

Octadecanamide, n-methyl-

Inchi: InChI=1S/C19H39NO/c1-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17-18-19(21)20-2/h3-18H2
InchiKey: HNUFCQUTJXHEPI-UHFFFAOYSA-N
Formula: C19H39NO
SMILES: CCCCCCCCCCCCCCCCC(O)=NC
Mol. weight [g/mol]: 297.52
CAS: 20198-92-9

Physical Properties

Property code	Value	Unit	Source
hf	-515.29	kJ/mol	Joback Method
hvap	77.96	kJ/mol	Joback Method
log10ws	-6.76		Crippen Method
logp	6.834		Crippen Method
mcvol	290.120	ml/mol	McGowan Method
pc	1074.98	kPa	Joback Method
tb	802.86	K	Joback Method
tc	984.91	K	Joback Method

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C20198929&Units=SI>
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>

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/89-303-7/Octadecanamide-n-methyl.pdf>

Generated by Cheméo on 2024-04-29 10:08:40.504915886 +0000 UTC m=+16674569.425493198.

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