

4,7-Methano-1H-indene, octahydro-5-(2-octyldecyl)-

Other names:	9-(5-exo-hexahydro-4,7-methanoindanyl)methyl)heptadecane
Inchi:	InChI=1S/C28H52/c1-3-5-7-9-11-13-16-23(17-14-12-10-8-6-4-2)20-24-21-25-22-28(24)2
InchiKey:	ZXUJNRWUNKDQTG-UHFFFAOYSA-N
Formula:	C28H52
SMILES:	CCCCCCCC(CCCCCC)CC1CC2CC1C1CCCC21
Mol. weight [g/mol]:	388.71
CAS:	55334-74-2

Physical Properties

Property code	Value	Unit	Source
gf	337.17	kJ/mol	Joback Method
hf	-454.97	kJ/mol	Joback Method
hfus	59.20	kJ/mol	Joback Method
hvap	76.83	kJ/mol	Joback Method
log10ws	-9.78		Crippen Method
logp	9.566		Crippen Method
mcvol	372.800	ml/mol	McGowan Method
pc	793.94	kPa	Joback Method
tb	854.75	K	Joback Method
tc	1048.87	K	Joback Method
tf	432.14	K	Joback Method
vc	1.458	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1314.51	J/molxK	854.75	Joback Method
cpg	1429.31	J/molxK	1016.52	Joback Method
cpg	1408.37	J/molxK	984.17	Joback Method
cpg	1386.54	J/molxK	951.81	Joback Method
cpg	1363.69	J/molxK	919.46	Joback Method
cpg	1339.71	J/molxK	887.10	Joback Method
cpg	1449.44	J/molxK	1048.87	Joback Method
dvisc	0.0010504	Paxs	854.75	Joback Method

dvisc	0.0011859	Paxs	784.32	Joback Method
dvisc	0.0013714	Paxs	713.88	Joback Method
dvisc	0.0016372	Paxs	643.44	Joback Method
dvisc	0.0020415	Paxs	573.01	Joback Method
dvisc	0.0027080	Paxs	502.57	Joback Method
dvisc	0.0039387	Paxs	432.14	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=C55334742&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	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
mc_{vol}:	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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