

1-(2-(Hexadecyloxy)ethoxy)octadecane

Inchi:	InChI=1S/C36H74O2/c1-3-5-7-9-11-13-15-17-19-20-22-24-26-28-30-32-34-38-36-35-37-
InchiKey:	RSHNRWZRUIITEEY-UHFFFAOYSA-N
Formula:	C36H74O2
SMILES:	CCCCCCCCCCCCCCCCCCOCCOCCCCCCCCCCCCCCCC
Mol. weight [g/mol]:	538.97

Physical Properties

Property code	Value	Unit	Source
gf	42.24	kJ/mol	Joback Method
hf	-1050.81	kJ/mol	Joback Method
hfus	91.37	kJ/mol	Joback Method
hvap	100.55	kJ/mol	Joback Method
log10ws	-13.07		Crippen Method
logp	12.762		Crippen Method
mvol	529.840	ml/mol	McGowan Method
pc	451.16	kPa	Joback Method
rinpol	412.60		NIST Webbook
rinpol	412.60		NIST Webbook
tb	1067.92	K	Joback Method
tc	1391.31	K	Joback Method
tf	539.94	K	Joback Method
vc	2.087	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1964.06	J/molxK	1067.92	Joback Method
cpg	1999.29	J/molxK	1121.82	Joback Method
cpg	2030.59	J/molxK	1175.72	Joback Method
cpg	2058.24	J/molxK	1229.61	Joback Method
cpg	2082.53	J/molxK	1283.51	Joback Method
cpg	2103.76	J/molxK	1337.41	Joback Method
cpg	2122.20	J/molxK	1391.31	Joback Method
dvisc	0.0001872	Paxs	539.94	Joback Method

dvisc	0.0000673	Paxs	627.94	Joback Method
dvisc	0.0000311	Paxs	715.93	Joback Method
dvisc	0.0000170	Paxs	803.93	Joback Method
dvisc	0.0000105	Paxs	891.93	Joback Method
dvisc	0.0000070	Paxs	979.92	Joback Method
dvisc	0.0000051	Paxs	1067.92	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=R516571&Units=SI

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion 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
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/49-794-7/1-2-Hexadecyloxy-ethoxyl-octadecane.pdf>

Generated by Cheméo on 2024-04-17 18:19:43.142503698 +0000 UTC m=+15667232.063081009.

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