1-Methyl-2-propylcyclohexane, trans

Other names:	Cyclohexane, 1-methyl-2-propyl-, trans- trans-1-Methyl-2-propylcyclohexane
Inchi:	InChI=1S/C10H20/c1-3-6-10-8-5-4-7-9(10)2/h9-10H,3-8H2,1-2H3/t9-,10-/m1/s1
InchiKey:	BVYJEKBXVYKYRA-NXEZZACHSA-N
Formula:	C10H20
SMILES:	CCCC1CCCCC1C
Mol. weight [g/mol]:	140.27
CAS:	42806-77-9

Physical Properties

Property code	Value	Unit	Source
gf	50.06	kJ/mol	Joback Method
hf	-215.75	kJ/mol	Joback Method
hfus	14.56	kJ/mol	Joback Method
hvap	37.97	kJ/mol	Joback Method
log10ws	-3.42		Crippen Method
logp	3.613		Crippen Method
mcvol	140.900	ml/mol	McGowan Method
pc	2485.07	kPa	Joback Method
rinpol	994.00		NIST Webbook
rinpol	993.40		NIST Webbook
rinpol	1016.00		NIST Webbook
rinpol	998.20		NIST Webbook
tb	443.08	K	Joback Method
tc	639.52	K	Joback Method
tf	205.60	K	Joback Method
vc	0.527	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	299.73	J/mol×K	443.08	Joback Method
cpg	389.89	J/mol×K	606.78	Joback Method
cpg	373.57	J/mol×K	574.04	Joback Method

cpg	356.42	J/mol×K	541.30	Joback Method
cpg	338.41	J/mol×K	508.56	Joback Method
cpg	319.51	J/mol×K	475.82	Joback Method
cpg	405.38	J/mol×K	639.52	Joback Method
dvisc	0.0002646	Pa×s	443.08	Joback Method
dvisc	0.0003397	Pa×s	403.50	Joback Method
dvisc	0.0004605	Pa×s	363.92	Joback Method
dvisc	0.0006724	Pa×s	324.34	Joback Method
dvisc	0.0010906	Pa×s	284.76	Joback Method
dvisc	0.0020679	Pa×s	245.18	Joback Method
dvisc	0.0050164	Pa×s	205.60	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=C42806779&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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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