

Glycine, 2-cyclohexyl-N-(2-ethylhexyl)oxycarbonyl-, hexyl ester

InChI: InChI=1S/C23H43NO4/c1-4-7-9-13-17-27-22(25)21(20-15-11-10-12-16-20)24-23(26)28-
InChIKey: SVTOHDHCKCCVMY-UHFFFAOYSA-N
Formula: C23H43NO4
SMILES: CCCCCCOC(=O)C(N=C(O)OCC(CC)CCCC)C1CCCCC1
Mol. weight [g/mol]: 397.59

Physical Properties

Property code	Value	Unit	Source
hf	-931.11	kJ/mol	Joback Method
hvap	98.08	kJ/mol	Joback Method
log10ws	-6.40		Crippen Method
logp	6.206		Crippen Method
mcpvol	348.930	ml/mol	McGowan Method
pc	986.40	kPa	Joback Method
rinpol	2597.00		NIST Webbook
tb	1011.76	K	Joback Method
tc	1239.76	K	Joback Method

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
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=U383150&Units=SI>

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

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