

2-Menthene, cis

Other names:	cis-3-isopropyl-6-methylcyclohexene
Inchi:	InChI=1S/C10H18/c1-8(2)10-6-4-9(3)5-7-10/h4,6,8-10H,5,7H2,1-3H3/t9-,10+/m0/s1
InchiKey:	WHNGPXQYYRWQAS-VHSXEESVSA-N
Formula:	C10H18
SMILES:	CC1C=CC(C(C)C)CC1
Mol. weight [g/mol]:	138.25
CAS:	2230-85-5

Physical Properties

Property code	Value	Unit	Source
gf	77.58	kJ/mol	Joback Method
hf	-163.25	kJ/mol	Joback Method
hfus	12.26	kJ/mol	Joback Method
hvap	37.88	kJ/mol	Joback Method
log10ws	-3.03		Crippen Method
logp	3.245		Crippen Method
mcvol	136.600	ml/mol	McGowan Method
pc	2597.78	kPa	Joback Method
rinpol	981.90		NIST Webbook
rinpol	976.00		NIST Webbook
rinpol	978.20		NIST Webbook
rinpol	976.00		NIST Webbook
ripol	1078.00		NIST Webbook
tb	441.80	K	Joback Method
tc	644.93	K	Joback Method
tf	191.36	K	Joback Method
vc	0.507	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	285.08	J/mol×K	441.80	Joback Method
cpg	371.86	J/mol×K	611.07	Joback Method
cpg	356.27	J/mol×K	577.22	Joback Method

cpg	339.82	J/mol×K	543.36	Joback Method
cpg	322.48	J/mol×K	509.51	Joback Method
cpg	304.25	J/mol×K	475.65	Joback Method
cpg	386.61	J/mol×K	644.93	Joback Method
dvisc	0.0002442	Pa×s	441.80	Joback Method
dvisc	0.0003145	Pa×s	400.06	Joback Method
dvisc	0.0004295	Pa×s	358.32	Joback Method
dvisc	0.0006369	Pa×s	316.58	Joback Method
dvisc	0.0010645	Pa×s	274.84	Joback Method
dvisc	0.0021384	Pa×s	233.10	Joback Method
dvisc	0.0058238	Pa×s	191.36	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C2230855&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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