

Tahsin F Kellici

List of Publications by Year in descending order

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Version: 2024-02-01

37
papers

556
citations

623574

14
h-index

642610

23
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38
all docs

38
docs citations

38
times ranked

848
citing authors

#	ARTICLE	IF	CITATIONS
1	Fighting COVID-19 with Artificial Intelligence. <i>Methods in Molecular Biology</i> , 2022, 2390, 103-112.	0.4	7
2	Structure–function analysis of naturally occurring apolipoprotein A-I L144R, A164S and L178P mutants provides insight on their role on HDL levels and cardiovascular risk. <i>Cellular and Molecular Life Sciences</i> , 2021, 78, 1523-1544.	2.4	8
3	Small-molecule modulators of serine protease inhibitor proteins (serpins). <i>Drug Discovery Today</i> , 2021, 26, 442-454.	3.2	13
4	Therapeutic Potential of Targeting Plasminogen Activator Inhibitor-1 in COVID-19. <i>Trends in Pharmacological Sciences</i> , 2021, 42, 431-433.	4.0	20
5	Synthesis, biology, computational studies and <i>in vitro</i> controlled release of new isoniazid-based adamantane derivatives. <i>Future Medicinal Chemistry</i> , 2019, 11, 2779-2802.	1.1	4
6	Host–Guest Interactions between Candesartan and Its Prodrug Candesartan Cilexetil in Complex with 2-Hydroxypropyl- β -cyclodextrin: On the Biological Potency for Angiotensin II Antagonism. <i>Molecular Pharmaceutics</i> , 2019, 16, 1255-1271.	2.3	17
7	Drug-Membrane Interactions in the Renin Angiotensin System. <i>Series in Bioengineering</i> , 2019, , 339-364.	0.3	1
8	Inclusion of Quercetin in Gold Nanoparticles Decorated with Supramolecular Hosts Amplifies Its Tumor Targeting Properties. <i>ACS Applied Bio Materials</i> , 2019, 2, 2715-2725.	2.3	30
9	The dynamic properties of angiotensin II type 1 receptor inverse agonists in solution and in the receptor site. <i>Arabian Journal of Chemistry</i> , 2019, 12, 5062-5078.	2.3	19
10	Molecular Dynamics Simulations on the Bioactive Molecule of hIAPP22–29 (NFGAILSS) and Rational Drug Design. <i>Methods in Molecular Biology</i> , 2018, 1824, 1-16.	0.4	0
11	Designing Natural Product Hybrids Bearing Triple Antiplatelet Profile and Evaluating Their Human Plasma Stability. <i>Methods in Molecular Biology</i> , 2018, 1824, 371-385.	0.4	4
12	Angiotensin II Type 1 Receptor Homology Models: A Comparison Between In Silico and the Crystal Structures. <i>Methods in Molecular Biology</i> , 2018, 1824, 449-460.	0.4	1
13	Three Regioselectively Acylated Flavonoid Aglycone Derivatives in Equimolar Yield at One Blow. <i>ChemistrySelect</i> , 2018, 3, 5207-5211.	0.7	2
14	Exploring the interactions of irbesartan and irbesartan–2-hydroxypropyl- β -cyclodextrin complex with model membranes. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2017, 1859, 1089-1098.	1.4	26
15	Molecular requirements involving the human platelet protease-activated receptor-4 mechanism of activation by peptide analogues of its tethered-ligand. <i>Platelets</i> , 2017, 28, 812-821.	1.1	13
16	In vitro Controlled Release from Solid Pharmaceutical Formulations of two new Adamantane Aminoethers with Antitubercular Activity (I).. <i>Drug Research</i> , 2017, 67, 447-450.	0.7	8
17	Crystal structure analysis, covalent docking, and molecular dynamics calculations reveal a conformational switch in PhaZ7 PHB depolymerase. <i>Proteins: Structure, Function and Bioinformatics</i> , 2017, 85, 1351-1361.	1.5	7
18	Tailoring naringenin conjugates with amplified and triple antiplatelet activity profile: Rational design, synthesis, human plasma stability and in vitro evaluation. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2017, 1861, 2609-2618.	1.1	13

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19	Rational design and structure-activity relationship studies of quercetin-amino acid hybrids targeting the anti-apoptotic protein Bcl-xL. <i>Organic and Biomolecular Chemistry</i> , 2017, 15, 7956-7976.	1.5	24
20	In vitro Controlled Release of two new Tuberculocidal Adamantane Aminoethers from Solid Pharmaceutical Formulations (II). <i>Drug Research</i> , 2017, 67, 653-660.	0.7	7
21	A combined NMR and molecular dynamics simulation study to determine the conformational properties of rat/mouse 35-55 myelin oligodendrocyte glycoprotein epitope implicated in the induction of experimental autoimmune encephalomyelitis. <i>Journal of Biomolecular Structure and Dynamics</i> , 2017, 35, 1559-1567.	2.0	3
22	The use of J-coupling as a sole criterion to assign the total absolute stereochemistry of new pyrrolidinone class synthetic analogs, derived from S-pyroglyutamic acid. <i>Journal of Molecular Structure</i> , 2017, 1129, 195-199.	1.8	0
23	Mapping the interactions and bioactivity of quercetin-(2-hydroxypropyl)- β -cyclodextrin complex. <i>International Journal of Pharmaceutics</i> , 2016, 511, 303-311.	2.6	48
24	PRESS: PRotEin S-Sulfenylation server. <i>Bioinformatics</i> , 2016, 32, 2710-2712.	1.8	19
25	New hydrazones of 5-nitro-2-furaldehyde with adamantanealkanohydrazides: synthesis and in vitro trypanocidal activity. <i>MedChemComm</i> , 2016, 7, 1229-1236.	3.5	12
26	Deconvoluting the Dual Antiplatelet Activity of a Plant Extract. <i>Journal of Agricultural and Food Chemistry</i> , 2016, 64, 4511-4521.	2.4	13
27	Antiplatelet effect of the main triterpenoids of an olive leaf extract. <i>Atherosclerosis</i> , 2016, 252, e98.	0.4	1
28	Calixarenes in Lipase Biocatalysis and Cancer Therapy. <i>Current Organic Chemistry</i> , 2016, 20, 1043-1057.	0.9	23
29	Leveraging NMR and X-ray Data of the Free Ligands to Build Better Drugs Targeting Angiotensin II Type 1 G-Protein Coupled Receptor. <i>Current Medicinal Chemistry</i> , 2015, 23, 36-59.	1.2	20
30	Rational Drug Design and Synthesis of Molecules Targeting the Angiotensin II Type 1 and Type 2 Receptors. <i>Molecules</i> , 2015, 20, 3868-3897.	1.7	36
31	Rosemary tea consumption results to anxiolytic- and anti-depressant-like behavior of adult male mice and inhibits all cerebral area and liver cholinesterase activity; phytochemical investigation and in silico studies. <i>Chemico-Biological Interactions</i> , 2015, 237, 47-57.	1.7	48
32	Investigation of the Interactions of Silibinin with 2-Hydroxypropyl- β -cyclodextrin through Biophysical Techniques and Computational Methods. <i>Molecular Pharmaceutics</i> , 2015, 12, 954-965.	2.3	55
33	Branched-chain sugar nucleosides: stereocontrolled synthesis and bioevaluation of novel 3-C-trifluoromethyl and 3-C-methyl pyranonucleosides. <i>Carbohydrate Research</i> , 2015, 407, 170-178.	1.1	2
34	A novel synthetic luteinizing hormone-releasing hormone (LHRH) analogue coupled with modified β -cyclodextrin: Insight into its intramolecular interactions. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2015, 1850, 159-168.	1.1	8
35	Pharmaceutical compositions for antihypertensive treatments: a patent review. <i>Expert Opinion on Therapeutic Patents</i> , 2015, 25, 1305-17.	2.4	20
36	Rational Drug Design Paradigms: The Odyssey for Designing Better Drugs. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2015, 18, 238-256.	0.6	8

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37	The application of solid-state NMR spectroscopy to study candesartan cilexetil (TCV-116) membrane interactions. Comparative study with the AT1R antagonist drug olmesartan. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2014, 1838, 2439-2450.	1.4	16