

Sayan Bagchi

List of Publications by Year in descending order

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papers

1,322
citations

516710

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477307

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docs citations

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times ranked

1436
citing authors

#	ARTICLE	IF	CITATIONS
1	Ligand Dynamics Time Scales Identify the Surfaceâ€“Ligand Interactions in Thiocyanate-Capped Cadmium Sulfide Nanocrystals. <i>Journal of Physical Chemistry Letters</i> , 2022, 13, 3059-3065.	4.6	6
2	Associationâ€“Dissociation Dynamics of Ionic Electrolytes in Low Dielectric Medium. <i>Journal of Physical Chemistry B</i> , 2022, 126, 239-248.	2.6	3
3	Predicting the Redox Potentials of Phenazine Derivatives Using DFT-Assisted Machine Learning. <i>ACS Omega</i> , 2022, 7, 11742-11755.	3.5	7
4	Onâ€“Off Infrared Absorption of the Sâ•O Vibrational Probe of Dimethyl Sulfoxide. <i>Journal of Physical Chemistry B</i> , 2022, 126, 4501-4508.	2.6	4
5	Reply to â€œComment on â€“Arresting an Unusual Amide Tautomer Using Divalent Cationsâ€™â€•. <i>Journal of Physical Chemistry B</i> , 2021, 125, 479-483.	2.6	1
6	Two Dimensional Infrared Spectroscopy: A Structure Sensitive Technique with Ultrafast Time Resolution. <i>Progress in Optical Science and Photonics</i> , 2021, , 39-56.	0.5	0
7	The Curious Case of Aqueous Warfarin: Structural Isomers or Distinct Excited States?. <i>Journal of Physical Chemistry B</i> , 2021, 125, 2871-2878.	2.6	4
8	Evidence of an nN(amide) â†’ Î€*Ar Interaction in N-Alkyl-N,Nâ€²-diacylhydrazines. <i>Organic Letters</i> , 2021, 23, 7003-7007.	4.6	7
9	Stabilization of Azapeptides by N_{amide}-Hâ€“N_{amide} Hydrogen Bonds. <i>Organic Letters</i> , 2021, 23, 4949-4954.	4.6	9
10	Transition of a Deep Eutectic Solution to Aqueous Solution: A Dynamical Perspective of the Dissolved Solute. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 8784-8789.	4.6	17
11	Electrostatic Manifestation of Micro-Heterogeneous Solvation Structures in Deep-Eutectic Solvents: A Spectroscopic Approach. <i>Journal of Physical Chemistry B</i> , 2020, 124, 3709-3715.	2.6	10
12	Hydrocarbon Chain-Length Dependence of Solvation Dynamics in Alcohol-Based Deep Eutectic Solvents: A Two-Dimensional Infrared Spectroscopic Investigation. <i>Journal of Physical Chemistry B</i> , 2019, 123, 9355-9363.	2.6	20
13	Arresting an Unusual Amide Tautomer Using Divalent Cations. <i>Journal of Physical Chemistry B</i> , 2019, 123, 8419-8424.	2.6	11
14	Interconverting Hydrogen-Bonding and Weak n â†’ Î€* Interactions in Aqueous Solution: A Direct Spectroscopic Evidence. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 5425-5429.	4.6	17
15	Two-Dimensional Infrared Spectroscopy Reveals Cosolvent-Composition-Dependent Crossover in Intermolecular Hydrogen-Bond Dynamics. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 1604-1609.	4.6	32
16	Correlating Nitrile IR Frequencies to Local Electrostatics Quantifies Noncovalent Interactions of Peptides and Proteins. <i>Journal of Physical Chemistry B</i> , 2016, 120, 4034-4046.	2.6	75
17	Electrostatic Interactions Are Key to Câ•O n-Î€* Shifts: An Experimental Proof. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2270-2275.	4.6	14
18	Pick and Choose the Spectroscopic Method to Calibrate the Local Electric Field inside Proteins. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2456-2460.	4.6	9

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19	Cosolvent Effects on Solute-Solvent Hydrogen-Bond Dynamics: Ultrafast 2D IR Investigations. <i>Journal of Physical Chemistry B</i> , 2015, 119, 15334-15343.	2.6	31
20	The non-uniform early structural response of globular proteins to cold denaturing conditions: A case study with Yfh1. <i>Journal of Chemical Physics</i> , 2014, 141, 205103.	3.0	20
21	Extreme electric fields power catalysis in the active site of ketosteroid isomerase. <i>Science</i> , 2014, 346, 1510-1514.	12.6	392
22	Experimental Determination of the Electrostatic Nature of Carbonyl Hydrogen-Bonding Interactions Using IR-NMR Correlations. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 3211-3215.	4.6	32
23	Measuring Electrostatic Fields in Both Hydrogen-Bonding and Non-Hydrogen-Bonding Environments Using Carbonyl Vibrational Probes. <i>Journal of the American Chemical Society</i> , 2013, 135, 11181-11192.	13.7	176
24	Ribonuclease S Dynamics Measured Using a Nitrile Label with 2D IR Vibrational Echo Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2012, 116, 4034-4042.	2.6	100
25	A Solvatochromic Model Calibrates Nitriles' Vibrational Frequencies to Electrostatic Fields. <i>Journal of the American Chemical Society</i> , 2012, 134, 10373-10376.	13.7	107
26	Dynamics of a Myoglobin Mutant Enzyme: 2D IR Vibrational Echo Experiments and Simulations. <i>Journal of the American Chemical Society</i> , 2010, 132, 18367-18376.	13.7	64
27	Conformational Switching between Protein Substates Studied with 2D IR Vibrational Echo Spectroscopy and Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2010, 114, 17187-17193.	2.6	33
28	Equilibrium Exchange Processes of the Aqueous Tryptophan Dipeptide. <i>Journal of Physical Chemistry B</i> , 2009, 113, 8412-8417.	2.6	20
29	2D-IR Experiments and Simulations of the Coupling between Amide-I and Ionizable Side Chains in Proteins: Application to the Villin Headpiece. <i>Journal of Physical Chemistry B</i> , 2009, 113, 11260-11273.	2.6	72
30	Two-Dimensional Infrared Investigation of N-Acetyl Tryptophan Methyl Amide in Solution. <i>Journal of Physical Chemistry B</i> , 2007, 111, 3010-3018.	2.6	29