

# Indene, 6-methyl-3a,4,7,7a-tetrahydro

<b>Other names:</b>	Indene, 3a,4,7,7a-tetrahydro-6-methyl
<b>Inchi:</b>	InChI=1S/C10H14/c1-8-5-6-9-3-2-4-10(9)7-8/h2-3,5,9-10H,4,6-7H2,1H3
<b>InchiKey:</b>	MGBYEAMVKYHVOF-UHFFFAOYSA-N
<b>Formula:</b>	C10H14
<b>SMILES:</b>	CC1=CCC2C=CCC2C1
<b>Mol. weight [g/mol]:</b>	134.22

## Physical Properties

Property code	Value	Unit	Source
gf	168.81	kJ/mol	Joback Method
hf	-18.52	kJ/mol	Joback Method
hfus	13.68	kJ/mol	Joback Method
hvap	39.44	kJ/mol	Joback Method
log10ws	-3.02		Crippen Method
logp	2.919		Crippen Method
mvol	121.440	ml/mol	McGowan Method
pc	3135.00	kPa	Joback Method
rinpol	1113.00		NIST Webbook
rinpol	1032.00		NIST Webbook
tb	457.79	K	Joback Method
tc	677.10	K	Joback Method
tf	241.82	K	Joback Method
vc	0.458	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	257.37	J/molxK	457.79	Joback Method
cpg	275.55	J/molxK	494.34	Joback Method
cpg	292.58	J/molxK	530.89	Joback Method
cpg	308.52	J/molxK	567.44	Joback Method
cpg	323.41	J/molxK	604.00	Joback Method
cpg	337.33	J/molxK	640.55	Joback Method
cpg	350.33	J/molxK	677.10	Joback Method

dvisc	0.0014256	Paxs	241.82	Joback Method
dvisc	0.0010239	Paxs	277.81	Joback Method
dvisc	0.0007934	Paxs	313.81	Joback Method
dvisc	0.0006480	Paxs	349.80	Joback Method
dvisc	0.0005495	Paxs	385.80	Joback Method
dvisc	0.0004794	Paxs	421.79	Joback Method
dvisc	0.0004272	Paxs	457.79	Joback Method

## Sources

<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=R3622&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R3622&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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</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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