

Boubaker Taoufik

List of Publications by Year in descending order

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33
papers

934
citations

566801

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h-index

454577

30
g-index

33
all docs

33
docs citations

33
times ranked

538
citing authors

#	ARTICLE	IF	CITATIONS
1	Nucleophilic Reactivities of Indoles. <i>Journal of Organic Chemistry</i> , 2006, 71, 9088-9095.	1.7	281
2	Ranking the Reactivity of Superelectrophilic Heteroaromatics on the Electrophilicity Scale. <i>Journal of Organic Chemistry</i> , 2005, 70, 6242-6253.	1.7	104
3	Nitrobenzoxadiazoles and related heterocycles: a relationship between aromaticity, superelectrophilicity and pericyclic reactivity. <i>Organic and Biomolecular Chemistry</i> , 2006, 4, 1910.	1.5	51
4	Mayr electrophilicity predicts the dual Diels-Alder and π -adduct formation behaviour of heteroaromatic super-electrophiles. <i>Organic and Biomolecular Chemistry</i> , 2007, 5, 1744-1751.	1.5	48
5	Electrophilicity of aromatic triflones in π -complexation processes. <i>Organic and Biomolecular Chemistry</i> , 2008, 6, 4041.	1.5	44
6	Steric effects on the intrinsic reactivity of nitrotriphenylmethanes. <i>Journal of Organic Chemistry</i> , 1992, 57, 3924-3929.	1.7	39
7	An extremely highly electrophilic heteroaromatic structure: 4,6-dinitrotetrazolo[1,5-a]pyridine Electronic supplementary information (ESI) available: Fig. S1 Effect of pH on the ionization of the OH group of C-4a to give the diadduct D-4a in aqueous solution. See http://www.rsc.org/suppdata/ob/b3/b306437a . <i>Organic and Biomolecular Chemistry</i> , 2003, 1, 2764.	1.5	38
8	Water and hydroxide ion pathways in the π -complexation of superelectrophilic 2-aryl-4,6-dinitrobenzotriazole 1-oxides in aqueous solution. A kinetic and thermodynamic study. <i>Perkin Transactions II RSC</i> , 2002, , 1627-1633.	1.1	32
9	The versatile reactivity of 2-aryl-4,6-dinitrobenzotriazole 1-oxides in Diels-Alder type condensations and in π -complexation - A relationship between superelectrophilicity and pericyclic reactivity. <i>Canadian Journal of Chemistry</i> , 2001, 79, 1617-1623.	0.6	29
10	Superelectrophilicity of the Nitroolefinic Fragment of 4-Nitrobenzodifuroxan in Michael-Type Reactions with Indoles: A Kinetic Study in Acetonitrile. <i>Chemistry - A European Journal</i> , 2007, 13, 8317-8324.	1.7	28
11	Kinetics of alkaline hydrolysis of p -substituted benzylidenemalononitriles in 50% aqueous acetonitrile: substituent effects and quantification of the electrophilic reactivity. <i>Journal of Physical Organic Chemistry</i> , 2014, 27, 484-489.	0.9	28
12	The Ambident Reactivity of 2,4,6-Tris(trifluoromethanesulfonyl)anisole in Methanol: Using the SO_2CF_3 Group as a Tool to Reach the Superelectrophilic Dimension in π -Complexation Processes. <i>Chemistry - A European Journal</i> , 2009, 15, 12018-12029.	1.7	27
13	Substituent effects on the reactions of aromatic triflones with p -anilines in methanol: Low intrinsic reactivity and transition state imbalances. <i>International Journal of Chemical Kinetics</i> , 2011, 43, 255-262.	1.0	24
14	Activation of the aromatic system by the SO_2CF_3 group: Kinetics study and structure-reactivity relationships. <i>International Journal of Chemical Kinetics</i> , 2010, 42, 203-210.	1.0	23
15	Ionization of nitrotriphenylmethanes. Remarkable kinetic evidence for steric inhibition to resonance and F-strain. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1990, , 1899.	0.9	20
16	Relationships between experimental and theoretical scales of electrophilicity of 7-L-4-nitrobenzofurazans. <i>Journal of Molecular Structure</i> , 2021, 1224, 128843.	1.8	14
17	The ambident electrophilic behavior of 5-nitro- β -thiophenes in π -complexation processes. <i>International Journal of Chemical Kinetics</i> , 2018, 50, 659-669.	1.0	11
18	Influence of solvent mixture on nucleophilicity parameters: the case of pyrrolidine in methanol-acetonitrile. <i>RSC Advances</i> , 2020, 10, 28635-28643.	1.7	11

#	ARTICLE	IF	CITATIONS
19	Kinetics and quantification of the electrophilic reactivities of substituted thiophenes and structure–reactivity relationships. <i>International Journal of Chemical Kinetics</i> , 2018, 50, 582-590.	1.0	9
20	Ambident electrophilicity of 4-nitrobenzochalcogenadiazoles: Kinetic studies and structure–reactivity relationships. <i>International Journal of Chemical Kinetics</i> , 2020, 52, 669-680.	1.0	9
21	Electrophilic reactivities of 7- and 4-nitrobenzofurazans in π -complexation processes: Kinetic studies and structure–reactivity relationships. <i>International Journal of Chemical Kinetics</i> , 2020, 52, 655-668.	1.0	9
22	Single-Electron Transfer in π -Complexation Reactions of 2,6-Dimethoxy-3,5-dinitropyridine with Para-X-Phenoxide Anions in Aqueous Solution. <i>International Journal of Chemical Kinetics</i> , 2016, 48, 523-530.	1.0	8
23	S _N Ar reactions of substituted pyridines with secondary amines in aqueous solution: Kinetic and reactivity indices in density functional theory. <i>International Journal of Chemical Kinetics</i> , 2019, 51, 249-257.	1.0	8
24	Regioselectivity and High Electrophilicity in the π Complexation of Aromatic Triflones. <i>European Journal of Organic Chemistry</i> , 2014, 2014, 6774-6786.	1.2	7
25	Structure–property relationships in para-substituted nitrobenzofurazans: electrochemical, optical, and theoretical analysis. <i>Chemical Papers</i> , 2022, 76, 4059-4080.	1.0	7
26	Electrophilicities of 4-nitrobenzochalcogenadiazoles. <i>ChemistrySelect</i> , 2020, 5, 7648-7657.	0.7	6
27	First-principles study of the reaction mechanism governing the S _N Ar of the dimethylamine on 2-methoxy-5-nitrothiophenes. <i>Theoretical Chemistry Accounts</i> , 2020, 139, 1.	0.5	5
28	Structure-Property Relationships in Benzofurazan Derivatives: A Combined Experimental and DFT/TD-DFT Investigation. , 0, , .		4
29	The N-alkylation of 4-nitrobenzochalcogenadiazoles: Synthesis and theoretical approach. <i>Journal of Molecular Structure</i> , 2019, 1197, 80-86.	1.8	3
30	Quantification of the Electrophilic Reactivities of Benzotriazoles and Structure–Reactivity Relationships. <i>ChemistrySelect</i> , 2021, 6, 4424-4431.	0.7	3
31	Exploring the reactivity of benzotriazole derivatives: Mayr's approach and density functional theory analysis. <i>Journal of Molecular Structure</i> , 2022, 1247, 131310.	1.8	3
32	Nucleophilicities of para-substituted aniline radical cations in acetonitrile: Kinetic investigation and structure–reactivity relationships. <i>International Journal of Chemical Kinetics</i> , 0, , .	1.0	1
33	Coordination Chemistry of a Bis(Tetrazine) Tweezer: A Case of Host-Guest Behavior with Silver Salts. <i>Molecules</i> , 2021, 26, 2705.	1.7	0