

Bis(ditrifluoromethyldiphosphino) sulfide

Other names:	Di[bis(trifluoromethyl)phosphino] sulfide
Inchi:	InChI=1S/C4F12P2S/c5-1(6,7)17(2(8,9)10)19-18(3(11,12)13)4(14,15)16
InchiKey:	ULASFWIFQPWNDN-UHFFFAOYSA-N
Formula:	C4F12P2S
SMILES:	FC(F)(F)P(SP(C(F)(F)F)C(F)(F)F)C(F)(F)F
Mol. weight [g/mol]:	370.04
CAS:	1486-20-0

Physical Properties

Property code	Value	Unit	Source
log10ws	0.30		Crippen Method
logp	6.593		Crippen Method
mccvol	145.730	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hvapt	42.20	kJ/mol	304.00	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=C1486200&Units=SI

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/80-691-6/Bis-ditrifluoromethyldiphosphino-sulfide.pdf>

Generated by Cheméo on 2024-05-04 07:54:15.108624723 +0000 UTC m=+17098504.029202037.

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