

Renato Pereira Orenha

List of Publications by Citations

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11
papers

35
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12
ext. papers

48
ext. citations

3.5
avg, IF

1.75
L-index

| # | Paper | IF | Citations |
|----|---|-----|-----------|
| 11 | How computational methods and relativistic effects influence the study of chemical reactions involving Ru-NO complexes?. <i>Journal of Computational Chemistry</i> , 2017 , 38, 883-891 | 3.5 | 16 |
| 10 | How does the total charge and isomerism influence the Ru-NO ammine complexes?. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 13348-13356 | 3.6 | 7 |
| 9 | How does the acidic milieu interfere in the capability of ruthenium nitrosyl complexes to release nitric oxide?. <i>New Journal of Chemistry</i> , 2020 , 44, 773-779 | 3.6 | 4 |
| 8 | On the recognition of chloride, bromide and nitrate anions by anthracene- π -quaramide conjugated compounds: a computational perspective. <i>New Journal of Chemistry</i> , 2020 , 44, 17831-17839 | 3.6 | 3 |
| 7 | Tracking the role of trans-ligands in ruthenium-NO bond lability: computational insight. <i>New Journal of Chemistry</i> , 2020 , 44, 11448-11456 | 3.6 | 2 |
| 6 | The anionic recognition mechanism based on polyol and boronic acid receptors. <i>New Journal of Chemistry</i> , 2020 , 44, 5564-5571 | 3.6 | 1 |
| 5 | Can the relative positions (cis/trans) of ligands really modulate the coordination of NO in ruthenium nitrosyl complexes?. <i>New Journal of Chemistry</i> , 2021 , 45, 1658-1666 | 3.6 | 1 |
| 4 | The design of anion- π interactions and hydrogen bonds for the recognition of chloride, bromide and nitrate anions. <i>Physical Chemistry Chemical Physics</i> , 2021 , 23, 11455-11465 | 3.6 | 1 |
| 3 | Double-bond elucidation for arsagermene with a tricoordinate germanium center: a theoretical survey. <i>New Journal of Chemistry</i> , 2019 , 43, 15681-15690 | 3.6 | 0 |
| 2 | Theoretical study of chloride complexes with hybrid macrocycles. <i>New Journal of Chemistry</i> , 2021 , 45, 463-470 | 3.6 | |
| 1 | The donor/acceptor trans effect on NO release in ruthenium nitrosyl complexes: a computational insight. <i>New Journal of Chemistry</i> , 2021 , 45, 8949-8957 | 3.6 | |