

# Cyclopentanecarboxylic acid, ethyl ester

Inchi:	InChI=1S/C8H14O2/c1-2-10-8(9)7-5-3-4-6-7/h7H,2-6H2,1H3
InchiKey:	UWSJCCUODNDXOT-UHFFFAOYSA-N
Formula:	C8H14O2
SMILES:	CCOC(=O)C1CCCC1
Mol. weight [g/mol]:	142.20

## Physical Properties

Property code	Value	Unit	Source
gf	-180.89	kJ/mol	Joback Method
hf	-392.77	kJ/mol	Joback Method
hfus	13.20	kJ/mol	Joback Method
hvap	42.81	kJ/mol	Joback Method
log10ws	-1.69		Crippen Method
logp	1.740		Crippen Method
mcvol	120.160	ml/mol	McGowan Method
pc	3228.31	kPa	Joback Method
rinsol	1036.00		NIST Webbook
tb	474.01	K	Joback Method
tc	677.64	K	Joback Method
tf	262.98	K	Joback Method
vc	0.449	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	263.68	J/molxK	474.01	Joback Method
cpg	332.13	J/molxK	643.70	Joback Method
cpg	319.85	J/molxK	609.76	Joback Method
cpg	306.88	J/molxK	575.83	Joback Method
cpg	293.21	J/molxK	541.89	Joback Method
cpg	278.81	J/molxK	507.95	Joback Method
cpg	343.74	J/molxK	677.64	Joback Method
dvisc	0.0003332	Paxs	474.01	Joback Method
dvisc	0.0004182	Paxs	438.84	Joback Method

dvisc	0.0005459	Paxs	403.67	Joback Method
dvisc	0.0007500	Paxs	368.50	Joback Method
dvisc	0.0011016	Paxs	333.32	Joback Method
dvisc	0.0017719	Paxs	298.15	Joback Method
dvisc	0.0032364	Paxs	262.98	Joback Method

## Sources

<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>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U375864&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U375864&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>

## 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>rinpol:</b>	Non-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/55-355-7/Cyclopentanecarboxylic-acid-ethyl-ester.pdf>

Generated by Cheméo on 2024-04-25 09:26:17.817506669 +0000 UTC m=+16326426.738083979.

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