

Dinaphtho[2,1,8,7-defg!2',1',8',7'-qrst]pentacene

Other names:	Dinaphtho[2,1,8,7-defg:2',1',8',7'-qrst]pentacene
Inchi:	InChI=1S/C32H16/c1-2-6-21-20(4-1)16-26-25-15-12-18-9-8-17-5-3-7-22-24-14-11-19-10
InchiKey:	FIVBPVSFZMYBBN-UHFFFAOYSA-N
Formula:	C32H16
SMILES:	c1ccc2c(c1)cc1c3ccc4ccc5cccc6c7ccc8ccc2c1c8c7c3c4c56
Mol. weight [g/mol]:	400.47
CAS:	120835-94-1

Physical Properties

Property code	Value	Unit	Source
gf	1099.48	kJ/mol	Joback Method
hf	844.61	kJ/mol	Joback Method
hfus	55.04	kJ/mol	Joback Method
hvap	104.95	kJ/mol	Joback Method
log10ws	-14.15		Crippen Method
logp	9.226		Crippen Method
mcvol	295.200	ml/mol	McGowan Method
pc	1792.42	kPa	Joback Method
tb	1121.84	K	Joback Method
tc	1399.15	K	Joback Method
tf	844.90	K	Joback Method
vc	1.185	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	953.39	J/molxK	1121.84	Joback Method
cpg	984.02	J/molxK	1168.06	Joback Method
cpg	1018.68	J/molxK	1214.28	Joback Method
cpg	1057.97	J/molxK	1260.50	Joback Method
cpg	1102.53	J/molxK	1306.71	Joback Method
cpg	1152.95	J/molxK	1352.93	Joback Method
cpg	1209.85	J/molxK	1399.15	Joback Method
dvisc	0.1009143	Paxs	844.90	Joback Method

dvisc	0.1032630	Paxs	891.06	Joback Method
dvisc	0.1054271	Paxs	937.21	Joback Method
dvisc	0.1074272	Paxs	983.37	Joback Method
dvisc	0.1092809	Paxs	1029.53	Joback Method
dvisc	0.1110035	Paxs	1075.68	Joback Method
dvisc	0.1126083	Paxs	1121.84	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C120835941&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cp_g:	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:	Log ₁₀ of Water solubility in mol/l
logp:	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

Latest version available from:

<https://www.chemeo.com/cid/77-213-0/Dinaphtho-2-1-8-7-defg-2-1-8-7-qrst-pentacene.pdf>

Generated by Cheméo on 2024-05-03 19:36:48.673776725 +0000 UTC m=+17054257.594354046.

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