## Congbao Kang

List of Publications by Year in descending order

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120 3,974 33 58 g-index

122 122 122 4473

times ranked

citing authors

docs citations

all docs

#	Article	IF	CITATIONS
1	Secondary structures, dynamics, and DNA binding of the homeodomain of human SIX1. Journal of Peptide Science, 2022, 28, e3376.	1.4	2
2	Structureâ€"activity relationship studies of allosteric inhibitors of <scp>EYA2</scp> tyrosine phosphatase. Protein Science, 2022, 31, 422-431.	7.6	4
3	Secondary Structures of the Transmembrane Domain of SARS-CoV-2 Spike Protein in Detergent Micelles. International Journal of Molecular Sciences, 2022, 23, 1040.	4.1	4
4	$1\mathrm{H},13\mathrm{C}$ and $15\mathrm{N}$ resonance assignments of the first BIR domain of cellular inhibitor of apoptosis protein $1.$ Biomolecular NMR Assignments, 2022, , $1.$	0.8	1
5	Structures and Dynamics of Dengue Virus Nonstructural Membrane Proteins. Membranes, 2022, 12, 231.	3.0	9
6	Structure and Dynamics of Zika Virus Protease and Its Insights into Inhibitor Design. Biomedicines, 2021, 9, 1044.	3.2	10
7	Perspectives on Fragment-based Drug Discovery: A Strategy Applicable to Diverse Targets. Current Topics in Medicinal Chemistry, 2021, 21, 1099-1112.	2.1	15
8	Targeting EYA2 tyrosine phosphatase activity in glioblastoma stem cells induces mitotic catastrophe. Journal of Experimental Medicine, 2021, 218, .	8.5	9
9	$1 {\rm \^AH}, 15 {\rm \^AN}$ and $13 {\rm \^AC}$ resonance assignments of the Q61H mutant of human KRAS bound to GDP. Biomolecular NMR Assignments, 2021, , .	0.8	1
10	Mechanisms of Action for Small Molecules Revealed by Structural Biology in Drug Discovery. International Journal of Molecular Sciences, 2020, 21, 5262.	4.1	34
11	Discovery of Covalent Inhibitors Targeting the Transcriptional Enhanced Associate Domain Central Pocket. Journal of Medicinal Chemistry, 2020, 63, 11972-11989.	6.4	36
12	Progress in Developing Inhibitors of SARS-CoV-2 3C-Like Protease. Microorganisms, 2020, 8, 1250.	3.6	90
13	A Practical Perspective on the Roles of Solution NMR Spectroscopy in Drug Discovery. Molecules, 2020, 25, 2974.	3.8	11
14	Breakthroughs in Medicinal Chemistry: New Targets and Mechanisms, New Drugs, New Hopes–7. Molecules, 2020, 25, 2968.	3.8	5
15	Identification and structural characterization of small molecule fragments targeting Zika virus NS2B-NS3 protease. Antiviral Research, 2020, 175, 104707.	4.1	15
16	Probing biological mechanisms with chemical tools. Pharmacological Research, 2020, 153, 104656.	7.1	4
17	Insights into Structures and Dynamics of Flavivirus Proteases from NMR Studies. International Journal of Molecular Sciences, 2020, 21, 2527.	4.1	11
18	PGE1 and PGA1 bind to Nurr1 and activate its transcriptional function. Nature Chemical Biology, 2020, 16, 876-886.	8.0	51

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19	Structural and Functional Analyses of an Allosteric EYA2 Phosphatase Inhibitor That Has On-Target Effects in Human Lung Cancer Cells. Molecular Cancer Therapeutics, 2019, 18, 1484-1496.	4.1	34
20	Dengue NS2A Protein Orchestrates Virus Assembly. Cell Host and Microbe, 2019, 26, 606-622.e8.	11.0	68
21	Targeting the Bacterial Epitranscriptome for Antibiotic Development: Discovery of Novel tRNA-(N <sup>1</sup> G37) Methyltransferase (TrmD) Inhibitors. ACS Infectious Diseases, 2019, 5, 326-335.	3.8	33
22	Backbone resonance assignment for the full length tRNA-(N1G37) methyltransferase of Pseudomonas aeruginosa. Biomolecular NMR Assignments, 2019, 13, 327-332.	0.8	0
23	Secondary structure and topology of the transmembrane domain of Syndecanâ€⊋ in detergent micelles. FEBS Letters, 2019, 593, 554-561.	2.8	6
24	Applications of In-Cell NMR in Structural Biology and Drug Discovery. International Journal of Molecular Sciences, 2019, 20, 139.	4.1	33
25	Backbone resonance assignment for the N-terminal region of bacterial tRNA-(N1G37) methyltransferase. Biomolecular NMR Assignments, 2019, 13, 49-53.	0.8	4
26	Expression, purification of Zika virus membrane protein-NS2B in detergent micelles for NMR studies. Protein Expression and Purification, 2019, 154, 1-6.	1.3	13
27	<sup>19</sup> F-NMR in Target-based Drug Discovery. Current Medicinal Chemistry, 2019, 26, 4964-4983.	2.4	22
28	Structural Insights into the Inhibition of Zika Virus NS2B-NS3 Protease by a Small-Molecule Inhibitor. Structure, 2018, 26, 555-564.e3.	3.3	70
29	Secondary structure and membrane topology of dengue virus NS4A protein in micelles. Biochimica Et Biophysica Acta - Biomembranes, 2018, 1860, 442-450.	2.6	21
30	Elucidating the bactericidal mechanism of action of the linear antimicrobial tetrapeptide BRBR-NH2. Biochimica Et Biophysica Acta - Biomembranes, 2018, 1860, 1517-1527.	2.6	12
31	The Dengue Virus Replication Complex: From RNA Replication to Protein-Protein Interactions to Evasion of Innate Immunity. Advances in Experimental Medicine and Biology, 2018, 1062, 115-129.	1.6	45
32	Structural and ligand-binding analysis of the YAP-binding domain of transcription factor TEAD4. Biochemical Journal, 2018, 475, 2043-2055.	3.7	35
33	Characterization of molecular interactions between Zika virus protease and peptides derived from the C-terminus of NS2B. Biochemical and Biophysical Research Communications, 2018, 503, 691-696.	2.1	15
34	Structural characterization of the linked <scp>NS</scp> 2Bâ€ <scp>NS</scp> 3 protease of Zika virus. FEBS Letters, 2017, 591, 2338-2347.	2.8	35
35	Backbone resonance assignments for the SET domain of human methyltransferase NSD3 in complex with its cofactor. Biomolecular NMR Assignments, 2017, 11, 225-229.	0.8	2
36	Zika Virus Protease: An Antiviral Drug Target. Trends in Microbiology, 2017, 25, 797-808.	7.7	80

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37	Structural Dynamics of Zika Virus NS2B-NS3 Protease Binding to Dipeptide Inhibitors. Structure, 2017, 25, 1242-1250.e3.	3.3	83
38	Solution NMR Spectroscopy in Target-Based Drug Discovery. Molecules, 2017, 22, 1399.	3.8	32
39	Erythropoietin Receptor Structural Domains. Vitamins and Hormones, 2017, 105, 1-17.	1.7	2
40	Structure of the transmembrane domain of human nicastrin-a component of $\hat{l}^3$ -secretase. Scientific Reports, 2016, 6, 19522.	3.3	11
41	Crystal structure of unlinked NS2B-NS3 protease from Zika virus. Science, 2016, 354, 1597-1600.	12.6	156
42	Structural basis for KCNE3 modulation of potassium recycling in epithelia. Science Advances, 2016, 2, e1501228.	10.3	45
43	Secondary Structure and Membrane Topology of the Full-Length Dengue Virus NS4B in Micelles. Angewandte Chemie, 2016, 128, 12247-12251.	2.0	5
44	Characterization of the Molecular Interactions between ParE/Gyrb and an Inhibitor and its Insight into Developing Antibacterial Agents. Biophysical Journal, 2016, 110, 542a.	0.5	0
45	Escherichia coli Topoisomerase IV E Subunit and an Inhibitor Binding Mode Revealed by NMR Spectroscopy. Journal of Biological Chemistry, 2016, 291, 17743-17753.	3.4	15
46	Secondary Structure and Membrane Topology of the Fullâ€Length Dengue Virus NS4B in Micelles. Angewandte Chemie - International Edition, 2016, 55, 12068-12072.	13.8	28
47	Structure of the NS2B-NS3 protease from Zika virus after self-cleavage. Nature Communications, 2016, 7, 13410.	12.8	169
48	Structure of the Cyclic Nucleotide-Binding Homology Domain of the hERG Channel and Its Insight into Type 2 Long QT Syndrome. Scientific Reports, 2016, 6, 23712.	3.3	9
49	Selection of suitable detergents for obtaining an active dengue protease in its natural form from E.Âcoli. Protein Expression and Purification, 2016, 121, 141-148.	1.3	6
50	Backbone assignment of the N-terminal 24-kDa fragment of Escherichia coli topoisomerase IV ParE subunit. Biomolecular NMR Assignments, 2016, 10, 135-138.	0.8	9
51	A Phenotypic Screen for Small-Molecule Inhibitors of Constitutively Active Mutant Thrombopoietin Receptor Implicated in Myeloproliferative Neoplasms. Combinatorial Chemistry and High Throughput Screening, 2016, 19, 824-833.	1.1	1
52	Structural transition in Bcl-xL and its potential association with mitochondrial calcium ion transport. Scientific Reports, 2015, 5, 10609.	3.3	17
53	Solution structure of the transmembrane domain of the mouse erythropoietin receptor in detergent micelles. Scientific Reports, 2015, 5, 13586.	3.3	21
54	NMR Structural Characterization for Proteases of Dengue and West Nile Viruses and its Insight into Drug Discovery. Biophysical Journal, 2015, 108, 375a.	0.5	0

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55	Application of Fragmentâ∈Based Drug Discovery against DNA Gyraseâ€B. ChemPlusChem, 2015, 80, 1250-12.	54.2.8	14
56	Identification of covalent active site inhibitors of dengue virus protease. Drug Design, Development and Therapy, 2015, 9, 6389.	4.3	25
57	Discovery of Dengue Virus NS4B Inhibitors. Journal of Virology, 2015, 89, 8233-8244.	3.4	77
58	Characterization of Dengue Virus NS4A and NS4B Protein Interaction. Journal of Virology, 2015, 89, 3455-3470.	3.4	116
59	Mapping the Interactions between the NS4B and NS3 Proteins of Dengue Virus. Journal of Virology, 2015, 89, 3471-3483.	3.4	83
60	Membrane topology of NS2B of dengue virus revealed by NMR spectroscopy. Biochimica Et Biophysica Acta - Biomembranes, 2015, 1848, 2244-2252.	2.6	63
61	Determinants of Dengue Virus NS4A Protein Oligomerization. Journal of Virology, 2015, 89, 6171-6183.	3.4	48
62	Secondary structure and membrane topology of dengue virus NS4B N-terminal 125 amino acids. Biochimica Et Biophysica Acta - Biomembranes, 2015, 1848, 3150-3157.	2.6	34
63	Targeting the Central Pocket in Human Transcription Factor TEAD as a Potential Cancer Therapeutic Strategy. Structure, 2015, 23, 2076-2086.	3.3	146
64	Characterization of the interaction between Escherichia coli topoisomerase IV E subunit and an ATP competitive inhibitor. Biochemical and Biophysical Research Communications, 2015, 467, 961-966.	2.1	7
65	Biophysical Studies of Bacterial Topoisomerases Substantiate Their Binding Modes to an Inhibitor. Biophysical Journal, 2015, 109, 1969-1977.	0.5	6
66	NMR structural characterization of the Nâ€terminal active domain of the gyrase B subunit from <i>Pseudomonas aeruginosa</i> and its complex with an inhibitor. FEBS Letters, 2015, 589, 2683-2689.	2.8	12
67	Structural analysis of the S4–S5 linker of the human KCNQ1 potassium channel. Biochemical and Biophysical Research Communications, 2015, 456, 410-414.	2.1	3
68	1H, 13C and 15N chemical shift assignments for the cyclic-nucleotide binding homology domain of a KCNH channel. Biomolecular NMR Assignments, 2015, 9, 55-58.	0.8	3
69	Structural insight into the transmembrane segments 3 and 4 of the hERG potassium channel. Journal of Peptide Science, 2014, 20, 935-944.	1.4	3
70	1H, 13C and 15N chemical shift assignments for the N-terminal PAS domain of the KCNH channel from Zebrafish. Biomolecular NMR Assignments, 2014, 8, 165-168.	0.8	4
71	Solution structure of the transmembrane domain of the insulin receptor in detergent micelles. Biochimica Et Biophysica Acta - Biomembranes, 2014, 1838, 1313-1321.	2.6	37
72	Dimerization of Flavivirus NS4B Protein. Journal of Virology, 2014, 88, 3379-3391.	3.4	77

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73	Structural Insight into the Transmembrane Domain and the Juxtamembrane Region of the Erythropoietin Receptor in Micelles. Biophysical Journal, 2014, 107, 2325-2336.	0.5	26
74	Insight into the molecular interaction between the cyclic nucleotideâ€binding homology domain and the eag domain of the hERG channel. FEBS Letters, 2014, 588, 2782-2788.	2.8	9
75	Solution structure of the cyclic-nucleotide binding homology domain of a KCNH channel. Journal of Structural Biology, 2014, 186, 68-74.	2.8	10
76	NMR Structural Study of the Domains of the KCNH Channels and its Insight into Channel Gating. Biophysical Journal, 2014, 106, 437a.	0.5	0
77	Drug Design For Flavivirus Proteases: What Are We Missing?. Current Pharmaceutical Design, 2014, 20, 3422-3427.	1.9	30
78	1H, 13C and 15N chemical shift assignments for an intracellular proteinase inhibitor of Bacillus subtilis. Biomolecular NMR Assignments, 2013, 7, 129-132.	0.8	7
79	Exploring the binding of peptidic West Nile virus NS2B–NS3 protease inhibitors by NMR. Antiviral Research, 2013, 97, 137-144.	4.1	33
80	West Nile virus protease activity in detergent solutions and application for affinity tag removal. Analytical Biochemistry, 2013, 435, 44-46.	2.4	10
81	Biomimetic membrane platform containing hERG potassium channel and its application to drug screening. Analyst, The, 2013, 138, 2007.	3.5	27
82	Membrane Topology and Function of Dengue Virus NS2A Protein. Journal of Virology, 2013, 87, 4609-4622.	3.4	162
83	Lyso-myristoyl phosphatidylcholine micelles sustain the activity of Dengue non-structural (NS) protein 3 protease domain fused with the full-length NS2B. Protein Expression and Purification, 2013, 92, 156-162.	1.3	17
84	NMR Analysis of a Novel Enzymatically Active Unlinked Dengue NS2B-NS3 Protease Complex. Journal of Biological Chemistry, 2013, 288, 12891-12900.	3.4	93
85	Functional role of the flexible N-terminal extension of FKBP38 in catalysis. Scientific Reports, 2013, 3, 2985.	3.3	10
86	Structure of the C-terminal Region of the Frizzled Receptor 1 in Detergent Micelles. Molecules, 2013, 18, 8579-8590.	3.8	12
87	Solution NMR Study of the Transmembrane Domain of Single-Span Membrane Proteins: Opportunities and Strategies. Current Protein and Peptide Science, 2012, 13, 585-600.	1.4	19
88	Purification and structural characterization of the voltage-sensor domain of the hERG potassium channel. Protein Expression and Purification, 2012, 86, 98-104.	1.3	9
89	Probing the Structural and Dynamic Properties of KCNE1 using Site-Directed Spin Labeling EPR Spectroscopy. Biophysical Journal, 2012, 102, 406a.	0.5	0
90	West Nile Virus (WNV) protease and membrane interactions revealed by NMR spectroscopy. Biochemical and Biophysical Research Communications, 2012, 423, 799-804.	2.1	7

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91	Expression and purification of the p75 neurotrophin receptor transmembrane domain using a ketosteroid isomerase tag. Microbial Cell Factories, 2012, 11, 45.	4.0	8
92	EPR Spectroscopic Distance Measurements of the KCNE1 Membrane Protein in Micelles and Lipid Bilayers. Biophysical Journal, 2012, 102, 406a.	0.5	0
93	The Natively Disordered Loop of Bcl-2 Undergoes Phosphorylation-Dependent Conformational Change and Interacts with Pin1. PLoS ONE, 2012, 7, e52047.	2.5	9
94	The solution structure of the S4–S5 linker of the hERG potassium channel. Journal of Peptide Science, 2012, 18, 140-145.	1.4	17
95	Reconstitution of KCNE1 into Lipid Bilayers: Comparing the Structural, Dynamic, and Activity Differences in Micelle and Vesicle Environments. Biochemistry, 2011, 50, 10851-10859.	2.5	31
96	Structural Studies on the Conformation of Human KCNEL1 Membrane Protein via Electron Paramagnetic Resonance Spectroscopy. Biophysical Journal, 2011, 100, 144a.	0.5	0
97	The Structural, Dynamic, and Functional Changes in the KCNE1 Membrane Protein Between Detergent Micelles and Lipid Bilayers. Biophysical Journal, 2011, 100, 144a.	0.5	0
98	NMR structural study of the intracellular loop 3 of the serotonin 5-HT1A receptor and its interaction with calmodulin. Biochimica Et Biophysica Acta - Biomembranes, 2011, 1808, 2224-2232.	2.6	22
99	Solution structure of a human minimembrane protein Ost4, a subunit of the oligosaccharyltransferase complex. Biochemical and Biophysical Research Communications, 2011, 409, 572-576.	2.1	22
100	Expression, purification, and initial structural characterization of nonstructural protein 2B, an integral membrane protein of Dengue-2 virus, in detergent micelles. Protein Expression and Purification, 2011, 80, 169-175.	1.3	30
101	An NMR study of the Nâ€terminal domain of wildâ€type hERG and a T65P trafficking deficient hERG mutant. Proteins: Structure, Function and Bioinformatics, 2011, 79, 2557-2565.	2.6	26
102	Solution NMR study of integral membrane proteins. Current Opinion in Chemical Biology, 2011, 15, 560-569.	6.1	71
103	NMR Solution Structure of Human Vaccinia-related Kinase 1 (VRK1) Reveals the C-terminal Tail Essential for Its Structural Stability and Autocatalytic Activity. Journal of Biological Chemistry, 2011, 286, 22131-22138.	3.4	40
104	1H, 13C and 15N chemical shift assignments for the N-terminal domain of the voltage-gated potassium channel-hERG. Biomolecular NMR Assignments, 2010, 4, 211-213.	0.8	16
105	Mechanistic basis for LQT1 caused by S3 mutations in the KCNQ1 subunit of <i>IKs</i> . Journal of General Physiology, 2010, 135, 433-448.	1.9	26
106	Functional Delivery of a Membrane Protein into Oocyte Membranes Using Bicelles. Biochemistry, 2010, 49, 653-655.	2.5	36
107	The impact of window functions on NMR-based paramagnetic relaxation enhancement measurements in membrane proteins. Biochimica Et Biophysica Acta - Biomembranes, 2010, 1798, 140-149.	2.6	14
108	NMR solution structure of the N-terminal domain of hERG and its interaction with the S4–S5 linker. Biochemical and Biophysical Research Communications, 2010, 403, 126-132.	2.1	62

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109	Mechanistic insights into non-immunosuppressive immunophilin ligands as potential antimalarial therapeutics. Malaria Journal, $2010, 9, .$	2.3	0
110	Structure of KCNE1 and Implications for How It Modulates the KCNQ1 Potassium Channel. Biochemistry, 2008, 47, 7999-8006.	2.5	183
111	FKBP Family Proteins: Immunophilins with Versatile Biological Functions. NeuroSignals, 2008, 16, 318-325.	0.9	278
112	Expression, purification, and molecular characterization of Plasmodium falciparum FK506-binding protein 35 (PfFKBP35). Protein Expression and Purification, 2007, 53, 179-185.	1.3	19
113	Preparation, Functional Characterization, and NMR Studies of Human KCNE1, a Voltage-Gated Potassium Channel Accessory Subunit Associated with Deafness and Long QT Syndrome <sup>,</sup> . Biochemistry, 2007, 46, 11459-11472.	2.5	61
114	1H, 13C, and 15N resonance assignments of FK506-binding domain of Plasmodium falciparum FKBP35. Biomolecular NMR Assignments, 2007, 1, 27-28.	0.8	1
115	The N-terminal domain of tumor suppressor p53 is involved in the molecular interaction with the anti-apoptotic protein Bcl-Xl. Biochemical and Biophysical Research Communications, 2006, 341, 938-944.	2.1	23
116	Molecular Characterization of the Recombinant A-chain of a Type II Ribosome-Inactivating Protein (RIP) from Viscum album coloratum and Structural Basis on its Ribosome-Inactivating Activity and the Sugar-binding Properties of the B-chain. BMB Reports, 2006, 39, 560-570.	2.4	18
117	Waste water produced from an oilfield and continuous treatment with an oil-degrading bacterium. Process Biochemistry, 2005, 40, 873-877.	3.7	96
118	Molecular characterization of FK-506 binding protein 38 and its potential regulatory role on the anti-apoptotic protein Bcl-2. Biochemical and Biophysical Research Communications, 2005, 337, 30-38.	2.1	48
119	The flexible loop of Bcl-2 is required for molecular interaction with immunosuppressant FK-506 binding protein 38 (FKBP38). FEBS Letters, 2005, 579, 1469-1476.	2.8	46
120	Application of microbial enhanced oil recovery technique to Daqing Oilfield. Biochemical Engineering Journal, 2002, 11, 197-199.	3.6	79