

# furfural ethyl propyl acetal

<b>Inchi:</b>	InChI=1S/C11H18O2/c1-3-9-13-11(12-4-2)10-7-5-6-8-10/h5-7,11H,3-4,8-9H2,1-2H3
<b>InchiKey:</b>	AASUBIVBGMLSMB-UHFFFAOYSA-N
<b>Formula:</b>	C11H18O2
<b>SMILES:</b>	CCCOC(OCC)C1=CC=CC1
<b>Mol. weight [g/mol]:</b>	182.26

## Physical Properties

Property code	Value	Unit	Source
gf	-76.15	kJ/mol	Joback Method
hf	-355.18	kJ/mol	Joback Method
hfus	18.02	kJ/mol	Joback Method
hvap	46.32	kJ/mol	Joback Method
log10ws	-2.81		Crippen Method
logp	2.662		Crippen Method
mcvol	158.130	ml/mol	McGowan Method
pc	2426.65	kPa	Joback Method
ripol	1550.00		NIST Webbook
tb	518.73	K	Joback Method
tc	712.95	K	Joback Method
tf	272.37	K	Joback Method
vc	0.596	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	368.75	J/molxK	518.73	Joback Method
cpg	384.61	J/molxK	551.10	Joback Method
cpg	399.74	J/molxK	583.47	Joback Method
cpg	414.17	J/molxK	615.84	Joback Method
cpg	427.89	J/molxK	648.21	Joback Method
cpg	440.93	J/molxK	680.58	Joback Method
cpg	453.29	J/molxK	712.95	Joback Method
dvisc	0.0026327	Paxs	272.37	Joback Method
dvisc	0.0012528	Paxs	313.43	Joback Method

dvisc	0.0007081	Paxs	354.49	Joback Method
dvisc	0.0004505	Paxs	395.55	Joback Method
dvisc	0.0003121	Paxs	436.61	Joback Method
dvisc	0.0002303	Paxs	477.67	Joback Method
dvisc	0.0001783	Paxs	518.73	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=R340879&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R340879&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>

## 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>ripol:</b>	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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