

Daisuke Kaneno

List of Publications by Year in descending order

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27
papers

1,036
citations

687220

13
h-index

501076

28
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34
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docs citations

34
times ranked

1335
citing authors

#	ARTICLE	IF	CITATIONS
1	Synthesis and photophysical properties of a new push-pull pyrene dye with green-to-far-red emission and its application to human cellular and skin tissue imaging. <i>Journal of Materials Chemistry B</i> , 2022, 10, 1641-1649.	2.9	9
2	Effects of the Monomeric Components of Poly-hydroxybutyrate-co-hydroxyhexanoate on the Growth of <i>Vibrio penaeicida</i> In Vitro and on the Survival of Infected Kuruma Shrimp (<i>Marsupenaeus japonicus</i>). <i>Animals</i> , 2021, 11, 567.	1.0	8
3	Protecting group-free method for synthesis of N-glycosyl carbamates and an assessment of the anomeric effect of nitrogen in the carbamate group. <i>Carbohydrate Research</i> , 2021, 505, 108280.	1.1	3
4	Bright and two-photon active red fluorescent dyes that selectively move back and forth between the mitochondria and nucleus upon changing the mitochondrial membrane potential. <i>Journal of Materials Chemistry B</i> , 2018, 6, 7396-7401.	2.9	15
5	Iriomoteolides-10a and 12a, Cytotoxic Macrolides from Marine Dinoflagellate <i>Amphidinium</i> Species. <i>Chemical and Pharmaceutical Bulletin</i> , 2016, 64, 1019-1023.	0.6	15
6	Synthesis of trans-2,6-Piperidinedicarboxamide Using the Ugi Reaction. A Plausible Model for the Biosynthesis of Halichonadin P. <i>Heterocycles</i> , 2016, 92, 857.	0.4	3
7	Synthesis and characterization of novel polyoxometalates with an inverted-Keggin structure as a new class of building unit. <i>Inorganic Chemistry Communication</i> , 2013, 38, 123-126.	1.8	3
8	Parr™s index to describe both electrophilicity and nucleophilicity. <i>Tetrahedron Letters</i> , 2013, 54, 339-342.	0.7	44
9	Intrinsic reactivity index as a single scale directed toward both electrophilicity and nucleophilicity using frontier molecular orbitals. <i>Tetrahedron</i> , 2013, 69, 4247-4258.	1.0	36
10	Intra- and Intermolecular Reaction Selectivities of $\hat{1}^3$ -Substituted Adamantanylidenes. <i>Journal of Organic Chemistry</i> , 2012, 77, 1340-1360.	1.7	17
11	Intramolecular titanium-promoted deoxygenative cyclization to 9-substituted-1,2,3,4-tetrahydrofluorene skeleton. <i>Tetrahedron Letters</i> , 2010, 51, 1651-1653.	0.7	3
12	The Danishefsky pyranone puzzle: an explanation based on the exterior frontier orbital extension model. <i>Tetrahedron Letters</i> , 2009, 50, 329-332.	0.7	1
13	Theoretical Study on the Mechanism and Diastereoselectivity of NaBH_4 Reduction. <i>Journal of Physical Chemistry A</i> , 2009, 113, 2578-2583.	1.1	29
14	Solvent effects on the diastereoselection in LiAlH_4 reduction of $\hat{1}^\pm$ -substituted ketones. <i>Tetrahedron Letters</i> , 2008, 49, 4223-4226.	0.7	13
15	The Importance of Lone Pair Delocalizations: Theoretical Investigations on the Stability of cis and trans Isomers in 1,2-Halodiazenes. <i>Journal of Organic Chemistry</i> , 2008, 73, 5429-5435.	1.7	17
16	The Origin of Cis Effect in 1,2-Dihaloethenes: The Quantitative Comparison of Electron Delocalizations and Steric Exchange Repulsions. <i>Bulletin of the Chemical Society of Japan</i> , 2008, 81, 1415-1422.	2.0	28
17	The Importance of Lone Pair Electron Delocalization in the cis-trans Isomers of 1,2-Dibromoethenes. <i>Chemistry Letters</i> , 2005, 34, 1190-1191.	0.7	15
18	Origin of Diastereoselection in the Hydrosilylation of Chiral N-Acyliminium Intermediates Derived from Pyroglutamic Acid. <i>Angewandte Chemie - International Edition</i> , 2004, 43, 2412-2415.	7.2	11

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19	Origin of facial diastereoselection in 2-adamantyl cations. Theoretical evidence against the Felkin-Anh and the Cieplak models. <i>Tetrahedron Letters</i> , 2004, 45, 4559-4562.	0.7	7
20	Origin of Facial Diastereoselection. Evidence for Negative Role of Antiperiplanar Hyperconjugation Effects in the Transition State of Carbene Insertion. <i>Organic Letters</i> , 2003, 5, 2947-2949.	2.4	21
21	Stereoelectronic and conformational effects on the stereochemical course of reduction of bicyclo[3.3.1]nonane 1,3-diketones. <i>Canadian Journal of Chemistry</i> , 2001, 79, 1598-1605.	0.6	5
22	Facial diastereoselection of hydride reduction of 1,3-diheteran-5-ones: application of the exterior frontier orbital extension model. <i>Heteroatom Chemistry</i> , 2001, 12, 358-368.	0.4	8
23	Prediction of Facial Diastereoselection with the Exterior Frontier Orbital Extension Model (EFOE) Tj ETQq1 1 0.784314 rgBT /Overlock 1	0.0	1
24	Reversal of fac-facial diastereoselection in the hydride reduction of selenanones. Further application of the exterior frontier orbital extension model. <i>Tetrahedron Letters</i> , 2000, 41, 4597-4601.	0.7	11
25	Origin of p-Facial Diastereoselection in Carbonyl Addition. Application of the Exterior Frontier Orbital Extension Model to 1,3-Diheteran-5-ones (Heteroatom = O, S). <i>Heterocycles</i> , 2000, 52, 1435.	0.4	13
26	Stereochemical Determination of Acyclic Structures Based on Carbon-13 Proton Spin-Coupling Constants. A Method of Configuration Analysis for Natural Products. <i>Journal of Organic Chemistry</i> , 1999, 64, 866-876.	1.7	697
27	Origin of fac-Facial Diastereoselection in Nucleophilic Addition to 1,3-Diheteran-5-ones (Heteroatom = O,) Tj ETQq1 1 0.784314 rgBT /Overlock 1161-1162.	0.7	2