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<b>C<sub>2</sub>H<sub>3</sub>Cl<sub>3</sub>Ge</b> CH <sub>2</sub> =CH-GeCl <sub>3</sub>	<b>Trichloroethenylgermane</b> GED structure, quantum chemical calculations on CH <sub>2</sub> =CH-MX <sub>3</sub> (M=C, Si, Ge, Sn and X=H, Cl), rotational barrier for the MX <sub>3</sub> group <i>S. Samdal, J.-C. Guillemin and S. Gundersen</i> J. Phys. Chem., submitted
<b>C<sub>3</sub>H<sub>5</sub>ClO</b> CHCICHCH <sub>2</sub> OH	<b>cis-3-Chloro-2-propen-1-ol</b> GED structure, quantum chemical calculations, conformations, large amplitude motion <i>T. G. Strand, S. Gundersen, H. Priebe, S. Samdal, R. Seip</i> J. Mol. Struct. <b>921</b> (2009) 72
<b>C<sub>3</sub>H<sub>6</sub>F<sub>2</sub>Si</b> CH <sub>2</sub> =CHCH <sub>2</sub> SiF <sub>2</sub> H	<b>Allyldifluorosilane</b> Microwave, quantum chemical calculations, conformations <i>H. Møllendal, S. Samdal, G. A. Guirgis and C. J. Wurrey</i> J. Phys. Chem., submitted
<b>C<sub>4</sub>Cl<sub>2</sub>N<sub>4</sub></b>	<b>Bis(chloroimino)butanedinitrile</b> GED structure, quantum chemical calculations, conformations <i>H. Thomassen, S. Gundersen and S. Samdal</i> J. Mol. Struct. <b>928</b> (2009) 182
<b>C<sub>6</sub>H<sub>5</sub>O<sub>2</sub></b>	<b>1,2-Cyclohexanedione</b> GED structure, quantum chemical calculations, conformations <i>Q. Shen, M. Traetteberg and S. Samdal</i> J. Mol. Struct. <b>923</b> (2009) 94
<b>C<sub>10</sub>Cl<sub>10</sub>Fe</b> Fe(C <sub>5</sub> Cl <sub>5</sub> ) <sub>2</sub>	<b>Decachloroferrocene</b> GED structure, quantum chemical calculations, barrier to internal rotation <i>L. Phillips, M. K. Cooper, A. Haaland, S. Samdal, N. I. Giricheva and G. V. Girichev</i> J. Chem. Soc. Dalton Trans., accepted