Marek Bajda

List of Publications by Year
in descending order

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83	2,376	236612	223531
papers	citations	25	46
		h-index	g-index

85	85	85	3328
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Multilevel virtual screening approach for discovery of thymidine phosphorylase inhibitors as potential anticancer agents. Journal of Molecular Structure, 2022, 1249, 131648.	1.8	0
2	New hybrids of tacrine and indomethacin as multifunctional acetylcholinesterase inhibitors. Chemical Papers, 2021, 75, 249-264.	1.0	7
3	Pyridinium-2-carbaldoximes with quinolinium carboxamide moiety are simultaneous reactivators of acetylcholinesterase and butyrylcholinesterase inhibited by nerve agent surrogates. Journal of Enzyme Inhibition and Medicinal Chemistry, 2021, 36, 437-449.	2.5	4
4	Biological evaluation and molecular docking of novel 1,3,4-thiadiazole-resorcinol conjugates as multifunctional cholinesterases inhibitors. Bioorganic Chemistry, 2021, 107, 104617.	2.0	19
5	Dual Inhibitors of Amyloid-Î² and Tau Aggregation with Amyloid-Î² Disaggregating Properties: Extended In Cellulo, In Silico, and Kinetic Studies of Multifunctional Anti-Alzheimerâ€™s Agents. ACS Chemical Neuroscience, 2021, 12, 2057-2068.	1.7	36
6	Guanidine Derivatives: How Simple Structural Modification of Histamine H₃R Antagonists Has Led to the Discovery of Potent Muscarinic M₂R/M₄R Antagonists. ACS Chemical Neuroscience, 2021, 12, 2503-2519.	1.7	7
7	Biphenylalkoxyamine Derivativesâ€“Histamine H3 Receptor Ligands with Butyrylcholinesterase Inhibitory Activity. Molecules, 2021, 26, 3580.	1.7	3
8	Design, Synthesis, and In Vitro Antiproliferative Activity of Hydantoin and Purine Derivatives with the 4-Acetylphenylpiperazinylalkyl Moiety. Materials, 2021, 14, 4156.	1.3	1
9	Novel Functionalized Amino Acids as Inhibitors of GABA Transporters with Analgesic Activity. ACS Chemical Neuroscience, 2021, 12, 3073-3100.	1.7	6
10	Cyanobiphenyls: Novel H3 receptor ligands with cholinesterase and MAO B inhibitory activity as multitarget compounds for potential treatment of Alzheimerâ€™s disease. Bioorganic Chemistry, 2021, 114, 105129.	2.0	8
11	Development of tricyclic N-benzyl-4-hydroxybutanamide derivatives as inhibitors of GABA transporters mGAT1-4 with anticonvulsant, antinociceptive, and antidepressant activity. European Journal of Medicinal Chemistry, 2021, 221, 113512.	2.6	6
12	Molecular Modeling Studies on the Multistep Reactivation Process of Organophosphate-Inhibited Acetylcholinesterase and Butyrylcholinesterase. Biomolecules, 2021, 11, 169.	1.8	6
13	Search for new multi-target compounds against Alzheimerâ€™s disease among histamine H3 receptor ligands. European Journal of Medicinal Chemistry, 2020, 185, 111785.	2.6	27
14	Rational design of new multitarget histamine H3 receptor ligands as potential candidates for treatment of Alzheimerâ€™s disease. European Journal of Medicinal Chemistry, 2020, 207, 112743.	2.6	17
15	Multidirectional inÂvitro and in cellulo studies as a tool for identification of multi-target-directed ligands aiming at symptoms and causes of Alzheimerâ€™s disease. Journal of Enzyme Inhibition and Medicinal Chemistry, 2020, 35, 1944-1952.	2.5	9
16	Structure modeling of Î³-aminobutyric acid transporters â€“ Molecular basics of ligand selectivity. International Journal of Biological Macromolecules, 2020, 158, 1380-1389.	3.6	10
17	Î³-Aminobutyric acid transporters as relevant biological target: Their function, structure, inhibitors and role in the therapy of different diseases. International Journal of Biological Macromolecules, 2020, 158, 750-772.	3.6	27
18	New Tetrahydroacridine Hybrids with Dichlorobenzoic Acid Moiety Demonstrating Multifunctional Potential for the Treatment of Alzheimerâ€™s Disease. International Journal of Molecular Sciences, 2020, 21, 3765.	1.8	6

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19	Structure Modeling of the Norepinephrine Transporter. Biomolecules, 2020, 10, 102.	1.8	12
20	Novel application of capillary electrophoresis with a liposome coated capillary for prediction of blood-brain barrier permeability. Talanta, 2020, 217, 121023.	2.9	13
21	Synthesis, physicochemical and biological evaluation of tacrine derivative labeled with technetium-99m and gallium-68 as a prospective diagnostic tool for early diagnosis of Alzheimerâ€™s disease. Bioorganic Chemistry, 2019, 91, 103136.	2.0	7
22	Discovery of New Cyclopentaquinoline Analogues as Multifunctional Agents for the Treatment of Alzheimerâ€™s Disease. International Journal of Molecular Sciences, 2019, 20, 498.	1.8	12
23	Search for multifunctional agents against Alzheimerâ€™s disease among non-imidazole histamine H3 receptor ligands. In vitro and in vivo pharmacological evaluation and computational studies of piperazine derivatives. Bioorganic Chemistry, 2019, 90, 103084.	2.0	13
24	EC359: A First-in-Class Small-Molecule Inhibitor for Targeting Oncogenic LIFR Signaling in Triple-Negative Breast Cancer. Molecular Cancer Therapeutics, 2019, 18, 1341-1354.	1.9	41
25	Biological Evaluation, Molecular Docking, and SAR Studies of Novel 2-(2,4-Dihydroxyphenyl)-1H-Benzimidazole Analogues. Biomolecules, 2019, 9, 870.	1.8	12
26	Characteristics of metabolic stability and the cell permeability of 2â€pyrimidinylâ€piperazinylâ€alkyl derivatives of 1Hâ€imidazo[2,1 â€f]purineâ€2,4(3 H ,8 H)â€dione with antidepressantâ€and anxiolyticâ€like activities. Chemical Biology and Drug Design, 2019, 93, 511-521.	1.5	8
27	Identification of 1,2,4-triazoles as new thymidine phosphorylase inhibitors: Future anti-tumor drugs. Bioorganic Chemistry, 2019, 85, 209-220.	2.0	41
28	Abstract 4716: Therapeutic utility of EC359 for targeting oncogenic LIFR signaling in triple negative breast cancer. , 2019, , .		0
29	Novel carbamate derivatives as selective butyrylcholinesterase inhibitors. Bioorganic Chemistry, 2018, 78, 29-38.	2.0	18
30	Design, Synthesis, and Biological Evaluation of 1-Benzylamino-2-hydroxyalkyl Derivatives as New Potential Disease-Modifying Multifunctional Anti-Alzheimerâ€™s Agents. ACS Chemical Neuroscience, 2018, 9, 1074-1094.	1.7	47
31	Tetrahydroacridine derivatives with dichloronicotinic acid moiety as attractive, multipotent agents for Alzheimer's disease treatment. European Journal of Medicinal Chemistry, 2018, 145, 760-769.	2.6	21
32	New cyclopentaquinoline hybrids with multifunctional capacities for the treatment of Alzheimerâ€™s disease. Journal of Enzyme Inhibition and Medicinal Chemistry, 2018, 33, 158-170.	2.5	17
33	Novel pyridineâ€2,4,6â€tricarbohydrazide thiourea compounds as small key organic molecules for the potential treatment of typeâ€2 diabetes mellitus: In vitro studies against yeast Î±â€and Î²â€glucosidase and in silico molecular modeling. Archiv Der Pharmazie, 2018, 351, 1700236.	2.1	3
34	4-tert-Pentylphenoxyalkyl derivatives â€“ Histamine H3 receptor ligands and monoamine oxidase B inhibitors. Bioorganic and Medicinal Chemistry Letters, 2018, 28, 3596-3600.	1.0	13
35	Determination of ligand efficiency indices in a group of 7Hâ€purineâ€2,6â€dione derivatives with psychotropic activity using micellar electrokinetic chromatography. Electrophoresis, 2018, 39, 2446-2453.	1.3	3
36	Design, Synthesis, and Biological Evaluation of 2-(Benzylamino-2-Hydroxyalkyl)Isoindoline-1,3-Diones Derivatives as Potential Disease-Modifying Multifunctional Anti-Alzheimer Agents. Molecules, 2018, 23, 347.	1.7	27

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37	Lipophilic properties of antiâ€Alzheimer's agents determined by micellar electrokinetic chromatography and reversedâ€phase thinâ€layer chromatography. Electrophoresis, 2017, 38, 1268-1275.	1.3	9
38	Tetrahydroacridine derivatives with fluorobenzoic acid moiety as multifunctional agents for Alzheimerâ€™s disease treatment. Bioorganic Chemistry, 2017, 72, 315-322.	2.0	17
39	Synthesis, physicochemical and biological studies of technetium-99m labeled tacrine derivative as a diagnostic tool for evaluation of cholinesterase level. Bioorganic and Medicinal Chemistry, 2017, 25, 912-920.	1.4	9
40	Design, synthesis and biological evaluation of new phthalimide and saccharin derivatives with alicyclic amines targeting cholinesterases, beta-secretase and amyloid beta aggregation. European Journal of Medicinal Chemistry, 2017, 125, 676-695.	2.6	85
41	Computational approach for the assessment of inhibitory potency against beta-amyloid aggregation. Bioorganic and Medicinal Chemistry Letters, 2017, 27, 212-216.	1.0	16
42	Hybrid approach to structure modeling of the histamine H3 receptor: Multi-level assessment as a tool for model verification. PLoS ONE, 2017, 12, e0186108.	1.1	14
43	Synthesis, Molecular Modelling and Biological Evaluation of Novel Heterodimeric, Multiple Ligands Targeting Cholinesterases and Amyloid Beta. Molecules, 2016, 21, 410.	1.7	29
44	Organocatalyzed Novel Synthetic Methodology for Highly Functionalized Piperidines as Potent Î±â€Glucosidase Inhibitors. Archiv Der Pharmazie, 2016, 349, 724-732.	2.1	9
45	Cholinesterase inhibitory activity of chlorophenoxy derivativesâ€”Histamine H3 receptor ligands. Bioorganic and Medicinal Chemistry Letters, 2016, 26, 4140-4145.	1.0	20
46	A Hybrid Approach to Structure and Function Modeling of G Protein-Coupled Receptors. Journal of Chemical Information and Modeling, 2016, 56, 630-641.	2.5	14
47	Novel biphenyl bis -sulfonamides as acetyl and butyrylcholinesterase inhibitors: Synthesis, biological evaluation and molecular modeling studies. Bioorganic Chemistry, 2016, 64, 13-20.	2.0	56
48	The Lipophilicity Estimation of Selected Dermatological Drugs Using Micellar Electrokinetic Chromatography Method. Journal of Liquid Chromatography and Related Technologies, 2015, 38, 1435-1438.	0.5	2
49	Multiple Ligands Targeting Cholinesterases and Î²â€Amyloid: Synthesis, Biological Evaluation of Heterodimeric Compounds with Benzylamine Pharmacophore. Archiv Der Pharmazie, 2015, 348, 556-563.	2.1	11
50	Design and synthesis of new barbituric- and thiobarbituric acid derivatives as potent urease inhibitors: Structure activity relationship and molecular modeling studies. Bioorganic and Medicinal Chemistry, 2015, 23, 6049-6058.	1.4	53
51	Isoindoline-1,3-dione derivatives targeting cholinesterases: Design, synthesis and biological evaluation of potential anti-Alzheimerâ€™s agents. Bioorganic and Medicinal Chemistry, 2015, 23, 1629-1637.	1.4	44
52	Development of multifunctional, heterodimeric isoindoline-1,3-dione derivatives as cholinesterase and Î²-amyloid aggregation inhibitors with neuroprotective properties. European Journal of Medicinal Chemistry, 2015, 92, 738-749.	2.6	60
53	Synthesis, biological evaluation and molecular modeling of new tetrahydroacridine derivatives as potential multifunctional agents for the treatment of Alzheimerâ€™s disease. Bioorganic and Medicinal Chemistry, 2015, 23, 5610-5618.	1.4	26
54	Synthesis, thymidine phosphorylase inhibition and molecular modeling studies of 1,3,4-oxadiazole-2-thione derivatives. Bioorganic Chemistry, 2015, 60, 37-41.	2.0	21

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55	Synthesis of new N-benzylpiperidine derivatives as cholinesterase inhibitors with Î²-amyloid anti-aggregation properties and beneficial effects on memory in vivo. Bioorganic and Medicinal Chemistry, 2015, 23, 2445-2457.	1.4	42
56	Study of early stages of amyloid AÎ²13-23 formation using molecular dynamics simulation in implicit environments. Computational Biology and Chemistry, 2015, 56, 13-18.	1.1	7
57	Pyridine sulfonamide as a small key organic molecule for the potential treatment of type-II diabetes mellitus and Alzheimerâ€™s disease: In vitro studies against yeast Î±-glucosidase, acetylcholinesterase and butyrylcholinesterase. Bioorganic Chemistry, 2015, 63, 64-71.	2.0	51
58	Discovery of butyrylcholinesterase inhibitors among derivatives of azaphenothiazines. Journal of Enzyme Inhibition and Medicinal Chemistry, 2015, 30, 98-106.	2.5	19
59	Organocatalyzed solvent free an efficient novel synthesis of 2,4,5-trisubstituted imidazoles for Î±-glucosidase inhibition to treat diabetes. Bioorganic Chemistry, 2015, 58, 65-71.	2.0	51
60	Design and Synthesis of New Dual Binding Site Cholinesterase Inhibitors: in vitro Inhibition Studies with in silico Docking. Letters in Drug Design and Discovery, 2014, 11, 331-338.	0.4	23
61	Application of Computational Methods for the Design of BACE-1 Inhibitors: Validation of in Silico Modelling. International Journal of Molecular Sciences, 2014, 15, 5128-5139.	1.8	17
62	Synthesis and biological evaluation of novel oxadiazole derivatives: A new class of thymidine phosphorylase inhibitors as potential anti-tumor agents. Bioorganic and Medicinal Chemistry, 2014, 22, 1008-1015.	1.4	51
63	Novel synthesis of dihydropyrimidines for Î±-glucosidase inhibition to treat type 2 diabetes: In vitro biological evaluation and in silico docking. Bioorganic Chemistry, 2014, 54, 96-104.	2.0	49
64	Advances in GPCR Modeling Evaluated by the GPCR Dock 2013 Assessment: Meeting New Challenges. Structure, 2014, 22, 1120-1139.	1.6	149
65	New Tacrine Analogs as Acetylcholinesterase Inhibitors â€" Theoretical Study with Chemometric Analysis. Molecules, 2013, 18, 2878-2894.	1.7	12
66	Determination of lipophilicity of Î³â€butyrolactone derivatives with anticonvulsant and analgesic activity using micellar electrokinetic chromatography. Electrophoresis, 2013, 34, 3079-3085.	1.3	6
67	Synthesis and biological evaluation of 1,3,4-thiadiazole analogues as novel AChE and BuChE inhibitors. European Journal of Medicinal Chemistry, 2013, 62, 311-319.	2.6	61
68	Structure-Based Search for New Inhibitors of Cholinesterases. International Journal of Molecular Sciences, 2013, 14, 5608-5632.	1.8	200
69	2,3-Dihydro-1H-cyclopenta[b]quinoline Derivatives as Acetylcholinesterase Inhibitorsâ€"Synthesis, Radiolabeling and Biodistribution. International Journal of Molecular Sciences, 2012, 13, 10067-10090.	1.8	24
70	Synthesis, Biological Activity and Molecular Modeling of 4-Fluoro-N-[ï‰-(1,2,3,4-tetrahydroacridin-9-ylamino)-alkyl]-benzamide Derivatives as Cholinesterase Inhibitors. Arzneimittelforschung, 2012, 62, 655-660.	0.5	7
71	Search for anticonvulsant and analgesic active derivatives of dihydrofuran-2(3H)-one. Bioorganic and Medicinal Chemistry, 2012, 20, 6533-6544.	1.4	12
72	Synthesis and Biological Activity of New 2,3-dihydro-1H-cyclopenta[b]- quinoline Derivatives as Acetylcholinesterase Inhibitors. Letters in Drug Design and Discovery, 2012, 9, 645-654.	0.4	10

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73	Design, Synthesis and Evaluation of Novel 2â€(Aminoalkyl)â€isoindolineâ€1,3â€dione Derivatives as Dualâ€Binding Site Acetylcholinesterase Inhibitors. Archiv Der Pharmazie, 2012, 345, 509-516.	2.1	48
74	Dualâ€Acting Diether Derivatives of Piperidine and Homopiperidine with Histamine H₃ Receptor Antagonistic and Anticholinesterase Activity. Archiv Der Pharmazie, 2012, 345, 591-597.	2.1	25
75	Antiarrhythmic properties of phenylpiperazine derivatives of phenytoin with Î±1-adrenoceptor affinities. Bioorganic and Medicinal Chemistry, 2012, 20, 2290-2303.	1.4	29
76	Multi-Target-Directed Ligands in Alzheimer's Disease Treatment. Current Medicinal Chemistry, 2011, 18, 4949-4975.	1.2	219
77	Chromatographic and computational studies of the physicochemical properties of cholinesterase inhibitors â€" alkyl- and arylcarbamate derivatives ofN-benzylpiperidine andN-benzylpiperazine. Journal of Planar Chromatography - Modern TLC, 2010, 23, 359-364.	0.6	2
78	Novel alkyl- and arylcarbamate derivatives with N-benzylpiperidine and N-benzylpiperazine moieties as cholinesterases inhibitors. European Journal of Medicinal Chemistry, 2010, 45, 5602-5611.	2.6	21
79	Application of a Library of Artificial Receptors Formed by Selfâ€Organization of Nâ€Lipidated Peptides Immobilized on Cellulose for Preliminary Studies of Binding of Nâ€Phenylpiperazines. QSAR and Combinatorial Science, 2009, 28, 728-736.	1.5	3
80	Electrophoretically mediated microanalysis technique as a tool for the rapid screening of novel acetylcholinesterase inhibitors. Acta Poloniae Pharmaceutica, 2009, 66, 357-62.	0.3	8
81	Determination of lipophilicity of Î±â€(4â€phenylpiperazine) derivatives of Nâ€benzylamides using chromatographic and computational methods. Biomedical Chromatography, 2008, 22, 428-432.	0.8	10
82	Recent Developments in Cholinesterases Inhibitors for Alzheimers Disease Treatment. Current Medicinal Chemistry, 2007, 14, 2654-2679.	1.2	186
83	Investigation of lipophilicity of anticancer-active thioquinoline derivatives. Biomedical Chromatography, 2007, 21, 123-131.	0.8	32