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Comparative conformational studies of 3,4,6-tri-*O*-acetyl-1,5-anhydro-2-deoxyhex-1-enitols at the DFT level

Andrzej Nowacki, Beata Liberek



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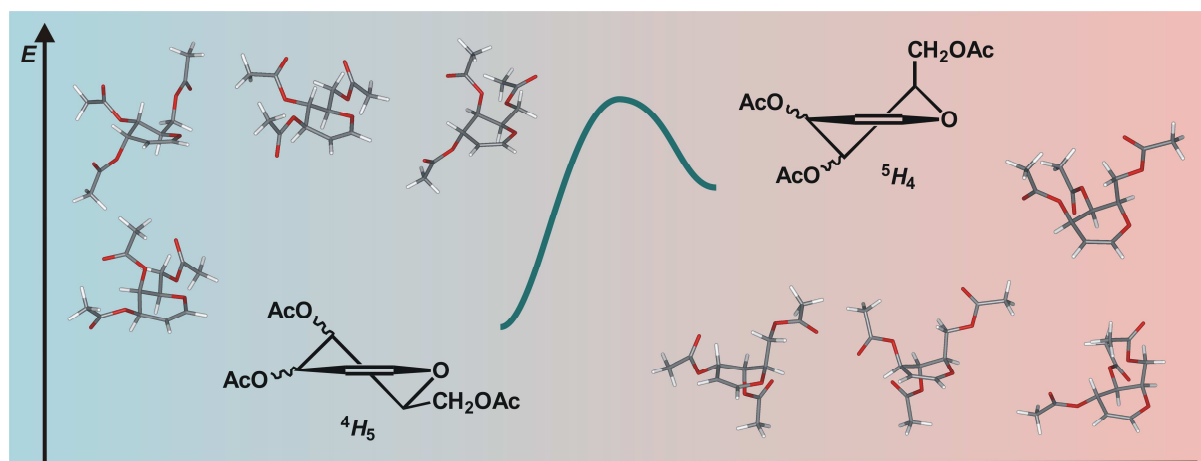
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