

Zhenyu Li

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179
papers

8,680
citations

48
h-index

89
g-index

189
ext. papers

9,776
ext. citations

6.7
avg, IF

6.37
L-index

#	Paper	IF	Citations
179	Half-metallicity in edge-modified zigzag graphene nanoribbons. <i>Journal of the American Chemical Society</i> , 2008 , 130, 4224-5	16.4	587
178	Epitaxial Growth of Single Layer Blue Phosphorus: A New Phase of Two-Dimensional Phosphorus. <i>Nano Letters</i> , 2016 , 16, 4903-8	11.5	490
177	Low-temperature growth of graphene by chemical vapor deposition using solid and liquid carbon sources. <i>ACS Nano</i> , 2011 , 5, 3385-90	16.7	304
176	How graphene is cut upon oxidation?. <i>Journal of the American Chemical Society</i> , 2009 , 131, 6320-1	16.4	289
175	Will zigzag graphene nanoribbon turn to half metal under electric field?. <i>Applied Physics Letters</i> , 2007 , 91, 243116	3.4	285
174	First-Principles Thermodynamics of Graphene Growth on Cu Surfaces. <i>Journal of Physical Chemistry C</i> , 2011 , 115, 17782-17787	3.8	276
173	Doping-induced structural phase transition in cobalt diselenide enables enhanced hydrogen evolution catalysis. <i>Nature Communications</i> , 2018 , 9, 2533	17.4	244
172	In Situ Thermal Atomization To Convert Supported Nickel Nanoparticles into Surface-Bound Nickel Single-Atom Catalysts. <i>Angewandte Chemie - International Edition</i> , 2018 , 57, 14095-14100	16.4	206
171	Obtaining two-dimensional electron gas in free space without resorting to electron doping: an electrified based design. <i>Journal of the American Chemical Society</i> , 2014 , 136, 13313-8	16.4	204
170	Electronic structures of SiC nanoribbons. <i>Journal of Chemical Physics</i> , 2008 , 129, 174114	3.9	199
169	Oxygen molecule dissociation on carbon nanostructures with different types of nitrogen doping. <i>Nanoscale</i> , 2012 , 4, 1184-9	7.7	195
168	Silicene as a highly sensitive molecule sensor for NH ₃ , NO and NO ₂ . <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 6957-62	3.6	173
167	A first-principles study of gas adsorption on germanene. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 22495-8	3.6	172
166	Proposed photosynthesis method for producing hydrogen from dissociated water molecules using incident near-infrared light. <i>Physical Review Letters</i> , 2014 , 112, 018301	7.4	163
165	Bipolar magnetic semiconductors: a new class of spintronics materials. <i>Nanoscale</i> , 2012 , 4, 5680-5	7.7	162
164	Adsorption energies of molecular oxygen on Au clusters. <i>Journal of Chemical Physics</i> , 2004 , 120, 9594-6009	3.9	149
163	Site-specific photocatalytic splitting of methanol on TiO ₂ (110). <i>Chemical Science</i> , 2010 , 1, 575	9.4	143

162	Half-metallicity in hybrid BCN nanoribbons. <i>Journal of Chemical Physics</i> , 2008 , 129, 084712	3.9	129
161	Mo2C nanoparticles embedded within bacterial cellulose-derived 3D N-doped carbon nanofiber networks for efficient hydrogen evolution. <i>NPG Asia Materials</i> , 2016 , 8, e288-e288	10.3	127
160	Electronic Structure Engineering via On-Plane Chemical Functionalization: A Comparison Study on Two-Dimensional Polysilane and Graphane. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 16741-16746	3.8	122
159	Electron-phonon coupling in a boron-doped diamond superconductor. <i>Physical Review B</i> , 2004 , 70,	3.3	122
158	Porous silicene as a hydrogen purification membrane. <i>Physical Chemistry Chemical Physics</i> , 2013 , 15, 5753-7	3.7	110
157	Remarkable chemical adsorption of manganese-doped titanate for direct carbon dioxide electrolysis. <i>Journal of Materials Chemistry A</i> , 2014 , 2, 6904-6915	13	101
156	Single-molecule chemistry of metal phthalocyanine on noble metal surfaces. <i>Accounts of Chemical Research</i> , 2010 , 43, 954-62	24.3	98
155	Growth of Quasi-Free-Standing Single-Layer Blue Phosphorus on Tellurium Monolayer Functionalized Au(111). <i>ACS Nano</i> , 2017 , 11, 4943-4949	16.7	92
154	Why the Band Gap of Graphene Is Tunable on Hexagonal Boron Nitride. <i>Journal of Physical Chemistry C</i> , 2012 , 116, 3142-3146	3.8	88
153	Oxidation states of graphene: insights from computational spectroscopy. <i>Journal of Chemical Physics</i> , 2009 , 131, 244505	3.9	84
152	Pt Single Atoms Embedded in the Surface of Ni Nanocrystals as Highly Active Catalysts for Selective Hydrogenation of Nitro Compounds. <i>Nano Letters</i> , 2018 , 18, 3785-3791	11.5	84
151	Structure of Graphene Oxide: Thermodynamics versus Kinetics. <i>Journal of Physical Chemistry C</i> , 2011 , 115, 11991-11995	3.8	83
150	Helium separation via porous silicene based ultimate membrane. <i>Nanoscale</i> , 2013 , 5, 9062-6	7.7	82
149	Electronic and optical properties of graphene and graphitic ZnO nanocomposite structures. <i>Journal of Chemical Physics</i> , 2013 , 138, 124706	3.9	82
148	Aerobic Oxidation of Cyclohexane on Catalysts Based on Twinned and Single-Crystal Au75Pd25 Bimetallic Nanocrystals. <i>Nano Letters</i> , 2015 , 15, 2875-80	11.5	81
147	Lattice mismatch induced nonlinear growth of graphene. <i>Journal of the American Chemical Society</i> , 2012 , 134, 6045-51	16.4	80
146	Structural, electronic, and optical properties of hybrid silicene and graphene nanocomposite. <i>Journal of Chemical Physics</i> , 2013 , 139, 154704	3.9	79
145	Theoretical study of nitric oxide adsorption on Au clusters. <i>Journal of Chemical Physics</i> , 2004 , 121, 2558-62	3.2	74

144	Dithiocarbamate anchoring in molecular wire junctions: a first principles study. <i>Journal of Physical Chemistry B</i> , 2006 , 110, 9893-8	3.4	72
143	Nature of well-defined conductance of amine-anchored molecular junctions: Density functional calculations. <i>Physical Review B</i> , 2007 , 76,	3.3	71
142	Precursor Triggering Synthesis of Self-Coupled Sulfide Polymorphs with Enhanced Photoelectrochemical Properties. <i>Journal of the American Chemical Society</i> , 2016 , 138, 12913-12919	16.4	69
141	Communication: Coalescence of carbon atoms on Cu (111) surface: Emergence of a stable bridging-metal structure motif. <i>Journal of Chemical Physics</i> , 2010 , 133, 071101	3.9	67
140	Is mayenite without clathrated oxygen an inorganic electride?. <i>Angewandte Chemie - International Edition</i> , 2004 , 43, 6479-82	16.4	63
139	Mechanisms of graphene growth on metal surfaces: theoretical perspectives. <i>Small</i> , 2014 , 10, 2136-50	11	62
138	Carbon dimers as the dominant feeding species in epitaxial growth and morphological phase transition of graphene on different Cu substrates. <i>Physical Review Letters</i> , 2015 , 114, 216102	7.4	61
137	MAGNETISM IN GRAPHENE SYSTEMS. <i>Nano</i> , 2008 , 03, 433-442	1.1	61
136	Contrasting Structural Reconstructions, Electronic Properties, and Magnetic Orderings along Different Edges of Zigzag Transition Metal Dichalcogenide Nanoribbons. <i>Nano Letters</i> , 2017 , 17, 1097-1101	11.5	60
135	Graphene Thickness Control via Gas-Phase Dynamics in Chemical Vapor Deposition. <i>Journal of Physical Chemistry C</i> , 2012 , 116, 10557-10562	3.8	59
134	Ratio-controlled synthesis of CuNi octahedra and nanocubes with enhanced catalytic activity. <i>Journal of the American Chemical Society</i> , 2015 , 137, 14027-30	16.4	57
133	A first-principles prediction of two-dimensional superconductivity in pristine B $\bar{1}$ single layers. <i>Nanoscale</i> , 2012 , 4, 3032-5	7.7	55
132	Hydrogenated bilayer wurtzite SiC nanofilms: a two-dimensional bipolar magnetic semiconductor material. <i>Physical Chemistry Chemical Physics</i> , 2013 , 15, 497-503	3.6	52
131	Diamondization of chemically functionalized graphene and graphene-BN bilayers. <i>Physical Chemistry Chemical Physics</i> , 2012 , 14, 8179-84	3.6	48
130	Dominant Kinetic Pathways of Graphene Growth in Chemical Vapor Deposition: The Role of Hydrogen. <i>Journal of Physical Chemistry C</i> , 2017 , 121, 25949-25955	3.8	47
129	Orbital interaction mechanisms of conductance enhancement and rectification by dithiocarboxylate anchoring group. <i>Journal of Physical Chemistry B</i> , 2006 , 110, 19116-20	3.4	47
128	A first-principles study of NO adsorption and oxidation on Au(111) surface. <i>Journal of Chemical Physics</i> , 2008 , 129, 134708	3.9	46
127	First principles nuclear magnetic resonance signatures of graphene oxide. <i>Journal of Chemical Physics</i> , 2010 , 133, 034502	3.9	45

126	Bilayer Graphene Growth via a Penetration Mechanism. <i>Journal of Physical Chemistry C</i> , 2014 , 118, 6201-6206	6.2	39
125	Diamond as an inert substrate of graphene. <i>Journal of Chemical Physics</i> , 2013 , 138, 054701	3.9	38
124	Discriminating early stage A{beta}42 monomer structures using chirality-induced 2DIR spectroscopy in a simulation study. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010 , 107, 15687-92	11.5	36
123	Inorganic electride: theoretical study on structural and electronic properties. <i>Journal of the American Chemical Society</i> , 2003 , 125, 6050-1	16.4	35
122	Gas-phase dynamics in graphene growth by chemical vapour deposition. <i>Physical Chemistry Chemical Physics</i> , 2015 , 17, 22832-6	3.6	34
121	Descriptor-Based Design Principle for Two-Dimensional Single-Atom Catalysts: Carbon Dioxide Electroreduction. <i>Journal of Physical Chemistry Letters</i> , 2020 , 11, 3481-3487	6.4	34
120	Single Faceted Two-Dimensional Mo2C Electrocatalyst for Highly Efficient Nitrogen Fixation. <i>ACS Catalysis</i> , 2020 , 10, 7864-7870	13.1	33
119	First Principles Study on the Geometric and Electronic Structures of the FeO/Pt(111) Surface. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 8302-8305	3.8	33
118	A density functional study on cationic Au(n)Cu(m)+ clusters and their monocarbonyls. <i>Physical Chemistry Chemical Physics</i> , 2009 , 11, 2329-34	3.6	32
117	Electride: from computational characterization to theoretical design. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2016 , 6, 430-440	7.9	31
116	Two-Dimensional Stoichiometric Boron Oxides as a Versatile Platform for Electronic Structure Engineering. <i>Journal of Physical Chemistry Letters</i> , 2017 , 8, 4347-4353	6.4	31
115	Hetero-site nucleation for growing twisted bilayer graphene with a wide range of twist angles. <i>Nature Communications</i> , 2021 , 12, 2391	17.4	31
114	In Situ Thermal Atomization To Convert Supported Nickel Nanoparticles into Surface-Bound Nickel Single-Atom Catalysts. <i>Angewandte Chemie</i> , 2018 , 130, 14291-14296	3.6	30
113	Synthesis of Monolayer Blue Phosphorus Enabled by Silicon Intercalation. <i>ACS Nano</i> , 2020 , 14, 3687-3695	6.7	28
112	Electronic Structures of Defective Boron Nitride Nanotubes under Transverse Electric Fields. <i>Journal of Physical Chemistry C</i> , 2008 , 112, 8424-8428	3.8	28
111	Probing electron correlations in molecules by two-dimensional coherent optical spectroscopy. <i>Journal of the American Chemical Society</i> , 2008 , 130, 3509-15	16.4	27
110	Two Dimensional Electronic Correlation Spectroscopy of the npi* and pipi* Protein Backbone Transitions: A Simulation Study. <i>Chemical Physics</i> , 2007 , 341, 29-36	2.3	27
109	Effects of discrete energy levels on single-electron tunneling in coupled metal particles. <i>Applied Physics Letters</i> , 2003 , 82, 3767-3769	3.4	27

108	Elementary process for CVD graphene on Cu(110): size-selective carbon clusters. <i>Scientific Reports</i> , 2014 , 4, 4431	4.9	26
107	Transition-Metal Diboride: A New Family of Two-Dimensional Materials Designed for Selective CO ₂ Electroreduction. <i>Journal of Physical Chemistry C</i> , 2019 , 123, 16294-16299	3.8	25
106	Rational Design of Two-dimensional Anode Materials: BS as a Strained Graphene. <i>Journal of Physical Chemistry Letters</i> , 2018 , 9, 4852-4856	6.4	25
105	Atomically thin semiconducting penta-PdP ₂ and PdAs ₂ with ultrahigh carrier mobility. <i>Journal of Materials Chemistry C</i> , 2018 , 6, 9055-9059	7.1	25
104	Boron K ₄ crystal: a stable chiral three-dimensional sp ² network. <i>Physical Chemistry Chemical Physics</i> , 2010 , 12, 12420-2	3.6	24
103	Coordinate activation in heterogeneous carbon dioxide reduction on Co-based molecular catalysts. <i>Applied Catalysis B: Environmental</i> , 2020 , 268, 118452	21.8	23
102	A first principles study on organic molecule encapsulated boron nitride nanotubes. <i>Journal of Chemical Physics</i> , 2008 , 128, 164701	3.9	22
101	Oxygen adsorption on Zr(0 0 1) surfaces: Density functional calculations and a multiple-layer adsorption model. <i>Surface Science</i> , 2008 , 602, 2212-2216	1.8	22
100	The Nanoparticle Size Effect in Graphene Cutting: A "Pac-Man" Mechanism. <i>Angewandte Chemie - International Edition</i> , 2016 , 55, 9918-21	16.4	22
99	Reversible Oxidation of Blue Phosphorus Monolayer on Au(111). <i>Nano Letters</i> , 2019 , 19, 5340-5346	11.5	21
98	Density Functional Study of Nonlinear Optical Properties of Grossly Warped Nanographene C ₈₀ H ₃₀ . <i>Journal of Physical Chemistry C</i> , 2014 , 118, 3313-3318	3.8	21
97	A Density Functional Study of the Nonlinear Optical Properties of Edge-Functionalized Nonplanar Nanographenes. <i>ChemPhysChem</i> , 2015 , 16, 2783-2788	3.2	21
96	Atomistic Simulations of Graphene Growth: From Kinetics to Mechanism. <i>Accounts of Chemical Research</i> , 2018 , 51, 728-735	24.3	20
95	Implementation of screened hybrid density functional for periodic systems with numerical atomic orbitals: basis function fitting and integral screening. <i>Journal of Chemical Physics</i> , 2011 , 135, 034110	3.9	20
94	Rectifying effect in polar conjugated molecular junctions: a first-principles study. <i>Journal of Nanoscience and Nanotechnology</i> , 2009 , 9, 774-8	1.3	20
93	Methanol-Selective Oxidation Pathways on Au Surfaces: A First-Principles Study. <i>Journal of Physical Chemistry C</i> , 2014 , 118, 17511-17520	3.8	19
92	Disorder and suppression of quantum confinement effects in Pd nanoparticles. <i>Physical Review Letters</i> , 2003 , 90, 246803	7.4	19
91	Structure of Blue Phosphorus Grown on Au(111) Surface Revisited. <i>Journal of Physical Chemistry C</i> , 2020 , 124, 2024-2029	3.8	19

90	Single-molecule imaging of activated nitrogen adsorption on individual manganese phthalocyanine. <i>Nano Letters</i> , 2015 , 15, 3181-8	11.5	18
89	Growth of boron nitride nanotubes from magnesium diboride catalysts. <i>Nanoscale</i> , 2018 , 10, 13895-13901	17	18
88	Density functional predictions of new silicon allotropes: Electronic properties and potential applications to Li-battery anode materials. <i>Solid State Communications</i> , 2011 , 151, 1228-1230	1.6	18
87	Geometrical, electronic, and magnetic properties of Na _{0.5} CoO ₂ from first principles. <i>Physical Review B</i> , 2005 , 71,	3.3	18
86	Distinct molecule adsorption behaviors on warped nanographene C ₈₀ H ₃₀ : A theoretical study. <i>Carbon</i> , 2016 , 100, 428-434	10.4	16
85	Surface and size effects on the charge state of NV center in nanodiamonds. <i>Computational and Theoretical Chemistry</i> , 2013 , 1021, 49-53	2	16
84	Diffusion and desorption of oxygen atoms on graphene. <i>Journal of Physics Condensed Matter</i> , 2013 , 25, 405301	1.8	16
83	Implementation of exact exchange with numerical atomic orbitals. <i>Journal of Physical Chemistry A</i> , 2010 , 114, 1039-43	2.8	16
82	Inorganic electrides. <i>Chemistry - A European Journal</i> , 2004 , 10, 1592-6	4.8	16
81	Interfacial properties of black phosphorus/transition metal carbide van der Waals heterostructures. <i>Frontiers of Physics</i> , 2018 , 13, 1	3.7	15
80	Proposal of a general scheme to obtain room-temperature spin polarization in asymmetric antiferromagnetic semiconductors. <i>Physical Review B</i> , 2015 , 92,	3.3	15
79	Linear scaling electronic structure calculations with numerical atomic basis set. <i>International Reviews in Physical Chemistry</i> , 2010 , 29, 665-691	7	15
78	Geometry and Excitation Energy Fluctuations of NMA in Aqueous Solution with CHARMM, AMBER, OPLS, and GROMOS Force Fields: Implications for Protein Ultraviolet Spectra Simulation. <i>Chemical Physics Letters</i> , 2008 , 452, 78-83	2.5	15
77	Water on silicene: A hydrogen bond-autocatalyzed physisorption-chemisorption-dissociation transition. <i>Nano Research</i> , 2017 , 10, 2223-2233	10	14
76	Detecting a Molecule-Surface Hybrid State by an Fe-Coated Tip with a Non-s-Like Orbital. <i>Journal of Physical Chemistry C</i> , 2008 , 112, 15603-15606	3.8	14
75	Electronic structures of organic molecule encapsulated BN nanotubes under transverse electric field. <i>Journal of Chemical Physics</i> , 2008 , 129, 024710	3.9	14
74	Phosphorus Nanostripe Arrays on Cu(110): A Case Study to Understand the Substrate Effect on the Phosphorus thin Film Growth. <i>Advanced Materials Interfaces</i> , 2017 , 4, 1601167	4.6	13
73	Adsorption and Diffusion of CO on Clean and CO ₂ -Precovered ZnO(101 0). <i>Journal of Physical Chemistry C</i> , 2018 , 122, 8919-8924	3.8	13

72	Electronic and transport properties of graphene with grain boundaries. <i>RSC Advances</i> , 2016 , 6, 1090-1093.7	3.7	13
71	Steric Hindrance Effect in High-Temperature Reactions. <i>CCS Chemistry</i> , 2020 , 2, 460-467	7.2	13
70	A Kinetic Pathway toward High-Density Ordered N Doping of Epitaxial Graphene on Cu(111) Using CNCl Precursors. <i>Journal of the American Chemical Society</i> , 2017 , 139, 7196-7202	16.4	12
69	Thickness-Dependent Adsorption of Melamine on Cu/Au(111) Films. <i>Journal of Physical Chemistry C</i> , 2017 , 121, 7977-7984	3.8	11
68	Reversible Tuning of Interfacial and Intramolecular Charge Transfer in Individual MnPc Molecules. <i>Nano Letters</i> , 2015 , 15, 8091-8	11.5	11
67	Single molecule tunneling spectroscopy investigation of reversibly switched dipolar vanadyl phthalocyanine on graphite. <i>Applied Physics Letters</i> , 2014 , 104, 113506	3.4	11
66	Are Azafullerene Encapsulated Single-Walled Carbon Nanotubes n-Type Semiconductors?. <i>Journal of Physical Chemistry C</i> , 2011 , 115, 12760-12762	3.8	11
65	Electron Transport in Butane Molecular Wires with Different Anchoring Groups Containing N, S, and P: A First Principles Study. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 21911-21914	3.8	11
64	First-principles simulation of amide and aromatic side chain ultraviolet spectroscopy of a cyclic dipeptide. <i>Journal of Physical Chemistry A</i> , 2007 , 111, 11579-83	2.8	11
63	Linear scaling calculation of maximally localized Wannier functions with atomic basis set. <i>Journal of Chemical Physics</i> , 2006 , 124, 234108	3.9	11
62	Electron-phonon interaction in a Ca ₂ N monolayer: Intrinsic mobility of electrene. <i>Physical Review B</i> , 2018 , 98,	3.3	11
61	Efficient Direct Band Gap Photovoltaic Material Predicted Via Doping Double Perovskites Cs ₂ AgBiX ₆ (X = Cl, Br). <i>Journal of Physical Chemistry C</i> , 2021 , 125, 10868-10875	3.8	10
60	First-Principles Study of Molecular Clusters Formed by Nitric Acid and Ammonia. <i>Journal of Physical Chemistry A</i> , 2017 , 121, 661-668	2.8	9
59	Ion Conductivity Enhancement in Anti-Spinel Li ₃ OBr with Intrinsic Vacancies. <i>Advanced Theory and Simulations</i> , 2019 , 2, 1800138	3.5	9
58	Designing Kagome Lattice from Potassium Atoms on Phosphorus-Gold Surface Alloy. <i>Nano Letters</i> , 2020 , 20, 5583-5589	11.5	9
57	Theoretical Insights into the Thermodynamics and Kinetics of Graphene Growth on Copper Surfaces. <i>Journal of Physical Chemistry C</i> , 2020 , 124, 16233-16247	3.8	9
56	Sandwich-Like Nanosheets of a N-Doped Porous Carbon/Graphene Composite with Enhanced Electrochemical Properties for Lithium and Sodium Storage. <i>ChemElectroChem</i> , 2018 , 5, 694-700	4.3	9
55	Single-stranded DNA adsorption on chiral molecule coated Au surface: a molecular dynamics study. <i>Physical Chemistry Chemical Physics</i> , 2010 , 12, 4431-4	3.6	9

54	Low-Temperature Heterolytic Adsorption of H ₂ on ZnO(101 0) Surface. <i>Journal of Physical Chemistry C</i> , 2019 , 123, 13283-13287	3.8	8
53	Theoretical Insights into Li-Ion Transport in LiTa ₂ PO ₈ . <i>Journal of Physical Chemistry C</i> , 2019 , 123, 19282-19287	3.8	8
52	Orientation-sensitive nonlinear growth of graphene: An epitaxial growth mechanism determined by geometry. <i>Physical Review B</i> , 2013 , 88,	3.3	8
51	STM studies of single molecules: molecular orbital aspects. <i>Chemical Communications</i> , 2011 , 47, 2747-625.8	5.8	8
50	First-principles calculations of conductance within a plane wave basis set via non-orthogonal Wannier-type atomic orbitals. <i>Journal of Physics Condensed Matter</i> , 2006 , 18, 1347-1358	1.8	8
49	A first-principles study on quasi-1D alkali metal chains within zeolite channels. <i>Journal of Chemical Physics</i> , 2004 , 120, 9725-8	3.9	7
48	Boosting ionic conductivity in antiperovskite Li ₃ OCl via defect engineering: Interstitial versus vacancy. <i>Physical Review Materials</i> , 2019 , 3,	3.2	7
47	Spin-phonon coupling in NiO nanoparticle. <i>Journal of Applied Physics</i> , 2020 , 128, 133905	2.5	7
46	The Nanoparticle Size Effect in Graphene Cutting: A Back-Man Mechanism. <i>Angewandte Chemie</i> , 2016 , 128, 10072-10075	3.6	7
45	Scanning tunneling microscopy and density functional theory combined studies of rutile TiO ₂ (1 1 0) surface chemistry: Watch surface processes at the atomic scale. <i>International Journal of Quantum Chemistry</i> , 2013 , 113, 89-95	2.1	6
44	A computational infrared spectroscopic study of graphene oxide. <i>Journal of Chemical Physics</i> , 2013 , 139, 084704	3.9	6
43	Nondecaying long range effect of surface decoration on the charge state of NV center in diamond. <i>Journal of Chemical Physics</i> , 2013 , 138, 034702	3.9	6
42	A first-principles study of ZnO polar surface growth: adsorption of Zn(x)O(y) clusters. <i>Journal of Chemical Physics</i> , 2013 , 139, 124704	3.9	6
41	Density functional study on mechanism of CO oxidation with activated water on O/Au (111) surface. <i>Science Bulletin</i> , 2009 , 54, 1973-1977	10.6	6
40	Graphene Oxide: Theoretical Perspectives 2012 , 69-84		6
39	Protonation effects on electron transport through diblock molecular junctions: A theoretical study. <i>Science in China Series B: Chemistry</i> , 2008 , 51, 1159-1165		6
38	Intrinsic ultra-wide completely spin-polarized state realized in a new CrO monolayer. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 17038-17041	3.6	6
37	One-Dimensional Magnetic Order Stabilized in Edge-Reconstructed MoS Nanoribbon via Bias Voltage. <i>Journal of Physical Chemistry Letters</i> , 2020 , 11, 7531-7535	6.4	6

- 36 Molecular Mechanism and Solvation Effect of Supramolecular Catalysis in a Synthetic Cavitand Receptor with an Inwardly Directed Carboxylic Acid for Ring-Opening Cyclization of Epoxy Alcohols. *ACS Catalysis*, **2018**, 8, 11910-11925 13.1 6
- 35 Molecular mechanism of heterogeneous supramolecular catalysis of metal-free cucurbituril solid for epoxide alcoholysis. *Molecular Catalysis*, **2019**, 467, 1-8 3.3 5
- 34 Blue Phosphorus Growth on Different Noble Metal Surfaces: From a 2D Alloy Network to an Extended Monolayer. *Journal of Physical Chemistry C*, **2021**, 125, 675-679 3.8 5
- 33 Dissociative adsorption and linear organization of formic acid on ZnO(100) surface. *Journal of Catalysis*, **2020**, 390, 109-116 7.3 5
- 32 Half-filled intermediate bands in doped inorganic perovskites for solar cells. *Physical Chemistry Chemical Physics*, **2020**, 22, 23804-23809 3.6 5
- 31 Rh Doping in Pd Nanocubes Optimizes the Adsorption of 3-Nitrostyrene towards Selective Hydrogenation of Vinyl Group. *ChemCatChem*, **2019**, 11, 2793-2798 5.2 4
- 30 Molecular Mechanisms and Atmospheric Implications of Criegee Intermediate-Alcohol Chemistry in the Gas Phase and Aqueous Surface Environments. *Journal of Physical Chemistry A*, **2020**, 124, 8585-8593^{2.8} 4
- 29 The mobility and solvation structure of a hydroxyl radical in a water nanodroplet: a Born-Oppenheimer molecular dynamics study. *Physical Chemistry Chemical Physics*, **2021**, 23, 14628-14635^{3.6} 4
- 28 Toward Epitaxial Growth of Misorientation-Free Graphene on Cu(111) Foils.. *ACS Nano*, **2021**, 16.7 4
- 27 Self-similarity of single-channel transmission for electron transport in nanowires. *Journal of Chemical Physics*, **2006**, 124, 104703 3.9 3
- 26 Is Mayenite without Clathrated Oxygen an Inorganic Electride?. *Angewandte Chemie*, **2004**, 116, 6641-6646 4.6 3
- 25 Simulating Periodic Systems on a Quantum Computer Using Molecular Orbitals. *Journal of Chemical Theory and Computation*, **2020**, 16, 6904-6914 6.4 3
- 24 On-Surface Synthesis of Nitrogen-Substituted Gold-Phosphorus Porous Network. *Chemistry of Materials*, **2020**, 32, 8561-8566 9.6 3
- 23 An efficient adaptive variational quantum solver of the Schrödinger equation based on reduced density matrices. *Journal of Chemical Physics*, **2021**, 154, 244112 3.9 3
- 22 Predictive design of intrinsic half-metallicity in zigzag tungsten dichalcogenide nanoribbons. *Physical Review B*, **2019**, 100, 3.3 3
- 21 Atom by Atom Condensation of Sn Single Clusters within Gold-Phosphorus Metal-Inorganic Porous Networks. *Journal of Physical Chemistry Letters*, **2021**, 12, 745-751 6.4 3
- 20 Obtaining Intrinsically Occupied Free-Space Superatom States in an Encapsulated CaN Nanotube. *ACS Omega*, **2018**, 3, 11966-11971 3.9 3
- 19 High-Pressure Phase Favored by a Symmetry-Recognized Nanoconfinement Effect. *Journal of Physical Chemistry Letters*, **2012**, 3, 2154-8 6.4 2

18	Understanding High-Temperature Chemical Reactions on Metal Surfaces: A Case Study on Equilibrium Concentration and Diffusivity of C H on a Cu(111) Surface.. <i>Jacs Au</i> , 2022 , 2, 443-452		2
17	Reaction Mechanism and Strategy for Optimizing the Hydrogen Evolution Reaction on Single-Layer 1T' WSe and WTe Based on Grand Canonical Potential Kinetics. <i>ACS Applied Materials & Interfaces</i> , 2021 , 13, 55611-55620	9.5	2
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