Cyclohexane, 1,1'-dodecylidenebis[4-methyl-

Other names:	1,1-Di(4-methylcyclohexyl)dodecane
	Cyclohexane, 1,1'-dodecylidenebis*4-methyl-
	1,1-bis(4-methylcyclohexyl)dodecane
Inchi:	InChI=1S/C26H50/c1-4-5-6-7-8-9-10-11-12-13-26(24-18-14-22(2)15-19-24)25-20-16-23(3
InchiKey:	ARDIJKGJYOGVEO-UHFFFAOYSA-N
Formula:	C26H50
SMILES:	CCCCCCCCCCCC(C1CCC(C)CC1)C1CCC(C)CC1
Mol. weight [g/mol]:	362.68
CAS:	55334-09-3

Physical Properties

Property code	Value	Unit	Source
gf	199.08	kJ/mol	Joback Method
hf	-517.29	kJ/mol	Joback Method
hfus	45.38	kJ/mol	Joback Method
hvap	73.32	kJ/mol	Joback Method
log10ws	-9.29		Crippen Method
logp	9.176		Crippen Method
mcvol	355.480	ml/mol	McGowan Method
pc	884.72	kPa	Joback Method
tb	823.60	K	Joback Method
tc	1022.38	K	Joback Method
tf	238.00 ± 3.00	K	NIST Webbook
vc	1.349	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1356.96	J/mol×K	1022.38	Joback Method
cpg	1223.41	J/mol×K	823.60	Joback Method
cpg	1249.73	J/mol×K	856.73	Joback Method
cpg	1274.35	J/mol×K	889.86	Joback Method
cpg	1297.33	J/mol×K	922.99	Joback Method
cpg	1318.72	J/mol×K	956.12	Joback Method

cpg	1338.58	J/mol×K	989.25	Joback Method
dvisc	0.0000710	Pa×s	823.60	Joback Method
dvisc	0.0029307	Pa×s	374.06	Joback Method
dvisc	0.0009397	Pa×s	448.98	Joback Method
dvisc	0.0004172	Pa×s	523.91	Joback Method
dvisc	0.0002269	Pa×s	598.83	Joback Method
dvisc	0.0001413	Pa×s	673.75	Joback Method
dvisc	0.0000968	Pa×s	748.68	Joback Method
hvapt	93.50	kJ/mol	502.00	NIST Webbook

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=C55334093&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
hvapt:	Enthalpy of vaporization at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	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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