloying conducting channels for reliable neuromorphic computing. <i>Nature Nanotechnology</i> , 2020, 15, 574-579.	31.5	160
5	Pocketlike Active Site of Rh ₁ /MoS ₂ Single-Atom Catalyst for Selective Crotonaldehyde Hydrogenation. <i>Journal of the American Chemical Society</i> , 2019, 141, 19289-19295.	13.7	141
6	Charge Mediated Reversible Metal-Insulator Transition in Monolayer MoTe ₂ and W _{1-x} Mo _x Te ₂ Alloy. <i>ACS Nano</i> , 2016, 10, 7370-7375.	14.6	133
7	Site-dependent multicomponent doping strategy for Ni-rich LiNi _{1-2y} Co _y Mn _y O ₂ ($y = 1/12$) cathode materials for Li-ion batteries. <i>Journal of Materials Chemistry A</i> , 2017, 5, 25303-25313.	10.3	119
8	Nucleation and growth of WSe ₂ : enabling large grain transition metal dichalcogenides. <i>2D Materials</i> , 2017, 4, 045019.	4.4	96
9	Quantum-Confined Electronic States Arising from the Moiré Pattern of MoS ₂ -WSe ₂ Heterobilayers. <i>Nano Letters</i> , 2018, 18, 1849-1855.	9.1	91
10	A kinetic Monte Carlo simulation method of van der Waals epitaxy for atomistic nucleation-growth processes of transition metal dichalcogenides. <i>Scientific Reports</i> , 2017, 7, 2977.	3.3	72
11	Graphene-assisted spontaneous relaxation towards dislocation-free heteroepitaxy. <i>Nature Nanotechnology</i> , 2020, 15, 272-276.	31.5	71
12	Stable and Active Oxidation Catalysis by Cooperative Lattice Oxygen Redox on SmMn ₂ O ₅ Mullite Surface. <i>Journal of the American Chemical Society</i> , 2019, 141, 10722-10728.	13.7	64
13	Quantum Transport and Band Structure Evolution under High Magnetic Field in Few-Layer Tellurene. <i>Nano Letters</i> , 2018, 18, 5760-5767.	9.1	60
14	First principles kinetic Monte Carlo study on the growth patterns of WSe ₂ monolayer. <i>2D Materials</i> , 2016, 3, 025029.	4.4	59
15	Obstacles toward unity efficiency of LiNi _{1-2x} Co _x Mn _x O ₂ ($x = 1/4, 1/3$) (NCM) cathode materials: Insights from ab initio calculations. <i>Journal of Power Sources</i> , 2017, 340, 217-228.	7.8	57
16	Ab Initio Study on Surface Segregation and Anisotropy of Ni-Rich LiNi _{1-2y} Co _y Mn _y O ₂ (NCM) ($y \approx 0.1$) Cathodes. <i>ACS Applied Materials & Interfaces</i> , 2018, 10, 6673-6680.	8.0	50
17	Controlling nucleation of monolayer WSe ₂ during metal-organic chemical vapor deposition growth. <i>2D Materials</i> , 2016, 3, 025015.	4.4	42
18	Higher superconducting transition temperature by breaking the universal pressure relation. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 2004-2008.	7.1	39

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19	Flat Bands and Mechanical Deformation Effects in the Moiré Superlattice of MoS ₂ -WSe ₂ Heterobilayers. ACS Nano, 2020, 14, 7564-7573.	14.6	38
20	Dislocation driven spiral and non-spiral growth in layered chalcogenides. Nanoscale, 2018, 10, 15023-15034.	5.6	24
21	Tuning electronic transport in epitaxial graphene-based van der Waals heterostructures. Nanoscale, 2016, 8, 8947-8954.	5.6	21
22	Defect-mediated ripening of core-shell nanostructures. Nature Communications, 2022, 13, 2211.	12.8	17
23	Ligand-induced reduction concerted with coating by atomic layer deposition on the example of TiO ₂ -coated magnetite nanoparticles. Chemical Science, 2019, 10, 2171-2178.	7.4	11
24	Theoretical Demonstration of the Ionic Barristor. Nano Letters, 2016, 16, 2090-2095.	9.1	9
25	WSe ₂ homojunctions and quantum dots created by patterned hydrogenation of epitaxial graphene substrates. 2D Materials, 2019, 6, 021001.	4.4	7
26	Chemical and physical adsorption of a H ₂ O molecule on a metal doped Zr (0001) surface. Journal of Nuclear Materials, 2014, 452, 493-499.	2.7	6
27	Chemisorption of a hydrogen adatom on metal doped $\hat{\pm}$ -Zr (0001) surfaces in a vacuum and an implicit solvation environment. Nuclear Materials and Energy, 2017, 13, 28-34.	1.3	6
28	Characteristics of Interlayer Tunneling Field-Effect Transistors Computed by a $\hat{\pm}$ DFT-Bardeen Method. Journal of Electronic Materials, 2017, 46, 1378-1389.	2.2	5
29	First principles study of the Mn-doping effect on the physical and chemical properties of mullite-family Al ₂ SiO ₅ . Physical Chemistry Chemical Physics, 2017, 19, 24991-25001.	2.8	5
30	Charge-transfer modified embedded atom method dynamic charge potential for Li-Co-O system. Journal of Physics Condensed Matter, 2017, 29, 475903.	1.8	3
31	A new route of synthesizing atomically thin 2D materials embedded in bulk oxides. Journal of Applied Physics, 2021, 130, 035302.	2.5	0