

Carol Beth Post

List of Publications by Year in descending order

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35
papers

1,121
citations

394421

19
h-index

395702

33
g-index

36
all docs

36
docs citations

36
times ranked

1611
citing authors

#	ARTICLE	IF	CITATIONS
1	Exchange-transferred NOE spectroscopy and bound ligand structure determination. <i>Current Opinion in Structural Biology</i> , 2003, 13, 581-588.	5.7	177
2	Shake, rattle, and roll: Impact of the dynamics of flavivirus particles on their interactions with the host. <i>Virology</i> , 2015, 479-480, 508-517.	2.4	103
3	Effects of impaired membrane interactions on α -synuclein aggregation and neurotoxicity. <i>Neurobiology of Disease</i> , 2015, 79, 150-163.	4.4	73
4	Src kinase activation: A switched electrostatic network. <i>Protein Science</i> , 2006, 15, 1051-1062.	7.6	71
5	Insights into Protein Compressibility from Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2001, 105, 715-724.	2.6	57
6	An electrostatic network and long-range regulation of Src kinases. <i>Protein Science</i> , 2008, 17, 1871-1880.	7.6	54
7	Tyr130 phosphorylation triggers Syk release from antigen receptor by long-distance conformational uncoupling. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 11760-11765.	7.1	53
8	α -Helix as a Switch in the Conformational Transition of Src/CDK-like Kinase Domains. <i>Journal of Physical Chemistry B</i> , 2012, 116, 4465-4475.	2.6	49
9	Theoretical studies of viral capsid proteins. <i>Current Opinion in Structural Biology</i> , 2000, 10, 170-173.	5.7	42
10	Constraining Binding Hot Spots: NMR and Molecular Dynamics Simulations Provide a Structural Explanation for Enthalpy-Entropy Compensation in SH2-Ligand Binding. <i>Journal of the American Chemical Society</i> , 2010, 132, 11058-11070.	13.7	42
11	Catalysis by Entropic Guidance from Enzymes. <i>Biochemistry</i> , 1996, 35, 15129-15133.	2.5	41
12	Long-distance correlations of rhinovirus capsid dynamics contribute to uncoating and antiviral activity. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2012, 109, 5271-5276.	7.1	35
13	Ubiquitin C-terminal Hydrolase L1: Biochemical and Cellular Characterization of a Covalent Cyanopyrrolidine-Based Inhibitor. <i>ChemBioChem</i> , 2020, 21, 712-722.	2.6	32
14	Unrestrained Computation of Free Energy along a Path. <i>Journal of Physical Chemistry B</i> , 2012, 116, 11046-11055.	2.6	28
15	Two Closely Spaced Tyrosines Regulate NFAT Signaling in B Cells via Syk Association with Vav. <i>Molecular and Cellular Biology</i> , 2011, 31, 2984-2996.	2.3	25
16	Docking multiple conformations of a flexible ligand into a protein binding site using NMR restraints. <i>Proteins: Structure, Function and Bioinformatics</i> , 2002, 46, 295-307.	2.6	24
17	Substrate Recognition by the Lyn Protein-tyrosine Kinase. <i>Journal of Biological Chemistry</i> , 2000, 275, 16174-16182.	3.4	23
18	Accuracy of bound peptide structures determined by exchange transferred nuclear Overhauser data: a simulation study. <i>Journal of Biomolecular NMR</i> , 2000, 17, 17-32.	2.8	20

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19	Molecular dynamics investigation of the effect of an antiviral compound on human rhinovirus. <i>Protein Science</i> , 1999, 8, 2281-2289.	7.6	20
20	A simple method for NMR t1 noise suppression. <i>Journal of Magnetic Resonance</i> , 2017, 276, 43-50.	2.1	18
21	Detection of Long-Range Concerted Motions in Protein by a Distance Covariance. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 3009-3014.	5.3	16
22	Relative Binding Enthalpies from Molecular Dynamics Simulations Using a Direct Method. <i>Journal of Chemical Theory and Computation</i> , 2014, 10, 2759-2768.	5.3	16
23	Insights into the allosteric regulation of Syk association with receptor ITAM, a multi-state equilibrium. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 5807-5818.	2.8	14
24	Application of a Substrate-Mediated Selection with c-Src Tyrosine Kinase to a DNA-Encoded Chemical Library. <i>Molecules</i> , 2019, 24, 2764.	3.8	14
25	Differential recognition of sykâ€binding sites by each of the two phosphotyrosineâ€binding pockets of the Vav SH2 domain. <i>Biopolymers</i> , 2013, 99, 897-907.	2.4	11
26	Entropic allostery dominates the phosphorylationâ€dependent regulation of Syk tyrosine kinase release from immunoreceptor tyrosineâ€based activation motifs. <i>Protein Science</i> , 2018, 27, 1780-1796.	7.6	11
27	Substrate binding to Src: A new perspective on tyrosine kinase substrate recognition from NMR and molecular dynamics. <i>Protein Science</i> , 2020, 29, 350-359.	7.6	9
28	Evaluating the dynamics and electrostatic interactions of folded proteins in implicit solvents. <i>Protein Science</i> , 2016, 25, 204-218.	7.6	7
29	NmrLineGuru: Standalone and User-Friendly GUIs for Fast 1D NMR Lineshape Simulation and Analysis of Multi-State Equilibrium Binding Models. <i>Scientific Reports</i> , 2019, 9, 16023.	3.3	7
30	All-atom adaptively biased path optimization of Src kinase conformational inactivation: Switched electrostatic network in the concerted motion of I±C helix and the activation loop. <i>Journal of Chemical Physics</i> , 2020, 153, 175101.	3.0	7
31	NMR structure of phospho-tyrosine signaling complexes. , 1999, 19, 295-305.		6
32	Intermolecular relaxation has little effect on intra-peptide exchange-transferred NOE intensities. <i>Journal of Biomolecular NMR</i> , 2002, 22, 303-315.	2.8	6
33	Protein Conformational Transitions from All-Atom Adaptively Biased Path Optimization. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 5372-5382.	5.3	6
34	A minimization principle for transition paths of maximum flux for collective variables. <i>Theoretical Chemistry Accounts</i> , 2017, 136, 1.	1.4	1
35	Editorial overview: Theory & computation. <i>Current Opinion in Structural Biology</i> , 2017, 43, iv-vi.	5.7	1