

# 2,6-Dimethyl-5-oxo-heptanoic acid

<b>Inchi:</b>	InChI=1S/C9H16O3/c1-6(2)8(10)5-4-7(3)9(11)12/h6-7H,4-5H2,1-3H3,(H,11,12)
<b>InchiKey:</b>	KFSWOOWOYKTCDK-UHFFFAOYSA-N
<b>Formula:</b>	C9H16O3
<b>SMILES:</b>	CC(C)C(=O)CCC(C)C(=O)O
<b>Mol. weight [g/mol]:</b>	172.22

## Physical Properties

Property code	Value	Unit	Source
gf	-374.64	kJ/mol	Joback Method
hf	-617.04	kJ/mol	Joback Method
hfus	19.31	kJ/mol	Joback Method
hvap	65.02	kJ/mol	Joback Method
log10ws	-1.49		Crippen Method
logp	1.712		Crippen Method
mcvol	146.680	ml/mol	McGowan Method
pc	2928.17	kPa	Joback Method
ripol	2515.00		NIST Webbook
tb	604.36	K	Joback Method
tc	786.29	K	Joback Method
tf	321.87	K	Joback Method
vc	0.558	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	373.45	J/molxK	604.36	Joback Method
cpg	384.73	J/molxK	634.68	Joback Method
cpg	395.46	J/molxK	665.00	Joback Method
cpg	405.65	J/molxK	695.32	Joback Method
cpg	415.32	J/molxK	725.64	Joback Method
cpg	424.48	J/molxK	755.97	Joback Method
cpg	433.14	J/molxK	786.29	Joback Method
dvisc	0.0137516	Paxs	321.87	Joback Method
dvisc	0.0035083	Paxs	368.95	Joback Method

dvisc	0.0012193	Paxs	416.03	Joback Method
dvisc	0.0005254	Paxs	463.12	Joback Method
dvisc	0.0002644	Paxs	510.20	Joback Method
dvisc	0.0001495	Paxs	557.28	Joback Method
dvisc	0.0000923	Paxs	604.36	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R326100&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R326100&amp;Units=SI</a>
<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>

## 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/92-117-0/2-6-Dimethyl-5-oxo-heptanoic-acid.pdf>

Generated by Cheméo on 2024-04-24 07:29:05.072421448 +0000 UTC m=+16232993.992998759.

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