1H,5H-Benzo[ij]quinolizin-8-ol, 2,3,6,7-tetrahydro-

Other names: 2,3,6,7-Tetrahydro-1H,5H-benzo(ij)quinolizin-8-ol; 2,3,6,7-Tetrahydro-1H,5H-pyrido[3,2,1-ij]quinolin-8-ol; 8-Hydroxyjulolidine.

InChI: InChI=1S/C12H15NO/c14-11-6-5-9-3-1-7-13-8-2-4-10(11)12(9)13/h5-6,14H,1-4,7-8H2

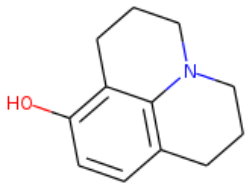
InChI Key: FOFUWJNBAQJABO-UHFFFAOYSA-N

Formula: C12H15NO

SMILES: Oc1ccc2c3c1CCCN3CCC2

Molecular Weight: 189.25

CAS: 41175-50-2



Physical Properties

Property	Value	Unit	Source
$logP_{oct/wat}$	2.09		Crippen Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/inchi/InChI=1S/C12H15NO/c14-11-6-5-9-3-1-7-13-8-2-4-10(11)12(9)13/h5-6,14H,1-4,7-8H2

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

Legend

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