

2-Furoic acid, 4-methoxyphenyl ester

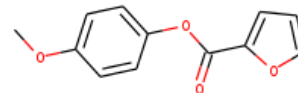
InChI: InChI=1S/C12H10O4/c1-14-9-4-6-10(7-5-9)16-12(13)11-3-2-8-15-11/h2-8H,1H3

InChI Key: PUIKCDWMCVPTGK-UHFFFAOYSA-N

Formula: C12H10O4

SMILES: COc1ccc(OC(=O)c2ccco2)cc1

Molecular Weight: 218.21



Physical Properties

Property	Value	Unit	Source
$\log P_{\text{oct/wat}}$	2.51		Crippen Method

Sources

NIST Webbook: [http://webbook.nist.gov/cgi/inchi/InChI=1S/C12H10O4/c1-14-9-4-6-10\(7-5-9\)16-12\(13\)11-3-2-8-15-11/h2-8H,1H3](http://webbook.nist.gov/cgi/inchi/InChI=1S/C12H10O4/c1-14-9-4-6-10(7-5-9)16-12(13)11-3-2-8-15-11/h2-8H,1H3)

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Legend

$\log P_{\text{oct/wat}}$: Octanol/Water partition coefficient .

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