

Benzquinamide M (N-des-Et), acetylated

Inchi: InChI=1S/C22H30N2O6/c1-6-24(13(2)25)22(27)17-12-23-8-7-15-9-20(28-4)21(29-5)10-1
InchiKey: KYXZHWVPMPALKB-UHFFFAOYSA-N
Formula: C22H30N2O6
SMILES: CCN(C(C)=O)C(=O)C1CN2CCc3cc(OC)c(OC)cc3C2CC1OC(C)=O
Mol. weight [g/mol]: 418.48

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.17		Crippen Method
logp	1.950		Crippen Method
mcvol	317.640	ml/mol	McGowan Method
rinpol	2960.00		NIST Webbook
rinpol	2960.00		NIST Webbook

Sources

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=R120497&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307I>

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/58-102-4/Benzquinamide-M-N-des-Et-acetylated.pdf>

Generated by Cheméo on 2025-12-05 14:50:15.630300602 +0000 UTC m=+4694413.160341256.

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