

# Cyclohexene, 6-iodo-1-methyl

Inchi:	InChI=1S/C7H11I/c1-6-4-2-3-5-7(6)8/h4,7H,2-3,5H2,1H3
InchiKey:	HOTMTBUTWQOXMD-UHFFFAOYSA-N
Formula:	C7H11I
SMILES:	CC1=CCCCC1I
Mol. weight [g/mol]:	222.07

## Physical Properties

Property code	Value	Unit	Source
gf	110.96	kJ/mol	Joback Method
hf	-10.31	kJ/mol	Joback Method
hfus	10.96	kJ/mol	Joback Method
hvap	41.93	kJ/mol	Joback Method
log10ws	-3.56		Crippen Method
logp	2.920		Crippen Method
mcvol	120.150	ml/mol	McGowan Method
pc	3517.91	kPa	Joback Method
rinpol	1181.00		NIST Webbook
tb	476.39	K	Joback Method
tc	722.59	K	Joback Method
tf	247.37	K	Joback Method
vc	0.434	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	216.17	J/molxK	476.39	Joback Method
cpg	279.33	J/molxK	681.56	Joback Method
cpg	268.44	J/molxK	640.53	Joback Method
cpg	256.72	J/molxK	599.49	Joback Method
cpg	244.12	J/molxK	558.46	Joback Method
cpg	230.62	J/molxK	517.42	Joback Method
cpg	289.44	J/molxK	722.59	Joback Method
dvisc	0.0003557	Paxs	476.39	Joback Method
dvisc	0.0004513	Paxs	438.22	Joback Method

dvisc	0.0005992	Paxs	400.05	Joback Method
dvisc	0.0008447	Paxs	361.88	Joback Method
dvisc	0.0012912	Paxs	323.71	Joback Method
dvisc	0.0022107	Paxs	285.54	Joback Method
dvisc	0.0044685	Paxs	247.37	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=R25508&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R25508&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>

## 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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