

# Margot Ernst

## List of Publications by Year in Descending Order

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**Version:** 2024-04-24

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

64  
papers

1,376  
citations

19  
h-index

36  
g-index

75  
ext. papers

1,663  
ext. citations

5.9  
avg, IF

4.45  
L-index

#	Paper	IF	Citations
64	6-Containing GABA Receptors: Functional Roles and Therapeutic Potentials.. <i>Pharmacological Reviews</i> , <b>2022</b> , 74, 238-270	22.5	1
63	A Benzodiazepine Ligand with Improved GABA Receptor 5-Subunit Selectivity Driven by Interactions with Loop C. <i>Molecular Pharmacology</i> , <b>2021</b> , 99, 39-48	4.3	0
62	Molecular basis of mood and cognitive adverse events elucidated via a combination of pharmacovigilance data mining and functional enrichment analysis. <i>Archives of Toxicology</i> , <b>2020</b> , 94, 2829-2845 <sup>0</sup>	5.8	5
61	Allosteric GABA Receptor Modulators-A Review on the Most Recent Heterocyclic Chemotypes and Their Synthetic Accessibility. <i>Molecules</i> , <b>2020</b> , 25,	4.8	13
60	GABA Receptor Ligands Often Interact with Binding Sites in the Transmembrane Domain and in the Extracellular Domain-Can the Promiscuity Code Be Cracked?. <i>International Journal of Molecular Sciences</i> , <b>2020</b> , 21,	6.3	6
59	A novel de novo variant of GABRA1 causes increased sensitivity for GABA in vitro. <i>Scientific Reports</i> , <b>2020</b> , 10, 2379	4.9	5
58	Two Distinct Populations of $\alpha 5$ -Containing GABA <sub>A</sub> -Receptors in Rat Cerebellum. <i>Frontiers in Synaptic Neuroscience</i> , <b>2020</b> , 12, 591129	3.5	6
57	Structure-Guided Computational Methods Predict Multiple Distinct Binding Modes for Pyrazoloquinolinones in GABA Receptors. <i>Frontiers in Neuroscience</i> , <b>2020</b> , 14, 611953	5.1	1
56	Defined concatenated $\alpha 1\beta 2$ GABA receptor constructs reveal dual action of pyrazoloquinolinone allosteric modulators. <i>Bioorganic and Medicinal Chemistry</i> , <b>2019</b> , 27, 3167-3178	3.4	6
55	Variations on a scaffold - Novel GABA receptor modulators. <i>European Journal of Medicinal Chemistry</i> , <b>2019</b> , 180, 340-349	6.8	1
54	Trigeminal neuropathic pain development and maintenance in rats are suppressed by a positive modulator of $\beta$ GABA receptors. <i>European Journal of Pain</i> , <b>2019</b> , 23, 973-984	3.7	18
53	Phenotypic variability of GABRA1-related epilepsy in monozygotic twins. <i>Annals of Clinical and Translational Neurology</i> , <b>2019</b> , 6, 2317-2322	5.3	3
52	A Novel Drug Target for Migraine: The GABA <sub>A</sub> Receptor $\beta$ Subtype in Trigeminal Ganglia. <i>FASEB Journal</i> , <b>2019</b> , 33, lb78	0.9	
51	Design and Synthesis of Novel Deuterated Ligands Functionally Selective for the $\beta$ Aminobutyric Acid Type A Receptor (GABA <sub>A</sub> ) $\beta$ Subtype with Improved Metabolic Stability and Enhanced Bioavailability. <i>Journal of Medicinal Chemistry</i> , <b>2018</b> , 61, 2422-2446	8.3	26
50	Cerebellar $\beta$ subunit-containing GABA receptors: a novel therapeutic target for disrupted prepulse inhibition in neuropsychiatric disorders. <i>British Journal of Pharmacology</i> , <b>2018</b> , 175, 2414-2427	8.6	20
49	Negative allosteric modulation of alpha 5-containing GABA receptors engenders antidepressant-like effects and selectively prevents age-associated hyperactivity in tau-depositing mice. <i>Psychopharmacology</i> , <b>2018</b> , 235, 1151-1161	4.7	13
48	Expedient synthesis of polyacetylenic water hemlock toxins and their effects on the major GABA receptor isoform. <i>Chemical Communications</i> , <b>2018</b> , 54, 2008-2011	5.8	7

47	The Benzodiazepine Binding Sites of GABA Receptors. <i>Trends in Pharmacological Sciences</i> , <b>2018</b> , 39, 659-671	6.1	89
46	Diversity matters: combinatorial information coding by GABA receptor subunits during spatial learning and its allosteric modulation. <i>Cellular Signalling</i> , <b>2018</b> , 50, 142-159	4.9	4
45	Engineered Flumazenil Recognition Site Provides Mechanistic Insight Governing Benzodiazepine Modulation in GABA Receptors. <i>ACS Chemical Biology</i> , <b>2018</b> , 13, 2040-2047	4.9	7
44	SAR-Guided Scoring Function and Mutational Validation Reveal the Binding Mode of CGS-8216 at the $\alpha 5/\alpha 1$ - Benzodiazepine Site. <i>Journal of Chemical Information and Modeling</i> , <b>2018</b> , 58, 1682-1696	6.1	3
43	The $\beta$ subunit-containing GABA receptor: A novel drug target for inhibition of trigeminal activation. <i>Neuropharmacology</i> , <b>2018</b> , 140, 1-13	5.5	14
42	Attaining in vivo selectivity of positive modulation of $\beta\gamma$ GABA receptors in rats: A hard task!. <i>European Neuropsychopharmacology</i> , <b>2018</b> , 28, 903-914	1.2	3
41	The cerebellar $\beta$ subunit-containing GABAA receptor: A novel therapeutic target for disrupted prepulse inhibition in neuropsychiatric disorders. <i>Proceedings for Annual Meeting of the Japanese Pharmacological Society</i> , <b>2018</b> , WCP2018, PO3-1-95	0	1
40	A Novel Target for Migraine Therapy: the $\beta$ Subunit-Containing GABAA Receptor. <i>Proceedings for Annual Meeting of the Japanese Pharmacological Society</i> , <b>2018</b> , WCP2018, PO1-1-136	0	
39	Towards functional selectivity for $\beta\gamma\delta$ GABA receptors: a series of novel pyrazoloquinolinones. <i>British Journal of Pharmacology</i> , <b>2018</b> , 175, 419-428	8.6	18
38	Stereoselective Synthesis of the Isomers of Notoincisol A: Assignment of the Absolute Configuration of this Natural Product and Biological Evaluation. <i>Journal of Natural Products</i> , <b>2018</b> , 81, 2419-2428	4.9	0
37	Different Benzodiazepines Bind with Distinct Binding Modes to GABA Receptors. <i>ACS Chemical Biology</i> , <b>2018</b> , 13, 2033-2039	4.9	12
36	Pharmacological and antihyperalgesic properties of the novel $\alpha 3$ preferring GABA receptor ligand MP-III-024. <i>Brain Research Bulletin</i> , <b>2017</b> , 131, 62-69	3.9	20
35	Molecular tools for GABA receptors: High affinity ligands for $\alpha 1$ -containing subtypes. <i>Scientific Reports</i> , <b>2017</b> , 7, 5674	4.9	19
34	$\beta$ subunits in GABA receptors are dispensable for GABA and diazepam action. <i>Scientific Reports</i> , <b>2017</b> , 7, 15498	4.9	15
33	Ester to amide substitution improves selectivity, efficacy and kinetic behavior of a benzodiazepine positive modulator of GABA receptors containing the $\beta$ subunit. <i>European Journal of Pharmacology</i> , <b>2016</b> , 791, 433-443	5.3	18
32	Synthesis and Characterization of a Novel $\gamma$ -Aminobutyric Acid Type A (GABA) Receptor Ligand That Combines Outstanding Metabolic Stability, Pharmacokinetics, and Anxiolytic Efficacy. <i>Journal of Medicinal Chemistry</i> , <b>2016</b> , 59, 10800-10806	8.3	34
31	First In Vivo Testing of Compounds Targeting Group 3 Medulloblastomas Using an Implantable Microdevice as a New Paradigm for Drug Development. <i>Journal of Biomedical Nanotechnology</i> , <b>2016</b> , 12, 1297-302	4	23
30	Targeting the $\gamma$ -Aminobutyric Acid A Receptor $\alpha 1$ Subunit in Airway Smooth Muscle to Alleviate Bronchoconstriction. <i>American Journal of Respiratory Cell and Molecular Biology</i> , <b>2016</b> , 54, 546-53	5.7	16

29	Structural Studies of GABAA Receptor Binding Sites: Which Experimental Structure Tells us What?. <i>Frontiers in Molecular Neuroscience</i> , <b>2016</b> , 9, 44	6.1	57
28	Mutagenesis and computational docking studies support the existence of a histamine binding site at the extracellular $\beta+\beta$ - interface of homooligomeric $\beta$ GABAA receptors. <i>Neuropharmacology</i> , <b>2016</b> , 108, 252-63	5.5	9
27	Development of GABAA Receptor Subtype-Selective Imidazobenzodiazepines as Novel Asthma Treatments. <i>Molecular Pharmaceutics</i> , <b>2016</b> , 13, 2026-38	5.6	18
26	Metal-assisted synthesis of unsymmetrical magnolol and honokiol analogs and their biological assessment as GABAA receptor ligands. <i>Bioorganic and Medicinal Chemistry Letters</i> , <b>2015</b> , 25, 400-3	2.9	7
25	GABAA receptor subtypes: structural variety raises hope for new therapy concepts. <i>E-Neuroforum</i> , <b>2015</b> , 6, 97-103		2
24	GABAA $\alpha$ Rezeptorsubtypen: Strukturelle Vielfalt gibt Hoffnung auf neue Therapiekonzepte. <i>Neuroforum</i> , <b>2015</b> , 21, 144-151	0.7	
23	Accelerated discovery of novel benzodiazepine ligands by experiment-guided virtual screening. <i>ACS Chemical Biology</i> , <b>2014</b> , 9, 1854-9	4.9	17
22	Unexpected Properties of $\alpha$ Containing GABAA Receptors in Response to Ligands Interacting with the $\beta$ Site. <i>Neurochemical Research</i> , <b>2014</b> , 39, 1057-1067	4.6	11
21	A pentasymmetric open channel blocker for Cys-loop receptor channels. <i>PLoS ONE</i> , <b>2014</b> , 9, e106688	3.7	6
20	Comparing the high affinity benzodiazepine binding site with the homologous $\alpha$ GS 9895 $\beta$ ite in GABA-A receptors (1059.1). <i>FASEB Journal</i> , <b>2014</b> , 28, 1059.1	0.9	
19	Subtype selectivity of $\beta$ Site ligands of GABAA receptors: identification of the first highly specific positive modulators at $\beta 2/3$ receptors. <i>British Journal of Pharmacology</i> , <b>2013</b> , 169, 384-99	8.6	41
18	Identification of novel positive allosteric modulators and null modulators at the GABAA receptor $\beta$ Site interface. <i>British Journal of Pharmacology</i> , <b>2013</b> , 169, 371-83	8.6	39
17	Molecular analysis of the site for 2-arachidonylglycerol (2-AG) on the $\beta$ ubunit of GABA(A) receptors. <i>Journal of Neurochemistry</i> , <b>2013</b> , 126, 29-36	6	19
16	A novel GABA(A) receptor pharmacology: drugs interacting with the $\beta$ (+) $\beta$ ) interface. <i>British Journal of Pharmacology</i> , <b>2012</b> , 166, 476-85	8.6	62
15	Understanding subtype-selective allosteric modulation of GABAareceptors. <i>BMC Pharmacology &amp; Toxicology</i> , <b>2012</b> , 13,	2.6	78
14	Diazepam-bound GABAA receptor models identify new benzodiazepine binding-site ligands. <i>Nature Chemical Biology</i> , <b>2012</b> , 8, 455-64	11.7	150
13	Pentameric ligand-gated ion channel ELIC is activated by GABA and modulated by benzodiazepines. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2012</b> , 109, E3028-34	11.5	106
12	The GABAA receptor alpha+beta- interface: a novel target for subtype selective drugs. <i>Journal of Neuroscience</i> , <b>2011</b> , 31, 870-7	6.6	91

11	Proximity-accelerated chemical coupling reaction in the benzodiazepine-binding site of gamma-aminobutyric acid type A receptors: superposition of different allosteric modulators. <i>Journal of Biological Chemistry</i> , <b>2007</b> , 282, 26316-25	5.4	28
10	Spontaneous cross-link of mutated alpha1 subunits during GABA(A) receptor assembly. <i>Journal of Biological Chemistry</i> , <b>2007</b> , 282, 4354-4363	5.4	9
9	Identification of amino acid residues important for assembly of GABA receptor alpha1 and gamma2 subunits. <i>Journal of Neurochemistry</i> , <b>2006</b> , 96, 983-95	6	15
8	Comparative models of GABAA receptor extracellular and transmembrane domains: important insights in pharmacology and function. <i>Molecular Pharmacology</i> , <b>2005</b> , 68, 1291-300	4.3	119
7	Affinity of various benzodiazepine site ligands in mice with a point mutation in the GABA(A) receptor gamma2 subunit. <i>Biochemical Pharmacology</i> , <b>2004</b> , 68, 1621-9	6	40
6	Raman and infrared vibrational spectra, ab initio calculations and normal coordinate analyses for 1,2-dimethyltetrachlorodisilane and 1,2-dimethyltetrachlorodisilane-d6. <i>Journal of Molecular Structure</i> , <b>1997</b> , 412, 83-95	3.4	12
5	Raman and infrared vibrational spectra, normal coordinate analysis and ab initio calculations of 1,1,2,2-tetrachlorodisilane and ab initio calculations of hexachlorodisilane. <i>Journal of Raman Spectroscopy</i> , <b>1997</b> , 28, 589-597	2.3	16
4	1,1,1,3,3,3-Hexabromotrisilane: structure and conformation determined by gas-phase electron diffraction, ab initio molecular orbital and molecular mechanics calculations, and vibrational spectroscopy. <i>Computational and Theoretical Chemistry</i> , <b>1995</b> , 372, 161-172		
3	GABAA Receptors1-12		
2	GABAA Receptor Modulators1-10		
1	Infrared and Raman Spectra, ab initio Calculations, and Rotational Isomerism of Methylated Disilanes241-247		