

Milan MladenoviÄ

List of Publications by Year in descending order

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Version: 2024-02-01

26
papers

540
citations

759190

12
h-index

642715

23
g-index

27
all docs

27
docs citations

27
times ranked

922
citing authors

#	ARTICLE	IF	CITATIONS
1	Human estrogen receptor \pm antagonists, part 2: Synthesis driven by rational design, in vitro antiproliferative, and in vivo anticancer evaluation of innovative coumarin-related antiestrogens as breast cancer suppressants. <i>European Journal of Medicinal Chemistry</i> , 2022, 227, 113869.	5.5	6
2	Human Estrogen Receptor Alpha Antagonists, Part 3: 3-D Pharmacophore and 3-D QSAR Guided Brefeldin A Hit-to-Lead Optimization toward New Breast Cancer Suppressants. <i>Molecules</i> , 2022, 27, 2823.	3.8	3
3	Novel non-covalent LSD1 inhibitors endowed with anticancer effects in leukemia and solid tumor cellular models. <i>European Journal of Medicinal Chemistry</i> , 2022, 237, 114410.	5.5	15
4	Ligand-based and structure-based studies to develop predictive models for SARS-CoV-2 main protease inhibitors through the 3d-qsar.com portal. <i>Journal of Computer-Aided Molecular Design</i> , 2022, 36, 483-505.	2.9	4
5	Low cytotoxic quinoline-4-carboxylic acids derived from vanillin precursors as potential human dihydroorotate dehydrogenase inhibitors. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2021, 46, 128194.	2.2	3
6	Pyrazole Derivatives of Medically Relevant Phenolic Acids: Insight into Antioxidative and Anti-LOX Activity. <i>Medicinal Chemistry</i> , 2021, 17, 807-819.	1.5	1
7	In vitro and in silico lipoxygenase inhibition studies and antimicrobial activity of pyrazolyl-phthalazine-diones. <i>Kragujevac Journal of Science</i> , 2021, , 35-52.	0.4	2
8	Human Estrogen Receptor \pm Antagonists. Part 1: 3-D QSAR-Driven Rational Design of Innovative Coumarin-Related Antiestrogens as Breast Cancer Suppressants through Structure-Based and Ligand-Based Studies. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 5028-5053.	5.4	5
9	Potent human dihydroorotate dehydrogenase inhibitory activity of new quinoline-4-carboxylic acids derived from phenolic aldehydes: Synthesis, cytotoxicity, lipophilicity and molecular docking studies. <i>Bioorganic Chemistry</i> , 2020, 105, 104373.	4.1	7
10	DFT calculations as an efficient tool for prediction of Raman and infra-red spectra and activities of newly synthesized cathinones. <i>Open Chemistry</i> , 2020, 18, 185-195.	1.9	11
11	16-membered macrolide antibiotics: a review. <i>International Journal of Antimicrobial Agents</i> , 2018, 51, 283-298.	2.5	118
12	The Targeted Pesticides as Acetylcholinesterase Inhibitors: Comprehensive Cross-Organism Molecular Modelling Studies Performed to Anticipate the Pharmacology of Harmfulness to Humans In Vitro. <i>Molecules</i> , 2018, 23, 2192.	3.8	36
13	Understanding the Molecular Determinant of Reversible Human Monoamine Oxidase B Inhibitors Containing 2- <i>H</i> -Chromen-2-One Core: Structure-Based and Ligand-Based Derived Three-Dimensional Quantitative Structure-Activity Relationships Predictive Models. <i>Journal of Chemical Information and Modeling</i> , 2017, 57, 787-814.	5.4	33
14	Mannich bases of 1,2,4-triazole-3-thione containing adamantane moiety: Synthesis, preliminary anticancer evaluation, and molecular modeling studies. <i>Chemical Biology and Drug Design</i> , 2017, 89, 943-952.	3.2	17
15	Filipendula ulmaria extracts attenuate cisplatin-induced liver and kidney oxidative stress in rats: In vivo investigation and LC-MS analysis. <i>Food and Chemical Toxicology</i> , 2017, 99, 86-102.	3.6	38
16	Serum albumin binding analysis and toxicological screening of novel chroman-2,4-diones as oral anticoagulants. <i>Chemico-Biological Interactions</i> , 2015, 227, 18-31.	4.0	5
17	Newly discovered chroman-2,4-diones neutralize the in vivo DNA damage induced by alkylation through the inhibition of Topoisomerase III \pm : A story behind the molecular modeling approach. <i>Biochemical Pharmacology</i> , 2015, 98, 243-266.	4.4	3
18	In vitro and in vivo assessment of the genotoxicity and antigenotoxicity of the Filipendula hexapetala and Filipendula ulmaria methanol extracts. <i>Journal of Ethnopharmacology</i> , 2015, 174, 287-292.	4.1	10

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19	Dropwort (<i>Filipendula hexapetala</i> Gilib.): potential role as antioxidant and antimicrobial agent. EXCLI Journal, 2015, 14, 1-20.	0.7	19
20	Synthesis and toxicological studies of in vivo anticoagulant activity of novel 3-(1-aminoethylidene)chroman-2,4-diones and 4-hydroxy-3-(1-iminoethyl)-2H-chromen-2-ones combined with a structure-based 3-D pharmacophore model. European Journal of Pharmaceutical Sciences, 2014, 55, 20-35.	4.0	10
21	Combining molecular docking and 3-D pharmacophore generation to enclose the in vivo antigenotoxic activity of naturally occurring aromatic compounds: Myricetin, quercetin, rutin, and rosmarinic acid. Biochemical Pharmacology, 2013, 86, 1376-1396.	4.4	27
22	Chemical composition, antioxidant and antigenotoxic activities of different fractions of <i>Gentiana asclepiadea</i> L. roots extract. EXCLI Journal, 2013, 12, 807-23.	0.7	13
23	Biochemical and pharmacological evaluation of 4-hydroxychromen-2-ones bearing polar C-3 substituents as anticoagulants. European Journal of Medicinal Chemistry, 2012, 54, 144-158.	5.5	12
24	In Vitro Antioxidant Activity of Selected 4-Hydroxy-chromene-2-one Derivativesâ€™ SAR, QSAR and DFT Studies. International Journal of Molecular Sciences, 2011, 12, 2822-2841.	4.1	78
25	Design of Novel 4-Hydroxy-chromene-2-one Derivatives as Antimicrobial Agents. Molecules, 2010, 15, 4294-4308.	3.8	24
26	Synthesis and Molecular Descriptor Characterization of Novel 4-Hydroxy-chromene-2-one Derivatives as Antimicrobial Agents. Molecules, 2009, 14, 1495-1512.	3.8	40