Jianfeng Jia

List of Publications by Year in descending order

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82	1,824	21	39
papers	citations	h-index	g-index
82	82	82	2241 citing authors
all docs	docs citations	times ranked	

#	Article	IF	CITATIONS
1	Carbon-coated BiVO4 prepared by molten salt method combined with ball milling for high-performance lithium-ion battery anode. Ionics, 2022, 28, 689-696.	2.4	3
2	The Investigation of Adsorption Behavior of Gas Molecules on FeN3-Doped Graphene. Journal of Sensors, 2022, 2022, 1-8.	1.1	2
3	Amorphous NiCo2O4 decorated Pd/C as electrocatalysts for boosting ethanol oxidation reaction in alkaline media. Electrochimica Acta, 2022, 411, 140048.	5.2	7
4	Facile in-situ electrochemical fabrication of highly efficient nickel hydroxide-iron hydroxide/graphene hybrid for oxygen evolution reaction. International Journal of Hydrogen Energy, 2022, 47, 12547-12558.	7.1	12
5	Adsorption Characteristics of Gas Molecules Adsorbed on Graphene Doped with Mn: A First Principle Study. Molecules, 2022, 27, 2315.	3.8	6
6	Immobilization of bismuth oxychloride on cellulose nanocrystal for sunlight-driven superior photosensitized degradation. International Journal of Biological Macromolecules, 2022, 206, 398-408.	7.5	7
7	Oxygen Vacancy-Mediated Selective C–N Coupling toward Electrocatalytic Urea Synthesis. Journal of the American Chemical Society, 2022, 144, 11530-11535.	13.7	142
8	A label-free electrochemical aptasensor based on the core–shell Cu-MOF@TpBD hybrid nanoarchitecture for the sensitive detection of PDGF-BB. Analyst, The, 2021, 146, 979-988.	3.5	28
9	Dative <i>versus</i> electron-sharing bonding in the isoelectronic argon compounds ArR ⁺ (R = CH ₃ , NH ₂ , OH, and F). New Journal of Chemistry, 2021, 45, 1363-1372.	2.8	1
10	Facile electrolytic synthesis of Pt and carbon quantum dots coloaded multiwall carbon nanotube as highly efficient electrocatalyst for hydrogen evolution and ethanol oxidation. Chemical Engineering Journal, 2021, 408, 127271.	12.7	27
11	Structural dependence of electrosynthesized cobalt phosphide/black phosphorus pre-catalyst for oxygen evolution in alkaline media. Nanoscale, 2021, 13, 7381-7388.	5.6	21
12	Highly sensitive and selective dopamine sensor uses three-dimensional cobalt phosphide nanowire array. Journal of Materials Science, 2021, 56, 6401-6410.	3.7	12
13	Hydroxyl-group-modified polymeric carbon nitride with the highly selective hydrogenation of nitrobenzene to <i>N</i> -phenylhydroxylamine under visible light. Green Chemistry, 2021, 23, 3612-3622.	9.0	22
14	Electrosynthesized CuO _x /graphene by a four-electrode electrolysis system for the oxygen reduction reaction to hydrogen peroxide. Chemical Communications, 2021, 57, 4118-4121.	4.1	12
15	Sc/Ti decorated novel C24N24 cage: Promising hydrogen storage materials. International Journal of Hydrogen Energy, 2021, 46, 7390-7401.	7.1	23
16	<scp>Ni₂P</scp> Nanosheets: A High Catalytic Activity Platform for Electrochemical Detection of Acetaminophen. Chinese Journal of Chemistry, 2021, 39, 1849-1854.	4.9	18
17	Tailoring phenol photomineralization pathway over polymeric carbon nitride with cyano group multifunctional active sites. Applied Catalysis B: Environmental, 2021, 284, 119710.	20.2	21
18	Li4Mo5O17 micron particles as new high-performance anode materials for lithium-ion batteries. Materials Letters, 2021, 305, 130803.	2.6	7

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19	Pd Nanoparticles Coupled to NiMoO ₄ –C Nanorods for Enhanced Electrocatalytic Ethanol Oxidation. ACS Applied Materials & Samp; Interfaces, 2021, 13, 53777-53786.	8.0	14
20	The Adsorption Behavior of Gas Molecules on Co/N Co–Doped Graphene. Molecules, 2021, 26, 7700.	3.8	6
21	Facile preparation of polyelectrolyte-functionalized reduced graphene oxide for significantly improving the performance of microbial fuel cells. Journal of Power Sources, 2020, 450, 227628.	7.8	16
22	The optimal adsorption pathway of H2 molecules on Ti-Acetylene/ ethylene compounds: A DFT study. International Journal of Hydrogen Energy, 2020, 45, 2105-2118.	7.1	3
23	Fe3O4 nanospheres decorated reduced graphene oxide as anode to promote extracellular electron transfer efficiency and power density in microbial fuel cells. Electrochimica Acta, 2020, 362, 137126.	5.2	30
24	A comparative study of electrocatalytic oxidation of glucose on conductive Ni-MOF nanosheet arrays with different ligands. New Journal of Chemistry, 2020, 44, 17849-17853.	2.8	26
25	Multicenter electron-sharing Ïf-bonding in the AgFe(CO) ₄ ^{â^'} complex. Dalton Transactions, 2020, 49, 15256-15266.	3.3	5
26	Cooperative physisorption and chemisorption of hydrogen on vanadium-decorated benzene. RSC Advances, 2020, 10, 37770-37778.	3.6	11
27	Triply Carbonyl-Bridged Ni ₂ (CO) ₅ Featuring Triple Three-Center Two-Electron Ni—C–Ni Bonds Instead of Ni≡Ni Triple Bond. Inorganic Chemistry, 2020, 59, 15365-15374.	4.0	3
28	Investigation of hydrogen storage on Sc/Ti-decorated novel B24N24. International Journal of Hydrogen Energy, 2020, 45, 33740-33750.	7.1	9
29	Defect Engineering in Pd/NiCo ₂ O _{4–<i>x</i>} for Selective Hydrogenation of α,β-Unsaturated Carbonyl Compounds under Ambient Conditions. ACS Sustainable Chemistry and Engineering, 2020, 8, 7851-7859.	6.7	29
30	Solvothermal synthesis of oxygen-incorporated MoS2-x nanosheets with abundant undercoordinated Mo for efficient hydrogen evolution. International Journal of Hydrogen Energy, 2020, 45, 19133-19143.	7.1	24
31	Thermodynamics and Kinetics of Gas-Phase CO Oxidation on the Scandium Monoxide Carbonyl Complexes. Journal of Physical Chemistry A, 2020, 124, 924-931.	2.5	4
32	Electric field-assisted synthesis of Pt, carbon quantum dots-coloaded graphene hybrid for hydrogen evolution reaction. Journal of Power Sources, 2020, 451, 227770.	7.8	32
33	A Sensitive Electrochemical MUC1 Sensing Platform Based on Electroactive Cu-MOFs Decorated by AuPt Nanoparticles. Journal of the Electrochemical Society, 2020, 167, 087502.	2.9	8
34	Molecular and dissociated adsorption of hydrogen on TiC6H6. International Journal of Hydrogen Energy, 2019, 44, 25800-25808.	7.1	8
35	The structures, stabilities and electronic properties of PdnB (n = 1–10) clusters. Computational and Theoretical Chemistry, 2019, 1164, 112554.	2.5	4
36	Facile <i>in situ</i> synthesis of a carbon quantum dot/graphene heterostructure as an efficient metal-free electrocatalyst for overall water splitting. Chemical Communications, 2019, 55, 1635-1638.	4.1	70

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37	Linear complex HC C-TMH (TM=Sc–Ni): A simple and efficient adsorbent for hydrogen molecules. International Journal of Hydrogen Energy, 2019, 44, 18145-18152.	7.1	8
38	Defining the optimal morphology of Rhn nanoparticles for efficient hydrazine adsorption: a DFT-D3 study. Journal of Materials Science, 2019, 54, 9533-9542.	3.7	5
39	High pressure behavior of crystal [2,2′-bi(1,3,4-oxadiazole)]-5,5′-dinitramide: A DFT investigation. Journal of Molecular Graphics and Modelling, 2019, 90, 87-93.	2.4	3
40	Light-assisted <i>O</i> -methylation of phenol with dimethyl carbonate over a layered double oxide catalyst. Catalysis Science and Technology, 2019, 9, 1774-1778.	4.1	12
41	The effect of interstitial boron on the mechanisms of acetylene hydrogenation catalyzed by Pd6: A DFT study. Computational and Theoretical Chemistry, 2019, 1170, 112636.	2.5	4
42	Theoretical Insights on the High Pressure Behavior of Pentazolate Anion Complex [Co(H2O)4(N5)2]·4H2O. Scientific Reports, 2019, 9, 15648.	3.3	3
43	Adsorption of multiple H2 molecules on the complex TiC6H6: An unusual combination of chemisorption and physisorption. Energy, 2019, 171, 315-325.	8.8	22
44	Geometries, stabilities, and magnetic properties of Co2Bn (n = 1–10) clusters. Journal of Molecular Modeling, 2019, 25, 27.	1.8	6
45	Facile synthesis of PdO-doped Co3O4 nanoparticles as an efficient bifunctional oxygen electrocatalyst. Applied Catalysis B: Environmental, 2019, 243, 175-182.	20.2	88
46	Hydrogenated graphene as support of Pd nanoparticles with improved electrocatalytic activity for ethanol oxidation reaction in alkaline media. Electrochimica Acta, 2019, 297, 856-863.	5.2	19
47	Electrochemical cathode exfoliation of bulky black phosphorus into few-layer phosphorene nanosheets. Electrochemistry Communications, 2018, 89, 10-13.	4.7	81
48	O 2 activation and CO oxidation on n-p codoped h-BN single-atom catalysts. Computational and Theoretical Chemistry, 2018, 1127, 31-36.	2.5	14
49	O2 adsorbed on Ptn clusters: Structure and optical absorption. AIP Advances, 2018, 8, 035307.	1.3	4
50	An insight into the structures, stabilities and magnetic properties of Fe2Bn (nÂ=Â1–10) clusters. Materials Chemistry and Physics, 2018, 205, 1-8.	4.0	15
51	Shape-control of super-branched Pd–Cu alloys with enhanced electrocatalytic performance for ethylene glycol oxidation. Chemical Communications, 2018, 54, 13363-13366.	4.1	33
52	Porous carbon-coated Li2MoO4 as high-performance anode materials for lithium-ion batteries. Materials Letters, 2018, 233, 302-305.	2.6	17
53	Molten salt synthesis of hexagonal tungsten trioxide nanoparticles for lithium-ion battery anode. Materials Letters, 2018, 233, 199-202.	2.6	10
54	A novel K2Ti8O17 nanorod photocatalyst rich in surface OH groups for efficient hydrogen production by water splitting. International Journal of Hydrogen Energy, 2018, 43, 18115-18124.	7.1	11

Digomerization of Vanadum-acetylene systems and its effect on hydrogen storage. International Journal of Hydrogen Energy, 2017, 42, 14188-14198. 7.1 16	#	Article	IF	CITATIONS
activity for ethylene glycol and glycerol oxidation. International Journal of Hydrogen Energy, 2017, 42, 25959. 259599. 259599. 259599. 259599. 259599. 259599. 259599. 259599. 2595999. 2595999. 2595999. 2595999. 2595999. 2595999. 2595999. 2595999. 25959999. 25959999. 25959999. 25959999. 25959999. 259599999. 25959999999. 259599999. 2595999999. 25959999999. 25959999999999	55	Oligomerization of Vanadium-acetylene systems and its effect on hydrogen storage. International Journal of Hydrogen Energy, 2017, 42, 14188-14198.	7.1	16
Density functional theory study of the interaction of hydrogen with TMC2H2(TM=Sc-Ni). International Journal of Hydrogen Energy, 2017, 42, 29384-29393. 7.1 9 Aucsub>nc/sub> (n = 1â€"16) clusters on the ZrO ⟨sub>2⟨sub>⟨slub⟩⟨111⟩ surface: a DFT+U investigation. Physical Chemistry Chemical Physics, 2016, 18, 30491-30497. Design and selection of triazole-based compounds with high energetic properties and stabilities. Design and selection of triazole-based compounds with high energetic properties and stabilities. Design and selection of triazole-based compounds with high energetic properties and stabilities. Design and selection of triazole-based compounds with high energetic properties and stabilities. Lis 6 Density functional theory study of Mo-doped M@ (BN)48 (MÅ=ÅSc, Ti, V, Cr, Mn, Fe, Co, Ni, and Cu) clusters. Journal of Molecular Structure, 2016, 1108, 92.95. Mechanistic investigation of palladium-catalyzed amidation of aryl halides. Journal of Molecular Modelling, 2016, 22, 53. Electronic structure, stability and magnetic properties of small M1â€"4(M = Fe, Co, Ni) clusters encapsulated inside a (BN)48 cage. Chemical Physics Letters, 2015, 622, 57-62. Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section At General, Atomic and Solid State Physics, 2015, 459, 131-136. Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section At General, Atomic and Solid State Physics, 2015, 459, 131-136. Structural and electronic properties of V2Bn (n=1â€"10) clusters. Chemical Physics, 2015, 459, 131-136. Computational investigation of hydrogen storage on scandiumā€"acetylene system. International Journal of Hydrogen Energy, 2015, 40, 420-428. Effect of COOH group on the performance of rhenium (I) tricarbonyl complexes with tetrathiafulvalene fused phenanthroline ligands as dyes in DSSC: DFT/TD-DFT t	56	activity for ethylene glycol and glycerol oxidation. International Journal of Hydrogen Energy, 2017,	7.1	33
International Journal of Hydrogen Energy, 2017, 42, 29384-29393. Ausubhn (subh (n = 1â€"16) clusters on the ZrO (subh 2c/subh (111) surface: a DFT+U investigation. Physical Chemistry Chemical Physics, 2016, 18, 30491-30497. Design and selection of triazole-based compounds with high energetic properties and stabilities. Journal of Chemical Sciences, 2016, 128, 1223-1236. Density functional theory study of Mo-doped M@ (BN)48 (MÂ=ÂSc, Ti, V, Cr, Mn, Fe, Co, Ni, and Cu) clusters. Journal of Molecular Structure, 2016, 1108, 92-95. Mechanistic investigation of palladium-catalyzed amidation of aryl halides. Journal of Molecular Modeling, 2016, 22, 53. Electronic structure, stability and magnetic properties of small M1â€"4(M = Fe, Co, Ni) clusters encapsulated inside a (BN)48 cage. Chemical Physics Letters, 2015, 622, 57-62. Computational investigation of hydrogen adsorption/desorption on Zrâ€"ſ-2 â€"(C 2 H 2) and its ion. Chemical Physics, 2015, 457, 57-62. Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section A: General, Atomic and Solid State Physics, 2.1 7 Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section A: General, Atomic and Solid State Physics, 2.1 7 Structural and electronic properties of V2Bn (n=1â€"10) clusters. Chemical Physics, 2015, 459, 131-136. 1.9 13 Computational investigation of hydrogen storage on scandiumâ€"acetylene system. International Journal of Hydrogen Energy, 2015, 40, 420-428. Effect of COOH group on the performance of rhenium (I) tricarbonyl complexes with tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/ID-DFT theoretical investigations. Structural Chemistry, 2015, 26, 421-430.	57		1.4	5
Physical Chemistry Chemical Physics, 2016, 18, 30491-30497. Design and selection of triazole-based compounds with high energetic properties and stabilities. 1.5 6 Density functional theory study of Mo-doped M@(BN)48 (MÂ=ÂSc, Ti, V, Cr, Mn, Fe, Co, Ni, and Cu) dusters. Journal of Molecular Structure, 2016, 1108, 92-95. Mechanistic investigation of palladium-catalyzed amidation of aryl halides. Journal of Molecular Modeling, 2016, 22, 53. Electronic structure, stability and magnetic properties of small M1â€"4(M = Fe, Co, Ni) clusters encapsulated inside a (BN)48 cage. Chemical Physics Letters, 2015, 622, 57-62. Computational investigation of hydrogen adsorption/desorption on Zrâ€"f. 2 â€" (C 2 H 2) and its ion. Chemical Physics, 2015, 457, 57-62. Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section At General, Atomic and Solid State Physics, 2.1 7 Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section At General, Atomic and Solid State Physics, 2.1 7 Structural and electronic properties of V2Bn (n=1â€"10) clusters. Chemical Physics, 2015, 459, 131-136. Computational investigation of hydrogen storage on scandiumâ€"acetylene system. International Journal of Hydrogen Energy, 2015, 40, 420-428. Effect of COOH group on the performance of rhenium (I) tricarbonyl complexes with tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/ITD-DFT theoretical investigations. Structural Chemistry, 2015, 26, 421-430.	58		7.1	9
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clusters. Journal of Molecular Structure, 2016, 1108, 92-95. Mechanistic investigation of palladium-catalyzed amidation of aryl halides. Journal of Molecular Modeling, 2016, 22, 53. Electronic structure, stability and magnetic properties of small M1â€"4(M = Fe, Co, Ni) clusters encapsulated inside a (BN)48 cage. Chemical Physics Letters, 2015, 622, 57-62. Computational investigation of hydrogen adsorption/desorption on Zrâ€"Î- 2 â€"(C 2 H 2) and its ion. Chemical Physics, 2015, 457, 57-62. Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section At General, Atomic and Solid State Physics, 2015, 379, 1715-1721. Structural and electronic properties of V2Bn (n=1â€"10) clusters. Chemical Physics, 2015, 459, 131-136. 1.9 13 Computational investigation of hydrogen storage on scandiumâ€"acetylene system. International Journal of Hydrogen Energy, 2015, 40, 420-428. Effect of COOH group on the performance of rhenium (I) tricarbonyl complexes with tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/ID-DFT theoretical investigations. Structural Chemistry, 2015, 26, 421-430.	60	Design and selection of triazole-based compounds with high energetic properties and stabilities. Journal of Chemical Sciences, 2016, 128, 1223-1236.	1.5	6
Modeling, 2016, 22, 53. Electronic structure, stability and magnetic properties of small M1â€"4(M = Fe, Co, Ni) clusters encapsulated inside a (BN)48 cage. Chemical Physics Letters, 2015, 622, 57-62. Computational investigation of hydrogen adsorption/desorption on Zrâ€". 2â€" (C 2 H 2) and its ion. Chemical Physics, 2015, 457, 57-62. Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section A: General, Atomic and Solid State Physics, 2015, 379, 1715-1721. Structural and electronic properties of V2Bn (n=1â€"10) clusters. Chemical Physics, 2015, 459, 131-136. 1.9 13 Computational investigation of hydrogen storage on scandiumâ€"acetylene system. International Journal of Hydrogen Energy, 2015, 40, 420-428. Effect of COOH group on the performance of rhenium (I) tricarbonyl complexes with tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/ID-DFT theoretical investigations. Structural Chemistry, 2015, 26, 421-430.	61	Density functional theory study of Mo-doped M@(BN)48 (MÂ=ÂSc, Ti, V, Cr, Mn, Fe, Co, Ni, and Cu) clusters. Journal of Molecular Structure, 2016, 1108, 92-95.	3.6	2
computational investigation of hydrogen adsorption/desorption on Zrâ€"η 2 â€"(C 2 H 2) and its ion. Chemical Physics, 2015, 457, 57-62. Electronic structure, stability and magnetic properties of small M1â€"2Cr (M = Fe, Co, and Ni) alloy encapsulated inside a (BN)48 cage. Physics Letters, Section A: General, Atomic and Solid State Physics, 2015, 379, 1715-1721. Structural and electronic properties of V2Bn (n=1â€"10) clusters. Chemical Physics, 2015, 459, 131-136. Computational investigation of hydrogen storage on scandiumâ€"acetylene system. International Journal of Hydrogen Energy, 2015, 40, 420-428. Effect of COOH group on the performance of rhenium (I) tricarbonyl complexes with tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/TD-DFT theoretical investigations. Structural Chemistry, 2015, 26, 421-430.	62		1.8	6
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Computational investigation of hydrogen storage on scandium–acetylene system. International Journal of Hydrogen Energy, 2015, 40, 420-428. Effect of COOH group on the performance of rhenium (I) tricarbonyl complexes with tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/TD-DFT theoretical investigations. Structural Chemistry, 2015, 26, 421-430.	65	encapsulated inside a (BN)48 cage. Physics Letters, Section A: General, Atomic and Solid State Physics,	2.1	7
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68 tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/TD-DFT theoretical 2.0 9 investigations. Structural Chemistry, 2015, 26, 421-430.	67		7.1	29
Viable Dhotocatalysts under Solarâf Spectrum Irradiation: Napplasmonic Motal Napoparticles	68	tetrathiafulvalene-fused phenanthroline ligands as dyes in DSSC: DFT/TD-DFT theoretical	2.0	9
Angewandte Chemie - International Edition, 2014, 53, 2935-2940.	69	Viable Photocatalysts under Solarâ€Spectrum Irradiation: Nonplasmonic Metal Nanoparticles. Angewandte Chemie - International Edition, 2014, 53, 2935-2940.	13.8	234
Efficient photocatalytic Suzuki cross-coupling reactions on Au–Pd alloy nanoparticles under visible light irradiation. Green Chemistry, 2014, 16, 4272.	70	Efficient photocatalytic Suzuki cross-coupling reactions on Au–Pd alloy nanoparticles under visible light irradiation. Green Chemistry, 2014, 16, 4272.	9.0	213
Density functional theory investigation on the structure and stability of Sc2B (n= 1–10) clusters. 2.5 20 Computational and Theoretical Chemistry, 2014, 1027, 128-134.	71	Density functional theory investigation on the structure and stability of Sc2B (n= $1\hat{a}$ e"10) clusters. Computational and Theoretical Chemistry, 2014, 1027, 128-134.	2.5	20

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73	Structures and stabilities of ScBn (n = 1–12) clusters: an ab initio investigation. Journal of Molecular Modeling, 2013, 19, 3255-3261.	1.8	26
74	Tiâ€"η2-(C2H2) and HC Câ€"TiH as high capacity hydrogen storage media. International Journal of Hydrogen Energy, 2013, 38, 16185-16192.	7.1	37
75	Elucidation of the Forces Governing the Stereochemistry of Biphenyl. European Journal of Organic Chemistry, 2013, 2013, 611-616.	2.4	26
76	Theoretical studies of COOH group effect on the performance of rhenium (I) tricarbonyl complexes with bispyridine sulfur-rich core ligand as dyes in DSSC. Theoretical Chemistry Accounts, 2012, 131, 1.	1.4	9
77	Structure and Stability of TiB <i>_n</i> (<i>n</i> i²i¼1—12) Clusters: An <i>abinitio</i> lnvestigation. Acta Chimica Sinica, 2012, 70, 1643.	1.4	9
78	Structure and Stability of Tube and Cage Ge ₆₀ H ₆₀ . Journal of Physical Chemistry A, 2010, 114, 12755-12758.	2.5	4
79	Tube and Cage C ₆₀ H ₆₀ : A Comparison with C ₆₀ F ₆₀ . Organic Letters, 2008, 10, 2573-2576.	4.6	12
80	Fused Five-Membered Rings Determine the Stability of C ₆₀ F ₆₀ . Journal of the American Chemical Society, 2008, 130, 3985-3988.	13.7	39
81	Structure and stability of boron nitride cages. Science Bulletin, 2003, 48, 1102-1107.	1.7	5
82	Photoelectron velocity-map imaging spectroscopy of nickel carbide: Examination of the low-lying electronic states. New Journal of Chemistry, 0, , .	2.8	1