Succinic acid, 8-chlorooctyl pentafluorobenzyl ester

Inchi: InChI=1S/C19H22ClF5O4/c20-9-5-3-1-2-4-6-10-28-13(26)7-8-14(27)29-11-12-15(21)17(2

InchiKey: YLQFBOCVIAEUSV-UHFFFAOYSA-N

Formula: C19H22ClF5O4

SMILES: O=C(CCC(=O)OCc1c(F)c(F)c(F)c(F)c1F)OCCCCCCCCCl

Mol. weight [g/mol]: 444.82

Physical Properties

Property code	Value	Unit	Source
gf	-1280.46	kJ/mol	Joback Method
hf	-1742.20	kJ/mol	Joback Method
hfus	62.23	kJ/mol	Joback Method
hvap	82.09	kJ/mol	Joback Method
log10ws	-6.91		Crippen Method
logp	5.328		Crippen Method
mcvol	290.780	ml/mol	McGowan Method
pc	1144.44	kPa	Joback Method
rinpol	2558.00		NIST Webbook
rinpol	2558.00		NIST Webbook
tb	872.06	K	Joback Method
tc	1067.87	K	Joback Method
tf	570.10	K	Joback Method
vc	1.179	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	879.10	J/mol×K	872.06	Joback Method
cpg	892.17	J/mol×K	904.70	Joback Method
cpg	904.21	J/mol×K	937.33	Joback Method
cpg	915.24	J/mol×K	969.97	Joback Method
cpg	925.25	J/mol×K	1002.60	Joback Method
cpg	934.26	J/mol×K	1035.24	Joback Method
cpg	942.26	J/mol×K	1067.87	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U389892&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
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