

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 | 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 |
| 2 | A Simple Preparation of Amides from Acids and Amines by Heating of Their Mixture. <i>Synthetic Communications</i> , 1993, 23, 2761-2770. | 2.1 | 125 |
| 3 | 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 |
| 4 | Rates of hydrolysis of N-acetylazoles: semiempirical calculations compared to experimental values. <i>Computational and Theoretical Chemistry</i> , 1994, 303, 177-183. | 1.5 | 54 |
| 5 | Molecular modeling methodology of β -cyclodextrin inclusion complexes. <i>Computational and Theoretical Chemistry</i> , 1996, 366, 113-117. | 1.5 | 42 |
| 6 | 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 |
| 7 | 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 |
| 8 | Comparison of AM1 and PM3 semiempirical to ab initio methods in the study of Diels-Alder reactions of butadiene and cyclopentadiene with cyanoethylenes. <i>Computational and Theoretical Chemistry</i> , 1994, 309, 249-257. | 1.5 | 32 |
| 9 | 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 |
| 10 | 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 |
| 11 | 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 |
| 12 | 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 |
| 13 | Theoretical Study of the Reactivity and Stereoselectivity of Heterodienophile Addition to 1,3,4-Oxadiazole. <i>Journal of Organic Chemistry</i> , 1994, 59, 3015-3019. | 3.2 | 19 |
| 14 | Semiempirical and ab initio study of 1,3-dipolar addition of azide anion to organic cyanides. <i>Computational and Theoretical Chemistry</i> , 1994, 312, 11-22. | 1.5 | 18 |
| 15 | 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 |
| 16 | Ab initio study of heterodienophile addition to oxazole. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1994, , 1877. | 0.9 | 16 |
| 17 | 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 |
| 18 | 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 |

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|----|--|-----|-----------|
| 19 | 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 |
| 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 | Bioluminescence determination of enzyme activity of firefly luciferase in the presence of pesticides. <i>Luminescence</i> , 2005, 20, 192-196. | 2.9 | 12 |
| 22 | 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 |
| 23 | 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 |
| 24 | 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 |
| 25 | 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 |
| 26 | Theoretical Study of Ethylene and Vinyl Alcohol Addition to 1,4-Dioxo-1,3-butadiene. <i>Journal of Organic Chemistry</i> , 1994, 59, 7732-7736. | 3.2 | 10 |
| 27 | 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 |
| 28 | 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 |
| 29 | 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 |
| 30 | 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 |
| 31 | 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 |
| 32 | 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 |
| 33 | Role of the Polymer Backbone in the Oxidative Properties of Polymer-Supported Dichromate. <i>Synthetic Communications</i> , 1989, 19, 1317-1324. | 2.1 | 5 |
| 34 | On the Formation of Peptide Bonds. <i>Journal of Chemical Education</i> , 1993, 70, 134. | 2.3 | 5 |
| 35 | 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. <i>Journal of Organic Chemistry</i> , 1993, 58, 5245-5250. | 3.2 | 5 |
| 36 | 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 |

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|----|--|-----|-----------|
| 37 | Theoretical study of azide anion addition to nonpolar and polar double and triple bonds. Computational and Theoretical Chemistry, 1995, 343, 149-155. | 1.5 | 3 |
| 38 | Tetrafluoroberyllate(2-) ions as Hydrogen-Bond Proton Acceptors: Spectroscopic Evidence. Spectroscopy Letters, 1996, 29, 867-875. | 1.0 | 2 |
| 39 | The theoretical study of a borane catalyzed azide anion addition to fluorinated acetonitriles. Computational and Theoretical Chemistry, 1997, 389, 69-74. | 1.5 | 2 |
| 40 | Theoretical study of the Diels-Alder reactions of benzene with fluorinated dienophiles. Computational and Theoretical Chemistry, 2004, 684, 99-102. | 1.5 | 2 |
| 41 | Theoretical Study of the Diastereofacial Isomers of Aldrin and Dieldrin. International Journal of Molecular Sciences, 2006, 7, 35-46. | 4.1 | 2 |
| 42 | Using MathCAD in chemistry calculations. Journal of Chemical Education, 1991, 68, A95. | 2.3 | 1 |
| 43 | Calculations of boron trifluoride catalyzed 1,3-dipolar additions of azide ion to organic nitriles. Computational and Theoretical Chemistry, 1995, 332, 127-135. | 1.5 | 1 |
| 44 | Development of a trap for fuel exhaust particulate matter under driving conditions and GC-MS method for their analysis. Macedonian Journal of Chemistry and Chemical Engineering, 2012, 30, 97. | 0.6 | 1 |
| 45 | Preparation and characterization of deuterated crystalhydrates: An undergraduate inorganic or spectroscopic experiment. Journal of Chemical Education, 1993, 70, 603. | 2.3 | 0 |
| 46 | Stopper with Interchangeable Septa for SPME. Journal of Chemical Education, 2009, 86, 924. | 2.3 | 0 |
| 47 | Macedonian journal of chemistry and chemical engineering: open journal systems--editor's perspective. Prilozi - Makedonska Akademija Na Naukite I Umetnostite Oddelenie Za Medicinski Nauki, 2014, 35, 51-5. | 0.5 | 0 |
| 48 | Elementary, my dear Watson! The making of a collection of the natural elements. Macedonian Journal of Chemistry and Chemical Engineering, 2022, 41, . | 0.6 | 0 |