Mire Zloh

List of Publications by Year
in descending order

130 papers	4,289 citations	126907 33 h-index	123424 61 g-index
135 all docs	135 docs citations	135 times ranked	6500 citing authors

#	Article	IF	Citations
1	Molecular recognition of acetylcholinesterase and its subnanomolar reversible inhibitor: a molecular simulations study. Journal of Biomolecular Structure and Dynamics, 2022, 40, 1671-1691.	3.5	2
2	Potential of Helicrysum italicum cultivated in urban environment: SCCO2 extract cytotoxicity & NF-kB activation in HeLa, MCF-7 and MRC-5Âcells. Sustainable Chemistry and Pharmacy, 2022, 26, 100622.	3.3	1
3	VEGFA, B, C: Implications of the C-Terminal Sequence Variations for the Interaction with Neuropilins. Biomolecules, 2022, 12, 372.	4.0	2
4	Modelling Approaches for Studies of Drug-Polymer Interactions in Drug Delivery Systems. , 2022, , 561-591.		1
5	Flavonoids from Artemisia rupestris and their synergistic antibacterial effects on drug-resistant Staphylococcus aureus. Natural Product Research, 2021, 35, 1881-1886.	1.8	29
6	Deep Learning for Novel Antimicrobial Peptide Design. Biomolecules, 2021, 11, 471.	4.0	44
7	Novel C-3-(N-alkyl-aryl)-aminomethyl rifamycin SV derivatives exhibit activity against rifampicin-resistant Mycobacterium tuberculosis RpoBS522L strain and display a different binding mode at the RNAP Î²-subunit site compared to rifampicin. European Journal of Medicinal Chemistry, 2021, 225, 113734.	5.5	4
8	Intrinsic acetamide brush-off by polyurea biodendrimers. Journal of Materials Chemistry B, 2021, 9, 3371-3376.	5.8	2
9	Protein modification by bis-alkylation. , 2020, , 351-385.		2
10	Synthesis and in Silico Modelling of the Potential Dual Mechanistic Activity of Small Cationic Peptides Potentiating the Antibiotic Novobiocin against Susceptible and Multi-Drug Resistant Escherichia coli. International Journal of Molecular Sciences, 2020, 21, 9134.	4.1	8
11	Early detection of metabolic changes in drug-induced steatosis using metabolomics approaches. RSC Advances, 2020, 10, 41047-41057.	3.6	3
12	An update on the use of molecular modeling in dendrimers design for biomedical applications: are we using its full potential?. Expert Opinion on Drug Discovery, 2020, 15, 1015-1024.	5.0	5
13	Use of near infrared spectroscopy and spectral databases to assess the quality of pharmaceutical products and aid characterization of unknown components. Journal of Near Infrared Spectroscopy, 2019, 27, 379-390.	1.5	2
14	Functionalized branched polymers: promising immunomodulatory tools for the treatment of cancer and immune disorders. Materials Horizons, 2019, 6, 1956-1973.	12.2	44
15	Inhibitory Effect of Berberine on Broiler P-glycoprotein Expression and Function: In Situ and In Vitro Studies. International Journal of Molecular Sciences, 2019, 20, 1966.	4.1	19
16	NMR spectroscopy in drug discovery and development: Evaluation of physico-chemical properties. ADMET and DMPK, 2019, 7, 242-251.	2.1	6
17	Use of quercetin in animal feed: effects on the P-gp expression and pharmacokinetics of orally administrated enrofloxacin in chicken. Scientific Reports, 2018, 8, 4400.	3.3	28
18	The benefits of in silico modeling to identify possible small-molecule drugs and their off-target interactions. Future Medicinal Chemistry, 2018, 10, 423-432.	2.3	36

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19	Tautomerism of 4-phenyl-2,4-dioxobutanoic acid. Insights from pH ramping NMR study and quantum chemical calculations. Structural Chemistry, 2018, 29, 423-434.	2.0	2
20	Design, synthesis and biological evaluation of novel aryldiketo acids with enhanced antibacterial activity against multidrug resistant bacterial strains. European Journal of Medicinal Chemistry, 2018, 143, 1474-1488.	5.5	13
21	An Evaluation of the Potential of NMR Spectroscopy and Computational Modelling Methods to Inform Biopharmaceutical Formulations. Pharmaceutics, 2018, 10, 165.	4.5	22
22	Detection of newly emerging psychoactive substances using Raman spectroscopy and chemometrics. RSC Advances, 2018, 8, 31924-31933.	3.6	21
23	Relevance of Breast Cancer Resistance Protein to Pharmacokinetics of Florfenicol in Chickens: A Perspective from In Vivo and In Vitro Studies. International Journal of Molecular Sciences, 2018, 19, 3165.	4.1	8
24	Naphthalene Periâ€…Annelated N,Nâ€•and N,Oâ€Heterocycles: The Effect of Heteroatomâ€Guided Periâ€Fusion on Their Structure and Reactivity Profilesâ€A Theoretical Endoscopy. ChemistrySelect, 2018, 3, 9743-9752.	1.5	6
25	Î±-Galactosylceramide and peptide-based nano-vaccine synergistically induced a strong tumor suppressive effect in melanoma. Acta Biomaterialia, 2018, 76, 193-207.	8.3	27
26	Octodrine: New Questions and Challenges in Sport Supplements. Brain Sciences, 2018, 8, 34.	2.3	22
27	In Silico Structural Evaluation of Short Cationic Antimicrobial Peptides. Pharmaceutics, 2018, 10, 72.	4.5	10
28	Practical computational toolkits for dendrimers and dendrons structure design. Journal of Computer-Aided Molecular Design, 2017, 31, 817-827.	2.9	8
29	Intended and unintended use of cathinone mixtures. Human Psychopharmacology, 2017, 32, e2598.	1.5	46
30	4,4â€²â€Dimethylaminorex (â€œ4,4â€²â€DMARâ€, â€œSerotoniâ€) misuse: A Webâ€based study. Human Psychopharmacology, 2017, 32, e2575.	1.5	16
31	Poly-glutamic dendrimer-based conjugates for cancer vaccination â€“ a computational design for targeted delivery of antigens. Journal of Drug Targeting, 2017, 25, 873-880.	4.4	9
32	Drowning in diversity? A systematic way of clustering and selecting a representative set of new psychoactive substances. RSC Advances, 2017, 7, 53181-53191.	3.6	13
33	Rational design of novel, fluorescent, tagged glutamic acid dendrimers with different terminal groups and in silico analysis of their properties. International Journal of Nanomedicine, 2017, Volume 12, 7053-7073.	6.7	15
34	Redox properties of alkyl-substituted 4-aryl-2,4-dioxobutanoic acids. Journal of the Serbian Chemical Society, 2017, 82, 303-316.	0.8	4
35	Identification of Proteinâ€“Excipient Interaction Hotspots Using Computational Approaches. International Journal of Molecular Sciences, 2016, 17, 853.	4.1	18
36	Small molecule recognition of mephedrone using an anthracene molecular clip. Chemical Communications, 2016, 52, 7474-7477.	4.1	10

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37	Optimization of protein loaded PLGA nanoparticle manufacturing parameters following a quality-by-design approach. RSC Advances, 2016, 6, 104502-104512.	3.6	7
38	Application of diffusionâ€edited and solvent suppression ¹Hâ€NMR to the direct analysis of markers in valerianâ€hop liquid herbal products. Phytochemical Analysis, 2016, 27, 100-106.	2.4	2
39	Evidence that diclofenac and celecoxib are thyroid hormone receptor beta antagonists. Life Sciences, 2016, 146, 66-72.	4.3	17
40	In silico modelling of prostacyclin and other lipid mediators to nuclear receptors reveal novel thyroid hormone receptor antagonist properties. Prostaglandins and Other Lipid Mediators, 2016, 122, 18-27.	1.9	6
41	Survey of knowledge of legal highs (novel psychoactive substances) amongst London pharmacists. Drugs and Alcohol Today, 2015, 15, 93-99.	0.7	49
42	Antagonistic effects of indoloquinazoline alkaloids on antimycobacterial activity of evocarpine. Journal of Applied Microbiology, 2015, 118, 864-872.	3.1	26
43	Intramolecular cyclization of Î²-nitroso-o-quinone methides. A theoretical endoscopy of a potentially useful innate â€˜reclusiveâ€™ reaction. Tetrahedron, 2015, 71, 359-369.	1.9	17
44	Molecular Modeling to Study Dendrimers for Biomedical Applications. Molecules, 2014, 19, 20424-20467.	3.8	66
45	Computational classification models for predicting the interaction of drugs with P-glycoprotein and breast cancer resistance protein. SAR and QSAR in Environmental Research, 2014, 25, 939-966.	2.2	28
46	Production of Water-Soluble Few-Layer Graphene Mesosheets by Dry Milling with Hydrophobic Drug. Langmuir, 2014, 30, 14999-15008.	3.5	10
47	Arene-fused 1,2-oxazole N-oxides and derivatives. The impact of the Nâ€“O dipole and substitution on their aromatic character and reactivity profile. Can it be a useful structure in synthesis? A theoretical insight. Structural Chemistry, 2014, 25, 1837-1846.	2.0	4
48	Rapid detection of sildenafil analogue in Eurycoma longifolia products using a new two-tier procedure of the near infrared (NIR) spectra database. Food Chemistry, 2014, 158, 296-301.	8.2	25
49	A phytochemical comparison of saw palmetto products using gas chromatography and 1H nuclear magnetic resonance spectroscopy metabolomic profiling. Journal of Pharmacy and Pharmacology, 2014, 66, 811-822.	2.4	40
50	Insights into mechanism of anticancer activity of pentacyclic oxindole alkaloids of Uncaria tomentosa by means of a computational reverse virtual screening and molecular docking approach. Monatshefte FÃ¼r Chemie, 2014, 145, 1201-1211.	1.8	11
51	Structural insights into binding of small molecule inhibitors to Enhancer of Zeste Homolog 2. Journal of Computer-Aided Molecular Design, 2014, 28, 1109-1128.	2.9	3
52	Î²-Nitroso-o-quinone methides: potent intermediates in organic chemistry and biology. The impact of the NO group on their structure and reactivity profile: a theoretical insight. Structural Chemistry, 2014, 25, 1711-1723.	2.0	7
53	Bioadhesive tablets containing cyclodextrin complex of itraconazole for the treatment of vaginal candidiasis. International Journal of Biological Macromolecules, 2014, 69, 124-136.	7.5	50
54	Cationic Poly-<scp>l</scp>-lysine Dendrimer Complexes Doxorubicin and Delays Tumor Growth in Vitro and in Vivo. ACS Nano, 2013, 7, 1905-1917.	14.6	124

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55	Molecular Dynamic Simulations of Ocular Tablet Dissolution. Journal of Chemical Information and Modeling, 2013, 53, 3000-3008.	5.4	11
56	Study of the selectivity of Î±1-adrenergic antagonists by molecular modeling of Î±1a-, Î±1b-, and Î±1d-adrenergic receptor subtypes and docking simulations. Monatshefte FÃ¼r Chemie, 2013, 144, 903-912.	1.8	1
57	Investigation of the protein alkylation sites of the STAT3:STAT3 inhibitor Stattic by mass spectrometry. Bioorganic and Medicinal Chemistry Letters, 2013, 23, 4719-4722.	2.2	45
58	Potential of Lichen Secondary Metabolites against Plasmodium Liver Stage Parasites with FAS-II as the Potential Target. Journal of Natural Products, 2013, 76, 1064-1070.	3.0	30
59	Crystal Structure of a Promoter Sequence in the B-raf Gene Reveals an Intertwined Dimer Quadruplex. Journal of the American Chemical Society, 2013, 135, 19319-19329.	13.7	45
60	Quality Control of Natural Product Legal High Materials in the UK Using NMR Based Metabolomic Profiling. Planta Medica, 2013, 79, .	1.3	0
61	Natural chalcones as dual inhibitors of HDACs and NF-ÎºB. Oncology Reports, 2012, 28, 797-805.	2.6	71
62	C9orf72 hexanucleotide repeat associated with amyotrophic lateral sclerosis and frontotemporal dementia forms RNA G-quadruplexes. Scientific Reports, 2012, 2, 1016.	3.3	275
63	Prediction of aqueous solubility of drug-like molecules using a novel algorithm for automatic adjustment of relative importance of descriptors implemented in counter-propagation artificial neural networks. International Journal of Pharmaceutics, 2012, 437, 232-241.	5.2	18
64	Preventing acute gut wall damage in infectious diarrhoeas with glycosylated dendrimers. EMBO Molecular Medicine, 2012, 4, 866-881.	6.9	34
65	Site-Specific PEGylation at Histidine Tags. Bioconjugate Chemistry, 2012, 23, 248-263.	3.6	68
66	Target fishing and docking studies of the novel derivatives of aryl-aminopyridines with potential anticancer activity. Bioorganic and Medicinal Chemistry, 2012, 20, 5220-5228.	3.0	32
67	Sequences in the HSP90 promoter form G-quadruplex structures with selectivity for disubstituted phenyl bis-oxazole derivatives. Bioorganic and Medicinal Chemistry Letters, 2012, 22, 5930-5935.	2.2	26
68	A Prodrug Nanoparticle Approach for the Oral Delivery of a Hydrophilic Peptide, Leucine⁵-enkephalin, to the Brain. Molecular Pharmaceutics, 2012, 9, 1665-1680.	4.6	64
69	Antibacterial Acylphloroglucinols from Hypericum olympicum. Journal of Natural Products, 2012, 75, 336-343.	3.0	62
70	Targeting glycolysis: a fragment based approach towards bifunctional inhibitors of hLDH-5. Chemical Communications, 2011, 47, 230-232.	4.1	24
71	Computational design principles for bioactive dendrimer based constructs as antagonists of the TLR4-MD-2-LPS complex. Biomaterials, 2011, 32, 8702-8711.	11.4	22
72	Structural studies of biologically active glycosylated polyamidoamine (PAMAM) dendrimers. Journal of Molecular Modeling, 2011, 17, 2051-2060.	1.8	23

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73	From sequence to 3D structure of hyperbranched molecules: application to surface modified PAMAM dendrimers. Journal of Molecular Modeling, 2011, 17, 2741-2749.	1.8	13
74	In silico screening for antibiotic escort molecules to overcome efflux. Journal of Molecular Modeling, 2011, 17, 2863-2872.	1.8	6
75	Relative quantification of polyethylene glycol 400 excreted in the urine of male and female volunteers by direct injection electrospray-selected ion monitoring mass spectrometry. International Journal of Pharmaceutics, 2011, 414, 35-41.	5.2	6
76	Near-infrared spectroscopy (NIRS) and chemometric analysis of Malaysian and UK paracetamol tablets: A spectral database study. International Journal of Pharmaceutics, 2011, 415, 102-109.	5.2	24
77	Partially Glycosylated Dendrimers Block MD-2 and Prevent TLR4-MD-2-LPS Complex Mediated Cytokine Responses. PLoS Computational Biology, 2011, 7, e1002095.	3.2	31
78	Metabolomic profiling of saw palmetto products using proton-NMR spectroscopy and multi-variate analysis. Planta Medica, 2011, 77, .	1.3	0
79	2-Hexadecynoic acid inhibits plasmodial FAS-II enzymes and arrests erythrocytic and liver stage Plasmodium infections. Bioorganic and Medicinal Chemistry, 2010, 18, 7475-7485.	3.0	29
80	An analysis of the â€˜legal highâ€™ mephedrone. Bioorganic and Medicinal Chemistry Letters, 2010, 20, 4135-4139.	2.2	141
81	Purification, characterisation and identification of acidocin LCHV, an antimicrobial peptide produced by Lactobacillus acidophilus n.v. Er 317/402 strain Narine. International Journal of Antimicrobial Agents, 2010, 35, 255-260.	2.5	31
82	Structureâ^'Activity Relationships of Monomeric C2-Aryl Pyrrolo[2,1-c][1,4]benzodiazepine (PBD) Antitumor Agents. Journal of Medicinal Chemistry, 2010, 53, 2927-2941.	6.4	39
83	Disruption of d-alanyl esterification of Staphylococcus aureus cell wall teichoic acid by the Î²-lactam resistance modifier (â^')-epicatechin gallate. Journal of Antimicrobial Chemotherapy, 2009, 63, 1156-1162.	3.0	54
84	Induction of the Cytoprotective Enzyme Heme Oxygenase-1 by Statins Is Enhanced in Vascular Endothelium Exposed to Laminar Shear Stress and Impaired by Disturbed Flow. Journal of Biological Chemistry, 2009, 284, 18882-18892.	3.4	96
85	Aryldiketo Acids Have Antibacterial Activity Against MDR Staphylococcus aureus Strains: Structural Insights Based on Similarity and Molecular Interaction Fields. ChemMedChem, 2009, 4, 1971-1975.	3.2	13
86	Direct metabolic fingerprinting of commercial herbal tinctures by nuclear magnetic resonance spectroscopy and mass spectrometry. Phytochemical Analysis, 2009, 20, 328-334.	2.4	30
87	Preparation and Characterisation of Natamycin: Î³-Cyclodextrin Inclusion Complex and its Evaluation in Vaginal Mucoadhesive Formulations. Journal of Pharmaceutical Sciences, 2008, 97, 4319-4335.	3.3	47
88	Direct NMR analysis of cannabis water extracts and tinctures and semi-quantitative data on Î”9-THC and Î”9-THC-acid. Phytochemistry, 2008, 69, 562-570.	2.9	42
89	2Î²-Acetoxyferruginolâ€”A new antibacterial abietane diterpene from the bark of Prumnopitys andina. Phytochemistry Letters, 2008, 1, 49-53.	1.2	18
90	New metabolites with antibacterial activity from the marine angiosperm Cymodocea nodosa. Tetrahedron, 2008, 64, 1696-1702.	1.9	55

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91	Role of complexes formation between drugs and penetration enhancers in transdermal delivery. International Journal of Pharmaceutics, 2008, 363, 40-49.	5.2	29
92	Disulfide bridge based PEGylation of proteins. Advanced Drug Delivery Reviews, 2008, 60, 3-12.	13.7	170
93	Solution Structure of a 2:1 C2-(2-Naphthyl) Pyrrolo[2,1-c][1,4]benzodiazepine DNA Adduct: Molecular Basis for Unexpectedly High DNA Helix Stabilization. Biochemistry, 2008, 47, 11818-11829.	2.5	24
94	The Effect of Phenyl Substituents on 13C NMR Shifts and Metal Ions Binding to 4-Phenyl-2,4-Dioxobutanoic Acid Derivatives. Letters in Organic Chemistry, 2008, 5, 692-699.	0.5	3
95	An LFER study of the protolytic equilibria of 4-aryl-2,4-dioxobutanoic acids in aqueous solutions. Journal of the Serbian Chemical Society, 2007, 72, 1201-1216.	0.8	16
96	Identification and insertion of 3-carbon bridges in protein disulfide bonds: a computational approach. Nature Protocols, 2007, 2, 1070-1083.	12.0	25
97	Site-Specific PEGylation of Protein Disulfide Bonds Using a Three-Carbon Bridge. Bioconjugate Chemistry, 2007, 18, 61-76.	3.6	152
98	Molecular Dynamics Simulations of Proteins with Chemically Modified Disulfide Bonds. Theoretical Chemistry Accounts, 2007, 117, 259-265.	1.4	12
99	Quantum Chemical Studies on Structure Activity Relationship of Natural Product Polyacetylenes. Theoretical Chemistry Accounts, 2007, 117, 247-252.	1.4	9
100	The Role of Small Moleculeâ€“small Molecule Interactions in Overcoming Biological Barriers for Antibacterial Drug Action. Theoretical Chemistry Accounts, 2007, 117, 231-238.	1.4	11
101	Synthesis of DNA-Directed Pyrrolidinyl and Piperidinyl Confined Alkylating Chloroalkylaminoanthraquinones:â€‰ Potential for Development of Tumor-Selective N-Oxides. Journal of Medicinal Chemistry, 2006, 49, 7013-7023.	6.4	43
102	Screening far red probes for use on optical biochip devices. , 2006, 6088, 122.		0
103	Site-specific PEGylation of native disulfide bonds in therapeutic proteins. Nature Chemical Biology, 2006, 2, 312-313.	8.0	246
104	PEGylation of native disulfide bonds in proteins. Nature Protocols, 2006, 1, 2241-2252.	12.0	110
105	Spectral analysis of the DNA targeting bisalkylaminoanthraquinone DRAQ5 in intact living cells. Cytometry Part A: the Journal of the International Society for Analytical Cytology, 2006, 69A, 805-814.	1.5	36
106	Amanicadol, a Pimarane-type Diterpene from Phlomis amanica Vierch.. Zeitschrift Fur Naturforschung - Section B Journal of Chemical Sciences, 2006, 61, 1433-1436.	0.7	8
107	Advanced microscopy solutions for monitoring the kinetics and dynamics of drug?DNA targeting in living cells. Advanced Drug Delivery Reviews, 2005, 57, 153-167.	13.7	47
108	Investigation of the association and flexibility of cationic lipidic peptide dendrons by NMR spectroscopy. Magnetic Resonance in Chemistry, 2005, 43, 47-52.	1.9	10

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109	Isopimaric acid fromPinus nigra shows activity against multidrug-resistant and EMRSA strains ofStaphylococcus aureus. Phytotherapy Research, 2005, 19, 538-542.	5.8	100
110	Putative DNA Quadruplex Formation within the Humanc-kitOncogene. Journal of the American Chemical Society, 2005, 127, 10584-10589.	13.7	526
111	Efficient Solid-Phase-Based Total Synthesis of the Bisintercalator TANDEM. Journal of Organic Chemistry, 2005, 70, 7654-7661.	3.2	20
112	Molecular Similarity of MDR Inhibitors. International Journal of Molecular Sciences, 2004, 5, 37-47.	4.1	12
113	Inhibitors of multidrug resistance (MDR) have affinity for MDR substrates. Bioorganic and Medicinal Chemistry Letters, 2004, 14, 881-885.	2.2	41
114	Acontyl-derived polymers for biomedical applications. Modeling study of cis-trans isomerisation. Theoretical Chemistry Accounts, 2003, 109, 206-212.	1.4	11
115	Solid-Phase Synthesis of the Cyclic Peptide Portion of Chlorofusin, an Inhibitor of p53-MDM2 Interactions. Organic Letters, 2003, 5, 5051-5054.	4.6	38
116	Competitive Reactions During Amine Addition to cis-Aconityl Anhydride. Australian Journal of Chemistry, 2002, 55, 467.	0.9	19
117	Docking of Transmembrane Helices Into Four Helix Bundles in the High Affinity IgE Receptor. , 2001, , 841-842.		0
118	Spectroscopy-Based Modelling of the 3D Structure of the Î² Subunit of the High Affinity IgE Receptor. Molecular Simulation, 2000, 24, 421-447.	2.0	3
119	NMR lipid profile of Agaricus bisporus. Phytochemistry, 1999, 50, 1311-1321.	2.9	24
120	44 NMR studies of the extracellular loop of the beta subunit of the high affinity IgE receptor. Biochemical Society Transactions, 1998, 26, S34-S34.	3.4	1
121	45 Molecular modelling of the IgE receptor loops - lipid interaction. Biochemical Society Transactions, 1998, 26, S35-S35.	3.4	1
122	Lipid mimetics: the design and properties of conformationally-restricted arachidonic acid lipidic and peptidic analogues. Biochemical Society Transactions, 1997, 25, 26S-26S.	3.4	5
123	NMR-Based Modelling Revealed an Alpha Helical Structure for Cytoplasmic Domain of the Alpha Subunit of FcÎµRI, the High Affinity IgE Receptor. Biochemical Society Transactions, 1997, 25, 55S-55S.	3.4	1
124	LIPID â€“ HELIX INTERACTIONS IN MEMBRANE RECEPTORS. Biochemical Society Transactions, 1996, 24, 305S-305S.	3.4	2
125	NMR LIPIDS PROFILES OF COMMON MUSHROOMS. Biochemical Society Transactions, 1995, 23, 613S-613S.	3.4	0
126	Conformational studies of the beta-subunit of the high affinity IgE receptor: circular dichroism and molecular modelling. Biomedical Peptides, Proteins & Nucleic Acids: Structure, Synthesis & Biological Activity, 1995, 1, 101-6.	0.1	0

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127	Mechanism of excitation and emission of papaverine molecule: fluorescence polarization spectroscopy study and MO calculation of transition moments. Journal of Luminescence, 1994, 59, 27-32.	3.1	1
128	N.m.r. studies of the cytoplasmic C-terminal Î²-subunit domain of the high-affinity IgE receptor. Biochemical Society Transactions, 1994, 22, 1027-1029.	3.4	4
129	Spectroscopic and conformational studies of the C-terminal cytoplasmic beta subunit 46-peptide of the high affinity IgE receptor. Biochemical Society Transactions, 1994, 22, 450S-450S.	3.4	2
130	Hydrophobic Core and Surface Charges of Human Î²2-Microglobulin Probed by CD Measurements. Collection of Czechoslovak Chemical Communications, 1992, 57, 1143-1148.	1.0	0