

Å^{1/2}eljko J Vitnik

List of Publications by Year in descending order

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papers

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citing authors

#	ARTICLE	IF	CITATIONS
1	Antioxidant properties of selected 4-phenyl hydroxycoumarins: Integrated in vitro and computational studies. <i>Chemico-Biological Interactions</i> , 2014, 214, 49-56.	4.0	40
2	Structure and stereochemistry of electrochemically synthesized poly-(l-naphthylamine) from neutral aceto- nitrile solution. <i>Journal of the Serbian Chemical Society</i> , 2002, 67, 867-877.	0.8	35
3	Quantum mechanical and spectroscopic (FT-IR, 13C, 1H NMR and UV) investigations of potent antiepileptic drug 1-(4-chloro-phenyl)-3-phenyl-succinimide. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2014, 117, 42-53.	3.9	27
4	Absorption and fluorescence spectral properties of azo dyes based on 3-amido-6-hydroxy-4-methyl-2-pyridone: Solvent and substituent effects. <i>Dyes and Pigments</i> , 2020, 175, 108139.	3.7	27
5	Synthesis, structure and solvatochromic properties of some novel 5-aryloxy-6-hydroxy-4-phenyl-3-cyano-2-pyridone dyes. <i>Chemistry Central Journal</i> , 2012, 6, 71.	2.6	25
6	On the structures of 5-(4-, 3- and 2-methoxyphenylazo)-3-cyano-1-ethyl-6-hydroxy-4-methyl-2-pyridone: An experimental and theoretical study. <i>Dyes and Pigments</i> , 2014, 104, 160-168.	3.7	25
7	Syntheses, characterization and antimicrobial activity of the first complexes of Zn(II), Cd(II) and Co(II) with N-benzoyloxycarbonylglycine. <i>Inorganica Chimica Acta</i> , 2008, 361, 86-94.	2.4	23
8	The spectroscopic (FT-IR, FT-Raman, 13C, 1H NMR and UV) and NBO analyses of 4-bromo-1-(ethoxycarbonyl)piperidine-4-carboxylic acid. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2015, 138, 1-12.	3.9	23
9	Novel azo pyridone dyes based on dihydropyrimidinone skeleton: Synthesis, DFT study and anticancer activity. <i>Dyes and Pigments</i> , 2021, 187, 109123.	3.7	23
10	Identification and Application of Enantiocomplementary Lactamases for Vince Lactam Derivatives. <i>ChemCatChem</i> , 2014, 6, 2517-2521.	3.7	18
11	New derivatives of hydantoin as potential antiproliferative agents: biological and structural characterization in combination with quantum chemical calculations. <i>Monatshefte für Chemie</i> , 2014, 145, 821-833.	1.8	16
12	Electrochemical characterization of phenytoin and its derivatives on bare gold electrode. <i>Electrochimica Acta</i> , 2015, 161, 378-387.	5.2	14
13	Synthesis, antimicrobial activity and quantum chemical investigation of novel succinimide derivatives. <i>Journal of Molecular Structure</i> , 2019, 1181, 148-156.	3.6	14
14	On the azo dyes derived from benzoic and cinnamic acids used as photosensitizers in dye-sensitized solar cells. <i>Turkish Journal of Chemistry</i> , 2019, 43, 1183-1203.	1.2	12
15	Experimental and theoretical insight into the electronic properties of 4-aryloxy-5-aryloxy-6-hydroxy-2-pyridone dyes. <i>Coloration Technology</i> , 2017, 133, 223-233.	1.5	10
16	Towards understanding intermolecular interactions in hydantoin derivatives: the case of cycloalkane-5-spirohydantoin derivatives tethered with a halogenated benzyl moiety. <i>CrystEngComm</i> , 2017, 19, 469-483.	2.6	10
17	N-H...O Interactions Stabilize the Most-Hindered Rotamer of the (S)-Tyrosinato Side Group in Bis[(S)-tyrosinato](diamine)cobalt(III) Complexes: An NMR Spectroscopic and DFT Study. <i>European Journal of Inorganic Chemistry</i> , 2005, 2005, 3172-3178.	2.0	8
18	Structural, spectroscopic and computational study of 5-(substituted) 1-phenylazo-2-pyridone dyes.	1.5	8

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19	Determination of all pK_a Values of Some di- and Tri-Carboxylic Unsaturated and Epoxy Acids and Their Polylinear Correlation with the Carboxylic Group Atomic Charges. <i>Journal of Chemical Research</i> , 2003, 2003, 247-248.	1.3	7
20	Structure-property relationship of 3-(4-substituted benzyl)-1,3-diazaspiro[4.4]nonane-2,4-diones as new potential anticonvulsant agents. An experimental and theoretical study. <i>Journal of Molecular Structure</i> , 2017, 1127, 88-98.	3.6	7
21	Experimental and theoretical study on the structure-property relationship of novel 1-aryl-3-methylsuccinimides. <i>Journal of Molecular Structure</i> , 2017, 1129, 271-282.	3.6	7
22	Synthesis, structural characterization, DFT calculations and antiproliferative evaluation of novel spirohydantoin derivatives containing a substituted benzyl moiety. <i>Journal of Molecular Structure</i> , 2019, 1180, 48-62.	3.6	7
23	Solvatochromic and quantum chemical investigations of newly synthesized succinimides: substituent effect on intramolecular charge transfer. <i>Monatshefte für Chemie</i> , 2013, 144, 1525-1535.	1.8	6
24	Spectroscopic and quantum mechanical investigation of N,N'-bisarylmalonamides: solvent and structural effects. <i>Journal of Molecular Modeling</i> , 2014, 20, 2384.	1.8	6
25	Solvatochromism and quantum mechanical investigation of disazo pyridone dye. <i>Coloration Technology</i> , 2018, 134, 478-490.	1.5	6
26	Acid-catalyzed rearrangement of some steroidal isoxazolidines. <i>Tetrahedron</i> , 1999, 55, 6681-6690.	1.9	5
27	Substituent and structural effects on the kinetics of the reaction of N-(substituted) Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 427 Chemical Society, 2007, 72, 1191-1200.	0.8	4
28	Improved synthesis and in vitro study of antimicrobial activity of α,β -unsaturated and α -bromo carboxylic acids. <i>Journal of the Serbian Chemical Society</i> , 2012, 77, 741-750.	0.8	3
29	Iodide Analogs of Arsenoplatinsâ€”Potential Drug Candidates for Triple Negative Breast Cancers. <i>Molecules</i> , 2021, 26, 5421.	3.8	3
30	Effect of substituents on the ^{13}C -NMR chemical shifts of 3-methylene-4-substituted-1, 4-pentadienes - Part I. <i>Journal of the Serbian Chemical Society</i> , 2003, 68, 67-76.	0.8	3
31	Kinetics and mechanism of the reaction of substituted 4-pyrimidine carboxylic acids with diazodiphenylmethane in dimethylformamide. <i>Journal of Chemical Research</i> , 2000, 2000, 506-507.	1.3	2
32	Computational and spectroscopic data correlation study of N,N'-bisarylmalonamides (Part II). <i>Journal of Molecular Modeling</i> , 2015, 21, 239.	1.8	2
33	Ring- and side-group conformational properties of di-O-acylated xylopyranosides: A computational study. <i>Computational and Theoretical Chemistry</i> , 2015, 1051, 104-109.	2.5	2
34	Structural and vibrational analyses of new potential anticancer drug 2-(phenylmethyl)-2-azaspiro[5.11]heptadecane-1,3,7-trione. <i>Journal of Molecular Structure</i> , 2017, 1137, 97-108.	3.6	2
35	Charge assisted assembly of zwitterionic pyridone hydrates. <i>Journal of Molecular Structure</i> , 2021, 1237, 130419.	3.6	2
36	Linear free energy relationships of the ^{13}C NMR chemical shifts in 5-(3- and 4-substituted) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 62 Td (0.5	2

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37	Molecular orbital investigation of various reaction pathways in reaction of ketones with bromoform. <i>International Journal of Quantum Chemistry</i> , 2006, 106, 1323-1329.	2.0	1
38	Carbenic vs. ionic mechanistic pathway in reaction of cyclohexanone with bromoform. <i>Journal of Molecular Modeling</i> , 2012, 18, 4721-4728.	1.8	1
39	A detailed UV-Vis spectral investigation of six azo dyes derived from benzoic- and cinnamic acids: experimental and theoretical insight. <i>Comptes Rendus Chimie</i> , 2021, 24, 267-280.	0.5	1
40	Spectroscopic and quantum chemical elucidation of newly synthesized 1-aryl-3-methyl-3-phenylpyrrolidine-2,5-diones as potential anticonvulsant agents. <i>Hemijaska Industrija</i> , 2019, 73, 125-137.	0.7	1
41	Improvement of theoretical UV-Vis spectra calculations by empirical solvatochromic parameters: Case study of 5-aryloxy-3-cyano-1-ethyl-6-hydroxy-4-methyl-2-pyridones. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2022, 272, 120978.	3.9	1
42	Linear free energy relationships of the ¹ H and ¹³ C NMR chemical shifts of 3-methylene-2-substituted-1,4-pentadienes. <i>Journal of Molecular Structure</i> , 2005, 744-747, 901-908.	3.6	0
43	Synthesis, NMR, DFT and antimicrobial studies of Zn(II) complexes with N-benzyloxycarbonyl-S-alanine. <i>Journal of the Serbian Chemical Society</i> , 2008, 73, 815-824.	0.8	0
44	Glutarimides: Biological activity, general synthetic methods and physicochemical properties. <i>Hemijaska Industrija</i> , 2015, 69, 523-536.	0.7	0