

cis-Bicyclo[3.1.0]hexane

Inchi:	InChI=1S/C6H10/c1-2-5-4-6(5)3-1/h5-6H,1-4H2/t5-,6+
InchiKey:	JAPMJSVZDUYFKL-OLQVQODUSA-N
Formula:	C6H10
SMILES:	C1CC2CC2C1
Mol. weight [g/mol]:	82.14
CAS:	285-58-5

Physical Properties

Property code	Value	Unit	Source
gf	121.14	kJ/mol	Joback Method
hf	-21.57	kJ/mol	Joback Method
hfus	7.57	kJ/mol	Joback Method
hvap	28.78	kJ/mol	Joback Method
log10ws	-1.64		Crippen Method
logp	1.806		Crippen Method
mvol	73.680	ml/mol	McGowan Method
pc	4249.61	kPa	Joback Method
rinpol	678.00		NIST Webbook
rinpol	678.00		NIST Webbook
tb	350.16	K	Joback Method
tc	545.42	K	Joback Method
tf	193.26	K	Joback Method
vc	0.285	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	123.56	J/mol×K	350.16	Joback Method
cpg	137.73	J/mol×K	382.70	Joback Method
cpg	150.96	J/mol×K	415.25	Joback Method
cpg	163.29	J/mol×K	447.79	Joback Method
cpg	174.79	J/mol×K	480.33	Joback Method
cpg	185.50	J/mol×K	512.88	Joback Method
cpg	195.48	J/mol×K	545.42	Joback Method

dvisc	0.0002505	Paxs	193.26	Joback Method
dvisc	0.0002758	Paxs	219.41	Joback Method
dvisc	0.0002976	Paxs	245.56	Joback Method
dvisc	0.0003163	Paxs	271.71	Joback Method
dvisc	0.0003327	Paxs	297.86	Joback Method
dvisc	0.0003471	Paxs	324.01	Joback Method
dvisc	0.0003599	Paxs	350.16	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.46227e+01
Coeff. B	-3.49263e+03
Coeff. C	-9.09100e+00
Temperature range (K), min.	252.73
Temperature range (K), max.	384.19

Sources

The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure
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=R294845&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

h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
m_{cvol}:	McGowan's characteristic volume
p_c:	Critical Pressure
p_{vap}:	Vapor pressure
r_{inpol}:	Non-polar retention indices
t_b:	Normal Boiling Point Temperature
t_c:	Critical Temperature
t_f:	Normal melting (fusion) point
v_c:	Critical Volume

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