Hamadeh Tarazi

List of Publications by Year
in descending order

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Version: 2024-02-01

28	513	687363 13	677142 22
papers	citations	h-index	g-index

28	28	28	826
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Safranal induces DNA double-strand breakage and ER-stress-mediated cell death in hepatocellular carcinoma cells. Scientific Reports, 2018, 8, 16951.	3.3	82
2	Camel whey protein hydrolysates induced G2/M cellcycle arrest in human colorectal carcinoma. Scientific Reports, 2021, 11, 7062.	3.3	47
3	Post-Ugi Cascade Transformations for Accessing Diverse Chromenopyrrole Collections. Organic Letters, 2018, 20, 836-839.	4.6	34
4	Inhibition of SHP2 by new compounds induces differential effects on RAS/RAF/ERK and PI3K/AKT pathways in different cancer cell types. Investigational New Drugs, 2019, 37, 252-261.	2.6	27
5	Rational design, synthesis, pharmacophore modeling, and docking studies for identification of novel potent DNA-PK inhibitors. Bioorganic Chemistry, 2017, 72, 234-247.	4.1	26
6	One-Pot Synthesis of Diverse Collections of Benzoxazepine and Indolopyrazine Fused to Heterocyclic Systems. Journal of Organic Chemistry, 2019, 84, 934-948.	3.2	25
7	A Review of HER4 (ErbB4) Kinase, Its Impact on Cancer, and Its Inhibitors. Molecules, 2021, 26, 7376.	3.8	25
8	Design and synthesis of new RAF kinase-inhibiting antiproliferative quinoline derivatives. Part 2: Diarylurea derivatives. European Journal of Medicinal Chemistry, 2017, 127, 413-423.	5.5	23
9	An Efficient Synthesis of biâ€Aryl Pyrimidine Heterocycles: Potential New Drug Candidates to Treat Alzheimer's Disease. Archiv Der Pharmazie, 2017, 350, 1600304.	4.1	22
10	Design, synthesis and SAR analysis of potent BACE1 inhibitors: Possible lead drug candidates for Alzheimer's disease. European Journal of Medicinal Chemistry, 2017, 125, 1213-1224.	5.5	21
11	Design, synthesis and biological evaluation of new pyrrolidine carboxamide analogues as potential chemotherapeutic agents for hepatocellular carcinoma. European Journal of Medicinal Chemistry, 2017, 139, 804-814.	5.5	18
12	Imidazothiazole-based potent inhibitors of V600E-B-RAF kinase with promising anti-melanoma activity: biological and computational studies. Journal of Enzyme Inhibition and Medicinal Chemistry, 2020, 35, 1712-1726.	5.2	17
13	Current Status of Baricitinib as a Repurposed Therapy for COVID-19. Pharmaceuticals, 2021, 14, 680.	3.8	15
14	Conjugation of 4-aminosalicylate with thiazolinones afforded non-cytotoxic potent in vitro and in vivo anti-inflammatory hybrids. Bioorganic Chemistry, 2020, 94, 103378.	4.1	14
15	Design, synthesis, biological evaluation, and modeling studies of novel conformationally-restricted analogues of sorafenib as selective kinase-inhibitory antiproliferative agents against hepatocellular carcinoma cells. European Journal of Medicinal Chemistry, 2021, 210, 113081.	5.5	13
16	Enhancing Student Communication Skills Through Arabic Language Competency and Simulated Patient Assessments. American Journal of Pharmaceutical Education, 2017, 81, 76.	2.1	12
17	Discovery of Novel Small-Molecule Inhibitors of SARS-CoV-2 Main Protease as Potential Leads for COVID-19 Treatment. Journal of Chemical Information and Modeling, 2021, 61, 4745-4757.	5.4	12
18	Discovery of first-in-class imidazothiazole-based potent and selective ErbB4 (HER4) kinase inhibitors. European Journal of Medicinal Chemistry, 2021, 224, 113674.	5.5	11

#	Article	IF	Citations
19	Bis-(5-substituted-2-thiono-1,3,5-thiadiazinan-3-yl) butane as a scaffold of anti-proliferative activity, blended by a multicomponent process. Medicinal Chemistry Research, 2018, 27, 1103-1110.	2.4	10
20	Discovery of a potent p38Î±/MAPK14 kinase inhibitor: Synthesis, inÂvitro/inÂvivo biological evaluation, and docking studies. European Journal of Medicinal Chemistry, 2019, 183, 111684.	5.5	10
21	Design, Synthesis and Qualitative Structure Activity Relationship Evaluations of Quinoline-Based Bisarylimidazoles as Antibacterial Motifs. Medicinal Chemistry, 2016, 12, 563-573.	1.5	10
22	Discovery of highly potent V600E-B-RAF kinase inhibitors: Molecular modeling study. Bioorganic and Medicinal Chemistry, 2019, 27, 655-663.	3.0	8
23	Molecular docking and glucosidase inhibition studies of novel N-arylthiazole-2-amines and Ethyl 2-[aryl(thiazol-2-yl)amino]acetates. Medicinal Chemistry Research, 2017, 26, 3247-3261.	2.4	7
24	Design and synthesis of new quinoline derivatives as selective C-RAF kinase inhibitors with potent anticancer activity. European Journal of Medicinal Chemistry, 2022, 238, 114434.	5.5	7
25	In-silico screening for DNA-dependent protein kinase (DNA-PK) inhibitors: Combined homology modeling, docking, molecular dynamic study followed by biological investigation. Biomedicine and Pharmacotherapy, 2016, 83, 693-703.	5.6	6
26	Design and synthesis of a new series of highly potent RAF kinase-inhibiting triarylpyrazole derivatives possessing antiproliferative activity against melanoma cells. Future Medicinal Chemistry, 2016, 8, 2197-2211.	2.3	6
27	Molecular properties prediction, synthesis, and diuretic activity of phenoxy acetic acid bearing pyrazolines. Medicinal Chemistry Research, 2013, 22, 916-926.	2.4	3
28	Peptide Derivatives of 1,2-Dihydro-3-Methyl-2-Oxoquinoxaline-6-Carboxylic Acid: Synthesis and Evaluation of Antimicrobial, Antifungal and Antiviral Potential. Pharmaceutical Chemistry Journal, 2016, 50, 331-338.	0.8	2