

# Carbonic acid, allyl 3,4-dichlorophenyl ester

<b>Inchi:</b>	InChI=1S/C10H8Cl2O3/c1-2-5-14-10(13)15-7-3-4-8(11)9(12)6-7/h2-4,6H,1,5H2
<b>InchiKey:</b>	XPYUFBCLCUBYAN-UHFFFAOYSA-N
<b>Formula:</b>	C10H8Cl2O3
<b>SMILES:</b>	C=CCOC(=O)Oc1ccc(Cl)c(Cl)c1
<b>Mol. weight [g/mol]:</b>	247.07

## Physical Properties

Property code	Value	Unit	Source
gf	-148.47	kJ/mol	Joback Method
hf	-319.21	kJ/mol	Joback Method
hfus	26.01	kJ/mol	Joback Method
hvap	61.12	kJ/mol	Joback Method
log10ws	-3.91		Crippen Method
logp	3.695		Crippen Method
mcvol	161.490	ml/mol	McGowan Method
pc	2829.33	kPa	Joback Method
rinqol	1670.00		NIST Webbook
tb	635.09	K	Joback Method
tc	859.63	K	Joback Method
tf	406.39	K	Joback Method
vc	0.609	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	350.72	J/molxK	635.09	Joback Method
cpg	361.24	J/molxK	672.51	Joback Method
cpg	371.09	J/molxK	709.94	Joback Method
cpg	380.27	J/molxK	747.36	Joback Method
cpg	388.78	J/molxK	784.78	Joback Method
cpg	396.62	J/molxK	822.20	Joback Method
cpg	403.78	J/molxK	859.63	Joback Method
dvisc	0.0009192	Paxs	406.39	Joback Method
dvisc	0.0006090	Paxs	444.51	Joback Method

dvisc	0.0004306	Paxs	482.62	Joback Method
dvisc	0.0003203	Paxs	520.74	Joback Method
dvisc	0.0002481	Paxs	558.86	Joback Method
dvisc	0.0001985	Paxs	596.97	Joback Method
dvisc	0.0001632	Paxs	635.09	Joback Method

## Sources

<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>
<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=U357812&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U357812&amp;Units=SI</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>rinpol:</b>	Non-polar retention indices
<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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