

# Muthuraja Perumal

## List of Publications by Year in descending order

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17  
papers

297  
citations

933447

10  
h-index

996975

15  
g-index

17  
all docs

17  
docs citations

17  
times ranked

166  
citing authors

#	ARTICLE	IF	CITATIONS
1	Hydrogen bonding interactions on molecular properties of pesticidal compound 4-nitrophthalic acid: experimental density functional theory computations, electron localized function, localized orbital locator analysis and molecular docking scrutiny. <i>Spectroscopy Letters</i> , 2022, 55, 362-372.	1.0	6
2	Analyzing integrated $\pi$ - $\pi$ , C-H $\cdots$ $\pi$ and hydrogen bonding interactions in N,N-Dimethyl-4-aminopyridinium benzilate. <i>Journal of Molecular Structure</i> , 2021, 1242, 130717.	3.6	3
3	Chlorine directed $\pi$ -Cl $\cdots$ $\pi$ and $\pi$ -H $\cdots$ Cl interactions in acridinium 3,5-dichlorosalicylate: Synthesis, X-ray diffraction and theoretical analysis. <i>Journal of Molecular Structure</i> , 2020, 1220, 128759.	3.6	2
4	Bioactivity of a radical scavenger bis(pyrazolium <i>p</i> -toluenesulphonate) on ctDNA and certain microbes: a combined experimental and theoretical analysis. <i>Toxicology Research</i> , 2019, 8, 421-431.	2.1	6
5	Molecular properties, crystal structure, Hirshfeld surface analysis and computational calculations of a new third order NLO organic crystal, 2-aminopyridinium benzilate. <i>Journal of Molecular Structure</i> , 2019, 1181, 118-130.	3.6	34
6	Supramolecular interactions in itaconic acid-2-amino-4,6-dimethylpyrimidine molecular adduct: Physicochemical characterisation and quantum chemical calculations for the molecular adduct. <i>Journal of Physics and Chemistry of Solids</i> , 2019, 129, 284-292.	4.0	2
7	Structural, spectral, physicochemical, computational studies and pharmacological screening of a new organic salt: 2, 6-diaminopyridinium-2-nitrobenzoate. <i>Journal of Molecular Structure</i> , 2019, 1178, 692-701.	3.6	0
8	Hydrogen bonding interactions on 1H-1,2,3-triazole based crystals: Featuring experimental and theoretical analysis. <i>Current Applied Physics</i> , 2018, 18, 774-784.	2.4	15
9	Influence of intramolecular hydrogen bonding interaction on the molecular properties of N-p-tolyl-5-oxo pyrrolidine-3-carboxylic acid: A theoretical and experimental study. <i>Chemical Physics Letters</i> , 2018, 691, 114-121.	2.6	22
10	Growth and characterization of metal halide perovskite crystals: Benzyltributyl ammonium tetrachloro manganate(II) monohydrate. <i>AIP Conference Proceedings</i> , 2018, , .	0.4	0
11	Hydrogen bonded $\pi$ - $\pi$ interactions in a new organic proton transfer crystal aminoguanidinium p-nitrobenzoate monohydrate for optical limiting applications. <i>Journal of Physics and Chemistry of Solids</i> , 2017, 111, 82-94.	3.6	42
12	Hydrogen bonding interactions and supramolecular assemblies in 2-amino guanidinium 4-methyl benzene sulphonate crystal structure: Hirshfeld surfaces and computational calculations. <i>Journal of Molecular Structure</i> , 2017, 1146, 723-734.	3.6	35
13	Theoretical and experimental evaluation of a new organic proton transfer crystal aminoguanidinium p-nitrobenzoate monohydrate for optical limiting applications. <i>Journal of Physics and Chemistry of Solids</i> , 2017, 111, 82-94.	4.0	31
14	Biological and spectral studies of O-Tolyl Biguanide: Experimental and theoretical approach. <i>Journal of Molecular Structure</i> , 2017, 1128, 290-299.	3.6	24
15	Single crystal X-ray diffraction and Hirshfeld surface analyses of supramolecular assemblies in certain hydrogen bonded heterocyclic organic crystals. <i>Journal of Molecular Structure</i> , 2016, 1122, 146-156.	3.6	35
16	Ligand based pharmacophoric modelling and docking of bioactive pyrazolium 3-nitrophthalate (P3NP) on <i>Bacillus subtilis</i> , <i>Aspergillus fumigatus</i> and <i>Aspergillus niger</i> – Computational and Hirshfeld surface analysis. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2016, 163, 352-365.	3.8	29
17	Biological applications and spectroscopic investigations of 4-nitrophenol-urea dimer: A DFT approach. <i>Chemical Physics Letters</i> , 2016, 645, 59-70.	2.6	11