

# 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  
g-index

24  
all docs

24  
docs citations

24  
times ranked

1206  
citing authors

#	ARTICLE	IF	CITATIONS
1	Ferroelectric polymers exhibiting behaviour reminiscent of a morphotropic phase boundary. <i>Nature</i> , 2018, 562, 96-100.	27.8	200
2	Chirality-induced relaxor properties in ferroelectric polymers. <i>Nature Materials</i> , 2020, 19, 1169-1174.	27.5	93
3	Structural Insight in the Interfacial Effect in Ferroelectric Polymer Nanocomposites. <i>Advanced Materials</i> , 2020, 32, e2005431.	21.0	84
4	Relaxor ferroelectric polymer exhibits ultrahigh electromechanical coupling at low electric field. <i>Science</i> , 2022, 375, 1418-1422.	12.6	74
5	Nonequilibrium Quantum Transport Properties of Organic Molecules on Silicon. <i>Physical Review Letters</i> , 2005, 95, 206805.	7.8	65
6	Seamless Staircase Electrical Contact to Semiconducting Graphene Nanoribbons. <i>Nano Letters</i> , 2017, 17, 6241-6247.	9.1	64
7	Controllable conversion of quasi-freestanding polymer chains to graphene nanoribbons. <i>Nature Communications</i> , 2017, 8, 14815.	12.8	58
8	Optical Anisotropy of the SiC(001)-(3Å <sup>-2</sup> ) Surface: Evidence for the Two-Adlayer Asymmetric-Dimer Model. <i>Physical Review Letters</i> , 2000, 85, 4381-4384.	7.8	37
9	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
10	Identification of Efficient Single-Atom Catalysts Based on V <sub>2</sub> CO <sub>2</sub> MXene by <i>ab Initio</i> Simulations. <i>Journal of Physical Chemistry C</i> , 2020, 124, 4090-4100.	3.1	31
11	Engineering Edge States of Graphene Nanoribbons for Narrow-Band Photoluminescence. <i>ACS Nano</i> , 2020, 14, 5090-5098.	14.6	27
12	Oxidization stability of atomically precise graphene nanoribbons. <i>Physical Review Materials</i> , 2018, 2, .	2.4	25
13	Implementation of ultrasoft pseudopotentials in large-scale grid-based electronic structure calculations. <i>Physical Review B</i> , 2007, 76, .	3.2	23
14	Design of Atomically Precise Nanoscale Negative Differential Resistance Devices. <i>Advanced Theory and Simulations</i> , 2019, 2, 1800172.	2.8	18
15	Direct writing of heterostructures in single atomically precise graphene nanoribbons. <i>Physical Review Materials</i> , 2019, 3, .	2.4	18
16	Insights into the Morphotropic Phase Boundary in Ferroelectric Polymers from the Molecular Perspective. <i>Journal of Physical Chemistry C</i> , 2019, 123, 8727-8730.	3.1	16
17	Step edge-mediated assembly of periodic arrays of long graphene nanoribbons on Au(111). <i>Chemical Communications</i> , 2019, 55, 11848-11851.	4.1	14
18	Local manifestations of thickness-dependent topology and edge states in the topological magnet $\text{MnBi}_2$ <i>Physical Review B</i> , 2022, 105, .	3.2	12

#	ARTICLE	IF	CITATIONS
19	On-surface cyclodehydrogenation reaction pathway determined by selective molecular deuterations. <i>Chemical Science</i> , 2021, 12, 15637-15644.	7.4	11
20	Large-Scale Phonon Calculations Using the Real-Space Multigrid Method. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 6859-6864.	5.3	9
21	Ab initio investigation of the cyclodehydrogenation process for polyanthrylene transformation to graphene nanoribbons. <i>Npj Computational Materials</i> , 2019, 5, .	8.7	9
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	Study of Anharmonicity in Zirconium Hydrides Using Inelastic Neutron Scattering and Ab-Initio Computer Modeling. <i>Inorganics</i> , 2021, 9, 29.	2.7	3
24	<i>Ab initio</i> simulations of metal contacts for graphene-based devices. <i>Journal of Applied Physics</i> , 2022, 131, 214301.	2.5	2