

Succinic acid, 2-chloro-6-fluorophenyl hexyl ester

Inchi:	InChI=1S/C16H20ClFO4/c1-2-3-4-5-11-21-14(19)9-10-15(20)22-16-12(17)7-6-8-13(16)1
InchiKey:	NLVUVUZIVQLEQR-UHFFFAOYSA-N
Formula:	C16H20ClFO4
SMILES:	CCCCCOC(=O)CCC(=O)Oc1c(F)cccc1Cl
Mol. weight [g/mol]:	330.78

Physical Properties

Property code	Value	Unit	Source
gf	-497.59	kJ/mol	Joback Method
hf	-861.43	kJ/mol	Joback Method
hfus	43.31	kJ/mol	Joback Method
hvap	76.69	kJ/mol	Joback Method
log10ws	-5.01		Crippen Method
logp	4.288		Crippen Method
mcvol	241.430	ml/mol	McGowan Method
pc	1675.53	kPa	Joback Method
rinpol	2142.00		NIST Webbook
rinpol	2142.00		NIST Webbook
tb	791.40	K	Joback Method
tc	991.82	K	Joback Method
tf	496.37	K	Joback Method
vc	0.939	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	681.78	J/mol×K	791.40	Joback Method
cpg	695.20	J/mol×K	824.80	Joback Method
cpg	707.70	J/mol×K	858.21	Joback Method
cpg	719.26	J/mol×K	891.61	Joback Method
cpg	729.92	J/mol×K	925.01	Joback Method
cpg	739.67	J/mol×K	958.42	Joback Method
cpg	748.52	J/mol×K	991.82	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=U389538&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
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
rinpola:	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.cheméo.com/cid/123-619-8/Succinic-acid-2-chloro-6-fluorophenyl-hexyl-ester.pdf>

Generated by Cheméo on 2025-12-22 15:56:26.926407909 +0000 UTC m=+6167184.456448564.

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