

Propanamide, 3-cyclopentyl-N-dodecyl-

Inchi: InChI=1S/C20H39NO/c1-2-3-4-5-6-7-8-9-10-13-18-21-20(22)17-16-19-14-11-12-15-19/h
InchiKey: WGIDJZTUXIVYTP-UHFFFAOYSA-N
Formula: C20H39NO
SMILES: CCCCCCCCCCCCN=C(O)CCC1CCCC1
Mol. weight [g/mol]: 309.53

Physical Properties

Property code	Value	Unit	Source
hf	-475.45	kJ/mol	Joback Method
hvap	80.44	kJ/mol	Joback Method
log10ws	-6.83		Crippen Method
logp	6.834		Crippen Method
mcvol	293.350	ml/mol	McGowan Method
pc	1136.73	kPa	Joback Method
rinpol	2558.00		NIST Webbook
rinpol	2558.00		NIST Webbook
tb	841.02	K	Joback Method
tc	1034.08	K	Joback Method

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
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=U407386&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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