

L-Proline, N-(hexanoyl)-, tetradecyl ester

Inchi: InChI=1S/C25H47NO3/c1-3-5-7-8-9-10-11-12-13-14-15-17-22-29-25(28)23-19-18-21-26
InchiKey: YNWVGLYLOGMERE-UHFFFAOYSA-N
Formula: C25H47NO3
SMILES: CCCCCCCCCCCCCOC(=O)C1CCCN1C(=O)CCCCC
Mol. weight [g/mol]: 409.65

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.50		Crippen Method
logp	6.802		Crippen Method
mcvol	371.240	ml/mol	McGowan Method
rinpol	3050.00		NIST Webbook
rinpol	3050.00		NIST Webbook

Sources

McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U346163&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws

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/95-452-5/L-Proline-N-hexanoyl-tetradecyl-ester.pdf>

Generated by Cheméo on 2025-05-21 21:45:55.749453539 +0000 UTC m=+3173001.249897765.

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