

# Methyl 3,3,3-trichloro-2-hydroxypropanoate

<b>Inchi:</b>	InChI=1S/C4H5Cl3O3/c1-10-3(9)2(8)4(5,6)7/h2,8H,1H3
<b>InchiKey:</b>	ZJUDLWAOAKTCQH-UHFFFAOYSA-N
<b>Formula:</b>	C4H5Cl3O3
<b>SMILES:</b>	COC(=O)C(O)C(Cl)(Cl)Cl
<b>Mol. weight [g/mol]:</b>	207.44

## Physical Properties

Property code	Value	Unit	Source
gf	-423.33	kJ/mol	Joback Method
hf	-584.17	kJ/mol	Joback Method
hfus	14.65	kJ/mol	Joback Method
hvap	61.80	kJ/mol	Joback Method
log10ws	-1.30		Crippen Method
logp	0.891		Crippen Method
mcvol	117.250	ml/mol	McGowan Method
pc	4173.09	kPa	Joback Method
tb	568.01	K	Joback Method
tc	772.13	K	Joback Method
tf	345.00	K	Joback Method
vc	0.432	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	229.94	J/molxK	568.01	Joback Method
cpg	235.89	J/molxK	602.03	Joback Method
cpg	241.40	J/molxK	636.05	Joback Method
cpg	246.51	J/molxK	670.07	Joback Method
cpg	251.23	J/molxK	704.09	Joback Method
cpg	255.58	J/molxK	738.11	Joback Method
cpg	259.57	J/molxK	772.13	Joback Method
dvisc	0.0068244	Paxs	345.00	Joback Method
dvisc	0.0024965	Paxs	382.17	Joback Method
dvisc	0.0010915	Paxs	419.34	Joback Method

dvisc	0.0005460	Paxs	456.50	Joback Method
dvisc	0.0003032	Paxs	493.67	Joback Method
dvisc	0.0001828	Paxs	530.84	Joback Method
dvisc	0.0001178	Paxs	568.01	Joback Method

## Sources

<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R80394&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R80394&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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