2-Phenoxy-3,6-dimethyl pyrazine

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

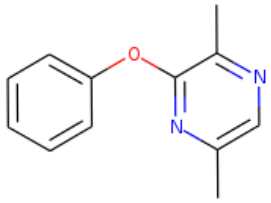
InChI Key: XHMYGFGJEPCOIE-UHFFFAOYSA-N

Formula: C12H12N2O

SMILES: Cc1cnc(C)c(Oc2ccccc2)n1

Molecular Weight: 200.24

CAS: 116660-24-3



Physical Properties

Property	Value	Unit	Source
$logP_{oct/wat}$	2.89		Crippen Method

Sources

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

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

***log*P$_{oct/wat}$**: Octanol/Water partition coefficient .