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<p>$\text{Ph-C}(\text{CH}=\text{CH})_3\text{C-X}$</p>	<p>4-Substituted 1-phenylbarrelene derivatives Structure by quantum chemical calculations, analysis of structural variation, transmission of field effects through the barrelene framework, field-induced p-polarization, electric dipole moments, comparison with bicyclo[2.2.2]octane derivatives <i>A. R. Campanelli, A. Domenicano, and F. Ramondo</i> Struct. Chem., in press</p>
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