

# Pentane, 1-chloro-4,4-dimethyl

Inchi:	InChI=1S/C7H15Cl/c1-7(2,3)5-4-6-8/h4-6H2,1-3H3
InchiKey:	LEQALULJXWQRKX-UHFFFAOYSA-N
Formula:	C7H15Cl
SMILES:	CC(C)(C)CCCCl
Mol. weight [g/mol]:	134.65

## Physical Properties

Property code	Value	Unit	Source
gf	-1.03	kJ/mol	Joback Method
hf	-212.30	kJ/mol	Joback Method
hfus	10.67	kJ/mol	Joback Method
hvap	34.27	kJ/mol	Joback Method
log10ws	-2.66		Crippen Method
logp	3.051		Crippen Method
mcvol	121.730	ml/mol	McGowan Method
pc	2746.90	kPa	Joback Method
rinpol	767.00		NIST Webbook
tb	393.76	K	Joback Method
tc	576.83	K	Joback Method
tf	200.99	K	Joback Method
vc	0.466	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	224.71	J/mol×K	393.76	Joback Method
cpg	283.83	J/mol×K	546.32	Joback Method
cpg	273.22	J/mol×K	515.80	Joback Method
cpg	262.03	J/mol×K	485.29	Joback Method
cpg	250.23	J/mol×K	454.78	Joback Method
cpg	237.80	J/mol×K	424.27	Joback Method
cpg	293.88	J/mol×K	576.83	Joback Method
dvisc	0.0003185	Paxs	393.76	Joback Method
dvisc	0.0004349	Paxs	361.63	Joback Method

dvisc	0.0006312	Paxs	329.50	Joback Method
dvisc	0.0009928	Paxs	297.38	Joback Method
dvisc	0.0017428	Paxs	265.25	Joback Method
dvisc	0.0035725	Paxs	233.12	Joback Method
dvisc	0.0092120	Paxs	200.99	Joback Method

## Sources

<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>
<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=R129701&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R129701&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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