

# Guochun Yang

## List of Publications by Year in Descending Order

**Source:** <https://exaly.com/author-pdf/8657032/guochun-yang-publications-by-year.pdf>

**Version:** 2024-04-20

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

203  
papers

3,969  
citations

34  
h-index

51  
g-index

218  
ext. papers

4,771  
ext. citations

4.9  
avg, IF

5.76  
L-index

#	Paper	IF	Citations
203	A superconducting boron allotrope featuring anticlinal pentapyramids. <i>Journal of Materials Chemistry C</i> , <b>2022</b> , 10, 672-679	7.1	1
202	Superconducting LaP2H2 with graphenelike phosphorus layers. <i>Physical Review B</i> , <b>2022</b> , 105,	3.3	2
201	Disproportionation of SO <sub>2</sub> at High Pressure and Temperature.. <i>Physical Review Letters</i> , <b>2022</b> , 128, 106001	7.4	0
200	Superconducting ScP4 with a novel phosphorus framework. <i>Applied Physics A: Materials Science and Processing</i> , <b>2022</b> , 128, 1	2.6	1
199	Theoretical considerations of superconducting HfBH2 and HfB2H under high pressure. <i>Journal of Applied Physics</i> , <b>2021</b> , 130, 153904	2.5	1
198	Janus MoP Monolayer as an Electrocatalyst for Hydrogen Evolution. <i>ACS Applied Materials &amp; Interfaces</i> , <b>2021</b> , 13, 57422-57429	9.5	0
197	Highly Photoluminescent Monolayer MoS2 and WS2 Achieved via Superacid Assisted Vacancy Reparation and Doping Strategy. <i>Laser and Photonics Reviews</i> , <b>2021</b> , 15, 2100104	8.3	0
196	Strain-Induced Magnetism in MSi2N4 (M = V, Cr): A First-Principles Study. <i>Annalen Der Physik</i> , <b>2021</b> , 533, 2100273	2.6	3
195	Pressure-induced hydride superconductors above 200 K. <i>Matter and Radiation at Extremes</i> , <b>2021</b> , 6, 068207	4.7	2
194	Anisotropic Janus SiP Monolayer as a Photocatalyst for Water Splitting. <i>Journal of Physical Chemistry Letters</i> , <b>2021</b> , 12, 2464-2470	6.4	16
193	Pressure-induced yttrium oxides with unconventional stoichiometries and novel properties. <i>Physical Review Materials</i> , <b>2021</b> , 5,	3.2	2
192	Ba with Unusual Oxidation States in Ba Chalcogenides under Pressure. <i>Journal of Physical Chemistry Letters</i> , <b>2021</b> , 12, 4203-4210	6.4	4
191	Tailoring p-type conductivity of aluminum nitride via transition metal and fluorine doping. <i>Journal of Alloys and Compounds</i> , <b>2021</b> , 862, 158017	5.7	0
190	Pressure-induced Na-Au compounds with novel structural units and unique charge transfer. <i>Physical Chemistry Chemical Physics</i> , <b>2021</b> , 23, 6455-6461	3.6	1
189	Hard and superconducting cubic boron phase via swarm-intelligence structural prediction driven by a machine-learning potential. <i>Physical Review B</i> , <b>2021</b> , 103,	3.3	7
188	IrN and IrN as potential high-energy-density materials. <i>Journal of Chemical Physics</i> , <b>2021</b> , 154, 054706	3.9	3
187	Anisotropic and High-Mobility CS Monolayer as a Photocatalyst for Water Splitting. <i>Journal of Physical Chemistry Letters</i> , <b>2021</b> , 12, 8320-8327	6.4	8

186	Unraveling the synergetic mechanism of physisorption and chemisorption in laser-irradiated monolayer WS <sub>2</sub> . <i>Nano Research</i> , <b>2021</b> , 14, 4274	10	1
185	Wide Band Gap PS Monolayer with Anisotropic and Ultrahigh Carrier Mobility. <i>Journal of Physical Chemistry Letters</i> , <b>2021</b> , 12, 8481-8488	6.4	4
184	SiCP Monolayer with a Direct Band Gap and High Carrier Mobility for Photocatalytic Water Splitting.. <i>Journal of Physical Chemistry Letters</i> , <b>2021</b> , 190-197	6.4	2
183	Au with sp Hybridization in LiAuP.. <i>Journal of Physical Chemistry Letters</i> , <b>2021</b> , 236-242	6.4	0
182	Superconducting boron allotropes. <i>Physical Review B</i> , <b>2020</b> , 101,	3.3	8
181	Unconventional stable stoichiometry of vanadium peroxide. <i>Physical Chemistry Chemical Physics</i> , <b>2020</b> , 22, 11460-11466	3.6	1
180	A controllable and reversible phase transformation between all-inorganic perovskites for white light emitting diodes. <i>Journal of Materials Chemistry C</i> , <b>2020</b> , 8, 8374-8379	7.1	27
179	Computational predictions of two-dimensional anode materials of metal-ion batteries. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , <b>2020</b> , 10, e1473	7.9	9
178	Exploring the Limits of Transition-Metal Fluorination at High Pressures. <i>Angewandte Chemie - International Edition</i> , <b>2020</b> , 59, 9155-9162	16.4	9
177	Exploring the Limits of Transition-Metal Fluorination at High Pressures. <i>Angewandte Chemie</i> , <b>2020</b> , 132, 9240-9247	3.6	
176	Achieving high conductivity p-type Ga <sub>2</sub> O <sub>3</sub> through Al-N and In-N co-doping. <i>Chemical Physics Letters</i> , <b>2020</b> , 746, 137308	2.5	7
175	The photophysical properties of cycloparaphenylene-based compounds with figure-eight configurations. <i>New Journal of Chemistry</i> , <b>2020</b> , 44, 12185-12193	3.6	3
174	Anisotropic PC6N Monolayer with Wide Band Gap and Ultrahigh Carrier Mobility. <i>Journal of Physical Chemistry C</i> , <b>2020</b> , 124, 4330-4337	3.8	7
173	Exploring the origin of electrochemical performance of Cr-doped LiNiMnO. <i>Physical Chemistry Chemical Physics</i> , <b>2020</b> , 22, 3831-3838	3.6	8
172	Prediction of new thermodynamically stable ZnNO at high pressure. <i>Physical Chemistry Chemical Physics</i> , <b>2020</b> , 22, 10941-10948	3.6	1
171	A Semiconducting Cationic Square-Grid Network with Fe <sup>III</sup> Centers Displaying Unusual Dynamic Behavior. <i>European Journal of Inorganic Chemistry</i> , <b>2020</b> , 2020, 1255-1259	2.3	1
170	Recent Advances and Applications of Inorganic Electrides. <i>Journal of Physical Chemistry Letters</i> , <b>2020</b> , 11, 3841-3852	6.4	16
169	Structural and electronic properties of tungsten oxides under high pressures. <i>Journal of Physics Condensed Matter</i> , <b>2020</b> , 32, 085403	1.8	2

168	Theoretical study on group III elements and F co-doped ZnO. <i>Journal of Alloys and Compounds</i> , <b>2020</b> , 819, 153012	5.7	5
167	Lithium lanthanum titanate perovskite as an anode for lithium ion batteries. <i>Nature Communications</i> , <b>2020</b> , 11, 3490	17.4	50
166	Achieving high hydrogen evolution reaction activity of a MoC monolayer. <i>Physical Chemistry Chemical Physics</i> , <b>2020</b> , 22, 26189-26199	3.6	4
165	Boron kagome-layer induced intrinsic superconductivity in a MnB <sub>3</sub> monolayer with a high critical temperature. <i>Physical Review B</i> , <b>2020</b> , 102,	3.3	6
164	O Adsorption Associated with Sulfur Vacancies on MoS Microspheres. <i>Inorganic Chemistry</i> , <b>2019</b> , 58, 21693-217625	3.2	5
163	Nickel Hydrides under High Pressure. <i>Journal of Physical Chemistry C</i> , <b>2019</b> , 123, 24243-24247	3.8	3
162	Unveiling the Role of Oxygen Vacancy in Li <sub>2</sub> MnO <sub>3</sub> upon Delithiation. <i>Journal of Physical Chemistry C</i> , <b>2019</b> , 123, 23403-23409	3.8	6
161	Metallic P3C monolayer as anode for sodium-ion batteries. <i>Journal of Materials Chemistry A</i> , <b>2019</b> , 7, 405-411	13	49
160	A hypervalent and cubically coordinated molecular phase of IF predicted at high pressure. <i>Chemical Science</i> , <b>2019</b> , 10, 2543-2550	9.4	20
159	High-Temperature Ferromagnetism in an FeP Monolayer with a Large Magnetic Anisotropy. <i>Journal of Physical Chemistry Letters</i> , <b>2019</b> , 10, 2733-2738	6.4	50
158	Predicted Pressure-Induced Superconducting Transition in Electride Li <sub>6</sub> P. <i>Physical Review Letters</i> , <b>2019</b> , 122, 097002	7.4	51
157	Nanotube-assembled pine-needle-like CuS as an effective energy booster for sodium-ion storage. <i>Journal of Materials Chemistry A</i> , <b>2019</b> , 7, 10619-10628	13	39
156	IrF Molecular Crystal under High Pressure. <i>Journal of the American Chemical Society</i> , <b>2019</b> , 141, 5409-5414	16.4	27
155	An Organic Emitter Displaying Dual Emissions and Efficient Delayed Fluorescence White OLEDs. <i>Advanced Optical Materials</i> , <b>2019</b> , 7, 1801667	8.1	16
154	Achieving high transparent Ga <sub>2</sub> O <sub>3</sub> through AlGa-InGa-VO. <i>Journal of Alloys and Compounds</i> , <b>2019</b> , 792, 405-410	5.7	3
153	Electronic structure and second-order nonlinear optical property of chiral peropyrenes. <i>Journal of Molecular Modeling</i> , <b>2019</b> , 25, 220	2	1
152	Pressure-induced new chemistry. <i>Chinese Physics B</i> , <b>2019</b> , 28, 106106	1.2	10
151	Phase diagrams and electronic properties of B-S and H-B-S systems under high pressure. <i>Physical Review B</i> , <b>2019</b> , 100,	3.3	7

150	FeP3 monolayer as a high-efficiency catalyst for hydrogen evolution reaction. <i>Journal of Materials Chemistry A</i> , <b>2019</b> , 7, 25665-25671	13	24
149	Isomer dependent molecular packing and carrier mobility of N-phenylcarbazolephenanthro[9,10-d]imidazole based materials as hosts for efficient electrophosphorescence devices. <i>Journal of Materials Chemistry C</i> , <b>2019</b> , 7, 13486-13492	7.1	14
148	Understanding photophysical properties of iridium complexes with N-(5-phenyl-1,3,4-oxadiazol-2-yl)-diphenylphosphinic amide as the ancillary ligand. <i>New Journal of Chemistry</i> , <b>2019</b> , 43, 16975-16980	3.6	
147	Two-Dimensional PC with Direct Band Gap and Anisotropic Carrier Mobility. <i>Journal of the American Chemical Society</i> , <b>2019</b> , 141, 1599-1605	16.4	76
146	Hexagonal BC2N with Remarkably High Hardness. <i>Journal of Physical Chemistry C</i> , <b>2018</b> , 122, 6801-6807	3.8	15
145	Complementary Resistive Switching Observed in Graphene Oxide-Based Memory Device. <i>IEEE Electron Device Letters</i> , <b>2018</b> , 39, 488-491	4.4	19
144	Photophysical Properties of Chiral Tetraphenylethylene Derivatives with the Fixed Propeller-Like Conformation. <i>Journal of Physical Chemistry C</i> , <b>2018</b> , 122, 5032-5039	3.8	11
143	TiC Monolayer with High Specific Capacity for Sodium-Ion Batteries. <i>Journal of the American Chemical Society</i> , <b>2018</b> , 140, 5962-5968	16.4	152
142	Accurate identification of layer number for few-layer WS and WSe via spectroscopic study. <i>Nanotechnology</i> , <b>2018</b> , 29, 124001	3.4	33
141	Photocatalytic Reduction of Graphene Oxide-TiO Nanocomposites for Improving Resistive-Switching Memory Behaviors. <i>Small</i> , <b>2018</b> , 14, e1801325	11	45
140	Gold with +4 and +6 Oxidation States in AuF and AuF. <i>Journal of the American Chemical Society</i> , <b>2018</b> , 140, 9545-9550	16.4	50
139	Unveiling the Photophysical Properties of Boron Heptaaryldipyromethene Derivatives. <i>ChemPhysChem</i> , <b>2018</b> , 19, 2751-2757	3.2	6
138	Theoretical Mechanistic Study of Nickel(0)/Lewis Acid Catalyzed Polyfluoroarylcyanation of Alkynes: Origin of Selectivity for C≡N Bond Activation. <i>Organometallics</i> , <b>2018</b> , 37, 2594-2601	3.8	10
137	Nonmetallic FeH6 under High Pressure. <i>Journal of Physical Chemistry C</i> , <b>2018</b> , 122, 12022-12028	3.8	21
136	Element substitution of kesterite CuZnSnS for efficient counter electrode of dye-sensitized solar cells. <i>Scientific Reports</i> , <b>2018</b> , 8, 8714	4.9	15
135	Mixed-valence Compounds: AuO2 and AuS. <i>ChemPhysChem</i> , <b>2018</b> , 19, 2971-2971	3.2	
134	Structural and Superconducting Properties of Tungsten Hydrides Under High Pressure. <i>Frontiers in Physics</i> , <b>2018</b> , 6,	3.9	5
133	Memory Devices: Photocatalytic Reduction of Graphene OxideTiO2 Nanocomposites for Improving Resistive-Switching Memory Behaviors (Small 29/2018). <i>Small</i> , <b>2018</b> , 14, 1870136	11	3

132	Mixed-valence Compounds: AuO and AuS. <i>ChemPhysChem</i> , <b>2018</b> , 19, 2989-2994	3.2	5
131	Barium in High Oxidation States in Pressure-Stabilized Barium Fluorides. <i>Journal of Physical Chemistry C</i> , <b>2018</b> , 122, 12448-12453	3.8	15
130	High-temperature driven inter-valley carrier transfer and significant fluorescence enhancement in multilayer WS. <i>Nanoscale Horizons</i> , <b>2018</b> , 3, 598-605	10.8	11
129	Chiral macrocyclic imine nickel(II) coordination complexes with diverse photophysical properties. <i>Dyes and Pigments</i> , <b>2017</b> , 140, 70-78	4.6	7
128	Understanding photophysical properties of chiral conjugated corrals for organic photovoltaics. <i>Journal of Materials Chemistry C</i> , <b>2017</b> , 5, 3495-3502	7.1	9
127	Enhancement of Exciton Emission from Multilayer MoS at High Temperatures: Intervalley Transfer versus Interlayer Decoupling. <i>Small</i> , <b>2017</b> , 13, 1700157	11	16
126	Highly Efficient Long-Wavelength Thermally Activated Delayed Fluorescence OLEDs Based on Dicyanopyrazino Phenanthrene Derivatives. <i>ACS Applied Materials &amp; Interfaces</i> , <b>2017</b> , 9, 9892-9901	9.5	128
125	Pressure-Induced Stable Beryllium Peroxide. <i>Inorganic Chemistry</i> , <b>2017</b> , 56, 5233-5238	5.1	15
124	Understanding the role of lithium sulfide clusters in lithium-sulfur batteries. <i>Journal of Materials Chemistry A</i> , <b>2017</b> , 5, 9293-9298	13	34
123	Tetragonal Structure BC <sub>4</sub> as a Superhard Material. <i>Journal of Physical Chemistry C</i> , <b>2017</b> , 121, 10119-10123	3.3	5
122	Adsorption Energy Optimization of Co <sub>3</sub> O <sub>4</sub> through Rapid Surface Sulfurization for Efficient Counter Electrode in Dye-Sensitized Solar Cells. <i>Journal of Physical Chemistry C</i> , <b>2017</b> , 121, 12524-12530	3.8	17
121	Phase diagram, stability and electronic properties of an FeB system under high pressure: a first principles study. <i>RSC Advances</i> , <b>2017</b> , 7, 15986-15991	3.7	18
120	Photophysical properties of chiral covalent organic cages. <i>Computational and Theoretical Chemistry</i> , <b>2017</b> , 1120, 1-7	2	2
119	Pressure-induced stable BeN <sub>4</sub> as a high-energy density material. <i>Journal of Power Sources</i> , <b>2017</b> , 365, 155-161	8.9	22
118	Pressure-Induced Stable Li <sub>5</sub> P for High-Performance Lithium-Ion Batteries. <i>Journal of Physical Chemistry C</i> , <b>2017</b> , 121, 21199-21205	3.8	27
117	Stable and metallic two-dimensional TaC <sub>2</sub> as an anode material for lithium-ion battery. <i>Journal of Materials Chemistry A</i> , <b>2017</b> , 5, 18698-18706	13	50
116	Unveiling photophysical properties of phenanthro[9,10-d]imidazole derivatives for organic light-emitting diodes. <i>Organic Electronics</i> , <b>2017</b> , 50, 220-227	3.5	5
115	Theoretical study on the photophysical properties of boron-fused double helicenes. <i>RSC Advances</i> , <b>2017</b> , 7, 56543-56549	3.7	2

114	Enhanced Electroluminescence from ZnO Quantum Dot Light-Emitting Diodes via Introducing Al <sub>2</sub> O <sub>3</sub> Retarding Layer and Ag@ZnO Hybrid Nanodots. <i>Advanced Optical Materials</i> , <b>2017</b> , 5, 1700493	8.1	16
113	Theoretical study on thiophene-based double helicenes with intrinsic large second-order nonlinear optical response. <i>RSC Advances</i> , <b>2016</b> , 6, 84705-84711	3.7	2
112	Tellurium Hydrides at High Pressures: High-Temperature Superconductors. <i>Physical Review Letters</i> , <b>2016</b> , 116, 057002	7.4	104
111	Structure and Electronic Properties of FeSH Compound under High Pressure. <i>Inorganic Chemistry</i> , <b>2016</b> , 55, 11434-11439	5.1	35
110	Silicon Framework-Based Lithium Silicides at High Pressures. <i>ACS Applied Materials &amp; Interfaces</i> , <b>2016</b> , 8, 16761-7	9.5	11
109	Pressure-induced reappearance of superconductivity in the oC24 phase of lithium. <i>Solid State Communications</i> , <b>2016</b> , 225, 7-11	1.6	2
108	Understanding the photophysical properties of chiral dinuclear Re(i) complexes and the role of Re(i) in their complexes. <i>Dalton Transactions</i> , <b>2016</b> , 45, 7285-93	4.3	12
107	Gold as a 6p-Element in Dense Lithium Aurides. <i>Journal of the American Chemical Society</i> , <b>2016</b> , 138, 4046-47	6.5	70
106	Dicyanomethylenated Acridone Based Crystals: Torsional Vibration Confinement Induced Emission with Supramolecular Structure Dependent and Stimuli Responsive Characteristics. <i>Journal of Physical Chemistry C</i> , <b>2016</b> , 120, 587-597	3.8	40
105	Structural transitions and electronic properties of sodium superoxide at high pressures. <i>RSC Advances</i> , <b>2016</b> , 6, 67910-67915	3.7	2
104	Ten-fold coordinated polymorph and metallization of TiO <sub>2</sub> under high pressure. <i>RSC Advances</i> , <b>2015</b> , 5, 54253-54257	3.7	11
103	Insight into the role of Li <sub>2</sub> S <sub>2</sub> in LiB batteries: a first-principles study. <i>Journal of Materials Chemistry A</i> , <b>2015</b> , 3, 8865-8869	13	57
102	Photophysical properties of quinoxaline-fused [7]carbohelicene derivatives. <i>RSC Advances</i> , <b>2015</b> , 5, 72907-72915	3.7	15
101	Pressure-induced structural changes and elemental dissociation of cadmium and mercury chalcogenides. <i>RSC Advances</i> , <b>2015</b> , 5, 104426-104432	3.7	4
100	Prediction of the XeHe binary phase diagram at high pressures. <i>Chemical Physics Letters</i> , <b>2015</b> , 640, 115-118	2.5	7
99	Phase Diagram and High-Temperature Superconductivity of Compressed Selenium Hydrides. <i>Scientific Reports</i> , <b>2015</b> , 5, 15433	4.9	56
98	High performance full color OLEDs based on a class of molecules with dual carrier transport channels and small singlet-triplet splitting. <i>Chemical Communications</i> , <b>2015</b> , 51, 10632-5	5.8	79
97	Structural and electronic properties of alkali metal peroxides at high pressures. <i>RSC Advances</i> , <b>2015</b> , 5, 104337-104342	3.7	10



96	Synthesis, crystal structures and DNA-binding properties of Cd(II), Cu(II) and Ni(II) complexes with 2-(2-pyridyl)benzothiazole. <i>Synthetic Metals</i> , <b>2015</b> , 200, 1-6	3.6	11
95	Nonplanar donor-acceptor chiral molecules with large second-order optical nonlinearities: 1,1,4,4-tetracyanobuta-1,3-diene derivatives. <i>Journal of Physical Chemistry A</i> , <b>2014</b> , 118, 1094-102	2.8	11
94	Electron transport via phenyl-perfluorophenyl interaction in crystals of fluorine-substituted dibenzalacetones. <i>RSC Advances</i> , <b>2014</b> , 4, 50188-50194	3.7	5
93	Organic polymorphs: one-compound-based crystals with molecular-conformation- and packing-dependent luminescent properties. <i>Advanced Materials</i> , <b>2014</b> , 26, 6168-73	2.4	224
92	The influence of the diphenylphosphoryl moiety on the phosphorescent properties of heteroleptic iridium(III) complexes and the OLED performance: a theoretical study. <i>Journal of Materials Chemistry C</i> , <b>2014</b> , 2, 2859	7.1	42
91	A Stable, Magnetic, and Metallic Li <sub>3</sub> O <sub>4</sub> Compound as a Discharge Product in a Li-Air Battery. <i>Journal of Physical Chemistry Letters</i> , <b>2014</b> , 5, 2516-21	6.4	48
90	The relationship between intermolecular interactions and charge transport properties of trifluoromethylated polycyclic aromatic hydrocarbons. <i>Organic Electronics</i> , <b>2014</b> , 15, 1896-1905	3.5	20
89	The effect of intermolecular interactions on the charge transport properties of thiazole/thiophene-based oligomers with trifluoromethylphenyl. <i>Journal of Molecular Graphics and Modelling</i> , <b>2014</b> , 51, 79-85	2.8	4
88	Theoretical study on photophysical properties of Pt(II) triarylborons with a 2,2-bpy core derivatives. <i>Journal of Molecular Graphics and Modelling</i> , <b>2013</b> , 44, 311-7	2.8	
87	Chirality recognition of the protonated serine dimer and octamer by infrared multiphoton dissociation spectroscopy. <i>Physical Chemistry Chemical Physics</i> , <b>2013</b> , 15, 1873-86	3.6	26
86	Theoretical study on the photophysical properties of chiral mononuclear and dinuclear zinc complexes. <i>RSC Advances</i> , <b>2013</b> , 3, 2241-2247	3.7	10
85	Chiroptical, linear, and second-order nonlinear optical properties of tetrathiafulvalenylallene: a multifunctional molecular material. <i>Journal of Materials Chemistry C</i> , <b>2013</b> , 1, 1399	7.1	20
84	Photophysical properties of azaboradibenzo[6]helicene derivatives. <i>Journal of Materials Chemistry C</i> , <b>2013</b> , 1, 2354	7.1	22
83	Theoretical insights into [PMo <sub>12</sub> O <sub>40</sub> ]( <sup>3-</sup> ) grafted on single-walled carbon nanotubes. <i>Physical Chemistry Chemical Physics</i> , <b>2013</b> , 15, 9177-85	3.6	23
82	Theoretical study on the rectifying performance of organoimido derivatives of hexamolybdates. <i>ChemPhysChem</i> , <b>2013</b> , 14, 610-7	3.2	15
81	Fluorescent photoswitching of a naphthopyran-benzimidazole dyad with high-degree fluorescent modulation within poly(methyl methacrylate) matrices. <i>Optical Materials</i> , <b>2013</b> , 35, 1504-1512	3.3	16
80	DFT characterization on the mechanism of water splitting catalyzed by single-Ru-substituted polyoxometalates. <i>Dalton Transactions</i> , <b>2013</b> , 42, 10617-25	4.3	25
79	Theoretical investigations on electronic spectra and the redox-switchable second-order nonlinear optical responses of rhodium(I)-9,10-phenanthrenediimine complexes. <i>Journal of Molecular Graphics and Modelling</i> , <b>2012</b> , 33, 19-25	2.8	7



78	TDDFT studies on the structures and ECD spectra of chiral bisarylimidos bearing different lengths of o-alkoxy chain-substituted polyoxomolybdates. <i>Journal of Molecular Graphics and Modelling</i> , <b>2012</b> , 35, 49-56	2.8	11
77	Density Functional Theory Investigation on the Second-Order Nonlinear Optical Properties of Chlorobenzyl-o-Carborane Derivatives. <i>Chinese Journal of Chemistry</i> , <b>2012</b> , 30, 2349-2355	4.9	8
76	Theoretical study on the charge transport property of Pt(CN(t)Bu) <sub>2</sub> (CN) <sub>2</sub> nanowires induced by Pt-Pt interactions. <i>Dalton Transactions</i> , <b>2012</b> , 41, 7272-7	4.3	11
75	TDDFT studies on chiral organophosphonate substituted divacant Keggin-type polyoxotungstate: diplex multistep-redox-triggered chiroptical and NLO switch. <i>Dalton Transactions</i> , <b>2012</b> , 41, 10097-104	4.3	11
74	THEORETICAL STUDY ON THE SECOND-ORDER NONLINEAR OPTICAL PROPERTIES OF C,B-SUBSTITUTED CARBORANE CONJUGATED DERIVATIVES. <i>Journal of Theoretical and Computational Chemistry</i> , <b>2012</b> , 11, 1121-1133	1.8	8
73	THEORETICAL INVESTIGATION ON SECOND-ORDER NONLINEAR OPTICAL PROPERTIES AND REDOX-SWITCHING OF PHENYL NITRONYL-NITROXIDE RADICAL DERIVATIVES. <i>Journal of Theoretical and Computational Chemistry</i> , <b>2012</b> , 11, 1075-1088	1.8	5
72	Redox-switching second-order nonlinear optical responses of N <sup>N</sup> N ruthenium complexes. <i>Computational and Theoretical Chemistry</i> , <b>2012</b> , 979, 112-118	2	16
71	Theoretical investigation on two-dimensional molecule-based second-order nonlinear optical materials of the disubstituted o-carborane derivatives. <i>Computational and Theoretical Chemistry</i> , <b>2012</b> , 992, 142-149	2	9
70	Chiroptical, linear, and second-order nonlinear optical properties of binaphthol derivatives. <i>Organic and Biomolecular Chemistry</i> , <b>2012</b> , 10, 8418-25	3.9	27
69	Computational study of chiral molecules with high intrinsic hyperpolarizabilities. <i>Molecular Physics</i> , <b>2012</b> , 110, 333-341	1.7	6
68	Theoretical study on photophysical properties of novel bis(BF <sub>2</sub> )-2,2'-bidipyrrins dyes: Effect of variation in monomer structure. <i>International Journal of Quantum Chemistry</i> , <b>2012</b> , 112, 440-452	2.1	9
67	Second-order nonlinear optical responses switching of N <sup>N</sup> N ruthenium carboxylate complexes with proton-electron transfer. <i>International Journal of Quantum Chemistry</i> , <b>2012</b> , 112, 779-788	2.1	6
66	Conformations of serine in aqueous solutions as revealed by vibrational circular dichroism. <i>ChemPhysChem</i> , <b>2012</b> , 13, 1272-81	3.2	30
65	Conformational distributions of N-acetyl-L-cysteine in aqueous solutions: a combined implicit and explicit solvation treatment of VA and VCD spectra. <i>ChemPhysChem</i> , <b>2012</b> , 13, 2310-21	3.2	38
64	Theoretical studies of the effect of electron-withdrawing dicyanovinyl group on the electronic and charge-transport properties of fluorene-thiophene oligomers. <i>Theoretical Chemistry Accounts</i> , <b>2012</b> , 131, 1	1.9	15
63	Effects of the substituting groups and proton abstraction on the nonlinear optical properties of heteroleptic bis-tridentate Ru(II) complexes. <i>Journal of Organometallic Chemistry</i> , <b>2011</b> , 696, 3384-3391	2.3	12
62	The effect of multiple weak interactions on the charge transport ability in polymorphs. <i>Synthetic Metals</i> , <b>2011</b> , 161, 1073-1078	3.6	5
61	Theoretical study on the second-order nonlinear optical properties and reorganization energy of silafluorenes and spirobisilafluorenes derivatives. <i>Theoretical Chemistry Accounts</i> , <b>2011</b> , 128, 249-256	1.9	8

60	Charge transport and electronic properties of N-heteroquinones: quadruple weak hydrogen bonds and strong $\pi$ -stacking interactions. <i>Theoretical Chemistry Accounts</i> , <b>2011</b> , 128, 257-264	1.9	19
59	Determination of the charge transport abilities of polymorphs [C <sub>6</sub> F <sub>5</sub> Cu] <sub>2</sub> (4,4'-bipy) with different interactions: a density functional theoretical investigation. <i>Theoretical Chemistry Accounts</i> , <b>2011</b> , 129, 45-51	1.9	3
58	Vibrational circular dichroism spectroscopy of chiral molecules. <i>Topics in Current Chemistry</i> , <b>2011</b> , 298, 189-236		57
57	Theoretical study on the chiroptical optical properties of chiral fullerene C <sub>60</sub> derivative. <i>Journal of Physical Chemistry A</i> , <b>2011</b> , 115, 13356-63	2.8	26
56	Theoretical discussions on electron transport properties of perylene bisimide derivatives with different molecular packings and intermolecular interactions. <i>Journal of Materials Chemistry</i> , <b>2011</b> , 21, 134-143		83
55	Quantum chemical study of structures, electronic spectrum, and nonlinear optical properties of polynuclear lithium compounds. <i>Computational and Theoretical Chemistry</i> , <b>2011</b> , 966, 14-19	2	17
54	Assignment of the absolute configuration of dinuclear zirconium complexes containing two homochiral N atoms using TDDFT calculations of ECD. <i>Chemical Physics Letters</i> , <b>2011</b> , 502, 266-270	2.5	11
53	A theoretical study of ambipolar organic transport material: 1,4-Bis(pentafluorobenzyl)[60]-fullerene. <i>Chemical Physics Letters</i> , <b>2011</b> , 506, 255-259	2.5	6
52	Computational study on second-order nonlinear optical (NLO) properties of a novel class of two-dimensional $\pi$ -and W-shaped sandwich metallocarborane-containing chromophores. <i>Journal of Organometallic Chemistry</i> , <b>2011</b> , 696, 2380-2387	2.3	28
51	Synthesis and characterizations of novel spindle-like terphenyl-type chromophores for non-linear optical materials. <i>Tetrahedron</i> , <b>2011</b> , 67, 4110-4117	2.4	12
50	Vibrational absorption, vibrational circular dichroism, and theoretical studies of methyl lactate self-aggregation and methyl lactate-methanol intermolecular interactions. <i>Journal of Chemical Physics</i> , <b>2010</b> , 132, 234513	3.9	29
49	Determination of the absolute configurations of bicyclo[3.1.0]hexane derivatives via electronic circular dichroism, optical rotation dispersion and vibrational circular dichroism spectroscopy and density functional theory calculations. <i>Organic and Biomolecular Chemistry</i> , <b>2010</b> , 8, 3777-83	3.9	20
48	Diastereomers of the pentacoordinate chiral phosphorus compounds in solution: absolute configurations and predominant conformations. <i>Dalton Transactions</i> , <b>2010</b> , 39, 6953-9	4.3	15
47	Theoretical study on dithieno[3,2-b:2',3'-d]phosphole derivatives: high-efficiency blue-emitting materials with ambipolar semiconductor behavior. <i>Theoretical Chemistry Accounts</i> , <b>2010</b> , 127, 419-427	1.9	11
46	Charge transport and luminescent properties of C <sub>6</sub> F <sub>5</sub> Cu(py) and their relationships with cuprophilic interactions: a density functional theory investigation. <i>Theoretical Chemistry Accounts</i> , <b>2010</b> , 127, 735-742	1.9	11
45	The modulation of electronic and optical properties of OXD-X through introduction of the electron-withdrawing groups: a DFT study. <i>Journal of Molecular Graphics and Modelling</i> , <b>2010</b> , 28, 427-34	2.8	3
44	Determination of the absolute configurations of synthetic daunorubicin analogues using vibrational circular dichroism spectroscopy and density functional theory. <i>Chirality</i> , <b>2010</b> , 22, 734-43	2.1	11
43	Determination of the absolute configuration of pentacoordinate chiral phosphorus compounds in solution by using vibrational circular dichroism spectroscopy and density functional theory. <i>Chemistry - A European Journal</i> , <b>2010</b> , 16, 2518-27	4.8	15

42	Theoretical Investigation on Electronic Transition of Tris(8-quinolinolate) Aluminum Grafted on Poly(p-phenylenevinylene) Units with the Localized-density-matrix Method. <i>Chinese Journal of Chemistry</i> , <b>2009</b> , 27, 1891-1896	4.9	1
41	Quantum chemical study of redox-switchable second-order optical nonlinearity in Keggin-type organoimido derivative $[PW_{11}O_{39}(ReNC_6H_5)]_n$ ( $n = 2, 3$ ). <i>Theoretical Chemistry Accounts</i> , <b>2009</b> , 122, 265-273	1.9	16
40	The influence of MM attraction on nonlinear optical properties of $(XMPH_3)_2$ ( $X = F, Cl$ ; and $M = Au, Ag$ and $Cu$ ): A theoretical study. <i>International Journal of Quantum Chemistry</i> , <b>2009</b> , 110, NA-NA	2.1	1
39	Theoretical study on the second-order nonlinear optical properties of nonconjugated D- $\pi$ A chromophores. <i>International Journal of Quantum Chemistry</i> , <b>2009</b> , 109, 1553-1559	2.1	12
38	Theoretical investigation of structures, electronic spectra and nonlinear optical properties of gold-pentacene ( $Au_2C_{22}H_{14}$ ) complexes. <i>Journal of Organometallic Chemistry</i> , <b>2009</b> , 694, 1266-1272	2.3	12
37	Theoretical study on the second-order nonlinear optical properties of gold (III) alkyl complexes. <i>Synthetic Metals</i> , <b>2009</b> , 159, 2406-2409	3.6	1
36	Triphenylamine-based pH chemosensor: Synthesis, crystal structure, photophysical properties and computational studies. <i>Synthetic Metals</i> , <b>2009</b> , 159, 2497-2501	3.6	13
35	Probing chiral solute-water hydrogen bonding networks by chirality transfer effects: a vibrational circular dichroism study of glycidol in water. <i>Journal of Chemical Physics</i> , <b>2009</b> , 130, 164506	3.9	53
34	Theoretical study on a novel series of fullerene-containing organometallics $Fe(\eta^5-C_5X_5)_2$ ( $X = CH, N, B$ ) and their large third-order nonlinear optical properties. <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 8086-92	2.8	21
33	The effects of self-aggregation on the vibrational circular dichroism and optical rotation measurements of glycidol. <i>Physical Chemistry Chemical Physics</i> , <b>2008</b> , 10, 6787-95	3.6	28
32	Effect of pi-conjugated length of bridging ligand on the optoelectronic properties of platinum(II) dimers. <i>Inorganic Chemistry</i> , <b>2008</b> , 47, 2347-55	5.1	48
31	Computational Study on Second-Order Nonlinear Response of a Series of Two-Dimensional Carbazole-Cored Chromophores. <i>Journal of Physical Chemistry C</i> , <b>2008</b> , 112, 7021-7028	3.8	29
30	Reversible redox-switchable second-order optical nonlinearity in polyoxometalate: a quantum chemical study of $[PW_{11}O_{39}(ReN)]_n$ ( $n = 3-7$ ). <i>Inorganic Chemistry</i> , <b>2008</b> , 47, 5245-52	5.1	43
29	Theoretical predication of third-order optical nonlinearities of $[Al_4MA_4]_n$ ( $n = 0, 1$ , $M = Ti, V$ and $Cr$ ) clusters. <i>Theoretical Chemistry Accounts</i> , <b>2008</b> , 119, 329-333	1.9	15
28	Density functional study of magnetic exchange of dinuclear manganese complexes with the heteropolymolyanion: $[MnII_2(Xn+Mo_9O_{33})_2]_2(n=0)$ ( $X = PV, AsV, SeVI$ ). <i>Science in China Series B: Chemistry</i> , <b>2008</b> , 51, 1174-1181		6
27	Theoretical study on the electronic structures and optical properties of oxadisilole-substituted acenes. <i>Chemical Physics Letters</i> , <b>2008</b> , 466, 37-43	2.5	5
26	Cooperative enhancement of two-photon absorption cross sections in three-branched oligofluorene with boron center. <i>Computational and Theoretical Chemistry</i> , <b>2008</b> , 855, 69-76		6
25	Quantum Chemical Study of Structures, Electronic Spectrum, and Nonlinear Optical Properties of GoldPentacene Complexes. <i>Organometallics</i> , <b>2007</b> , 26, 2082-2087	3.8	33

24	Theoretical study on photophysical properties of phenolpyridyl boron complexes. <i>Journal of Physical Chemistry A</i> , <b>2007</b> , 111, 2739-44	2.8	35
23	Theoretical Study on One- and Two-Photon Absorption Properties of PPV Derivative with Electron-Donor Phenylanthracene as Pendent Group. <i>Chinese Journal of Chemistry</i> , <b>2007</b> , 25, 465-471	4.9	2
22	DFT Studies on second-order nonlinear optical properties of mono (salicylaldiminato) Nickel(II) polyenyl Schiff base metal complexes. <i>Chemical Physics Letters</i> , <b>2007</b> , 443, 163-168	2.5	29
21	Theoretical studies on the electronic structure and spectral properties of versatile diarylethene-containing 1,10-phenanthroline ligands and their rhenium(I) complexes. <i>Journal of Organometallic Chemistry</i> , <b>2007</b> , 692, 5368-5374	2.3	24
20	THE INVESTIGATION ON ELECTRONIC STRUCTURE AND SECOND-ORDER NONLINEAR OPTICAL PROPERTIES OF III-V SEMICONDUCTOR CLUSTERS BY TIME-DEPENDENT DENSITY FUNCTIONAL THEORY. <i>Journal of Theoretical and Computational Chemistry</i> , <b>2007</b> , 06, 585-594	1.8	1
19	How Do the Different Defect Structures and Element Substitutions Affect the Nonlinear Optical Properties of Lacunary Keggin Polyoxometalates? A DFT Study. <i>European Journal of Inorganic Chemistry</i> , <b>2006</b> , 2006, 4179-4183	2.3	30
18	Prediction of second-order optical nonlinearity of trisorganotin-substituted beta-Keggin polyoxotungstate. <i>Inorganic Chemistry</i> , <b>2006</b> , 45, 7864-8	5.1	42
17	DFT Study on Second-order Nonlinear Optical Properties of a Series of Three-dimensional Cobalt(II) Metal Complexes. <i>Acta Physico-chimica Sinica</i> , <b>2006</b> , 22, 836-839		3
16	Theoretical study on photophysical and charge transport properties of 1,6-bis(2-hydroxyphenol)pyridylboron bis(4-n-butylphenyl)phenyleneamine compound. <i>Journal of Physical Chemistry A</i> , <b>2006</b> , 110, 8758-62	2.8	40
15	Theoretical study on the optoelectronic properties of electron-withdrawing substituted diethynylfluorenyl gold(I) complexes. <i>Journal of Physical Chemistry A</i> , <b>2006</b> , 110, 13036-44	2.8	21
14	Theoretical study on the electronic spectrum and the origin of remarkably large third-order nonlinear optical properties of organoimide derivatives of hexamolybdates. <i>Journal of Physical Chemistry B</i> , <b>2006</b> , 110, 23092-8	3.4	81
13	Theoretical studies on one- and two-photon absorption properties of rubrene and its derivatives. <i>Synthetic Metals</i> , <b>2006</b> , 156, 1218-1224	3.6	7
12	Theoretical study on characteristics of structure and vibrational frequency of spiro-linked complex Zn(PyIm) <sub>2</sub> (PyIm = 2(2?-pyridine)-imidazole) in excited state. <i>Chemical Physics Letters</i> , <b>2006</b> , 418, 302-306	2.5	5
11	TD-DFT investigation on the low-lying excited states of spiro-bithiophene. <i>Chemical Physics Letters</i> , <b>2006</b> , 429, 180-184	2.5	21
10	Quantum chemistry studies on the Ru-III interactions and the 31P NMR in [Ru(CO) <sub>3</sub> (Ph <sub>2</sub> Ppy) <sub>2</sub> (MCl <sub>2</sub> )] (M=Zn, Cd, Hg). <i>Journal of Organometallic Chemistry</i> , <b>2006</b> , 691, 1927-1933	2.3	8
9	Hyperpolarizabilities of para-nitroaniline and bis[4-(dimethylamino)phenyl] squaraine: The effects of functional/basis set based on TDDFT/BOS method. <i>Computational and Theoretical Chemistry</i> , <b>2006</b> , 773, 9-14		16
8	Theoretical study on the second-order nonlinear optical properties of asymmetric spiro-silabifluorene derivatives. <i>Journal of Physical Chemistry A</i> , <b>2006</b> , 110, 4817-21	2.8	55
7	Density functional theory study on the first hyperpolarizabilities of organoimido derivatives of hexamolybdates. <i>Journal of Physical Chemistry B</i> , <b>2005</b> , 109, 22332-6	3.4	57

6	Predication of second-order optical nonlinearity of [(Bu <sub>2</sub> t Im)AuX] (X=halogen) using time-dependent density-functional theory combined with sum-over-states method. <i>Journal of Chemical Physics</i> , <b>2005</b> , 123, 134302	3.9	22
5	DFT/FF study on electronic structure and second-order NLO property of dinuclear gold complex [Au(SeC <sub>2</sub> B <sub>10</sub> H <sub>11</sub> )(PPh <sub>3</sub> )] <sub>2</sub> . <i>Synthetic Metals</i> , <b>2005</b> , 152, 273-276	3.6	12
4	Calculations of two-photon absorption cross-sections of stilbene and bis (styryl) benzene derivatives by means of TDDFT-SOS method. <i>Computational and Theoretical Chemistry</i> , <b>2005</b> , 726, 61-65		18
3	Fabrication of Alkali Metal Boride: Honeycomb-Like Structured NaB <sub>4</sub> with High Hardness and Excellent Electrical Conductivity. <i>Advanced Functional Materials</i> , 2110872	15.6	2
2	BC6P Monolayer: Isostructural and Isoelectronic Analogues of Graphene with Desirable Properties for K-Ion Batteries. <i>Chemistry of Materials</i> ,	9.6	1
1	Superconducting ternary hydrides under high pressure. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> ,	7.9	4