

Magdalena J Ålusarż

List of Publications by Year in descending order

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Version: 2024-02-01

20
papers

253
citations

933447

10
h-index

940533

16
g-index

21
all docs

21
docs citations

21
times ranked

305
citing authors

#	ARTICLE	IF	CITATIONS
1	Analysis of interactions responsible for vasopressin binding to human neurohypophyseal hormone receptorsâ€”molecular dynamics study of the activated receptorâ€”vasopressinâ€”G1± systems. <i>Journal of Peptide Science</i> , 2006, 12, 180-189.	1.4	33
2	Molecular Docking-Based Study of Vasopressin Analogues Modified at Positions 2 and 3 with N-Methylphenylalanine: Influence on Receptor-Bound Conformations and Interactions with Vasopressin and Oxytocin Receptors. <i>Journal of Medicinal Chemistry</i> , 2006, 49, 2463-2469.	6.4	31
3	Molecular dynamics simulation of human neurohypophyseal hormone receptors complexed with oxytocinâ€”modeling of an activated state. <i>Journal of Peptide Science</i> , 2006, 12, 171-179.	1.4	27
4	Investigation of mechanism of desmopressin binding in vasopressin V2 receptor versus vasopressin V1a and oxytocin receptors: Molecular dynamics simulation of the agonist-bound state in the membraneâ€”aqueous system. <i>Biopolymers</i> , 2006, 81, 321-338.	2.4	24
5	Oxytocin-Gly-Lys-Arg: A Novel Cardiomyogenic Peptide. <i>PLoS ONE</i> , 2010, 5, e13643.	2.5	23
6	Modeling protein structures with the coarse-grained UNRES force field in the CASP14 experiment. <i>Journal of Molecular Graphics and Modelling</i> , 2021, 108, 108008.	2.4	17
7	Molecular Modeling of the Neurohypophyseal Receptor/Atosiban Complexes. <i>Protein and Peptide Letters</i> , 2003, 10, 295-302.	0.9	15
8	Conformational stability of the fullâ€”atom hexameric model of the ClpB chaperone from <i>Escherichia coli</i> . <i>Biopolymers</i> , 2010, 93, 47-60.	2.4	14
9	Conformational studies of vasopressin analogues modified with N-methylphenylalanine enantiomers in dimethyl sulfoxide solution. <i>Biopolymers</i> , 2006, 82, 603-614.	2.4	10
10	Exploring the Ligand Recognition Properties of the Human Vasopressin V1a Receptor Using QSAR and Molecular Modeling Studies. <i>Chemical Biology and Drug Design</i> , 2014, 83, 207-223.	3.2	10
11	Influence of bulky 3,3â€”diphenylalanine enantiomers replacing position 2 of AVP analogues on their conformations: NMR and molecular modeling studies. <i>European Journal of Medicinal Chemistry</i> , 2010, 45, 4065-4073.	5.5	8
12	Structure determination of UL49.5 transmembrane protein from bovine herpesvirus 1 by NMR spectroscopy and molecular dynamics. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2019, 1861, 926-938.	2.6	8
13	Investigation of cis/trans ratios of peptide bonds in AVP analogues containing N-methylphenylalanine enantiomers. <i>Journal of Peptide Science</i> , 2006, 12, 13-24.	1.4	7
14	Interactions of vasopressin and oxytocin receptors with vasopressin analogues substituted in position 2 with 3,3â€”diphenylalanine â€” a molecular docking study. <i>Journal of Peptide Science</i> , 2013, 19, 118-126.	1.4	6
15	Vasopressin V1a and V1b receptor modulators: a patent review (2012 â€” 2014). <i>Expert Opinion on Therapeutic Patents</i> , 2015, 25, 711-722.	5.0	6
16	PTD4 Peptide Increases Neural Viability in an In Vitro Model of Acute Ischemic Stroke. <i>International Journal of Molecular Sciences</i> , 2021, 22, 6086.	4.1	5
17	Molecular Dynamics Study of the Internal Water Molecules in Vasopressin and Oxytocin Receptors. <i>Protein and Peptide Letters</i> , 2009, 16, 342-350.	0.9	4
18	Theoretical studies, synthesis, and biological activity of 1-[(4-methylphenyl)sulfonyl]-5-oxo-2,3,4,5-tetrahydro-1H-1-benzazepine-4-carbonitrile (C9) as a non-peptide antagonist of the arginine vasopressin V1a and V2 receptors. <i>Medicinal Chemistry Research</i> , 2014, 23, 1581-1590.	2.4	4

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19	Investigation of the Effects of Primary Structure Modifications within the RRE Motif on the Conformation of Synthetic Bovine Herpesvirus 1â€€Encoded UL49.5 Protein Fragments. Chemistry and Biodiversity, 2021, 18, e2000883.	2.1	1
20	Molecular modeling study of the opioid receptor interactions with series of cyclic deltorphin analogues. Journal of Peptide Science, 2011, 17, 554-564.	1.4	0