

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

Inchi:	InChI=1S/C8H10Cl4O4/c9-5(10)3-15-7(13)1-2-8(14)16-4-6(11)12/h5-6H,1-4H2
InchiKey:	APJYUYGGYUZQET-UHFFFAOYSA-N
Formula:	C8H10Cl4O4
SMILES:	O=C(CCC(=O)OCC(Cl)Cl)OCC(Cl)Cl
Mol. weight [g/mol]:	311.98

Physical Properties

Property code	Value	Unit	Source
gf	-503.96	kJ/mol	Joback Method
hf	-771.57	kJ/mol	Joback Method
hfus	31.79	kJ/mol	Joback Method
hvap	68.48	kJ/mol	Joback Method
log10ws	-2.72		Crippen Method
logp	2.460		Crippen Method
mcvol	187.420	ml/mol	McGowan Method
pc	2436.25	kPa	Joback Method
rinpola	1886.00		NIST Webbook
tb	683.86	K	Joback Method
tc	891.26	K	Joback Method
tf	413.92	K	Joback Method
vc	0.716	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	418.79	J/molxK	683.86	Joback Method
cpg	459.43	J/molxK	856.69	Joback Method
cpg	452.54	J/molxK	822.12	Joback Method
cpg	445.04	J/molxK	787.56	Joback Method
cpg	436.91	J/molxK	752.99	Joback Method
cpg	428.15	J/molxK	718.43	Joback Method
cpg	465.68	J/molxK	891.26	Joback Method
dvisc	0.0001414	Paxs	683.86	Joback Method
dvisc	0.0001841	Paxs	638.87	Joback Method

dvisc	0.0002494	Paxs	593.88	Joback Method
dvisc	0.0003552	Paxs	548.89	Joback Method
dvisc	0.0005387	Paxs	503.90	Joback Method
dvisc	0.0008867	Paxs	458.91	Joback Method
dvisc	0.0016264	Paxs	413.92	Joback Method

Sources

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=U349420&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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

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