

Heini W Dirr

List of Publications by Year in Descending Order

Source: <https://exaly.com/author-pdf/7877691/heini-w-dirr-publications-by-year.pdf>

Version: 2024-04-25

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

98
papers

2,960
citations

28
h-index

52
g-index

98
ext. papers

3,133
ext. citations

3.7
avg. IF

4.58
L-index

#	Paper	IF	Citations
98	An empirical and theoretical description of <i>Schistosoma japonicum</i> glutathione transferase inhibition by bromosulphophthalein and indanyloxyacetic acid 94. <i>Journal of Molecular Structure</i> , 2021 , 1223, 128892	3.4	0
97	Molecular basis of inhibition of <i>Schistosoma japonicum</i> glutathione transferase by ellagic acid: Insights into biophysical and structural studies. <i>Molecular and Biochemical Parasitology</i> , 2020 , 240, 111319 ⁹	1.9	2
96	Drug susceptibility and replication capacity of a rare HIV-1 subtype C protease hinge region variant. <i>Antiviral Therapy</i> , 2019 , 24, 333-342	1.6	1
95	Double trouble? Gag in conjunction with double insert in HIV protease contributes to reduced DRV susceptibility. <i>Biochemical Journal</i> , 2019 , 476, 375-384	3.8	2
94	An update on the biophysical character of the human eukaryotic elongation factor 1 beta: Perspectives from interaction with elongation factor 1 gamma. <i>Journal of Molecular Recognition</i> , 2018 , 31, e2708	2.6	3
93	Molecular dynamics and ligand docking of a hinge region variant of South African HIV-1 subtype C protease. <i>Journal of Molecular Graphics and Modelling</i> , 2018 , 82, 1-11	2.8	10
92	A Phosphomimetic Study Implicates Ser557 in Regulation of FOXP2 DNA Binding. <i>Protein Journal</i> , 2018 , 37, 311-323	3.9	3
91	Overexpression, Purification and Functional Characterisation of Wild-Type HIV-1 Subtype C Protease and Two Variants Using a Thioredoxin and His-Tag Protein Fusion System. <i>Protein Journal</i> , 2018 , 37, 369-379	3.9	3
90	The forkhead domain hinge-loop plays a pivotal role in DNA binding and transcriptional activity of FOXP2. <i>Biological Chemistry</i> , 2018 , 399, 881-893	4.5	3
89	The study of degradation mechanisms of glyco-engineered plant produced anti-rabies monoclonal antibodies E559 and 62-71-3. <i>PLoS ONE</i> , 2018 , 13, e0209373	3.7	2
88	Structural and biochemical characterization of <i>Plasmodium falciparum</i> Hsp70-x reveals functional versatility of its C-terminal EEVN motif. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018 , 86, 1189-1201	4.2	26
87	A conserved cation binding site in the DNA binding domain of forkhead box transcription factors regulates DNA binding by FOXP2. <i>Archives of Biochemistry and Biophysics</i> , 2018 , 657, 56-64	4.1	2
86	The effects of mutating Tyr9 and Arg15 on the structure, stability, conformational dynamics and mechanism of GSTA3-3. <i>Biophysical Chemistry</i> , 2017 , 224, 40-48	3.5	3
85	The FOXP2 forkhead domain binds to a variety of DNA sequences with different rates and affinities. <i>Journal of Biochemistry</i> , 2017 , 162, 45-54	3.1	5
84	(-)-Epigallocatechin-3-Gallate Inhibits the Chaperone Activity of <i>Plasmodium falciparum</i> Hsp70 Chaperones and Abrogates Their Association with Functional Partners. <i>Molecules</i> , 2017 , 22,	4.8	23
83	Energetics of Glutathione Binding to Human Eukaryotic Elongation Factor 1 Gamma: Isothermal Titration Calorimetry and Molecular Dynamics Studies. <i>Protein Journal</i> , 2016 , 35, 448-458	3.9	2
82	The Lyssavirus glycoprotein: A key to cross-immunity. <i>Virology</i> , 2016 , 498, 250-256	3.6	6

81	Plasmodium falciparum Hsp70-z, an Hsp110 homologue, exhibits independent chaperone activity and interacts with Hsp70-1 in a nucleotide-dependent fashion. <i>Cell Stress and Chaperones</i> , 2016 , 21, 499-513	4.13	29
80	A Single Amino Acid in the Hinge Loop Region of the FOXP Forkhead Domain is Significant for Dimerisation. <i>Protein Journal</i> , 2015 , 34, 111-21	3.9	10
79	The kinetics of the substitution of coordinated H ₂ O on Co(III) by cyanide in aquacobalamin (vitamin B12a) and in a corrole analogue. <i>Inorganic Chemistry Communication</i> , 2015 , 57, 15-17	3.1	3
78	Overexpression, Purification and Characterisation of the Plasmodium falciparum Hsp70-z (PfHsp70-z) Protein. <i>PLoS ONE</i> , 2015 , 10, e0129445	3.7	28
77	Purification and characterisation of recombinant human eukaryotic elongation factor 1 gamma. <i>Protein Expression and Purification</i> , 2014 , 99, 70-7	2	12
76	Phosphorylation- and nucleotide-binding-induced changes to the stability and hydrogen exchange patterns of JNK1 β provide insight into its mechanisms of activation. <i>Journal of Molecular Biology</i> , 2014 , 426, 3569-89	6.5	6
75	A Lys-Trp cation- π interaction mediates the dimerization and function of the chloride intracellular channel protein 1 transmembrane domain. <i>Biochemistry</i> , 2014 , 53, 57-67	3.2	19
74	A conserved cationic motif enhances membrane binding and insertion of the chloride intracellular channel protein 1 transmembrane domain. <i>European Biophysics Journal</i> , 2014 , 43, 405-14	1.9	5
73	The isomerization of β -androstene-3,17-dione by the human glutathione transferase A3-3 proceeds via a conjugated heteroannular diene intermediate. <i>Journal of Biological Chemistry</i> , 2014 , 289, 32243-32252	5.4	4
72	Amide hydrogen exchange in HIV-1 subtype B and C proteases--insights into reduced drug susceptibility and dimer stability. <i>FEBS Journal</i> , 2014 , 281, 5395-410	5.7	5
71	Polymorphic Diversity: N-Phenylbenzamide as a Possible Polymorphophore. <i>Crystal Growth and Design</i> , 2013 , 13, 3463-3474	3.5	12
70	F99 is critical for dimerization and activation of South African HIV-1 subtype C protease. <i>Protein Journal</i> , 2013 , 32, 560-7	3.9	2
69	JNK1 β is phosphorylated during expression in E. coli and in vitro by MKK4 at three identical novel sites. <i>Biochemical and Biophysical Research Communications</i> , 2013 , 432, 683-8	3.4	4
68	High yield purification of JNK1 β and activation by in vitro reconstitution of the MEKK1-MKK4-JNK MAPK phosphorylation cascade. <i>Protein Expression and Purification</i> , 2013 , 87, 87-99	2	13
67	S-Nitrosation destabilizes glutathione transferase P1-1. <i>Biochemistry</i> , 2013 , 52, 9394-402	3.2	7
66	Membrane mimetics induce helix formation and oligomerization of the chloride intracellular channel protein 1 transmembrane domain. <i>Biochemistry</i> , 2013 , 52, 2739-49	3.2	6
65	Structural insights into the South African HIV-1 subtype C protease: impact of hinge region dynamics and flap flexibility in drug resistance. <i>Journal of Biomolecular Structure and Dynamics</i> , 2013 , 31, 1370-80	3.6	29
64	S-nitrosation of glutathione transferase p1-1 is controlled by the conformation of a dynamic active site helix. <i>Journal of Biological Chemistry</i> , 2013 , 288, 14973-84	5.4	8

63	Role of arginine 29 and glutamic acid 81 interactions in the conformational stability of human chloride intracellular channel 1. <i>Biochemistry</i> , 2012 , 51, 7854-62	3.2	13
62	Role of individual histidines in the pH-dependent global stability of human chloride intracellular channel 1. <i>Biochemistry</i> , 2012 , 51, 995-1004	3.2	19
61	A conserved interdomain interaction is a determinant of folding cooperativity in the GST fold. <i>Biochemistry</i> , 2011 , 50, 7067-75	3.2	4
60	Energetics of ligand binding to human glutathione transferase A1-1: Tyr-9 associated localisation of the C-terminal helix is ligand-dependent. <i>Biophysical Chemistry</i> , 2011 , 156, 153-8	3.5	3
59	Class Pi glutathione transferase unfolds via a dimeric and not monomeric intermediate: functional implications for an unstable monomer. <i>Biochemistry</i> , 2010 , 49, 5074-81	3.2	17
58	Stability of the domain interface contributes towards the catalytic function at the H-site of class alpha glutathione transferase A1-1. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2010 , 1804, 2228-33	4	14
57	Arginine 15 stabilizes an S(N)Ar reaction transition state and the binding of anionic ligands at the active site of human glutathione transferase A1-1. <i>Biophysical Chemistry</i> , 2010 , 146, 118-25	3.5	17
56	The role of a topologically conserved isoleucine in glutathione transferase structure, stability and function. <i>Acta Crystallographica Section F: Structural Biology Communications</i> , 2010 , 66, 776-80		3
55	Structural dynamics of soluble chloride intracellular channel protein CLIC1 examined by amide hydrogen-deuterium exchange mass spectrometry. <i>Biochemistry</i> , 2009 , 48, 8413-21	3.2	30
54	Formation of an unfolding intermediate state of soluble chloride intracellular channel protein CLIC1 at acidic pH. <i>Biochemistry</i> , 2008 , 47, 11674-81	3.2	38
53	Stability and unfolding of reduced Escherichia coli glutaredoxin 2: a monomeric structural homologue of the glutathione transferase family. <i>Biochemistry</i> , 2008 , 47, 10801-8	3.2	2
52	Active-site mutations in the South african human immunodeficiency virus type 1 subtype C protease have a significant impact on clinical inhibitor binding: kinetic and thermodynamic study. <i>Journal of Virology</i> , 2008 , 82, 11476-9	6.6	30
51	Characterization of the binding of 8-anilino-naphthalene sulfonate to rat class Mu GST M1-1. <i>Biophysical Chemistry</i> , 2008 , 137, 100-4	3.5	11
50	Double mutation at the subunit interface of glutathione transferase rGSTM1-1 results in a stable, folded monomer. <i>Biochemistry</i> , 2006 , 45, 2267-73	3.2	26
49	The intersubunit lock-and-key motif in human glutathione transferase A1-1: role of the key residues Met51 and Phe52 in function and dimer stability. <i>Biochemical Journal</i> , 2006 , 393, 523-8	3.8	18
48	Influence of the dimer interface on glutathione transferase structure and dynamics revealed by amide H/D exchange mass spectrometry. <i>Biochemistry</i> , 2005 , 44, 10605-12	3.2	18
47	Tertiary interactions stabilise the C-terminal region of human glutathione transferase A1-1: a crystallographic and calorimetric study. <i>Journal of Molecular Biology</i> , 2005 , 349, 825-38	6.5	27
46	A conserved N-capping motif contributes significantly to the stabilization and dynamics of the C-terminal region of class Alpha glutathione S-transferases. <i>Journal of Biological Chemistry</i> , 2005 , 280, 19480-7	5.4	16

45	Characterization of bromosulphophthalein binding to human glutathione S-transferase A1-1: thermodynamics and inhibition kinetics. <i>Biochemical Journal</i> , 2004 , 382, 703-9	3.8	23
44	The role of an evolutionarily conserved cis-proline in the thioredoxin-like domain of human class Alpha glutathione transferase A1-1. <i>Biochemical Journal</i> , 2003 , 372, 241-6	3.8	30
43	Residue 219 impacts on the dynamics of the C-terminal region in glutathione transferase A1-1: implications for stability and catalytic and ligandin functions. <i>Biochemistry</i> , 2003 , 42, 15326-32	3.2	24
42	Protease inhibition in African subtypes of HIV-1. <i>AIDS Reviews</i> , 2003 , 5, 165-71	1.5	32
41	Impact of domain interchange on conformational stability and equilibrium folding of chimeric class micro glutathione transferases. <i>Protein Science</i> , 2002 , 11, 2208-17	6.3	16
40	Thermodynamics of the ligandin function of human class Alpha glutathione transferase A1-1: energetics of organic anion ligand binding. <i>Biochemical Journal</i> , 2002 , 363, 341-6	3.8	24
39	Thermodynamics of the ligandin function of human class Alpha glutathione transferase A1-1: energetics of organic anion ligand binding. <i>Biochemical Journal</i> , 2002 , 363, 341-346	3.8	34
38	Molecular recognition at the dimer interface of a class mu glutathione transferase: role of a hydrophobic interaction motif in dimer stability and protein function. <i>Biochemistry</i> , 2002 , 41, 14238-47	3.2	37
37	The cochaperone murine stress-inducible protein 1: overexpression, purification, and characterization. <i>Protein Expression and Purification</i> , 2001 , 21, 462-9	2	14
36	Heat shock cognate protein 70 chaperone-binding site in the co-chaperone murine stress-inducible protein 1 maps to within three consecutive tetratricopeptide repeat motifs. <i>Biochemical Journal</i> , 2000 , 345 Pt 3, 645-51	3.8	22
35	Electrostatic interactions affecting the active site of class Sigma glutathione S-transferase. <i>Biochemical Journal</i> , 2000 , 347, 193	3.8	8
34	Heat shock cognate protein 70 chaperone-binding site in the co-chaperone murine stress-inducible protein 1 maps to within three consecutive tetratricopeptide repeat motifs. <i>Biochemical Journal</i> , 2000 , 345, 645-651	3.8	38
33	Electrostatic interactions affecting the active site of class Sigma glutathione S-transferase. <i>Biochemical Journal</i> , 2000 , 347, 193-197	3.8	26
32	Domain-domain interface packing at conserved Trp-20 in class alpha glutathione transferase impacts on protein stability. <i>BBA - Proteins and Proteomics</i> , 2000 , 1478, 325-32		20
31	Analysis of the levels of conservation of the J domain among the various types of DnaJ-like proteins. <i>Cell Stress and Chaperones</i> , 2000 , 5, 347-58	4	57
30	The hydrophobic lock-and-key intersubunit motif of glutathione transferase A1-1: implications for catalysis, ligandin function and stability. <i>FEBS Letters</i> , 2000 , 465, 169-72	3.8	48
29	Equilibrium folding of dimeric class mu glutathione transferases involves a stable monomeric intermediate. <i>Biochemistry</i> , 2000 , 39, 12336-44	3.2	55
28	Folding and assembly of dimeric human glutathione transferase A1-1. <i>Biochemistry</i> , 1999 , 38, 16686-94	3.2	48

27	Role of the C-terminal helix 9 in the stability and ligandin function of class alpha glutathione transferase A1-1. <i>Biochemistry</i> , 1999 , 38, 15631-40	3.2	52
26	Aflatoxin B1 and sulphobromophthalein binding to the dimeric human glutathione S-transferase A1-1: a fluorescence spectroscopic analysis. <i>FEBS Journal</i> , 1998 , 257, 434-42		15
25	Class sigma glutathione transferase unfolds via a dimeric and a monomeric intermediate: impact of subunit interface on conformational stability in the superfamily. <i>Biochemistry</i> , 1998 , 37, 15534-41	3.2	38
24	Equilibrium and kinetic unfolding properties of dimeric human glutathione transferase A1-1. <i>Biochemistry</i> , 1998 , 37, 5320-8	3.2	71
23	A topologically conserved aliphatic residue in alpha-helix 6 stabilizes the hydrophobic core in domain II of glutathione transferases and is a structural determinant for the unfolding pathway. <i>Biochemical Journal</i> , 1998 , 336 (Pt 2), 413-8	3.8	10
22	Conformational stability of pGEX-expressed <i>Schistosoma japonicum</i> glutathione S-transferase: a detoxification enzyme and fusion-protein affinity tag. <i>Protein Science</i> , 1997 , 6, 399-406	6.3	113
21	Effect of glutathione, glutathione sulphonate and S-hexylglutathione on the conformational stability of class pi glutathione S-transferase. <i>FEBS Letters</i> , 1996 , 391, 313-6	3.8	13
20	Determination of a binding site for a non-substrate ligand in mammalian cytosolic glutathione S-transferases by means of fluorescence-resonance energy transfer. <i>FEBS Journal</i> , 1996 , 241, 484-8		28
19	Class-pi glutathione S-transferase is unable to regain its native conformation after oxidative inactivation by hydrogen peroxide. <i>FEBS Journal</i> , 1996 , 242, 301-7		12
18	Porcine class pi glutathione S-transferase: anionic ligand binding and conformational analysis. <i>BBA - Proteins and Proteomics</i> , 1995 , 1247, 225-30		26
17	X-ray structure methods for glutathione binding. <i>Methods in Enzymology</i> , 1995 , 251, 243-54	1.7	4
16	Conformational stability of Cys45-alkylated and hydrogen peroxide-oxidised glutathione S-transferase. <i>FEBS Letters</i> , 1995 , 371, 94-8	3.8	9
15	Native Dimer Stabilizes the Subunit Tertiary Structure of Porcine Class pi Glutathione S-transferase. <i>FEBS Journal</i> , 1995 , 230, 614-620		46
14	X-ray crystal structures of cytosolic glutathione S-transferases. Implications for protein architecture, substrate recognition and catalytic function. <i>FEBS Journal</i> , 1994 , 220, 645-61		361
13	Refined crystal structure of porcine class Pi glutathione S-transferase (pGST P1-1) at 2.1 A resolution. <i>Journal of Molecular Biology</i> , 1994 , 243, 72-92	6.5	86
12	X-ray crystal structures of cytosolic glutathione S-transferases 1994 , 57-73		
11	Structure determination and refinement of human alpha class glutathione transferase A1-1, and a comparison with the Mu and Pi class enzymes. <i>Journal of Molecular Biology</i> , 1993 , 232, 192-212	6.5	428
10	Three-dimensional structure of class pi glutathione S-transferase from human placenta in complex with S-hexylglutathione at 2.8 A resolution. <i>Journal of Molecular Biology</i> , 1992 , 227, 214-26	6.5	261

9	Mutational substitution of residues implicated by crystal structure in binding the substrate glutathione to human glutathione S-transferase pi. <i>Journal of Molecular Biology</i> , 1992 , 226, 319-22	6.5	57
8	Class pi glutathione S-transferase from pig lung. Purification, biochemical characterization, primary structure and crystallization. <i>FEBS Journal</i> , 1991 , 196, 693-8		62
7	Equilibrium unfolding of class pi glutathione S-transferase. <i>Biochemical and Biophysical Research Communications</i> , 1991 , 180, 294-300	3.4	58
6	Purification and properties of an esterase from Cucurbita maxima fruit tissue. <i>Phytochemistry</i> , 1989 , 28, 379-383	4	11
5	Accessibility of the tryptophan residues and flavin prosthetic group of Cyclopiazionate oxidocyclase to solvent studied by fluorescence quenching. <i>International Journal of Biochemistry & Cell Biology</i> , 1988 , 20, 109-111		1
4	Purification and partial characterization of the glutathione S-transferase of rat erythrocytes. <i>BBA - Proteins and Proteomics</i> , 1988 , 957, 173-7		6
3	Solvent effects on the spectroscopic properties of aflatoxin B1. <i>International Journal of Biochemistry & Cell Biology</i> , 1987 , 19, 1137-40		9
2	Characterization of the aflatoxin B1-binding site of rat albumin. <i>BBA - Proteins and Proteomics</i> , 1987 , 913, 300-7		3
1	Aflatoxin B1 transport in rat blood plasma. Binding to albumin in vivo and in vitro and spectrofluorimetric studies into the nature of the interaction. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 1986 , 881, 383-90	4	20