Alexander E Pogonin

List of Publications by Year in descending order

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1684188 1474206 19 90 5 9 citations g-index h-index papers 19 19 19 73 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	Molecular structure and bonding in octamethylporphyrin tin(ii), SnN4C28H28. Dalton Transactions, 2012, 41, 7550.	3.3	23
2	Conformation analysis of copper(II) etioporphyrin-II by combined gas electron diffraction/mass-spectrometry methods and DFT calculations. Journal of Molecular Structure, 2015, 1085, 276-285.	3.6	11
3	Vibrational Spectra of Cobalt (II), Nickel(II), Copper(II), Zinc(II) Etioporphyrins-II, MN4C32H36. Macroheterocycles, 2014, 7, 60-72.	0.5	8
4	Combined gas-phase electron diffraction/mass spectrometry and DFT study of the molecular structure of zinc(II) etioporphyrin-II. Structural Chemistry, 2015, 26, 1521-1530.	2.0	8
5	Mass-Spectrometric Study of Cobalt, Nickel, Copper and Zinc Etioporphyrin-II Sublimation. Macroheterocycles, 2012, 5, 315-320.	0.5	8
6	Gas-phase structures of hemiporphyrazine and dicarbahemiporphyrazine: Key role of interactions inside coordination cavity. Journal of Molecular Structure, 2019, 1184, 576-582.	3. 6	5
7	Hydrogen bonds determine the signal arrangement in 13 C NMR spectra of nicotinate. Journal of Molecular Structure, 2018, 1154, 565-569.	3.6	4
8	Quantum Chemical Study Aimed at Modeling Efficient Aza-BODIPY NIR Dyes: Molecular and Electronic Structure, Absorption, and Emission Spectra. Molecules, 2020, 25, 5361.	3.8	4
9	Molecular structure of cobalt(II) etioporphyrin-II determined by combined gas-phase electron diffraction/mass-spectrometry and quantum chemical calculations: Searching a ruffling and saddling effects. Journal of Molecular Structure, 2020, 1216, 128319.	3.6	4
10	CONFORMATIONAL BEHAVIOR OF HYDRAZONE DERIVED FROM PYRIDOXAL 5'-PHOSPHATE AND ISONIAZID. ChemChemTech, 2018, 61, 101-107.	0.3	4
11	Molecular Structure of Nickel Octamethylporphyrinâ€"Rare Experimental Evidence of a Ruffling Effect in Gas Phase. International Journal of Molecular Sciences, 2022, 23, 320.	4.1	4
12	Structure and energetical properties of metal pivalate chelates M(piv)3 (MÂ=ÂAl, Ga, In, Tl) by DFT calculations. Structural Chemistry, 2015, 26, 1443-1450.	2.0	2
13	An alternative approach to structural analysis for gas electron diffraction method. Computational and Theoretical Chemistry, 2018, 1123, 149-153.	2.5	2
14	Molecular structure of 5,10,15,20-tetrakis(4′-fluorophenyl)porphyrin by combined gas-phase electron diffraction/mass spectrometry experiment and DFT calculations. Journal of Molecular Structure, 2020, 1221, 128662.	3.6	2
15	Molecular and electronic structure of substituted BODIPY dyes: Quantum chemical study. Computational and Theoretical Chemistry, 2022, 1212, 113719.	2.5	1
16	Energy of Solvation and a Quantum-Chemical Model of the Structure of 18-Crown-6 Ether in Nonaqueous Solvents. Russian Journal of Physical Chemistry A, 2018, 92, 1494-1498.	0.6	0
17	Intramolecular Hydrogen Bonding and Electronic Structure of Thiadiazole Annulated Hemihexaphyrazine. Macroheterocycles, 2018, 11, 67-72.	0.5	0
18	MOLECULAR STRUCTURE AND VIBRATION SPECTRA OF PIVALIC ACID. ChemChemTech, 2018, 59, 17.	0.3	0

#	Article	IF	CITATIONS
19	DFT STUDY OF MOLECULAR STRUCTURE OF 5,10,15,20-TETRAKIS(4'-HALOGENOPHENYL)PORPHYRINS AND THEIR ISOMERS. ChemChemTech, 2019, 63, 51-57.	0.3	0