

Kunpeng Yuan

List of Publications by Year in descending order

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Version: 2024-02-01

23
papers

469
citations

687363

13
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713466

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23
docs citations

23
times ranked

495
citing authors

#	ARTICLE	IF	CITATIONS
1	Ultra-low thermal conductivity and high thermoelectric performance of two-dimensional triphosphides (InP ₃ , GaP ₃ , SbP ₃ and SnP ₃): a comprehensive first-principles study. <i>Nanoscale</i> , 2020, 12, 3330-3342.	5.6	68
2	Effects of tensile strain and finite size on thermal conductivity in monolayer WSe ₂ . <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 468-477.	2.8	60
3	Anomalous pressure effect on the thermal conductivity of ZnO, GaN, and AlN from first-principles calculations. <i>Physical Review B</i> , 2018, 98, .	3.2	42
4	A first-principles study of the thermoelectric properties of rhombohedral GeSe. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 1911-1922.	2.8	32
5	Lattice Thermal Conductivity Prediction Using Symbolic Regression and Machine Learning. <i>Journal of Physical Chemistry A</i> , 2021, 125, 435-450.	2.5	32
6	Disparate strain response of the thermal transport properties of bilayer penta-graphene as compared to that of monolayer penta-graphene. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 15647-15655.	2.8	28
7	Self-Assembled Monolayers for the Polymer/Semiconductor Interface with Improved Interfacial Thermal Management. <i>ACS Applied Materials & Interfaces</i> , 2019, 11, 42708-42714.	8.0	27
8	Predicting Elastic Properties of Materials from Electronic Charge Density Using 3D Deep Convolutional Neural Networks. <i>Journal of Physical Chemistry C</i> , 2020, 124, 17262-17273.	3.1	23
9	Tailoring phononic, electronic, and thermoelectric properties of orthorhombic GeSe through hydrostatic pressure. <i>Scientific Reports</i> , 2019, 9, 9490.	3.3	21
10	Efficient thermal conductivity modulation by manipulating interlayer interactions: A comparative study of bilayer graphene and graphite. <i>Journal of Applied Physics</i> , 2019, 126, .	2.5	21
11	Pressure tuning of the thermal conductivity of gallium arsenide from first-principles calculations. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 30331-30339.	2.8	16
12	Electric field tuned anisotropic to isotropic thermal transport transition in monolayer borophene without altering its atomic structure. <i>Nanoscale</i> , 2020, 12, 19178-19190.	5.6	15
13	First-principles calculations of interfacial thermal transport properties between SiC/Si substrates and compounds of boron with selected group V elements. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 6011-6020.	2.8	14
14	First-principles analysis of phonon thermal transport properties of two-dimensional WS ₂ /WSe ₂ heterostructures*. <i>Chinese Physics B</i> , 2021, 30, 034401.	1.4	13
15	Why thermal conductivity of CaO is lower than that of CaS: a study from the perspective of phonon splitting of optical mode. <i>Nanotechnology</i> , 2021, 32, 025709.	2.6	13
16	Ultralow lattice thermal conductivity and dramatically enhanced thermoelectric properties of monolayer InSe induced by an external electric field. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 13633-13646.	2.8	10
17	Zintl Phase Compounds Mg ₃ Sb ₂ ~xBix (x = 0, 1, and 2) Monolayers: Electronic, Phonon and Thermoelectric Properties From ab Initio Calculations. <i>Frontiers in Mechanical Engineering</i> , 2022, 8, .	1.8	7
18	The Abnormally Excellent Figure of Merit of 14,14,18-Graphyne at Room Temperature: A Study on the Thermoelectric Characteristic of Graphyne. <i>ACS Applied Energy Materials</i> , 2022, 5, 6363-6372.	5.1	6

#	ARTICLE	IF	CITATIONS
19	Thermal conductivity of SrTiO ₃ under high-pressure. Applied Physics Letters, 2022, 120, .	3.3	6
20	Alkaline treatment of used carbon-brush anodes for restoring power generation of microbial fuel cells. RSC Advances, 2018, 8, 36754-36760.	3.6	5
21	Low thermal conductivity of peanut-shaped carbon nanotube and its insensitive response to uniaxial strain. Nanotechnology, 2020, 31, 115701.	2.6	4
22	Giant Manipulation of Phonon Hydrodynamics in Ferroelectric Bilayer Boron Nitride at Room Temperature and Beyond. ACS Applied Energy Materials, 2022, 5, 8781-8790.	5.1	4
23	Potential thermoelectric materials: first-principles prediction of low lattice thermal conductivity of two-dimensional (2D) orthogonal ScX ₂ (X = C and N) compounds. Physical Chemistry Chemical Physics, 2021, 23, 23718-23729.	2.8	2