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<b>C<sub>18</sub>H<sub>16</sub>Si</b> HSiPh <sub>3</sub>	<b>Triphenylsilane</b> Structure by ED and quantum chemical calculations; structural variations in H <sub>4</sub> -nSiPh <sub>n</sub> molecules (n = 1-4) <i>A. R. Campanelli, A. Domenicano, F. Ramondo and I. Hargittai</i> Struct. Chem. (2011) published online
Ph-C <sub>8</sub> H <sub>6</sub> -X	<b>Phenylcubane derivatives</b> Structure by quantum chemical calculations, analysis of structural variations, effect of substituents on molecular geometry <i>A. R. Campanelli, A. Domenicano and F. Ramondo</i> Struct. Chem. (2011) submitted
	<b>Benzene derivatives</b> Comparison of the structures of five benzene derivatives determined by ED in two different laboratories <i>A. R. Campanelli, A. Domenicano and I. Hargittai</i> Struct. Chem. <b>21</b> (2010), 803
Ph-[C(CH <sub>2</sub> ) <sub>3</sub> C] <sub>n</sub> -X	<b>[n]Staffane derivatives, n = 1-5</b> Structure by quantum chemical calculations, analysis of structural variations, effect of substituents on molecular geometry <i>A. R. Campanelli, A. Domenicano, G. Piacente and F. Ramondo</i> J. Phys. Chem. A <b>114</b> (2010), 5162