Ademir J Camargo

List of Publications by Year
in descending order

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68	900	535685 17	563245 28
papers	citations	h-index	g-index

68	68	68	1309
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	An Update on the Synthesis and Pharmacological Properties of Pyrazoles Obtained from Chalcone. Current Organic Chemistry, 2022, 26, 81-90.	0.9	4
2	Cyclohexanone-Based Chalcones as Alternatives for Fuel Additives. ACS Omega, 2022, 7, 11871-11886.	1.6	6
3	Molecular modeling and nonlinear optical properties of new isostructural halogenated dihydroquinolinones. New Journal of Chemistry, 2022, 46, 14192-14204.	1.4	2
4	Antioxidant effects of polyphenolic compounds and structure-activity relationship predicted by multivariate regression tree. LWT - Food Science and Technology, 2021, 137, 110366.	2.5	20
5	A new isostructural halogenated chalcone with optical properties. Journal of Molecular Modeling, 2021, 27, 52.	0.8	3
6	Effect of ortho- and para-chlorine substitution on hydroxychlorochalcone. Journal of Molecular Modeling, 2021, 27, 65.	0.8	6
7	Synthesis and Structural Studies of Two New Anthracene Derivatives. Crystals, 2021, 11, 934.	1.0	1
8	Aqueous solvation study of melatonin using ab initio molecular dynamics. Journal of Molecular Liquids, 2021, 343, 117451.	2.3	4
9	Ab Initio Molecular Dynamics Simulations of Aqueous Glucosamine Solutions: Solvation Structure and Mechanism of Proton Transfer from Water to Amino Group. Journal of Physical Chemistry B, 2020, 124, 6986-6997.	1.2	6
10	Halogen bonds on substituted dibromonitrobenzene derivatives. Journal of Molecular Modeling, 2020, 26, 319.	0.8	0
11	Synthesis, characterization, and computational study of a new heteroaryl chalcone. Journal of Molecular Modeling, 2020, 26, 243.	0.8	2
12	Synthesis and structural studies on (E)-3-(2,6-difluorophenyl)-1-(4-fluorophenyl)prop-2-en-1-one: a promising nonlinear optical material. RSC Advances, 2020, 10, 22542-22555.	1.7	15
13	Structural studies on dihydropyrimidine derivatives as Mycobacterium tuberculosis coenzyme-A carboxylase inhibitors. Zeitschrift Fur Kristallographie - Crystalline Materials, 2019, 234, 657-669.	0.4	3
14	Structureâ€“activity relationship of tacrine and its analogues in relation to inhibitory activity against Alzheimerâ€™s disease. Journal of Molecular Modeling, 2019, 25, 116.	0.8	2
15	Molecular modeling of cytotoxic activity of a new terpenoid-like bischalcone. New Journal of Chemistry, 2019, 43, 18451-18460.	1.4	6
16	A Comprehensive Topological Analysis of a Novel Flavonoid Extracted from Brazilian Cerrado Plants. ChemistrySelect, 2019, 4, 14012-14020.	0.7	3
17	Theoretical Studies on the Glucosamine: A Systematic Review. Revista Virtual De Quimica, 2019, 11, 1835-1852.	0.1	0
18	Uma RevisÃ£o SistemÃ¡tica sobre InteraÃ§Ãµes de HalogÃªnio em Derivados de Nitrobenzeno SubstituÃdos. Revista Processos QuÃmicos, 2019, 13, 23-30.	0.0	0

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19	Explicit Aqueous Solvation Treatment of Epinephrine from Carâ€“Parrinello Molecular Dynamics: Effect of Hydrogen Bonding on the Electronic Absorption Spectrum. Journal of Physical Chemistry B, 2018, 122, 8439-8450.	1.2	7
20	Synthesis, characterization, and computational study of the supramolecular arrangement of a novel cinnamic acid derivative. Journal of Molecular Modeling, 2017, 23, 35.	0.8	7
21	Substitution effect on a hydroxylated chalcone: Conformational, topological and theoretical studies. Journal of Molecular Structure, 2017, 1136, 69-79.	1.8	16
22	Conformation analysis of a novel fluorinated chalcone. Journal of Molecular Modeling, 2017, 23, 97.	0.8	6
23	A novel dihydrocoumarin under experimental and theoretical characterization. Journal of Molecular Modeling, 2017, 23, 315.	0.8	5
24	Contribution of Directional Dihydrogen Interactions in the Supramolecular Assembly of Single Crystals: Quantum Chemical and Structural Investigation of C₁₇H₁₇N₃O₂ Azine. Crystal Growth and Design, 2017, 17, 5145-5153.	1.4	22
25	Estudo TeÃ³rico dos ParÂ¢metros Estruturais da CafeÃna no VÃ¡cuo Usando DinÂ¢mica Molecular de Car-Parrinello. Revista Processos QuÃmicos, 2017, 11, 17-24.	0.0	0
26	Structural and Theoretical Investigation of Anhydrous 3,4,5-Triacetoxybenzoic Acid. PLoS ONE, 2016, 11, e0158029.	1.1	3
27	Stereodirectional Origin of anti-Arrhenius Kinetics for a Tetraatomic Hydrogen Exchange Reaction: Bornâ€“Oppenheimer Molecular Dynamics for OH + HBr. Journal of Physical Chemistry A, 2016, 120, 5408-5417.	1.1	30
28	Methanol Solvation Effect on the Proton Rearrangement of Curcuminâ€™s Enol Forms: An Ab Initio Molecular Dynamics and Electronic Structure Viewpoint. Journal of Physical Chemistry C, 2016, 120, 19923-19931.	1.5	27
29	A Quantum Chemical and Chemometrical Study of Styrylbenzylsulfones and their Analogues with Citotoxic Activity against Prostate Cancer Cells. Revista Virtual De Quimica, 2016, 8, 506-514.	0.1	0
30	Stereodynamical Origin of Anti-Arrhenius Kinetics: Negative Activation Energy and Roaming for a Four-Atom Reaction. Journal of Physical Chemistry Letters, 2015, 6, 1553-1558.	2.1	63
31	Synthesis, characterization, and computational study of a new dimethoxy-chalcone. Journal of Molecular Modeling, 2014, 20, 2526.	0.8	42
32	Effect of the Methanol Molecule on the Stabilization of C₁₈H₁₈O₄ Crystal: Combined Theoretical and Structural Investigation. Journal of Physical Chemistry A, 2014, 118, 10048-10056.	1.1	5
33	Theoretical investigation on ruthenium tetraazaporphyrin as potential nitric oxide carrier in biological systems. Journal of Molecular Modeling, 2013, 19, 1727-1737.	0.8	1
34	Conformational variability in a new terpenoid-like bischalcone: Structure and theoretical studies. Journal of Structural Chemistry, 2013, 54, 1112-1121.	0.3	4
35	Biological and structure-activity evaluation of chalcone derivatives against bacteria and fungi. Journal of the Brazilian Chemical Society, 2013, 24, 133-144.	0.6	29
36	Structure-activity relationship study of rutaecarpine analogous active against central nervous system cancer. Journal of the Brazilian Chemical Society, 2012, 23, 2183-2190.	0.6	5

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37	X-ray diffraction and theoretical investigation of the Gedunin crystal structure. Journal of Molecular Structure, 2012, 1008, 83-87.	1.8	6
38	Theoretical investigation of nitric oxide interaction with aluminum phthalocyanine. Journal of Molecular Graphics and Modelling, 2011, 29, 777-783.	1.3	9
39	Theoretical investigation of the interaction of glycerol with aluminum and magnesium phthalocyanines. Journal of Molecular Graphics and Modelling, 2010, 29, 206-213.	1.3	7
40	A quantum chemical and chemometrical study of indolo[2,1-b]quinazoline and their analogues with cytotoxic activity against breast cancer cells. SAR and QSAR in Environmental Research, 2009, 20, 537-549.	1.0	7
41	AproximaÃ§Ãµes da MecÃ¢nica QuÃ¢ntica no Estudo de Propriedades Moleculares. Revista Processos QuÃmicos, 2009, 3, 9-16.	0.0	2
42	DinÃ¢mica Molecular de Car-Parrinello. Revista Processos QuÃmicos, 2009, 3, 59-72.	0.0	0
43	Estudo TeÃ³rico da RelaÃ§Ã£o Estrutura Atividade da Indolo [2,1b] Quinazolina e seus Derivados AnÃ¡logos Contra o CÃ¢ncer de OvÃ¡rio. Revista Processos QuÃmicos, 2009, 3, 24-30.	0.0	0
44	Estudo QuÃmico QuÃ¢ntico da AdsorÃ§Ã£o dos Gases O2 e H2 sobre a Ftalocianina de AlumÃnio. Revista Processos QuÃmicos, 2008, 2, 23-30.	0.0	1
45	Estudo QuÃmico QuÃ¢ntico da Atividade da Indolo [2,1b] Quinazolina e seus Derivados AnÃ¡logos Contra o CÃ¢ncer de Mama. Revista Processos QuÃmicos, 2008, 2, 51-61.	0.0	0
46	Theoretical investigation of the intramolecular hydrogen bond formation, non-linear optic properties, and electronic absorption spectra of the 8-hydroxiquinoline. Computational and Theoretical Chemistry, 2007, 816, 145-151.	1.5	46
47	Estudo teÃ³rico quÃmico-quÃ¢ntico das propriedades geomÃ©tricas e fÃsico-quÃmicas das Ftalocianinas de Co, Cr, Cu, Mn, Ni, Fe, Sc, Ti, VO. Revista Processos QuÃmicos, 2007, 1, 21-34.	0.0	0
48	AnÃ¡lise da difraÃ§Ã£o dos Raios X. Revista Processos QuÃmicos, 2007, 1, 35-45.	0.0	1
49	AnÃ¡lise Quantitativa do Solvente em Cristais. Revista Processos QuÃmicos, 2007, 1, 42-50.	0.0	0
50	Structural insights of a potential inhibition against leishmania major. Acta Crystallographica Section A: Foundations and Advances, 2005, 61, c290-c290.	0.3	0
51	Ruthenium Tetraammines as a Model of Nitric Oxide Donor Compounds. European Journal of Inorganic Chemistry, 2004, 2004, 1879-1885.	1.0	48
52	Study of the Oâ€“Ruâ€“N bonding in trans-[Ru(NH3)4(SO4)L]+ complexes (L=imidazole, histidine and) Tj ETQq0 0 0 rgBT /Overlock 10 T 2004, 357, 3147-3154.	1.2	16
53	A quantum chemical and photophysical study of acridine-9-N-methacrylamide. Computational and Theoretical Chemistry, 2004, 674, 213-225.	1.5	15
54	Reactivity of Radicals Generated on Irradiation oftrans-[Ru(NH3)4(NO2)P(OEt)3](PF6). Journal of the American Chemical Society, 2004, 126, 2546-2555.	6.6	25

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55	A quantum chemical and statistical study of flavonoid compounds (flavones) with anti-HIV activity. European Journal of Medicinal Chemistry, 2003, 38, 929-938.	2.6	22
56	Structure characterization of molecular complexes for non-linear optical materials I. X-ray analysis and AM1 calculations of 1 : 1 complexes of 8-hydroxiquinoline (1) and isonicotinamide (2) with 2,4,6-trinitrophenol. Zeitschrift Fur Kristallographie - Crystalline Materials, 2003, 218, 575-580.	0.4	12
57	A study of neolignan compounds with biological activity against Paracoccidioides brasiliensis by using quantum chemical and chemometric methods. Journal of the Brazilian Chemical Society, 2003, 14, 809-814.	0.6	19
58	A Structure-Activity Relationship (SAR) Study of Neolignan Compounds with Anti-schistosomiasis Activity. Journal of the Brazilian Chemical Society, 2002, 13, 300-307.	0.6	18
59	Synthesis, structure, electronic and vibrational spectra of 9-(Diethylamino)-benzo(a)phenoxazin-7-ium-5-N-methacrylamide. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2002, 58, 3103-3111.	2.0	15
60	A structureâ€“activity relationship (SAR) study of synthetic neolignans and related compounds with biological activity against Escherichia coli. Computational and Theoretical Chemistry, 2002, 583, 105-116.	1.5	19
61	Molecular orbital calculations, experimental and theoretical UV spectra of granulatimides and didemnimides, biologically active polycyclic heteroaromatic alkaloids from the ascidian Didemnum granulatum. Journal of Molecular Structure, 2001, 559, 67-77.	1.8	6
62	On-line mass spectrometry investigation of the reduction of carbon dioxide in acidic media on polycrystalline Pt. Electrochemistry Communications, 2001, 3, 603-607.	2.3	40
63	A multiple linear regression and partial least squares study of flavonoid compounds with anti-HIV activity. Computational and Theoretical Chemistry, 2001, 541, 81-88.	1.5	24
64	A structureâ€“activity relationship study of HEPT-analog compounds with anti-HIV activity. Computational and Theoretical Chemistry, 2000, 530, 39-47.	1.5	26
65	Title is missing!. Journal of Solution Chemistry, 2000, 29, 1047-1060.	0.6	126
66	A quantum chemical and statistical study of flavonoid compounds with anti-HIV activity. Computational and Theoretical Chemistry, 1999, 491, 123-131.	1.5	24
67	Crystal, Molecular, and Electronic Structure of 1-Acetyl-indoline and Derivatives. Structural Chemistry, 1998, 9, 365-373.	1.0	3
68	Synthesis, Antimicrobial Activity and Structure-Activity Relationship of Some 5-Arylidene-thiazolidine-2,4-dione Derivatives. Journal of the Brazilian Chemical Society, 0, , .	0.6	8