

Sobhan Sen

List of Publications by Year in descending order

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54
papers

1,868
citations

218677

26
h-index

265206

42
g-index

58
all docs

58
docs citations

58
times ranked

1676
citing authors

#	ARTICLE	IF	CITATIONS
1	Molecular Picture of the Effect of Cosolvent Crowding on Ligand Binding and Dispersed Solvation Dynamics in G-Quadruplex DNA. <i>Journal of Physical Chemistry B</i> , 2022, 126, 1668-1681.	2.6	6
2	Graphene Quantum Dot-Based Optical Sensing Platform for Aflatoxin B1 Detection <i>via</i> the Resonance Energy Transfer Phenomenon. <i>ACS Applied Bio Materials</i> , 2022, 5, 1179-1186.	4.6	24
3	DNA damage, cell cycle perturbation and cell death by naphthalene diimide derivative in gastric cancer cells. <i>Chemico-Biological Interactions</i> , 2022, 358, 109881.	4.0	4
4	Cdr1p highlights the role of the non-hydrolytic ATP-binding site in driving drug translocation in asymmetric ABC pumps. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2020, 1862, 183131.	2.6	12
5	Structure of an Unfolding Intermediate of an RRM Domain of ETR-3 Reveals Its Native-like Fold. <i>Biophysical Journal</i> , 2020, 118, 352-365.	0.5	1
6	Origin of Slow Solvation Dynamics in DNA: DAPI in Minor Groove of Dickerson-Drew DNA. <i>Journal of Physical Chemistry B</i> , 2019, 123, 10202-10216.	2.6	15
7	Role of Ser65, His148 and Thr203 in the Organic Solvent-Dependent Spectral Shift in Green Fluorescent Protein. <i>Photochemistry and Photobiology</i> , 2019, 95, 543-555.	2.5	9
8	Ras hyperactivation versus overexpression: Lessons from Ras dynamics in <i>Candida albicans</i> . <i>Scientific Reports</i> , 2018, 8, 5248.	3.3	5
9	Dynamics of Water and Ions Near DNA: Perspective from Time-Resolved Fluorescence Stokes Shift Experiments and Molecular Dynamics Simulation. <i>Reviews in Fluorescence</i> , 2018, , 231-279.	0.5	2
10	Dynamics of water and ions around DNA: What is so special about them?. <i>Journal of Biosciences</i> , 2018, 43, 499-518.	1.1	14
11	Dynamics of water and ions around DNA: What is so special about them?. <i>Journal of Biosciences</i> , 2018, 43, 499-518.	1.1	4
12	Multidrug ABC transporter Cdr1 of <i>Candida albicans</i> harbors specific and overlapping binding sites for human steroid hormones transport. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2017, 1859, 1778-1789.	2.6	9
13	Effect of T•T Mismatch on DNA Dynamics Probed by Minor Groove Binders: Comparison of Dynamic Stokes Shifts of Hoechst and DAPI. <i>Journal of Physical Chemistry B</i> , 2017, 121, 10735-10748.	2.6	12
14	Probe-location dependent resonance energy transfer at lipid/water interfaces: comparison between the gel- and fluid-phase of lipid bilayer. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 25870-25885.	2.8	9
15	New insight into probe-location dependent polarity and hydration at lipid/water interfaces: comparison between gel- and fluid-phases of lipid bilayers. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 24185-24197.	2.8	14
16	Dispersed dynamics of solvation in G-quadruplex DNA: comparison of dynamic Stokes shifts of probes in parallel and antiparallel quadruplex structures. <i>Methods and Applications in Fluorescence</i> , 2016, 4, 034009.	2.3	8
17	Measuring Size, Size Distribution, and Polydispersity of Water-in-Oil Microemulsion Droplets using Fluorescence Correlation Spectroscopy: Comparison to Dynamic Light Scattering. <i>Journal of Physical Chemistry B</i> , 2016, 120, 1008-1020.	2.6	57
18	Newly identified motifs in <i>Candida albicans</i> Cdr1 protein nucleotide binding domains are pleiotropic drug resistance subfamily-specific and functionally asymmetric. <i>Scientific Reports</i> , 2016, 6, 27132.	3.3	6

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19	Sequence-Dependent Solvation Dynamics of Minor-Groove Bound Ligand Inside Duplex-DNA. <i>Journal of Physical Chemistry B</i> , 2015, 119, 11019-11029.	2.6	23
20	Power-Law Solvation Dynamics in G-Quadruplex DNA: Role of Hydration Dynamics on Ligand Solvation inside DNA. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 1754-1760.	4.6	26
21	Potent Antimalarial Activity of Acriflavine <i>in Vitro</i> and <i>in Vivo</i> . <i>ACS Chemical Biology</i> , 2014, 9, 2366-2373.	3.4	44
22	Rationally Designed Transmembrane Peptide Mimics of the Multidrug Transporter Protein Cdr1 Act as Antagonists to Selectively Block Drug Efflux and Chemosensitize Azole-resistant Clinical Isolates of <i>Candida albicans</i> . <i>Journal of Biological Chemistry</i> , 2013, 288, 16775-16787.	3.4	31
23	Insight into Pleiotropic Drug Resistance ATP-binding Cassette Pump Drug Transport through Mutagenesis of Cdr1p Transmembrane Domains*. <i>Journal of Biological Chemistry</i> , 2013, 288, 24480-24493.	3.4	42
24	Understanding Growth Kinetics of Nanorods in Microemulsion: A Combined Fluorescence Correlation Spectroscopy, Dynamic Light Scattering, and Electron Microscopy Study. <i>Journal of the American Chemical Society</i> , 2012, 134, 19677-19684.	13.7	54
25	Probe Position-Dependent Counterion Dynamics in DNA: Comparison of Time-Resolved Stokes Shift of Groove-Bound to Base-Stacked Probes in the Presence of Different Monovalent Counterions. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2621-2626.	4.6	23
26	Understanding Ligand Interaction with Different Structures of G-Quadruplex DNA: Evidence of Kinetically Controlled Ligand Binding and Binding-Mode Assisted Quadruplex Structure Alteration. <i>Analytical Chemistry</i> , 2012, 84, 7218-7226.	6.5	40
27	Fluorescence Correlation Spectroscopy: An Efficient Tool for Measuring Size, Size-Distribution and Polydispersity of Microemulsion Droplets in Solution. <i>Analytical Chemistry</i> , 2011, 83, 7736-7744.	6.5	113
28	Probe Position Dependence of DNA Dynamics: Comparison of the Time-Resolved Stokes Shift of Groove-Bound to Base-Stacked Probes. <i>Journal of the American Chemical Society</i> , 2010, 132, 9277-9279.	13.7	52
29	“Half-hydration” at the air/water interface revealed by heterodyne-detected electronic sum frequency generation spectroscopy, polarization second harmonic generation, and molecular dynamics simulation. <i>Journal of Chemical Physics</i> , 2010, 132, 144701.	3.0	23
30	Different Molecules Experience Different Polarities at the Air/Water Interface. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 6439-6442.	13.8	33
31	Dynamics of Water and Ions Near DNA: Comparison of Simulation to Time-Resolved Stokes-Shift Experiments. <i>Journal of the American Chemical Society</i> , 2009, 131, 1724-1735.	13.7	86
32	Role of Monovalent Counterions in the Ultrafast Dynamics of DNA. <i>Journal of Physical Chemistry B</i> , 2006, 110, 13248-13255.	2.6	30
33	Ultrafast Dynamics in DNA: “Fraying” at the End of the Helix. <i>Journal of the American Chemical Society</i> , 2006, 128, 6885-6892.	13.7	130
34	Effect of Protein Binding on Ultrafast DNA Dynamics: Characterization of a DNA:APE1 Complex. <i>Biophysical Journal</i> , 2005, 89, 4129-4138.	0.5	32
35	Solvation Dynamics in Dimyristoyl-Phosphatidylcholine Entrapped Inside a Sol [~] Gel Matrix. <i>Journal of Physical Chemistry B</i> , 2004, 108, 2309-2312.	2.6	18
36	Solvation dynamics in a protein “surfactant” aggregate. TNS in HSA “SDS. <i>Chemical Physics Letters</i> , 2003, 379, 471-478.	2.6	20

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37	Solvation dynamics in a protein-surfactant complex. <i>Chemical Physics Letters</i> , 2003, 377, 229-235.	2.6	46
38	Solvation dynamics of 4-aminophthalimide in a polymer (PVP)-surfactant (SDS) aggregate. <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 4875-4879.	2.8	10
39	Solvation Dynamics of a Probe Covalently Bound to a Protein and in an AOT Microemulsion: 4-(N-Bromoacetyl-amino)-Phthalimide. <i>Journal of Physical Chemistry B</i> , 2002, 106, 10741-10747.	2.6	63
40	Excited State Proton Transfer of 1-Naphthol in a Hydroxypropylcellulose/Sodium Dodecyl Sulfate System. <i>Langmuir</i> , 2002, 18, 7867-7871.	3.5	25
41	Solvation Dynamics in the Water Pool of Aerosol Sodium Dioctylsulfosuccinate Microemulsion: Effect of Polymer. <i>Journal of Physical Chemistry A</i> , 2002, 106, 6017-6023.	2.5	34
42	Solvation Dynamics in Bile Salt Aggregates. <i>Journal of Physical Chemistry B</i> , 2002, 106, 7745-7750.	2.6	49
43	Solvation Dynamics in Aqueous Polymer Solution and in Polymer-Surfactant Aggregate. <i>Journal of Physical Chemistry B</i> , 2002, 106, 3763-3769.	2.6	76
44	Isomerization and fluorescence depolarization of merocyanine 540 in polyacrylic acid. Effect of pH. <i>Journal of Chemical Sciences</i> , 2002, 114, 501-511.	1.5	2
45	Photoisomerization of merocyanine 540 in polymer-surfactant aggregate. <i>Journal of Chemical Sciences</i> , 2002, 114, 83-91.	1.5	2
46	Solvation dynamics of TNS in polymer (PEG)-surfactant (SDS) aggregate. <i>Chemical Physics Letters</i> , 2002, 359, 15-21.	2.6	24
47	Femtosecond study of solvation dynamics of DCM in micelles. <i>Chemical Physics Letters</i> , 2002, 359, 77-82.	2.6	59
48	Fluorescence Anisotropy Decay in Polymer-Surfactant Aggregates. <i>Journal of Physical Chemistry A</i> , 2001, 105, 7495-7500.	2.5	79
49	Slow Solvation Dynamics of Dimethylformamide in a Nanocavity. 4-Aminophthalimide in β -Cyclodextrin. <i>Journal of Physical Chemistry A</i> , 2001, 105, 10635-10639.	2.5	56
50	Solvation Dynamics of DCM in Human Serum Albumin. <i>Journal of Physical Chemistry B</i> , 2001, 105, 1438-1441.	2.6	103
51	Solvation dynamics of DCM in micelles. <i>Chemical Physics Letters</i> , 2000, 327, 91-96.	2.6	74
52	Solvation Dynamics of DCM in Dipalmitoyl Phosphatidylcholine Lipid. <i>Tetrahedron</i> , 2000, 56, 6999-7002.	1.9	20
53	Solvation Dynamics of Coumarin 480 in Sol-Gel Matrix. <i>Journal of Physical Chemistry B</i> , 2000, 104, 2613-2616.	2.6	68
54	Excited State Proton Transfer as a Probe for Polymer-Surfactant Interaction. <i>Journal of Physical Chemistry B</i> , 2000, 104, 6128-6132.	2.6	47