

Isonipectic acid, N-(3-cyclopentylpropionyl)-, tridecyl ester

Inchi: InChI=1S/C27H49NO3/c1-2-3-4-5-6-7-8-9-10-11-14-23-31-27(30)25-19-21-28(22-20-25)
InchiKey: QQUOTIOEANKACW-UHFFFAOYSA-N
Formula: C27H49NO3
SMILES: CCCCCCCCCCCCCOC(=O)C1CCN(C(=O)CCC2CCCC2)CC1
Mol. weight [g/mol]: 435.68

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|--------|----------------|
| log10ws | -7.64 | | Crippen Method |
| logp | 7.050 | | Crippen Method |
| mcvol | 388.560 | ml/mol | McGowan Method |
| rinpol | 3539.00 | | NIST Webbook |
| rinpol | 3539.00 | | NIST Webbook |

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U361093&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.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/20-795-7/Isonipectic-acid-N-3-cyclopentylpropionyl-tridecyl-ester.pdf>

Generated by Cheméo on 2024-05-22 14:06:55.097241484 +0000 UTC m=+18676064.017818797.

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