

# Goran Å inko

## List of Publications by Year in descending order

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Version: 2024-02-01

38  
papers

1,424  
citations

361413

20  
h-index

315739

38  
g-index

43  
all docs

43  
docs citations

43  
times ranked

1810  
citing authors

#	ARTICLE	IF	CITATIONS
1	Use of connectivity index and simple topological parameters for estimating the inhibition potency of acetylcholinesterase. Saudi Pharmaceutical Journal, 2022, 30, 369-376.	2.7	3
2	Interaction of silver nanoparticles with plasma transport proteins: A systematic study on impacts of particle size, shape and surface functionalization. Chemico-Biological Interactions, 2021, 335, 109364.	4.0	13
3	Assessment of four organophosphorus pesticides as inhibitors of human acetylcholinesterase and butyrylcholinesterase. Scientific Reports, 2021, 11, 21486.	3.3	27
4	Interactions of Paraoxonase-1 with Pharmacologically Relevant Carbamates. Molecules, 2020, 25, 211.	3.8	8
5	Evaluation of high-affinity phenyltetrahydroisoquinoline aldoximes, linked through anti-triazoles, as reactivators of phosphorylated cholinesterases. Toxicology Letters, 2020, 321, 83-89.	0.8	13
6	Targeting organophosphorus compounds poisoning by novel quinuclidine-3 oximes: development of butyrylcholinesterase-based bioscavengers. Archives of Toxicology, 2020, 94, 3157-3171.	4.2	21
7	Enantioseparation, <i>in vitro</i> testing, and structural characterization of triple-binding reactivators of organophosphate-inhibited cholinesterases. Biochemical Journal, 2020, 477, 2771-2790.	3.7	12
8	Synthesis and In Vitro Screening of Novel Heterocyclic $\beta$ -D-Gluco- and $\beta$ -D-Galactoconjugates as Butyrylcholinesterase Inhibitors. Molecules, 2019, 24, 2833.	3.8	4
9	Assessment of scoring functions and <i>in silico</i> parameters for AChE-ligand interactions as a tool for predicting inhibition potency. Chemico-Biological Interactions, 2019, 308, 216-223.	4.0	23
10	Structural aspects of 4-aminoquinolines as reversible inhibitors of human acetylcholinesterase and butyrylcholinesterase. Chemico-Biological Interactions, 2019, 308, 101-109.	4.0	26
11	The estimation of oxime efficiency is affected by the experimental design of phosphorylated acetylcholinesterase reactivation. Toxicology Letters, 2018, 293, 222-228.	0.8	25
12	Oxime-assisted reactivation of tabun-inhibited acetylcholinesterase analysed by active site mutations. Toxicology, 2018, 406-407, 104-113.	4.2	28
13	Resorcinol-, catechol- and saligenin-based bronchodilating $\beta$ -2-agonists as inhibitors of human cholinesterase activity. Journal of Enzyme Inhibition and Medicinal Chemistry, 2017, 32, 789-797.	5.2	14
14	Interaction between the zebrafish ( <i>Danio rerio</i> ) organic cation transporter 1 (Oct1) and endo- and xenobiotics. Aquatic Toxicology, 2017, 187, 18-28.	4.0	17
15	The Lock is the Key: Development of Novel Drugs through Receptor Based Combinatorial Chemistry. Acta Chimica Slovenica, 2017, 64, 15-39.	0.6	9
16	Design and synthesis of N-substituted-2-hydroxyiminoacetamides and interactions with cholinesterases. Chemico-Biological Interactions, 2016, 259, 122-132.	4.0	27
17	Pyridoxal oxime derivative potency to reactivate cholinesterases inhibited by organophosphorus compounds. Toxicology Letters, 2016, 262, 114-122.	0.8	14
18	Response of biochemical biomarkers in the aquatic crustacean <i>Daphnia magna</i> exposed to silver nanoparticles. Environmental Science and Pollution Research, 2015, 22, 19990-19999.	5.3	59

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19	Alteration of cholinesterase activity as possible mechanism of silver nanoparticle toxicity. <i>Environmental Science and Pollution Research</i> , 2014, 21, 1391-1400.	5.3	41
20	Enzyme-catalyzed cascade synthesis of hydroxyiminoacetamides. <i>Tetrahedron Letters</i> , 2014, 55, 4338-4341.	1.4	5
21	Inactivation of cholinesterases by silver and gold ions in vitro. <i>Open Chemistry</i> , 2013, 11, 935-944.	1.9	3
22	Peripheral site and acyl pocket define selective inhibition of mouse butyrylcholinesterase by two bis carbamates. <i>Archives of Biochemistry and Biophysics</i> , 2013, 529, 140-145.	3.0	13
23	<i>In vitro</i> inhibition of blood cholinesterase activities from cattle by triazole fungicides. <i>Caryologia</i> , 2013, 66, 346-350.	0.3	9
24	Metaproterenol, Isoproterenol, and Their Bisdimethylcarbamate Derivatives as Human Cholinesterase Inhibitors. <i>Journal of Medicinal Chemistry</i> , 2012, 55, 6716-6723.	6.4	37
25	Mechanism of stereoselective interaction between butyrylcholinesterase and ethopropazine enantiomers. <i>Biochimie</i> , 2011, 93, 1797-1807.	2.6	26
26	Structural aspects of flavonoids as inhibitors of human butyrylcholinesterase. <i>European Journal of Medicinal Chemistry</i> , 2010, 45, 186-192.	5.5	154
27	Pseudo-catalytic scavenging: Searching for a suitable reactivator of phosphorylated butyrylcholinesterase. <i>Chemico-Biological Interactions</i> , 2010, 187, 167-171.	4.0	53
28	Interactions of pyridinium oximes with acetylcholinesterase. <i>Chemico-Biological Interactions</i> , 2010, 187, 172-176.	4.0	28
29	Cholinesterase Interactions with Oximes. <i>Current Bioactive Compounds</i> , 2010, 6, 9-15.	0.5	7
30	Oximes: Reactivators of phosphorylated acetylcholinesterase and antidotes in therapy against tabun poisoning. <i>Chemico-Biological Interactions</i> , 2008, 175, 173-179.	4.0	37
31	Structure-Activity Approach in the Reactivation of Tabun-Phosphorylated Human Acetylcholinesterase with Bispyridinium para-Aldoximes. <i>Arhiv Za Higijenu Rada I Toksikologiju</i> , 2007, 58, 201-209.	0.7	26
32	Limitation of the Ellman method: Cholinesterase activity measurement in the presence of oximes. <i>Analytical Biochemistry</i> , 2007, 370, 223-227.	2.4	103
33	para- and ortho-Pyridinium aldoximes in reaction with acetylthiocholine. <i>FEBS Letters</i> , 2006, 580, 3167-3172.	2.8	44
34	Preparative HPLC separation of bambuterol enantiomers and stereoselective inhibition of human cholinesterases. <i>Analytical and Bioanalytical Chemistry</i> , 2006, 385, 1513-1519.	3.7	32
35	(42) Structure-inhibition relationships in the interaction of butyrylcholinesterase with bambuterol, haloxon and their leaving groups. <i>Chemico-Biological Interactions</i> , 2005, 157-158, 421-423.	4.0	3
36	Molar absorption coefficients for the reduced Ellman reagent: reassessment. <i>Analytical Biochemistry</i> , 2003, 312, 224-227.	2.4	435

#	ARTICLE	IF	CITATIONS
37	Separation, Conformation in Solution and Absolute Configuration of Ethopropazine Enantiomers. Enantiomer, 2002, 7, 149-156.	0.5	3
38	Kinetic Model of Ethopropazine Interaction with Horse Serum Butyrylcholinesterase and Its Docking into the Active Site. Archives of Biochemistry and Biophysics, 2002, 398, 23-31.	3.0	20