

Myristamide, N-(3-methylbutyl)-

Inchi:	InChI=1S/C19H39NO/c1-4-5-6-7-8-9-10-11-12-13-14-15-19(21)20-17-16-18(2)3/h18H,4-
InchiKey:	IYZXUMQBCKGGBU-UHFFFAOYSA-N
Formula:	C19H39NO
SMILES:	CCCCCCCCCCCC(O)=NCCC(C)C
Mol. weight [g/mol]:	297.52

Physical Properties

Property code	Value	Unit	Source
hf	-520.57	kJ/mol	Joback Method
hvap	77.57	kJ/mol	Joback Method
log10ws	-6.52		Crippen Method
logp	6.690		Crippen Method
mcvol	290.120	ml/mol	McGowan Method
pc	1080.64	kPa	Joback Method
rinpol	2327.00		NIST Webbook
rinpol	2327.00		NIST Webbook
tb	802.42	K	Joback Method
tc	985.15	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=U408022&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

Latest version available from:

<https://www.cheméo.com/cid/89-300-0/Myristamide-N-3-methylbutyl.pdf>

Generated by Cheméo on 2024-04-30 12:29:05.469567077 +0000 UTC m=+16769394.390144389.

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