

# Davi Brasil

## List of Publications by Year in descending order

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Version: 2024-02-01

52  
papers

832  
citations

516561

16  
h-index

501076

28  
g-index

52  
all docs

52  
docs citations

52  
times ranked

1241  
citing authors

#	ARTICLE	IF	CITATIONS
1	Adsorption of basic dyes onto activated carbon: Experimental and theoretical investigation of chemical reactivity of basic dyes using DFT-based descriptors. <i>Applied Surface Science</i> , 2018, 448, 662-670.	3.1	124
2	Obtaining extracts rich in antioxidant polysaccharides from the edible mushroom <i>Pleurotus ostreatus</i> using binary system with hot water and supercritical CO <sub>2</sub> . <i>Food Chemistry</i> , 2020, 330, 127173.	4.2	62
3	Gastroprotective effect of aparisthman, a diterpene isolated from <i>Aparisthmium cordatum</i> , on experimental gastric ulcer models in rats and mice. <i>Phytomedicine</i> , 2001, 8, 94-100.	2.3	59
4	Mercury levels assessment in hair of riverside inhabitants of the Tapaj�s River, Par� State, Amazon, Brazil: Fish consumption as a possible route of exposure. <i>Journal of Trace Elements in Medicine and Biology</i> , 2015, 30, 66-76.	1.5	46
5	Naphthoquinones isolated from <i>Eleutherine plicata</i> herb: in vitro antimalarial activity and molecular modeling to investigate their binding modes. <i>Medicinal Chemistry Research</i> , 2020, 29, 487-494.	1.1	43
6	Combined Kinetic Studies and Computational Analysis on Kojic Acid Analogs as Tyrosinase Inhibitors. <i>Molecules</i> , 2014, 19, 9591-9605.	1.7	41
7	A SAR and QSAR Study of New Artemisinin Compounds with Antimalarial Activity. <i>Molecules</i> , 2014, 19, 367-399.	1.7	38
8	An In Silico Study of the Antioxidant Ability for Two Caffeine Analogs Using Molecular Docking and Quantum Chemical Methods. <i>Molecules</i> , 2018, 23, 2801.	1.7	38
9	Chemical profile of <i>Lippia thymoides</i> , evaluation of the acetylcholinesterase inhibitory activity of its essential oil, and molecular docking and molecular dynamics simulations. <i>PLoS ONE</i> , 2019, 14, e0213393.	1.1	34
10	Identification of Novel Protein Kinase Receptor Type 2 Inhibitors Using Pharmacophore and Structure-Based Virtual Screening. <i>Molecules</i> , 2018, 23, 453.	1.7	30
11	Evaluation of the Gastroprotective Activity of Cordatin, a Diterpene Isolated from <i>Aparisthmium cordatum</i> (Euphorbiaceae).. <i>Biological and Pharmaceutical Bulletin</i> , 2000, 23, 1465-1469.	0.6	29
12	Biotransformation of chalcones by the endophytic fungus <i>Aspergillus flavus</i> isolated from <i>Paspalum maritimum</i> trin. <i>Journal of the Brazilian Chemical Society</i> , 2011, 22, 1333-1338.	0.6	28
13	Virtual Screening and Statistical Analysis in the Design of New Caffeine Analogues Molecules with Potential Epithelial Anticancer Activity. <i>Current Pharmaceutical Design</i> , 2018, 24, 576-594.	0.9	28
14	H <sub>3</sub> PO <sub>4</sub> -activated carbons produced from a�sai stones and Brazil nut shells: removal of basic blue 26 dye from aqueous solutions by adsorption. <i>Environmental Science and Pollution Research</i> , 2019, 26, 28533-28547.	2.7	25
15	A QSAR, Pharmacokinetic and Toxicological Study of New Artemisinin Compounds with Anticancer Activity. <i>Molecules</i> , 2014, 19, 10670-10697.	1.7	24
16	Leishmanicidal Activity of (+)-Phyllanthidine and the Phytochemical Profile of <i>Margaritaria nobilis</i> (Phyllanthaceae). <i>Molecules</i> , 2015, 20, 22157-22169.	1.7	24
17	Studies of NMR, molecular docking, and molecular dynamics simulation of new promising inhibitors of cruzaine from the parasite <i>Trypanosoma cruzi</i> . <i>Medicinal Chemistry Research</i> , 2019, 28, 246-259.	1.1	15
18	Rational Design of Antimalarial Drugs Using Molecular Modeling and Statistical Analysis. <i>Current Pharmaceutical Design</i> , 2015, 21, 4112-4127.	0.9	13

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19	Ligand- and structure-based virtual screening from 16-((diisobutylamino)methyl)-6 $\beta$ -hydroxyivouacapane-7 $\beta$ ,17 $\beta$ -lactone a compound with potential anti-prostate cancer activity. <i>Journal of the Serbian Chemical Society</i> , 2019, 84, 153-174.	0.4	13
20	Crystal structure and theoretical calculations of Julocrotine, a natural product with antileishmanial activity. <i>International Journal of Quantum Chemistry</i> , 2008, 108, 513-520.	1.0	12
21	Molecular Modeling of the Major Compounds of Sesquiterpenes Class in Copaiba Oil-resin. <i>British Journal of Pharmaceutical Research</i> , 2015, 7, 247-263.	0.4	12
22	Essential oil composition of <i>Croton palanostigma</i> Klotzsch from north Brazil. <i>Journal of the Brazilian Chemical Society</i> , 2009, 20, 1188-1192.	0.6	11
23	Volatiles, A Glutarimide Alkaloid and Antimicrobial Effects of <i>Croton pullei</i> (Euphorbiaceae). <i>Molecules</i> , 2013, 18, 3195-3205.	1.7	11
24	Crystal structure and theoretical study of IR and $^1\text{H}$ and $^{13}\text{C}$ NMR spectra of cordatin, a natural product with antiulcerogenic activity. <i>International Journal of Quantum Chemistry</i> , 2008, 108, 2564-2575.	1.0	10
25	Theoretical and experimental study of aparisthman: A natural product with anti-ulcer activity. <i>International Journal of Quantum Chemistry</i> , 2006, 106, 2706-2713.	1.0	8
26	Isolation, X-ray crystal structure and theoretical calculations of the new compound 8-Eepicordatin and identification of others terpenes and steroids from the bark and leaves of <i>Croton palanostigma</i> Klotzsch. <i>Journal of the Brazilian Chemical Society</i> , 2010, 21, 731-739.	0.6	8
27	Glycerol and fatty acid influences on the rheological and technological properties of composite films from residues of <i>Cynoscion acoupa</i> . <i>Food Bioscience</i> , 2020, 38, 100773.	2.0	6
28	Study of molecular interactions between Chitosan and Vi Antigen. <i>Journal of Molecular Graphics and Modelling</i> , 2017, 72, 148-155.	1.3	5
29	Theoretical study via DFT for prediction of $^{13}\text{C}$ and $^1\text{H}$ NMR data of two diterpenoids derived from the root of <i>salvia grandifolia</i> . <i>Journal of the Serbian Chemical Society</i> , 2019, 84, 591-598.	0.4	5
30	Identification of ( $\hat{\alpha}$ )-(E)-N-[2(S)-Hydroxy-2-(4-hydroxyphenyl) ethyl]ferulamide, a Natural Product Isolated from <i>Croton Pullei</i> : Theoretical and Experimental Analysis. <i>International Journal of Molecular Sciences</i> , 2011, 12, 9389-9403.	1.8	4
31	Antioxidant Activity of an Industrial Cupuassu Seed By- $\text{\textcircled{e}}$ product: Molecular Modeling of Phenolic Compounds. <i>Chemical Engineering and Technology</i> , 2019, 42, 397-406.	0.9	3
32	Biological activities of <i>Croton palanostigma</i> Klotzsch. <i>Pharmacognosy Magazine</i> , 2015, 11, 601.	0.3	3
33	<i>Aspidosperma excelsum</i> and its pharmacological potential: in silico studies of pharmacokinetic prediction, toxicological and biological activity. <i>Research, Society and Development</i> , 2020, 9, e3629108635.	0.0	3
34	In silico identification of novel allosteric inhibitors of Dengue virus NS2B/NS3 serine protease. <i>Journal of the Serbian Chemical Society</i> , 2022, 87, 693-706.	0.4	3
35	Density Functional Theory Calculations of the Nuclear Magnetic Resonance Parameters for Two Dihydrochalcones. <i>Journal of Computational and Theoretical Nanoscience</i> , 2012, 9, 953-956.	0.4	2
36	Structure modeling of a metalloendopeptidase from <i>Corynebacterium pseudotuberculosis</i> . <i>Computers in Biology and Medicine</i> , 2012, 42, 538-541.	3.9	2

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37	Processos de extração e usos industriais de óleos de andiroba e açafrão: uma revisão. <i>Research, Society and Development</i> , 2021, 10, e229101220227.	0.0	2
38	Psicologia ambiental e problemas ambientais: uma revisão de literatura. <i>Doxa Revista Brasileira De Psicologia E Educação</i> , 2020, 22, 108-122.	0.1	2
39	Síntese e caracterização de zeólita tipo sodalita obtida a partir de resíduo de caulim. <i>Ceramica</i> , 2020, 66, 404-412.	0.3	2
40	Óleo de Castanha-do-Brasil: métodos de extração e aplicações na indústria. <i>Research, Society and Development</i> , 2022, 11, e29511427256.	0.0	2
41	Pharmaceutical and Biological Potential of the Croton palanostigma Isolated Compounds. <i>Journal of Computational and Theoretical Nanoscience</i> , 2019, 16, 1773-1782.	0.4	1
42	Estudo de quantificação dos compostos fenólicos, avaliação das atividades antioxidante e antimicrobiana da casca do caule do Croton cajucara BENTH. <i>Research, Society and Development</i> , 2020, 9, e1929119742.	0.0	1
43	Isolation, Identification, predictions of biological activity and molecular docking of the diterpene jateorin obtained from the stem of odontocarya tamoides (D.C.) miers. <i>Chemical Data Collections</i> , 2019, 21, 100215.	1.1	0
44	Abordagens de modelagem para investigar a interação entre Fullereno e $\beta$ -Amiloide. <i>Research, Society and Development</i> , 2021, 10, e261101119585.	0.0	0
45	Crystal Structure of Limonoid 6-Acetylsvietephragmin and Theoretical Study of Nuclear Magnetic Resonance Spectra of Phragmalin Limonoids. <i>Advanced Science Letters</i> , 2012, 18, 150-157.	0.2	0
46	Saúde ambiental: uma análise situacional e a possibilidade de prevenção de doenças ambientais em Redenção-PA. <i>Research, Society and Development</i> , 2020, 9, e217996580.	0.0	0
47	Computational simulation of nanostructured lipid carrier containing lipids from Cupuassu ( <i>Theobroma grandiflorum</i> ) seed fat: Design, interaction and molecular dynamic study. <i>Research, Society and Development</i> , 2020, 9, e92191110433.	0.0	0
48	IMPLEMENTAÇÃO DO CAMPO DE FORÇA CLAYFF NO GROMACS: UMA APLICAÇÃO EM ESTRUTURA DE CAULINITA. <i>Química Nova</i> , 0, , .	0.3	0
49	Análise isotérmica da atividade de água (aw) de sementes de pimenta-do-reino ( <i>Piper nigrum</i> , L.) em câmara de secagem. <i>Revista Brasileira De Tecnologia Agroindustrial</i> , 2020, 14, .	0.1	0
50	Atividades biológicas da espécie <i>spilanthus acmella</i> : uma revisão. <i>Research, Society and Development</i> , 2021, 10, e404101422035.	0.0	0
51	Extinção de copos descartáveis: análise ambiental e econômica em uma instituição de ensino superior do sul do Pará. <i>Research, Society and Development</i> , 2020, 9, e2189108321.	0.0	0
52	Guaranã ( <i>Paullinia cupana</i> Kunth), marapuama ( <i>Ptychopetalum olacoides</i> Benth.), genciana ( <i>Gentiana</i> ) Tj ETQq0 0 0 rgBT /Overlock 10 <i>Development</i> , 2022, 11, e21711224592.	0.0	0