Phenol, 2,4,6-trichloro-, acetate

Other names:	2,4,6-Trichlorophenol acetate
	2,4,6-Trichlorophenyl acetate
	Acetic acid, (2,4,6-trichlorophenyl) ester
	2,4,6-Trichlorfenylester kyseliny octove
Inchi:	InChI=1S/C8H5Cl3O2/c1-4(12)13-8-6(10)2-5(9)3-7(8)11/h2-3H,1H3
InchiKey:	RFOCPJZGJNKBOI-UHFFFAOYSA-N
Formula:	C8H5Cl3O2
SMILES:	CC(=O)Oc1c(Cl)cc(Cl)cc1Cl
Mol. weight [g/mol]:	239.48
CAS:	23399-90-8

Physical Properties

Property code	Value	Unit	Source
gf	-169.71	kJ/mol	Joback Method
hf	-298.35	kJ/mol	Joback Method
hfus	24.73	kJ/mol	Joback Method
hvap	59.97	kJ/mol	Joback Method
log10ws	-3.84		Crippen Method
logp	3.572		Crippen Method
mcvol	143.980	ml/mol	McGowan Method
pc	3210.04	kPa	Joback Method
rinpol	1420.00		NIST Webbook
rinpol	1454.00		NIST Webbook
rinpol	1424.00		NIST Webbook
rinpol	1420.00		NIST Webbook
ripol	2035.00		NIST Webbook
tb	612.64	K	Joback Method
tc	848.96	K	Joback Method
tf	405.82	K	Joback Method
vc	0.546	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source

cpg	277.62	J/mol×K	612.64	Joback Method
cpg	314.35	J/mol×K	809.58	Joback Method
cpg	308.11	J/mol×K	770.19	Joback Method
cpg	301.31	J/mol×K	730.80	Joback Method
cpg	293.97	J/mol×K	691.41	Joback Method
cpg	286.07	J/mol×K	652.03	Joback Method
cpg	320.04	J/mol×K	848.96	Joback Method
dvisc	0.0002256	Pa×s	612.64	Joback Method
dvisc	0.0002695	Pa×s	578.17	Joback Method
dvisc	0.0003292	Pa×s	543.70	Joback Method
dvisc	0.0004132	Pa×s	509.23	Joback Method
dvisc	0.0005361	Pa×s	474.76	Joback Method
dvisc	0.0007244	Pa×s	440.29	Joback Method
dvisc	0.0010302	Pa×s	405.82	Joback Method

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C23399908&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

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
ripol: Polar retention indices
tb: Normal Boiling Point Temperature

tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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