

Succinic acid, 3-chlorophenyl 2-chloro-4-methylphenyl ester

Inchi:	InChI=1S/C17H14Cl2O4/c1-11-5-6-15(14(19)9-11)23-17(21)8-7-16(20)22-13-4-2-3-12(18)
InchiKey:	OQRSAGPENRFXTR-UHFFFAOYSA-N
Formula:	C17H14Cl2O4
SMILES:	<chem>Cc1ccc(OC(=O)CCC(=O)Oc2cccc(Cl)c2)c(Cl)c1</chem>
Mol. weight [g/mol]:	353.20

Physical Properties

Property code	Value	Unit	Source
gf	-203.51	kJ/mol	Joback Method
hf	-476.64	kJ/mol	Joback Method
hfus	40.67	kJ/mol	Joback Method
hvap	87.06	kJ/mol	Joback Method
log10ws	-5.59		Crippen Method
logp	4.593		Crippen Method
mcvol	242.230	ml/mol	McGowan Method
pc	2038.23	kPa	Joback Method
rinpol	2664.00		NIST Webbook
rinpol	2664.00		NIST Webbook
tb	884.10	K	Joback Method
tc	1122.62	K	Joback Method
tf	575.91	K	Joback Method
vc	0.917	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	653.59	J/molxK	884.10	Joback Method
cpg	664.42	J/molxK	923.85	Joback Method
cpg	674.03	J/molxK	963.61	Joback Method
cpg	682.45	J/molxK	1003.36	Joback Method
cpg	689.70	J/molxK	1043.11	Joback Method
cpg	695.80	J/molxK	1082.87	Joback Method
cpg	700.78	J/molxK	1122.62	Joback Method
dvisc	0.0003841	Paxs	575.91	Joback Method

dvisc	0.0002542	Paxs	627.28	Joback Method
dvisc	0.0001791	Paxs	678.64	Joback Method
dvisc	0.0001325	Paxs	730.00	Joback Method
dvisc	0.0001020	Paxs	781.37	Joback Method
dvisc	0.0000811	Paxs	832.73	Joback Method
dvisc	0.0000663	Paxs	884.10	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U390221&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/115-765-5/Succinic-acid-3-chlorophenyl-2-chloro-4-methylphenyl-ester.pdf>

Generated by Cheméo on 2024-04-29 09:06:31.328414369 +0000 UTC m=+16670840.248991696.

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