

Renan Vidal Viesser

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Stereoelectronic interactions: A booster for $\langle sup \rangle 4 \langle /sup \rangle \langle i \rangle J \langle /i \rangle \langle sub \rangle HF \langle /sub \rangle$ transmission. Magnetic Resonance in Chemistry, 2022, 60, 481-488.	1.9	1
2	Inverse halogen dependence in anion $\langle sup \rangle 13 \langle /sup \rangle C$ NMR. Physical Chemistry Chemical Physics, 2021, 23, 3019-3030.	2.8	3
3	Counterintuitive deshielding on the $\langle sup \rangle 13 \langle /sup \rangle C$ NMR chemical shift for the trifluoromethyl anion. Magnetic Resonance in Chemistry, 2020, 58, 540-547.	1.9	2
4	Dealing with Hydrogen Bonding on the Conformational Preference of 1,3-Aminopropanols: Experimental and Molecular Dynamics Approaches. Journal of Physical Chemistry A, 2019, 123, 8583-8594.	2.5	9
5	$\langle sup \rangle 1 \langle /sup \rangle \langle i \rangle J \langle /i \rangle \langle sub \rangle CH \langle /sub \rangle$ Coupling in Benzaldehyde Derivatives: Ortho Substitution Effect. ACS Omega, 2019, 4, 1494-1503.	3.5	5
6	The Antagonist Effect of Nitrogen Lone Pair: 3 J HF versus 5 J HF. ChemPhysChem, 2018, 19, 1358-1362.	2.1	7
7	The halogen effect on the $\langle sup \rangle 13 \langle /sup \rangle C$ NMR chemical shift in substituted benzenes. Physical Chemistry Chemical Physics, 2018, 20, 11247-11259.	2.8	34
8	The unexpected roles of \tilde{J}_f and \tilde{J}_C orbitals in electron donor and acceptor group effects on the $\langle sup \rangle 13 \langle /sup \rangle C$ NMR chemical shifts in substituted benzenes. Chemical Science, 2017, 8, 6570-6576.	7.4	39
9	Trends of intramolecular hydrogen bonding in substituted alcohols: a deeper investigation. Physical Chemistry Chemical Physics, 2017, 19, 16904-16913.	2.8	30
10	NMR spin-spin coupling constants: bond angle dependence of the sign and magnitude of the vicinal $\langle sup \rangle 3 \langle /sup \rangle J \langle sub \rangle HF \langle /sub \rangle$ coupling. Physical Chemistry Chemical Physics, 2016, 18, 24119-24128.	2.8	14
11	Experimental and theoretical evaluation of trans-3-halo-2-hydroxy-tetrahydropyran conformational preferences. Beyond anomeric interaction. RSC Advances, 2015, 5, 35412-35420.	3.6	14
12	Effects of stereoelectronic interactions on the relativistic spin-orbit and paramagnetic components of the $\langle sup \rangle 13 \langle /sup \rangle C$ NMR shielding tensors of dihaloethenes. Physical Chemistry Chemical Physics, 2015, 17, 19315-19324.	2.8	18
13	Experimental and Theoretical Studies of Intramolecular Hydrogen Bonding in 3-Hydroxytetrahydropyran: Beyond AIM Analysis. Journal of Physical Chemistry A, 2014, 118, 2794-2800.	2.5	18
14	Influence of OH-N and NH-O inter- and intramolecular hydrogen bonds in the conformational equilibrium of some 1,3-disubstituted cyclohexanes through NMR spectroscopy and theoretical calculations. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2011, 78, 1599-1605.	3.9	12