Xiaojie Liu

List of Publications by Year in descending order

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		687363	454955
35	904	13	30
papers	citations	h-index	g-index
2.5	2.5	0.5	1076
35	35	35	1276
all docs	docs citations	times ranked	citing authors

#	Article	IF	CITATIONS
1	Effect of the co-adsorption of small molecules from air on the properties of penta-graphene and their proton transfer calculation. Physical Chemistry Chemical Physics, 2022, , .	2.8	O
2	Comparative Study of Proton Exchange in Tri- and Hexatitanates: Correlations between Stability and Electronic Properties. Inorganic Chemistry, 2022, 61, 3918-3930.	4.0	6
3	Hybrid heterostructure of transition metal dichalcogenides as potential photocatalyst for hydrogen evolution. Applied Surface Science, 2022, 599, 154057.	6.1	7
4	g-C ₃ N ₄ /TiO ₂ -B{100} heterostructures used as promising photocatalysts for water splitting from a hybrid density functional study. Physical Chemistry Chemical Physics, 2022, 24, 17703-17715.	2.8	6
5	Strain-driven anisotropic Agl–S chains and semiconductor-to-metal transition in monoclinic Ag2S. Journal of Physics and Chemistry of Solids, 2022, 169, 110879.	4.0	0
6	Mechanism of Metal Intercalation under Graphene through Small Vacancy Defects. Journal of Physical Chemistry C, 2021, 125, 6954-6962.	3.1	13
7	Interactions and electronic properties of adatom/Gra/adatom sandwich complexes. Materials Chemistry and Physics, 2021, 272, 125013.	4.0	0
8	Structures and stabilities of glycine and water complexes. Chemical Physics, 2020, 528, 110528.	1.9	5
9	Amorphous Ag2-xCuxS quantum dots: "all-in-one―theranostic nanomedicines for near-infrared fluorescence/photoacoustics dual-modal-imaging-guided photothermal therapy. Chemical Engineering Journal, 2020, 399, 125777.	12.7	19
10	Effect of intrinsic vacancy defects on the electronic properties of monoclinic Ag2S. Materials Chemistry and Physics, 2020, 249, 122961.	4.0	14
11	Electronic and optical properties of gold-doped endohedral fullerenes. Journal of Materials Science, 2020, 55, 12980-12994.	3.7	2
12	Defect-mediated intercalation of dysprosium on buffer layer graphene supported by SiC(0001) substrate. Chemical Physics Letters, 2020, 742, 137162.	2.6	3
13	Theoretical investigation of dissociative and non-dissociative acetic-acid on TiO2-B surfaces. Applied Surface Science, 2019, 494, 850-858.	6.1	8
14	The effect of oxidation on the electronic properties of penta-graphene: first-principles calculation. RSC Advances, 2019, 9, 8253-8261.	3.6	14
15	Ringâ€Stacking Water Clusters: Morphology and Stabilities. ChemistryOpen, 2019, 8, 210-218.	1.9	15
16	Niobium-Doped TiO ₂ : Effect of an Interstitial Oxygen Atom on the Charge State of Niobium. Inorganic Chemistry, 2019, 58, 3090-3098.	4.0	14
17	Dy adsorption and penetration on defected graphene by first-principles calculations. Materials Research Express, 2018, 5, 025022.	1.6	8
18	First-principles study of electronic properties of Cu doped Ag ₂ S. Journal of Physics Condensed Matter, 2018, 30, 425502.	1.8	10

#	Article	IF	Citations
19	Interplay between quantum confinement and surface effects in thickness selective stability of thin Ag and Eu films. Journal of Physics Condensed Matter, 2017, 29, 185504.	1.8	4
20	Oscillatory electrostatic potential on graphene induced by group IV element decoration. Scientific Reports, 2017, 7, 13152.	3.3	4
21	Transition metal partially supported graphene: Magnetism and oscillatory electrostatic potentials. Journal of Applied Physics, 2017, 122, .	2.5	2
22	Growth mode and structures of magnetic Mn clusters on graphene. RSC Advances, 2016, 6, 64595-64604.	3.6	7
23	Metal intercalation-induced selective adatom mass transport on graphene. Nano Research, 2016, 9, 1434-1441.	10.4	7
24	Interplay between surface and surface resonance states on height selective stability of fcc Dy(111) film at nanoscale. Physical Chemistry Chemical Physics, 2016, 18, 31238-31243.	2.8	2
25	Magnetic Moment Enhancement for Mn ₇ Cluster on Graphene. Journal of Physical Chemistry C, 2014, 118, 19123-19128.	3.1	12
26	Electronic and spin transport properties of graphene nanoribbon mediated by metal adatoms: a study by the QUAMBO–NEGF approach. Journal of Physics Condensed Matter, 2013, 25, 105302.	1.8	7
27	Metals on Graphene: Interactions, Growth Morphology, and Thermal Stability. Crystals, 2013, 3, 79-111.	2.2	135
28	Directed assembly of Ru nanoclusters on Ru(0001)-supported graphene: STM studies and atomistic modeling. Physical Review B, 2012, 86, .	3.2	27
29	Growth morphology and thermal stability of metal islands on graphene. Physical Review B, 2012, 86, .	3.2	38
30	Metals on graphene: correlation between adatom adsorption behavior and growth morphology. Physical Chemistry Chemical Physics, 2012, 14, 9157.	2.8	145
31	Fe-Fe adatom interaction and growth morphology on graphene. Physical Review B, 2011, 84, .	3.2	23
32	Bonding and charge transfer by metal adatom adsorption on graphene. Physical Review B, 2011, 83, .	3.2	167
33	Energetic and fragmentation stability of water clusters (H2O)n, n=2–30. Chemical Physics Letters, 2011, 508, 270-275.	2.6	45
34	Metal Nanostructure Formation on Graphene: Weak versus Strong Bonding. Advanced Materials, 2011, 23, 2082-2087.	21.0	69
35	Adsorption and growth morphology of rare-earth metals on graphene studied by (i>ab initio (i) calculations and scanning tunneling microscopy. Physical Review B, 2010, 82, .	3.2	66