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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 GED 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>6</sub>HCl<sub>5</sub></b> C <sub>6</sub> HCl <sub>5</sub>	<b>Pentachlorobenzene</b> Structure by ED and liquid crystal NMR spectroscopy <i>R. Blom, D. A. Wann, D. W. H. Rankin et al.</i> Manuscript in preparation.
<b>C<sub>6</sub>H<sub>2</sub>Cl<sub>4</sub></b> 1,2,3,4-C <sub>6</sub> H <sub>2</sub> Cl <sub>4</sub>	<b>1,2,3,4-Tetrachlorobenzene</b> Structure by ED and liquid crystal NMR spectroscopy <i>R. Blom, D. A. Wann, D. W. H. Rankin et al.</i> Manuscript in preparation.
<b>C<sub>6</sub>H<sub>12</sub>F<sub>6</sub>Si<sub>2</sub></b> CF <sub>3</sub> Me <sub>2</sub> SiSiMe <sub>2</sub> CF <sub>3</sub>	<b>1,2-Trifluoromethyl-1,1,2,2-tetramethyldisilane</b> Structure by ED, X-ray diffraction and ab initio calculations, interpretation of Raman spectra <i>S. L. Masters, H. E. Robertson, D. A. Wann, M. Hölbling, K. Hassler, R. Björnsson, S. Ó. Wallevik and I. Arnason</i> <i>J. Phys. Chem. A, in press</i>
<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 ED and ab initio calculations <i>E. Seppälä, W.-W. du Mont, S. L. Masters, D. W. H. Rankin, and H. E. Robertson</i> Manuscript in preparation
<b>C<sub>10</sub>H<sub>20</sub>N<sub>2</sub></b> ButN=CHCH=NBut	<b>N,N'-Di-tert-butyl-1,4-diaza-1,3-butadiene</b> Structure by ED and computational methods <i>C. Jones, D. A. Wann, H. E. Robertson, and D. W. H. Rankin</i> Manuscript in preparation.
<b>C<sub>11</sub>H<sub>30</sub>Br<sub>2</sub>Si<sub>4</sub></b> C(SiMe <sub>3</sub> ) <sub>2</sub> (SiMe <sub>2</sub> Br) <sub>2</sub>	<b>Bis(bromodimethylsilyl)bis(trimethylsilyl)methane</b> Structure by ED and computational methods <i>D. A. Wann, K. Bätz, M. S. Robinson, S. L. Masters, H. E. Robertson and P. D. Lickiss</i> <i>J. Phys. Chem. A, in press</i>
<b>C<sub>11</sub>H<sub>30</sub>Cl<sub>2</sub>Si<sub>4</sub></b> (Me <sub>3</sub> Si) <sub>2</sub> C(SiClMe <sub>2</sub> ) <sub>2</sub>	<b>Bis(chlorodimethylsilyl)bis(trimethylsilyl)methane</b> Structure by ED and computational methods <i>D. A. Wann, K. Bätz, M. S. Robinson, S. L. Masters, H. E. Robertson and P. D. Lickiss</i> <i>J. Phys. Chem. A, in press</i>
<b>C<sub>11</sub>H<sub>32</sub>Si<sub>4</sub></b> C(SiMe <sub>3</sub> ) <sub>2</sub> (SiMe <sub>2</sub> H) <sub>2</sub>	<b>Bis(dimethylsilyl)bis(trimethylsilyl)methane</b> Structure by ED and computational methods <i>D. A. Wann, K. Bätz, M. S. Robinson, S. L. Masters, H. E. Robertson and P. D. Lickiss</i> <i>J. Phys. Chem. A, in press</i>
<b>C<sub>12</sub>Fe<sub>3</sub>O<sub>12</sub></b>	<b>Dodecacarbonyltriiron</b> Structure by ED and ab initio calculations

Fe3(CO)12

*G. R. Kafka, S. L. Masters, D. W. H. Rankin et al.*  
Manuscript in preparation.

**Apparatus development**

York time-averaged electron diffractometer

*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*  
Manuscript in preparation