

Naphthalene, 1,2,4-trimethyl-

Other names:	1,2,4-Trimethylnaphthalene
Inchi:	InChI=1S/C13H14/c1-9-8-10(2)12-6-4-5-7-13(12)11(9)3/h4-8H,1-3H3
InchiKey:	JCNGSJUYPCVGAM-UHFFFAOYSA-N
Formula:	C13H14
SMILES:	<chem>Cc1cc(C)c2cccc2c1C</chem>
Mol. weight [g/mol]:	170.25
CAS:	2717-42-2

Physical Properties

Property code	Value	Unit	Source
gf	248.75	kJ/mol	Joback Method
hf	81.54	kJ/mol	Joback Method
hfus	19.32	kJ/mol	Joback Method
hvap	50.43	kJ/mol	Joback Method
log10ws	-4.70		Crippen Method
logp	3.765		Crippen Method
mcvol	150.810	ml/mol	McGowan Method
pc	2670.78	kPa	Joback Method
rinpol	268.29		NIST Webbook
rinpol	268.54		NIST Webbook
rinpol	269.30		NIST Webbook
rinpol	269.23		NIST Webbook
rinpol	268.50		NIST Webbook
rinpol	268.30		NIST Webbook
rinpol	269.23		NIST Webbook
rinpol	1575.00		NIST Webbook
rinpol	268.50		NIST Webbook
rinpol	1575.00		NIST Webbook
tb	557.44	K	Joback Method
tc	785.77	K	Joback Method
tf	328.00 ± 4.00	K	NIST Webbook
vc	0.578	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	342.00	J/molxK	557.44	Joback Method
cpg	409.52	J/molxK	747.71	Joback Method
cpg	397.73	J/molxK	709.66	Joback Method
cpg	385.15	J/molxK	671.60	Joback Method
cpg	371.70	J/molxK	633.55	Joback Method
cpg	357.34	J/molxK	595.49	Joback Method
cpg	420.56	J/molxK	785.77	Joback Method
dvisc	0.0002834	Paxs	557.44	Joback Method
dvisc	0.0003273	Paxs	520.03	Joback Method
dvisc	0.0003866	Paxs	482.61	Joback Method
dvisc	0.0004696	Paxs	445.20	Joback Method
dvisc	0.0005911	Paxs	407.78	Joback Method
dvisc	0.0007795	Paxs	370.37	Joback Method
dvisc	0.0010939	Paxs	332.95	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.50877e+01
Coeff. B	-4.43327e+03
Coeff. C	-1.19607e+02
Temperature range (K), min.	419.15
Temperature range (K), max.	573.08

Sources

The Yaws Handbook of Vapor

Pressure:

Crippen Method:

Crippen Method:

Joback Method:

McGowan Method:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

https://www.chemeo.com/doc/models/crippen_log10ws

https://en.wikipedia.org/wiki/Joback_method

<http://link.springer.com/article/10.1007/BF02311772>

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
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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