

Dinaphtho[1,2-g:1',2'-g']naphtho[2,1-c:7,8-c']diphenyl

Inchi:	InChI=1S/C50H28/c1-3-7-40-29(5-1)9-11-31-13-15-33-17-19-35-21-23-37-25-27-39-28-2
InchiKey:	WRIBLYWAWHZTKJ-UHFFFAOYSA-N
Formula:	C50H28
SMILES:	c1ccc2c(c1)ccc1ccc3ccc4ccc5ccc6ccc7ccc8ccc9ccc%10ccc%11cccc%11c%10c9c8c7c
Mol. weight [g/mol]:	628.76
CAS:	57468-46-9

Physical Properties

Property code	Value	Unit	Source
gf	1559.38	kJ/mol	Joback Method
hf	1148.27	kJ/mol	Joback Method
hfus	82.62	kJ/mol	Joback Method
hvap	153.83	kJ/mol	Joback Method
ie	6.93	eV	NIST Webbook
log10ws	-21.32		Crippen Method
logp	14.372		Crippen Method
mcvol	477.540	ml/mol	McGowan Method
pc	1012.95	kPa	Joback Method
tb	1628.66	K	Joback Method
tc	2003.27	K	Joback Method
tf	1164.58	K	Joback Method
vc	1.869	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	2264.14	J/molxK	1628.66	Joback Method
cpg	2424.65	J/molxK	1691.10	Joback Method
cpg	2606.94	J/molxK	1753.53	Joback Method
cpg	2812.83	J/molxK	1815.97	Joback Method
cpg	3044.15	J/molxK	1878.40	Joback Method
cpg	3302.73	J/molxK	1940.84	Joback Method
cpg	3590.40	J/molxK	2003.27	Joback Method
dvisc	0.0382713	Paxs	1164.58	Joback Method

dvisc	0.0357658	Paxs	1241.93	Joback Method
dvisc	0.0336907	Paxs	1319.27	Joback Method
dvisc	0.0319468	Paxs	1396.62	Joback Method
dvisc	0.0304627	Paxs	1473.97	Joback Method
dvisc	0.0291856	Paxs	1551.31	Joback Method
dvisc	0.0280760	Paxs	1628.66	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=C57468469&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
ie:	Ionization energy
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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