

# 3-chloro-2-butoxy-tetrahydro-pyran

<b>Inchi:</b>	InChI=1S/C9H17ClO2/c1-2-3-6-11-9-8(10)5-4-7-12-9/h8-9H,2-7H2,1H3
<b>InchiKey:</b>	YVJCTCVZSCULEB-UHFFFAOYSA-N
<b>Formula:</b>	C9H17ClO2
<b>SMILES:</b>	CCCCOC1OCCCC1Cl
<b>Mol. weight [g/mol]:</b>	192.68

## Physical Properties

Property code	Value	Unit	Source
gf	-161.41	kJ/mol	Joback Method
hf	-475.07	kJ/mol	Joback Method
hfus	25.34	kJ/mol	Joback Method
hvap	47.05	kJ/mol	Joback Method
log10ws	-2.53		Crippen Method
logp	2.547		Crippen Method
mcvol	150.790	ml/mol	McGowan Method
pc	2568.89	kPa	Joback Method
rinpol	1275.00		NIST Webbook
rinpol	1275.00		NIST Webbook
tb	507.00	K	Joback Method
tc	709.81	K	Joback Method
tf	273.05	K	Joback Method
vc	0.559	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	348.05	J/molxK	507.00	Joback Method
cpg	426.03	J/molxK	676.01	Joback Method
cpg	412.01	J/molxK	642.20	Joback Method
cpg	397.21	J/molxK	608.40	Joback Method
cpg	381.62	J/molxK	574.60	Joback Method
cpg	365.23	J/molxK	540.80	Joback Method
cpg	439.27	J/molxK	709.81	Joback Method
dvisc	0.0002651	Paxs	507.00	Joback Method

dvisc	0.0003423	Paxs	468.01	Joback Method
dvisc	0.0004629	Paxs	429.02	Joback Method
dvisc	0.0006650	Paxs	390.02	Joback Method
dvisc	0.0010355	Paxs	351.03	Joback Method
dvisc	0.0018009	Paxs	312.04	Joback Method
dvisc	0.0036685	Paxs	273.05	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=R132920&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R132920&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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