

D-Proline, N-methoxycarbonyl-, nonyl ester

Inchi: InChI=1S/C16H29NO4/c1-3-4-5-6-7-8-9-13-21-15(18)14-11-10-12-17(14)16(19)20-2/h14
InchiKey: VBYSRQOYCHVUDBK-UHFFFAOYSA-N
Formula: C16H29NO4
SMILES: CCCCCCCCCOC(=O)C1CCCN1C(=O)OC
Mol. weight [g/mol]: 299.41

Physical Properties

Property code	Value	Unit	Source
logp	3.511		Crippen Method
mcvol	250.300	ml/mol	McGowan Method
tf	308.25	K	Cheméo Relay (1.0)

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U320793&Units=SI>
McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>

Legend

logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
tf: Normal melting (fusion) point

Latest version available from:

<https://www.chemeo.com/cid/63-994-9/D-Proline-N-methoxycarbonyl-nonyl-ester.pdf>

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