«beta»-Alanine, N-caproyl-, decyl ester

Inchi: InChI=1S/C19H37NO3/c1-3-5-7-8-9-10-11-13-17-23-19(22)15-16-20-18(21)14-12-6-4-2/h
InchiKey: OBYFDWIWPUJARY-UHFFFAOYSA-N
Formula: C19H37NO3
SMILES: CCCCCCCCCCOC(=O)CCN=C(O)CCCCC
Mol. weight [g/mol]: 327.50

Physical Properties

Property code	Value	Unit	Source
hf	-760.09	kJ/mol	Joback Method
hvap	87.12	kJ/mol	Joback Method
log10ws	-5.62		Crippen Method
logp	5.597		Crippen Method
mcvol	297.560	ml/mol	McGowan Method
pc	1108.15	kPa	Joback Method
rinpol	2478.00		NIST Webbook
rinpol	2478.00		NIST Webbook
tb	879.15	K	Joback Method
tc	1076.36	K	Joback Method

Sources

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=U321787&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

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