

# Methyl 4-oxooctanoate

<b>Inchi:</b>	InChI=1S/C9H16O3/c1-3-4-5-8(10)6-7-9(11)12-2/h3-7H2,1-2H3
<b>InchiKey:</b>	PONRPCIHEGPNCW-UHFFFAOYSA-N
<b>Formula:</b>	C9H16O3
<b>SMILES:</b>	CCCCC(=O)CCC(=O)OC
<b>Mol. weight [g/mol]:</b>	172.22

## Physical Properties

Property code	Value	Unit	Source
gf	-337.94	kJ/mol	Joback Method
hf	-586.47	kJ/mol	Joback Method
hfus	23.45	kJ/mol	Joback Method
hvap	51.53	kJ/mol	Joback Method
log10ws	-1.73		Crippen Method
logp	1.699		Crippen Method
mcvol	146.680	ml/mol	McGowan Method
pc	2568.89	kPa	Joback Method
ripol	1797.00		NIST Webbook
tb	535.48	K	Joback Method
tc	718.27	K	Joback Method
tf	313.28	K	Joback Method
vc	0.570	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	343.51	J/molxK	535.48	Joback Method
cpg	356.03	J/molxK	565.94	Joback Method
cpg	368.04	J/molxK	596.41	Joback Method
cpg	379.53	J/molxK	626.87	Joback Method
cpg	390.51	J/molxK	657.34	Joback Method
cpg	400.99	J/molxK	687.80	Joback Method
cpg	410.97	J/molxK	718.27	Joback Method
dvisc	0.0026992	Paxs	313.28	Joback Method
dvisc	0.0014842	Paxs	350.31	Joback Method

dvisc	0.0009150	Paxs	387.35	Joback Method
dvisc	0.0006138	Paxs	424.38	Joback Method
dvisc	0.0004390	Paxs	461.41	Joback Method
dvisc	0.0003300	Paxs	498.45	Joback Method
dvisc	0.0002580	Paxs	535.48	Joback Method

## 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=R289239&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R289239&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>

## 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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