Giuseppe Zagotto

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53 681 16 22 g-index

54 787 4 4.12 ext. papers ext. citations avg, IF L-index

#	Paper	IF	Citations
53	Aminoacyl-anthraquinone conjugates as telomerase inhibitors: synthesis, biophysical and biological evaluation. <i>Journal of Medicinal Chemistry</i> , 2008 , 51, 5566-74	8.3	50
52	Mechanistic Insight into the Oxidation of Organic Phenylselenides by H O. <i>Chemistry - A European Journal</i> , 2017 , 23, 2405-2422	4.8	41
51	Peptidyl anthraquinones as potential antineoplastic drugs: synthesis, DNA binding, redox cycling, and biological activity. <i>Journal of Medicinal Chemistry</i> , 1996 , 39, 3114-22	8.3	39
50	Tuning G-quadruplex vs double-stranded DNA recognition in regioisomeric lysyl-peptidyl-anthraquinone conjugates. <i>Bioconjugate Chemistry</i> , 2011 , 22, 2126-35	6.3	34
49	New Therapeutic Applications of Phosphodiesterase 5 Inhibitors (PDE5-Is). <i>Current Medicinal Chemistry</i> , 2016 , 23, 1239-49	4.3	27
48	Mapping drug interactions at the covalent topoisomerase II-DNA complex by bisantrene/amsacrine congeners. <i>Journal of Biological Chemistry</i> , 1998 , 273, 12732-9	5.4	26
47	Sequence-specific interactions of drugs interfering with the topoisomerase-DNA cleavage complex. <i>Biochimica Et Biophysica Acta - Molecular Basis of Disease</i> , 2002 , 1587, 145-54	6.9	23
46	Development of DNA topoisomerase-related therapeutics: a short perspective of new challenges. <i>Anti-Cancer Agents in Medicinal Chemistry</i> , 2004 , 4, 335-45		23
45	The Old Made New: Natural Compounds against Erectile Dysfunction. <i>Archiv Der Pharmazie</i> , 2015 , 348, 607-14	4.3	22
44	An Overview of New Possible Treatments of Alzheimer Disease, Based on Natural Products and Semi-Synthetic Compounds. <i>Current Medicinal Chemistry</i> , 2017 , 24, 3749-3773	4.3	21
43	The Medicinal Chemistry of Natural and Semisynthetic Compounds against Parkinson's and Huntington's Diseases. <i>ACS Chemical Neuroscience</i> , 2017 , 8, 2356-2368	5.7	19
42	Antioxidant Potential of Psychotropic Drugs: From Clinical Evidence to In Vitro and In Vivo Assessment and toward a New Challenge for in Silico Molecular Design. <i>Antioxidants</i> , 2020 , 9,	7.1	19
41	Major Depressive Disorder and Oxidative Stress: In Silico Investigation of Fluoxetine Activity against ROS. <i>Applied Sciences (Switzerland)</i> , 2019 , 9, 3631	2.6	18
40	Psychiatric Disorders and Oxidative Injury: Antioxidant Effects of Zolpidem Therapy disclosed. Computational and Structural Biotechnology Journal, 2019, 17, 311-318	6.8	18
39	Semi-synthetic derivatives of natural isoflavones from Maclura pomifera as a novel class of PDE-5A inhibitors. <i>Floterap</i> []2015 , 105, 132-8	3.2	16
38	New naphthoquinone derivatives against glioma cells. <i>European Journal of Medicinal Chemistry</i> , 2015 , 96, 458-66	6.8	16
37	Remarkable interference with telomeric function by a G-quadruplex selective bisantrene regioisomer. <i>Biochemical Pharmacology</i> , 2010 , 79, 1781-90	6	16

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36	Preliminary studies of berberine and its semi-synthetic derivatives as a promising class of multi-target anti-parkinson agents. <i>Natural Product Research</i> , 2018 , 32, 1395-1401	2.3	15
35	Semi-synthetic isoflavones as BACE-1 inhibitors against Alzheimer u disease. <i>Bioorganic Chemistry</i> , 2019 , 87, 474-483	5.1	14
34	Rational design, synthesis, and DNA binding properties of novel sequence-selective peptidyl congeners of ametantrone. <i>ChemMedChem</i> , 2010 , 5, 1080-91	3.7	13
33	Therapeutic Potential of Phosphodiesterase Inhibitors against Neurodegeneration: The Perspective of the Medicinal Chemist. <i>ACS Chemical Neuroscience</i> , 2020 , 11, 1726-1739	5.7	12
32	Antiproliferative activity of Juglone derivatives on rat glioma. <i>Natural Product Research</i> , 2017 , 31, 632-	63 83	12
31	Electron paramagnetic resonance (EPR) study of spin-labeled camptothecin derivatives: a different look of the ternary complex. <i>Journal of Medicinal Chemistry</i> , 2011 , 54, 1003-9	8.3	12
30	DNA-binding preferences of bisantrene analogues: relevance to the sequence specificity of drug-mediated topoisomerase II poisoning. <i>Molecular Pharmacology</i> , 1998 , 54, 1036-45	4.3	12
29	Pharmacophore-guided discovery of CDC25 inhibitors causing cell cycle arrest and tumor regression. <i>Scientific Reports</i> , 2019 , 9, 1335	4.9	11
28	Fluoxetine scaffold to design tandem molecular antioxidants and green catalysts <i>RSC Advances</i> , 2020 , 10, 18583-18593	3.7	11
27	Constrained bisantrene derivatives as G-quadruplex binders. <i>Arkivoc</i> , 2016 , 2016, 145-160	0.9	10
26	Natural phosphodiesterase 5 (PDE5) inhibitors: a computational approach. <i>Natural Product Research</i> , 2021 , 35, 1648-1653	2.3	10
25	Plant natural products with anti-thyroid cancer activity. Floterap[]2020, 146, 104640	3.2	9
24	A novel class of selective CK2 inhibitors targeting its open hinge conformation. <i>European Journal of Medicinal Chemistry</i> , 2020 , 195, 112267	6.8	8
23	Synthesis and Evaluation of New Naphthalene and Naphthoquinone Derivatives as Anticancer Agents. <i>Archiv Der Pharmazie</i> , 2017 , 350, e1600286	4.3	7
22	Overcoming resistance in non-small-cell lung cancer: A practical lesson for the medicinal chemist. <i>Archiv Der Pharmazie</i> , 2018 , 351, e1800037	4.3	7
21	Evidence on selective binding to G-quadruplex DNA of isoflavones from by mass spectrometry and molecular docking. <i>Natural Product Research</i> , 2021 , 35, 2583-2587	2.3	7
20	Novel ametantrone-amsacrine related hybrids as topoisomerase IIIpoisons and cytotoxic agents. <i>Archiv Der Pharmazie</i> , 2014 , 347, 728-37	4.3	7
19	Synthesis, DNA-damaging and cytotoxic properties of novel topoisomerase II-directed bisantrene analogues. <i>Bioorganic and Medicinal Chemistry Letters</i> , 1998 , 8, 121-6	2.9	7

A versatile synthesis of the 1,4-dihydroxynaphthoquinone nucleus. Tetrahedron Letters, 2000, 41, 6631-6634 18 Synthesis via A3 Coupling Reaction of Anthracene-Propargylamine as a New Scaffold for the 1.8 17 7 Interaction with DNA. ChemistrySelect, 2019, 4, 13138-13142 5-Hydroxy-3-(4-hydroxyphenyl)-8,8-dimethyl-6-(3-methylbut-2-enyl)pyrano[2,3-h]chromen-4-one. 16 0.5 7 MolBank, 2018, 2018, M1004 Investigation of the molecular reactivity of bioactive oxiranylmethyloxy anthraquinones. Archiv Der 6 15 4.3 Pharmazie, 2019, 352, e1900030 Preferred interaction of D-peptidyl-anthraquinones with double-stranded B-DNA. International 6 7.9 14 Journal of Biological Macromolecules, 1997, 21, 319-26 Scouting Novel Protein Kinase A (PKA) Inhibitors by Using a Consensus Docking-Based Virtual 0.8 6 13 Screening Approach. Letters in Drug Design and Discovery, 2009, 6, 327-336 2-(3,4-Dihydroxyphenyl)-4-(2-(4-nitrophenyl)hydrazono)-4H-chromene-3,5,7-triol. MolBank, 2020, 12 0.5 5 2020, M1144 ROS-Scavenging Selenofluoxetine Derivatives Inhibit Serotonin Reuptake.. ACS Omega, 2022, 7, 8314-8322 11 8-Hydroxynaphthalene-1,4-dione derivative as novel compound for glioma treatment. Bioorganic 10 2.9 4 and Medicinal Chemistry Letters, 2011, 21, 2079-82 Natural Compounds Promoting Weight Loss: Mechanistic Insights from the Point of View of the 0.6 9 4 Medicinal Chemist. Natural Products Journal, 2019, 9, 78-85 Photoactivated semi-synthetic derivative of osajin selectively interacts with G-quadruplex DNA. 8 2.3 3 Natural Product Research, 2020, 1-6 Combinatorial library generation, molecular docking and molecular dynamics simulations for enhancing the isoflavone scaffold in phosphodiesterase inhibition. New Journal of Chemistry, 2020, 3.6 44, 19472-19488 Enhanced G-quadruplex selectivity of flavonoid glycoside rutin over quercetin. Natural Product 2.3 2 Research, 2020, 1-5 Design and synthesis of a peptide derivative of ametantrone targeting the major groove of the 3.6 2 d(GGCGCC)2 palindromic sequence. New Journal of Chemistry, 2020, 44, 3624-3631 Selenoxide Elimination Triggers Enamine Hydrolysis to Primary and Secondary Amines: A Combined 4.8 1 Experimental and Theoretical Investigation. Molecules, 2021, 26, Photo-induced spin switching in a modified anthraguinone modulated by DNA binding. 1 4.2 Photochemical and Photobiological Sciences, 2019, 18, 2199-2207 Synthesis and biological evaluation of heteroalicyclic cyanoguanidines at histamine receptors. 4.3 O *Archiv Der Pharmazie*, **2019**, 352, e1900107 A new sensitive and subunit-selective molecular tool for investigating protein kinase A in the brain. 4.3 *Archiv Der Pharmazie*, **2020**, 353, e1900326