

# Tanveer Hasan

## List of Publications by Year in descending order

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13  
papers

86  
citations

1937685  
4  
h-index

1474206  
9  
g-index

13  
all docs

13  
docs citations

13  
times ranked

120  
citing authors

| #  | ARTICLE   | IF  | CITATIONS |
|----|---|-----|-----------|
| 1  | VIBRATIONAL DYNAMICS AND POTENTIAL ENERGY DISTRIBUTION OF TWO WELL-KNOWN NEUROTRANSMITTER RECEPTORS: TYRAMINE AND DOPAMINE HYDROCHLORIDE. <i>Journal of Theoretical and Computational Chemistry</i> , 2009, 08, 433-450.            | 1.8 | 29        |
| 2  | Molecular structure, vibrational spectra and potential energy distribution of protopine using ab initio and density functional theory. <i>Journal of Structural Chemistry</i> , 2009, 50, 411-420.                                  | 1.0 | 15        |
| 3  | FT-IR spectra and vibrational spectroscopy of Andrographolide. <i>Spectroscopy</i> , 2006, 20, 275-283.   | 0.8 | 11        |
| 4  | Molecular Structure, Vibrational Spectra and Potential Energy Distribution of Colchicine Using ab Initio and Density Functional Theory. <i>Journal of Computer Chemistry Japan</i> , 2009, 8, 59-72.                                | 0.1 | 9         |
| 5  | An investigation on structural, vibrational and nonlinear optical behavior of 4b,9b-dihydroxy-7,8-dihydro-4bH-Indeno[1,2-b] Benzofuran-9,10(6H,9bH)-dione: A DFT study. <i>Journal of Chemical Sciences</i> , 2015, 127, 2217-2223. | 1.5 | 4         |
| 6  | Vibrational Spectra, NBO and NLO Analyses, and A Molecular Docking Study of 3a,8a-Dihydroxy-2-Thioxo-1,3,3a,8a-Tetrahydroindeno[1,2-d]Imidazol-8(2H)-One Using DFT. <i>Journal of Structural Chemistry</i> , 2018, 59, 1078-1087.   | 1.0 | 4         |
| 7  | Normal coordinate analysis and quantum chemical study of tris(p-fluorophenyl)antimony di(N-phenylglycinate) [(p-FC6H4)3Sb(O2CCH2NHC6H5)2]. <i>Pramana - Journal of Physics</i> , 2007, 69, 675-680.                                 | 1.8 | 3         |
| 8  | Efficiently-mapped adaptive hierarchical QAM for reliable scalable video communication. , 2011, , .   |     | 3         |
| 9  | Vibrational dynamics of the organometallic compound triarylorganoantimony (V) SbPh3[O2CC(OH)Ph2]2. <i>Pramana - Journal of Physics</i> , 2007, 68, 875-881.   | 1.8 | 2         |
| 10 | Vibrational and Quantum Chemical Study of Triphenylantimony(V)-o-salicylate. <i>E-Journal of Chemistry</i> , 2008, 5, 723-735.  | 0.5 | 2         |
| 11 | Vibrational Dynamics of the Diterpene-Neoandrographolide. <i>E-Journal of Chemistry</i> , 2009, 6, 183-188.   | 0.5 | 2         |
| 12 | Vibrational analysis of deoxy-andrographolide using MM/QM methods. <i>Spectroscopy</i> , 2007, 21, 279-292.   | 0.8 | 1         |
| 13 | 2,2â€²-(2-Oxopropane-1,3-diyl)bis-(2-hydroxy-1H-indene-1,3(2H)-dione): synthesis, crystal, electronic and molecular docking studies. <i>SN Applied Sciences</i> , 2019, 1, 1.   | 2.9 | 1         |