

Yeonjoon Kim

List of Publications by Year in descending order

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Version: 2024-02-01

20
papers

653
citations

687363

13
h-index

713466

21
g-index

22
all docs

22
docs citations

22
times ranked

817
citing authors

#	ARTICLE	IF	CITATIONS
1	Predicting Catalytic Pyrolysis Aromatic Selectivity from Pyrolysis Vapor Composition Using Mass Spectra Coupled with Statistical Analysis. <i>ACS Sustainable Chemistry and Engineering</i> , 2022, 10, 234-244.	6.7	3
2	Understanding how chemical structure affects ignition-delay-time $\langle \text{mml:math} \text{xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si2.svg"} \rangle$ -sensitivity. <i>Combustion and Flame</i> , 2021, 225, 377-387.	5.2	7
3	Investigation of structural effects of aromatic compounds on sooting tendency with mechanistic insight into ethylphenol isomers. <i>Proceedings of the Combustion Institute</i> , 2021, 38, 1143-1151.	3.9	10
4	Quantum chemical calculations for over 200,000 organic radical species and 40,000 associated closed-shell molecules. <i>Scientific Data</i> , 2020, 7, 244.	5.3	49
5	Prediction of Hydroxymethylfurfural Yield in Glucose Conversion through Investigation of Lewis Acid and Organic Solvent Effects. <i>ACS Catalysis</i> , 2020, 10, 14707-14721.	11.2	41
6	Prediction of organic homolytic bond dissociation enthalpies at near chemical accuracy with sub-second computational cost. <i>Nature Communications</i> , 2020, 11, 2328.	12.8	128
7	A perspective on biomass-derived biofuels: From catalyst design principles to fuel properties. <i>Journal of Hazardous Materials</i> , 2020, 400, 123198.	12.4	23
8	Graph theory-based reaction pathway searches and DFT calculations for the mechanism studies of free radical-initiated peptide sequencing mass spectrometry (FRIPS MS): a model gas-phase reaction of GGR tri-peptide. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 5057-5069.	2.8	5
9	Isotopic Studies for Tracking Biogenic Carbon during Co-processing of Biomass and Vacuum Gas Oil. <i>ACS Sustainable Chemistry and Engineering</i> , 2020, 8, 2652-2664.	6.7	14
10	Ga/ZSM-5 catalyst improves hydrocarbon yields and increases alkene selectivity during catalytic fast pyrolysis of biomass with co-fed hydrogen. <i>Green Chemistry</i> , 2020, 22, 2403-2418.	9.0	26
11	Performance of ACE-Reaction on 26 Organic Reactions for Fully Automated Reaction Network Construction and Microkinetic Analysis. <i>Journal of Physical Chemistry A</i> , 2019, 123, 4796-4805.	2.5	11
12	Feasibility of Activation Energy Prediction of Gas-Phase Reactions by Machine Learning. <i>Chemistry - A European Journal</i> , 2018, 24, 12354-12358.	3.3	35
13	Efficient structural elucidation of microhydrated biomolecules through the interrogation of hydrogen bond networks. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 8185-8191.	2.8	2
14	Efficient prediction of reaction paths through molecular graph and reaction network analysis. <i>Chemical Science</i> , 2018, 9, 825-835.	7.4	82
15	Poly(amide-imide) materials for transparent and flexible displays. <i>Science Advances</i> , 2018, 4, eaau1956.	10.3	57
16	Enhancing the Activity of Platinum-Based Nanocrystal Catalysts for Organic Synthesis through Electronic Structure Modification. <i>ChemCatChem</i> , 2016, 8, 2450-2454.	3.7	3
17	Computational searching for new stable graphyne structures and their electronic properties. <i>Carbon</i> , 2016, 98, 404-410.	10.3	37
18	Universal Structure Conversion Method for Organic Molecules: From Atomic Connectivity to Three-Dimensional Geometry. <i>Bulletin of the Korean Chemical Society</i> , 2015, 36, 1769-1777.	1.9	36

#	ARTICLE	IF	CITATIONS
19	Bicyclic Bridgehead Phosphoramidite (Briphos) Ligands with Tunable σ -Acceptor Ability and Catalytic Activity in the Rhodium-Catalyzed Conjugate Additions. <i>Organic Letters</i> , 2014, 16, 5490-5493.	4.6	32
20	Efficient Basin-Hopping Sampling of Reaction Intermediates through Molecular Fragmentation and Graph Theory. <i>Journal of Chemical Theory and Computation</i> , 2014, 10, 2419-2426.	5.3	46