

N-Formylalbine

Inchi: InChI=1S/C15H20N2O2/c1-2-3-14-13-6-11(9-17(14)10-18)8-16-5-4-12(19)7-15(13)16/h2
InchiKey: WASGYVQSAZPDIT-AXVKAWJUSA-N
Formula: C15H20N2O2
SMILES: C=CCC1C2CC(CN1C=O)CN1C=CC(=O)CC21
Mol. weight [g/mol]: 260.33

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.92		Crippen Method
logp	1.196		Crippen Method
mcvol	204.130	ml/mol	McGowan Method
rinpol	2543.00		NIST Webbook
rinpol	2530.00		NIST Webbook
rinpol	2543.00		NIST Webbook

Sources

McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R319969&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
rinpol: Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/18-779-8/N-Formylalbine.pdf>

Generated by Cheméo on 2025-12-22 13:11:18.178310811 +0000 UTC m=+6157275.708351477.

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