

Naphthalene, 1,7-diphenyl-

Other names:	1,7-Diphenyl-naphthalene
Inchi:	InChI=1S/C22H16/c1-3-8-17(9-4-1)20-15-14-19-12-7-13-21(22(19)16-20)18-10-5-2-6-11
InchiKey:	GQHLUGQTYDTQDG-UHFFFAOYSA-N
Formula:	C22H16
SMILES:	<chem>c1ccc(-c2ccc3ccccc(-c4ccccc4)c3c2)cc1</chem>
Mol. weight [g/mol]:	280.36
CAS:	970-06-9

Physical Properties

Property code	Value	Unit	Source
gf	558.98	kJ/mol	Joback Method
hf	380.31	kJ/mol	Joback Method
hfus	31.10	kJ/mol	Joback Method
hvap	74.36	kJ/mol	Joback Method
log10ws	-8.48		Crippen Method
logp	6.174		Crippen Method
mcvol	230.100	ml/mol	McGowan Method
pc	2220.80	kPa	Joback Method
tb	811.74	K	Joback Method
tc	1087.66	K	Joback Method
tf	474.70	K	Joback Method
vc	0.866	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	641.58	J/molxK	811.74	Joback Method
cpg	712.88	J/molxK	1041.68	Joback Method
cpg	700.72	J/molxK	995.69	Joback Method
cpg	687.72	J/molxK	949.70	Joback Method
cpg	673.67	J/molxK	903.71	Joback Method
cpg	658.36	J/molxK	857.73	Joback Method
cpg	724.41	J/molxK	1087.66	Joback Method
dvisc	0.0001568	Paxs	811.74	Joback Method

dvisc	0.0001897	Paxs	755.57	Joback Method
dvisc	0.0002367	Paxs	699.39	Joback Method
dvisc	0.0003069	Paxs	643.22	Joback Method
dvisc	0.0004184	Paxs	587.05	Joback Method
dvisc	0.0006089	Paxs	530.87	Joback Method
dvisc	0.0009686	Paxs	474.70	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=C970069&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
log₁₀w_s:	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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