

Adipic acid, di(2-chloro-5-methylphenyl) ester

Inchi:

InChI=1S/C20H20Cl2O4/c1-13-7-9-15(21)17(11-13)25-19(23)5-3-4-6-20(24)26-18-12-14

InchiKey:

SOJKHWZPVNYKQD-UHFFFAOYSA-N

Formula:

C20H20Cl2O4

SMILES:

Cc1ccc(Cl)c(OC(=O)CCCCC(=O)Oc2cc(C)ccc2Cl)c1

Mol. weight [g/mol]:

395.28

Physical Properties

Property code	Value	Unit	Source
gf	-187.88	kJ/mol	Joback Method
hf	-550.03	kJ/mol	Joback Method
hfus	48.05	kJ/mol	Joback Method
hvap	94.40	kJ/mol	Joback Method
log10ws	-6.91		Crippen Method
logp	5.682		Crippen Method
mcvol	284.500	ml/mol	McGowan Method
pc	1577.21	kPa	Joback Method
rinpol	2981.00		NIST Webbook
tb	957.72	K	Joback Method
tc	1191.17	K	Joback Method
tf	622.24	K	Joback Method
vc	1.085	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	821.52	J/mol×K	957.72	Joback Method
cpg	832.62	J/mol×K	996.63	Joback Method
cpg	842.39	J/mol×K	1035.54	Joback Method
cpg	850.87	J/mol×K	1074.45	Joback Method
cpg	858.09	J/mol×K	1113.35	Joback Method
cpg	864.06	J/mol×K	1152.26	Joback Method
cpg	868.80	J/mol×K	1191.17	Joback Method
dvisc	0.0002590	Pa×s	622.24	Joback Method
dvisc	0.0001710	Pa×s	678.15	Joback Method

dvisc	0.0001202	Pa×s	734.07	Joback Method
dvisc	0.0000889	Pa×s	789.98	Joback Method
dvisc	0.0000684	Pa×s	845.89	Joback Method
dvisc	0.0000543	Pa×s	901.81	Joback Method
dvisc	0.0000444	Pa×s	957.72	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U353968&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
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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