

Christopher M Cheatum

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9039325/publications.pdf>

Version: 2024-02-01

37
papers

983
citations

430874

18
h-index

434195

31
g-index

37
all docs

37
docs citations

37
times ranked

962
citing authors

#	ARTICLE	IF	CITATIONS
1	Evolution of the Chemical Step in Enzyme Catalysis. <i>ACS Catalysis</i> , 2021, 11, 6726-6732.	11.2	14
2	Least-Squares Fitting of Multidimensional Spectra to Kubo Line-Shape Models. <i>Journal of Physical Chemistry B</i> , 2021, 125, 12876-12891.	2.6	7
3	Evolution of Optimized Hydride Transfer Reaction and Overall Enzyme Turnover in Human Dihydrofolate Reductase. <i>Biochemistry</i> , 2021, 60, 3822-3828.	2.5	3
4	Edge-pixel referencing suppresses correlated baseline noise in heterodyned spectroscopies. <i>Journal of Chemical Physics</i> , 2020, 152, 094201.	3.0	18
5	Low-Frequency Protein Motions Coupled to Catalytic Sites. <i>Annual Review of Physical Chemistry</i> , 2020, 71, 267-288.	10.8	20
6	Optimized reconstructions of compressively sampled two-dimensional infrared spectra. <i>Journal of Chemical Physics</i> , 2019, 150, 234202.	3.0	4
7	Isotopic Labeling of Formate Dehydrogenase Perturbs the Protein Dynamics. <i>Journal of Physical Chemistry B</i> , 2019, 123, 10403-10409.	2.6	14
8	Oscillatory Active-Site Motions Correlate with Kinetic Isotope Effects in Formate Dehydrogenase. <i>ACS Catalysis</i> , 2019, 9, 11199-11206.	11.2	29
9	Evolution Conserves the Network of Coupled Residues in Dihydrofolate Reductase. <i>Biochemistry</i> , 2019, 58, 3861-3868.	2.5	3
10	Evolutionary Effects on Bound Substrate pK_a in Dihydrofolate Reductase. <i>Journal of the American Chemical Society</i> , 2018, 140, 16650-16660.	13.7	17
11	Two-dimensional infrared study of the C D and C O stretching vibrations in strongly hydrogen-bonded complexes. <i>Chemical Physics</i> , 2018, 512, 3-12.	1.9	3
12	Effect of Asp122 Mutation on the Hydride Transfer in <i>E. coli</i> DHFR Demonstrates the Goldilocks of Enzyme Flexibility. <i>Journal of Physical Chemistry B</i> , 2018, 122, 8006-8017.	2.6	11
13	Picosecond Active-Site Dynamics Correlate with the Temperature Dependence of KIEs in Enzyme-Catalyzed Hydride Transfer. <i>FASEB Journal</i> , 2018, 32, .	0.5	0
14	Compressively Sampled Two-Dimensional Infrared Spectroscopy That Preserves Line Shape Information. <i>Journal of Physical Chemistry A</i> , 2017, 121, 3088-3093.	2.5	6
15	Protein Mass Effects on Formate Dehydrogenase. <i>Journal of the American Chemical Society</i> , 2017, 139, 17405-17413.	13.7	17
16	Accelerating two-dimensional infrared spectroscopy while preserving lineshapes using GIRAF. <i>Optics Letters</i> , 2017, 42, 4573.	3.3	5
17	Effects of Isotopic Substitution in Enzyme and Co-factor on Enzyme Catalyzed Hydride Transfer. <i>FASEB Journal</i> , 2017, 31, 764.1.	0.5	0
18	Structural and Kinetic Studies of Formate Dehydrogenase from <i>Candida boidinii</i> . <i>Biochemistry</i> , 2016, 55, 2760-2771.	2.5	76

#	ARTICLE	IF	CITATIONS
19	Oscillatory Enzyme Dynamics Revealed by Two-Dimensional Infrared Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2507-2511.	4.6	33
20	Improved Parametrization for Extended Derjaguin, Landau, Verwey, and Overbeek Predictions of Functionalized Gold Nanosphere Stability. <i>Journal of Physical Chemistry C</i> , 2015, 119, 10064-10075.	3.1	59
21	Line shape analysis of two-dimensional infrared spectra. <i>Journal of Chemical Physics</i> , 2015, 142, 212427.	3.0	76
22	Characterization of Catalytically Relevant Fast Dynamics at the Active Site of Formate Dehydrogenase. <i>FASEB Journal</i> , 2015, 29, 891.1.	0.5	0
23	Relationship of Femtosecond-Picosecond Dynamics to Enzyme-Catalyzed H-Transfer. <i>Topics in Current Chemistry</i> , 2013, 337, 1-39.	4.0	20
24	2D IR Spectroscopy using Four-Wave Mixing, Pulse Shaping, and IR Upconversion: A Quantitative Comparison. <i>Journal of Physical Chemistry A</i> , 2013, 117, 6073-6083.	2.5	56
25	3-Picolyl Azide Adenine Dinucleotide as a Probe of Femtosecond to Picosecond Enzyme Dynamics. <i>Journal of Physical Chemistry B</i> , 2012, 116, 542-548.	2.6	36
26	Hydrogen Donor-Acceptor Fluctuations from Kinetic Isotope Effects: A Phenomenological Model. <i>Biochemistry</i> , 2012, 51, 6860-6870.	2.5	53
27	2D IR Spectroscopy of the C-D stretching vibration of the deuterated formic acid dimer. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 6098.	2.8	19
28	Two-dimensional infrared spectroscopy of azido-nicotinamide adenine dinucleotide in water. <i>Journal of Chemical Physics</i> , 2011, 135, 055106.	3.0	27
29	Characterization of azido-NAD ⁺ to assess its potential as a two-dimensional infrared probe of enzyme dynamics. <i>Analytical Biochemistry</i> , 2010, 407, 241-246.	2.4	19
30	Characterizing the dynamics of functionally relevant complexes of formate dehydrogenase. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010, 107, 17974-17979.	7.1	72
31	Two-dimensional infrared study of 3-azidopyridine as a potential spectroscopic reporter of protonation state. <i>Journal of Chemical Physics</i> , 2010, 133, 134506.	3.0	50
32	Exploring the Molecular Origins of Protein Dynamics in the Active Site of Human Carbonic Anhydrase II. <i>Journal of Physical Chemistry B</i> , 2009, 113, 11505-11510.	2.6	30
33	Efforts toward Developing Probes of Protein Dynamics: Vibrational Dephasing and Relaxation of Carbon-Deuterium Stretching Modes in Deuterated Leucine. <i>Journal of Physical Chemistry B</i> , 2009, 113, 7991-7994.	2.6	26
34	Examination of Enzymatic H-Tunneling through Kinetics and Dynamics. <i>Journal of the American Chemical Society</i> , 2009, 131, 10151-10155.	13.7	42
35	Fast Enzyme Dynamics at the Active Site of Formate Dehydrogenase. <i>Journal of the American Chemical Society</i> , 2008, 130, 22-23.	13.7	80
36	Relaxation and anharmonic couplings of the O-H stretching vibration of asymmetric strongly hydrogen-bonded complexes. <i>Journal of Chemical Physics</i> , 2007, 127, 044501.	3.0	23

#	ARTICLE	IF	CITATIONS
37	Vibrational relaxation of C–D stretching vibrations in CDCl ₃ , CDBr ₃ , and CDI ₃ . <i>Journal of Chemical Physics</i> , 2006, 125, 174503.	3.0	15