(2-Phenylethyl)pyrazine

Other names: (2-Phenylethyl)pyrazine; Phenylethyl-pyrazine; Pyrazine, phenethyl-.

InChI: InChI=1S/C12H12N2/c1-2-4-11(5-3-1)6-7-12-10-13-8-9-14-12/h1-5,8-10H,6-7H2

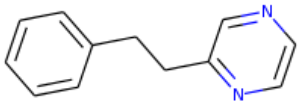
InChI Key: LQWPAMNCTNMMDB-UHFFFAOYSA-N

Formula: C12H12N2

SMILES: c1ccc(CCc2cnccn2)cc1

Molecular Weight: 184.24

CAS: 91391-83-2



Physical Properties

Property	Value	Unit	Source
*log*P$_{oct/wat}$	2.26		Crippen Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/inchi/InChI=1S/C12H12N2/c1-2-4-11(5-3-1)6-7-12-10-13-8-9-14-12/h1-5,8-10H,6-7H2

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

Legend

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