

# Cyclopropane, 1-propenyl-

<b>Other names:</b>	Propene, 1-cyclopropyl- Cyclopropane, propenyl- 1-Propene, 1-cyclopropyl-
<b>Inchi:</b>	InChI=1S/C6H10/c1-2-3-6-4-5-6/h2-3,6H,4-5H2,1H3/b3-2+
<b>InchiKey:</b>	TWAJIXJFYMNJJO-NSCUHMNNSA-N
<b>Formula:</b>	C6H10
<b>SMILES:</b>	CC=CC1CC1
<b>Mol. weight [g/mol]:</b>	82.14
<b>CAS:</b>	4663-21-2

## Physical Properties

Property code	Value	Unit	Source
gf	140.61	kJ/mol	Joback Method
hf	22.85	kJ/mol	Joback Method
hfus	9.63	kJ/mol	Joback Method
hvap	28.82	kJ/mol	Joback Method
log10ws	-1.84		Crippen Method
logp	1.973		Crippen Method
mcvol	80.240	ml/mol	McGowan Method
pc	3848.31	kPa	Joback Method
rinpol	684.00		NIST Webbook
tb	347.58	K	Joback Method
tc	536.08	K	Joback Method
tf	170.24	K	Joback Method
vc	0.308	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	127.26	J/molxK	347.58	Joback Method
cpg	180.94	J/molxK	504.66	Joback Method
cpg	171.52	J/molxK	473.24	Joback Method
cpg	161.48	J/molxK	441.83	Joback Method
cpg	150.78	J/molxK	410.41	Joback Method

cpg	139.39	J/molxK	379.00	Joback Method
cpg	189.79	J/molxK	536.08	Joback Method
dvisc	0.0002305	Paxs	347.58	Joback Method
dvisc	0.0002521	Paxs	318.02	Joback Method
dvisc	0.0002809	Paxs	288.47	Joback Method
dvisc	0.0003209	Paxs	258.91	Joback Method
dvisc	0.0003792	Paxs	229.35	Joback Method
dvisc	0.0004709	Paxs	199.80	Joback Method
dvisc	0.0006305	Paxs	170.24	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=C4663212&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C4663212&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>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

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