

1,4,6,7-tetramethylnaphthalene

Inchi:	InChI=1S/C14H16/c1-9-5-6-10(2)14-8-12(4)11(3)7-13(9)14/h5-8H,1-4H3
InchiKey:	VPSPONOBZCLIU-UHFFFAOYSA-N
Formula:	C14H16
SMILES:	Cc1cc2c(C)ccc(C)c2cc1C
Mol. weight [g/mol]:	184.28
CAS:	13764-18-6

Physical Properties

Property code	Value	Unit	Source
af	0.5131		Cheméo Relay (1.0)
gf	247.54	kJ/mol	Joback Method
hf	35.53	kJ/mol	Cheméo Relay (1.0)
hfus	21.52	kJ/mol	Joback Method
hvap	76.17	kJ/mol	Cheméo Relay (1.0)
ie	7.47	eV	Cheméo Relay (1.0)
log10ws	-5.63		Cheméo Relay (1.0)
logp	4.073		Crippen Method
mcvol	164.900	ml/mol	McGowan Method
pc	2388.85	kPa	Joback Method
tb	571.41	K	Cheméo Relay (1.0)
tc	802.12	K	Cheméo Relay (1.0)
tf	401.82	K	Cheméo Relay (1.0)
vc	0.624	m3/kmol	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	389.23	J/mol×K	585.30	Joback Method
cpg	470.88	J/mol×K	810.38	Joback Method
cpg	459.28	J/mol×K	772.87	Joback Method
cpg	446.95	J/mol×K	735.36	Joback Method
cpg	433.83	J/mol×K	697.84	Joback Method
cpg	419.87	J/mol×K	660.33	Joback Method
cpg	405.02	J/mol×K	622.81	Joback Method

dvisc	0.0004369	Pa×s	471.02	Joback Method
dvisc	0.0002702	Pa×s	585.30	Joback Method
dvisc	0.0003101	Pa×s	547.21	Joback Method
dvisc	0.0003634	Pa×s	509.11	Joback Method
dvisc	0.0009613	Pa×s	356.74	Joback Method
dvisc	0.0005425	Pa×s	432.93	Joback Method
dvisc	0.0007025	Pa×s	394.83	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.63870e+01
Coeff. B	-5.38420e+03
Coeff. C	-1.03895e+02
Temperature range (K), min.	438.33
Temperature range (K), max.	590.03

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C13764186&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

af:	Acentric Factor
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
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.cheméo.com/cid/35-610-5/1-4-6-7-tetramethylnaphthalene.pdf>

Generated by Cheméo on 2026-07-21 20:21:55.400636078 +0000 UTC m=+176411.242521453.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.