Benzo[2,1-a!3,4-a']dianthracene

Other names: Benzo[2,1-a:3,4-a']dianthracene
Inchi: InChI=1S/C30H18/c1-3-7-23-17-27-25(15-21(23)5-1)13-11-19-9-10-20-12-14-26-16-22-6
InchiKey: IYYZTTJANRCFKN-UHFFFAOYSA-N
Formula: C30H18
SMILES: c1ccc2cc3c(ccc4ccc5ccc6cc7ccccc7cc6c5c43)cc2c1
Mol. weight [g/mol]: 378.46
CAS: 188-51-2

Physical Properties

Property code	Value	Unit	Source
gf	905.88	kJ/mol	Joback Method
hf	663.07	kJ/mol	Joback Method
hfus	47.67	kJ/mol	Joback Method
hvap	97.80	kJ/mol	Joback Method
log10ws	-12.30		Crippen Method
logp	8.606		Crippen Method
mcvol	293.040	ml/mol	McGowan Method
pc	1777.34	kPa	Joback Method
tb	1051.26	K	Joback Method
tc	1330.47	K	Joback Method
tf	713.08	K	Joback Method
vc	1.139	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	909.76	J/mol×K	1051.26	Joback Method
cpg	929.78	J/mol×K	1097.80	Joback Method
cpg	951.06	J/mol×K	1144.33	Joback Method
cpg	974.05	J/mol×K	1190.87	Joback Method
cpg	999.18	J/mol×K	1237.40	Joback Method
cpg	1026.90	J/mol×K	1283.94	Joback Method
cpg	1057.66	J/mol×K	1330.47	Joback Method
dvisc	0.0046199	Pa×s	713.08	Joback Method

dvisc	0.0041148	Pa×s	769.44	Joback Method
dvisc	0.0037234	Pa×s	825.81	Joback Method
dvisc	0.0034125	Pa×s	882.17	Joback Method
dvisc	0.0031604	Pa×s	938.53	Joback Method
dvisc	0.0029526	Pa×s	994.90	Joback Method
dvisc	0.0027786	Pa×s	1051.26	Joback Method

Sources

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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C188512&Units=SI

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume