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List of Publications by Year in descending order

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713013 623188 39 541 14 21 citations g-index h-index papers 43 43 43 618 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	Novel scaffold hopping of potent benzothiazole and isatin analogues linked to 1,2,3-triazole fragment that mimic quinazoline epidermal growth factor receptor inhibitors: Synthesis, antitumor and mechanistic analyses. Bioorganic Chemistry, 2020, 103, 104133.	2.0	53
2	Design, synthesis, molecular docking of new lipophilic acetamide derivatives affording potential anticancer and antimicrobial agents. Bioorganic Chemistry, 2018, 76, 332-342.	2.0	38
3	Design, Synthesis and in Vivo Anti-inflammatory Activities of 2,4-Diaryl-5-4H-imidazolone Derivatives. Molecules, 2012, 17, 12262-12275.	1.7	37
4	The rational design, synthesis, and antimicrobial investigation of 2-Amino-4-Methylthiazole analogues inhibitors of GlcN-6-P synthase. Bioorganic Chemistry, 2020, 99, 103781.	2.0	31
5	Crystal structure of carbonmonoxy sickle hemoglobin in R-state conformation. Journal of Structural Biology, 2016, 194, 446-450.	1.3	30
6	Novel molecular discovery of promising amidine-based thiazole analogues as potent dual Matrix Metalloproteinase-2 and 9 inhibitors: Anticancer activity data with prominent cell cycle arrest and DNA fragmentation analysis effects. Bioorganic Chemistry, 2020, 101, 103992.	2.0	26
7	Design, synthesis, and biological evaluation studies of novel quinazolinone derivatives as anticonvulsant agents. Medicinal Chemistry Research, 2013, 22, 5823-5831.	1.1	25
8	Molecular Mimics of Classic P-Glycoprotein Inhibitors as Multidrug Resistance Suppressors and Their Synergistic Effect on Paclitaxel. PLoS ONE, 2017, 12, e0168938.	1.1	22
9	Antimicrobial screening and pharmacokinetic profiling of novel phenyl-[1,2,4]triazolo[4,3-a]quinoxaline analogues targeting DHFR and E. coli DNA gyrase B. Bioorganic Chemistry, 2020, 96, 103656.	2.0	22
10	Synthesis, Modelling, and Anticonvulsant Studies of New Quinazolines Showing Three Highly Active Compounds with Low Toxicity and High Affinity to the GABA-A Receptor. Molecules, 2017, 22, 188.	1.7	19
11	Thiophenesâ€"Naturally Occurring Plant Metabolites: Biological Activities and In Silico Evaluation of Their Potential as Cathepsin D Inhibitors. Plants, 2022, 11, 539.	1.6	19
12	Synthesis and screening of some new fluorinated quinazolinone–sulphonamideÂhybrids as anticancer agents. Journal of Taibah University Medical Sciences, 2015, 10, 333-339.	0.5	17
13	Identification of a novel class of covalent modifiers of hemoglobin as potential antisickling agents. Organic and Biomolecular Chemistry, 2015, 13, 6353-6370.	1.5	16
14	Inborn errors in the vitamin B6 salvage enzymes associated with neonatal epileptic encephalopathy and other pathologies. Biochimie, 2021, 183, 18-29.	1.3	16
15	Discovery of novel quinoline-based analogues of combretastatin A-4 as tubulin polymerisation inhibitors with apoptosis inducing activity and potent anticancer effect. Journal of Enzyme Inhibition and Medicinal Chemistry, 2021, 36, 802-818.	2.5	15
16	VZHE-039, a novel antisickling agent that prevents erythrocyte sickling under both hypoxic and anoxic conditions. Scientific Reports, 2020, 10, 20277.	1.6	14
17	Chaetomugilins and Chaetoviridinsâ€"Promising Natural Metabolites: Structures, Separation, Characterization, Biosynthesis, Bioactivities, Molecular Docking, and Molecular Dynamics. Journal of Fungi (Basel, Switzerland), 2022, 8, 127.	1.5	14
18	Potent Quinoline-Containing Combretastatin A-4 Analogues: Design, Synthesis, Antiproliferative, and Anti-Tubulin Activity. Pharmaceuticals, 2020, 13, 393.	1.7	12

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19	Targeted potent antimicrobial benzochromene-based analogues: Synthesis, computational studies, and inhibitory effect against 14î±-Demethylase and DNA Gyrase. Bioorganic Chemistry, 2020, 105, 104387.	2.0	10
20	Design, Synthesis, and Investigation of Novel Nitric Oxide (NO)-Releasing Prodrugs as Drug Candidates for the Treatment of Ischemic Disorders: Insights into NO-Releasing Prodrug Biotransformation and Hemoglobin–NO Biochemistry. Biochemistry, 2015, 54, 7178-7192.	1.2	9
21	Phytoconstituents, In Vitro Anti-Infective Activity of Buddleja indica Lam., and In Silico Evaluation of its SARS-CoV-2 Inhibitory Potential. Frontiers in Pharmacology, 2021, 12, 619373.	1.6	9
22	Modulating hemoglobin allostery for treatment of sickle cell disease: current progress and intellectual property. Expert Opinion on Therapeutic Patents, 2022, 32, 115-130.	2.4	9
23	Design, Synthesis and Antiproliferative Activities of Oxidative Stress Inducers Based on 2-Styryl-3,5-dihydro-4 <i>H</i> -imidazol-4-one Scaffold. Chemical and Pharmaceutical Bulletin, 2018, 66, 967-975.	0.6	8
24	Discovery and SAR of Novel Disubstituted Quinazolines as Dual PI3Kalpha/mTOR Inhibitors Targeting Breast Cancer. ACS Medicinal Chemistry Letters, 2020, 11, 2156-2164.	1.3	8
25	<p>Zein-alpha lipoic acid-loaded nanoparticles to enhance the oral bioavailability of dapoxetine: optimization and clinical pharmacokinetic evaluation</p> . International Journal of Nanomedicine, 2019, Volume 14, 7461-7473.	3.3	7
26	Exploring the Activity of Fungal Phenalenone Derivatives as Potential CK2 Inhibitors Using Computational Methods. Journal of Fungi (Basel, Switzerland), 2022, 8, 443.	1.5	7
27	Structural modification of azolylacryloyl derivatives yields a novel class of covalent modifiers of hemoglobin as potential antisickling agents. MedChemComm, 2019, 10, 1900-1906.	3.5	6
28	An Investigation of Structure-Activity Relationships of Azolylacryloyl Derivatives Yielded Potent and Long-Acting Hemoglobin Modulators for Reversing Erythrocyte Sickling. Biomolecules, 2020, 10, 1508.	1.8	6
29	1H-Imidazole-2,5-Dicarboxamides as NS4A Peptidomimetics: Identification of a New Approach to Inhibit HCV-NS3 Protease. Biomolecules, 2020, 10, 479.	1.8	5
30	Aryloxyalkanoic Acids as Non-Covalent Modifiers of the Allosteric Properties of Hemoglobin. Molecules, 2016, 21, 1057.	1.7	4
31	Introducing of potent cytotoxic novel 2-(aroylamino)cinnamamide derivatives against colon cancer mediated by dual apoptotic signal activation and oxidative stress. Bioorganic Chemistry, 2020, 101, 103953.	2.0	4
32	Improving the Solubility and Oral Bioavailability of a Novel Aromatic Aldehyde Antisickling Agent (PP10) for the Treatment of Sickle Cell Disease. Pharmaceutics, 2021, 13, 1148.	2.0	4
33	Molecular insight into 2-phosphoglycolate activation of the phosphatase activity of bisphosphoglycerate mutase. Acta Crystallographica Section D: Structural Biology, 2022, 78, 472-482.	1.1	4
34	Pairing 3D-Printing with Nanotechnology to Manage Metabolic Syndrome. International Journal of Nanomedicine, 2022, Volume 17, 1783-1801.	3.3	4
35	Synthetic bulky NS4A peptide variants bind to and inhibit HCV NS3 protease. Journal of Advanced Research, 2020, 24, 251-259.	4.4	3
36	Insights on Cancer Cell Inhibition, Subcellular Activities, and Kinase Profile of Phenylacetamides Pending 1H-Imidazol-5-One Variants. Frontiers in Pharmacology, 2021, 12, 794325.	1.6	3

#	Article	IF	CITATIONS
37	Clinical Pharmacokinetic Evaluation of Optimized Liquisolid Tablets as a Potential Therapy for Male Sexual Dysfunction. Pharmaceutics, 2020, 12, 1187.	2.0	2
38	A <scp>singleâ€step</scp> synthesis of 1,3,4, <scp>6â€tetraaryl</scp> â€5â€aryliminopiperazinâ€2â€one. Journ Heterocyclic Chemistry, 2021, 58, 442-449.	al of 1.4	2
39	Reexamining Povarov Reaction's Scope and Limitation in the Generation of HCV-NS4A Peptidomimetics. Heteroatom Chemistry, 2022, 2022, 1-12.	0.4	0