Succinic acid, 2-chloroethyl 3,5-dimethylphenyl ester

Inchi: InChI=1S/C14H17ClO4/c1-10-7-11(2)9-12(8-10)19-14(17)4-3-13(16)18-6-5-15/h7-9H,3-6

InchiKey: ZJVOYYYZPOJQKF-UHFFFAOYSA-N

Formula: C14H17ClO4

SMILES: Cc1cc(C)cc(OC(=O)CCC(=O)OCCCl)c1

Mol. weight [g/mol]: 284.74

Physical Properties

Property code	Value	Unit	Source
gf	-319.62	kJ/mol	Joback Method
hf	-624.04	kJ/mol	Joback Method
hfus	35.05	kJ/mol	Joback Method
hvap	73.05	kJ/mol	Joback Method
log10ws	-3.43		Crippen Method
logp	2.771		Crippen Method
mcvol	211.480	ml/mol	McGowan Method
pc	2054.89	kPa	Joback Method
rinpol	2123.00		NIST Webbook
rinpol	2123.00		NIST Webbook
tb	746.37	K	Joback Method
tc	956.53	K	Joback Method
tf	473.24	K	Joback Method
vc	0.808	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	564.84	J/mol×K	746.37	Joback Method
cpg	621.94	J/mol×K	921.51	Joback Method
cpg	612.31	J/mol×K	886.48	Joback Method
cpg	601.78	J/mol×K	851.45	Joback Method
cpg	590.37	J/mol×K	816.42	Joback Method
cpg	578.06	J/mol×K	781.40	Joback Method
cpg	630.68	J/mol×K	956.53	Joback Method
dvisc	0.0001079	Pa×s	746.37	Joback Method

dvisc	0.0001337	Pa×s	700.85	Joback Method
dvisc	0.0001708	Pa×s	655.33	Joback Method
dvisc	0.0002262	Pa×s	609.81	Joback Method
dvisc	0.0003135	Pa×s	564.28	Joback Method
dvisc	0.0004601	Pa×s	518.76	Joback Method
dvisc	0.0007271	Pa×s	473.24	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U360727&Units=SI

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
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

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