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C₈H₄N₂ C ₆ H ₄ (NC) ₂	1,4-Diisocyanobenzene Structure by ED and quantum chemical calculations, comparison with solid-state results <i>A. R. Campanelli, A. Domenicano, F. Ramondo and I. Hargittai</i> Struct. Chem., 23 (2012), 287
C₁₈H₁₆Si HSiPh ₃	Triphenylsilane Structure by ED and quantum chemical calculations; structural variations in H ₄ -nSiPh _n molecules (n = 1-4) <i>A. R. Campanelli, A. Domenicano, F. Ramondo and I. Hargittai</i> Struct. Chem., 22 (2011), 361
Ph-C ₈ H ₆ -X	Phenylcubane derivatives 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., 22 (2011), 449
Ph-C ₆ H ₄ -X	4-Substituted biphenyl derivatives Structure by quantum chemical calculations, analysis of structural variation, field and resonance effects <i>A. R. Campanelli, A. Domenicano and F. Ramondo</i> 14th European Symposium on Gas Phase Electron Diffraction, Moscow, June 24-28, 2011. Book of Abstracts, p. 10
Ph-X	Monosubstituted benzene derivatives Structure by quantum chemical calculations at different levels of theory, effect of substituents on molecular geometry, electronegativity and resonance parameters <i>A. R. Campanelli, A. Domenicano, M. Macchiagodena and F. Ramondo</i> Struct. Chem., 22 (2011), 1131