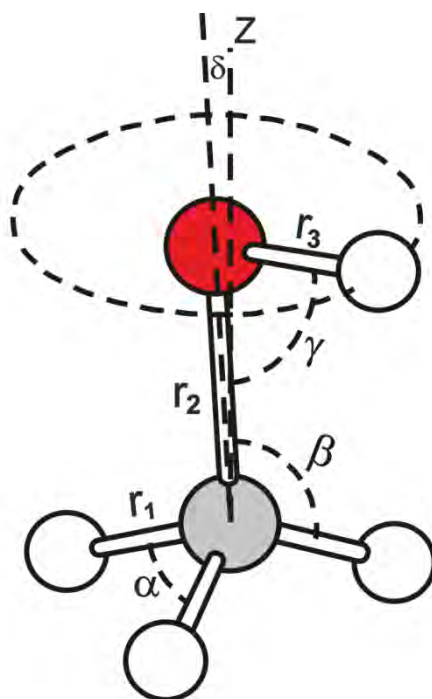


GEDIS Letter

Gas Electron Diffraction Information Service



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Issue February 2020

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CH₂OS H ₂ C=S=O	Sulfine r _e by MW and ab initio <i>J. Demaison, N. Vogt, and D. N. Ksenafontov</i> J. Mol. Struct. (2020), in press
CH₅As CH ₃ AsH ₂	Methylarsine MW spectrum in the ground, first, and second excited torsional state <i>R. A. Motiyenko, V. V. Ilyushin, J. Demaison, N. Vogt, L. Margulès, H. Møllendal, and J.-C. Guillemin</i> J. Mol. Struct., submitted
C₄H₃FN₂O₂	5-Fluorouracil Molecular structure by ED and coupled cluster calculations <i>N. Vogt, D. N. Ksenafontov, and A. N. Rykov</i> Manuscript in preparation
C₄H₄O₃	Succinic anhydride r _e by MW and ab initio <i>M. K. Jahn, D. A. Obenchain, K. P. Rajappan Nair, J.-U. Grabow, N. Vogt, J. Demaison, P. D. Godfrey, and D. McNaughton</i> Phys. Chem. Chem. Phys., (2020) in press
C₅H₄N₂O₄	Orotic acid (1,2,3,6-tetrahydro-2,6-dioxo-4-pyrimidinecarboxylic acid) Molecular structure and conformations by ED and QC <i>N. Vogt, E. P. Altova, A. N. Rykov, and I. F. Shishkov</i> Manuscript in preparation
C₆H₄N₂	2-Pyridinecarbonitrile r _e structure by (ED+MW) and ab initio <i>N. Vogt, L. S. Khaikin, A. N. Rykov, O. E. Grikina, A. A. Batiukov, J. Vogt, I. V. Kochikov, and I. F. Shishkov</i> Struct. Chem. 30 (2019) 1699-1706
C₆H₅NO₂ C ₅ H ₄ NCOOH	Pyridine-3-carboxylic acid (nicotinic acid) Molecular structure and conformations by ED and coupled cluster calculations <i>N. Vogt, I. I. Marochkin, and A. N. Rykov</i> Work in progress
C₆H₁₀S₂ (H ₂ C=CHCH ₂) ₂ S ₂	Diallyl disulfide r _e by MW and ab initio <i>J. Demaison, N. Vogt, R. Tama Saragi, M. Juanes, H.-D. Rudolph, and A. Lesarri</i> Phys. Chem. Chem. Phys. 21 (2019) 19732-19736
C₆H₁₀S₂	Diallyl disulfide Cover Story r _e by MW and ab initio

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$(\text{H}_2\text{C}=\text{CHCH}_2)_2\text{S}_2$	<i>J. Demaison, N. Vogt, R. Tama Saragi, M. Juanes, H.-D. Rudolph, and A. Lesarri</i> Phys. Chem. Chem. Phys. 21 (2019) 19732-19736
$\text{C}_8\text{H}_4\text{FNO}_2$	5-Fluoroisatin r_e from CCSD(T) computations <i>A. V. Belyakov, K. O. Nikolaenko, A. A. Oskorbin, N. Vogt, A. Rykov, and I. F. Shishkov</i> Mol. Phys. 117(14) (2019) 1850-1857
$\text{C}_8\text{H}_4\text{O}_3$	Phthalic anhydride r_e from MW data and CCSD(T) computations <i>A. V. Belyakov, R. Yu. Kulishenko, N. Vogt, and J. Demaison</i> Manuscript in preparation
$\text{C}_8\text{H}_5\text{NO}_2$	Isatin Equilibrium structure from CCSD(T) computations <i>A. V. Belyakov, K. O. Nikolaenko, A. A. Oskorbin, N. Vogt, A. Rykov, and I. F. Shishkov</i> Mol. Phys. 117(14) (2019) 1850-1857
$\text{C}_8\text{H}_6\text{N}_2\text{O}_2$	3-Aminophthalimide r_e structure from ED and coupled cluster computations <i>N. Vogt, D. Savelev, N. I. Giricheva, and G. V. Girichev</i> Manuscript in preparation
$\text{C}_9\text{H}_7\text{NO}_2$	1-Methylisatin r_e from ED data and CCSD(T) computations <i>A. V. Belyakov, K. O. Nikolaenko, A. A. Oskorbin, N. Vogt, A. Rykov, I. F. Shishkov</i> Mol. Phys. 117(14) (2019) 1850-1857
$\text{C}_{12}\text{H}_{10}\text{S}_2$ $\text{C}_6\text{H}_5\text{-SS-C}_6\text{H}_5$	Diphenyl disulfide r_e by MW and ab initio <i>J. Demaison, N. Vogt, R. Tama Saragi, M. Juanes, H.-D. Rudolph, and A. Lesarri</i> ChemPhysChem 20 (2019) 366-373
$\text{C}_{12}\text{H}_{10}\text{S}_2$ $\text{C}_6\text{H}_5\text{-SS-C}_6\text{H}_5$	Diphenyl disulfide Cover Feature r_e by MW and ab initio <i>J. Demaison, N. Vogt, R. Tama Saragi, M. Juanes, H.-D. Rudolph, and A. Lesarri</i> ChemPhysChem 20 (2019) 344
	MOGADOC update <i>J. Vogt, N. Vogt, R. Rudert, K. Deutzmann, and S. Schlagenhauf</i> update in preparation
	Structure Data of Free Polyatomic Molecules <i>N. Vogt and J. Vogt</i> Springer Nature Switzerland, in press
	Multidimensional hypersurfaces Visualization and minima finding of multidimensional hypersurface <i>E. V. Popov, A. A. Batiukov, N. Vogt, T. P. Popova, and J. Vogt</i> IGI Global Disseminator of Knowledge, Chapter 15, in press