

2-Chloroethyl ethyl carbonate

Inchi:	InChI=1S/C5H9ClO3/c1-2-8-5(7)9-4-3-6/h2-4H2,1H3
InchiKey:	SEYHRCMNMPHDG-UHFFFAOYSA-N
Formula:	C5H9ClO3
SMILES:	CCOC(=O)OCCCl
Mol. weight [g/mol]:	152.58

Physical Properties

Property code	Value	Unit	Source
gf	-359.63	kJ/mol	Joback Method
hf	-539.29	kJ/mol	Joback Method
hfus	16.88	kJ/mol	Joback Method
hvap	42.67	kJ/mol	Joback Method
log10ws	-1.00		Crippen Method
logp	1.398		Crippen Method
mvol	106.860	ml/mol	McGowan Method
pc	3435.91	kPa	Joback Method
rinpol	1004.00		NIST Webbook
rinpol	1004.00		NIST Webbook
tb	449.94	K	Joback Method
tc	635.36	K	Joback Method
tf	270.42	K	Joback Method
vc	0.406	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	207.39	J/molxK	449.94	Joback Method
cpg	215.59	J/molxK	480.84	Joback Method
cpg	223.57	J/molxK	511.75	Joback Method
cpg	231.31	J/molxK	542.65	Joback Method
cpg	238.81	J/molxK	573.55	Joback Method
cpg	246.06	J/molxK	604.45	Joback Method
cpg	253.04	J/molxK	635.36	Joback Method
dvisc	0.0022644	Paxs	270.42	Joback Method

dvisc	0.0013308	Paxs	300.34	Joback Method
dvisc	0.0008611	Paxs	330.26	Joback Method
dvisc	0.0005990	Paxs	360.18	Joback Method
dvisc	0.0004406	Paxs	390.10	Joback Method
dvisc	0.0003385	Paxs	420.02	Joback Method
dvisc	0.0002694	Paxs	449.94	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U373774&Units=SI
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

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/54-359-4/2-Chloroethyl-ethyl-carbonate.pdf>

Generated by Cheméo on 2022-10-05 12:28:19.492234185 +0000 UTC m=+262606.427092020.

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