

Saeed Jamehbozorgi

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

59
papers

493
citations

12
h-index

18
g-index

65
ext. papers

609
ext. citations

2.1
avg, IF

4.25
L-index

#	Paper	IF	Citations
59	Understanding delivery and adsorption of Flutamide drug with ZnONS based on: Dispersion-corrected DFT calculations and MD simulations. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2022 , 135, 114937	3	5
58	Hydrothermal synthesis of a dinuclear copper(II) complex with 2-aminopyrimidine and bridging oxalate ligands: in situ degradation of tartaric acid to oxalate and thermal decomposition to CuO nanoparticles. <i>Inorganic and Nano-Metal Chemistry</i> , 2020 , 50, 1353-1357	1.2	0
57	Selective and Sensitive Two New Macroacyclic Schiff base Fluorescent Turn-Off Receptors for Fe ³⁺ , DFT Calculation and Their Antibacterial Activity. <i>Applied Organometallic Chemistry</i> , 2020 , 34, e5430	3.1	2
56	Synthesis of 1-amidoalkyl-2-naphthol derivatives using a magnetic nano-Fe ₃ O ₄ @SiO ₂ @Hexamethylenetetramine-supported ionic liquid as a catalyst under solvent-free conditions. <i>Journal of the Chinese Chemical Society</i> , 2020 , 67, 603-609	1.5	7
55	The effect of encapsulation of lithium atom on supramolecular triad complexes performance in solar cell by using theoretical approach. <i>Adsorption</i> , 2020 , 26, 471-489	2.6	1
54	Graphyne and its boron nitride analogue as nanocarriers for anti-cancer drug delivery. <i>Molecular Physics</i> , 2020 , 118, e1691748	1.7	6
53	A new type of heteroborospherene as a versatile carrier for drug delivery: A theoretical study. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2020 , 117, 113852	3	10
52	Preparation of CoFe ₂ O ₄ /sawdust and NiFe ₂ O ₄ /sawdust magnetic nanocomposites for removal of oil from the water surface. <i>Journal of the Chinese Chemical Society</i> , 2020 , 67, 288-297	1.5	3
51	The interaction of 5-fluorouracil with graphene in presence of external electric field: a theoretical investigation. <i>Adsorption</i> , 2020 , 26, 905-911	2.6	4
50	Theoretical Study of Substituent Effect on the Electronic and Optical Properties of 4-Substituted Ferrocenylethynylbenzenes. <i>Russian Journal of Physical Chemistry A</i> , 2019 , 93, 1747-1753	0.7	2
49	Adsorption of Cd and Ni from water by graphene oxide and graphene oxide-almond shell composite. <i>Water Environment Research</i> , 2019 , 91, 475-482	2.8	9
48	Tandem dispersive liquid-liquid microextraction coupled with micro-sampling flame atomic absorption spectrometry for rapid determination of lead(II) and cadmium(II) ions in environmental water samples. <i>International Journal of Environmental Analytical Chemistry</i> , 2019 , 99, 1235-1246	1.8	6
47	Enhancing Photocatalytic Activity of Cu ₂ O in Degradation of Sulphonic Acid-Based Dye. <i>Russian Journal of Applied Chemistry</i> , 2019 , 92, 141-149	0.8	4
46	Effect of the External Electric Field on the Electronic Structure and Aromaticity of Iridabenzene: A DFT Study. <i>Journal of Structural Chemistry</i> , 2019 , 60, 547-555	0.9	5
45	The Role of Polyethylene Glycol Size in Chemical Spectra, Cytotoxicity, and Release of PEGylated Nanoliposomal Cisplatin. <i>Assay and Drug Development Technologies</i> , 2019 , 17, 231-239	2.1	13
44	Microfunnel-filter-based emulsification microextraction followed by gas chromatography for simple determination of organophosphorus pesticides in environmental water samples. <i>Journal of Separation Science</i> , 2019 , 42, 2418-2425	3.4	9
43	The Role of Pseudo Jahn-Teller Effect in Geometry and Electronic Parameters of Tetrahalodigermene Ge ₂ X ₄ (X= Cl, Br, I). <i>Russian Journal of Physical Chemistry A</i> , 2019 , 93, 1297-1304	0.7	1

42	Synthesis of Tetrazoles Catalyzed by Novel Cobalt Magnetic Nanoparticles. <i>Russian Journal of Organic Chemistry</i> , 2019 , 55, 1777-1784	0.7	1
41	Synthesis and characterization of a new poly(N-heterocyclic carbene Cu complex) immobilized on nano-silica, (CuII-NHCs) _n @nSiO ₂ , and its application as an efficient and reusable catalyst in the synthesis of benzimidazoles, benzothiazoles, 1,2,3-triazoles, bis-triazoles and Sonogashira-Bagchi reactions. <i>Inorganica Chimica Acta</i> , 2019 , 485, 173-189	2.7	11
40	The effect of water on the electronic and field emission properties of inorganic AlN nanocones: Computational study. <i>Inorganic Chemistry Communication</i> , 2018 , 90, 86-91	3.1	9
39	DFT/TD-semiempirical study on the structural and electronic properties and absorption spectra of supramolecular fullerene-porphyrine-metalloporphyrine triads based dye-sensitized solar cells. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2018 , 194, 57-66	4.4	12
38	Preparation of nanostructured Co/Mo alloy electrodes and investigation of their electrocatalytic activity for hydrazine oxidation in alkaline medium. <i>Journal of the Chinese Chemical Society</i> , 2018 , 65, 970-976	1.5	4
37	Solvent Effects on the Structure And Spectroscopic Properties of the Second-Generation Anticancer Drug Carboplatin: A Theoretical Insight. <i>Journal of Structural Chemistry</i> , 2018 , 59, 245-251	0.9	18
36	Influence of Solvent and Electric Field on the Structure and IR, 31P NMR Spectroscopic Properties of a Titanocene-Benzyne Complex. <i>Journal of Applied Spectroscopy</i> , 2018 , 85, 526-534	0.7	13
35	A Computational Approach for Hydrolysis of the Third-Generation Anticancer Drug: Trans-Platinum(II) Complex of 3-Aminoflavone. <i>Journal of Structural Chemistry</i> , 2018 , 59, 1791-1796	0.9	3
34	Analysis of the Interaction Between the C ₂₀ Cage and cis-PtCl ₂ (NH ₃) ₂ : A DFT Investigation of the Solvent Effect, Structures, Properties, and Topologies. <i>Journal of Structural Chemistry</i> , 2018 , 59, 1044-1051	0.9	4
33	Quantum Chemical Study of the Solvent Effect on the Anticancer Active Molecule of Iproplatin: Structural, Electronic, and Spectroscopic Properties (IR, 1H NMR, UV). <i>Journal of Applied Spectroscopy</i> , 2017 , 83, 909-916	0.7	19
32	Theoretical Study of Substituent Effects on Geometric and Spectroscopic Parameters (IR, 13C, 29Si NMR) and Energy Decomposition Analysis of the Bonding in Molybdenum Silylidyne Complexes CpMo(CO) ₂ (Si-para-C ₆ H ₄ X). <i>Journal of the Chinese Chemical Society</i> , 2017 , 64, 522-530	1.5	14
31	Warped C ₈₀ H ₃₀ nanographene as a chemical sensor for CO gas: DFT studies. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2017 , 381, 646-651	2.3	25
30	Assessing of Removal Efficiency of Indigo Carmine from Wastewater Using MWCNTs 2017 , 41, 1047-1053		8
29	Facile, low-cost, and organic-free fabrication of diverse nanoporous alumina as support for drug release; on the salt effect, calcination temperature, and reaction time dependence. <i>Journal of Sol-Gel Science and Technology</i> , 2017 , 83, 627-639	2.3	1
28	Solvent effect on the linkage isomerism in [Fe(CO) ₄ (NCS)] ⁻ and [Fe(CO) ₄ (SCN)] ⁻ anions: A theoretical investigation. <i>Physics and Chemistry of Liquids</i> , 2017 , 55, 444-456	1.5	20
27	Computational study of osmabenzyne: The solvent effects on the structure and spectroscopic properties (IR, NMR). <i>Journal of Structural Chemistry</i> , 2017 , 58, 1324-1331	0.9	17
26	Structural and electronic properties of metalloporphyrin (MP, M = Fe, Co and Zn) adsorbed on single walled BNNT and SiCNT. <i>Applied Surface Science</i> , 2016 , 360, 69-76	6.7	11
25	Cytotoxicity of Nanoliposomal Cisplatin Coated with Synthesized Methoxypolyethylene Glycol Propionaldehyde in Human Ovarian Cancer Cell Line A2780CP. <i>Tropical Journal of Pharmaceutical Research</i> , 2016 , 15, 563	0.8	2

24	A comparative study of structural and electronic properties of formaldehyde molecule on monolayer honeycomb structures based on vdW-DF prospective. <i>Applied Surface Science</i> , 2016 , 384, 175-181	6.7	18
23	Structural, energetic and electrical properties of encapsulation of penicillamine drug into the CNTs based on vdW-DF perspective. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2015 , 72, 120-127		6
22	Nickel(II) catalyzed oxidation of aldehyde derivatives to their carboxylic acid or ester analogs. <i>Research on Chemical Intermediates</i> , 2013 , 39, 3319-3325	2.8	6
21	Hybrid-DFT Study and NBO Interpretation of the Configurational Behavior of 2-Halotetrahydrothiopyran S-Oxides. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2013 , 188, 839-849	1	0
20	Synthesis, characterization, spectroscopic study, and catalytic behavior of ruthenium(II) o-phenanthroline complex with dipicolinate. <i>Journal of Coordination Chemistry</i> , 2012 , 65, 994-1003	1.6	10
19	Theoretical study on the electronic, structural, properties and reactivity of a series of mono-, di-, tri- and tetrachlorothiophenes as well as corresponding radical cation forms as monomers for conducting polymers. <i>Chemistry Central Journal</i> , 2011 , 5, 13		4
18	Theoretical study on the electronic, structural, properties and reactivity of a series of mono-, di-, tri- and tetrafluorothiophenes as monomers for new conducting polymers. <i>Journal of Fluorine Chemistry</i> , 2011 , 132, 190-195	2.1	7
17	Simultaneous Voltammetric Determination of Lead and Tin by Adsorptive Differential Pulse Stripping Method and Orthogonal Signal Correction-Partial Least Squares in Water Samples. <i>Journal of the Chinese Chemical Society</i> , 2008 , 55, 276-285	1.5	4
16	Electrochemical synthesis of 4-(dihydroxyphenylthio)-2H-chromen-2-one derivatives. <i>Chemical and Pharmaceutical Bulletin</i> , 2008 , 56, 1562-6	1.9	8
15	Prediction of toxicity of nitrobenzenes using ab initio and least squares support vector machines. <i>Journal of Hazardous Materials</i> , 2008 , 151, 603-9	12.8	59
14	Quantitative structure-property relationship study of the solubility of thiazolidine-4-carboxylic acid derivatives using ab initio and genetic algorithm-partial least squares. <i>Chinese Chemical Letters</i> , 2007 , 18, 621-624	8.1	4
13	DFT study and NBO analysis of the mutual interconversion of cumulene compounds. <i>Journal of Physical Organic Chemistry</i> , 2007 , 20, 355-364	2.1	20
12	Ab Initio and NBO Studies of the Decomposition ([2+2] and [2+4]Elimination) Mechanisms of Alkylisothiocyanates (alkyl = ethyl-, iso-propyl- and tert-butyl). <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2006 , 181, 75-85	1	6
11	An Ab Initio Study and NBO Analysis of the Stability and Conformational Properties of Hexakis(trimethylelementhyl)benzene (Element = C, Si, Ge, and Sn). <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2006 , 181, 2419-2434	1	4
10	Ab Initio Study of Geminal Steric Hindrance Effects on the Stability of Conformations of Cyclohexane Derivatives. <i>Journal of Chemical Research</i> , 2005 , 2005, 508-515	0.6	5
9	Ab initio study of the conformational properties of (Z,Z)-cyclodeca-1,6-diene and its heterocyclic analogue containing oxygen. <i>Computational and Theoretical Chemistry</i> , 2005 , 716, 211-215		
8	Ab initio study of the structures and dynamic stereochemistry of biaryls. <i>Computational and Theoretical Chemistry</i> , 2005 , 717, 41-51		38
7	Ab Initio Study of the Various Pathways of the Decomposition ([2+2]elimination) of 2-Chloroethyltrichlorosilane. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2005 , 180, 1611-1619		19

6	Synthesis, Characterization, and an ab initio Study of a Manganese(II) Macrocyclic Schiff-Base Complex with Two 2-Aminoethyl Pendant Arms. <i>Transition Metal Chemistry</i> , 2004 , 29, 523-527	2.1	2
5	Ab initio Study of Configurational and Conformational Properties of Cyclodeca-1,2,4,6,7,9-Hexaene and Cyclodeca-1,2,4,6,8,9-Hexaene. <i>Journal of Chemical Research</i> , 2003 , 2003, 384-385	0.6	2
4	Ab Initio Study of Conformational Properties of (Z,Z,Z)-Cyclonona-1,3,6-Triene. <i>Journal of Chemical Research</i> , 2003 , 2003, 546-548	0.6	
3	Ab Initio Study of Ring Flipping of the Overcrowded Peri-Substituted Naphthalenes. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2003 , 178, 2529-2537	1	3
2	Ab Initio Study of Structures, Metallootropic 1,2-Shifts and Prototropic 1,2-Shifts of Cyclopentadienyl(trimethyl)silane, -germane and -stannane. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 2003 , 178, 341-351	1	5
1	Ab initio Study of Configurations of Cycloundeca-1,2,4,5,7,8,10-heptaene. <i>Journal of Chemical Research</i> , 2002 , 2002, 544-546	0.6	