Richie R Bhandare

List of Publications by Year in descending order

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759233 713466 34 485 12 21 citations h-index g-index papers 35 35 35 349 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	An overview of <i>in silico</i> methods used in the design of VEGFR-2 inhibitors as anticancer agents. ChemistrySelect, 2023, 8, 2441-2457.	1.5	O
2	Forging of nicotine for the effective management of diabetic wounds: A hybrid of scaffold hopping and molecular dynamics simulation approaches. Arabian Journal of Chemistry, 2022, 15, 103585.	4.9	1
3	Antitubercular and antioxidant activities of hydroxy and chloro substituted chalcone analogues: Synthesis, biological and computational studies. Arabian Journal of Chemistry, 2022, 15, 103581.	4.9	19
4	A critical review on phytopharmacology, spectral and computational analysis of phytoconstituents from Streblus asper Lour. Phytomedicine Plus, 2022, 2, 100177.	2.0	2
5	Heterogeneous graphitic carbon nitrides in visible-light-initiated organic transformations. Green Chemistry, 2022, 24, 438-479.	9.0	47
6	Quinoline conjugated 2-azetidinone derivatives as prospective anti-breast cancer agents: In vitro antiproliferative and anti-EGFR activities, molecular docking and in-silico drug likeliness studies. Journal of Saudi Chemical Society, 2022, 26, 101471.	5.2	15
7	Multistep synthesis and screening of heterocyclic tetrads containing furan, pyrazoline, thiazole and triazole (or oxadiazole) as antimicrobial and anticancer agents. Journal of Saudi Chemical Society, 2022, 26, 101447.	5.2	20
8	L-Glutamic acid loaded collagen chitosan composite scaffold as regenerative medicine for the accelerated healing of diabetic wounds. Arabian Journal of Chemistry, 2022, 15, 103841.	4.9	3
9	Antitubercular activity assessment of fluorinated chalcones, 2-aminopyridine-3-carbonitrile and 2-amino-4H-pyran-3-carbonitrile derivatives: In vitro, molecular docking and in-silico drug likeliness studies. PLoS ONE, 2022, 17, e0265068.	2.5	4
10	Selectivity profile comparison for certain \hat{I}^3 -butyrolactone and oxazolidinone-based ligands on a sigma 2 receptor over sigma 1: a molecular docking approach. RSC Advances, 2022, 12, 20096-20109.	3.6	3
11	Synthesis, and biological screening of chloropyrazine conjugated benzothiazepine derivatives as potential antimicrobial, antitubercular and cytotoxic agents. Arabian Journal of Chemistry, 2021, 14, 102915.	4.9	15
12	Antimicrobial Hexaaquacopper(II) Complexes with Novel Polyiodide Chains. Polymers, 2021, 13, 1005.	4.5	7
13	Design, synthesis, and biological evaluation of novel bromo-pyrimidine analogues as tyrosine kinase inhibitors. Arabian Journal of Chemistry, 2021, 14, 103054.	4.9	7
14	Thiazole–Chalcone Hybrids as Prospective Antitubercular and Antiproliferative Agents: Design, Synthesis, Biological, Molecular Docking Studies and In Silico ADME Evaluation. Molecules, 2021, 26, 2847.	3.8	26
15	Design, multistep synthesis and in-vitro antimicrobial and antioxidant screening of coumarin clubbed chalcone hybrids through molecular hybridization approach. Arabian Journal of Chemistry, 2021, 14, 103154.	4.9	24
16	A key review on oxadiazole analogs as potential methicillin-resistant Staphylococcus aureus (MRSA) activity: Structure-activity relationship studies. European Journal of Medicinal Chemistry, 2021, 219, 113442.	5.5	58
17	2D-Quantitative structure activity relationship (QSAR) modeling, docking studies, synthesis and in-vitro evaluation of 1,3,4-thiadiazole tethered coumarin derivatives as antiproliferative agents. Journal of Saudi Chemical Society, 2021, 25, 101279.	5.2	7
18	Synthetic Strategies of Pyrimidine-Based Scaffolds as Aurora Kinase and Polo-like Kinase Inhibitors. Molecules, 2021, 26, 5170.	3.8	8

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19	Assessment of the Antimicrobial and Antiproliferative Activities of Chloropyrazine-Tethered Pyrimidine Derivatives: In Vitro, Molecular Docking, and In-Silico Drug-Likeness Studies. Applied Sciences (Switzerland), 2021, 11, 10734.	2.5	10
20	Design, Synthesis, and Antibacterial and Antifungal Activities of Novel Trifluoromethyl and Trifluoromethoxy Substituted Chalcone Derivatives. Pharmaceuticals, 2020, 13, 375.	3.8	29
21	Design, Facile Synthesis and Characterization of Dichloro Substituted Chalcones and Dihydropyrazole Derivatives for Their Antifungal, Antitubercular and Antiproliferative Activities. Molecules, 2020, 25, 3188.	3.8	26
22	Antimicrobial, Antioxidant, and Anticancer Activities of Some Novel Isoxazole Ring Containing Chalcone and Dihydropyrazole Derivatives. Molecules, 2020, 25, 1047.	3.8	71
23	Enhanced solubility of microwave-assisted synthesized acyclovir Co-crystals. Research Journal of Pharmacy and Technology, 2020, 13, 5979-5986.	0.8	6
24	Green Synthesis of Potent Antimicrobial Silver Nanoparticles Using Different Plant Extracts and Their Mixtures. Processes, 2019, 7, 510.	2.8	41
25	Emerging Phytochemicals and Bioactive Compounds from a Desert Plant Prosopis cineraria (L.) Druce and Future Prospects., 2019,, 19-51.		2
26	Public perception toward e-commerce of medicines and comparative pharmaceutical quality assessment study of two different products of furosemide tablets from community and illicit online pharmacies. Journal of Pharmacy and Bioallied Sciences, 2019, 11, 284.	0.6	14
27	In Vitro Comparative Quality Attributes of Selected Brands of Fexofenadine Hydrochloride Tablets Marketed in UAE and India., 2019,, 85-96.		O
28	Cardiovascular Concern of 5-HT2B Receptor and Recent Vistas in the Development of Its Antagonists. Cardiovascular & Hematological Disorders Drug Targets, 2017, 17, 86-104.	0.7	4
29	Novel \hat{I}^3 -Butyrolactone Derivatives as Muscarinic Receptor Antagonists: Pharmacophore Elucidation and Docking Analyses. , 2016, , 155-179.		0
30	Homologation as a lead modification approach en route to a series of lactone-based muscarinic ligands. Medicinal Chemistry Research, 2014, 23, 1023-1030.	2.4	5
31	Bioisosteric Replacement and Related Analogs in the Design, Synthesis and Evaluation of Ligands for Muscarinic Acetylcholine Receptors. Medicinal Chemistry, 2014, 10, 361-375.	1.5	3
32	Synthesis and preliminary evaluation of affinity to retinoic acid receptors for new organosilicon-based retinoids. Pharmaceutical Chemistry Journal, 2012, 45, 612-621.	0.8	2
33	Modifications to five-substituted 3,3-diethyl-4,5-dihydro-2(3H)-furanones en route to novel muscarinic receptor ligands. Medicinal Chemistry Research, 2011, 20, 558-565.	2.4	6
34	Identification of N-substituted-oxazolidinones as subtype selective 5-HT _{2b} ligands. , 0, , .		O