

# 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

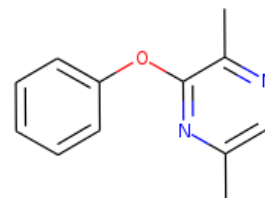
**InChI Key:** XHMYGFGJEP COIE-UHFFFAOYSA-N

**Formula:** C<sub>12</sub>H<sub>12</sub>N<sub>2</sub>O

**SMILES:** Cc1cnc(C)c(Oc2ccccc2)n1

**Molecular Weight:** 200.24

**CAS:** 116660-24-3



## 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/c1-9-8-13-10\(2\)12\(14-9\)15-11-6-4-3-5-7-11/h3-8H,1-2H3](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/ci9903071>

## Legend

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

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