

# Hexyl chloroformate

<b>Other names:</b>	Carbonochloridic acid, hexyl ester Hexyl chloridocarbonate
<b>Inchi:</b>	InChI=1S/C7H13ClO2/c1-2-3-4-5-6-10-7(8)9/h2-6H2,1H3
<b>InchiKey:</b>	KIWBRXCOTCXSSZ-UHFFFAOYSA-N
<b>Formula:</b>	C7H13ClO2
<b>SMILES:</b>	CCCCCCOC(=O)Cl
<b>Mol. weight [g/mol]:</b>	164.63
<b>CAS:</b>	6092-54-2

## Physical Properties

Property code	Value	Unit	Source
gf	-237.79	kJ/mol	Joback Method
hf	-448.35	kJ/mol	Joback Method
hfus	20.87	kJ/mol	Joback Method
hvap	44.72	kJ/mol	Joback Method
log10ws	-2.75		Crippen Method
logp	2.942		Crippen Method
mvol	129.170	ml/mol	McGowan Method
pc	2826.33	kPa	Joback Method
rinpol	1055.00		NIST Webbook
rinpol	1055.00		NIST Webbook
tb	473.28	K	Joback Method
tc	655.90	K	Joback Method
tf	270.73	K	Joback Method
vc	0.500	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	265.22	J/mol×K	473.28	Joback Method
cpg	276.01	J/mol×K	503.72	Joback Method
cpg	286.40	J/mol×K	534.15	Joback Method
cpg	296.38	J/mol×K	564.59	Joback Method
cpg	305.96	J/mol×K	595.03	Joback Method

cpg	315.13	J/molxK	625.46	Joback Method
cpg	323.90	J/molxK	655.90	Joback Method
dvisc	0.0030645	Paxs	270.73	Joback Method
dvisc	0.0016518	Paxs	304.49	Joback Method
dvisc	0.0010072	Paxs	338.25	Joback Method
dvisc	0.0006719	Paxs	372.00	Joback Method
dvisc	0.0004794	Paxs	405.76	Joback Method
dvisc	0.0003603	Paxs	439.52	Joback Method
dvisc	0.0002820	Paxs	473.28	Joback Method

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	333.70	K	0.90	NIST Webbook

## Sources

<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=C6092542&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C6092542&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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>

## 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>tbrp:</b>	Boiling point at reduced pressure
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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