

Phosphonothioic difluoride

Other names:	Phosphine sulfide, difluoro- Hydrothiophosphoryl difluoride
Inchi:	InChI=1S/F2HPS/c1-3(2)4/h3H
InchiKey:	MVVUNUURHWXOFX-UHFFFAOYSA-N
Formula:	F2HPS
SMILES:	F[PH](F)=S
Mol. weight [g/mol]:	102.04
CAS:	13780-63-7

Physical Properties

Property code	Value	Unit	Source
log10ws	3.02		Crippen Method
logp	1.431		Crippen Method
mcvol	51.210	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hvapt	29.10	kJ/mol	223.00	NIST Webbook
hvapt	29.10	kJ/mol	223.00	NIST Webbook

Sources

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=C13780637&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

hvapt:	Enthalpy of vaporization at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume

Latest version available from:

<https://www.cheméo.com/cid/74-883-0/Phosphonothioic-difluoride.pdf>

Generated by Cheméo on 2024-07-20 09:42:45.921727263 +0000 UTC m=+158435.168832621.

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