

1-Aminocyclopentanecarboxylic acid, N-(propargyloxycarbonyl)-, tridecyl ester

Inchi:	lnChI=1S/C23H39NO4/c1-3-5-6-7-8-9-10-11-12-13-16-20-27-21(25)23(17-14-15-18-23)2
InchiKey:	IQAKGOGTFPNXGU-UHFFFAOYSA-N
Formula:	C23H39NO4
SMILES:	C#CCOC(O)=NC1(C(=O)OCCCCCCCCCCC)CCCC1
Mol. weight [g/mol]:	393.56

Physical Properties

Property code	Value	Unit	Source
hf	-607.25	kJ/mol	Joback Method
hvap	97.39	kJ/mol	Joback Method
log10ws	-6.68		Crippen Method
logp	5.717		Crippen Method
mcvol	340.330	ml/mol	McGowan Method
pc	1087.78	kPa	Joback Method
tb	998.73	K	Joback Method
tc	1222.85	K	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U392471&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307i
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

hf:	Enthalpy of formation at standard conditions
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
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

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