

Succinic acid, butyl 2,2-dichloroethyl ester

Inchi:

InChI=1S/C10H16Cl2O4/c1-2-3-6-15-9(13)4-5-10(14)16-7-8(11)12/h8H,2-7H2,1H3

InchiKey:

CBJUDFLVVQADHN-UHFFFAOYSA-N

Formula:

C10H16Cl2O4

SMILES:

CCCCOC(=O)CCC(=O)OCC(Cl)Cl

Mol. weight [g/mol]:

271.14

Physical Properties

Property code	Value	Unit	Source
gf	-460.82	kJ/mol	Joback Method
hf	-776.09	kJ/mol	Joback Method
hfus	32.10	kJ/mol	Joback Method
hvap	64.55	kJ/mol	Joback Method
log10ws	-2.65		Crippen Method
logp	2.457		Crippen Method
mcvol	191.120	ml/mol	McGowan Method
pc	2159.31	kPa	Joback Method
rinpol	1698.00		NIST Webbook
rinpol	1698.00		NIST Webbook
tb	655.20	K	Joback Method
tc	847.01	K	Joback Method
tf	391.62	K	Joback Method
vc	0.736	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	470.31	J/molxK	655.20	Joback Method
cpg	482.41	J/molxK	687.17	Joback Method
cpg	493.88	J/molxK	719.14	Joback Method
cpg	504.72	J/molxK	751.11	Joback Method
cpg	514.92	J/molxK	783.08	Joback Method
cpg	524.48	J/molxK	815.04	Joback Method
cpg	533.42	J/molxK	847.01	Joback Method
dvisc	0.0017020	Paxs	391.62	Joback Method

dvisc	0.0009246	Paxs	435.55	Joback Method
dvisc	0.0005617	Paxs	479.48	Joback Method
dvisc	0.0003711	Paxs	523.41	Joback Method
dvisc	0.0002614	Paxs	567.34	Joback Method
dvisc	0.0001936	Paxs	611.27	Joback Method
dvisc	0.0001493	Paxs	655.20	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U349405&Units=SI
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

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/37-322-3/Succinic-acid-butyl-2-2-dichloroethyl-ester.pdf>

Generated by Cheméo on 2024-05-01 04:05:16.117170264 +0000 UTC m=+16825565.037747575.

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