

Ethanol, 2-(tetradecyloxy)-

Other names:

Myristyl monoethoxylate
2-Tetradecyloxyethanol
Ethylene glycol monotetradecyl ether

Inchi: InChI=1S/C16H34O2/c1-2-3-4-5-6-7-8-9-10-11-12-13-15-18-16-14-17/h17H,2-16H2,1H3**InchiKey:** QHERURXKGFABNZ-UHFFFAOYSA-N**Formula:** C16H34O2**SMILES:** CCCCCCCCCCCCCOCCO**Mol. weight [g/mol]:** 258.44**CAS:** 2136-70-1

Physical Properties

Property code	Value	Unit	Source
gf	-157.98	kJ/mol	Joback Method
hf	-658.02	kJ/mol	Joback Method
hfus	42.47	kJ/mol	Joback Method
hvap	70.30	kJ/mol	Joback Method
log10ws	-4.87		Crippen Method
logp	4.696		Crippen Method
mcvol	248.040	ml/mol	McGowan Method
pc	1393.33	kPa	Joback Method
tb	680.08	K	Joback Method
tc	841.71	K	Joback Method
tf	353.13	K	Joback Method
vc	0.969	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	723.53	J/molxK	680.08	Joback Method
cpg	800.99	J/molxK	814.78	Joback Method
cpg	786.83	J/molxK	787.84	Joback Method
cpg	772.02	J/molxK	760.90	Joback Method
cpg	756.54	J/molxK	733.96	Joback Method
cpg	740.38	J/molxK	707.02	Joback Method

cpg	814.51	J/molxK	841.71	Joback Method
dvisc	0.0000300	Paxs	680.08	Joback Method
dvisc	0.0000479	Paxs	625.59	Joback Method
dvisc	0.0000837	Paxs	571.10	Joback Method
dvisc	0.0001646	Paxs	516.61	Joback Method
dvisc	0.0003794	Paxs	462.11	Joback Method
dvisc	0.0010932	Paxs	407.62	Joback Method
dvisc	0.0043674	Paxs	353.13	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=C2136701&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
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/61-905-9/Ethanol-2-tetradecyloxy.pdf>

Generated by Cheméo on 2024-04-30 01:39:00.146335424 +0000 UTC m=+16730389.066912737.

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