

Zoran Zdravkovski

List of Publications by Year in descending order

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48
papers

934
citations

516710

16
h-index

454955

30
g-index

48
all docs

48
docs citations

48
times ranked

650
citing authors

#	ARTICLE	IF	CITATIONS
1	Elementary, my dear Watson! The making of a collection of the natural elements. <i>Macedonian Journal of Chemistry and Chemical Engineering</i> , 2022, 41, .	0.6	0
2	Application of the electro-Fenton process to mesotrione aqueous solutions: Kinetics, degradation pathways, mineralization and evolution of the toxicity. <i>Macedonian Journal of Chemistry and Chemical Engineering</i> , 2014, 33, 121.	0.6	6
3	Macedonian journal of chemistry and chemical engineering: open journal systems--editor's perspective. <i>Prilozi - Makedonska Akademija Na Naukite I Umetnostite Oddelenie Za Medicinski Nauki</i> , 2014, 35, 51-5.	0.5	0
4	Determination of dimethoate, 2,4-dichlorophenoxy acetic acid, mecoprop and linuron pesticides in environmental waters in Republic of Macedonia by high performance liquid chromatography. <i>Macedonian Journal of Chemistry and Chemical Engineering</i> , 2013, 27, 25.	0.6	6
5	Degradation and mineralization of sulcotrione and mesotrione in aqueous medium by the electro-Fenton process: a kinetic study. <i>Environmental Science and Pollution Research</i> , 2012, 19, 1563-1573.	5.3	42
6	Development of a trap for fuel exhaust particulate matter under driving conditions and GC-MS method for their analysis. <i>Macedonian Journal of Chemistry and Chemical Engineering</i> , 2012, 30, 97.	0.6	1
7	Stopper with Interchangeable Septa for SPME. <i>Journal of Chemical Education</i> , 2009, 86, 924.	2.3	0
8	Theoretical Study of the Diastereofacial Isomers of Aldrin and Dieldrin. <i>International Journal of Molecular Sciences</i> , 2006, 7, 35-46.	4.1	2
9	Bioluminescence determination of enzyme activity of firefly luciferase in the presence of pesticides. <i>Luminescence</i> , 2005, 20, 192-196.	2.9	12
10	Rapid and simple method for direct determination of several amphetamines in seized tablets by GC-FID. <i>Forensic Science International</i> , 2005, 152, 199-203.	2.2	15
11	Theoretical study of the Diels-Alder reactions of benzene with fluorinated dienophiles. <i>Computational and Theoretical Chemistry</i> , 2004, 684, 99-102.	1.5	2
12	Supercritical Fluid Extraction of Quercetin and Rutin from Hyperici Herba. <i>Journal of Liquid Chromatography and Related Technologies</i> , 2003, 26, 2517-2533.	1.0	22
13	The theoretical study of a borane catalyzed azide anion addition to fluorinated acetonitriles. <i>Computational and Theoretical Chemistry</i> , 1997, 389, 69-74.	1.5	2
14	Tetrafluoroberyllate(2-) Ions as Hydrogen-Bond Proton Acceptors: Spectroscopic Evidence. <i>Spectroscopy Letters</i> , 1996, 29, 867-875.	1.0	2
15	Theoretical study of the Diels-Alder reaction between the S-methylthiophenium ion and ethene. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1996, , 455-459.	0.9	11
16	Molecular modeling methodology of β -cyclodextrin inclusion complexes. <i>Computational and Theoretical Chemistry</i> , 1996, 366, 113-117.	1.5	42
17	An ab initio study of heterodienophiles addition to 2,3-diaza-1,3-butadiene: An example of endo-lone-pair effect on the reaction energy barrier. <i>Journal of Computational Chemistry</i> , 1996, 17, 298-305.	3.3	8
18	Theoretical investigation of cis- and trans-nitric oxide dimers with ab initio and density functional Gaussian-type orbital approach. <i>International Journal of Quantum Chemistry</i> , 1995, 54, 161-166.	2.0	98

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19	Theoretical investigation of the conrotatory ring opening of cyclobutene and 1, 2-dihydro-1, 2-diazacyclobutadienes with ab initio and density functional Gaussian-type-orbital approach. <i>International Journal of Quantum Chemistry</i> , 1995, 56, 115-123.	2.0	41
20	Ab initio transition structures for hetero Diels-Alder cycloadditions to furan. <i>Computational and Theoretical Chemistry</i> , 1995, 331, 215-221.	1.5	12
21	Diels-Alder transition structures of hetero-dienophile addition to 4H-pyrazole calculated by ab initio methods. <i>Computational and Theoretical Chemistry</i> , 1995, 331, 229-234.	1.5	9
22	PM3 study of the stereochemistry of heterodienophile cycloadditions to pyrrole: endo lone pair effect. <i>Computational and Theoretical Chemistry</i> , 1995, 332, 39-45.	1.5	13
23	Calculations of boron trifluoride catalyzed 1,3-dipolar additions of azide ion to organic nitriles. <i>Computational and Theoretical Chemistry</i> , 1995, 332, 127-135.	1.5	1
24	Ab initio study of the 1,3-cycloaddition of methyl azide to fluorinated acetonitriles. <i>Computational and Theoretical Chemistry</i> , 1995, 333, 209-214.	1.5	5
25	Why is tetrazole not practical as a diene in Diels-Alder reactions? An ab initio theoretical study. <i>Computational and Theoretical Chemistry</i> , 1995, 337, 9-16.	1.5	11
26	Theoretical study of azide anion addition to nonpolar and polar double and triple bonds. <i>Computational and Theoretical Chemistry</i> , 1995, 343, 149-155.	1.5	3
27	Reaction of Imidazoles with Ethylene and Singlet Oxygen. An ab Initio Theoretical Study. <i>Journal of Organic Chemistry</i> , 1995, 60, 2865-2869.	3.2	16
28	Theoretical Study of the BH ₃ -Catalyzed Hetero Diels-Alder Reaction between Ethylene and Nitrosoethylene. <i>Journal of Organic Chemistry</i> , 1995, 60, 3163-3168.	3.2	23
29	DFT study of the Diels-Alder reactions between ethylene with buta-1,3-diene and cyclopentadiene. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1995, , 1223-1226.	0.9	145
30	PM3 calculations of Diels-Alder reactions of α -pull-push-activated isoprenes with aceto- and acrylonitrile. <i>Tetrahedron</i> , 1994, 50, 10379-10390.	1.9	7
31	AB initio study of Diels-Alder reactions of 1,3,4-oxadiazole with ethylene, acrylonitrile, maleonitrile, fumaronitrile and 1,1-dicyanoethylene. Inverse order and ratio of endo/exo reactivity. <i>Journal of Physical Organic Chemistry</i> , 1994, 7, 634-640.	1.9	7
32	AB initio calculations of Diels-Alder transition structures for hetero-dienophile additions to cyclopentadiene. <i>Journal of Physical Organic Chemistry</i> , 1994, 7, 641-645.	1.9	19
33	Determination of enantiomeric composition of 2-phenyl-2-(2-piperidyl)acetamide. A routine method for evaluation of enantiomeric purity of primary amides. <i>Tetrahedron: Asymmetry</i> , 1994, 5, 1711-1716.	1.8	10
34	The utility of the PM3 method for predicting the reactivities of cyanoethenes in diels-Alder reactions with pyrrole. <i>Journal of Heterocyclic Chemistry</i> , 1994, 31, 1429-1432.	2.6	24
35	PM3-CI calculation of Diels-Alder transition structures of hetero dienophile addition to butadiene: comparison with PM3 and ab initio generated transition structures. <i>Computational and Theoretical Chemistry</i> , 1994, 315, 85-90.	1.5	10
36	Semiempirical and ab initio transition state calculations for the transformation of N-acetyltetrazoles into corresponding 1,3,4-oxadiazoles. <i>Computational and Theoretical Chemistry</i> , 1994, 309, 241-247.	1.5	17

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37	Comparison of AM1 and PM3 semiempirical to ab initio methods in the study of Diels-Alder reactions of butadiene and cyclopentadiene with cyanoethylenes. Computational and Theoretical Chemistry, 1994, 309, 249-257.	1.5	32
38	Rates of hydrolysis of N-acetylazoles: semiempirical calculations compared to experimental values. Computational and Theoretical Chemistry, 1994, 303, 177-183.	1.5	54
39	Semiempirical and ab initio study of 1,3-dipolar addition of azide anion to organic cyanides. Computational and Theoretical Chemistry, 1994, 312, 11-22.	1.5	18
40	Ab initio study of heterodienophile addition to oxazole. Journal of the Chemical Society Perkin Transactions II, 1994, , 1877.	0.9	16
41	Theoretical Study of the Reactivity and Stereoselectivity of Heterodienophile Addition to 1,3,4-Oxadiazole. Journal of Organic Chemistry, 1994, 59, 3015-3019.	3.2	19
42	Theoretical Study of Ethylene and Vinyl Alcohol Addition to 1,4-Dioxo-1,3-butadiene. Journal of Organic Chemistry, 1994, 59, 7732-7736.	3.2	10
43	On the Formation of Peptide Bonds. Journal of Chemical Education, 1993, 70, 134.	2.3	5
44	Preparation and characterization of deuterated crystalhydrates: An undergraduate inorganic or spectroscopic experiment. Journal of Chemical Education, 1993, 70, 603.	2.3	0
45	Molecular mechanics calculations and comparison of proton, fluorine, and carbon NMR diastereomer discrimination via nonbonding interactions between fluorine-labeled enantiomeric amides and enantiomerically pure chiral solvating agents. Journal of Organic Chemistry, 1993, 58, 5245-5250.	3.2	5
46	A Simple Preparation of Amides from Acids and Amines by Heating of Their Mixture. Synthetic Communications, 1993, 23, 2761-2770.	2.1	125
47	Using MathCAD in chemistry calculations. Journal of Chemical Education, 1991, 68, A95.	2.3	1
48	Role of the Polymer Backbone in the Oxidative Properties of Polymer-Supported Dichromate. Synthetic Communications, 1989, 19, 1317-1324.	2.1	5