11H-Pyrido[2,1-b]quinazolin-11-one, 6,7,8,9-tetrahydro, 8-methyl

Inchi: InChI=1S/C13H14N2O/c1-9-6-7-12-14-11-5-3-2-4-10(11)13(16)15(12)8-9/h2-5,9H,6-8H2

InchiKey: JEGAIHUNMXTWSA-UHFFFAOYSA-N

Formula: C13H14N2O

SMILES: CC1CCc2nc3ccccc3c(=O)n2C1

Mol. weight [g/mol]: 214.26

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.60		Crippen Method
logp	1.979		Crippen Method
mcvol	165.780	ml/mol	McGowan Method
rinpol	2190.00		NIST Webbook
rinpol	2190.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R64176&Units=SI

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

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

Legend

log10ws: Log10 of Water solubility in mol/l

logp: Octanol/Water partition coefficient

mcvol: McGowan's characteristic volume

rinpol: Non-polar retention indices

Cheméo (https://www.chemeo.com) is the biggest free database of chemical and physical data for the process industry.