

Andrea

List of Publications by Year in descending order

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Version: 2024-02-01

24
papers

396
citations

840776

11
h-index

752698

20
g-index

25
all docs

25
docs citations

25
times ranked

588
citing authors

#	ARTICLE	IF	CITATIONS
1	Synthesis, Biological Evaluation, and Computational Studies of Tri- and Tetracyclic Nitrogen-Bridgehead Compounds as Potent Dual-Acting AChE Inhibitors and H_3 Receptor Antagonists. <i>ACS Chemical Neuroscience</i> , 2014, 5, 225-242.	3.5	67
2	Structure-Activity Relationships and Computational Investigations into the Development of Potent and Balanced Dual-Acting Butyrylcholinesterase Inhibitors and Human Cannabinoid Receptor 2 Ligands with Pro-Cognitive in Vivo Profiles. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 1646-1663.	6.4	50
3	Pharmacological Profile of Histaprodifens at Four Recombinant Histamine H_1 Receptor Species Isoforms. <i>Journal of Pharmacology and Experimental Therapeutics</i> , 2008, 324, 60-71.	2.5	47
4	Aminobenzimidazoles and Structural Isomers as Templates for Dual-Acting Butyrylcholinesterase Inhibitors and CB_2 Receptor Ligands To Combat Neurodegenerative Disorders. <i>ChemMedChem</i> , 2016, 11, 1270-1283.	3.2	28
5	Comparison of the pharmacological properties of human and rat histamine H_3 -receptors. <i>Biochemical Pharmacology</i> , 2010, 80, 1437-1449.	4.4	27
6	Striking differences of action of lisuride stereoisomers at histamine H_1 receptors. <i>Naunyn-Schmiedeberg's Archives of Pharmacology</i> , 2006, 374, 215-222.	3.0	20
7	The First Photochromic Affinity Switch for the Human Cannabinoid Receptor 2. <i>Advanced Therapeutics</i> , 2018, 1, 1700032.	3.2	20
8	A Dynamic, Split-Luciferase-Based Mini-G Protein Sensor to Functionally Characterize Ligands at All Four Histamine Receptor Subtypes. <i>International Journal of Molecular Sciences</i> , 2020, 21, 8440.	4.1	20
9	Mathematical analysis of the sodium sensitivity of the human histamine H_3 receptor. <i>In Silico Pharmacology</i> , 2014, 2, 1.	3.3	16
10	Analysis of the activation mechanism of the guinea-pig Histamine H_1 -receptor. <i>Journal of Computer-Aided Molecular Design</i> , 2007, 21, 499-509.	2.9	13
11	Binding pathway of histamine to the hH_4R , observed by unconstrained molecular dynamics. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2015, 25, 1259-1268.	2.2	12
12	Dibenzo[b , f][1,4]oxazepines and dibenzo[b , e]oxepines: Influence of the chlorine substitution pattern on the pharmacology at the H_1 R, H_4 R, 5-HT $2A$ R and other selected GPCRs. <i>Pharmacological Research</i> , 2016, 113, 610-625.	7.1	11
13	2,4-Diaminopyrimidines as dual ligands at the histamine H_1 and H_4 receptor H_1/H_4 -receptor selectivity. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2016, 26, 292-300.	2.2	11
14	Conformational Restriction and Enantioseparation Increase Potency and Selectivity of Cyanoguanidine-Type Histamine H_4 Receptor Agonists. <i>Journal of Medicinal Chemistry</i> , 2016, 59, 3452-3470.	6.4	9
15	Abolishing Dopamine D_{2L} / D_3 Receptor Affinity of Subtype-Selective Carbamoylguanidine-Type Histamine H_2 Receptor Agonists. <i>Journal of Medicinal Chemistry</i> , 2021, 64, 8684-8709.	6.4	8
16	Label-Free Investigations on the G Protein Dependent Signaling Pathways of Histamine Receptors. <i>International Journal of Molecular Sciences</i> , 2021, 22, 9739.	4.1	7
17	LigPath: a module for predictive calculation of a ligand's pathway into a receptor-application to the gpH_1 - receptor. <i>Journal of Molecular Modeling</i> , 2006, 13, 209-218.	1.8	6
18	Distinct interactions between the human adrenergic β_2 receptor and $G_{i/s}$ in silico study. <i>Journal of Molecular Modeling</i> , 2010, 16, 1307-1318.	1.8	6

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19	Competitive association binding kinetic assays: a new tool to detect two different binding orientations of a ligand to its target protein under distinct conditions?. <i>Naunyn-Schmiedeberg's Archives of Pharmacology</i> , 2017, 390, 595-612.	3.0	5
20	Specific Engineered G Protein Coupling to Histamine Receptors Revealed from Cellular Assay Experiments and Accelerated Molecular Dynamics Simulations. <i>International Journal of Molecular Sciences</i> , 2021, 22, 10047.	4.1	4
21	In silico analysis of the histaprodifen induced activation pathway of the guinea-pig histamine H1-receptor. <i>Journal of Computer-Aided Molecular Design</i> , 2010, 24, 759-769.	2.9	3
22	Structure-Dependent Deconjugation of Flavonoid Glucuronides by Human β -Glucuronidase " In Vitro and In Silico Analyses. <i>Planta Medica</i> , 2015, 81, 1182-1189.	1.3	3
23	Molecular Modelling Approaches for the Analysis of Histamine Receptors and Their Interaction with Ligands. <i>Handbook of Experimental Pharmacology</i> , 2017, 241, 31-61.	1.8	3
24	Shining light on the histamine H2 receptor: Synthesis of carbamoylguanidine-type agonists as a pharmacological tool to study internalization. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2021, 52, 128388.	2.2	0