

Brahim Houari

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Interactions in inclusion complex of β -cyclodextrin/I-Methionine: DFT computational studies. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 2020, 96, 43-54.	1.6	19
2	A quantum chemical study of encapsulation and stabilization of gallic acid in β -cyclodextrin as a drug delivery system. Canadian Journal of Chemistry, 2020, 98, 204-214.	1.1	8
3	Computational studies of Ni(II) photosensitizers complexes containing 1,1'-bis(diphenylphosphino)ferrocene and dithio ligands. Canadian Journal of Chemistry, 2020, 98, 194-203.	1.1	4
4	Theoretical study of geometric, optical, nonlinear optical, UV-Vis spectra and phosphorescence properties of iridium(III) complexes based on 5-nitro-2-(2,4-difluorophenyl)pyridyl. Theoretical Chemistry Accounts, 2020, 139, 1.	1.4	19
5	Investigation of intermolecular interactions in inclusion complexes of pyroquilon with cucurbit[n]urils (n=7,8) using DFT-D3 correction dispersion. Journal of Molecular Liquids, 2020, 309, 113233.	4.9	23
6	Theoretical investigation on green emitting heteroleptic cyclometalated iridium(III) complexes with fluorinated 2-phenylpyridine ligands. Journal of Photochemistry and Photobiology A: Chemistry, 2020, 398, 112624.	3.9	5
7	DFT/TD-DFT investigation on the UV-Vis absorption and phosphorescence spectra of platinum(II) and palladium(II) complexes with Schiff-base ligands. Journal of Luminescence, 2019, 210, 96-103.	3.1	13
8	Structural, optical and nonlinear optical properties and TD-DFT analysis of heteroleptic bis-cyclometalated iridium(III) complex containing 2-phenylpyridine and picolinate ligands. Theoretical Chemistry Accounts, 2018, 137, 1.	1.4	38
9	Structure, electronic properties, and NBO and TD-DFT analyses of nickel(II), zinc(II), and palladium(II) complexes based on Schiff-base ligands. Journal of Molecular Modeling, 2018, 24, 301.	1.8	10
10	Effects of C(2) Methylation on Thermal Behavior and Interionic Interactions in Imidazolium-Based Ionic Liquids with Highly Symmetric Anions. Applied Sciences (Switzerland), 2018, 8, 1043.	2.5	24
11	Theoretical study of geometric structures and electronic absorption spectra of Iridium(III) complexes based on 2-phenyl-5-nitropyridyl with different ancillary ligands. Computational and Theoretical Chemistry, 2017, 1101, 8-19.	2.5	9
12	Quantitative conformational stability host-guest complex of Carvacrol and Thymol with β -cyclodextrin: a theoretical investigation. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 2017, 89, 143-155.	1.6	10
13	DFT/TDDFT computational study of the structural, electronic and optical properties of rhodium (III) and iridium (III) complexes based on tris-picolinate bidentate ligands. Journal of Molecular Modeling, 2017, 23, 344.	1.8	14
14	Structural and spectroscopic properties of Ir(III) complexes with phenylpyridine ligands: Absorption spectra without and with spin-orbit-coupling. Computational and Theoretical Chemistry, 2014, 1040-1041, 219-229.	2.5	31
15	Spin-orbit absorption spectroscopy of transition metal hydrides: A TD-DFT and MS-CASPT2 study of HM(CO) ₅ (M = Mn, Re). International Journal of Quantum Chemistry, 2012, 112, 2085-2097.	2.0	22