

# trans-1-Methyl-2-propylcyclopropane

<b>Other names:</b>	1-methyl-trans-2-propyl-cyclopropane
<b>Inchi:</b>	InChI=1S/C7H14/c1-3-4-7-5-6(7)2/h6-7H,3-5H2,1-2H3/t6-,7-/m1/s1
<b>InchiKey:</b>	SSOHKKMWPDYMSB-RNFRBKRXSA-N
<b>Formula:</b>	C7H14
<b>SMILES:</b>	CCCC1CC1C
<b>Mol. weight [g/mol]:</b>	98.19

## Physical Properties

Property code	Value	Unit	Source
gf	61.10	kJ/mol	Joback Method
hf	-135.35	kJ/mol	Joback Method
hfus	13.09	kJ/mol	Joback Method
hvap	30.78	kJ/mol	Joback Method
log10ws	-2.16		Crippen Method
logp	2.442		Crippen Method
mcvol	98.630	ml/mol	McGowan Method
pc	3110.57	kPa	Joback Method
rinpol	668.00		NIST Webbook
rinpol	668.00		NIST Webbook
tb	361.63	K	Joback Method
tc	539.10	K	Joback Method
tf	182.35	K	Joback Method
vc	0.384	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	177.61	J/mol×K	361.63	Joback Method
cpg	191.26	J/mol×K	391.21	Joback Method
cpg	204.28	J/mol×K	420.79	Joback Method
cpg	216.71	J/mol×K	450.36	Joback Method
cpg	228.57	J/mol×K	479.94	Joback Method
cpg	239.87	J/mol×K	509.52	Joback Method
cpg	250.64	J/mol×K	539.10	Joback Method

dvisc	0.0004919	Paxs	182.35	Joback Method
dvisc	0.0004196	Paxs	212.23	Joback Method
dvisc	0.0003722	Paxs	242.11	Joback Method
dvisc	0.0003390	Paxs	271.99	Joback Method
dvisc	0.0003145	Paxs	301.87	Joback Method
dvisc	0.0002957	Paxs	331.75	Joback Method
dvisc	0.0002809	Paxs	361.63	Joback Method

## Sources

<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>
<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=R137549&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R137549&amp;Units=SI</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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