

4H-Benz[de]anthracene, 5,6-dihydro-

Other names:	5,6-Dihydro-4H-benz[de]anthracene
Inchi:	InChI=1S/C17H14/c1-2-9-15-13(5-1)11-14-8-3-6-12-7-4-10-16(15)17(12)14/h1-2,4-5,7,9-
InchiKey:	VFZDNSWGNPSQLY-UHFFFAOYSA-N
Formula:	C17H14
SMILES:	<chem>c1ccc2c(c1)cc1c3c(cccc32)CCC1</chem>
Mol. weight [g/mol]:	218.29
CAS:	4389-09-7

Physical Properties

Property code	Value	Unit	Source
gf	457.54	kJ/mol	Joback Method
hf	283.19	kJ/mol	Joback Method
hfus	23.76	kJ/mol	Joback Method
hvap	61.20	kJ/mol	Joback Method
log10ws	-6.16		Crippen Method
logp	4.482		Crippen Method
mcvol	176.850	ml/mol	McGowan Method
pc	2738.28	kPa	Joback Method
tb	679.35	K	Joback Method
tc	935.56	K	Joback Method
tf	353.60 ± 1.50	K	NIST Webbook
vc	0.681	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	462.24	J/mol×K	679.35	Joback Method
cpg	478.18	J/mol×K	722.05	Joback Method
cpg	492.87	J/mol×K	764.75	Joback Method
cpg	506.49	J/mol×K	807.45	Joback Method
cpg	519.26	J/mol×K	850.15	Joback Method
cpg	531.38	J/mol×K	892.86	Joback Method
cpg	543.05	J/mol×K	935.56	Joback Method
dvisc	0.0020061	Paxs	432.91	Joback Method

dvisc	0.0016661	Paxs	473.98	Joback Method
dvisc	0.0014254	Paxs	515.06	Joback Method
dvisc	0.0012479	Paxs	556.13	Joback Method
dvisc	0.0011126	Paxs	597.20	Joback Method
dvisc	0.0010068	Paxs	638.28	Joback Method
dvisc	0.0009221	Paxs	679.35	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C4389097&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
h_{vap}:	Enthalpy of vaporization at standard conditions
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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