

trans-Chlorfenvinphos

Other names:

Phosphoric acid, 2-chloro-1-(2,4-dichlorophenyl)ethenyl diethyl ester, (E)-
Phosphoric acid, 2-chloro-1-(2,4-dichlorophenyl)vinyl diethyl ester, (E)-
trans-Isomer of chlorfenvinphos
Chlorfenvinphos (trans)
(E)-Chlorfenvinfos
2-Chloro-1-(2,4-dichlorophenyl)ethenyl diethyl phosphate, (E)-
«alpha»-Chlorfenvinphos
(E)-2-Chloro-1-(2,4-dichlorophenyl)ethenyl diethyl phosphate
Chlorfenvinphos E
Chlorfenvinphos (E), alpha
(E)-Chlorfenvinphos

Inchi: InChI=1S/C12H14Cl3O4P/c1-3-17-20(16,18-4-2)19-12(8-13)10-6-5-9(14)7-11(10)15/h5-8**InchiKey:** FSAVDKDHPDSTO-XYOKQWHBSA-N**Formula:** C12H14Cl3O4P**SMILES:** CCOP(=O)(OCC)OC(=CCl)c1ccc(Cl)cc1Cl**Mol. weight [g/mol]:** 359.57**CAS:** 18708-86-6

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.01		Crippen Method
logp	5.728		Crippen Method
mcvol	232.540	ml/mol	McGowan Method
rinpol	2046.00		NIST Webbook
rinpol	2048.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C18708866&Units=SI>**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>**Crippen Method:** https://www.chemeo.com/doc/models/crippen_log10ws**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.cheméo.com/cid/66-898-3/trans-Chlorfenvinphos.pdf>

Generated by Cheméo on 2024-05-06 17:13:30.005420276 +0000 UTC m=+17304858.925997591.

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