

Naphthalene, 6-(1,1-dimethylethyl)-1,2,3,4-tetrahydro-

Other names:	6-tert-Butyltetralin 6-tert-Butyl-(1,2,3,4-tetrahydronaphthalene)
Inchi:	InChI=1S/C14H20/c1-14(2,3)13-9-8-11-6-4-5-7-12(11)10-13/h8-10H,4-7H2,1-3H3
InchiKey:	XCYVLZXWLBNID-UHFFFAOYSA-N
Formula:	C14H20
SMILES:	CC(C)(C)c1ccc2c(c1)CCCC2
Mol. weight [g/mol]:	188.31
CAS:	42044-26-8

Physical Properties

Property code	Value	Unit	Source
gf	219.35	kJ/mol	Joback Method
hf	-40.47	kJ/mol	Joback Method
hfus	12.83	kJ/mol	Joback Method
hvap	49.46	kJ/mol	Joback Method
log10ws	-4.29		Crippen Method
logp	3.863		Crippen Method
mvol	173.500	ml/mol	McGowan Method
pc	2370.28	kPa	Joback Method
tb	568.81	K	Joback Method
tc	800.54	K	Joback Method
tf	320.08	K	Joback Method
vc	0.650	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	432.38	J/molxK	568.81	Joback Method
cpg	452.59	J/molxK	607.43	Joback Method
cpg	471.33	J/molxK	646.05	Joback Method
cpg	488.73	J/molxK	684.68	Joback Method
cpg	504.87	J/molxK	723.30	Joback Method
cpg	519.88	J/molxK	761.92	Joback Method
cpg	533.84	J/molxK	800.54	Joback Method

dvisc	0.0026770	Paxs	320.08	Joback Method
dvisc	0.0014325	Paxs	361.54	Joback Method
dvisc	0.0008718	Paxs	402.99	Joback Method
dvisc	0.0005821	Paxs	444.44	Joback Method
dvisc	0.0004164	Paxs	485.90	Joback Method
dvisc	0.0003139	Paxs	527.36	Joback Method
dvisc	0.0002467	Paxs	568.81	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C42044268&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
log₁₀ws:	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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