

Diethylene glycol, decyl ether

Other names:	2-(2-decyloxy-ethoxy)-ethanol
Inchi:	InChI=1S/C14H30O3/c1-2-3-4-5-6-7-8-9-11-16-13-14-17-12-10-15/h15H,2-14H2,1H3
InchiKey:	UVNSFSOKRSZVEZ-UHFFFAOYSA-N
Formula:	C14H30O3
SMILES:	CCCCCCCCCOCCOCCO
Mol. weight [g/mol]:	246.39

Physical Properties

Property code	Value	Unit	Source
gf	-279.82	kJ/mol	Joback Method
hf	-748.96	kJ/mol	Joback Method
hfus	38.48	kJ/mol	Joback Method
hvap	68.26	kJ/mol	Joback Method
log10ws	-3.12		Crippen Method
logp	3.153		Crippen Method
mvol	225.730	ml/mol	McGowan Method
pc	1594.89	kPa	Joback Method
rinpol	1864.10		NIST Webbook
tb	656.74	K	Joback Method
tc	817.27	K	Joback Method
tf	352.82	K	Joback Method
vc	0.875	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	641.91	J/molxK	656.74	Joback Method
cpg	657.54	J/molxK	683.49	Joback Method
cpg	672.56	J/molxK	710.25	Joback Method
cpg	686.96	J/molxK	737.00	Joback Method
cpg	700.76	J/molxK	763.76	Joback Method
cpg	713.95	J/molxK	790.51	Joback Method
cpg	726.55	J/molxK	817.27	Joback Method
dvisc	0.0035246	Paxs	352.82	Joback Method

dvisc	0.0009817	Paxs	403.47	Joback Method
dvisc	0.0003636	Paxs	454.13	Joback Method
dvisc	0.0001644	Paxs	504.78	Joback Method
dvisc	0.0000859	Paxs	555.43	Joback Method
dvisc	0.0000500	Paxs	606.09	Joback Method
dvisc	0.0000317	Paxs	656.74	Joback Method

Sources

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=R184362&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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/68-356-2/Diethylene-glycol-decyl-ether.pdf>

Generated by Cheméo on 2024-04-27 04:14:31.311840405 +0000 UTC m=+16480520.232417754.

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