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<b>B<sub>7</sub>H<sub>9</sub>S<sub>2</sub></b> arachno-B <sub>7</sub> H <sub>9</sub> S <sub>2</sub>	<b>arachno-Dithianonaborane(9)</b> Structure by ED and ab initio calculations <i>D. A. Wann, P. D. Lane, H. E. Robertson, J. Holub, and D. Hnyk</i> Inorg. Chem., <b>52</b> (2013), 4502
<b>BrNa</b> NaBr	<b>Sodium bromide</b> Structure by ED and ab initio calculations <i>D. A. Wann, D. W. H. Rankin, P. D. McCaffrey, and R. J. Mawhorter</i> Manuscript submitted
<b>Br<sub>2</sub>Na<sub>2</sub></b> Na <sub>2</sub> Br <sub>2</sub>	<b>Sodium bromide dimer</b> Structure by ED and ab initio calculations <i>D. A. Wann, D. W. H. Rankin, P. D. McCaffrey, and R. J. Mawhorter</i> Manuscript submitted
<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>4</sub>O<sub>4</sub>S<sub>2</sub></b> O <sub>2</sub> -cyclo-S(CH <sub>2</sub> ) <sub>2</sub> S-O <sub>2</sub>	<b>1,3-Dithietane-1,1,3,3-tetraoxide</b> Structure by ED and ab initio calculations <i>D. A. Wann, A. Bil, P. D. Lane, H. E. Robertson, D. W. H. Rankin, and E. Block</i> Struct. Chem., <b>24</b> (2013), 827
<b>C<sub>2</sub>H<sub>12</sub>B<sub>10</sub>S<sub>2</sub></b> 9,12-(SH) <sub>2</sub> -closo- <sup>-1,2-</sup> C <sub>2</sub> B <sub>10</sub> H <sub>10</sub>	<b>9,12-Dimercapto-closo-1,2-dicarbado-decarborane(12)</b> Structure by ED and ab initio calculations <i>D. A. Wann, P. D. Lane, H. E. Robertson, T. Base, and D. Hnyk</i> Dalton Trans., <b>42</b> (2013), 12015
<b>C<sub>2</sub>H<sub>13</sub>B<sub>7</sub></b> arachno-B <sub>7</sub> H <sub>9</sub> (CH <sub>2</sub> ) <sub>2</sub>	<b>arachno-6,9-Dicarbanoaborane(13)</b> Structure by ED and ab initio calculations <i>D. A. Wann, P. D. Lane, H. E. Robertson, J. Holub, and D. Hnyk</i> Inorg. Chem., <b>52</b> (2013), 4502
<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, D. A. Wann, H. E. Robertson, F. Lennox, D. W. H. Rankin, I. Arnason, K. Hassler et al.</i> Manuscript in preparation
<b>C<sub>9</sub>H<sub>24</sub>Br<sub>4</sub>Si<sub>4</sub></b> C(SiMe <sub>2</sub> Br) <sub>4</sub>	<b>Tetrakis(bromodimethylsilyl)methane</b> Structure by ED and computational methods <i>D. A. Wann, K. Bätz, S. Young, S. L. Masters, H. E. Robertson, and P. D. Lickiss</i> Manuscript completed
<b>C<sub>9</sub>H<sub>24</sub>Cl<sub>4</sub>Si<sub>4</sub></b>	<b>Tetrakis(chlorodimethylsilyl)methane</b> Structure by ED and computational methods

$C(SiMe_2Cl)_4$	<i>D. A. Wann, K. Bätz, S. Young, S. L. Masters, H. E. Robertson, and P. D. Lickiss</i> Manuscript completed
$C_9H_{24}F_4Si_4$ $C(SiMe_2F)_4$	<b>Tetrakis(fluorodimethylsilyl)methane</b> Structure by ED and computational methods <i>D. A. Wann, K. Bätz, S. Young, S. L. Masters, H. E. Robertson, and P. D. Lickiss</i> Manuscript completed
$C_9H_{28}Si_4$ $C(SiMe_2H)_4$	<b>Tetrakis(dimethylsilyl)methane</b> Structure by ED and computational methods <i>D. A. Wann, K. Bätz, S. Young, S. L. Masters, H. E. Robertson, and P. D. Lickiss</i> Manuscript completed
$C_{11}H_{30}Br_2Si_4$ $C(SiMe_3)_2(SiMe_2Br)_2$	<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> Manuscript completed
$C_{11}H_{30}Cl_2Si_4$ $(Me_3Si)_2C(SiClMe_2)_2$	<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> Manuscript completed
$C_{11}H_{32}Si_4$ $C(SiMe_3)_2(SiMe_2H)_2$	<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> Manuscript completed
$C_{12}H_{38}Si_6$ $Si_2H_2(SiMe_3)_4$	<b>1,1,2,2-Tetrakis(trimethylsilyl)disilane</b> Structure by ED and ab initio calculations <i>J. Schwabedissen, P. D. Lane, S. L. Masters, K. Hassler, and D. A. Wann</i> Manuscript completed
$C_{14}H_{42}Si_6$ $(SiMe_3)_2MeSiSiMe(SiMe_3)_2$	<b>1,1,2,2-Tetrakis-trimethylsilyl-1,2-dimethyldisilane</b> Structure by ED and ab initio calculations <i>J. Schwabedissen, P. D. Lane, S. L. Masters, K. Hassler, and D. A. Wann</i> Manuscript completed
<b>FNa</b> NaF	<b>Sodium fluoride</b> Structure by ED and ab initio calculations <i>D. A. Wann, D. W. H. Rankin, P. D. McCaffrey, and R. J. Mawhorter</i> Manuscript submitted
<b>F<sub>2</sub>Na<sub>2</sub></b> Na <sub>2</sub> F <sub>2</sub>	<b>Sodium fluoride dimer</b> Structure by ED and ab initio calculations <i>D. A. Wann, D. W. H. Rankin, P. D. McCaffrey, and R. J. Mawhorter</i> Manuscript submitted
<b>INa</b> NaI	<b>Sodium iodide</b> Structure by ED and ab initio calculations <i>D. A. Wann, D. W. H. Rankin, P. D. McCaffrey, and R. J. Mawhorter</i> Manuscript submitted
<b>I<sub>2</sub>Na<sub>2</sub></b> Na <sub>2</sub> I <sub>2</sub>	<b>Sodium iodide dimer</b> Structure by ED and ab initio calculations <i>D. A. Wann, D. W. H. Rankin, P. D. McCaffrey, and R. J. Mawhorter</i> Manuscript submitted