Rui Qin

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

39 2,579 19 41 g-index

41 2,810 5.6 4.63 ext. papers ext. citations avg, IF L-index

#	Paper	IF	Citations
39	High harmonic studies of structural phase transitions in silicon. <i>Computational Materials Science</i> , 2021 , 197, 110621	3.2	1
38	Probing structural chirality of crystals using high-order harmonic generation in solids. <i>Physical Review A</i> , 2020 , 101,	2.6	3
37	High harmonic generation in grapheneBoron nitride heterostructures. <i>Journal of Materials Chemistry C</i> , 2020 , 8, 12085-12091	7.1	6
36	Strong-field nonlinear optical properties of monolayer black phosphorus. <i>Nanoscale</i> , 2019 , 11, 16377-16	6 38 3	9
35	Circularly polarized extreme ultraviolet high harmonic generation in graphene. <i>Optics Express</i> , 2019 , 27, 3761-3770	3.3	26
34	Strain-controlled high harmonic generation with Dirac fermions in silicene. <i>Nanoscale</i> , 2018 , 10, 22593-2	2 ≱6 00	12
33	Sign-tunable Poissonঙ ratio in semi-fluorinated graphene. <i>Nanoscale</i> , 2017 , 9, 128-133	7.7	23
32	Uniaxial strain-induced mechanical and electronic property modulation of silicene. <i>Nanoscale Research Letters</i> , 2014 , 9, 521	5	59
31	First-principles calculations of mechanical and electronic properties of silicene under strain. <i>AIP Advances</i> , 2012 , 2, 022159	1.5	124
30	Tunable bandgap in silicene and germanene. <i>Nano Letters</i> , 2012 , 12, 113-8	11.5	1043
29	Tunable and sizable band gap of single-layer graphene sandwiched between hexagonal boron nitride. <i>NPG Asia Materials</i> , 2012 , 4, e6-e6	10.3	140
28	Polarized Nonresonant Raman Spectra of Graphene Nanoribbons. <i>Journal of Physical Chemistry C</i> , 2011 , 115, 24463-24468	3.8	9
27	Functionalized graphene for high-performance two-dimensional spintronics devices. <i>ACS Nano</i> , 2011 , 5, 2601-10	16.7	101
26	Structure and Electronic and Transport Properties of Transition Metal Intercalated Graphene and Graphene-Hexagonal-Boron-Nitride Bilayer. <i>Journal of Physical Chemistry C</i> , 2011 , 115, 25273-25280	3.8	21
25	Electric-Field-Induced Energy Gap in Few-Layer Graphene. <i>Journal of Physical Chemistry C</i> , 2011 , 115, 9458-9464	3.8	66
24	Family-Dependent Rectification Characteristics in Ultra-Short Graphene NanoribbonpfiJunctions. Journal of Physical Chemistry C, 2011 , 115, 8547-8554	3.8	27
23	Negative differential resistance in parallel single-walled carbon nanotube contacts. <i>Physical Review B</i> , 2011 , 83,	3.3	30

(2009-2011)

22	Quasiparticle energies and excitonic effects of the two-dimensional carbon allotrope graphdiyne: Theory and experiment. <i>Physical Review B</i> , 2011 , 84,	3.3	246
21	All-Metallic High-Performance Field Effect Transistor Based on Telescoping Carbon Nanotubes: An ab Initio Study. <i>Journal of Physical Chemistry C</i> , 2011 , 115, 6933-6938	3.8	20
20	Negative rectification and negative differential resistance in nanoscale single-walled carbon nanotube p-n junctions. <i>Theoretical Chemistry Accounts</i> , 2011 , 130, 353-359	1.9	9
19	Tuning graphene nanoribbon field effect transistors via controlling doping level. <i>Theoretical Chemistry Accounts</i> , 2011 , 130, 483-489	1.9	4
18	Room-temperature giant magnetoresistance over one billion percent in a bare graphene nanoribbon device. <i>Physical Review B</i> , 2010 , 81,	3.3	43
17	Electronic Structure and Stability of Ultranarrow Single-Layer SnS2 Nanoribbons: A First-Principles Study. <i>Journal of Computational and Theoretical Nanoscience</i> , 2010 , 7, 2100-2103	0.3	10
16	Polarized Vibrational Infrared Absorption of Graphene Nanoribbons. <i>Journal of Physical Chemistry C</i> , 2010 , 114, 6959-6965	3.8	10
15	Structural, Electronic, and Transport Properties of Gd/Eu Atomic Chains Encapsulated in Single-Walled Carbon Nanotubes. <i>Journal of Physical Chemistry C</i> , 2010 , 114, 15347-15353	3.8	15
14	Functionalized Metallic Single-Walled Carbon Nanotubes as a High-Performance Single-Molecule Organic Field Effect Transistor: An ab Initio Study. <i>Journal of Physical Chemistry C</i> , 2010 , 114, 15816-158	822 ⁸	12
13	Selection of single-walled carbon nanotubes according to both their diameter and chirality via nanotweezers. <i>Nano Research</i> , 2010 , 3, 296-306	10	12
12	First-principles study of the formation mechanisms of nitrogen molecule in annealed ZnO. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2010 , 374, 3546-3550	2.3	19
11	Electron localization and emission mechanism in wurtzite (Al, In, Ga)N alloys. <i>Physica Status Solidi</i> (B): Basic Research, 2010 , 247, 109-114	1.3	14
10	Electronic and Mechanical Coupling in Bent ZnO Nanowires. Advanced Materials, 2009, 21, 4937-4941	24	128
9	Magnetism in carbon nanoscrolls: Quasi-half-metals and half-metals in pristine hydrocarbons. <i>Nano Research</i> , 2009 , 2, 844-850	10	10
8	Optical Absorption Spectra of Charge-Doped Single-Walled Carbon Nanotubes from First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 7058-7064	3.8	2
7	Magnetic Properties of Fully Bare and Half-Bare Boron Nitride Nanoribbons. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 2273-2276	3.8	96
6	Origin of p-Type Doping in Zinc Oxide Nanowires Induced by Phosphorus Doping: A First Principles Study. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 9541-9545	3.8	26
5	Study on the dispersion of charged single-wall carbon nanotube bundles by first principles calculation. <i>Journal of Nanoscience and Nanotechnology</i> , 2009 , 9, 5170-2	1.3	2

4	Electronic-type- and diameter-dependent reduction of single-walled carbon nanotubes induced by adsorption of electron-donor molecules. <i>Small</i> , 2009 , 5, 244-55	11	10
3	Novel one-dimensional organometallic half metals: vanadium-cyclopentadienyl, vanadium-cyclopentadienyl-benzene, and vanadium-anthracene wires. <i>Nano Letters</i> , 2008 , 8, 3640-4	11.5	116
2	Optical Absorption Spectra and Polarizabilities of Silicon Carbide Nanotubes: A First Principles Study. <i>Journal of Physical Chemistry C</i> , 2007 , 111, 18864-18870	3.8	11
1	Why semiconducting single-walled carbon nanotubes are separated from their metallic counterparts. <i>Small</i> , 2007 , 3, 1566-76	11	62