Succinic acid, 2-chloro-6-fluorophenyl cis-pent-2-en-1-yl ester

Inchi: InChI=1S/C15H16ClFO4/c1-2-3-4-10-20-13(18)8-9-14(19)21-15-11(16)6-5-7-12(15)17/h3

InchiKey: BCNOMFZMJHTICI-ARJAWSKDSA-N

Formula: C15H16ClFO4

SMILES: CCC=CCOC(=O)CCC(=O)Oc1c(F)cccc1Cl

Mol. weight [g/mol]: 314.74

Physical Properties

Property code	Value	Unit	Source
gf	-425.79	kJ/mol	Joback Method
hf	-723.57	kJ/mol	Joback Method
hfus	40.92	kJ/mol	Joback Method
hvap	74.42	kJ/mol	Joback Method
log10ws	-4.45		Crippen Method
logp	3.674		Crippen Method
mcvol	223.040	ml/mol	McGowan Method
pc	1898.60	kPa	Joback Method
rinpol	2097.00		NIST Webbook
rinpol	2097.00		NIST Webbook
tb	772.68	K	Joback Method
tc	979.96	K	Joback Method
tf	480.02	K	Joback Method
vc	0.863	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	601.45	J/mol×K	772.68	Joback Method
cpg	613.96	J/mol×K	807.23	Joback Method
cpg	625.60	J/mol×K	841.77	Joback Method
cpg	636.40	J/mol×K	876.32	Joback Method
cpg	646.36	J/mol×K	910.87	Joback Method
cpg	655.52	J/mol×K	945.41	Joback Method
cpg	663.89	J/mol×K	979.96	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=U391264&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume