

# Bruno N C Tenorio

## List of Publications by Citations

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20  
papers

127  
citations

7  
h-index

10  
g-index

26  
ext. papers

211  
ext. citations

3  
avg, IF

2.93  
L-index

#	Paper	IF	Citations
20	Dalton Project: A Python platform for molecular- and electronic-structure simulations of complex systems. <i>Journal of Chemical Physics</i> , <b>2020</b> , 152, 214115	3.9	24
19	Molecular inner-shell photoabsorption/photoionization cross sections at core-valence-separated coupled cluster level: Theory and examples. <i>Journal of Chemical Physics</i> , <b>2019</b> , 150, 224104	3.9	22
18	Coupled Cluster Study of Photoionization and Photodetachment Cross Sections. <i>Journal of Chemical Theory and Computation</i> , <b>2016</b> , 12, 4440-59	6.4	14
17	Time-dependent density functional theory description of total photoabsorption cross sections. <i>Journal of Chemical Physics</i> , <b>2018</b> , 148, 074104	3.9	11
16	Atomic versus molecular Auger decay in CHCl and CDCl molecules. <i>Journal of Chemical Physics</i> , <b>2018</b> , 149, 054303	3.9	10
15	Coupled Cluster and Time-Dependent Density Functional Theory Description of Inner Shell Photoabsorption Cross Sections of Molecules. <i>Journal of Chemical Theory and Computation</i> , <b>2018</b> , 14, 5324-5338	6.4	10
14	Transitions energies, optical oscillator strengths and partial potential energy surfaces of inner-shell states of water clusters. <i>Chemical Physics</i> , <b>2018</b> , 508, 26-33	2.3	8
13	Lanczos-based equation-of-motion coupled-cluster singles-and-doubles approach to the total photoionization cross section of valence excited states. <i>Journal of Chemical Physics</i> , <b>2019</b> , 151, 184106	3.9	5
12	Theoretical study of the absolute inner-shell photoionization cross sections of the formic acid and some of its hydrogen-bonded clusters. <i>Journal of Chemical Physics</i> , <b>2019</b> , 150, 154308	3.9	4
11	Soft X-ray Chlorine Photolysis on Chlorobenzene Ice: An Experimental and Theoretical Study. <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 1389-1398	2.8	3
10	Ab initio study of diffusion of hydrogen, silver and lithium in PbS and Ag <sub>2</sub> S. <i>Computational Materials Science</i> , <b>2019</b> , 166, 75-81	3.2	3
9	Experimental and theoretical results of resonant and normal Auger decay in dichloromethane. <i>European Physical Journal D</i> , <b>2019</b> , 73, 1	1.3	3
8	Fragmentation of Valence and Carbon Core Excited and Ionized CHF <sub>2</sub> CF Molecule. <i>Journal of Physical Chemistry A</i> , <b>2018</b> , 122, 9755-9760	2.8	3
7	Inner-valence Auger decay in chloroform after Cl 2p ionization. <i>Nuclear Instruments &amp; Methods in Physics Research B</i> , <b>2019</b> , 461, 133-136	1.2	2
6	X-ray Photoionization Cross Section Spectra of Water and Ammonia Bonded on Polycyclic Aromatic Hydrocarbons: A Quantum Mechanical Interpretation to the Absorption Spectra on Graphene. <i>Journal of Physical Chemistry A</i> , <b>2020</b> , 124, 2591-2600	2.8	1
5	Evidence of ultrafast dissociation in the CHCl <sub>3</sub> molecule. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , <b>2021</b> , 54, 015202	1.3	1
4	Molecular Photoionization and Photodetachment Cross Sections Based on L <sup>(2)</sup> Basis Sets: Theory and Selected Examples. <i>Progress in Theoretical Chemistry and Physics</i> , <b>2021</b> , 151-179	0.6	1

3	Multi-reference approach to the computation of double core-hole spectra. <i>Journal of Chemical Physics</i> , <b>2021</b> , 155, 131101	3.9	0
2	Insights on the site-selective fragmentation of CF <sub>2</sub> Cl <sub>2</sub> and CH <sub>2</sub> Cl <sub>2</sub> at the chlorine K-edge from ab initio calculations. <i>Chemical Physics</i> , <b>2021</b> , 548, 111226	2.3	0
1	Atomic and molecular Auger decay in CHCl <sub>3</sub> molecule. <i>Journal of Physics: Conference Series</i> , <b>2020</b> , 1412, 122013	0.3	