

Acta Crystallographica Section E

# Structure Reports Online

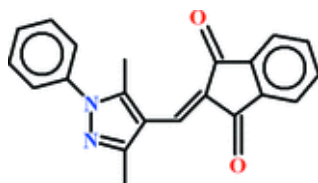
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**organic compounds**

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<a href="#">html</a>	<a href="#">pdf</a>	<a href="#">cif</a>	<a href="#">3d view</a>	<a href="#">structure factors</a>	<a href="#">supplementary materials</a>	<a href="#">checkCIF</a>	<a href="#">similar papers</a>	<a href="#">contents of issue</a>	<a href="#">open access</a>
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## 2-[(3,5-Dimethyl-1-phenyl-1*H*-pyrazol-4-yl)methylidene]indan-1,3-dione

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**Abstract:** In the title compound, C<sub>21</sub>H<sub>16</sub>N<sub>2</sub>O<sub>2</sub>, the five-membered heterocyclic ring makes a dihedral angle of 47.06 (6)° with the attached benzene ring, whereas the indan-1,3-dione ring system and the benzene ring are oriented at a dihedral angle of 21.92 (7)°. In the crystal, inversion dimers linked by pairs of C-H...O hydrogen bonds generate R<sub>2</sub><sup>2</sup>(22) loops. Aromatic  $\pi$ - $\pi$  stacking interactions [centroid-centroid distances = 3.8325 (12)-3.8600 (12) Å] also occur.