

Acetylseneciphylline

Inchi: InChI=1S/C20H25NO6/c1-5-14-10-12(2)20(4,27-13(3)22)19(24)25-11-15-6-8-21-9-7-16(18)
InchiKey: CTCKXBIRQMSUIU-BYVYPXNKSA-N
Formula: C20H25NO6
SMILES: C=C1CC(=CC)C(=O)OC2CCN3CC=C(COC(=O)C1(C)OC(C)=O)C23
Mol. weight [g/mol]: 375.42

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.93		Crippen Method
logp	1.684		Crippen Method
mcvol	279.480	ml/mol	McGowan Method
rinpol	2483.00		NIST Webbook
rinpol	2460.00		NIST Webbook
rinpol	2483.00		NIST Webbook
rinpol	2460.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R178362&Units=SI>
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>

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.cheméo.com/cid/44-940-9/Acetylseneciphylline.pdf>

Generated by Cheméo on 2025-12-20 05:49:47.778263595 +0000 UTC m=+5957985.308304262.

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