

Naphthalene, 1,2,3,4-tetrahydro-1,4,6-trimethyl-

Other names:	1,2,3,4-tetrahydro-1,4,6-trimethylnaphthalene
Inchi:	InChI=1S/C13H18/c1-9-4-7-12-10(2)5-6-11(3)13(12)8-9/h4,7-8,10-11H,5-6H2,1-3H3
InchiKey:	HIACZWPSGFMDRF-UHFFFAOYSA-N
Formula:	C13H18
SMILES:	Cc1ccc2c(c1)C(C)CCC2C
Mol. weight [g/mol]:	174.28
CAS:	22824-32-4

Physical Properties

Property code	Value	Unit	Source
gf	192.67	kJ/mol	Joback Method
hf	-51.76	kJ/mol	Joback Method
hfus	19.79	kJ/mol	Joback Method
hvap	47.91	kJ/mol	Joback Method
log10ws	-4.21		Crippen Method
logp	3.996		Crippen Method
mcvol	159.410	ml/mol	McGowan Method
pc	2395.87	kPa	Joback Method
tb	539.82	K	Joback Method
tc	761.35	K	Joback Method
tf	297.91	K	Joback Method
vc	0.604	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	381.81	J/molxK	539.82	Joback Method
cpg	467.80	J/molxK	724.43	Joback Method
cpg	452.70	J/molxK	687.51	Joback Method
cpg	436.60	J/molxK	650.59	Joback Method
cpg	419.45	J/molxK	613.66	Joback Method
cpg	401.20	J/molxK	576.74	Joback Method
cpg	481.94	J/molxK	761.35	Joback Method
dvisc	0.0003468	Paxs	539.82	Joback Method

dvisc	0.0003976	Paxs	499.50	Joback Method
dvisc	0.0004671	Paxs	459.18	Joback Method
dvisc	0.0005658	Paxs	418.86	Joback Method
dvisc	0.0007141	Paxs	378.55	Joback Method
dvisc	0.0009527	Paxs	338.23	Joback Method
dvisc	0.0013740	Paxs	297.91	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=C22824324&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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
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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