

13«alpha»-Propanyloxylupanine

Inchi: InChI=1S/C18H28N2O3/c1-2-18(22)23-14-6-7-15-13-8-12(9-19(15)11-14)16-4-3-5-17(21)
InchiKey: DVXPHLIHSQFWMH-GMOYYMRRSA-N
Formula: C18H28N2O3
SMILES: CCC(=O)OC1CCC2C3CC(CN2C1)C1CCCC(=O)N1C3
Mol. weight [g/mol]: 320.43

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.57		Crippen Method
logp	1.803		Crippen Method
mcvol	250.010	ml/mol	McGowan Method
rinpol	2530.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R263851&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

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