

Chloroacetic acid, 2,3-dimethylphenyl ester

Inchi:	InChI=1S/C10H11ClO2/c1-7-4-3-5-9(8(7)2)13-10(12)6-11/h3-5H,6H2,1-2H3
InchiKey:	QGXOVVMPIWLEDT-UHFFFAOYSA-N
Formula:	C10H11ClO2
SMILES:	<chem>Cc1cccc(OC(=O)CCl)c1C</chem>
Mol. weight [g/mol]:	198.65

Physical Properties

Property code	Value	Unit	Source
gf	-119.38	kJ/mol	Joback Method
hf	-296.68	kJ/mol	Joback Method
hfus	21.90	kJ/mol	Joback Method
hvap	55.00	kJ/mol	Joback Method
log10ws	-2.89		Crippen Method
logp	2.448		Crippen Method
mvol	147.680	ml/mol	McGowan Method
pc	2865.80	kPa	Joback Method
rinpol	1513.00		NIST Webbook
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tb	578.56	K	Joback Method
tc	797.75	K	Joback Method
tf	356.00	K	Joback Method
vc	0.560	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	325.21	J/molxK	578.56	Joback Method
cpg	337.43	J/molxK	615.09	Joback Method
cpg	348.97	J/molxK	651.62	Joback Method
cpg	359.84	J/molxK	688.16	Joback Method
cpg	370.04	J/molxK	724.69	Joback Method
cpg	379.59	J/molxK	761.22	Joback Method
cpg	388.49	J/molxK	797.75	Joback Method
dvisc	0.0013217	Paxs	356.00	Joback Method

dvisc	0.0008372	Paxs	393.09	Joback Method
dvisc	0.0005738	Paxs	430.19	Joback Method
dvisc	0.0004175	Paxs	467.28	Joback Method
dvisc	0.0003184	Paxs	504.37	Joback Method
dvisc	0.0002520	Paxs	541.47	Joback Method
dvisc	0.0002055	Paxs	578.56	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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=U354611&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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