Sotirios Katsamakas

List of Publications by Year in descending order

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933447 677142 26 504 10 22 citations g-index h-index papers 27 27 27 935 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	Discovery of Benzothiazole Scaffold-Based DNA Gyrase B Inhibitors. Journal of Medicinal Chemistry, 2016, 59, 8941-8954.	6.4	99
2	Discovery of 4,5,6,7-Tetrahydrobenzo[1,2- <i>d</i>]thiazoles as Novel DNA Gyrase Inhibitors Targeting the ATP-Binding Site. Journal of Medicinal Chemistry, 2015, 58, 5501-5521.	6.4	92
3	RGD-mediated delivery of small-molecule drugs. Future Medicinal Chemistry, 2017, 9, 579-604.	2.3	61
4	Probing the Inhibition of Microtubule Affinity Regulating Kinase 4 by N-Substituted Acridones. Scientific Reports, 2019, 9, 1676.	3.3	49
5	Synthesis and biological evaluation of a Platinum(II)-c(RGDyK) conjugate for integrin-targeted photodynamic therapy. European Journal of Medicinal Chemistry, 2017, 141, 221-231.	5.5	38
6	Advances of Phenoxazines: Synthesis, Reactivity and Their Medicinal Applications. Current Medicinal Chemistry, 2016, 23, 2972-2999.	2.4	18
7	Synthesis and evaluation of gallocyanine dyes as potential agents for the treatment of Alzheimer's disease and related neurodegenerative tauopathies. European Journal of Medicinal Chemistry, 2016, 108, 28-38.	5.5	17
8	Synthesis, structure elucidation and biological evaluation of triple bridged dinuclear copper(II) complexes as anticancer and antioxidant/anti-inflammatory agents. Materials Science and Engineering C, 2017, 76, 1026-1040.	7.3	16
9	Interaction of Fe(III) with herbicide-carboxylato ligands – Di-, tri- and tetra-nuclear compounds: Structure and magnetic behavior. Polyhedron, 2007, 26, 763-772.	2.2	15
10	Design and synthesis of gallocyanine inhibitors of DKK1/LRP6 interactions for treatment of Alzheimer's disease. Bioorganic Chemistry, 2018, 80, 230-244.	4.1	15
11	Boronic Acid Group: A Cumbersome False Negative Case in the Process of Drug Design. Molecules, 2016, 21, 1185.	3.8	10
12	Novel c(RGDyK)-based conjugates of POPAM and 5-fluorouracil for integrin-targeted cancer therapy. Future Medicinal Chemistry, 2017, 9, 2181-2196.	2.3	10
13	Mononuclear copper(II) complexes with 2-thiophene carboxylate and N-N donors; DNA interaction, antioxidant/anti-inflammatory and antitumor activity. Materials Science and Engineering C, 2019, 94, 493-508.	7.3	9
14	Sulfanyl 5H-dihydro-pyrrole derivatives via 1,3-dipolar cycloaddition, their further chemical manipulation and antioxidant activity. Arkivoc, 2015, 2015, 214-231.	0.5	9
15	Examining barbiturate scaffold for the synthesis of new agents with biological interest. Future Medicinal Chemistry, 2019, 11, 2063-2079.	2.3	8
16	Considering Autotaxin Inhibitors in Terms of 2D-QSAR and 3D-Mapping-Review and Evaluation. Current Medicinal Chemistry, 2015, 22, 1428-1461.	2.4	8
17	Targeting on poly(ADPâ€ribose) polymerase activity with DNAâ€damaging hybrid lactamâ€steroid alkylators in wildâ€type and BRCA1â€mutated ovarian cancer cells. Chemical Biology and Drug Design, 2017, 90, 854-866.	3.2	7
18	Discovery of novel phenoxazinone derivatives as DKK1/LRP6 interaction inhibitors: Synthesis, biological evaluation and structure–activity relationships. Bioorganic and Medicinal Chemistry, 2016, 24, 1014-1022.	3.0	6

#	Article	IF	CITATIONS
19	\hat{l} ±-Amination and the 5-exo-trig cyclization reaction of sulfur-containing Schiff bases with N-phenyltriazolinedione and their anti-lipid peroxidation activity. Comptes Rendus Chimie, 2017, 20, 424-434.	0.5	6
20	Computational Design of Multitarget Drugs Against Alzheimer's Disease. Methods in Pharmacology and Toxicology, 2018, , 203-253.	0.2	2
21	Boronic Acid Based Inhibitors of Autotaxin: Understanding their Biological Role in Terms of Quantitative Structure Activity Relationships (QSAR). Letters in Drug Design and Discovery, 2013, 10, 11-18.	0.7	2
22	A novel synthetic protocol for the synthesis of <i>pulvinones,</i> and naturally occurring <i>Aspulvinone E</i> , molecules of medicinal interest. Synthetic Communications, 2022, 52, 117-128.	2.1	2
23	Design and Synthesis of Novel Bis-Imidazolyl Phenyl Butadiyne Derivatives as HCV NS5A Inhibitors. Pharmaceuticals, 2022, 15, 632.	3.8	2
24	1-Methyl-8-phenyl-1,3-diazaspiro[4.5]decane-2,4-dione. MolBank, 2021, 2021, M1228.	0.5	1
25	Redesigning of the cap conformation and symmetry of the diphenylethyne core to yield highly potent pan-genotypic NS5A inhibitors with high potency and high resistance barrier. European Journal of Medicinal Chemistry, 2022, 229, 114034.	5.5	1
26	Deep learning-assisted pipeline for Virtual Screening of ligand compound databases: Application on inhibiting the entry of SARS-CoV-2 into human cells. , 2020, , .		1