

Tricyclohexylmethanol

Other names:	Cyclohexanemethanol, «alpha»,«alpha»-dicyclohexyl-
Inchi:	InChI=1S/C19H34O/c20-19(16-10-4-1-5-11-16,17-12-6-2-7-13-17)18-14-8-3-9-15-18/h16
InchiKey:	MRGVKQSSSOCONZ-UHFFFAOYSA-N
Formula:	C19H34O
SMILES:	OC(C1CCCCC1)(C1CCCCC1)C1CCCCC1
Mol. weight [g/mol]:	278.47
CAS:	17687-74-0

Physical Properties

Property code	Value	Unit	Source
gf	48.47	kJ/mol	Joback Method
hf	-433.51	kJ/mol	Joback Method
hfus	17.14	kJ/mol	Joback Method
hvap	74.56	kJ/mol	Joback Method
log10ws	-6.11		Crippen Method
logp	5.458		Crippen Method
mcvol	251.860	ml/mol	McGowan Method
pc	1864.33	kPa	Joback Method
tb	781.72	K	Joback Method
tc	1014.51	K	Joback Method
tf	366.00 ± 2.00	K	NIST Webbook
vc	0.906	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	861.17	J/mol×K	781.72	Joback Method
cpg	964.89	J/mol×K	975.71	Joback Method
cpg	947.63	J/mol×K	936.91	Joback Method
cpg	928.74	J/mol×K	898.12	Joback Method
cpg	908.11	J/mol×K	859.32	Joback Method
cpg	885.62	J/mol×K	820.52	Joback Method
cpg	980.64	J/mol×K	1014.51	Joback Method
dvisc	0.0000190	Paxs	781.72	Joback Method

dvisc	0.0000319	Paxs	716.31	Joback Method
dvisc	0.0000595	Paxs	650.90	Joback Method
dvisc	0.0001273	Paxs	585.50	Joback Method
dvisc	0.0003297	Paxs	520.09	Joback Method
dvisc	0.0011236	Paxs	454.68	Joback Method
dvisc	0.0057811	Paxs	389.27	Joback Method

Sources

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=C17687740&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cp_g:	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:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
m_cvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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