

Venelin Enchev

List of Publications by Year in descending order

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docs citations

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times ranked

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citing authors

#	ARTICLE	IF	CITATIONS
1	»;Potential of hydroxybenzoic acids from <i>Graptopetalum paraguayense</i> for inhibiting of herpes simplex virus DNA polymerase – metabolome profiling, molecular docking and quantum-chemical analysis. <i>Pharmacia</i> , 2022, 69, 113-123.	1.2	1
2	Tautomerism of cytosine, cytidine, and deoxycytidine: Proton transfer through water bridges. <i>International Journal of Quantum Chemistry</i> , 2022, 122, .	2.0	3
3	Chemical evolution: from formamide to nucleobases and amino acids without the presence of catalyst. <i>Journal of Biomolecular Structure and Dynamics</i> , 2021, 39, 5563-5578.	3.5	11
4	Self-catalytic mechanism of prebiotic reactions: II. From urea and glycinate to hypoxanthine. <i>International Journal of Quantum Chemistry</i> , 2021, 121, e26508.	2.0	1
5	Binding Expedient of 2-carbamido-1,3-indandione to Nucleic Acids: Potential Fluorescent Probe. <i>Photochemistry and Photobiology</i> , 2021, 97, 710-717.	2.5	1
6	Self-catalytic mechanism of prebiotic reactions: from formamide to pterins and guanine. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 19043-19053.	2.8	4
7	Effect of external electric field on the tautomeric equilibrium and structure of 2-carbamido-1,3-indandione. <i>International Journal of Quantum Chemistry</i> , 2021, 121, e26760.	2.0	1
8	Self-catalytic mechanism of prebiotic reactions: From formamide to purine bases. <i>International Journal of Quantum Chemistry</i> , 2020, 120, e26362.	2.0	6
9	Anti-Herpes Simplex virus and antibacterial activities of <i>< i>Graptopetalum paraguayense</i></i> E. Walther leaf extract: a pilot study. <i>Biotechnology and Biotechnological Equipment</i> , 2019, 33, 1251-1259.	1.3	6
10	Tautomerism of Inosine in Water: Is It Possible?. <i>Journal of Physical Chemistry B</i> , 2019, 123, 622-630.	2.6	10
11	Hybrid MC/QC simulations of water-assisted proton transfer in nucleosides. Guanosine and its analog acyclovir. <i>Journal of Biomolecular Structure and Dynamics</i> , 2017, 35, 1168-1188.	3.5	6
12	Ultrasound-assisted green bromination of N-cinnamoyl amino acid amides – Structural characterization and antimicrobial evaluation. <i>Journal of Molecular Structure</i> , 2017, 1135, 144-152.	3.6	2
13	Computational insight on the chalcone formation mechanism by the Claisen-Schmidt reaction. <i>International Journal of Quantum Chemistry</i> , 2017, 117, e25365.	2.0	8
14	Green synthesis, structure and fluorescence spectra of new azacyanine dyes. <i>Journal of Molecular Structure</i> , 2017, 1147, 380-387.	3.6	8
15	2-Methylthio-imidazolines: a rare case of different tautomeric forms in solid state and in solution. <i>Structural Chemistry</i> , 2017, 28, 757-772.	2.0	4
16	A Hybrid statistical mechanics–quantum chemical model for proton transfer in 5-azauracil and 6-azauracil in water solution. <i>International Journal of Quantum Chemistry</i> , 2015, 115, 477-485.	2.0	9
17	2-Carbamido-1,3-indandione – a Fluorescent Molecular Probe and Sunscreen Candidate. <i>Journal of Fluorescence</i> , 2015, 25, 1601-1614.	2.5	7
18	Anti-conjunctivitis effect of fresh juice of <i>xGraptovaria</i> (<i>Crassulaceae</i>) - a phytochemical and ethnobotanical study. <i>Journal of Intercultural Ethnopharmacology</i> , 2015, 4, 24.	0.9	3

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19	Solid-state structures of 2-(4-hydroxyphenyl)-substituted phenalene-1,3-dione and indan-1,3-dione. Journal of Structural Chemistry, 2014, 55, 446-455.	1.0	1
20	Synthesis of 3',4'-dihydro-2H,2'H,5H-spiro [imidazolidine-4,1'-naphthalene]-2,5-dione and its derivatives. Acta Chimica Slovenica, 2014, 61, 420-4.	0.6	2
21	A model system with intramolecular hydrogen bonding: Effect of external electric field on the tautomeric conversion and electronic structures. Computational and Theoretical Chemistry, 2013, 1006, 113-122.	2.5	12
22	Excited state proton transfer in 3,6-bis(4,5-dihydroxyoxazo-2-yl)benzene-1,2-diol. Chemical Physics Letters, 2013, 563, 43-49.	2.6	12
23	Antiradical and antioxidant activities of new bio-antioxidants. Biochimie, 2012, 94, 403-415.	2.6	26
24	Tacticity of poly(butyl- \pm -cyanoacrylate) chains in nanoparticles: NMR spectroscopy and DFT calculations. Structural Chemistry, 2012, 23, 815-824.	2.0	9
25	Solid-State Tautomerism in 2-Carboxyindan-1,3-dione. Journal of Physical Chemistry A, 2011, 115, 2026-2034.	2.5	4
26	Ab initio study of the tautomerism of 2,5-substituted diazoles. Structural Chemistry, 2010, 21, 1053-1060.	2.0	7
27	Influence of pH on the cis-“trans isomerization of Valine-Proline dipeptide: An integrated NMR and theoretical investigation. Journal of Molecular Structure, 2010, 975, 330-334.	3.6	9
28	Tautomeric Equilibria of 5-Fluorouracil Anionic Species in Water. Journal of Physical Chemistry A, 2010, 114, 13154-13162.	2.5	49
29	Does tautomeric equilibrium exist in 4-nitroso-5-pyrazolones?. Computational and Theoretical Chemistry, 2009, 897, 55-60.	1.5	14
30	Physicochemical characterization and in vitro behavior of daunorubicin-loaded poly(butylcyanoacrylate) nanoparticles. Acta Biomaterialia, 2009, 5, 2109-2121.	8.3	46
31	Theoretical and Spectroscopic Study of 2-Substituted Indan-1,3-diones: A Coherent Picture of the Tautomeric Equilibrium. Journal of Physical Chemistry A, 2007, 111, 9901-9913.	2.5	9
32	Intramolecular Hydrogen-Bonding Interactions in 2-Nitrosophenol and Nitrosonaphthols: Ab Initio, Density Functional, and Nuclear Magnetic Resonance Theoretical Study. Journal of Physical Chemistry A, 2007, 111, 7112-7123.	2.5	28
33	The chemistry of the processes involved in the production of lanthanide titanates by the polymerized-complex method. Canadian Journal of Chemistry, 2007, 85, 547-559.	1.1	12
34	Ab Initio Investigation on the Second-Order Nonlinear Optical Responses in Keto-Enol Equilibria of Salicylideneanilines. Journal of Physical Chemistry A, 2007, 111, 9914-9923.	2.5	49
35	Synthesis of trans/cis 4-substituted 3-furyl-2-phenethyltetrahydroisoquinolin-1-ones: conformation of the trans-4-(pyrrolidinylcarbonyl) derivative. Tetrahedron Letters, 2006, 47, 2119-2123.	1.4	4
36	Fast intramolecular proton transfer in 2-(hydroxyaminomethylidene)-indan-1,3-dione. Computational and Theoretical Chemistry, 2005, 719, 169-175.	1.5	15

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37	Ab Initio Study of 2,4-Substituted Azolidines. II. Amino α 'Imino Tautomerism of 2-Aminothiazolidine-4-one and 4-Aminothiazolidine-2-one in Water Solution. <i>Journal of Physical Chemistry A</i> , 2005, 109, 8904-8913.	2.5	14
38	Oxo β 'Hydroxy Tautomerism of 5-Fluorouracil: Water-Assisted Proton Transfer. <i>Journal of Physical Chemistry A</i> , 2005, 109, 1981-1988.	2.5	79
39	Experimental and computational studies of the structure and vibrational spectra of 4-dimethylamino pyridinium-betaine of squaric acid. <i>Journal of Molecular Structure</i> , 2004, 691, 241-248.	3.6	23
40	Water-assisted proton transfer in formamide, thioformamide and selenoformamide. <i>Computational and Theoretical Chemistry</i> , 2004, 679, 195-205.	1.5	21
41	Ab initio study of 2,4-substituted azolidines. I. Tautomerism. <i>Computational and Theoretical Chemistry</i> , 2004, 711, 201-207.	1.5	25
42	Ab initio investigation of the structure and nonlinear optical properties of five-membered heterocycles containing sulfur. <i>Chemical Physics</i> , 2004, 298, 29-36.	1.9	36
43	Study on the Role of 5-fluorouracil in the Polymerization of Butylcyanoacrylate during the Formation of Nanoparticles. <i>Journal of Drug Targeting</i> , 2004, 12, 49-56.	4.4	18
44	On the chemical nature of lanthanum α titanium citric complexes, precursors of La ₂ Ti ₂ O ₇ . <i>Materials Letters</i> , 2004, 58, 3559-3563.	2.6	13
45	Tautomeric and conformational equilibrium of 2-nitrosophenol and 9,10-phenanthrenequinonemonooxime: ab initio and NMR study. <i>Computational and Theoretical Chemistry</i> , 2003, 640, 149-162.	1.5	15
46	Reaction of 2-acetyl-indane-1,3-dione with aniline - Schiff base or enamine?. <i>Journal of Molecular Structure</i> , 2003, 654, 11-20.	3.6	10
47	Poly(butylcyanoacrylate) nanoparticles for topical delivery of 5-fluorouracil. <i>International Journal of Pharmaceutics</i> , 2003, 263, 133-140.	5.2	64
48	New Stable Complexes of Au(III) with Biuret: X-ray Structure of cis-[Au(Biu)Br ₂]PPh ₄ and Ab Initio Investigation of cis-[Au(Biu)X ₂] β . <i>Journal of Coordination Chemistry</i> , 2003, 56, 299-305.	2.2	6
49	2-Acetylindan-1,3-dione and its Cu ²⁺ and Zn ²⁺ complexes as promising sunscreen agents. <i>International Journal of Cosmetic Science</i> , 2002, 24, 103-110.	2.6	20
50	Synthesis, cytotoxicity, antibacterial and antitumor activity of platinum(II) complexes of 3-aminocyclohexanespiro-5-hydantoin. <i>Journal of Inorganic Biochemistry</i> , 2002, 89, 203-211.	3.5	28
51	Spectroscopic and quantum chemical study of the structure of a new paramagnetic dimeric palladium(II,III) complex with creatine. <i>Journal of Molecular Structure</i> , 2002, 609, 61-65.	3.6	5
52	Structure of the symmetric monooxime of 1,2,3-indantrione in gas, solution and solid states. <i>Journal of Molecular Structure</i> , 2002, 608, 193-200.	3.6	8
53	Title is missing!. <i>Chemistry of Heterocyclic Compounds</i> , 2002, 38, 1110-1120.	1.2	31
54	Does tautomeric equilibrium exist in ortho-nitrosonaphthols?. <i>Chemical Physics</i> , 2001, 264, 235-244.	1.9	30

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55	Quantum chemical and spectroscopic study of the structure of 2-acetylindan-1,3-dione complexes with metal(II) ions. <i>Journal of Molecular Structure</i> , 2001, 595, 67-76.	3.6	15
56	Intramolecular proton transfer reactions in internally hydrogen-bonded Schiff bases: ab initio and semiempirical study. <i>Computational and Theoretical Chemistry</i> , 2000, 530, 223-235.	1.5	18
57	Tautomeric and conformational equilibrium of acenaphthenequinonemonooxime. <i>Journal of Molecular Structure</i> , 1999, 508, 149-161.	3.6	20
58	Excited state intramolecular proton transfer in 2-acetylindan-1,3-dione. <i>Chemical Physics Letters</i> , 1999, 314, 234-238.	2.6	35
59	Copper(II) Complexes of Spirohydantoins. Synthesis, Quantum-Chemical, and Spectroscopic Study. <i>Structural Chemistry</i> , 1999, 10, 381-385.	2.0	4
60	Ab initio quantum chemical and NMR study of the symmetric monooximes of 1,2,3-phenalenetrione and 1,2,3-indantrione. <i>Journal of Molecular Structure</i> , 1998, 440, 227-235.	3.6	15
61	Structure of six- and seven-membered cyclic- β -diketones and their metal(II) complexes. <i>Polyhedron</i> , 1997, 16, 1693-1699.	2.2	4
62	Tautomerism in 2,2'-bipyridyl-3,3'-diol. <i>International Journal of Quantum Chemistry</i> , 1996, 57, 721-728.	2.0	8
63	Electronic structure and polarizabilities of some heterocycles: I. Hydroxypyrazoles and related systems. <i>Molecular Engineering</i> , 1995, 5, 347-361.	0.2	3
64	Tautomerism of rhodanine. <i>Structural Chemistry</i> , 1994, 5, 225-231.	2.0	22
65	Comparative theoretical study of intramolecular proton transfer in the photochemical cycles of 2-(2-hydroxyphenyl)benzoxazole and 5,8-dimethyl-1-tetralone. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 1994, 80, 135-144.	3.9	3
66	AM1 study of ground state intramolecular proton transfer reaction in 2-(2-hydroxy-5-methylphenyl)benzotriazole and 2-(2-hydroxyphenyl)benzotriazole. <i>Computational and Theoretical Chemistry</i> , 1993, 288, 63-66.	1.5	4
67	A semiempirical and ab initio MO study of the tautomers of N-unsubstituted pyrazolones [hydroxypyrazoles]. <i>Structural Chemistry</i> , 1992, 3, 231-238.	2.0	10
68	Tautomerism of N-unsubstituted pyrazolones (hydroxypyrazoles): MNDO and MNDO + CI study of C-substituted tautomers. <i>Computational and Theoretical Chemistry</i> , 1992, 258, 217-234.	1.5	9
69	Methodology for deriving quantitative structure-retention relationships in gas chromatography. <i>Analytica Chimica Acta</i> , 1992, 260, 69-74.	5.4	23
70	MNDO study of intermediates by anionic polymerization of ethylene oxide. <i>Die Makromolekulare Chemie Rapid Communications</i> , 1990, 11, 423-426.	1.1	0
71	Syntheses and Structures of 2(<i>i</i> O <i>j</i>)Acylhydroxyimino)1,3-indandiones. <i>Liebigs Annalen Der Chemie</i> , 1987, 1987, 375-376.	0.8	4
72	A classification of polyenes into 4L + 2- and 4L-classes on the basis of Coulson's bond orders and information theory and its application to the interpretation of electrocyclic reactions. <i>International Journal of Quantum Chemistry</i> , 1984, 26, 993-1015.	2.0	3

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73	Subchromophore Recognition in Some New Luminophores. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 1984, 39, 1143-1144.		1.5	1
74	A theoreticalâ€”information study on the electron delocalization (aromaticity) of annulenes with and without bond alternation. Computational and Theoretical Chemistry, 1982, 88, 105-118.		1.5	8
75	Electronic Structure and Aromaticity of (AB) _N -Heteroannulenes. Zeitschrift Fur Physikalische Chemie, 1981, 128, 169-178.		2.8	2
76	Some contributions and generalizations to the electronic theory of even polyenes and annulenes. Chemical Physics Letters, 1981, 78, 560-565.		2.6	7
77	Electronic structure of mÃ¶bius annulenes. Chemical Physics Letters, 1981, 83, 529-532.		2.6	11