

# Ranjan Das

## List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Decoding the Kinetic Pathways toward a Lipid/DNA Complex of Alkyl Alcohol Cationic Lipids Formed in a Microfluidic Channel. <i>Journal of Physical Chemistry B</i> , 2022, 126, 588-600.	2.6	5
2	Essential Loop Dynamics Modulates Catalytic Activity in $\beta$ -Chymotrypsin. <i>ChemistrySelect</i> , 2022, 7, .	1.5	3
3	Host-Assisted Delivery of a Model Drug to Genomic DNA: Key Information from Ultrafast Spectroscopy and <i>in silico</i> Study. <i>ChemBioChem</i> , 2022, , .	2.6	1
4	Oral drug delivery using a polymeric nanocarrier: chitosan nanoparticles in the delivery of rifampicin. <i>Materials Advances</i> , 2022, 3, 4622-4628.	5.4	20
5	Reversible photoswitching of spiropyran in biomolecular interfaces: A combined spectroscopy and computational study. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2022, 430, 113958.	3.9	3
6	Pico-second resolved Förster resonance energy transfer (FRET) differentiates self-assembled biological macromolecules in aqueous medium. <i>Chemical Physics Impact</i> , 2022, 4, 100081.	3.5	1
7	Host assisted molecular recognition by human serum albumin: Study of molecular recognition controlled protein/drug mimic binding in a microfluidic channel. <i>International Journal of Biological Macromolecules</i> , 2021, 176, 137-144.	7.5	4
8	Polyethylene Glycol-Mediated Fusion of Extracellular Vesicles with Cationic Liposomes for the Design of Hybrid Delivery Systems. <i>ACS Applied Bio Materials</i> , 2021, 4, 8259-8266.	4.6	16
9	Deciphering the response of asymmetry in the hydrophobic chains of novel cationic lipids towards biological function. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 1738-1746.	2.8	15
10	Differential flexibility leading to crucial microelastic properties of asymmetric lipid vesicles for cellular transfection: A combined spectroscopic and atomic force microscopy studies. <i>Colloids and Surfaces B: Biointerfaces</i> , 2020, 196, 111363.	5.0	10
11	Flexibility modulates the catalytic activity of a thermostable enzyme: key information from optical spectroscopy and molecular dynamics simulation. <i>Soft Matter</i> , 2020, 16, 3050-3062.	2.7	15
12	Selective and Fast Responsive Sensitized Micelle for Detection of Fluoride Level in Drinking Water. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 16355-16363.	6.7	15
13	Modulation of Kinetic Pathways of Enzyme-Substrate Interaction in a Microfluidic Channel: Nanoscopic Water Dynamics as a Switch. <i>Chemistry - A European Journal</i> , 2019, 25, 9728-9736.	3.3	4
14	Unraveling the Role of Monoolein in Fluidity and Dynamical Response of a Mixed Cationic Lipid Bilayer. <i>Langmuir</i> , 2019, 35, 4682-4692.	3.5	33
15	Probing relaxation dynamics of a cationic lipid based non-viral carrier: a time-resolved fluorescence study. <i>RSC Advances</i> , 2019, 9, 35549-35558.	3.6	12
16	Picosecond Solvation Dynamics in Nanoconfinement: Role of Water and Host-Guest Complexation. <i>Journal of Physical Chemistry B</i> , 2018, 122, 3996-4005.	2.6	7
17	Modulation of Solvation and Molecular Recognition of a Lipid Bilayer under Dynamical Phase Transition. <i>ChemPhysChem</i> , 2018, 19, 2709-2716.	2.1	12
18	A sensitive fluorescent probe for the polar solvation dynamics at protein-surfactant interfaces. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 12237-12245.	2.8	20

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19	Slow solvation dynamics in supramolecular systems based on bile salts: Role of structural rigidity of bile salt aggregates. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2017, 346, 17-23.	3.9	5
20	Environment sensitive fluorescent analogue of biologically active oxazoles differentially recognizes human serum albumin and bovine serum albumin: Photophysical and molecular modeling studies. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2017, 175, 191-199.	3.9	14
21	An Efficient Two-Step Synthesis of 2,5-Disubstituted Oxazole Derivatives Involving Cu-Promoted Carbon-Carbon Single Bond Formation. <i>Asian Journal of Chemistry</i> , 2016, 28, 1519-1522.	0.3	1
22	Conquering 2-Aminopurine's Deficiencies: Highly Emissive Isomorphous Guanosine Surrogate Faithfully Monitors Guanosine Conformation and Dynamics in DNA. <i>Journal of the American Chemical Society</i> , 2015, 137, 3185-3188.	13.7	60
23	Fluorescent Amino Acid Undergoing Excited State Intramolecular Proton Transfer for Site-Specific Probing and Imaging of Peptide Interactions. <i>Journal of Physical Chemistry B</i> , 2015, 119, 2585-2595.	2.6	54
24	Photophysical study of a charge transfer oxazole dye in micelles: Role of surfactant headgroups. <i>Journal of Luminescence</i> , 2015, 163, 21-27.	3.1	7
25	Tuning excited-state proton transfer dynamics of a 3-hydroxychromone dye in supramolecular complexes via host-guest steric compatibility. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 776-784.	2.8	25
26	Sensing Micelle Hydration by Proton-Transfer Dynamics of a 3-Hydroxychromone Dye: Role of the Surfactant Headgroup and Chain Length. <i>Langmuir</i> , 2012, 28, 7147-7159.	3.5	26
27	Unusually slow proton transfer dynamics of a 3-hydroxychromone dye in protic solvents. <i>Photochemical and Photobiological Sciences</i> , 2009, 8, 1583-1589.	2.9	41
28	Excited State Proton Transfer and Solvent Relaxation of a 3-Hydroxyflavone Probe in Lipid Bilayers. <i>Journal of Physical Chemistry B</i> , 2008, 112, 11929-11935.	2.6	43
29	Proton transfer reaction of 4-methyl-2,6-dicarbomethoxyphenol in nonpolar and weakly polar solvents. <i>Journal of Chemical Physics</i> , 2002, 117, 5280-5289.	3.0	9
30	Proton transfer reaction of 4-methyl-2,6-diacetylphenol and an analysis with AM1 potential-energy surfaces. <i>Journal of Chemical Physics</i> , 2001, 114, 1336-1343.	3.0	13
31	Photoinduced proton transfer in 3-methyl-6-hydroxy-m-phthalic acid. <i>Journal of Luminescence</i> , 1999, 81, 61-70.	3.1	13
32	Ground- and Excited-State Proton Transfer in Some <i>o</i> -Hydroxy Aromatic Compounds and Solvent Effect. <i>Israel Journal of Chemistry</i> , 1999, 39, 375-385.	2.3	9
33	Study of proton transfer reactions in binary solvent mixtures by steady state and nanosecond spectroscopy. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 1998, 54, 1073-1081.	3.9	11
34	Energetic and Dynamic Aspects of Intramolecular Proton Transfer in 4-Methyl-2,6-diformylphenol: A Detailed Analysis with AM1 Potential Energy Surfaces. <i>Journal of Physical Chemistry A</i> , 1997, 101, 293-298.	2.5	55
35	Intramolecular Charge Transfer as Probing Reaction: A Fluorescence Monitoring of Protein-Surfactant Interaction. <i>Journal of Physical Chemistry A</i> , 1997, 101, 4042-4047.	2.5	76
36	Dynamical solvent effect on the intramolecular proton transfer reaction of 4-methyl-2,6-diformylphenol in mixed solvents. <i>Journal of Molecular Liquids</i> , 1996, 68, 65-79.	4.9	5

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37	Excited State Proton Transfer Reaction as a Probe for the Microenvironment of a Binding Site of Bovine Serum Albumin: Effect of Urea. <i>The Journal of Physical Chemistry</i> , 1996, 100, 14514-14519.	2.9	33
38	Intramolecular proton transfer in the first excited singlet state of 4-methyl-2,6-diformylphenol: effect of non-polar and weakly polar aprotic solvents. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 1995, 51, 363-369.	3.9	11
39	Complex formation and photophysical properties of luminol: solvent effects. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 1995, 87, 225-230.	3.9	14
40	Effect of the Alkyl Group of Some Alcohols on the Proton-Transfer Reaction of 4-Methyl-2,6-diformylphenol and its Interaction with Triethylamine.. <i>Acta Chemica Scandinavica</i> , 1995, 49, 469-474.	0.7	4
41	Proton transfer and excitation-wavelength-dependent fluorescence and phosphorescence spectra of 4-methyl-2,6-diformyl phenol: interaction with triethylamine. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 1994, 79, 49-54.	3.9	14
42	Ground and excited state proton transfer of 4-methyl-2,6-diformylphenol in different solvent mixtures. <i>Chemical Physics Letters</i> , 1994, 221, 368-372.	2.6	11
43	Emission spectra of 4-methyl-2,6-diacetylphenol in 3-methylpentane at room temperature and 77 K. <i>Chemical Physics Letters</i> , 1994, 228, 393-397.	2.6	20
44	Excitation wavelength and solvent dependent emission spectra in weakly polar aprotic solvents at room temperature and 77 K. <i>Spectrochimica Acta Part A: Molecular Spectroscopy</i> , 1994, 50, 1301-1308.	0.1	16
45	Intramolecular proton transfer in the first-excited electronic states of 4-methyl-2,6-diformyl phenol in some hydrocarbon solvents. <i>Chemical Physics Letters</i> , 1993, 202, 549-554.	2.6	38
46	Proton Transfer in the Ground and Excited Electronic States of 4-Methyl-2,6-diformylphenol : Role of Alcoholic Solvents. <i>Bulletin of the Chemical Society of Japan</i> , 1993, 66, 2492-2497.	3.2	33