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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
C₄N₄S	1,2,5-Thiadiazole-3,4-dicarbonitrile Molecular structure by ED <i>N. V. Tverdova, N. I. Giricheva, D. S. Savelyev, M. S. Mikhailov, N. Vogt, O. I. Koifman, P. A. Stuzhin, and G. V. Girichev</i> Macroheterocyclics, 10 (2017) 27
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
C₅H₉NO₂	Proline <i>r_e</i> structure from MW data and CCSD(T) computations <i>N. Vogt, J. Demaison, S. Krasnoshchekov, N. Stepanov, and H. D. Rudolf</i> Mol. Phys., 115 (2017) 942
C₆H₄N₂	4-Pyridinecarbonitrile <i>r_e</i> structure by (ED+MW) and ab initio <i>L. S. Khaikin, N. Vogt, A. N. Rykov, O. E. Grikina, J. Demaison, J. Vogt, I. V. Kochikov, Y. D. Shishova, E. S. Ageeva, and I. F. Shishkov</i> Russ. J. Phys. Chem., submitted
C₆H₄N₂	3-Pyridinecarbonitrile <i>r_e</i> structure by (ED+MW) and ab initio <i>L. S. Khaikin, N. Vogt, A. N. Rykov, O. E. Grikina, J. Vogt, I. V. Kochikov, E. S. Ageeva, and I. F. Shishkov</i> Mendeleev Comm., submitted
C₆H₄N₂	4-Pyridinecarbonitrile <i>r_e</i> structure by MW and ab initio <i>N. Vogt, K. P. R. Nair, J.-U. Grabow, and J. Demaison</i> Mol. Phys., submitted
C₆H₄N₂	2-Pyridinecarbonitrile <i>r_e</i> structure by (ED+MW) and ab initio <i>N. Vogt, L. S. Khaikin, A. N. Rykov, O. E. Grikina, et al</i> Work in progress
C₆H₅NO₂ C₅H₄NCOOH	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
C₆H₅NO₂ C₅H₄NCOOH	Pyridine-2-carboxylic acid (picolinic acid) Molecular structure and conformations by ED and coupled cluster calculations <i>N. Vogt, I. I. Marochkin, and A. N. Rykov</i> Manuscript submitted
C₆H₁₁F	Fluorocyclohexane r_e , r_s structure by MW and ab initio <i>M. Juanes, N. Vogt, J. Demaison, I. Leon, A. Lesarri, and H. D. Rudolph</i> Phys. Chem. Chem. Phys., 19 (2017) 29162
C₈H₆N₂O₂	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
C₈H₁₂	Ethynylcyclohexane r_e , r_s , r_m structures by MW <i>N. Vogt, J. Demaison, H. D. Rudolph, M. Juanes, J. Fernandez, and A. Lesarri</i> J. Chem. Phys., in press
C₉H₁₅NO	Pseudopelletierine Cover story <i>M. Vallejo-Lopez, P. Ecija, N. Vogt, J. Demaison, A. Lesarri, F. J. Basterretxea, and E. J. Cocinero</i> Chem. Eur. J., 23 (2017) 16412
C₉H₁₅NO	Pseudopelletierine r_e structure by MW and CCSD(T) calculation <i>M. Vallejo-Lopez, P. Ecija, N. Vogt, J. Demaison, A. Lesarri, F. J. Basterretxea, and E. J. Cocinero</i> Chem. Eur. J., 23 (2017) 16491
C₁₆N₁₆S₄Zn	Tetrakis(1,2,5-thiadiazolo)-porphyrazinatozinc(II) r_{h1} structure by ED <i>N. V. Tverdova, N. I. Giricheva, D. S. Savelyev, M. S. Mikhailov, N. Vogt, O. I. Koifman, P. A. Stuzhin, and G. V. Girichev</i> Macroheterocyclics, 10 (2017) 27
	MOGADOC update 2016/2017 <i>J. Vogt, N. Vogt, R. Rudert, K. Deutzmann, and S. Schlagenhauf</i> update in preparation