Ethane, 1,2-dichloro-1-ethoxy-

Other names:	Ether, 1,2-dichloroethyl ethyl
	1,2-Dichloroethyl ethyl ether
	1,2-Dichlorodiethyl ether
	Ethyl 1,2-dichloroethyl ether
	1,2-dichloro-2-ethoxyethane
Inchi:	InChI=1S/C4H8Cl2O/c1-2-7-4(6)3-5/h4H,2-3H2,1H3
InchiKey:	NNBUKAPOVBEMNI-UHFFFAOYSA-N
Formula:	C4H8Cl2O
SMILES:	CCOC(Cl)CCl
Mol. weight [g/mol]:	143.01
CAS:	623-46-1

Physical Properties

Property code	Value	Unit	Source
gf	-148.50	kJ/mol	Joback Method
hf	-294.87	kJ/mol	Joback Method
hfus	12.18	kJ/mol	Joback Method
hvap	35.29	kJ/mol	Joback Method
log10ws	-1.50		Crippen Method
logp	1.827		Crippen Method
mcvol	97.570	ml/mol	McGowan Method
pc	3522.09	kPa	Joback Method
rinpol	865.00		NIST Webbook
rinpol	856.00		NIST Webbook
rinpol	856.00		NIST Webbook
tb	414.70 ± 0.60	K	NIST Webbook
tb	418.20	K	NIST Webbook
tb	415.70	K	NIST Webbook
tc	574.35	K	Joback Method
tf	201.91	K	Joback Method
vc	0.369	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	161.62	J/mol×K	387.76	Joback Method
cpg	169.00	J/mol×K	418.86	Joback Method
cpg	176.14	J/mol×K	449.96	Joback Method
cpg	183.02	J/mol×K	481.05	Joback Method
cpg	189.66	J/mol×K	512.15	Joback Method
cpg	196.06	J/mol×K	543.25	Joback Method
cpg	202.20	J/mol×K	574.35	Joback Method
dvisc	0.0050358	Pa×s	201.91	Joback Method
dvisc	0.0023000	Pa×s	232.89	Joback Method
dvisc	0.0012627	Pa×s	263.86	Joback Method
dvisc	0.0007863	Pa×s	294.84	Joback Method
dvisc	0.0005358	Pa×s	325.81	Joback Method
dvisc	0.0003902	Pa×s	356.78	Joback Method
dvisc	0.0002990	Pa×s	387.76	Joback Method

Sources

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Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
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

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