

D-Proline, N-butoxycarbonyl-, hexyl ester

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

Physical Properties

Property code	Value	Unit	Source
hvap	91.49	kJ/mol	Cheméo Relay (1.0)
log10ws	-3.37		Cheméo Relay (1.0)
logp	3.511		Crippen Method
mcvol	250.300	ml/mol	McGowan Method
rinpol	1979.00		NIST Webbook
rinpol	1979.00		NIST Webbook
tb	607.13	K	Cheméo Relay (1.0)
tf	271.68		Cheméo Relay (1.0)

Sources

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=U321084&Units=SI
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient

mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tf:	Normal melting (fusion) point

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