

Wenchang Lu

List of Publications by Year in descending order

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Version: 2024-02-01

24
papers

930
citations

567281

15
h-index

610901

24
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24
all docs

24
docs citations

24
times ranked

1206
citing authors

#	ARTICLE	IF	CITATIONS
1	Local manifestations of thickness-dependent topology and edge states in the topological magnet <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:msub><mml:mi>MnBi</mml:mi><mml:mn>2</mml:mn></mml:msub></mml:mrow></mml:math> Physical Review B, 2022, 105, .	3.2	12
2	Relaxor ferroelectric polymer exhibits ultrahigh electromechanical coupling at low electric field. Science, 2022, 375, 1418-1422.	12.6	74
3	<i>Ab initio</i> simulations of metal contacts for graphene-based devices. Journal of Applied Physics, 2022, 131, 214301.	2.5	2
4	Study of Anharmonicity in Zirconium Hydrides Using Inelastic Neutron Scattering and Ab-Initio Computer Modeling. Inorganics, 2021, 9, 29.	2.7	3
5	On-surface cyclodehydrogenation reaction pathway determined by selective molecular deuterations. Chemical Science, 2021, 12, 15637-15644.	7.4	11
6	Structural Insight in the Interfacial Effect in Ferroelectric Polymer Nanocomposites. Advanced Materials, 2020, 32, e2005431.	21.0	84
7	Chirality-induced relaxor properties in ferroelectric polymers. Nature Materials, 2020, 19, 1169-1174.	27.5	93
8	Identification of Efficient Single-Atom Catalysts Based on V₂CO₂ MXene by <i>ab Initio</i> Simulations. Journal of Physical Chemistry C, 2020, 124, 4090-4100.	3.1	31
9	Engineering Edge States of Graphene Nanoribbons for Narrow-Band Photoluminescence. ACS Nano, 2020, 14, 5090-5098.	14.6	27
10	Large-Scale Phonon Calculations Using the Real-Space Multigrid Method. Journal of Chemical Theory and Computation, 2019, 15, 6859-6864.	5.3	9
11	Ab initio investigation of the cyclodehydrogenation process for polyanthrylene transformation to graphene nanoribbons. Npj Computational Materials, 2019, 5, .	8.7	9
12	Step edge-mediated assembly of periodic arrays of long graphene nanoribbons on Au(111). Chemical Communications, 2019, 55, 11848-11851.	4.1	14
13	Insights into the Morphotropic Phase Boundary in Ferroelectric Polymers from the Molecular Perspective. Journal of Physical Chemistry C, 2019, 123, 8727-8730.	3.1	16
14	Design of Atomically Precise Nanoscale Negative Differential Resistance Devices. Advanced Theory and Simulations, 2019, 2, 1800172.	2.8	18
15	Direct writing of heterostructures in single atomically precise graphene nanoribbons. Physical Review Materials, 2019, 3, .	2.4	18
16	Ferroelectric polymers exhibiting behaviour reminiscent of a morphotropic phase boundary. Nature, 2018, 562, 96-100.	27.8	200
17	Oxidization stability of atomically precise graphene nanoribbons. Physical Review Materials, 2018, 2, .	2.4	25
18	Controllable conversion of quasi-freestanding polymer chains to graphene nanoribbons. Nature Communications, 2017, 8, 14815.	12.8	58

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19	Selective sensing of ethylene and glucose using carbon-nanotube-based sensors: an ab initio investigation. <i>Nanoscale</i> , 2017, 9, 1687-1698.	5.6	33
20	Seamless Staircase Electrical Contact to Semiconducting Graphene Nanoribbons. <i>Nano Letters</i> , 2017, 17, 6241-6247.	9.1	64
21	Implementation of ultrasoft pseudopotentials in large-scale grid-based electronic structure calculations. <i>Physical Review B</i> , 2007, 76, .	3.2	23
22	Density functional theory studies of quantum transport in molecular systems. <i>International Journal of Quantum Chemistry</i> , 2006, 106, 3334-3342.	2.0	4
23	Nonequilibrium Quantum Transport Properties of Organic Molecules on Silicon. <i>Physical Review Letters</i> , 2005, 95, 206805.	7.8	65
24	Optical Anisotropy of the SiC(001)-(3 \times 3) Surface: Evidence for the Two-Adlayer Asymmetric-Dimer Model. <i>Physical Review Letters</i> , 2000, 85, 4381-4384.	7.8	37