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

Inchi: InChI=1S/C23H41NO4/c1-3-5-6-7-8-9-10-11-12-13-16-20-27-21(25)23(17-14-15-18-23)2

InchiKey: DOZYWTDWARGJOC-UHFFFAOYSA-N

Formula: C23H41NO4

SMILES: C=CCOC(O)=NC1(C(=O)OCCCCCCCCCCCCC)CCCC1

Mol. weight [g/mol]: 395.58

Physical Properties

Property code	Value	Unit	Source
hf	-773.72	kJ/mol	Joback Method
hvap	96.87	kJ/mol	Joback Method
log10ws	-6.74		Crippen Method
logp	6.270		Crippen Method
mcvol	344.630	ml/mol	McGowan Method
pc	1015.53	kPa	Joback Method
rinpol	2767.00		NIST Webbook
tb	1005.29	K	Joback Method
tc	1231.99	K	Joback Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U392482&Units=SI

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Joback Method: https://en.wikipedia.org/wiki/Joback_method

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature