1,1'-Biphenyl, 2,3',4',5-tetrachloro-

Other names:	1,2-dichloro-4-(2,5-dichlorophenyl)benzene
	2,3',4',5-PCB
	2,3',4',5-Tetrachloro-1,1'-biphenyl
	2,3',4',5-Tetrachlorobiphenyl
	2,5,3',4'-Tetrachlorobiphenyl
	Biphenyl, 2,3',4',5-tetrachloro-
	PCB-70
Inchi:	InChI=1S/C12H6Cl4/c13-8-2-4-10(14)9(6-8)7-1-3-11(15)12(16)5-7/h1-6H
InchiKey:	KENZYIHFBRWMOD-UHFFFAOYSA-N
Formula:	C12H6Cl4
SMILES:	Clc1ccc(Cl)c(-c2ccc(Cl)c(Cl)c2)c1
Mol. weight [g/mol]:	291.99
CAS:	32598-11-1

Physical Properties

Property code	Value	Unit	Source
gf	188.74	kJ/mol	Joback Method
hf	73.21	kJ/mol	Joback Method
hfus	30.15	kJ/mol	Joback Method
hvap	67.05	kJ/mol	Joback Method
log10ws	-7.25		Aqueous Solubility Prediction Method
log10ws	-7.25		Estimated Solubility Method
logp	5.967		Crippen Method
mcvol	181.380	ml/mol	McGowan Method
pc	2724.01	kPa	Joback Method
rinpol	2020.60		NIST Webbook
rinpol	1989.00		NIST Webbook
rinpol	2009.90		NIST Webbook
rinpol	2019.00		NIST Webbook
rinpol	2031.60		NIST Webbook
rinpol	2031.30		NIST Webbook
rinpol	2019.00		NIST Webbook
rinpol	1988.00		NIST Webbook
rinpol	2000.00		NIST Webbook
rinpol	2031.30		NIST Webbook

rinpol	1989.00		NIST Webbook
tb	696.96	K	Joback Method
tc	963.35	K	Joback Method
tf	447.60	K	Joback Method
vc	0.688	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	417.60	J/mol×K	963.35	Joback Method
cpg	411.33	J/mol×K	918.95	Joback Method
cpg	404.36	J/mol×K	874.55	Joback Method
cpg	396.63	J/mol×K	830.15	Joback Method
cpg	388.08	J/mol×K	785.76	Joback Method
cpg	378.65	J/mol×K	741.36	Joback Method
cpg	368.27	J/mol×K	696.96	Joback Method
dvisc	0.0008409	Pa×s	447.60	Joback Method
dvisc	0.0001717	Pa×s	696.96	Joback Method
dvisc	0.0002057	Pa×s	655.40	Joback Method
dvisc	0.0002526	Pa×s	613.84	Joback Method
dvisc	0.0003196	Pa×s	572.28	Joback Method
dvisc	0.0004195	Pa×s	530.72	Joback Method
dvisc	0.0005766	Pa×s	489.16	Joback Method
hvapt	84.80	kJ/mol	370.50	NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C32598111&Units=SI

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Joback Method: https://en.wikipedia.org/wiki/Joback_method

Aqueous Solubility Prediction Method: http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa

Estimated Solubility Method: http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

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
hvapt:	Enthalpy of vaporization at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume