

L-Proline, N-(3-cyclopentylpropionyl)-, undecyl ester

Inchi: InChI=1S/C24H43NO3/c1-2-3-4-5-6-7-8-9-12-20-28-24(27)22-16-13-19-25(22)23(26)18-
InchiKey: SWNPLDKJETUMOW-UHFFFAOYSA-N
Formula: C24H43NO3
SMILES: CCCCCCCCCCOC(=O)C1CCCN1C(=O)CCC1CCCC1
Mol. weight [g/mol]: 393.60

Physical Properties

Property code	Value	Unit	Source
log10ws	-6.74		Crippen Method
logp	6.022		Crippen Method
mcvol	346.290	ml/mol	McGowan Method
rinpol	3035.00		NIST Webbook
rinpol	3035.00		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=U346416&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

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<https://www.chemeo.com/cid/84-116-0/L-Proline-N-3-cyclopentylpropionyl-undecyl-ester.pdf>

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