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

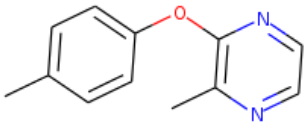
InChI Key: VFXFVQJLZMZESR-UHFFFAOYSA-N

Formula: C12H12N2O

SMILES: Cc1ccc(Oc2nccnc2C)cc1

Molecular Weight: 200.24

CAS: 91392-06-2



Physical Properties

Property	Value	Unit	Source
*log*P$_{oct/wat}$	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

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

Legend

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