

DL-Alanine, N-methyl-N-(2-ethylhexyloxycarbonyl)-, tetradecyl ester

InChI: C27H53NO4
InChIKey: LMBZGTWDZUFXTE-UHFFFAOYSA-N

Formula: C27H53NO4

SMILES: CCCCCCCCCCCCCOC(=O)C(C)N(C)C(=O)OCC(CC)CCCC

Mol. weight [g/mol]: 455.71

Physical Properties

Property code	Value	Unit	Source
gf	-185.48	kJ/mol	Joback Method
hf	-1033.24	kJ/mol	Joback Method
hfus	67.24	kJ/mol	Joback Method
hvap	95.28	kJ/mol	Joback Method
log10ws	-8.27		Crippen Method
logp	7.904		Crippen Method
mvol	416.150	ml/mol	McGowan Method
pc	728.88	kPa	Joback Method
rinpol	2870.00		NIST Webbook
rinpol	2870.00		NIST Webbook
tb	981.30	K	Joback Method
tc	1212.75	K	Joback Method
tf	540.84	K	Joback Method
vc	1.601	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1455.55	J/molxK	981.30	Joback Method
cpg	1477.49	J/molxK	1019.88	Joback Method
cpg	1497.54	J/molxK	1058.45	Joback Method
cpg	1515.78	J/molxK	1097.03	Joback Method
cpg	1532.28	J/molxK	1135.60	Joback Method
cpg	1547.11	J/molxK	1174.18	Joback Method
cpg	1560.35	J/molxK	1212.75	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U392665&Units=SI

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion 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
rinpola:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/94-933-2/DL-Alanine-N-methyl-N-2-ethylhexyloxycarbonyl-tetradecyl-ester.pdf>

Generated by Cheméo on 2024-04-23 16:15:51.736580981 +0000 UTC m=+16178200.657158292.

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