

Phosphoric acid, 2-bromo-1-(2,4-dichlorophenyl)vinyl diethyl ester

Other names:
ester

Bromfenvinfos
Bromfenvinphos
Bromfenwinfos
Bromphenvinphos
O-1-(2,4-Dichlorophenyl)-2-bromovinyl O,O-diethyl phosphate
O,O-Dwuetylo-O-1-(2,4-dwuchlorofenylo)-2-bromowinylofosforan
IPO 62
lpofos
lpophos
Phosphoric acid, 2-bromo-1-(2,4-dichlorophenyl)ethenyl diethyl ester
SD 8989
[2-Bromo-1-(2,4-dichlorophenyl)vinyl] diethyl phosphate
InChI=1S/C12H14BrCl2O4P/c1-3-17-20(16,18-4-2)19-12(8-13)10-6-5-9(14)7-11(10)15/h
ORDKAVSHIKNMAN-XYOKQWHBSA-N
C12H14BrCl2O4P
CCOP(=O)(OCC)OC(=CBr)c1ccc(Cl)cc1Cl
Mol. weight [g/mol]: 404.02
CAS: 33399-00-7

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.29		Crippen Method
logp	5.884		Crippen Method
mcvol	237.800	ml/mol	McGowan Method
rmpol	2089.00		NIST Webbook

Sources

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>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C33399007&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

Latest version available from:

<https://www.cheméo.com/cid/32-327-3/Phosphoric-acid-2-bromo-1-2-4-dichlorophenyl-vinyl-diethyl-ester.pdf>

Generated by Cheméo on 2024-05-06 21:30:07.44173385 +0000 UTC m=+17320256.362311162.

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