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<b>C<sub>7</sub>H<sub>5</sub>N</b> C <sub>6</sub> H <sub>5</sub> CN	<b>Cyanobenzene</b> Structure by ED and quantum chemical calculations <i>A. R. Campanelli, A. Domenicano, F. Ramondo and I. Hargittai</i> <i>J. Phys. Chem. A</i> <b>112</b> (2008) 10998-11008
<b>C<sub>8</sub>H<sub>4</sub>N<sub>2</sub></b> C <sub>6</sub> H <sub>4</sub> (CN) <sub>2</sub>	<b>1,4-Dicyanobenzene</b> Structure by ED and quantum chemical calculations <i>A. R. Campanelli, A. Domenicano, F. Ramondo and I. Hargittai</i> <i>J. Phys. Chem. A</i> <b>112</b> (2008) 10998-11008
<b>C<sub>10</sub>H<sub>2</sub>N<sub>4</sub></b> C <sub>6</sub> H <sub>2</sub> (CN) <sub>4</sub>	<b>1,2,4,5-Tetracyanobenzene</b> Structure by ED and quantum chemical calculations <i>A. R. Campanelli, A. Domenicano, F. Ramondo and I. Hargittai</i> <i>J. Phys. Chem. A</i> <b>112</b> (2008) 10998-11008
	<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., submitted
Ph-[C(CH <sub>2</sub> ) <sub>3</sub> C]n-X	<b>[n]Staffane derivatives</b> Structure by quantum chemical calculations, analysis of structural variation, effect of substituents on molecular geometry <i>A. R. Campanelli, A. Domenicano, G. Piacente and F. Ramondo</i> <i>J. Phys. Chem. A</i> <b>114</b> (2010) published in the website