

Ju Xie

List of Publications by Year in descending order

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papers

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#	ARTICLE	IF	CITATIONS
1	Self-Assembled Polyoxometalate Nanodots as Bidirectional Cluster Catalysts for Polysulfide/Sulfide Redox Conversion in Lithium–Sulfur Batteries. <i>ACS Nano</i> , 2021, 15, 12222-12236.	7.3	77
2	Controllable Supramolecular Chiral Twisted Nanoribbons from Achiral Conjugated Oligoaniline Derivatives. <i>Journal of the American Chemical Society</i> , 2018, 140, 9417-9425.	6.6	62
3	Theoretical insight into the degradation of p-nitrophenol by OH radicals synergized with other active oxidants in aqueous solution. <i>Journal of Hazardous Materials</i> , 2020, 389, 121901.	6.5	62
4	Embedding Cobalt Atom Clusters in CNT@Wired MoS ₂ Tube@CNT Tube Nanostructures with Enhanced Sulfur Immobilization and Catalyzation for Li–S Batteries. <i>Small</i> , 2021, 17, e2102710.	5.2	52
5	Formation of a series of stable pillar[5]arene-based pseudo[1]-rotaxanes and their [1]rotaxanes in the crystal state. <i>Scientific Reports</i> , 2016, 6, 28748.	1.6	40
6	Self-Assembled Supramolecular Polyoxometalate Hybrid Architecture as a Multifunctional Oxidation Catalyst. <i>ACS Applied Materials & Interfaces</i> , 2019, 11, 38708-38718.	4.0	38
7	Development of the Visible–Light Response of CeO ₂ with a high Ce ³⁺ Content and Its Photocatalytic Properties. <i>ChemCatChem</i> , 2018, 10, 1267-1271.	1.8	37
8	A new precursor to synthesize g-C ₃ N ₄ with superior visible light absorption for photocatalytic application. <i>Catalysis Science and Technology</i> , 2017, 7, 1826-1830.	2.1	35
9	The roles of HO•, ClO• and BrO• radicals in caffeine degradation: A theoretical study. <i>Science of the Total Environment</i> , 2021, 768, 144733.	3.9	31
10	Gaseous and heterogeneous reactions of low-molecular-weight (LMW) unsaturated ketones with O ₃ : Mechanisms, kinetics, and effects of mineral dust in tropospheric chemical processes. <i>Chemical Engineering Journal</i> , 2020, 395, 125083.	6.6	27
11	Molecular diversity of the three-component reaction of α -amino acids, dialkyl acetylenedicarboxylates and N-substituted maleimides. <i>Organic and Biomolecular Chemistry</i> , 2016, 14, 6497-6507.	1.5	25
12	Specific binding and inhibition of 6-benzylaminopurine to catalase: Multiple spectroscopic methods combined with molecular docking study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2014, 123, 327-335.	2.0	23
13	Full insights into the roles of pH on hydroxylation of aromatic acids/bases and toxicity evaluation. <i>Water Research</i> , 2021, 190, 116689.	5.3	22
14	Ozonation of diclofenac in the aqueous solution: Mechanism, kinetics and ecotoxicity assessment. <i>Environmental Research</i> , 2020, 188, 109713.	3.7	19
15	Water determines the products: an unexpected Brønsted acid-catalyzed PO–R cleavage of P(<i>sc</i>) esters selectively producing P(O)–H and P(O)–R compounds. <i>Green Chemistry</i> , 2019, 21, 2916-2922.	4.6	18
16	Pillar[5]arene-based [3]rotaxanes: Convenient construction via multicomponent reaction and pH responsive self-assembly in water. <i>Chinese Chemical Letters</i> , 2020, 31, 1550-1553.	4.8	18
17	Theoretical investigation on the contribution of HO•, SO ₄ ⁻ and CO ₃ ⁻ radicals to the degradation of phenacetin in water: Mechanisms, kinetics, and toxicity evaluation. <i>Ecotoxicology and Environmental Safety</i> , 2020, 204, 110977.	2.9	18
18	Highly Efficient Dehydrogenation of 2,3-Butanediol Induced by Metal–Support Interface over Cu–SiO ₂ Catalysts. <i>ACS Sustainable Chemistry and Engineering</i> , 2020, 8, 15716-15731.	3.2	18

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19	Composite polymerized molecular imprinting membrane-based electrochemical sensor for sensitive determination of curcumin by using 4-pentenoyl-aminoacyl-chitosan oligosaccharide as functional monomer oligomer. <i>Journal of Electroanalytical Chemistry</i> , 2020, 879, 114793.	1.9	18
20	Computational electrochemistry of Pillar[5]quinone cathode material for lithium-ion batteries. <i>Computational Materials Science</i> , 2017, 137, 233-242.	1.4	17
21	Highly Selective and Efficient Reduction of CO ₂ to Methane by Activated Alkaline Earth Metal Hydrides without a Catalyst. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 4831-4841.	3.2	17
22	Acetaminophen degradation by hydroxyl and organic radicals in the peracetic acid-based advanced oxidation processes: Theoretical calculation and toxicity assessment. <i>Journal of Hazardous Materials</i> , 2021, 416, 126250.	6.5	17
23	Effect of pH on H^+ -induced degradation progress of syringol/syringaldehyde and health effect. <i>Chemosphere</i> , 2020, 255, 126893.	4.2	15
24	Bimetallic sulfide Co ₈ FeS ₈ /N-C dodecahedral nanocages via cation exchange as counter electrode for dye-sensitized solar cells. <i>Journal of Alloys and Compounds</i> , 2020, 829, 154526.	2.8	15
25	Degradation of prosulfocarb by hydroxyl radicals in gas and aqueous phase: Mechanisms, kinetics and toxicity. <i>Ecotoxicology and Environmental Safety</i> , 2020, 191, 110175.	2.9	15
26	Theoretical investigation of pillar[4]quinone as a cathode active material for lithium-ion batteries. <i>Journal of Molecular Modeling</i> , 2017, 23, 105.	0.8	14
27	Inorganic-organic hybrid supramolecular architectures based on Keggin polyoxometalates and crown ether: synthesis, crystal structure and electrochemical properties. <i>CrystEngComm</i> , 2021, 23, 8482-8489.	1.3	14
28	Electronic and optical properties of germanene/MoS ₂ heterobilayers: first principles study. <i>Journal of Molecular Modeling</i> , 2018, 24, 333.	0.8	12
29	Construction of Polyfunctionalized 2,4-Dioxo-8-azaspiro[5.5]undec-9-enes and 2,4,8-Triazaspiro[5.5]undec-9-enes via a Domino [2+2+2] Cycloaddition Reaction. <i>Journal of Organic Chemistry</i> , 2021, 86, 1827-1842.	1.7	12
30	Theoretical study of solvent effects on the structures and isomerization of silylenoid H ₂ SiLiF. <i>Structural Chemistry</i> , 2007, 18, 65-70.	1.0	11
31	Mechanisms and Kinetic Parameters for the Gas-Phase Reactions of 3-Methyl-3-buten-2-one and 3-Methyl-3-penten-2-one with Ozone. <i>Journal of Physical Chemistry A</i> , 2019, 123, 2745-2755.	1.1	11
32	Density functional theory study of selective aerobic oxidation of cyclohexane: the roles of acetic acid and cobalt ion. <i>Journal of Molecular Modeling</i> , 2019, 25, 71.	0.8	11
33	Heteroatom-bridged pillar[4]quinone: evolutionary active cathode material for lithium-ion battery using density functional theory. <i>Journal of Chemical Sciences</i> , 2021, 133, 1.	0.7	11
34	Ionic S _N 2 Nucleophilic Substitution in N-Methylaniline-Induced Si-Si Bond Cleavages of Si ₂ Cl ₆ . <i>Chemistry - A European Journal</i> , 2016, 22, 5010-5016.	1.7	10
35	Theoretical study of a novel imino bridged pillar[5]arene derivative. <i>Chemical Physics Letters</i> , 2016, 662, 25-30.	1.2	10
36	New cytotoxic trichothecene macrolide epimers from endophytic <i>Myrothecium roridum</i> IFB-E012. <i>Journal of Antibiotics</i> , 2016, 69, 652-655.	1.0	10

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37	Theoretical study on host-guest interaction between pillar[4]arene and molecules or ions. <i>Journal of Molecular Modeling</i> , 2018, 24, 199.	0.8	10
38	Theoretical prediction of structures and inclusion properties of heteroatom-bridged pillar[n]arenes. <i>Structural Chemistry</i> , 2020, 31, 329-337.	1.0	9
39	Reactivity of aromatic contaminants towards nitrate radical in tropospheric gas and aqueous phase. <i>Journal of Hazardous Materials</i> , 2021, 401, 123396.	6.5	9
40	Ozonolysis of Permethrin in the Atmosphere: Mechanism, Kinetics, and Evaluation of Toxicity. <i>Journal of Physical Chemistry A</i> , 2021, 125, 7705-7715.	1.1	9
41	Unravelling the effects of complexation of transition metal ions on the hydroxylation of catechol over the whole pH region. <i>Journal of Environmental Sciences</i> , 2022, 115, 392-402.	3.2	7
42	Template-assisted synthesis of LiNi _{0.8} Co _{0.15} Al _{0.05} O ₂ hollow nanospheres as cathode material for lithium ion batteries. <i>Journal of Materials Science</i> , 2020, 55, 9493-9503.	1.7	6
43	Degradation mechanisms, kinetics and eco-toxicity assessment of 2,4-Dinitrophenol by oxygen-containing free radicals in aqueous solution. <i>Molecular Physics</i> , 2021, 119, e1886365.	0.8	6
44	[Mo ₃ S ₁₃] ²⁺ as bidirectional cluster catalysts for high-performance Li-S batteries. <i>Catalysis Science and Technology</i> , 2022, 12, 3431-3435.	2.1	6
45	4-Pentenoyl-oleucyl-chitosan oligosaccharide and acrylamide functional monomer-dependent hybrid bilayer molecularly imprinted membrane for sensitive electrochemical sensing of bisphenol A. <i>RSC Advances</i> , 2021, 11, 36769-36776.	1.7	5
46	Computational simulation study on the anion recognition properties of functionalized tetraphenyl porphyrins. <i>Journal of Molecular Modeling</i> , 2015, 21, 140.	0.8	4
47	Bidirectional heterostructures consisting of graphene and lateral MoS ₂ /WS ₂ composites: a first-principles study. <i>RSC Advances</i> , 2019, 9, 34986-34994.	1.7	4
48	Atmospheric ozonolysis of crotonaldehyde in the absence and presence of hydroxylated silica oligomer cluster adsorption. <i>Chemosphere</i> , 2021, 281, 130996.	4.2	4
49	Theoretical study on halide anion recognitions by pyrrole- and benzene-strapped calix[4]pyrroles. <i>Computational and Theoretical Chemistry</i> , 2013, 1007, 1-8.	1.1	3
50	Theoretical Study of Ozonation of Methylparaben and Ethylparaben in Aqueous Solution. <i>Journal of Physical Chemistry A</i> , 2020, 124, 10967-10976.	1.1	3
51	Adsorption of [BF ₄] ⁻ anion-based ionic liquids on phosphorene, arsenene, and antimonene: A density functional theory study. <i>International Journal of Quantum Chemistry</i> , 2021, 121, e26668.	1.0	3
52	Theoretical insights into the degradation of swep by hydroxyl radicals in atmosphere and water environment: Mechanisms, kinetics and toxicity. <i>Science of the Total Environment</i> , 2022, 816, 151651.	3.9	3
53	Cycloaddition reactions of N-heterocyclic stable silylenes with ethylene and formaldehyde. <i>Journal of Organometallic Chemistry</i> , 2011, 696, 841-845.	0.8	2
54	Theoretical study on ion-pair recognition of M+X- (M = Li, Na, K and X = F, Cl, Br) by formylaminocalix[4]arene derivatives. <i>Journal of Molecular Modeling</i> , 2012, 18, 4985-4993.	0.8	2

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55	Theoretical study on the ion-pair recognition of Na ⁺ /X ⁻ (X = F ⁻ , Cl ⁻ , Br ⁻) by urea calix[4]bis crown-3 derivative. <i>Molecular Simulation</i> , 2013, 39, 621-628.	0.9	2
56	Cyclodimerization of 3-phenacylideneoxindolines with amino esters for the synthesis of dispiro[indoline-3,1 ⁻² -cyclopentane-3 ⁻² ,3 ⁻³ -indolines]. <i>Heterocyclic Communications</i> , 2017, 23, 297-303.	0.6	2
57	Solubility of Acetoguanamine in Twelve Neat Solvents from 283.15 to 323.15 K. <i>Journal of Chemical & Engineering Data</i> , 2019, 64, 4546-4550.	1.0	2
58	Theoretical study on a kind of cyclization mechanism of boron trichloride and hexamethyldisilazane to form the borazine ring for SiBCN ceramics precursor. <i>Ceramics International</i> , 2021, 47, 6068-6076.	2.3	2
59	Theoretical prediction on photoelectric and supramolecular properties of benzoquinone-tetrathiafulvalene macrocyclic molecules. <i>Journal of Molecular Modeling</i> , 2021, 27, 157.	0.8	1
60	H-bond catalytic mechanism of aromatic electrophilic substitution between phenol and formaldehyde. <i>Tetrahedron</i> , 2020, 76, 131055.	1.0	0
61	Theoretical Study on the Mechanism of the Initial Ammonolysis Reactions of the Main Raw Materials for the Formation of SiBCN Ceramics. <i>ChemistrySelect</i> , 2021, 6, 12233-12245.	0.7	0