

Marino Convertino

List of Publications by Year in descending order

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Version: 2024-02-01

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papers

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516710

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1703
citing authors

#	ARTICLE	IF	CITATIONS
1	Identification and characterization of novel candidate compounds targeting μ - and δ -transmembrane μ -opioid receptor isoforms. <i>British Journal of Pharmacology</i> , 2021, 178, 2709-2726.	5.4	4
2	Stabilization of μ -opioid receptor facilitates its cellular translocation and signaling. <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 878-884.	2.6	6
3	Discovery of Novel Adenosine Receptor Antagonists through a Combined Structure- and Ligand-Based Approach Followed by Molecular Dynamics Investigation of Ligand Binding Mode. <i>Journal of Chemical Information and Modeling</i> , 2018, 58, 794-815.	5.4	22
4	Reversible and Tunable Photoswitching of Protein Function through Genetic Encoding of Azobenzene Amino Acids in Mammalian Cells. <i>ChemBioChem</i> , 2018, 19, 2178-2185.	2.6	40
5	miRNA-711 Binds and Activates TRPA1 Extracellularly to Evoke Acute and Chronic Pruritus. <i>Neuron</i> , 2018, 99, 449-463.e6.	8.1	79
6	Molecular Mechanisms of the R61T Mutation in Apolipoprotein E4: A Dynamic Rescue. <i>Biophysical Journal</i> , 2017, 113, 2192-2198.	0.5	2
7	Reducing the Flexibility of Type II Dehydroquinase for Inhibition: A Fragment-Based Approach and Molecular Dynamics Study. <i>ChemMedChem</i> , 2017, 12, 1512-1524.	3.2	4
8	Structural and functional interactions between six-transmembrane μ -opioid receptors and β 2-adrenoreceptors modulate opioid signaling. <i>Scientific Reports</i> , 2016, 5, 18198.	3.3	34
9	Docking and Scoring with Target-Specific Pose Classifier Succeeds in Native-Like Pose Identification But Not Binding Affinity Prediction in the CSAR 2014 Benchmark Exercise. <i>Journal of Chemical Information and Modeling</i> , 2016, 56, 1032-1041.	5.4	15
10	Pharmacological Chaperones: Design and Development of New Therapeutic Strategies for the Treatment of Conformational Diseases. <i>ACS Chemical Biology</i> , 2016, 11, 1471-1489.	3.4	85
11	Computational Modeling of Small Molecule Ligand Binding Interactions and Affinities. <i>Methods in Molecular Biology</i> , 2016, 1414, 23-32.	0.9	4
12	The Molecular Basis for Dual Fatty Acid Amide Hydrolase (FAAH)/Cyclooxygenase (COX) Inhibition. <i>ChemMedChem</i> , 2016, 11, 1252-1258.	3.2	33
13	ApoE4-specific Misfolded Intermediate Identified by Molecular Dynamics Simulations. <i>PLoS Computational Biology</i> , 2015, 11, e1004359.	3.2	21
14	Differential Regulation of 6- and 7-Transmembrane Helix Variants of μ -Opioid Receptor in Response to Morphine Stimulation. <i>PLoS ONE</i> , 2015, 10, e0142826.	2.5	14
15	Differences in the Antinociceptive Effects and Binding Properties of Propranolol and Bupranolol Enantiomers. <i>Journal of Pain</i> , 2015, 16, 1321-1333.	1.4	27
16	An optimized polyamine moiety boosts the potency of human type II topoisomerase poisons as quantified by comparative analysis centered on the clinical candidate F14512. <i>Chemical Communications</i> , 2015, 51, 14310-14313.	4.1	32
17	μ -Opioid receptor 6-transmembrane isoform: A potential therapeutic target for new effective opioids. <i>Progress in Neuro-Psychopharmacology and Biological Psychiatry</i> , 2015, 62, 61-67.	4.8	26
18	Carnosine Inhibits Al^{2+} Aggregation by Perturbing the H-Bond Network in and around the Central Hydrophobic Cluster. <i>ChemBioChem</i> , 2013, 14, 583-592.	2.6	76

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19	Methylations of Tryptophan-Modified Naphthoquinone Affect Its Inhibitory Potential toward A β Aggregation. <i>Journal of Physical Chemistry B</i> , 2013, 117, 1780-1789.	2.6	16
20	Structural Basis for Inhibiting β -Amyloid Oligomerization by a Non-coded β -Breaker-Substituted Endomorphin Analogue. <i>ACS Chemical Biology</i> , 2011, 6, 1265-1276.	3.4	32
21	Disordered Binding of Small Molecules to A β (12-28). <i>Journal of Biological Chemistry</i> , 2011, 286, 41578-41588.	3.4	46
22	Complete Phenotypic Recovery of an Alzheimer's Disease Model by a Quinone-Tryptophan Hybrid Aggregation Inhibitor. <i>PLoS ONE</i> , 2010, 5, e11101.	2.5	129
23	9,10-Anthraquinone hinders β aggregation: How does a small molecule interfere with A β peptide amyloid fibrillation?. <i>Protein Science</i> , 2009, 18, 792-800.	7.6	91