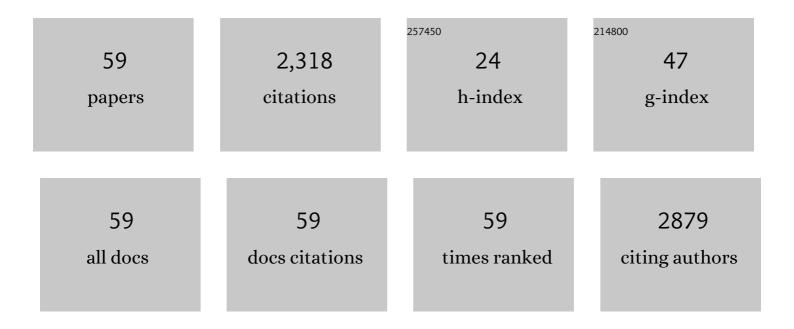
Xiao-Fei Li

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	First-Principles Observation of Bonded 2D B4C3 Bilayers. ACS Omega, 2021, 6, 13218-13224.	3.5	0
2	Rectification with controllable directions in sulfur-doped armchair graphene nanoribbon heterojunctions. Chemical Physics, 2021, 546, 111140.	1.9	1
3	A 4×4 Planar Dual-Polarization Retrodirective Array. , 2021, , .		1
4	Spin Polarization-Induced Facile Dioxygen Activation in Boron-Doped Graphitic Carbon Nitride. ACS Applied Materials & Interfaces, 2020, 12, 52741-52748.	8.0	15
5	Interface Magnetism in Topological Armchair/Cove-Edged Graphene Nanoribbons. Journal of Physical Chemistry C, 2020, 124, 15448-15453.	3.1	9
6	Folded C-Type SIW Butler Matrix. , 2019, , .		1
7	Improving the Oxygen Reduction Reaction Activity of FeN ₄ –Graphene via Tuning Electronic Characteristics. ACS Applied Energy Materials, 2019, 2, 6634-6641.	5.1	37
8	Ultrahigh conductivity of graphene nanoribbons doped with ordered nitrogen. Nanoscale Advances, 2019, 1, 4359-4364.	4.6	4
9	Rationally designed 2D/2D SiC/g-C ₃ N ₄ photocatalysts for hydrogen production. Catalysis Science and Technology, 2019, 9, 3896-3906.	4.1	35
10	Family-dependent magnetism in atomic boron adsorbed armchair graphene nanoribbons. Journal of Materials Chemistry C, 2019, 7, 6241-6245.	5.5	16
11	Perfect spin-filtering in p-aminophenol functionalized zigzag graphene nanoribbons: The role of sp3 hybridized nitrogen. Physics Letters, Section A: General, Atomic and Solid State Physics, 2019, 383, 2093-2096.	2.1	2
12	Industrial wastes applications for alkalinity regulation in bauxite residue: A comprehensive review. Journal of Central South University, 2019, 26, 268-288.	3.0	114
13	Alkalinity stabilization behavior of bauxite residue: Ca-driving regulation characteristics of gypsum. Journal of Central South University, 2019, 26, 383-392.	3.0	11
14	Migration and distribution of saline ions in bauxite residue during water leaching. Transactions of Nonferrous Metals Society of China, 2018, 28, 534-541.	4.2	37
15	Effect of ammonium chloride on leaching behavior of alkaline anion and sodium ion in bauxite residue. Transactions of Nonferrous Metals Society of China, 2018, 28, 2125-2134.	4.2	25
16	Leaching optimization and dissolution behavior of alkaline anions in bauxite residue. Transactions of Nonferrous Metals Society of China, 2018, 28, 1248-1255.	4.2	39
17	Cooperative Spin Transition of Monodispersed FeN ₃ Sites within Graphene Induced by CO Adsorption. Journal of the American Chemical Society, 2018, 140, 15149-15152.	13.7	108
18	Tunable Electronic and Magnetic Properties of Grapheneâ€Embedded Transition Metalâ€N ₄ Complexes: Insight From Firstâ€Principles Calculations. Chemistry - an Asian Journal, 2018, 13, 3239-3245.	3.3	18

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19	SiC/MoS2 layered heterostructures: Promising photocatalysts revealed by a first-principles study. Materials Chemistry and Physics, 2018, 216, 64-71.	4.0	63
20	Strong current-polarization and negative differential resistance in FeN3-embedded armchair graphene nanoribbons. Chinese Journal of Chemical Physics, 2018, 31, 756-760.	1.3	4
21	Superlight and Superflexible Threeâ€Dimensional Semiconductor Frameworks A(X≡Y) ₄ (A=Si,) Tj Chemistry - an Asian Journal, 2017, 12, 804-810.	ETQq1 1 3.3	0.784314 rg 3
22	Uniform and perfectly linear current–voltage characteristics of nitrogen-doped armchair graphene nanoribbons for nanowires. Physical Chemistry Chemical Physics, 2017, 19, 44-48.	2.8	13
23	Electric-field-induced widely tunable direct and indirect band gaps in hBN/MoS ₂ van der Waals heterostructures. Journal of Materials Chemistry C, 2017, 5, 4426-4434.	5.5	29
24	Strong current polarization and perfect negative differential resistance in few-FeN ₄ -embedded zigzag graphene nanoribbons. Physical Chemistry Chemical Physics, 2017, 19, 2674-2678.	2.8	19
25	Conversion of Dinitrogen to Ammonia by FeN ₃ -Embedded Graphene. Journal of the American Chemical Society, 2016, 138, 8706-8709.	13.7	562
26	Isolated pentagons induced enhancement of conductance in ultra-narrow armchair graphene nanoribbon junctions. Journal of Applied Physics, 2016, 120, 164303.	2.5	5
27	Insights into enhanced visible-light photocatalytic activity of C ₆₀ modified g-C ₃ N ₄ hybrids: the role of nitrogen. Physical Chemistry Chemical Physics, 2016, 18, 33094-33102.	2.8	31
28	Unraveling the formation mechanism of graphitic nitrogen-doping in thermally treated graphene with ammonia. Scientific Reports, 2016, 6, 23495.	3.3	111
29	Realizing Fano-like resonance in a one terminal closed T-shaped waveguide. European Physical Journal B, 2015, 88, 1.	1.5	14
30	Half-filled energy bands induced negative differential resistance in nitrogen-doped graphene. Nanoscale, 2015, 7, 4156-4162.	5.6	32
31	Resonance induced spin-selective transport behavior in carbon nanoribbon/nanotube/nanoribbon heterojunctions. Physics Letters, Section A: General, Atomic and Solid State Physics, 2015, 379, 1722-1725.	2.1	4
32	Semiconductor to metal transition by tuning the location of N2AA in armchair graphene nanoribbons. Journal of Applied Physics, 2014, 115, 053707.	2.5	12
33	Tunable bandwidth of the terahertz metamaterial absorber. Optics Communications, 2014, 325, 78-83.	2.1	33
34	A simple design of ultra-broadband and polarization insensitive terahertz metamaterial absorber. Applied Physics A: Materials Science and Processing, 2014, 115, 1187-1192.	2.3	67
35	Tuning electronic and magnetic properties of armchair zigzag hybrid graphene nanoribbons by the choice of supercell model of grain boundaries. Journal of Applied Physics, 2014, 115, 104303.	2.5	10
36	Frequency Continuous Tunable Terahertz Metamaterial Absorber. Journal of Lightwave Technology, 2014, 32, 1183-1189.	4.6	102

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#	Article	IF	CITATIONS
37	Realistic-contact-induced enhancement of rectifying in carbon-nanotube/graphene-nanoribbon junctions. Applied Physics Letters, 2014, 104, 103107.	3.3	12
38	Theoretical Investigation of Broadband and Wide-Angle Terahertz Metamaterial Absorber. IEEE Photonics Technology Letters, 2014, 26, 111-114.	2.5	176
39	A simple design of a broadband, polarization-insensitive, and low-conductivity alloy metamaterial absorber. Applied Physics Express, 2014, 7, 082601.	2.4	40
40	Conductivity of carbon-based molecular junctions from ab-initio methods. Frontiers of Physics, 2014, 9, 748-759.	5.0	5
41	A broadband, polarisation-insensitive and wide-angle coplanar terahertz metamaterial absorber. European Physical Journal B, 2014, 87, 1.	1.5	13
42	Tuning electron transport through a single molecular junction by bridge modification. Journal of Applied Physics, 2014, 116, .	2.5	5
43	Metamaterial-Based Low-Conductivity Alloy Perfect Absorber. Journal of Lightwave Technology, 2014, 32, 2293-2298.	4.6	49
44	A simple nested metamaterial structure with enhanced bandwidth performance. Optics Communications, 2013, 303, 13-14.	2.1	23
45	Big Bandgap in Highly Reduced Graphene Oxides. Journal of Physical Chemistry C, 2013, 117, 6049-6054.	3.1	52
46	First principles study on magnetic properties in ZnS doped with palladium. European Physical Journal B, 2013, 86, 1.	1.5	18
47	Strong current polarization and negative differential resistance in chiral graphene nanoribbons with reconstructed (2,1)-edges. Applied Physics Letters, 2012, 101, 073101.	3.3	15
48	Electronic transport through zigzag/armchair graphene nanoribbon heterojunctions. Journal of Physics Condensed Matter, 2012, 24, 095801.	1.8	18
49	A triangular shaped channel MIM waveguide filter. Journal of Modern Optics, 2012, 59, 1686-1689.	1.3	18
50	Ferromagnetic coupling in Mgâ€doped passivated AlN nanowires: A firstâ€principles study. Physica Status Solidi (B): Basic Research, 2012, 249, 185-189.	1.5	10
51	Design of Graphene-Nanoribbon Heterojunctions from First Principles. Journal of Physical Chemistry C, 2011, 115, 12616-12624.	3.1	49
52	Tuning the Electronic Transport Properties of Zigzag Graphene Nanoribbons via Hydrogenation Separators. Journal of Physical Chemistry C, 2011, 115, 24366-24372.	3.1	10
53	Firstâ€principles study of magnetic properties in Agâ€doped SnO ₂ . Physica Status Solidi (B): Basic Research, 2011, 248, 1961-1966.	1.5	26
54	Important Structural Factors Controlling the Conductance of DNA Pairs in Molecular Junctions. Journal of Physical Chemistry C, 2010, 114, 14240-14242.	3.1	11

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55	Effects of Interface Roughness on Electronic Transport Properties of Nanotubeâ^'Moleculeâ^'Nanotube Junctions. Journal of Physical Chemistry C, 2010, 114, 12335-12340.	3.1	11
56	Nanomechanically induced molecular conductance switch. Applied Physics Letters, 2009, 95, 232118.	3.3	14
57	Effect of intertube interaction on the transport properties of a carbon double-nanotube device. Journal of Applied Physics, 2007, 101, 064514.	2.5	25
58	Effect of length and size of heterojunction on the transport properties of carbon-nanotube devices. Applied Physics Letters, 2007, 91, 133511.	3.3	109
59	Coupling effect on the electronic transport through dimolecular junctions. Physics Letters, Section A: General, Atomic and Solid State Physics, 2007, 365, 489-494.	2.1	22