

Harish Jangra

List of Publications by Year in descending order

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Version: 2024-02-01

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papers

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1163065

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citing authors

#	ARTICLE	IF	CITATIONS
1	Reactivities of allenic and olefinic Michael acceptors towards phosphines. <i>Chemical Communications</i> , 2022, 58, 3358-3361.	4.1	10
2	Stereoselective and Stereospecific Triflate-Mediated Intramolecular Schmidt Reaction: Ready Access to Alkaloid Skeletons**. <i>Angewandte Chemie</i> , 2021, 133, 10267-10273.	2.0	2
3	Stereoselective and Stereospecific Triflate-Mediated Intramolecular Schmidt Reaction: Ready Access to Alkaloid Skeletons**. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 10179-10185.	13.8	9
4	Radical chain monoalkylation of pyridines. <i>Chemical Science</i> , 2021, 12, 15362-15373.	7.4	7
5	Molecule-Induced Radical Formation (MIRF) Reactions: A Reappraisal. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 6318-6329.	13.8	8
6	A Predictive Model Towards Site-Selective Metalations of Functionalized Heterocycles, Arenes, Olefins, and Alkanes using TMPZnCl ₂ ·LiCl. <i>Angewandte Chemie</i> , 2020, 132, 15102-15109.	2.0	8
7	A Predictive Model Towards Site-Selective Metalations of Functionalized Heterocycles, Arenes, Olefins, and Alkanes using TMPZnCl ₂ ·LiCl. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 14992-14999.	13.8	20
8	Phosphine-catalyzed [3 + 2] annulation of 2-aminoacrylates with allenates and mechanistic studies. <i>Catalysis Science and Technology</i> , 2020, 10, 3959-3964.	4.1	6
9	Construction of α,β -disubstituted α -Amino Acid Derivatives via aza-Morita-Baylis-Hillman Reactions of α -Aminoacrylates with Activated Olefins. <i>ChemCatChem</i> , 2020, 12, 1143-1147.	3.7	2
10	Ferrocenes with a Persulfurated Cyclopentadienyl Ring: Synthesis, Structural Studies, and Optoelectronic Properties. <i>Chemistry - A European Journal</i> , 2019, 25, 12684-12688.	3.3	6
11	Fast Microsecond Dynamics of the Protein-Water Network in the Active Site of Human Carbonic Anhydrase II Studied by Solid-State NMR Spectroscopy. <i>Journal of the American Chemical Society</i> , 2019, 141, 19276-19288.	13.7	46
12	Lewis Acid Directed Regioselective Metalations of Pyridazine. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 9244-9247.	13.8	20
13	Kinetics and Mechanism of Oxirane Formation by Darzens Condensation of Ketones: Quantification of the Electrophilicities of Ketones. <i>Journal of the American Chemical Society</i> , 2018, 140, 5500-5515.	13.7	34
14	Nucleophilicity and Electrophilicity Parameters for Predicting Absolute Rate Constants of Highly Asynchronous 1,3-Dipolar Cycloadditions of Aryldiazomethanes. <i>Journal of the American Chemical Society</i> , 2018, 140, 16758-16772.	13.7	52
15	A third generation of radical fluorinating agents based on N-fluoro-N-arylsulfonamides. <i>Nature Communications</i> , 2018, 9, 4888.	12.8	58
16	Electrostatic Effects on the Stability of Peptide Radicals. <i>Journal of Physical Chemistry B</i> , 2018, 122, 8880-8890.	2.6	5
17	Unique Stereoselective Homolytic C=O Bond Activation in Diketopiperazine-Derived Alkoxyamines by Adjacent Amide Pyramidalization. <i>Chemistry - A European Journal</i> , 2018, 24, 15336-15345.	3.3	7
18	Quantification and Theoretical Analysis of the Electrophilicities of Michael Acceptors. <i>Journal of the American Chemical Society</i> , 2017, 139, 13318-13329.	13.7	168

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19	Conformational Preferences in Small Peptide Models: The Relevance of <i>cis</i> / <i>trans</i> Conformations. Chemistry - A European Journal, 2016, 22, 13328-13335.	3.3	4
20	Detour matrix-based adjacent path eccentric distance sum indices for QSAR/QSPR. Part I: development and evaluation. International Journal of Computational Biology and Drug Design, 2014, 7, 295.	0.3	1
21	Detour Cum Distance Matrix Based Topological Descriptors for QSAR/QSPR Part-I: Development and Evaluation. Letters in Drug Design and Discovery, 2014, 11, 844-863.	0.7	0
22	Formation/Fate of Reactive Metabolites from General Anesthetics and A Comparison of Toxic and Non-Toxic Analogues: A DFT Study. Drug Metabolism Letters, 2013, 6, 221-234.	0.8	2
23	Fourth generation detour matrix-based topological indices for QSAR/QSPR - Part-1: development and evaluation. International Journal of Computational Biology and Drug Design, 2012, 5, 335.	0.3	3