

Naphthalene, 1,2,3,4,6-pentachloro

Inchi:	InChI=1S/C10H3Cl5/c11-4-1-2-5-6(3-4)8(13)10(15)9(14)7(5)12/h1-3H
InchiKey:	BAOLNVSMVTYGDA-UHFFFAOYSA-N
Formula:	C10H3Cl5
SMILES:	Clc1ccc2c(Cl)c(Cl)c(Cl)c(Cl)c2c1
Mol. weight [g/mol]:	300.40

Physical Properties

Property code	Value	Unit	Source
gf	144.58	kJ/mol	Joback Method
hf	41.82	kJ/mol	Joback Method
hfus	31.76	kJ/mol	Joback Method
hvap	67.00	kJ/mol	Joback Method
log10ws	-6.70		Crippen Method
logp	6.107		Crippen Method
mcvol	169.740	ml/mol	McGowan Method
pc	2884.30	kPa	Joback Method
rinpol	2186.00		NIST Webbook
rinpol	2186.00		NIST Webbook
tb	685.91	K	Joback Method
tc	947.92	K	Joback Method
tf	473.78	K	Joback Method
vc	0.654	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	310.73	J/molxK	685.91	Joback Method
cpg	318.01	J/molxK	729.58	Joback Method
cpg	324.67	J/molxK	773.25	Joback Method
cpg	330.79	J/molxK	816.91	Joback Method
cpg	336.44	J/molxK	860.58	Joback Method
cpg	341.68	J/molxK	904.25	Joback Method
cpg	346.58	J/molxK	947.92	Joback Method
dvisc	0.0009715	Paxs	473.78	Joback Method

dvisc	0.0007697	Paxs	509.14	Joback Method
dvisc	0.0006286	Paxs	544.49	Joback Method
dvisc	0.0005262	Paxs	579.85	Joback Method
dvisc	0.0004495	Paxs	615.20	Joback Method
dvisc	0.0003907	Paxs	650.56	Joback Method
dvisc	0.0003445	Paxs	685.91	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R128446&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/77-010-5/Naphthalene-1-2-3-4-6-pentachloro.pdf>

Generated by Cheméo on 2024-04-24 22:02:18.905304984 +0000 UTC m=+16285387.825882300.

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