Marcelo Zaldini Hernandes

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Streptomyces hygroscopicus UFPEDA 3370: A valuable source of the potent cytotoxic agent nigericin and its evaluation against human colorectal cancer cells. Chemico-Biological Interactions, 2021, 333, 109316.	1.7	2
2	NPCdc, a synthetic natriuretic peptide, is a substrate to neprilysin and enhances blood pressure-lowering induced by enalapril in 5/6 nephrectomized rats. Toxicon, 2021, 203, 30-39.	0.8	2
3	Chagas disease: Immunology of the disease at a glance. Cytokine and Growth Factor Reviews, 2021, 62, 15-22.	3.2	18
4	Immunogenicity of Potential CD4+ and CD8+ T Cell Epitopes Derived From the Proteome of Leishmania braziliensis. Frontiers in Immunology, 2020, 10, 3145.	2.2	4
5	Chagas Disease Treatment and Rational Drug Discovery: A Challenge That Remains. Frontiers in Pharmacology, 2019, 10, 873.	1.6	9
6	A docking-based structural analysis of geldanamycin-derived inhibitor binding to human or Leishmania Hsp90. Scientific Reports, 2019, 9, 14756.	1.6	15
7	Biological Evaluation of Arylsemicarbazone Derivatives as Potential Anticancer Agents. Pharmaceuticals, 2019, 12, 169.	1.7	9
8	Design, Synthesis and In Vitro Trypanocidal and Leishmanicidal Activities of 2â€(2â€Arylidene)hydrazonoâ€4â€oxothiazolidineâ€5â€acetic Acid Derivatives. ChemistrySelect, 2019, 4, 13163	-13772.	7
9	CycloMolder software: building theoretical cyclodextrin derivatives models and evaluating their host:guest interactions. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 2019, 93, 301-308.	0.9	9
10	Elucidation of the mechanism of complexation between oncocalyxone A and cyclodextrins by isothermal titration calorimetry and molecular modeling. Journal of Molecular Liquids, 2019, 274, 165-172.	2.3	17
11	Synthesis, antitrypanosomal activity and molecular docking studies of pyrimidine derivatives. Medicinal Chemistry Research, 2018, 27, 2512-2522.	1.1	5
12	Anti-hypersensitivity effects of the phthalimide derivative N-(4methyl-phenyl)-4-methylphthalimide in different pain models in mice. Biomedicine and Pharmacotherapy, 2017, 96, 503-512.	2.5	5
13	Supramolecular interactions between βâ€lapachone with cyclodextrins studied using isothermal titration calorimetry and molecular modeling. Journal of Molecular Recognition, 2017, 30, e2646.	1.1	10
14	Combination of In Silico Methods in the Search for Potential CD4+ and CD8+ T Cell Epitopes in the Proteome of Leishmania braziliensis. Frontiers in Immunology, 2016, 7, 327.	2.2	47
15	New 1,3-thiazole derivatives and their biological and ultrastructural effects on Trypanosoma cruzi. European Journal of Medicinal Chemistry, 2016, 121, 387-398.	2.6	46
16	Molecular modeling and cytotoxicity of diffractaic acid: HP-β-CD inclusion complex encapsulated in microspheres. International Journal of Biological Macromolecules, 2016, 92, 494-503.	3.6	15
17	New PPARÎ ³ partial agonist improves obesity-induced metabolic alterations and atherosclerosis in LDLrâ^'/â^' mice. Pharmacological Research, 2016, 104, 49-60.	3.1	26
18	Design, synthesis and structure–activity relationship of phthalimides endowed with dual antiproliferative and immunomodulatory activities. European Journal of Medicinal Chemistry, 2015, 96, 491-503.	2.6	34

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19	Thiosemicarbazones as Aedes aegypti larvicidal. European Journal of Medicinal Chemistry, 2015, 100, 162-175.	2.6	36
20	Synthesis, in vitro anticancer activity and in silico study of new disubstituted thiazolidinedione derivatives. Medicinal Chemistry Research, 2014, 23, 3220-3226.	1.1	18
21	Structural Design, Synthesis and Structure–Activity Relationships of Thiazolidinones with Enhanced Antiâ€ <i>Trypanosoma cruzi</i> Activity. ChemMedChem, 2014, 9, 177-188.	1.6	39
22	2-Pyridyl thiazoles as novel anti-Trypanosoma cruzi agents: Structural design, synthesis and pharmacological evaluation. European Journal of Medicinal Chemistry, 2014, 86, 48-59.	2.6	86
23	Sulfonamide–metal complexes endowed with potent anti-Trypanosoma cruzi activity. Journal of Enzyme Inhibition and Medicinal Chemistry, 2014, 29, 230-236.	2.5	13
24	Isolation, homology modeling and renal effects of a C-type natriuretic peptide from the venom of the Brazilian yellow scorpion (Tityus serrulatus). Toxicon, 2013, 74, 19-26.	0.8	14
25	Synthesis and in vitro anticancer activity of novel thiazacridine derivatives. Medicinal Chemistry Research, 2013, 22, 2421-2429.	1.1	23
26	Chemical synthesis, docking studies and biological effects of a pan peroxisome proliferator-activated receptor agonist and cyclooxygenase inhibitor. European Journal of Pharmaceutical Sciences, 2013, 48, 689-697.	1.9	18
27	Inhibition of Neurotoxic Secretory Phospholipases A2Enzymatic, Edematogenic, and Myotoxic Activities by Harpalycin 2, an Isoflavone Isolated fromHarpalyce brasilianaBenth. Evidence-based Complementary and Alternative Medicine, 2012, 2012, 1-9.	0.5	12
28	Structural Investigation of Anti- <i>Trypanosoma cruzi</i> 2-Iminothiazolidin-4-ones Allows the Identification of Agents with Efficacy in Infected Mice. Journal of Medicinal Chemistry, 2012, 55, 10918-10936.	2.9	55
29	Enhanced Antiproliferative Activity of the New Anticancer Candidate LPSF/AC04 in Cyclodextrin Inclusion Complexes Encapsulated into Liposomes. AAPS PharmSciTech, 2012, 13, 1355-1366.	1.5	26
30	The encapsulation of β-lapachone in 2-hydroxypropyl-β-cyclodextrin inclusion complex into liposomes: A physicochemical evaluation and molecular modeling approach. European Journal of Pharmaceutical Sciences, 2011, 44, 332-340.	1.9	59
31	Quercetin as an inhibitor of snake venom secretory phospholipase A2. Chemico-Biological Interactions, 2011, 189, 9-16.	1.7	62
32	Solid Dispersions of Imidazolidinedione by PEG and PVP Polymers with Potential Antischistosomal Activities. AAPS PharmSciTech, 2011, 12, 401-410.	1.5	22
33	The Use of Solid Dispersion Systems in Hydrophilic Carriers to Increase Benznidazole Solubility. Journal of Pharmaceutical Sciences, 2011, 100, 2443-2451.	1.6	53
34	A theoretical study of red-shifting and blue-shifting hydrogen bonds occurring between imidazolidine derivatives and PEG/PVP polymers. Journal of Molecular Modeling, 2010, 16, 119-127.	0.8	38
35	Discovery of Phthalimides as Immunomodulatory and Antitumor Drug Prototypes. ChemMedChem, 2010, 5, 523-528.	1.6	35
36	Synthesis and anti-inflammatory activity of new arylidene-thiazolidine-2,4-diones as PPARÎ ³ ligands. Bioorganic and Medicinal Chemistry, 2010, 18, 3805-3811.	1.4	86

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37	Studies toward the structural optimization of novel thiazolylhydrazone-based potent antitrypanosomal agents. Bioorganic and Medicinal Chemistry, 2010, 18, 7826-7835.	1.4	46
38	Halogen Atoms in the Modern Medicinal Chemistry: Hints for the Drug Design. Current Drug Targets, 2010, 11, 303-314.	1.0	528
39	Interaction of Morphine With a New α2-Adrenoceptor Agonist in Mice. Journal of Pain, 2010, 11, 71-78.	0.7	31
40	Novel Nitrofurazone Derivatives Endowed with Antimicrobial Activity. Archiv Der Pharmazie, 2008, 341, 655-660.	2.1	5
41	Estrutura, reatividade e propriedades biológicas de hidantoÃnas. Quimica Nova, 2008, 31, 614-622.	0.3	15
42	Synthesis, Cruzain Docking, and in vitro Studies of Arylâ€4â€Oxothiazolylhydrazones Against <i>Trypanosoma cruzi</i> . ChemMedChem, 2007, 2, 1339-1345.	1.6	50
43	A theoretical study of the solvent effects in ethylene oxide: Hydrofluoric acid complex using continuum and new discrete models. Computational and Theoretical Chemistry, 2007, 802, 91-97.	1.5	27
44	Synthesis, biological evaluation and molecular modeling studies of arylidene-thiazolidinediones with potential hypoglycemic and hypolipidemic activities. European Journal of Medicinal Chemistry, 2007, 42, 1263-1271.	2.6	46
45	Synthesis, docking, and in vitro activity of thiosemicarbazones, aminoacyl-thiosemicarbazides and acyl-thiazolidones against Trypanosoma cruzi. Bioorganic and Medicinal Chemistry, 2006, 14, 3749-3757.	1.4	98
46	AIPAR: ab initio parametrization of intermolecular potentials for computer simulations. Journal of Molecular Modeling, 2005, 11, 61-68.	0.8	7
47	Principal component analysis of the effects of wavefunction modification on the electrostatic potential of indole. International Journal of Quantum Chemistry, 2005, 102, 379-386.	1.0	1
48	Solute relaxation on the solvatochromism of ortho-betaine dyes. A sequential Monte Carlo/quantum mechanics study. Physical Chemistry Chemical Physics, 2004, 6, 2088.	1.3	27
49	Chemometric study of liquid water simulations. I. The parameters of the TIP4P model potential. Journal of Computational Chemistry, 2003, 24, 973-981.	1.5	19
50	The use of the generator coordinate method for designing basis set. Application to oxo-diperoxo molybdenum complexes. Computational and Theoretical Chemistry, 2002, 589-590, 251-264.	1.5	10
51	AGOA: A Hydration Procedure and Its Application to the 1-Phenyl-beta-Carboline Molecule. Journal of the Brazilian Chemical Society, 2002, 13, 36-42.	0.6	7
52	Molecular Structure of the Molybdenum Oxo-Diperoxo Compound MoO(O2)2(OPy)(H2O):Â A Computational and X-ray Study. Inorganic Chemistry, 2001, 40, 6022-6025.	1.9	28
53	Selective and mild oxidation of sulfides to sulfoxides by oxodiperoxo molybdenum complexes adsorbed onto silica gel. Tetrahedron, 2001, 57, 9669-9676.	1.0	60
54	Preparation, crystal structure determination, and properties of adducts of halogenomethyl compounds of indium with Group 16 donors. Journal of Organometallic Chemistry, 2001, 626, 68-75.	0.8	12

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55	Preparation, crystal structure determination and properties of adducts of indium methylene compounds with Group 15 donors. Journal of Organometallic Chemistry, 2000, 603, 203-212.	0.8	10
56	Coordination Chemistry of Br2InCH2Br:  Coordination at the Metal Center. Organometallics, 1999, 18, 99-105.	1.1	13
57	Rotational features of carbon-nitrogen bonds in axially chiral o-tert-butyl anilides and related molecules. Potential substrates for the â€~prochiral auxiliary' approach to asymmetric synthesis. Tetrahedron: Asymmetry, 1997, 8, 3955-3975.	1.8	112