

I-Norvaline, N-isobutoxycarbonyl-, octyl ester

Inchi:	InChI=1S/C18H35NO4/c1-5-7-8-9-10-11-13-22-17(20)16(12-6-2)19-18(21)23-14-15(3)4/
InchiKey:	PNJMFOCHCUQIEG-UHFFFAOYSA-N
Formula:	C18H35NO4
SMILES:	CCCCCCCCCOC(=O)C(CCC)N=C(O)OCC(C)C
Mol. weight [g/mol]:	329.47

Physical Properties

Property code	Value	Unit	Source
hf	-882.23	kJ/mol	Joback Method
hvap	86.52	kJ/mol	Joback Method
log10ws	-4.66		Crippen Method
logp	4.645		Crippen Method
mcvol	289.340	ml/mol	McGowan Method
pc	1181.71	kPa	Joback Method
rinpol	1946.00		NIST Webbook
rinpol	1946.00		NIST Webbook
tb	877.81	K	Joback Method
tc	1075.29	K	Joback Method

Sources

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=U320718&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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
mvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

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