Kenneth A Jacobson

List of Publications by Year
in descending order

774	42,097	2963	6282
papers	citations	93 h-index	158 g-index

817	817	817	23289
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Adenosine receptors as therapeutic targets. Nature Reviews Drug Discovery, 2006, 5, 247-264.	21.5	1,243
2	International Union of Pharmacology LVIII: Update on the P2Y G Protein-Coupled Nucleotide Receptors: From Molecular Mechanisms and Pathophysiology to Therapy. Pharmacological Reviews, 2006, 58, 281-341.	7.1	1,147
3	International Union of Basic and Clinical Pharmacology. LXXXI. Nomenclature and Classification of Adenosine Receptorsâ€"An Update. Pharmacological Reviews, 2011, 63, 1-34.	7.1	1,135
4	Structure of an Agonist-Bound Human A _{2A} Adenosine Receptor. Science, 2011, 332, 322-327.	6.0	783
5	UDP acting at P2Y6 receptors is a mediator of microglial phagocytosis. Nature, 2007, 446, 1091-1095.	13.7	698
6	THE CONCISE GUIDE TO PHARMACOLOGY 2019/20: G proteinâ€coupled receptors. British Journal of Pharmacology, 2019, 176, S21-S141.	2.7	519
7	The Concise Guide to PHARMACOLOGY 2015/16: G proteinâ€coupled receptors. British Journal of Pharmacology, 2015, 172, 5744-5869.	2.7	507
8	Adenosine receptors: pharmacology, structure-activity relationships, and therapeutic potential. Journal of Medicinal Chemistry, 1992, 35, 407-422.	2.9	488
9	Coordinated Adenine Nucleotide Phosphohydrolysis and Nucleoside Signaling in Posthypoxic Endothelium. Journal of Experimental Medicine, 2003, 198, 783-796.	4.2	444
10	Characterization of the UDP-glucose receptor (re-named here the P2Y14 receptor) adds diversity to the P2Y receptor family. Trends in Pharmacological Sciences, 2003, 24, 52-55.	4.0	382
11	Recent developments in adenosine receptor ligands and their potential as novel drugs. Biochimica Et Biophysica Acta - Biomembranes, 2011, 1808, 1290-1308.	1.4	375
12	THE CONCISE GUIDE TO PHARMACOLOGY 2021/22: G proteinâ€coupled receptors. British Journal of Pharmacology, 2021, 178, S27-S156.	2.7	337
13	Purine and Pyrimidine (P2) Receptors as Drug Targets. Journal of Medicinal Chemistry, 2002, 45, 4057-4093.	2.9	334
14	Structure of the human P2Y12 receptor in complex with an antithrombotic drug. Nature, 2014, 509, 115-118.	13.7	330
15	Towards a revised nomenclature for P1 and P2 receptors. Trends in Pharmacological Sciences, 1997, 18, 79-82.	4.0	315
16	Two disparate ligand-binding sites in the human P2Y1 receptor. Nature, 2015, 520, 317-321.	13.7	305
17	Adenosine A3 receptors: novel ligands and paradoxical effects. Trends in Pharmacological Sciences, 1998, 19, 184-191.	4.0	292
18	Agonist-bound structure of the human P2Y12 receptor. Nature, 2014, 509, 119-122.	13.7	279

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19	Adenosine A3 receptor stimulation and cerebral ischemia. European Journal of Pharmacology, 1994, 263, 59-67.	1.7	266
20	Structure-Activity Relationships of N6-Benzyladenosine-5'-uronamides as A3-Selective Adenosine Agonists. Journal of Medicinal Chemistry, 1994, 37, 636-646.	2.9	248
21	Adenosine receptor ligands: differences with acute versus chronic treatment. Trends in Pharmacological Sciences, 1996, 17, 108-113.	4.0	248
22	The Concise Guide to PHARMACOLOGY 2015/16: Overview. British Journal of Pharmacology, 2015, 172, 5729-5743.	2.7	220
23	2-Substitution of N6-Benzyladenosine-5'-uronamides Enhances Selectivity for A3 Adenosine Receptors. Journal of Medicinal Chemistry, 1994, 37, 3614-3621.	2.9	213
24	Medicinal chemistry of adenosine, P2Y and P2X receptors. Neuropharmacology, 2016, 104, 31-49.	2.0	213
25	Structure-Based Discovery of A_{2A}Adenosine Receptor Ligands. Journal of Medicinal Chemistry, 2010, 53, 3748-3755.	2.9	212
26	A physiological role of the adenosine A3 receptor: Sustained cardioprotection. Proceedings of the National Academy of Sciences of the United States of America, 1998, 95, 6995-6999.	3.3	202
27	Pharmacological and therapeutic effects of A3 adenosine receptor agonists. Drug Discovery Today, 2012, 17, 359-366.	3.2	193
28	Anilide Derivatives of an 8-Phenylxanthine Carboxylic Congener Are Highly Potent and Selective Antagonists at Human A2BAdenosine Receptors. Journal of Medicinal Chemistry, 2000, 43, 1165-1172.	2.9	192
29	Site-directed Mutagenesis Identifies Residues Involved in Ligand Recognition in the Human A2a Adenosine Receptor. Journal of Biological Chemistry, 1995, 270, 13987-13997.	1.6	191
30	Competitive and selective antagonism of P2Y1 receptors by N 6 -methyl 2â€²-deoxyadenosine 3â€²,5â€²-bisphosphate. British Journal of Pharmacology, 1998, 124, 1-3.	2.7	188
31	Synthesis, CoMFA Analysis, and Receptor Docking of 3,5-Diacyl-2,4-Dialkylpyridine Derivatives as Selective A3Adenosine Receptor Antagonists. Journal of Medicinal Chemistry, 1999, 42, 706-721.	2.9	187
32	Molecular architecture of G protein-coupled receptors. Drug Development Research, 1996, 37, 1-38.	1.4	180
33	Allosteric Coupling of Drug Binding and Intracellular Signaling in the A2A Adenosine Receptor. Cell, 2018, 172, 68-80.e12.	13.5	173
34	Increased Signaling via Adenosine A1 Receptors, Sleep Deprivation, Imipramine, and Ketamine Inhibit Depressive-like Behavior via Induction of Homer1a. Neuron, 2015, 87, 549-562.	3.8	168
35	Induction of Apoptosis in HL-60 Human Promyelocytic Leukemia Cells by Adenosine A3Receptor Agonists. Biochemical and Biophysical Research Communications, 1996, 219, 904-910.	1.0	166
36	Update of P2X receptor properties and their pharmacology: IUPHAR Review 30. British Journal of Pharmacology, 2021, 178, 489-514.	2.7	165

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37	Activation of Hippocampal Adenosine A₃Receptors Produces a Desensitization of A₁Receptor-Mediated Responses in Rat Hippocampus. Journal of Neuroscience, 1997, 17, 607-614.	1.7	159
38	8-(3-Chlorostyryl)caffeine (CSC) is a selective A2 -adenosine antagonist in vitro and in vivo. FEBS Letters, 1993, 323, 141-144.	1.3	158
39	Derivatives of the Triazoloquinazoline Adenosine Antagonist (CGS15943) Are Selective for the Human A3Receptor Subtype. Journal of Medicinal Chemistry, 1996, 39, 4142-4148.	2.9	154
40	Progress in the pursuit of therapeutic adenosine receptor antagonists. Medicinal Research Reviews, 2006, 26, 131-159.	5.0	154
41	Human P2Y1Receptor:Â Molecular Modeling and Site-Directed Mutagenesis as Tools To Identify Agonist and Antagonist Recognition Sites. Journal of Medicinal Chemistry, 1998, 41, 1456-1466.	2.9	153
42	New paradigms in GPCR drug discovery. Biochemical Pharmacology, 2015, 98, 541-555.	2.0	152
43	Structure-activity relationships of 8-styrylxanthines as A2-selective adenosine antagonists. Journal of Medicinal Chemistry, 1993, 36, 1333-1342.	2.9	151
44	Structural Determinants of A3Adenosine Receptor Activation:Â Nucleoside Ligands at the Agonist/Antagonist Boundary. Journal of Medicinal Chemistry, 2002, 45, 4471-4484.	2.9	151
45	Systematic Investigation of Polyamidoamine Dendrimers Surface-Modified with Poly(ethylene glycol) for Drug Delivery Applications: Synthesis, Characterization, and Evaluation of Cytotoxicity. Bioconjugate Chemistry, 2008, 19, 1660-1672.	1.8	151
46	Update of P2Y receptor pharmacology: IUPHAR Review 27. British Journal of Pharmacology, 2020, 177, 2413-2433.	2.7	151
47	Pharmacological characterization of novel A3 adenosine receptor-selective antagonists. Neuropharmacology, 1997, 36, 1157-1165.	2.0	150
48	Chronic caffeine alters the density of adenosine, adrenergic, cholinergic, GABA, and serotonin receptors and calcium channels in mouse brain. Cellular and Molecular Neurobiology, 1993, 13, 247-261.	1.7	149
49	Historical and Current Adenosine Receptor Agonists in Preclinical and Clinical Development. Frontiers in Cellular Neuroscience, 2019, 13, 124.	1.8	146
50	A role for central A3-adenosine receptors. FEBS Letters, 1993, 336, 57-60.	1.3	145
51	Adenosine A1 and A2 receptors: Structureâ€“function relationships. Medicinal Research Reviews, 1992, 12, 423-471.	5.0	144
52	Diisothiocyanate derivatives as potent, insurmountable antagonists of P2Y6 nucleotide receptors. Biochemical Pharmacology, 2004, 67, 1763-1770.	2.0	142
53	Architecture of P2Y Nucleotide Receptors:â€‰ Structural Comparison Based on Sequence Analysis, Mutagenesis, and Homology Modeling. Journal of Medicinal Chemistry, 2004, 47, 5393-5404.	2.9	139
54	N6-Substituted adenosine derivatives: selectivity, efficacy, and species differences at A3 adenosine receptors. Biochemical Pharmacology, 2003, 65, 1675-1684.	2.0	136

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55	MRS2500 [2-Iodo-N6-methyl-(N)-methanocarba-2â€²-deoxyadenosine-3â€²,5â€²-bisphosphate], a Potent, Selective, and Stable Antagonist of the Platelet P2Y1 Receptor with Strong Antithrombotic Activity in Mice. Journal of Pharmacology and Experimental Therapeutics, 2006, 316, 556-563.	1.3	135
56	Identification by Site-directed Mutagenesis of Residues Involved in Ligand Recognition and Activation of the Human A3 Adenosine Receptor. Journal of Biological Chemistry, 2002, 277, 19056-19063.	1.6	134
57	Structural Connection between Activation Microswitch and Allosteric Sodium Site in GPCR Signaling. Structure, 2018, 26, 259-269.e5.	1.6	134
58	Cerebral ischemia in gerbils: effects of acute and chronic treatment with adenosine A2A receptor agonist and antagonist. European Journal of Pharmacology, 1995, 287, 295-302.	1.7	133
59	Deoxyadenosine Bisphosphate Derivatives as Potent Antagonists at P2Y1Receptors. Journal of Medicinal Chemistry, 1998, 41, 183-190.	2.9	133
60	Activation of Th1 and Tc1 cell adenosine A2A receptors directly inhibits IL-2 secretion in vitro and IL-2-driven expansion in vivo. Blood, 2005, 105, 4707-4714.	0.6	133
61	The Role of Amino Acids in Extracellular Loops of the Human P2Y1 Receptor in Surface Expression and Activation Processes. Journal of Biological Chemistry, 1999, 274, 14639-14647.	1.6	132
62	P2Y nucleotide receptors: promise of therapeutic applications. Drug Discovery Today, 2010, 15, 570-578.	3.2	132
63	A3-adenosine receptors: Design of selective ligands and therapeutic prospects. Drugs of the Future, 1995, 20, 689.	0.0	132
64	Differential effects of P₂â€purinoceptor antagonists on phospholipase Câ€•and adenylyl cyclaseâ€coupled P_{2Y}â€purinoceptors. British Journal of Pharmacology, 1994, 113, 614-620.	2.7	129
65	Synthesis, Biological Activity, and Molecular Modeling of Ribose-Modified Deoxyadenosine Bisphosphate Analogues as P2Y1Receptor Ligands. Journal of Medicinal Chemistry, 2000, 43, 829-842.	2.9	129
66	Methanocarba Analogues of Purine Nucleosides as Potent and Selective Adenosine Receptor Agonists. Journal of Medicinal Chemistry, 2000, 43, 2196-2203.	2.9	127
67	Structureâ^'Activity Relationships and Molecular Modeling of 3,5-Diacyl-2,4-dialkylpyridine Derivatives as Selective A3 Adenosine Receptor Antagonists. Journal of Medicinal Chemistry, 1998, 41, 3186-3201.	2.9	126
68	Role of the Extracellular Loops of G Protein-Coupled Receptors in Ligand Recognition:Â A Molecular Modeling Study of the Human P2Y1Receptor. Biochemistry, 1999, 38, 3498-3507.	1.2	125
69	2-Substitution of Adenine Nucleotide Analogues Containing a Bicyclo[3.1.0]hexane Ring System Locked in a Northern Conformation:Â Enhanced Potency as P2Y1Receptor Antagonists. Journal of Medicinal Chemistry, 2003, 46, 4974-4987.	2.9	125
70	Modeling the Adenosine Receptors:Â Comparison of the Binding Domains of A2AAgonists and Antagonists. Journal of Medicinal Chemistry, 2003, 46, 4847-4859.	2.9	125
71	Direct preconditioning of cultured chick ventricular myocytes. Novel functions of cardiac adenosine A2a and A3 receptors.. Journal of Clinical Investigation, 1996, 98, 1773-1779.	3.9	124
72	A Mutational Analysis of Residues Essential for Ligand Recognition at the Human P2Y₁Receptor. Molecular Pharmacology, 1997, 52, 499-507.	1.0	123

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73	Endogenous adenosine A3 receptor activation selectively alleviates persistent pain states. Brain, 2015, 138, 28-35.	3.7	120
74	Identification of potent, selective P2Y-purinoceptor agonists: structure-activity relationships for 2-thioether derivatives of adenosine 5'-triphosphate. Journal of Medicinal Chemistry, 1993, 36, 3937-3946.	2.9	116
75	Structureâ€“activity relationships of thiazole and thiadiazole derivatives as potent and selective human adenosine A3 receptor antagonists. Bioorganic and Medicinal Chemistry, 2004, 12, 613-623.	1.4	115
76	Deficiency of adenosine deaminase 2 triggers adenosine-mediated NETosis and TNF production in patients with DADA2. Blood, 2019, 134, 395-406.	0.6	115
77	Functionalized congeners of 1,3-dialkylxanthines: preparation of analogs with high affinity for adenosine receptors. Journal of Medicinal Chemistry, 1985, 28, 1334-1340.	2.9	114
78	Spinal neuroimmune activation is independent of T-cell infiltration and attenuated by A3 adenosine receptor agonists in a model of oxaliplatin-induced peripheral neuropathy. Brain, Behavior, and Immunity, 2015, 44, 91-99.	2.0	114
79	[3H]xanthine amine congener of 1,3-dipropyl-8-phenylxanthine: an antagonist radioligand for adenosine receptors.. Proceedings of the National Academy of Sciences of the United States of America, 1986, 83, 4089-4093.	3.3	111
80	Synthesis and Biological Activities of Flavonoid Derivatives as A3 Adenosine Receptor Antagonists. Journal of Medicinal Chemistry, 1996, 39, 2293-2301.	2.9	111
81	A₃ Adenosine Receptors as Modulators of Inflammation: From Medicinal Chemistry to Therapy. Medicinal Research Reviews, 2018, 38, 1031-1072.	5.0	111
82	Interactions of Flavonoids and Other Phytochemicals with Adenosine Receptors. Journal of Medicinal Chemistry, 1996, 39, 781-788.	2.9	110
83	Interaction of 1,4-Dihydropyridine and Pyridine Derivatives with Adenosine Receptors:â€‰ Selectivity for A3 Receptors. Journal of Medicinal Chemistry, 1996, 39, 2980-2989.	2.9	108
84	Small molecule blockers of the Alzheimer AÎ² calcium channel potently protect neurons from AÎ² cytotoxicity. Proceedings of the National Academy of Sciences of the United States of America, 2009, 106, 3348-3353.	3.3	108
85	Xanthines as Adenosine Receptor Antagonists. Handbook of Experimental Pharmacology, 2011, , 151-199.	0.9	107
86	Derivatives of the Triazoloquinazoline Adenosine Antagonist (CGS 15943) Having High Potency at the Human A2Band A3Receptor Subtypes. Journal of Medicinal Chemistry, 1998, 41, 2835-2845.	2.9	106
87	Chronic administration of selective adenosine A1 receptor agonist or antagonist in cerebral ischemia. European Journal of Pharmacology, 1994, 256, 161-167.	1.7	104
88	Dihydropyridines as inhibitors of capacitative calcium entry in leukemic HL-60 cells. Biochemical Pharmacology, 2003, 65, 329-338.	2.0	103
89	Structure activity relationships for derivatives of adenosine-5?-triphosphate as agonists at P2 purinoceptors: Heterogeneity within P2x and P2y subtypes. Drug Development Research, 1994, 31, 206-219.	1.4	101
90	Adenosine A3Receptor Agonists Protect HL-60 and U-937 Cells from Apoptosis Induced by A3Antagonists. Biochemical and Biophysical Research Communications, 1997, 232, 317-322.	1.0	101

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91	Development of selective agonists and antagonists of P2Y receptors. Purinergic Signalling, 2009, 5, 75-89.	1.1	101
92	Emerging adenosine receptor agonists. Expert Opinion on Emerging Drugs, 2007, 12, 479-492.	1.0	100
93	Controlling murine and rat chronic pain through A ₃ adenosine receptor activation. FASEB Journal, 2012, 26, 1855-1865.	0.2	99
94	Adenosine-induced cell death: evidence for receptor-mediated signalling. Apoptosis: an International Journal on Programmed Cell Death, 1999, 4, 197-211.	2.2	98
95	5â€²-Phosphate and 5â€²-Phosphonate Ester Derivatives of (N)-Methanocarba Adenosine with in Vivo Cardioprotective Activity. Journal of Medicinal Chemistry, 2013, 56, 902-914.	2.9	98
96	Search for New Purine- and Ribose-Modified Adenosine Analogs as Selective Agonists and Antagonists at Adenosine Receptors. Journal of Medicinal Chemistry, 1995, 38, 1174-1188.	2.9	97
97	Quantitation of the P2Y1Receptor with a High Affinity Radiolabeled Antagonist. Molecular Pharmacology, 2002, 62, 1249-1257.	1.0	95
98	Induction of Novel Agonist Selectivity for the ADP-Activated P2Y1 Receptor Versus the ADP-Activated P2Y12 and P2Y13 Receptors by Conformational Constraint of an ADP Analog. Journal of Pharmacology and Experimental Therapeutics, 2004, 311, 1038-1043.	1.3	95
99	Synthesis of pyridoxal phosphate derivatives with antagonist activity at the P2Y13 receptor. Biochemical Pharmacology, 2005, 70, 266-274.	2.0	95
100	Positive Inotropic Effects by Uridine Triphosphate (UTP) and Uridine Diphosphate (UDP) via P2Y 2 and P2Y 6 Receptors on Cardiomyocytes and Release of UTP in Man During Myocardial Infarction. Circulation Research, 2006, 98, 970-976.	2.0	95
101	Behavioral characterization of mice lacking the A3 adenosine receptor: sensitivity to hypoxic neurodegeneration. Cellular and Molecular Neurobiology, 2003, 23, 431-447.	1.7	94
102	Stimulation of the P2X7 receptor kills rat retinal ganglion cells in vivo. Experimental Eye Research, 2010, 91, 425-432.	1.2	93
103	A3 adenosine receptor agonist prevents the development of paclitaxel-induced neuropathic pain by modulating spinal glial-restricted redox-dependent signaling pathways. Pain, 2014, 155, 2560-2567.	2.0	93
104	Species differences in structure-activity relationships of adenosine agonists and xanthine antagonists at brain A1 adenosine receptors. FEBS Letters, 1986, 209, 122-128.	1.3	92
105	Induction of Apoptosis in Cardiac Myocytes by an A3Adenosine Receptor Agonist. Experimental Cell Research, 1998, 243, 383-397.	1.2	91
106	Identification of the A2 adenosine receptor binding subunit by photoaffinity crosslinking. Proceedings of the National Academy of Sciences of the United States of America, 1989, 86, 6572-6576.	3.3	90
107	Methanocarba Modification of Uracil and Adenine Nucleotides:Â High Potency of Northern Ring Conformation at P2Y1, P2Y2, P2Y4, and P2Y11but Not P2Y6Receptors. Journal of Medicinal Chemistry, 2002, 45, 208-218.	2.9	90
108	Evaluation of Homology Modeling of G-Protein-Coupled Receptors in Light of the A_{2A} Adenosine Receptor Crystallographic Structure. Journal of Medicinal Chemistry, 2009, 52, 3284-3292.	2.9	90

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109	[3H]MRS 1754, a selective antagonist radioligand for A2B adenosine receptors. Biochemical Pharmacology, 2001, 61, 657-663.	2.0	89
110	2-Chloro N6 -methyl-(N)-methanocarba-2â€²-deoxyadenosine-3â€²,5â€²-bisphosphate is a selective high affinity P2Y1 receptor antagonist. British Journal of Pharmacology, 2002, 135, 2004-2010.	2.7	89
111	Introduction to Adenosine Receptors as Therapeutic Targets. Handbook of Experimental Pharmacology, 2009, , 1-24.	0.9	89
112	A Selective High-Affinity Antagonist of the P2Y₁₄ Receptor Inhibits UDP-Glucoseâ€“Stimulated Chemotaxis of Human Neutrophils. Molecular Pharmacology, 2013, 84, 41-49.	1.0	89
113	6-Phenyl-1,4-dihydropyridine Derivatives as Potent and Selective A3Adenosine Receptor Antagonists. Journal of Medicinal Chemistry, 1996, 39, 4667-4675.	2.9	88
114	The A3Adenosine Receptor Mediates Cell Spreading, Reorganization of Actin Cytoskeleton, and Distribution of Bcl-xL: Studies in Human Astroglioma Cells. Biochemical and Biophysical Research Communications, 1997, 241, 297-304.	1.0	88
115	(N)-Methanocarba 2,N6-Disubstituted Adenine Nucleosides as Highly Potent and Selective A3Adenosine Receptor Agonists. Journal of Medicinal Chemistry, 2005, 48, 1745-1758.	2.9	88
116	Digitoxin mimics gene therapy with CFTR and suppresses hypersecretion of IL-8 from cystic fibrosis lung epithelial cells. Proceedings of the National Academy of Sciences of the United States of America, 2004, 101, 7693-7698.	3.3	87
117	Locomotor activity in mice during chronic treatment with caffeine and withdrawal. Pharmacology Biochemistry and Behavior, 1993, 44, 199-216.	1.3	86
118	The effects of adenosine A3 receptor stimulation on seizures in mice. European Journal of Pharmacology, 1995, 275, 23-29.	1.7	86
119	Structureâ^'Activity Relationships of 4-(Phenylethynyl)-6-phenyl-1,4- dihydropyridines as Highly Selective A3 Adenosine Receptor Antagonists. Journal of Medicinal Chemistry, 1997, 40, 2596-2608.	2.9	86
120	Structureâ^'Activity Relationships of Pyridoxal Phosphate Derivatives as Potent and Selective Antagonists of P2X1Receptors. Journal of Medicinal Chemistry, 2001, 44, 340-349.	2.9	86
121	Cardioprotective effects of adenosine A1 and A3 receptor activation during hypoxia in isolated rat cardiac myocytes. Molecular and Cellular Biochemistry, 2001, 217, 143-152.	1.4	86
122	A2B adenosine receptor blockade inhibits growth of prostate cancer cells. Purinergic Signalling, 2013, 9, 271-280.	1.1	86
123	Nucleotides Acting at P2Y Receptors: Connecting Structure and Function. Molecular Pharmacology, 2015, 88, 220-230.	1.0	86
124	Antiaggregatory activity in human platelets of potent antagonists of the P2Y1 receptor. Biochemical Pharmacology, 2004, 68, 1995-2002.	2.0	85
125	N6-Substituted D-4â€˜-Thioadenosine-5â€˜-methyluronamides:â€‰ Potent and Selective Agonists at the Human A3 Adenosine Receptor. Journal of Medicinal Chemistry, 2003, 46, 3775-3777.	2.9	83
126	Allosteric Modulation of A3 Adenosine Receptors by a Series of 3-(2-Pyridinyl)isoquinoline Derivatives. Molecular Pharmacology, 2001, 60, 1057-1063.	1.0	82

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127	Imidazo[2,1-i]purin-5-ones and Related Tricyclic Water-Soluble Purine Derivatives:Â Potent A2A- and A3-Adenosine Receptor Antagonistsâ€. Journal of Medicinal Chemistry, 2002, 45, 3440-3450.	2.9	81
128	Identification of Essential Residues Involved in the Allosteric Modulation of the Human A3Adenosine Receptor. Molecular Pharmacology, 2003, 63, 1021-1031.	1.0	81
129	Discovery of a New Nucleoside Template for Human A3 Adenosine Receptor Ligands:â€‰ d-4â€˜-Thioadenosine Derivatives without 4â€˜-Hydroxymethyl Group as Highly Potent and Selective Antagonists. Journal of Medicinal Chemistry, 2007, 50, 3159-3162.	2.9	81
130	Effects of chronic administration of adenosine A1 receptor agonist and antagonist on spatial learning and memory. European Journal of Pharmacology, 1993, 249, 271-280.	1.7	80
131	Neoceptor Concept Based on Molecular Complementarity in GPCRs:â€‰ A Mutant Adenosine A3 Receptor with Selectively Enhanced Affinity for Amine-Modified Nucleosides. Journal of Medicinal Chemistry, 2001, 44, 4125-4136.	2.9	80
132	Activation of the A 3 adenosine receptor affects cell cycle progression and cell growth. Naunyn-Schmiedeberg's Archives of Pharmacology, 2000, 361, 225-234.	1.4	79
133	Identification of Acidic Residues in the Extracellular Loops of the Seven-transmembrane Domain of the Human Ca2+ Receptor Critical for Response to Ca2+ and a Positive Allosteric Modulator. Journal of Biological Chemistry, 2002, 277, 46622-46631.	1.6	79
134	Acyclic Analogues of Adenosine Bisphosphates as P2Y Receptor Antagonists:Â Phosphate Substitution Leads to Multiple Pathways of Inhibition of Platelet Aggregation. Journal of Medicinal Chemistry, 2002, 45, 5694-5709.	2.9	79
135	Functionalized congeners of adenosine: preparation of analogs with high affinity for A1-adenosine receptors. Journal of Medicinal Chemistry, 1985, 28, 1341-1346.	2.9	78
136	Heteromultimeric P2X1/2 Receptors Show a Novel Sensitivity to Extracellular pH. Journal of Pharmacology and Experimental Therapeutics, 2002, 300, 673-680.	1.3	78
137	Treatment of Dry Eye Syndrome with Orally Administered CF101. Ophthalmology, 2010, 117, 1287-1293.	2.5	78
138	G protein-coupled adenosine (P1) and P2Y receptors: ligand design and receptor interactions. Purinergic Signalling, 2012, 8, 419-436.	1.1	78
139	Modulation of Apoptosis by Adenosine in the Central Nervous System: a Possible Role for the A3Receptor.. Annals of the New York Academy of Sciences, 1997, 825, 11-22.	1.8	77
140	New Insights for Drug Design from the X-Ray Crystallographic Structures of G-Protein-Coupled Receptors. Molecular Pharmacology, 2012, 82, 361-371.	1.0	77
141	Induction of Apoptosis in Rat Cardiocytes by A3 Adenosine Receptor Activation and Its Suppression by Isoproterenol. Experimental Cell Research, 2000, 257, 111-126.	1.2	76
142	Structureâ^'Activity Relationships at Human and Rat A2BAdenosine Receptors of Xanthine Derivatives Substituted at the 1-, 3-, 7-, and 8-Positions. Journal of Medicinal Chemistry, 2002, 45, 2131-2138.	2.9	76
143	A Region in the Seven-transmembrane Domain of the Human Ca2+ Receptor Critical for Response to Ca2+. Journal of Biological Chemistry, 2005, 280, 5113-5120.	1.6	76
144	Quantification of G_i-Mediated Inhibition of Adenylyl Cyclase Activity Reveals That UDP Is a Potent Agonist of the Human P2Y₁₄ Receptor. Molecular Pharmacology, 2009, 76, 1341-1348.	1.0	76

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145	Structure-Guided Design of A₃ Adenosine Receptor-Selective Nucleosides: Combination of 2-Arylethynyl and Bicyclo[3.1.0]hexane Substitutions. Journal of Medicinal Chemistry, 2012, 55, 4847-4860.	2.9	76
146	Effects of a Calcimimetic Compound and Naturally Activating Mutations on the Human Ca²⁺Receptor and on Ca²⁺Receptor/Metabotropic Glutamate Chimeric Receptors. Endocrinology, 2000, 141, 4156-4163.	1.4	75
147	Extracellular nucleotides induce vasodilatation in human arteries via prostaglandins, nitric oxide and endothelium-derived hyperpolarising factor. British Journal of Pharmacology, 2003, 138, 1451-1458.	2.7	75
148	Human P2Y6 Receptor:â€‰Molecular Modeling Leads to the Rational Design of a Novel Agonist Based on a Unique Conformational Preference. Journal of Medicinal Chemistry, 2005, 48, 8108-8111.	2.9	75
149	Structureâ^'Activity Relationships of Uridine 5â€˜-Diphosphate Analogues at the Human P2Y6Receptor. Journal of Medicinal Chemistry, 2006, 49, 5532-5543.	2.9	75
150	Structureâ^'Activity Relationship of (N)-Methanocarba Phosphonate Analogues of 5â€²-AMP as Cardioprotective Agents Acting Through a Cardiac P2X Receptor. Journal of Medicinal Chemistry, 2010, 53, 2562-2576.	2.9	75
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