Succinic acid, 2-(pentafluorophenoxy)ethyl pentyl ester

Inchi: InChI=1S/C17H19F5O5/c1-2-3-4-7-25-10(23)5-6-11(24)26-8-9-27-17-15(21)13(19)12(18

InchiKey: NKAQQIMFEPQUMR-UHFFFAOYSA-N

Formula: C17H19F5O5

SMILES: CCCCCOC(=O)CCC(=O)OCCOc1c(F)c(F)c(F)c(F)c1F

Mol. weight [g/mol]: 398.32

Physical Properties

Property code	Value	Unit	Source
gf	-1390.37	kJ/mol	Joback Method
hf	-1817.40	kJ/mol	Joback Method
hfus	54.04	kJ/mol	Joback Method
hvap	75.66	kJ/mol	Joback Method
log10ws	-5.16		Crippen Method
logp	3.818		Crippen Method
mcvol	256.230	ml/mol	McGowan Method
pc	1328.10	kPa	Joback Method
rinpol	2138.00		NIST Webbook
rinpol	2138.00		NIST Webbook
tb	811.29	K	Joback Method
tc	996.24	K	Joback Method
tf	539.87	K	Joback Method
vc	1.036	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	766.85	J/mol×K	811.29	Joback Method
cpg	779.72	J/mol×K	842.11	Joback Method
cpg	791.71	J/mol×K	872.94	Joback Method
cpg	802.80	J/mol×K	903.76	Joback Method
cpg	812.98	J/mol×K	934.59	Joback Method
cpg	822.25	J/mol×K	965.41	Joback Method
cpg	830.58	J/mol×K	996.24	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U381550&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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