

Triazophos

Other names:

Phosphorothioic acid, O,O-diethyl O-(1-phenyl-1H-1,2,4-triazol-3-yl) ester
Hostathion
Hostation
HOE 2960
Triazofos
Phosphorothioic acid, O,O-diethyl O-(1-phenyl-1,2,4-triazolyl) ester
O,O-Diethyl O-(1-phenyl-1H-1,2,4-triazol-3-yl)phosphorothioate
HOE 2960 OJ
1-Phenyl-3-(O,O-diethyl-thionophosphoryl)-1,2,4-triazole
1-Phenyl-1,2,4-triazolyl-3-(O,O-diethylthionophosphate)
Triazofosz
Trazophos
Thiasophos

Inchi: InChI=1S/C12H16N3O3PS/c1-3-16-19(20,17-4-2)18-12-13-10-15(14-12)11-8-6-5-7-9-11**InchiKey:** AMFGTOFWMRQMEM-UHFFFAOYSA-N**Formula:** C12H16N3O3PS**SMILES:** CCOP(=S)(OCC)Oc1ncn(-c2ccccc2)n1**Mol. weight [g/mol]:** 313.31**CAS:** 24017-47-8

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.46		Crippen Method
logp	2.944		Crippen Method
mcvol	221.080	ml/mol	McGowan Method
rinpol	2299.00		NIST Webbook
rinpol	2309.00		NIST Webbook
rinpol	2310.00		NIST Webbook
rinpol	2310.00		NIST Webbook
rinpol	2322.20		NIST Webbook
rinpol	2310.00		NIST Webbook
rinpol	2299.00		NIST Webbook
rinpol	2322.20		NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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=C24017478&Units=SI

Legend

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

Latest version available from:

<https://www.chemeo.com/cid/114-459-6/Triazophos.pdf>

Generated by Cheméo on 2024-04-28 03:50:21.624083393 +0000 UTC m=+16565470.544660707.

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