

# 2-(Cyclopenten-1-yl)acetic acid

<b>Other names:</b>	cyclopent-2-enylacetic acid
<b>Inchi:</b>	InChI=1S/C7H10O2/c8-7(9)5-6-3-1-2-4-6/h1,3,6H,2,4-5H2,(H,8,9)
<b>InchiKey:</b>	NNRZTJAACCRFRV-UHFFFAOYSA-N
<b>Formula:</b>	C7H10O2
<b>SMILES:</b>	O=C(O)CC1C=CCC1
<b>Mol. weight [g/mol]:</b>	126.15
<b>CAS:</b>	13668-61-6

## Physical Properties

Property code	Value	Unit	Source
gf	-191.17	kJ/mol	Joback Method
hf	-334.36	kJ/mol	Joback Method
hfus	14.73	kJ/mol	Joback Method
hvap	55.15	kJ/mol	Joback Method
log10ws	-1.36		Crippen Method
logp	1.427		Crippen Method
mcvol	101.770	ml/mol	McGowan Method
pc	4333.96	kPa	Joback Method
tb	520.05	K	Joback Method
tc	717.25	K	Joback Method
tf	291.06	K	Joback Method
vc	0.380	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	232.43	J/mol×K	520.05	Joback Method
cpg	281.06	J/mol×K	684.39	Joback Method
cpg	272.49	J/mol×K	651.52	Joback Method
cpg	263.36	J/mol×K	618.65	Joback Method
cpg	253.66	J/mol×K	585.78	Joback Method
cpg	243.36	J/mol×K	552.92	Joback Method
cpg	289.10	J/mol×K	717.25	Joback Method
dvisc	0.0002181	Paxs	520.05	Joback Method

dvisc	0.0003315	Paxs	481.88	Joback Method
dvisc	0.0005417	Paxs	443.72	Joback Method
dvisc	0.0009707	Paxs	405.55	Joback Method
dvisc	0.0019637	Paxs	367.39	Joback Method
dvisc	0.0046771	Paxs	329.23	Joback Method
dvisc	0.0139872	Paxs	291.06	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=C13668616&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C13668616&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>cp<sub>g</sub>:</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>h<sub>vap</sub>:</b>	Enthalpy of vaporization at standard conditions
<b>log<sub>10</sub>ws:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<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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