

# Folpet

<b>Other names:</b>	1H-Isoindole-1,3(2H)-dione, 2-[(trichloromethyl)thio]- Phthalimide, N-[(trichloromethyl)thio]- Faltan Ftalan Fungitrol 11 N-[(Trichloromethyl)thio]phthalimide Orthophaltan Phaltan Phaltane Phthaltan Spolacid 2-[(Trichloromethyl)thio]-1H-isoindole-1,3(2H)-dione Folpan Folpel Folpex N-(Trichloromethylmercapto)phthalimide Trichloromethyl(thio)phthalimide Acryptan Folnit Intercide TMP Vinicoll
<b>Inchi:</b>	InChI=1S/C9H4Cl3NO2S/c10-9(11,12)16-13-7(14)5-3-1-2-4-6(5)8(13)15/h1-4H
<b>InchiKey:</b>	HKIOYBQGHSTUDB-UHFFFAOYSA-N
<b>Formula:</b>	C9H4Cl3NO2S
<b>SMILES:</b>	O=C1c2ccccc2C(=O)N1SC(Cl)(Cl)Cl
<b>Mol. weight [g/mol]:</b>	296.56
<b>CAS:</b>	133-07-3

## Physical Properties

Property code	Value	Unit	Source
log10ws	-4.76		Crippen Method
logp	3.259		Crippen Method
mcvol	169.240	ml/mol	McGowan Method
rinpol	2077.00		NIST Webbook
rinpol	2015.00		NIST Webbook
rinpol	2086.00		NIST Webbook

rropol	2086.00		NIST Webbook
rropol	2015.00		NIST Webbook
rropol	2015.00		NIST Webbook
rropol	2015.00		NIST Webbook
tf	454.52 ± 0.20	K	NIST Webbook
tf	454.10 ± 0.20	K	NIST Webbook

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hfust	35.49	kJ/mol	454.20	NIST Webbook

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C133073&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C133073&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>

## Legend

<b>hfust:</b>	Enthalpy of fusion at a given temperature
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>rropol:</b>	Non-polar retention indices
<b>tf:</b>	Normal melting (fusion) point

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