

Succinic acid, 2,4-dichloronaphth-1-yl ethyl ester

Inchi:	InChI=1S/C16H14Cl2O4/c1-2-21-14(19)7-8-15(20)22-16-11-6-4-3-5-10(11)12(17)9-13(18)
InchiKey:	FSJWXEZRUNIYAU-UHFFFAOYSA-N
Formula:	C16H14Cl2O4
SMILES:	CCOC(=O)CCC(=O)Oc1c(Cl)cc(Cl)c2ccccc12
Mol. weight [g/mol]:	341.19

Physical Properties

Property code	Value	Unit	Source
gf	-217.69	kJ/mol	Joback Method
hf	-501.46	kJ/mol	Joback Method
hfus	41.06	kJ/mol	Joback Method
hvap	84.19	kJ/mol	Joback Method
log10ws	-5.50		Crippen Method
logp	4.395		Crippen Method
mvol	232.440	ml/mol	McGowan Method
pc	2056.76	kPa	Joback Method
rinpol	2421.00		NIST Webbook
rinpol	2421.00		NIST Webbook
tb	853.52	K	Joback Method
tc	1082.62	K	Joback Method
tf	570.92	K	Joback Method
vc	0.891	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	620.83	J/molxK	853.52	Joback Method
cpg	666.64	J/molxK	1044.44	Joback Method
cpg	659.23	J/molxK	1006.26	Joback Method
cpg	650.97	J/molxK	968.07	Joback Method
cpg	641.85	J/molxK	929.89	Joback Method
cpg	631.81	J/molxK	891.70	Joback Method
cpg	673.24	J/molxK	1082.62	Joback Method
dvisc	0.0001460	Paxs	853.52	Joback Method

dvisc	0.0001730	Paxs	806.42	Joback Method
dvisc	0.0002093	Paxs	759.32	Joback Method
dvisc	0.0002597	Paxs	712.22	Joback Method
dvisc	0.0003323	Paxs	665.12	Joback Method
dvisc	0.0004414	Paxs	618.02	Joback Method
dvisc	0.0006145	Paxs	570.92	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U389915&Units=SI

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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