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<b>C<sub>2</sub>F<sub>6</sub>O<sub>2</sub>S<sub>2</sub></b> CF <sub>3</sub> SO <sub>2</sub> SCF <sub>3</sub>	<b>Trifluoromethanesulfonothioic acid trifluoromethyl ester</b> Structure by ED and ab initio calculations <i>S. L. Masters, D. A. Wann, H. E. Robertson, D. W. H. Rankin, A. Ben Altabef et al.</i> Manuscript in preparation
<b>C<sub>2</sub>H<sub>2</sub>O</b> C(H <sub>2</sub> )CO	<b>Ketene</b> Structure by VHT-GED generation of pyrolysis products <i>S. J. Atkinson, R. Noble-Eddy, and S. L. Masters</i> <i>J. Phys. Chem. A, <b>120</b> (2016) 2041</i>
<b>C<sub>2</sub>H<sub>4</sub>O<sub>2</sub></b> CH <sub>3</sub> C(O)OH	<b>Acetic acid</b> Structure by VHT-GED generation of pyrolysis products <i>S. J. Atkinson, R. Noble-Eddy, and S. L. Masters</i> <i>J. Phys. Chem. A, <b>120</b> (2016) 2041</i>
<b>C<sub>3</sub>H<sub>6</sub>NCl<sub>3</sub></b> N(CH <sub>2</sub> Cl) <sub>3</sub>	<b>Tris(chloromethyl)amine</b> Structure by ED and ab initio calculations, interpretation of Raman spectra <i>C. D. Rankine, S. J. Atkinson, M. R. Waterland, S. L. Masters, and D. A. Wann</i> Manuscript in preparation
<b>C<sub>4</sub>H<sub>6</sub>O<sub>3</sub></b> CH <sub>3</sub> C(O)OC(O)CH <sub>3</sub>	<b>Acetic anhydride</b> Structure by VHT-GED generation of pyrolysis products <i>S. J. Atkinson, R. Noble-Eddy, and S. L. Masters</i> <i>J. Phys. Chem. A, <b>120</b> (2016) 2041</i>
<b>C<sub>6</sub>H<sub>8</sub>O<sub>4</sub></b>	<b>2,2-Dimethyl-1,3-dioxane-4,6-dione (Meldrum's Acid)</b> Structure by GED and ab initio calculations <i>S. J. Atkinson and S. L. Masters</i> Manuscript in preparation
<b>C<sub>7</sub>H<sub>7</sub></b> C <sub>6</sub> H <sub>5</sub> CH <sub>2</sub> .	<b>Benzyl radical</b> Structure by VHT-GED generation of pyrolysis products <i>R. Noble-Eddy, B. H. C. Wilson, and S. L. Masters</i> Manuscript in preparation
<b>C<sub>7</sub>H<sub>16</sub>Cl<sub>3</sub>PSi</b> (tBu)(iPr)PSiCl <sub>3</sub>	<b>(Tert-butyl)(iso-propyl)(trichlorosilyl)phosphine</b> Structure by GED, UCONGA and ab initio / DFT methods <i>I. Wagner, E. Seppälä, H. E. Robertson W.-W. du Mont, and S. L. Masters</i> Manuscript in preparation
<b>C<sub>12</sub>H<sub>36</sub>P<sub>2</sub>Si<sub>4</sub></b> P <sub>2</sub> (SiMe <sub>3</sub> ) <sub>4</sub>	<b>1,1,2,2-Tetrakis(trimethylsilyl)diphosphane</b> Structure by GED, UCONGA and ab initio / DFT methods <i>A. P. Flanagan, H. Humphrey-Taylor, N. R. Gunby, H. E. Robertson, and S. L. Masters</i> Manuscript complete
<b>C<sub>14</sub>H<sub>14</sub>O<sub>2</sub>S</b> (C <sub>6</sub> H <sub>5</sub> CH <sub>2</sub> ) <sub>2</sub> SO <sub>2</sub>	<b>Dibenzyl sulfone</b> Structure by GED, UCONGA and ab initio / DFT methods <i>R. Noble-Eddy, B. H. C. Wilson, and S. L. Masters</i> Manuscript in preparation
	<b>Hexakis(trimethylsilyl)disilane</b>

<b>C<sub>18</sub>H<sub>54</sub>Si<sub>8</sub></b> Si <sub>2</sub> (SiMe <sub>3</sub> ) <sub>6</sub>	Vibrational spectra and structure by GED and ab initio calculations <i>K. Hassler, H. E. Robertson, S. L. Masters et al.</i> Manuscript in preparation
	<b>Apparatus development</b> York time-averaged electron diffractometer  <i>M. H. P. Ardebili, R. S. Fender, M. A. D. Fluendy, S. A. Hayes, P. D. Lane, S. L. Masters, R. J. Mawhorter, J. P. F. Nunes, P. Papathomas, D. W. H. Rankin, C. D. Rankine, D. A. Wann, and S. Young</i> Manuscript in preparation
	<b>Method development</b> Universal Conformer Generation and Analysis (UCONGA) <i>N. R. Gunby, S. L. Masters, and D. L. Crittenden</i> Manuscript complete