

L-Norvaline, N-hexyloxycarbonyl-, butyl ester

Inchi:	InChI=1S/C16H31NO4/c1-4-7-9-10-13-21-16(19)17-14(11-6-3)15(18)20-12-8-5-2/h14H,4
InchiKey:	GTKAYWRBSUKNLE-CQSZACIVSA-N
Formula:	C16H31NO4
SMILES:	CCCCCOC(=O)NC(CCC)C(=O)OCCCC
Mol. weight [g/mol]:	301.42

Physical Properties

Property code	Value	Unit	Source
gf	-297.05	kJ/mol	Joback Method
hf	-814.98	kJ/mol	Joback Method
hfus	44.35	kJ/mol	Joback Method
hvap	75.57	kJ/mol	Joback Method
log10ws	-4.52		Crippen Method
logp	3.805		Crippen Method
mvol	261.160	ml/mol	McGowan Method
pc	1433.72	kPa	Joback Method
rinpol	2023.00		NIST Webbook
rinpol	2023.00		NIST Webbook
tb	767.79	K	Joback Method
tc	950.65	K	Joback Method
tf	452.06	K	Joback Method
vc	1.008	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	795.32	J/molxK	767.79	Joback Method
cpg	811.65	J/molxK	798.27	Joback Method
cpg	827.08	J/molxK	828.74	Joback Method
cpg	841.60	J/molxK	859.22	Joback Method
cpg	855.22	J/molxK	889.70	Joback Method
cpg	867.97	J/molxK	920.17	Joback Method
cpg	879.84	J/molxK	950.65	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U392832&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
h vap:	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
r in pol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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