

# 2-propoxy-tetrahydro-furan

<b>Other names:</b>	Tetrahydrofuran, 2-propyloxy
<b>Inchi:</b>	InChI=1S/C7H14O2/c1-2-5-8-7-4-3-6-9-7/h7H,2-6H2,1H3
<b>InchiKey:</b>	OHNPPRNQKABHPI-UHFFFAOYSA-N
<b>Formula:</b>	C7H14O2
<b>SMILES:</b>	CCCOC1CCCO1
<b>Mol. weight [g/mol]:</b>	130.18

## Physical Properties

Property code	Value	Unit	Source
gf	-146.51	kJ/mol	Joback Method
hf	-391.55	kJ/mol	Joback Method
hfus	16.99	kJ/mol	Joback Method
hvap	38.35	kJ/mol	Joback Method
log10ws	-1.43		Crippen Method
logp	1.550		Crippen Method
mcvol	110.370	ml/mol	McGowan Method
pc	3318.18	kPa	Joback Method
rinpol	900.00		NIST Webbook
rinpol	900.00		NIST Webbook
tb	424.21	K	Joback Method
tc	618.47	K	Joback Method
tf	228.35	K	Joback Method
vc	0.407	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	229.35	J/molxK	424.21	Joback Method
cpg	295.43	J/molxK	586.09	Joback Method
cpg	283.41	J/molxK	553.71	Joback Method
cpg	270.80	J/molxK	521.34	Joback Method
cpg	257.59	J/molxK	488.96	Joback Method
cpg	243.78	J/molxK	456.59	Joback Method
cpg	306.88	J/molxK	618.47	Joback Method

dvisc	0.0003243	Paxs	424.21	Joback Method
dvisc	0.0004129	Paxs	391.57	Joback Method
dvisc	0.0005494	Paxs	358.92	Joback Method
dvisc	0.0007739	Paxs	326.28	Joback Method
dvisc	0.0011766	Paxs	293.64	Joback Method
dvisc	0.0019863	Paxs	260.99	Joback Method
dvisc	0.0038949	Paxs	228.35	Joback Method

## Sources

<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>
<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=R91141&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R91141&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>mvol:</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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