

# 1-Ethyl-3-methylcyclohexane (c,t)

<b>Other names:</b>	1-Ethyl-3-methylcyclohexane
<b>Inchi:</b>	InChI=1S/C9H18/c1-3-9-6-4-5-8(2)7-9/h8-9H,3-7H2,1-2H3
<b>InchiKey:</b>	UDDVMPHNQKRNS-UHFFFAOYSA-N
<b>Formula:</b>	C9H18
<b>SMILES:</b>	CCC1CCCC(C)C1
<b>Mol. weight [g/mol]:</b>	126.24
<b>CAS:</b>	3728-55-0

## Physical Properties

Property code	Value	Unit	Source
gf	41.64	kJ/mol	Joback Method
hf	-195.11	kJ/mol	Joback Method
hfus	11.97	kJ/mol	Joback Method
hvap	35.75	kJ/mol	Joback Method
log10ws	-3.00		Crippen Method
logp	3.223		Crippen Method
mcvol	126.810	ml/mol	McGowan Method
pc	2741.15	kPa	Joback Method
rinpol	931.00		NIST Webbook
tb	420.20	K	Joback Method
tc	618.70	K	Joback Method
tf	194.33	K	Joback Method
vc	0.471	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	255.07	J/mol×K	420.20	Joback Method
cpg	273.91	J/mol×K	453.28	Joback Method
cpg	291.90	J/mol×K	486.37	Joback Method
cpg	309.07	J/mol×K	519.45	Joback Method
cpg	325.42	J/mol×K	552.53	Joback Method
cpg	340.98	J/mol×K	585.62	Joback Method
cpg	355.76	J/mol×K	618.70	Joback Method

dvisc	0.0046654	Paxs	194.33	Joback Method
dvisc	0.0019666	Paxs	231.97	Joback Method
dvisc	0.0010552	Paxs	269.62	Joback Method
dvisc	0.0006594	Paxs	307.26	Joback Method
dvisc	0.0004567	Paxs	344.91	Joback Method
dvisc	0.0003400	Paxs	382.56	Joback Method
dvisc	0.0002668	Paxs	420.20	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=C3728550&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3728550&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>
<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>mvol:</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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