

# Rui Qin

## List of Publications by Citations

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39  
papers

2,579  
citations

19  
h-index

41  
g-index

41  
ext. papers

2,810  
ext. citations

5.6  
avg, IF

4.63  
L-index

#	Paper	IF	Citations
39	Tunable bandgap in silicene and germanene. <i>Nano Letters</i> , <b>2012</b> , 12, 113-8	11.5	1043
38	Quasiparticle energies and excitonic effects of the two-dimensional carbon allotrope graphdiyne: Theory and experiment. <i>Physical Review B</i> , <b>2011</b> , 84,	3.3	246
37	Tunable and sizable band gap of single-layer graphene sandwiched between hexagonal boron nitride. <i>NPG Asia Materials</i> , <b>2012</b> , 4, e6-e6	10.3	140
36	Electronic and Mechanical Coupling in Bent ZnO Nanowires. <i>Advanced Materials</i> , <b>2009</b> , 21, 4937-4941	24	128
35	First-principles calculations of mechanical and electronic properties of silicene under strain. <i>AIP Advances</i> , <b>2012</b> , 2, 022159	1.5	124
34	Novel one-dimensional organometallic half metals: vanadium-cyclopentadienyl, vanadium-cyclopentadienyl-benzene, and vanadium-anthracene wires. <i>Nano Letters</i> , <b>2008</b> , 8, 3640-4	11.5	116
33	Functionalized graphene for high-performance two-dimensional spintronics devices. <i>ACS Nano</i> , <b>2011</b> , 5, 2601-10	16.7	101
32	Magnetic Properties of Fully Bare and Half-Bare Boron Nitride Nanoribbons. <i>Journal of Physical Chemistry C</i> , <b>2009</b> , 113, 2273-2276	3.8	96
31	Electric-Field-Induced Energy Gap in Few-Layer Graphene. <i>Journal of Physical Chemistry C</i> , <b>2011</b> , 115, 9458-9464	3.8	66
30	Why semiconducting single-walled carbon nanotubes are separated from their metallic counterparts. <i>Small</i> , <b>2007</b> , 3, 1566-76	11	62
29	Uniaxial strain-induced mechanical and electronic property modulation of silicene. <i>Nanoscale Research Letters</i> , <b>2014</b> , 9, 521	5	59
28	Room-temperature giant magnetoresistance over one billion percent in a bare graphene nanoribbon device. <i>Physical Review B</i> , <b>2010</b> , 81,	3.3	43
27	Negative differential resistance in parallel single-walled carbon nanotube contacts. <i>Physical Review B</i> , <b>2011</b> , 83,	3.3	30
26	Family-Dependent Rectification Characteristics in Ultra-Short Graphene Nanoribbon p-n Junctions. <i>Journal of Physical Chemistry C</i> , <b>2011</b> , 115, 8547-8554	3.8	27
25	Origin of p-Type Doping in Zinc Oxide Nanowires Induced by Phosphorus Doping: A First Principles Study. <i>Journal of Physical Chemistry C</i> , <b>2009</b> , 113, 9541-9545	3.8	26
24	Circularly polarized extreme ultraviolet high harmonic generation in graphene. <i>Optics Express</i> , <b>2019</b> , 27, 3761-3770	3.3	26
23	Sign-tunable Poisson's ratio in semi-fluorinated graphene. <i>Nanoscale</i> , <b>2017</b> , 9, 128-133	7.7	23

22	Structure and Electronic and Transport Properties of Transition Metal Intercalated Graphene and Graphene-Hexagonal-Boron-Nitride Bilayer. <i>Journal of Physical Chemistry C</i> , <b>2011</b> , 115, 25273-25280	3.8	21
21	All-Metallic High-Performance Field Effect Transistor Based on Telescoping Carbon Nanotubes: An ab Initio Study. <i>Journal of Physical Chemistry C</i> , <b>2011</b> , 115, 6933-6938	3.8	20
20	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> , <b>2010</b> , 374, 3546-3550	2.3	19
19	Structural, Electronic, and Transport Properties of Gd/Eu Atomic Chains Encapsulated in Single-Walled Carbon Nanotubes. <i>Journal of Physical Chemistry C</i> , <b>2010</b> , 114, 15347-15353	3.8	15
18	Electron localization and emission mechanism in wurtzite (Al, In, Ga)N alloys. <i>Physica Status Solidi (B): Basic Research</i> , <b>2010</b> , 247, 109-114	1.3	14
17	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> , <b>2010</b> , 114, 15816-15822	3.8	12
16	Selection of single-walled carbon nanotubes according to both their diameter and chirality via nanotweezers. <i>Nano Research</i> , <b>2010</b> , 3, 296-306	10	12
15	Strain-controlled high harmonic generation with Dirac fermions in silicene. <i>Nanoscale</i> , <b>2018</b> , 10, 22593-22600	7.6	12
14	Optical Absorption Spectra and Polarizabilities of Silicon Carbide Nanotubes: A First Principles Study. <i>Journal of Physical Chemistry C</i> , <b>2007</b> , 111, 18864-18870	3.8	11
13	Electronic Structure and Stability of Ultranarrow Single-Layer SnS <sub>2</sub> Nanoribbons: A First-Principles Study. <i>Journal of Computational and Theoretical Nanoscience</i> , <b>2010</b> , 7, 2100-2103	0.3	10
12	Polarized Vibrational Infrared Absorption of Graphene Nanoribbons. <i>Journal of Physical Chemistry C</i> , <b>2010</b> , 114, 6959-6965	3.8	10
11	Magnetism in carbon nanoscrolls: Quasi-half-metals and half-metals in pristine hydrocarbons. <i>Nano Research</i> , <b>2009</b> , 2, 844-850	10	10
10	Electronic-type- and diameter-dependent reduction of single-walled carbon nanotubes induced by adsorption of electron-donor molecules. <i>Small</i> , <b>2009</b> , 5, 244-55	11	10
9	Strong-field nonlinear optical properties of monolayer black phosphorus. <i>Nanoscale</i> , <b>2019</b> , 11, 16377-16383	7.3	9
8	Polarized Nonresonant Raman Spectra of Graphene Nanoribbons. <i>Journal of Physical Chemistry C</i> , <b>2011</b> , 115, 24463-24468	3.8	9
7	Negative rectification and negative differential resistance in nanoscale single-walled carbon nanotube p-n junctions. <i>Theoretical Chemistry Accounts</i> , <b>2011</b> , 130, 353-359	1.9	9
6	High harmonic generation in grapheneBoron nitride heterostructures. <i>Journal of Materials Chemistry C</i> , <b>2020</b> , 8, 12085-12091	7.1	6
5	Tuning graphene nanoribbon field effect transistors via controlling doping level. <i>Theoretical Chemistry Accounts</i> , <b>2011</b> , 130, 483-489	1.9	4

4	Probing structural chirality of crystals using high-order harmonic generation in solids. <i>Physical Review A</i> , <b>2020</b> , 101,	2.6	3
3	Optical Absorption Spectra of Charge-Doped Single-Walled Carbon Nanotubes from First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , <b>2009</b> , 113, 7058-7064	3.8	2
2	Study on the dispersion of charged single-wall carbon nanotube bundles by first principles calculation. <i>Journal of Nanoscience and Nanotechnology</i> , <b>2009</b> , 9, 5170-2	1.3	2
1	High harmonic studies of structural phase transitions in silicon. <i>Computational Materials Science</i> , <b>2021</b> , 197, 110621	3.2	1