

Succinic acid, cyclohexylmethyl 3-fluorophenyl ester

Inchi:	InChI=1S/C17H21FO4/c18-14-7-4-8-15(11-14)22-17(20)10-9-16(19)21-12-13-5-2-1-3-6-
InchiKey:	GNGPJCYNFYNJH-UHFFFAOYSA-N
Formula:	C17H21FO4
SMILES:	O=C(CCC(=O)Oc1cccc(F)c1)OCC1CCCCC1
Mol. weight [g/mol]:	308.34

Physical Properties

Property code	Value	Unit	Source
gf	-443.16	kJ/mol	Joback Method
hf	-800.54	kJ/mol	Joback Method
hfus	33.93	kJ/mol	Joback Method
hvap	74.30	kJ/mol	Joback Method
log10ws	-4.40		Crippen Method
logp	3.635		Crippen Method
mcvol	232.420	ml/mol	McGowan Method
pc	1920.30	kPa	Joback Method
rinpol	2210.00		NIST Webbook
rinpol	2210.00		NIST Webbook
tb	791.42	K	Joback Method
tc	1009.64	K	Joback Method
tf	472.58	K	Joback Method
vc	0.878	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	706.12	J/molxK	791.42	Joback Method
cpg	722.45	J/molxK	827.79	Joback Method
cpg	737.45	J/molxK	864.16	Joback Method
cpg	751.12	J/molxK	900.53	Joback Method
cpg	763.51	J/molxK	936.90	Joback Method
cpg	774.62	J/molxK	973.27	Joback Method
cpg	784.49	J/molxK	1009.64	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U390332&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

cp_g:	Ideal gas heat capacity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
rin_{pol}:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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