

# Zhong-Jun Zhou

## List of Publications by Year in descending order

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30  
papers

686  
citations

687363

13  
h-index

677142

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docs citations

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times ranked

416  
citing authors

#	ARTICLE	IF	CITATIONS
1	Lacking of ESIPT Band of Aromatic ortho-Aminoaldehyde Derivatives Triggered by the N-H Vibration. <i>Physical Chemistry Chemical Physics</i> , 2022, , .	2.8	0
2	Generation of singlet oxygen catalyzed by the room-temperature-stable anthraquinone anion radical. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 14165-14171.	2.8	2
3	Dearomatization of Benzenoid Arenes Triggered by Triplet Excited State Intramolecular Proton Transfer. <i>Journal of Physical Chemistry A</i> , 2022, 126, 4424-4431.	2.5	1
4	Metalloradical complex $\text{Co}^{\text{I}}\text{Cp}^*\text{Ph}_3$ catalyzes the $\text{CO}_2$ reduction in gas phase: a theoretical study. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 1392-1400.	2.8	0
5	Electride-Sponsored Radical-Controlled $\text{CO}_2$ Reduction to Organic Acids: A Computational Design. <i>Chemistry - A European Journal</i> , 2020, 26, 6234-6239.	3.3	3
6	Copper(I) catalyzed $\text{CO}_2$ transformation: A density functional theory investigation. <i>Computational and Theoretical Chemistry</i> , 2020, 1175, 112745.	2.5	0
7	Theoretical study of a novel organic electride with large nonlinear optical responses. <i>International Journal of Quantum Chemistry</i> , 2020, 120, e26235.	2.0	3
8	Small Janus dimer as electric field manipulated molecular clam switch and electric information storage unit. <i>International Journal of Quantum Chemistry</i> , 2019, 119, e26005.	2.0	1
9	Theoretical study of the photochemical isomerization process of perfluoroaryltetrahedrane to perfluoroarylcyclobutadiene mediated by 9,10-dicyanoanthracene. <i>Theoretical Chemistry Accounts</i> , 2019, 138, 1.	1.4	0
10	Theoretical study of substituent effects on electride characteristics and the nonlinear optical properties of $\text{Li}@\text{calix}[4]\text{pyrrole}$ . <i>RSC Advances</i> , 2019, 9, 37919-37925.	3.6	5
11	Finding all-Nonmetal transition-metal-like superatom and its magnetic building block. <i>International Journal of Quantum Chemistry</i> , 2018, 118, e25570.	2.0	2
12	Efficient External Electric Field Manipulated Nonlinear Optical Switches of All-Metal Electride Molecules with Infrared Transparency: Nonbonding Electron Transfer Forms an Excess Electron Lone Pair. <i>Journal of Physical Chemistry C</i> , 2017, 121, 958-968.	3.1	53
13	Theoretical investigation of boron-doped lithium clusters, $\text{BLi}_n$ ( $n = 3-6$ ), activating $\text{CO}_2$ : an example of the carboxylation of $\text{C-H}$ bonds. <i>RSC Advances</i> , 2016, 6, 84042-84049.	3.6	8
14	An External Electric Field Manipulated Second-Order Nonlinear Optical Switch of an Electride Molecule: A Long-Range Electron Transfer Forms a Lone Excess Electron Pair and Quenches Singlet Diradical. <i>Journal of Physical Chemistry C</i> , 2016, 120, 13656-13666.	3.1	50
15	Excess-electron-induced $\text{C-C}$ bond formation in transformation of carbon dioxide. <i>RSC Advances</i> , 2016, 6, 851-858.	3.6	2
16	Nonlinear optical response and transparency of hexagonal boron nitride hybrid graphene nanoribbons. <i>Chemical Physics Letters</i> , 2014, 614, 57-61.	2.6	14
17	Theoretical investigation on nonlinear optical properties of carbon nanotubes with Stone-Wales defect rings. <i>Journal of Materials Chemistry C</i> , 2014, 2, 306-311.	5.5	42
18	The effects of external electric field: creating non-zero first hyperpolarizability for centrosymmetric benzene and strongly enhancing first hyperpolarizability for non-centrosymmetric edge-modified graphene ribbon $\text{H}_2\text{N}-(3,3)\text{ZGNR-NO}_2$ . <i>Journal of Molecular Modeling</i> , 2013, 19, 3983-3991.	1.8	14

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19	The structure and large nonlinear optical properties of a novel octupolar electride Li@36Adz. Computational and Theoretical Chemistry, 2013, 1023, 99-103.	2.5	29
20	Design of Lewis acid–base complex: enhancing the stability and first hyperpolarizability of large excess electron compound. Journal of Molecular Modeling, 2013, 19, 4805-4813.	1.8	7
21	New Acceptor–Bridge–Donor Strategy for Enhancing NLO Response with Long-Range Excess Electron Transfer from the NH <sub>2</sub> ...M/M <sub>3</sub> O Donor (M = Li, Na, K) to Inside the Electron Hole Cage C <sub>20</sub> F <sub>19</sub> Acceptor through the Unusual If Chain Bridge (CH <sub>2</sub> ) <sub>4</sub> . Journal of Physical Chemistry A, 2013, 117, 2835-2843.	2.5	78
22	The interaction between superalkalis (M <sub>3</sub> O, M = Na, K) and a C <sub>20</sub> F <sub>20</sub> cage forming superalkali electride salt molecules with excess electrons inside the C <sub>20</sub> F <sub>20</sub> cage: dramatic superalkali effect on the nonlinear optical property. Journal of Materials Chemistry, 2012, 22, 9652.	6.7	97
23	Theoretical study on the ion–molecule reaction of NH <sup>+</sup> with CH <sub>2</sub> O. International Journal of Quantum Chemistry, 2012, 112, 1654-1666.	2.0	0
24	What is the role of defects in single-walled carbon nanotubes for nonlinear optical property?. Journal of Materials Chemistry, 2011, 21, 8905.	6.7	16
25	Electric Field-Driven Acid–Base Chemistry: Proton Transfer from Acid (HCl) to Base (NH <sub>3</sub> /H <sub>2</sub> O). Journal of Physical Chemistry A, 2011, 115, 1418-1422.	2.5	43
26	Exceptionally Large Second-Order Nonlinear Optical Response in Donor–Graphene Nanoribbon–Acceptor Systems. Chemistry - A European Journal, 2011, 17, 2414-2419.	3.3	59
27	Modulated Nonlinear Optical Responses and Charge Transfer Transition in Endohedral Fullerene Dimers Na@C <sub>60</sub> C <sub>60</sub> @F with <i>n</i> -Fold Covalent Bond ( <i>n</i> = 1, 2, 5, and) Tj ETC, 1 0.7843 14 rg	1.7	11
28	Push–pull electron effects of the complexant in a Li atom doped molecule with electride character: a new strategy to enhance the first hyperpolarizability. Physical Chemistry Chemical Physics, 2010, 12, 10562.	2.8	66
29	Effect of substitution and cooperativity on the Cl–F blue shift in single-electron halogen-bonded H <sub>3</sub> C–F complex. Molecular Physics, 2010, 108, 2021-2026.	1.7	11
30	Bonding and correlation analysis of various Si <sub>2</sub> CO isomers on the potential energy surface. International Journal of Quantum Chemistry, 2009, 109, 907-919.	2.0	0