

Succinic acid, 2-fluorophenyl 4-chlorobenzyl ester

Inchi:	InChI=1S/C17H14ClFO4/c18-13-7-5-12(6-8-13)11-22-16(20)9-10-17(21)23-15-4-2-1-3-14
InchiKey:	CTFFCTSXVAKIGY-UHFFFAOYSA-N
Formula:	C17H14ClFO4
SMILES:	O=C(CCC(=O)Oc1ccccc1F)OCc1ccc(Cl)cc1
Mol. weight [g/mol]:	336.74

Physical Properties

Property code	Value	Unit	Source
gf	-376.76	kJ/mol	Joback Method
hf	-645.54	kJ/mol	Joback Method
hfus	39.94	kJ/mol	Joback Method
hvap	81.19	kJ/mol	Joback Method
log10ws	-5.03		Crippen Method
logp	3.908		Crippen Method
mvol	231.760	ml/mol	McGowan Method
pc	2051.17	kPa	Joback Method
rinpol	2505.00		NIST Webbook
rinpol	2505.00		NIST Webbook
tb	840.96	K	Joback Method
tc	1068.17	K	Joback Method
tf	534.06	K	Joback Method
vc	0.886	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	641.78	J/molxK	840.96	Joback Method
cpg	653.54	J/molxK	878.83	Joback Method
cpg	664.15	J/molxK	916.70	Joback Method
cpg	673.66	J/molxK	954.56	Joback Method
cpg	682.07	J/molxK	992.43	Joback Method
cpg	689.42	J/molxK	1030.30	Joback Method
cpg	695.73	J/molxK	1068.17	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U389675&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.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
hvp:	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
rinp:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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