

Asmaa M Aboulmagd

List of Publications by Year in descending order

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Version: 2024-02-01

20
papers

275
citations

1040056

9
h-index

940533

16
g-index

20
all docs

20
docs citations

20
times ranked

409
citing authors

#	ARTICLE	IF	CITATIONS
1	Development and Greenness Assessment of HPLC Method for Studying the Pharmacokinetics of Co-Administered Metformin and Papaya Extract. <i>Molecules</i> , 2022, 27, 375.	3.8	1
2	Pioglitazone Synthetic Analogue Ameliorates Streptozotocin-Induced Diabetes Mellitus through Modulation of ACE 2/Angiotensin 1 α 7 via PI3K/AKT/mTOR Signaling Pathway. <i>Pharmaceuticals</i> , 2022, 15, 341.	3.8	2
3	Radiolabeling, biological distribution, docking and ADME studies of 99mTc-Ros as a promising natural tumor tracer. <i>Applied Radiation and Isotopes</i> , 2022, 184, 110196.	1.5	3
4	Design, synthesis and mechanistic study of new benzenesulfonamide derivatives as anticancer and antimicrobial agents <i>via</i> carbonic anhydrase IX inhibition. <i>RSC Advances</i> , 2021, 11, 26241-26257.	3.6	30
5	Metabolomic profiling, biological evaluation of <i>Aspergillus awamori</i> , the river Nile-derived fungus using epigenetic and OSMAC approaches. <i>RSC Advances</i> , 2021, 11, 6709-6719.	3.6	7
6	Ligand-based design, molecular dynamics and ADMET studies of suggested SARS-CoV-2 Mpro inhibitors. <i>RSC Advances</i> , 2021, 11, 4523-4538.	3.6	9
7	Ecological HPLC method for analyzing an antidiabetic drug in real rat plasma samples and studying the effects of concurrently administered fenugreek extract on its pharmacokinetics. <i>RSC Advances</i> , 2021, 11, 4740-4750.	3.6	6
8	Holospiniferoside: A New Antitumor Cerebroside from The Red Sea Cucumber <i>Holothuria spinifera</i> : In Vitro and In Silico Studies. <i>Molecules</i> , 2021, 26, 1555.	3.8	4
9	Analysis of sunitinib malate, a multi-targeted tyrosine kinase inhibitor: A critical review. <i>Microchemical Journal</i> , 2021, 163, 105926.	4.5	3
10	Design, microwave assisted synthesis, and molecular modeling study of some new 1,3,4-thiadiazole derivatives as potent anticancer agents and potential VEGFR-2 inhibitors. <i>Bioorganic Chemistry</i> , 2021, 112, 104923.	4.1	13
11	Antiulcer secondary metabolites from <i>Elaeocarpus grandis</i> , family Elaeocarpaceae, supported by <i>in silico</i> studies. <i>RSC Advances</i> , 2020, 10, 34788-34799.	3.6	6
12	New Cytotoxic Cerebrosides from the Red Sea Cucumber <i>Holothuria spinifera</i> Supported by In-Silico Studies. <i>Marine Drugs</i> , 2020, 18, 405.	4.6	7
13	New fused pyrimidine derivatives with anticancer activity: Synthesis, topoisomerase II inhibition, apoptotic inducing activity and molecular modeling study. <i>Bioorganic Chemistry</i> , 2020, 103, 104134.	4.1	25
14	Quinazoline-Schiff base conjugates: <i>in silico</i> study and ADMET predictions as multi-target inhibitors of coronavirus (SARS-CoV-2) proteins. <i>RSC Advances</i> , 2020, 10, 34033-34045.	3.6	37
15	Radiolabeling, docking studies, <i>in silico</i> ADME and biological evaluation of serotonin with 125I for 5-HTRs imaging. <i>Nuclear Science and Techniques/Hewuli</i> , 2020, 31, 1.	3.4	1
16	Nature as a treasure trove of potential anti-SARS-CoV drug leads: a structural/mechanistic rationale. <i>RSC Advances</i> , 2020, 10, 19790-19802.	3.6	71
17	Triple-negative breast cancer suppressive activities, antioxidants and pharmacophore model of new acylated rhamnopyranoses from <i>Premna odorata</i> . <i>RSC Advances</i> , 2020, 10, 10584-10598.	3.6	16
18	Testicular Caspase-3 and β -Catenin Regulators Predicted via Comparative Metabolomics and Docking Studies. <i>Metabolites</i> , 2020, 10, 31.	2.9	14

#	ARTICLE	IF	CITATIONS
19	Saccharomonosporine A inspiration; synthesis of potent analogues as potential PIM kinase inhibitors. RSC Advances, 2020, 10, 6752-6762.	3.6	8
20	Bioassay-Guided Isolation, Metabolic Profiling, and Docking Studies of Hyaluronidase Inhibitors from <i>Ravenala madagascariensis</i> . Molecules, 2020, 25, 1714.	3.8	12