

C N Ramachandran

List of Publications by Year in descending order

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50
papers

802
citations

516710

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526287

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docs citations

50
times ranked

863
citing authors

#	ARTICLE	IF	CITATIONS
1	Clustering of Auro-acetylenes via C-Au π Interactions: Gold-Hydrogen Analogy. <i>Chemical Physics Letters</i> , 2022, , 139387.	2.6	1
2	Atoms and molecules in a confined environment. <i>European Physical Journal D</i> , 2021, 75, 1.	1.3	9
3	Density functional studies of the catalytic oxidation of CO using small aurocarbons. <i>Chemical Physics Impact</i> , 2021, 2, 100023.	3.5	4
4	Stacked dimers of Fe-porphyrin with hydrazine and pyrazine as linkers. <i>Chemical Physics Letters</i> , 2021, 779, 138826.	2.6	0
5	Density functional studies on the conversion of hydrogen cyanide to vinyl isocyanide using carbon-supported platinum catalysts. <i>Computational and Theoretical Chemistry</i> , 2021, 1205, 113442.	2.5	2
6	Hydracyanation of acetylene on carbon intercalated gold clusters: Co-operativity and site specificity. <i>Chemical Physics Letters</i> , 2020, 761, 138059.	2.6	5
7	Optoelectronic and charge transport properties of the complex of carbon nanotube with perylene bisimide. <i>International Journal of Quantum Chemistry</i> , 2019, 119, e26026.	2.0	2
8	Isatin π -Triazole π -Functionalized Rhodamine: A Dual Sensor for Cu ²⁺ and Fe ³⁺ Ions and Its Application to Cell Imaging. <i>ChemistrySelect</i> , 2019, 4, 7532-7540.	1.5	11
9	Effect of confinement on structure, energy and vibrational spectra of (HF) _n , n = 1-4. <i>Chemical Physics Letters</i> , 2019, 733, 136670.	2.6	7
10	A multifunctional triazine-based nanoporous polymer as a versatile organocatalyst for CO ₂ utilization and C-C bond formation. <i>Chemical Communications</i> , 2019, 55, 11607-11610.	4.1	24
11	Triazole-appended pyrano[2,3- <i>c</i>]pyrazolone based colorimetric chemosensors for recognition of Fe ³⁺ ions and their molecular logic gate behavior. <i>Analytical Methods</i> , 2019, 11, 3230-3243.	2.7	22
12	Synthesis of a Highly Efficient Multifunctional Copper (II) π -Pyridyl Complex for Adsorption and Photocatalytic Degradation of Organic Dyes. <i>ChemistrySelect</i> , 2019, 4, 4952-4961.	1.5	7
13	Catalytic reduction of SO ₂ by CO using carbon intercalated gold clusters. <i>Chemical Physics Letters</i> , 2019, 726, 111-116.	2.6	9
14	Nitrogen Amelioration-Driven Carbon Dioxide Capture by Nanoporous Polytriazine. <i>Langmuir</i> , 2019, 35, 4893-4901.	3.5	21
15	Spin density transfer from guest to host in endohedral heterofullerene dimers. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 7605-7612.	2.8	1
16	Switching the charge transfer characteristics of quaterthiophene from p-type to n-type <i>via</i> interactions with carbon nanotubes. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 24820-24827.	2.8	0
17	Formation of a nanobubble and its effect on the structural ordering of water in a CH ₄ π -N ₂ π -CO ₂ π -H ₂ O mixture. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 9157-9166.	2.8	13
18	A highly sensitive pyridine-dicarbohydrazide based chemosensor for colorimetric recognition of Cu ²⁺ , AMP ²⁺ , F ⁺ and AcO ⁺ ions. <i>New Journal of Chemistry</i> , 2018, 42, 8567-8576.	2.8	35

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19	Structural, optoelectronic and charge transport properties of the complexes of indigo encapsulated in carbon nanotubes. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 15158-15167.	2.8	6
20	Host-guest and guest-guest interactions in noble gas hydrates. <i>Molecular Physics</i> , 2018, 116, 54-63.	1.7	11
21	Hydrogen-tetrahydrofuran mixed hydrates: A computational study. <i>International Journal of Hydrogen Energy</i> , 2018, 43, 19559-19566.	7.1	14
22	Adsorption and dissolution of methane at the surface of the methanol-water mixture. <i>Fluid Phase Equilibria</i> , 2018, 473, 310-317.	2.5	3
23	Effect of surface roughness on adsorption and distribution of methane at the water-methane interface. <i>Journal of Molecular Liquids</i> , 2018, 266, 856-863.	4.9	9
24	Optoelectronic Properties of Cycloparaphenylene-Carbon Nanotube Based Molecular Architectures. <i>Journal of Physical Chemistry C</i> , 2018, 122, 19904-19912.	3.1	12
25	Structure, Stability, and Properties of Boron Encapsulated Complexes of C ₆₀ , C ₅₉ B, and C ₅₉ N. <i>Journal of Physical Chemistry A</i> , 2017, 121, 1708-1714.	2.5	12
26	Self-Assembling and Luminescent Properties of Chiral Bisoxadiazole Derivatives in Solution and Liquid-Crystalline Phases. <i>Journal of Physical Chemistry B</i> , 2017, 121, 1922-1929.	2.6	12
27	A comprehensive study of the optoelectronic properties of donor-acceptor based derivatives of 1,3,4-oxadiazole. <i>Chemical Physics Letters</i> , 2017, 679, 102-111.	2.6	5
28	Natural Gas Evolution in a Gas Hydrate Melt: Effect of Thermodynamic Hydrate Inhibitors. <i>Journal of Physical Chemistry B</i> , 2017, 121, 153-163.	2.6	26
29	Effect of multiple and adjacent cage occupancies on host-guest interaction and NMR chemical shifts in gas hydrates. <i>Computational and Theoretical Chemistry</i> , 2016, 1092, 57-67.	2.5	8
30	Spin-Spin Coupling in Nitrogen Atom Encapsulated C ₆₀ , C ₅₉ N, and Their Respective Dimers. <i>Journal of Physical Chemistry A</i> , 2016, 120, 6990-6997.	2.5	7
31	Electronic and optical absorption properties of the derivatives of 1,3,4-Oxadiazole. <i>Chemical Data Collections</i> , 2016, 5-6, 88-95.	2.3	4
32	Charge transport and optical properties of the complexes of indigo wrapped over carbon nanotubes. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 14040-14045.	2.8	11
33	Carbon dioxide induced bubble formation in a CH ₄ -CO ₂ -H ₂ O ternary system: a molecular dynamics simulation study. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 3746-3754.	2.8	19
34	Revisiting the structural pattern and the stability of (H ₂ O) ₂₀ clusters using the dispersion corrected density functional method. <i>Chemical Physics Letters</i> , 2015, 626, 39-42.	2.6	4
35	Density functional studies of fused dodecahedral and irregular-dodecahedral water cages. <i>RSC Advances</i> , 2015, 5, 74270-74273.	3.6	4
36	Interaction of rare gas dimers in the confines of a carbon nanotube. <i>Chemical Physics Letters</i> , 2015, 618, 42-45.	2.6	19

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37	Encapsulation of paramagnetic diatomic molecules B ₂ , O ₂ and Ge ₂ inside C ₆₀ . Chemical Physics Letters, 2014, 610-611, 251-255.	2.6	17
38	Structure and Properties of Small Aurocarbons: A Selective Study. Journal of Physical Chemistry A, 2013, 117, 6803-6808.	2.5	9
39	Density functional studies of endosulphan and its interaction with glycine and GABA#. Journal of Chemical Sciences, 2012, 124, 203-207.	1.5	1
40	Encapsulation of the interstellar abundant H ₃ ⁺ in a C ₆₀ fullerene. International Journal of Quantum Chemistry, 2011, 111, 3695-3700.	2.0	3
41	Density functional theoretical studies of the isomers of croconic acid and their dimers. Computational and Theoretical Chemistry, 2011, 973, 28-32.	2.5	20
42	Water clustering in the presence of a CO ₂ molecule. Computational and Theoretical Chemistry, 2011, 966, 84-90.	2.5	16
43	Revisiting the potential energy surface for the He^+H^+ complex. Chemical Physics Letters, 2009, 469, 26-30.	2.6	46
44	Guest species trapped inside carbon nanotubes. Chemical Physics Letters, 2009, 473, 146-150.	2.6	33
45	Host-guest interaction in endohedral fullerenes. Chemical Physics Letters, 2008, 461, 87-92.	2.6	54
46	Time-Dependent Density Functional Theoretical Study of the Absorption Properties of BN-Substituted C ₆₀ Fullerenes. Journal of Physical Chemistry A, 2007, 111, 6901-6903.	2.5	21
47	Hydrogen Peroxide Clusters: The Role of Open Book Motif in Cage and Helical Structures. Journal of Physical Chemistry A, 2006, 110, 6294-6300.	2.5	33
48	Blue Shift in X-H Stretching Frequency of Molecules Due to Confinement. Journal of Physical Chemistry A, 2006, 110, 2-4.	2.5	96
49	Water clusters in a confined nonpolar environment. Chemical Physics Letters, 2005, 410, 348-351.	2.6	94
50	Cis-trans isomerisation and absorption properties of the ring-extended azobenzene. Molecular Physics, 0, , Å.	1.7	0