

Sebacic acid, di(2,2-dichloroethyl) ester

Inchi:	InChI=1S/C14H22Cl4O4/c15-11(16)9-21-13(19)7-5-3-1-2-4-6-8-14(20)22-10-12(17)18/h
InchiKey:	UUYXFUXOPDQHKU-UHFFFAOYSA-N
Formula:	C14H22Cl4O4
SMILES:	O=C(CCCCCCCC(=O)OCC(Cl)Cl)OCC(Cl)Cl
Mol. weight [g/mol]:	396.13

Physical Properties

Property code	Value	Unit	Source
gf	-453.44	kJ/mol	Joback Method
hf	-895.41	kJ/mol	Joback Method
hfus	47.33	kJ/mol	Joback Method
hvap	81.83	kJ/mol	Joback Method
log10ws	-5.23		Crippen Method
logp	4.801		Crippen Method
mcvol	271.960	ml/mol	McGowan Method
pc	1477.02	kPa	Joback Method
rinpol	2516.00		NIST Webbook
rinpol	2516.00		NIST Webbook
tb	821.14	K	Joback Method
tc	1020.17	K	Joback Method
tf	481.54	K	Joback Method
vc	1.052	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	734.38	J/molxK	821.14	Joback Method
cpg	787.54	J/molxK	987.00	Joback Method
cpg	778.64	J/molxK	953.83	Joback Method
cpg	768.89	J/molxK	920.66	Joback Method
cpg	758.27	J/molxK	887.48	Joback Method
cpg	746.77	J/molxK	854.31	Joback Method
cpg	795.60	J/molxK	1020.17	Joback Method
dvisc	0.0000579	Paxs	821.14	Joback Method

dvisc	0.0000767	Paxs	764.54	Joback Method
dvisc	0.0001064	Paxs	707.94	Joback Method
dvisc	0.0001562	Paxs	651.34	Joback Method
dvisc	0.0002466	Paxs	594.74	Joback Method
dvisc	0.0004286	Paxs	538.14	Joback Method
dvisc	0.0008483	Paxs	481.54	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=U355479&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/63-011-9/Sebacic-acid-di-2-2-dichloroethyl-ester.pdf>

Generated by Cheméo on 2024-04-24 10:18:30.40646443 +0000 UTC m=+16243159.327041741.

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