

cis,cis-Bicyclo[6.1.0]non-3-ene

Inchi:	InChI=1S/C9H14/c1-2-4-6-9-7-8(9)5-3-1/h1,3,8-9H,2,4-7H2/b3-1-/t8-,9+/m1/s1
InchiKey:	HTNGHCFQBNOUQK-LGSOUNPISA-N
Formula:	C9H14
SMILES:	C1=CCC2CC2CCC1
Mol. weight [g/mol]:	122.21

Physical Properties

Property code	Value	Unit	Source
gf	140.06	kJ/mol	Joback Method
hf	-44.19	kJ/mol	Joback Method
hfus	10.26	kJ/mol	Joback Method
hvap	36.26	kJ/mol	Joback Method
log10ws	-2.75		Crippen Method
logp	2.753		Crippen Method
mcvol	111.650	ml/mol	McGowan Method
pc	3415.86	kPa	Joback Method
rinsol	1012.00		NIST Webbook
tb	430.77	K	Joback Method
tc	649.36	K	Joback Method
tf	217.27	K	Joback Method
vc	0.415	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	226.80	J/molxK	430.77	Joback Method
cpg	310.78	J/molxK	612.93	Joback Method
cpg	296.19	J/molxK	576.50	Joback Method
cpg	280.56	J/molxK	540.06	Joback Method
cpg	263.83	J/molxK	503.63	Joback Method
cpg	245.93	J/molxK	467.20	Joback Method
cpg	324.39	J/molxK	649.36	Joback Method
dvisc	0.0004387	Paxs	430.77	Joback Method
dvisc	0.0005047	Paxs	395.19	Joback Method

dvisc	0.0005970	Paxs	359.60	Joback Method
dvisc	0.0007327	Paxs	324.02	Joback Method
dvisc	0.0009458	Paxs	288.44	Joback Method
dvisc	0.0013118	Paxs	252.85	Joback Method
dvisc	0.0020254	Paxs	217.27	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=R293486&Units=SI

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/15-599-1/cis-cis-Bicyclo-6-1-0-non-3-ene.pdf>

Generated by Cheméo on 2024-04-19 22:24:18.365346524 +0000 UTC m=+15854707.285923840.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.