

2-(p-methylphenoxy)-3-methyl pyrazine

InChI: InChI=1S/C12H12N2O/c1-9-3-5-11(6-4-9)15-12-10(2)13-7-8-14-12/h3-8H,1-2H3

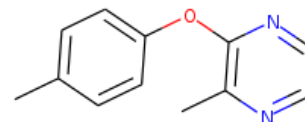
InChI Key: VFXFVQJLZMZESR-UHFFFAOYSA-N

Formula: C12H12N2O

SMILES: Cc1ccc(Oc2nccnc2C)cc1

Molecular Weight: 200.24

CAS: 91392-06-2



Physical Properties

| Property | Value | Unit | Source |
|---------------------------|-------|------|----------------|
| $\log P_{\text{oct/wat}}$ | 2.89 | | Crippen Method |

Sources

NIST Webbook: [http://webbook.nist.gov/cgi/inchi/InChI=1S/C12H12N2O/c1-9-3-5-11\(6-4-9\)15-12-10\(2\)13-7-8-14-12/h3-8H,1-2H3](http://webbook.nist.gov/cgi/inchi/InChI=1S/C12H12N2O/c1-9-3-5-11(6-4-9)15-12-10(2)13-7-8-14-12/h3-8H,1-2H3)

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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