

1,1'-Biphenyl, 2,3',4,4',6-pentachloro-

Other names:	2,3',4,4',6-Pentachlorobiphenyl PCB 119
Inchi:	InChI=1S/C12H5Cl5/c13-7-4-10(16)12(11(17)5-7)6-1-2-8(14)9(15)3-6/h1-5H
InchiKey:	OAEQTHQGPZKTQP-UHFFFAOYSA-N
Formula:	C12H5Cl5
SMILES:	Clc1cc(Cl)c(-c2ccc(Cl)c(Cl)c2)c(Cl)c1
Mol. weight [g/mol]:	326.44

Physical Properties

Property code	Value	Unit	Source
hf	73.79	kJ/mol	Cheméo Relay (1.0)
hfus	33.96	kJ/mol	Joback Method
hvap	95.38	kJ/mol	Cheméo Relay (1.0)
ie	8.52	eV	Cheméo Relay (1.0)
log10ws	-7.59		Cheméo Relay (1.0)
logp	6.621		Crippen Method
mcvol	193.620	ml/mol	McGowan Method
rinpol	2182.00		NIST Webbook
rinpol	2185.00		NIST Webbook
rinpol	2182.00		NIST Webbook
rinpol	2185.00		NIST Webbook
rinpol	2182.00		NIST Webbook
rinpol	2182.00		NIST Webbook
tb	609.04	K	Cheméo Relay (1.0)
tf	365.45	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	385.50	J/mol×K	739.37	Joback Method
cpg	394.64	J/mol×K	784.18	Joback Method
cpg	402.91	J/mol×K	828.99	Joback Method
cpg	410.38	J/mol×K	873.79	Joback Method
cpg	417.08	J/mol×K	918.60	Joback Method

cpg	423.09	J/mol×K	963.41	Joback Method
cpg	428.46	J/mol×K	1008.22	Joback Method
dvisc	0.0006694	Pa×s	490.04	Joback Method
dvisc	0.0004783	Pa×s	531.60	Joback Method
dvisc	0.0003588	Pa×s	573.15	Joback Method
dvisc	0.0002798	Pa×s	614.71	Joback Method
dvisc	0.0002252	Pa×s	656.26	Joback Method
dvisc	0.0001860	Pa×s	697.82	Joback Method
dvisc	0.0001570	Pa×s	739.37	Joback Method

Sources

Cheméo Relay (1.0, beta):

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C56558179&Units=SI>

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

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

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

Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws

Cheméo Relay (1.0):

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tf:	Normal melting (fusion) point

Latest version available from:

<https://www.cheméo.com/cid/39-514-8/1-1-Biphenyl-2-3-4-4-6-pentachloro.pdf>

Generated by Cheméo on 2026-07-21 12:52:08.665566944 +0000 UTC m=+149424.507452374.

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