

I-Norvaline, N-allyloxycarbonyl-, undecyl ester

Inchi:	InChI=1S/C20H37NO4/c1-4-7-8-9-10-11-12-13-14-17-24-19(22)18(15-5-2)21-20(23)25-1
InchiKey:	NATQURDTERXJCY-UHFFFAOYSA-N
Formula:	C20H37NO4
SMILES:	C=CCOC(=O)NC(CCC)C(=O)OCCCCCCCCCCCC
Mol. weight [g/mol]:	355.51

Physical Properties

Property code	Value	Unit	Source
gf	-175.53	kJ/mol	Joback Method
hf	-772.11	kJ/mol	Joback Method
hfus	53.43	kJ/mol	Joback Method
hvap	83.80	kJ/mol	Joback Method
log10ws	-6.05		Crippen Method
logp	5.141		Crippen Method
mcvol	313.220	ml/mol	McGowan Method
pc	1123.81	kPa	Joback Method
rinpol	2162.00		NIST Webbook
rinpol	2162.00		NIST Webbook
tb	855.99	K	Joback Method
tc	1049.04	K	Joback Method
tf	495.38	K	Joback Method
vc	1.214	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1006.16	J/mol×K	855.99	Joback Method
cpg	1023.48	J/mol×K	888.17	Joback Method
cpg	1039.68	J/mol×K	920.34	Joback Method
cpg	1054.77	J/mol×K	952.52	Joback Method
cpg	1068.80	J/mol×K	984.69	Joback Method
cpg	1081.78	J/mol×K	1016.87	Joback Method
cpg	1093.73	J/mol×K	1049.04	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U320750&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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
rlnpol:	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/29-653-5/l-Norvaline-N-allyloxycarbonyl-undecyl-ester.pdf>

Generated by Cheméo on 2025-12-23 06:11:39.706486473 +0000 UTC m=+6218497.236527127.

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