

GÃ¼nter A-zpÄ±nar

List of Publications by Year in descending order

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Version: 2024-02-01

15
papers

270
citations

1163117

8
h-index

1058476

14
g-index

15
all docs

15
docs citations

15
times ranked

461
citing authors

#	ARTICLE	IF	CITATIONS
1	An improved generalized AMBER force field (GAFF) for urea. <i>Journal of Molecular Modeling</i> , 2010, 16, 1427-1440.	1.8	77
2	7-Oxy-3-(3,4,5-trimethoxyphenyl)coumarin substituted phthalonitrile derivatives as fluorescent sensors for detection of Fe ³⁺ ions: Experimental and theoretical study. <i>Sensors and Actuators B: Chemical</i> , 2014, 194, 377-388.	7.8	35
3	Intramolecular excimer formation in hexakis(pyrenyloxy)cyclotriphosphazene: photophysical properties, crystal structure, and theoretical investigation. <i>Dalton Transactions</i> , 2014, 43, 3428-3433.	3.3	34
4	QSPR Study on the Bioconcentration Factors of Nonionic Organic Compounds in Fish by Characteristic Root Index and Semiempirical Molecular Descriptors. <i>Journal of Chemical Information and Computer Sciences</i> , 2004, 44, 985-992.	2.8	27
5	Effects of position (\hat{I}_1 or \hat{I}_2) and linker heteroatom (O or S) of substituent on the photophysicochemical behavior of poly(oxyethylene) substituted ZnPcs and assessment of J-aggregation or protonation using TD-DFT computations. <i>Dalton Transactions</i> , 2013, 42, 14892.	3.3	23
6	The effect of α -on/off β -molecular switching on the photophysical and photochemical properties of axially calixarene substituted activatable silicon($\langle scp \rangle iv \langle /scp \rangle$)phthalocyanine photosensitizers. <i>Dalton Transactions</i> , 2016, 45, 7634-7641.	3.3	18
7	Investigation on the aromaticity of 1,3,4-thiadiazole-2-thione and its oxygen analogs including their tautomeric forms. <i>Computational and Theoretical Chemistry</i> , 2005, 726, 233-243.	1.5	17
8	Quantum chemical modeling of the inhibition mechanism of monoamine oxidase by oxazolidinone and analogous heterocyclic compounds. <i>Journal of Enzyme Inhibition and Medicinal Chemistry</i> , 2014, 29, 81-86.	5.2	10
9	A test of improved force field parameters for urea: molecular-dynamics simulations of urea crystals. <i>Journal of Molecular Modeling</i> , 2012, 18, 3455-3466.	1.8	8
10	The Combination of Cross β -Hyperconjugation and γ δ -Conjugation in 2,5 β -Oligosilylanyl Substituted Siloles. <i>Chemistry - A European Journal</i> , 2020, 26, 17252-17260.	3.3	8
11	Reductive Elimination at Pb(II) Center of an (Amino)plumbylene β -Substituted Phosphaketene: New Pathway for Phosphinidene Synthesis. <i>Chemistry - A European Journal</i> , 2022, 28, .	3.3	4
12	Formation of the Vilsmeier-Haack complex: the performance of different levels of theory. <i>Journal of Molecular Modeling</i> , 2011, 17, 3209-3217.	1.8	3
13	Norbornene based-sulfide-stabilized silylium ions: synthesis, structure and application in catalysis. <i>Dalton Transactions</i> , 2022, 51, 1407-1414.	3.3	3
14	NH bond activation of ammonia and amines by ditetrelenes: key insights into the stereochemistry of nucleophilic addition. <i>Dalton Transactions</i> , 2021, 50, 17734-17750.	3.3	2
15	Synthesis of (3,6-dihydro-2H-pyran-2-yl)phosphonate derivatives and investigation of catalyst effect on frontier molecular orbitals using DFT method. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2016, 191, 1262-1267.	1.6	1