

# Futing Xia

## List of Publications by Year in descending order

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Version: 2024-02-01

9  
papers

93  
citations

1684188  
5  
h-index

1474206  
9  
g-index

9  
all docs

9  
docs citations

9  
times ranked

95  
citing authors

#	ARTICLE	IF	CITATIONS
1	Improved catalytic activity and N <sub>2</sub> selectivity of Fe-Mn-O x catalyst for selective catalytic reduction of NO by NH <sub>3</sub> at low temperature. <i>Research on Chemical Intermediates</i> , 2018, 44, 2703-2717.	2.7	27
2	Removal of toluene over bi-metallic Pt-Pd-SBA-15 catalysts: Kinetic and mechanistic study. <i>Microporous and Mesoporous Materials</i> , 2020, 302, 110111.	4.4	23
3	Alkaline hydrolysis of ethylene phosphate: An <i>ab initio</i> study by supermolecule model and polarizable continuum approach. <i>Journal of Computational Chemistry</i> , 2011, 32, 2545-2554.	3.3	11
4	Density functional calculations on alcoholysis and thiolysis of phosphate triesters: Stepwise or concerted?. <i>Computational and Theoretical Chemistry</i> , 2013, 1017, 60-71.	2.5	10
5	Catalytic hydrolysis of HCN over Fe-Ti-O catalysts prepared by different calcination temperatures: Effect of Fe chemical valence and Ti phase. <i>Microporous and Mesoporous Materials</i> , 2020, 292, 109753.	4.4	10
6	Effect of sulfur substitution for methanolysis of paraoxon: CO vs. PO bond cleavage from density-functional theory. <i>Computational and Theoretical Chemistry</i> , 2012, 982, 8-16.	2.5	5
7	The promotional effect of SO <sub>4</sub> <sup>2-</sup> on N <sub>2</sub> selectivity for selective catalytic oxidation of ammonia over RuO <sub>2</sub> /ZrO <sub>2</sub> catalyst. <i>Research on Chemical Intermediates</i> , 2020, 46, 803-820.	2.7	3
8	Density functional calculations on the effect of sulfur substitution for 2-hydroxypropyl-p-nitrophenyl phosphate: C-O vs. P-O bond cleavage. <i>Bioorganic Chemistry</i> , 2012, 40, 99-107.	4.1	2
9	Theoretical studies on the effect of sulfur substitution for the methanolysis of cyclic and acyclic phosphate esters. <i>Computational and Theoretical Chemistry</i> , 2014, 1048, 35-45.	2.5	2