

Desacetyldoronine

Inchi: InChI=1S/C18H28ClNO6/c1-11-9-18(24,12(2)19)16(22)26-14-6-8-20(4)7-5-13(15(14)21)
InchiKey: QAISIHLCOPYMAEI-ZVPVIJKSA-N
Formula: C18H28ClNO6
SMILES: CC1CC(O)(C(C)Cl)C(=O)OC2CCN(C)CC=C(OCC1(C)O)C2=O
Mol. weight [g/mol]: 389.87

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.18		Crippen Method
logp	0.853		Crippen Method
mcvol	287.300	ml/mol	McGowan Method
rinpol	2737.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R422703&Units=SI>

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/37-012-7/Desacetyldoronine.pdf>

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