

Supplementary Material

Striking difference in reactivity of trienes and oximes derived from glucose and xylose

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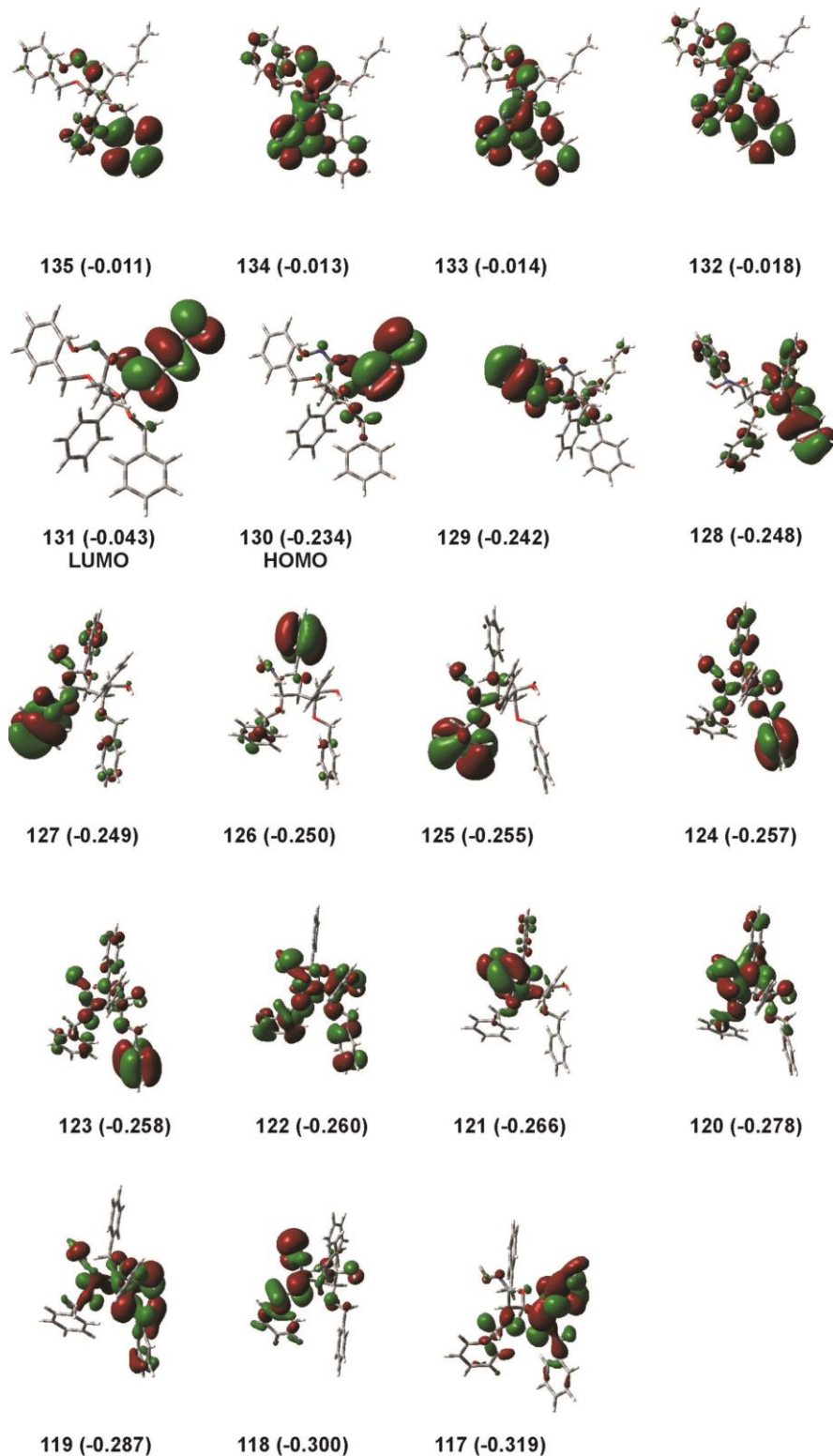
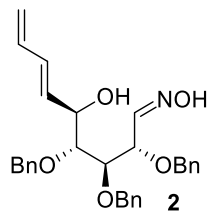
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Calculations of molecular orbitals (MO) were performed by the Gaussian 09 program suite, at the B3LYP/6-311G(2d,p) theory level, using B3LYP/6-31G(2d,p) optimized compound structures. Conformations of compounds were selected arbitrarily. Numbering plan of orbitals, as well as their energy (a.u.), was given as provided by the Gaussian software.

Figure 1S. Selected MO for **2**.

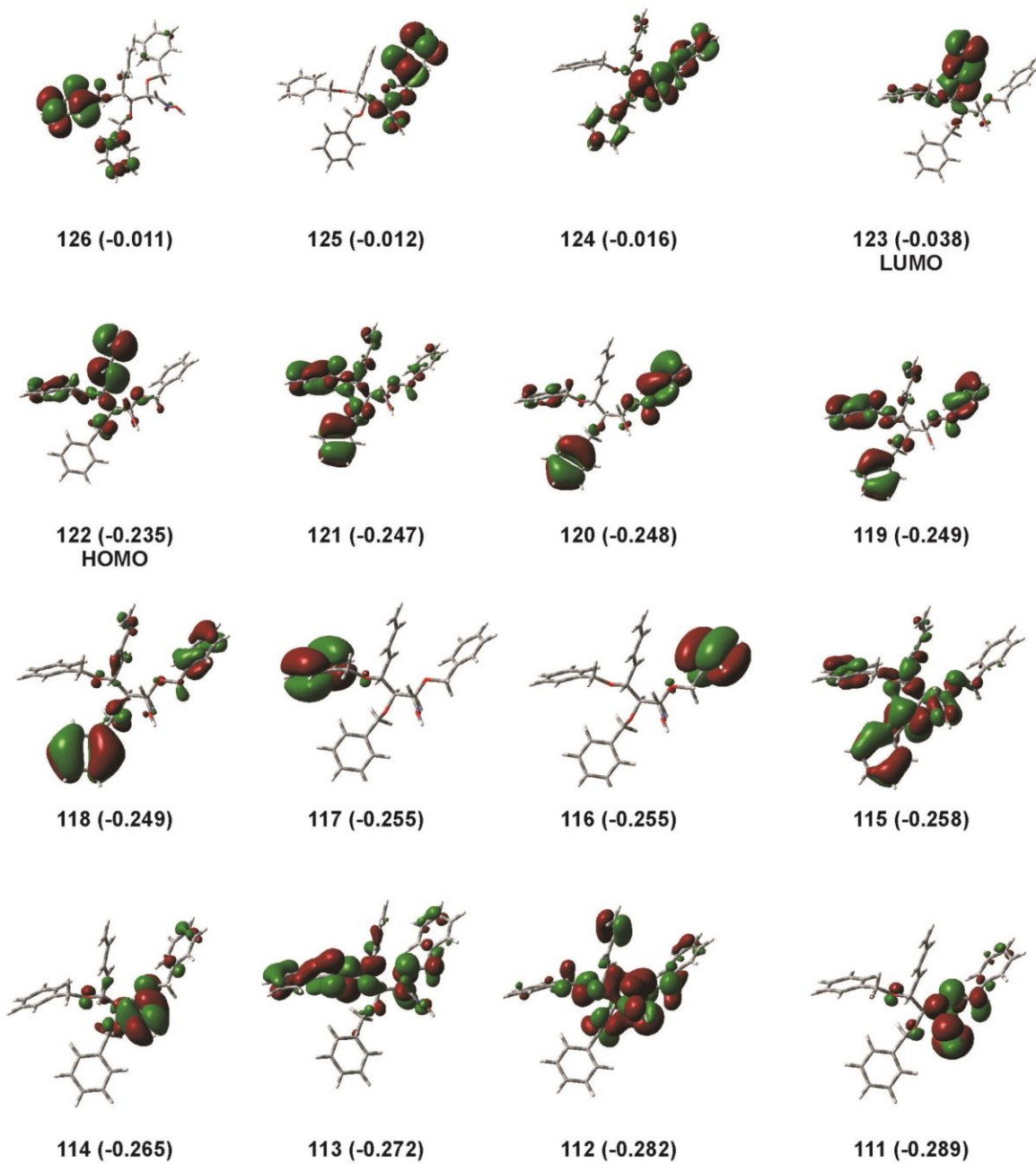
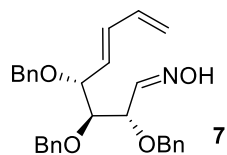


Figure 2S. Selected MO for 7.

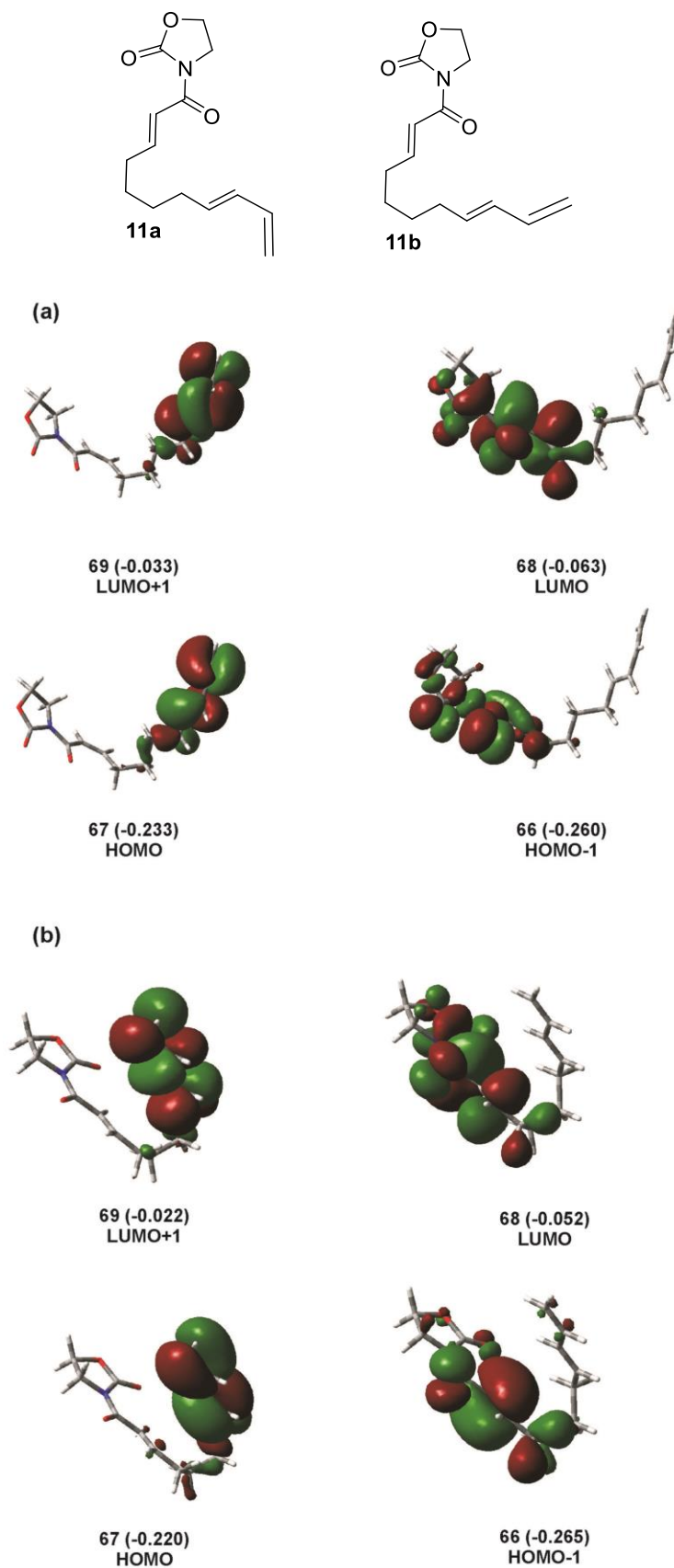


Figure 3S. Selected MO for two conformers of **11**, *s-cis* (a) and *s-trans* (b).

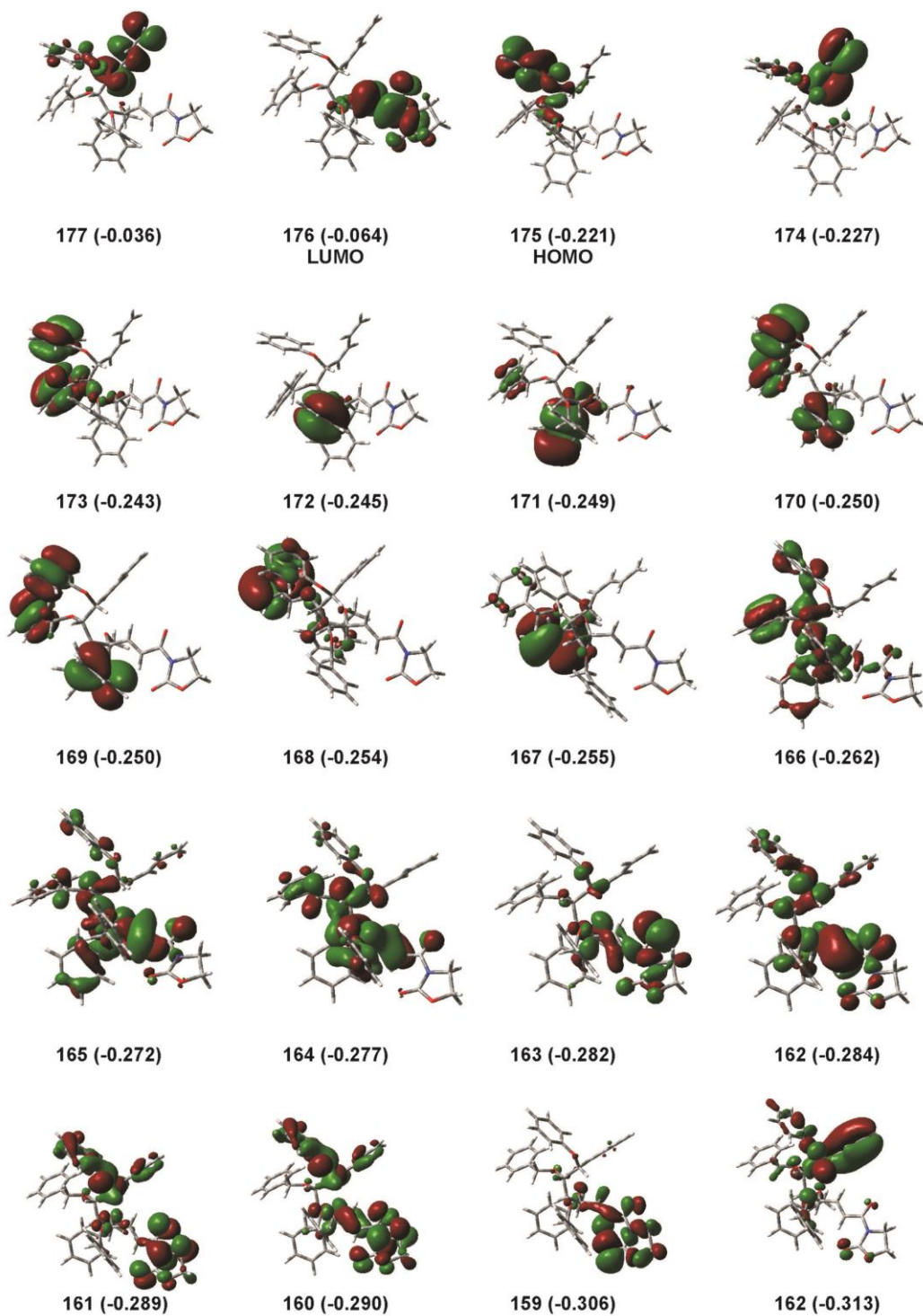
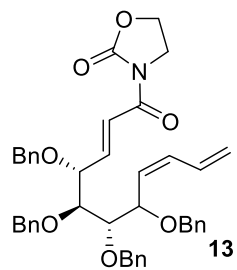


Figure 4S. Some MO for 13.

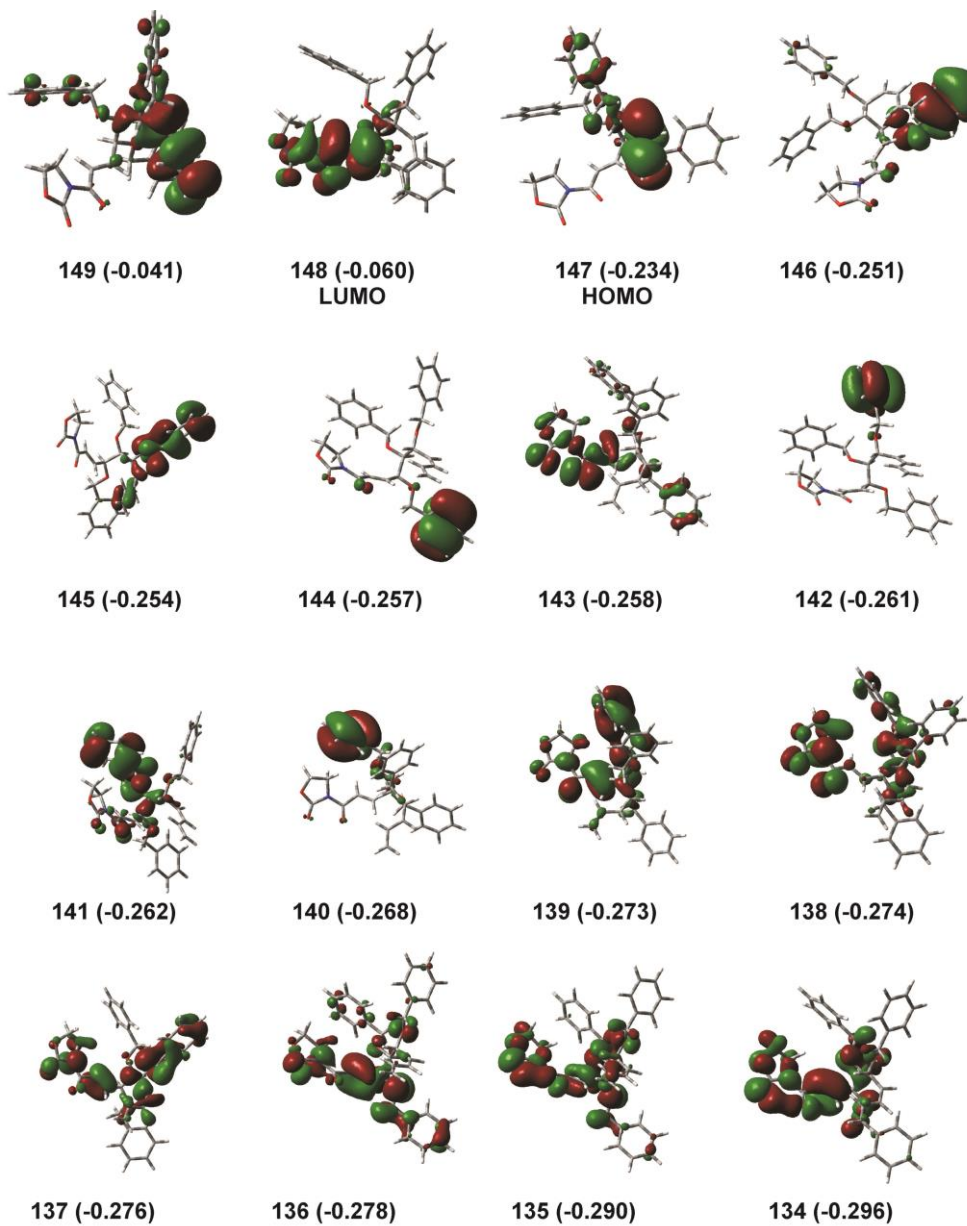
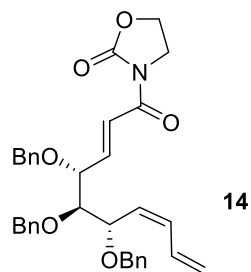


Figure S5. Some molecular orbitals for 14.