

# 3-heptanon-2-ol

<b>Inchi:</b>	InChI=1S/C7H14O2/c1-3-4-5-7(9)6(2)8/h6,8H,3-5H2,1-2H3
<b>InchiKey:</b>	KTKQJWVLAMSBTM-UHFFFAOYSA-N
<b>Formula:</b>	C7H14O2
<b>SMILES:</b>	CCCCC(=O)C(C)O
<b>Mol. weight [g/mol]:</b>	130.18

## Physical Properties

Property code	Value	Unit	Source
gf	-260.12	kJ/mol	Joback Method
hf	-457.90	kJ/mol	Joback Method
hfus	16.05	kJ/mol	Joback Method
hvap	54.21	kJ/mol	Joback Method
log10ws	-1.41		Crippen Method
logp	1.127		Crippen Method
mcvol	116.930	ml/mol	McGowan Method
pc	3352.86	kPa	Joback Method
ripol	1551.00		NIST Webbook
ripol	1551.00		NIST Webbook
tb	505.17	K	Joback Method
tc	678.63	K	Joback Method
tf	264.40	K	Joback Method
vc	0.447	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	265.80	J/molxK	505.17	Joback Method
cpg	312.64	J/molxK	649.72	Joback Method
cpg	304.09	J/molxK	620.81	Joback Method
cpg	295.14	J/molxK	591.90	Joback Method
cpg	285.78	J/molxK	562.99	Joback Method
cpg	276.01	J/molxK	534.08	Joback Method
cpg	320.81	J/molxK	678.63	Joback Method
dvisc	0.0001662	Paxs	505.17	Joback Method

dvisc	0.0002761	Paxs	465.04	Joback Method
dvisc	0.0005050	Paxs	424.91	Joback Method
dvisc	0.0010476	Paxs	384.78	Joback Method
dvisc	0.0025758	Paxs	344.66	Joback Method
dvisc	0.0080276	Paxs	304.53	Joback Method
dvisc	0.0353274	Paxs	264.40	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=R241059&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R241059&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

Latest version available from:

<https://www.chemeo.com/cid/92-650-8/3-heptanon-2-ol.pdf>

Generated by Cheméo on 2024-04-24 19:31:23.833027162 +0000 UTC m=+16276332.753604479.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.