

Nasrin Masnabadi

List of Publications by Year in descending order

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Version: 2024-02-01

18
papers

70
citations

1937685

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all docs

18
docs citations

18
times ranked

102
citing authors

#	ARTICLE	IF	CITATIONS
1	Oxidative dimerization of thiols to disulfide using recyclable magnetic nanoparticles. <i>Research on Chemical Intermediates</i> , 2017, 43, 1609-1618.	2.7	11
2	Volatile Constituents of <i>Alococarpum erianthum</i> (DC) H. Riedl & Kuber. <i>Ferula ovina</i> (Boiss.) Boiss. and <i>Pimpinella affinis</i> Ledeb. Three Umbelliferae Herbs Growing Wild in Iran. <i>Journal of Essential Oil Research</i> , 2008, 20, 232-235.	2.7	10
3	Chemical Composition of the Essential Oil from Flower, Stem and Leaves of <i>Astragalus schahrudensis</i> Bge. from Iran. <i>Journal of Essential Oil Research</i> , 2007, 19, 269-270.	2.7	6
4	Conformational Behaviors of 2-Substituted Cyclohexanone Oximes: An Ab Initio, Hybrid Dft Study, and NBO Interpretation. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2012, 187, 276-293.	1.6	5
5	Ab Initio Calculations of the Conformational Preferences of 1,3-Oxathiane S-Oxide and its Analogs Containing S and SE Atoms: Evidence for Stereoelectronic Interactions Associated with the Anomeric Effects. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2013, 188, 1053-1063.	1.6	5
6	Potentiality of carbon nanotube to encapsulate some alkylating agent anticancer drugs: a molecular simulation study. <i>Structural Chemistry</i> , 2021, 32, 869-877.	2.0	5
7	Conformational behaviours of 2-substituted cyclohexanones: a complete basis set, hybrid-DFT study and NBO interpretation. <i>Molecular Simulation</i> , 2011, 37, 1207-1220.	2.0	4
8	Hybrid-Density Functional Theory, MO Study, and NBO Interpretation of Conformational Behaviors of 2-Halo-1,3-Dioxanes and Their Dithiane and Diselenane Analogs. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2012, 187, 305-320.	1.6	4
9	PHYTOCHEMICAL ANALYSIS AND ANTIBACTERIAL ACTIVITY OF <i>ERYNGIUM PYRAMIDALE</i> BOISS. & HAUSSKN. <i>Journal of the Chilean Chemical Society</i> , 2021, 66, 5230-5236.	1.2	4
10	Composition of the Essential Oils of <i>Bupleurum falcatum</i> L. and <i>Bupleurum gerardi</i> All. from Iran. <i>Journal of Essential Oil-bearing Plants: JEOP</i> , 2010, 13, 727-731.	1.9	3
11	A Computational Study of the Conformational Behavior of 2,5-Dimethyl-1,4-dithiane-2,5-diol and Analogous S and Se: DFT and NBO Study. <i>Letters in Organic Chemistry</i> , 2020, 17, 749-759.	0.5	3
12	Structural, Electronic, Reactivity, and Conformational Features of 2,5,5-Trimethyl-1,3,2-diheterophosphinane-2-sulfide, and Its Derivatives: DFT, MEP, and NBO Calculations. <i>Molecules</i> , 2022, 27, 4011.	3.8	3
13	Conformational Stability, FMO, NMR, MEP and NBO Analysis of 2,5-Di-methyl-2,5-bis(methylthio)-1,4-dithiane and Dimethoxy compounds by DFT Approach. <i>Letters in Organic Chemistry</i> , 2021, 17, .	0.5	2
14	Functionalized GO@ZIF-90-supported sulfuric acid and its application in the catalytic synthesis of sulfonamides. <i>Research on Chemical Intermediates</i> , 2022, 48, 457-470.	2.7	2
15	Theoretical study of encapsulation of diethylstilbestrol drug into the inner surface of BNNT toward designing a new nanocarrier for drug delivery systems. <i>Materials Research Express</i> , 2022, 9, 045002.	1.6	2
16	Variational RRKM theory calculation of thermal rate constant for carbon-hydrogen bond fission reaction of nitro benzene. <i>Russian Journal of Physical Chemistry A</i> , 2013, 87, 1175-1179.	0.6	1
17	Calculation of Arrhenius Parameters by the Rrkm Method for the Carbon-Hydrogen Bond Fission Reaction of Fluorobenzene. <i>Progress in Reaction Kinetics and Mechanism</i> , 2017, 42, 182-190.	2.1	0
18	The study of Letrozole adsorption upon CCT nanotube: A DFT/TD-DFT and spectroscopic (excited states) Tj ETQq0 0,0 rgBT /Qverlock 10	0.8	0