

2-Chlorobenzoic acid, 4-chlorophenyl ester

Inchi:	InChI=1S/C13H8Cl2O2/c14-9-5-7-10(8-6-9)17-13(16)11-3-1-2-4-12(11)15/h1-8H
InchiKey:	FAWUTQLFEATJSE-UHFFFAOYSA-N
Formula:	C13H8Cl2O2
SMILES:	O=C(Oc1ccc(Cl)cc1)c1ccccc1Cl
Mol. weight [g/mol]:	267.11

Physical Properties

Property code	Value	Unit	Source
gf	6.36	kJ/mol	Joback Method
hf	-137.81	kJ/mol	Joback Method
hfus	27.91	kJ/mol	Joback Method
hvap	68.33	kJ/mol	Joback Method
log10ws	-4.93		Crippen Method
logp	4.213		Crippen Method
mcvol	178.430	ml/mol	McGowan Method
pc	2884.30	kPa	Joback Method
rinpola	1992.00		NIST Webbook
tb	711.31	K	Joback Method
tc	966.18	K	Joback Method
tf	446.15	K	Joback Method
vc	0.669	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	410.63	J/molxK	711.31	Joback Method
cpg	422.43	J/molxK	753.79	Joback Method
cpg	433.14	J/molxK	796.27	Joback Method
cpg	442.82	J/molxK	838.75	Joback Method
cpg	451.51	J/molxK	881.22	Joback Method
cpg	459.25	J/molxK	923.70	Joback Method
cpg	466.10	J/molxK	966.18	Joback Method
dvisc	0.0009179	Paxs	446.15	Joback Method
dvisc	0.0005874	Paxs	490.34	Joback Method

dvisc	0.0004047	Paxs	534.54	Joback Method
dvisc	0.0002951	Paxs	578.73	Joback Method
dvisc	0.0002251	Paxs	622.92	Joback Method
dvisc	0.0001780	Paxs	667.12	Joback Method
dvisc	0.0001449	Paxs	711.31	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=U307815&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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