

Dibenzofuran, 1,4,7,8-tetrachloro

Other names:	1,4,7,8-tetrachlorodibenzofuran 2,3,6,9-Tetrachlorodibenzofuran
Inchi:	InChI=1S/C12H4Cl4O/c13-6-1-2-7(14)12-11(6)5-3-8(15)9(16)4-10(5)17-12/h1-4H
InchiKey:	IVRDRRWQABCXLY-UHFFFAOYSA-N
Formula:	C12H4Cl4O
SMILES:	Clc1cc2oc3c(Cl)ccc(Cl)c3c2cc1Cl
Mol. weight [g/mol]:	305.97

Physical Properties

Property code	Value	Unit	Source
log10ws	-11.55		Crippen Method
logp	6.200		Crippen Method
mcvol	176.390	ml/mol	McGowan Method
rinpol	2338.00		NIST Webbook
rinpol	2338.00		NIST Webbook
rinpol	2290.00		NIST Webbook
rinpol	2290.00		NIST Webbook
rinpol	2290.00		NIST Webbook
rinpol	2290.00		NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R29613&Units=SI

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient

mcvol: McGowan's characteristic volume

rinpol: Non-polar retention indices

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