

Diethyl phosphite

Other names:	Phosphonic acid, diethyl ester Diethyl acid phosphite Diethyl hydrogen phosphite Diethyl phosphonate Hydrogen diethyl phosphite O,O-Diethyl phosphonate OPH(OC2H5)2 Diethoxyphosphine oxide Ethyl phosphonate ((EtO)2HPO) Phosphorous acid, diethyl ester NSC 2665
Inchi:	InChI=1S/C4H11O3P/c1-3-6-8(5)7-4-2/h8H,3-4H2,1-2H3
InchiKey:	MJUJXFBTEFXVKU-UHFFFAOYSA-N
Formula:	C4H11O3P
SMILES:	CCO[PH](=O)OCC
Mol. weight [g/mol]:	138.10
CAS:	762-04-9

Physical Properties

Property code	Value	Unit	Source
hvap	49.30	kJ/mol	NIST Webbook
ie	10.31	eV	NIST Webbook
ie	10.31	eV	NIST Webbook
log10ws	-2.13		Crippen Method
logp	1.449		Crippen Method
mcvol	105.290	ml/mol	McGowan Method
rinpol	953.00		NIST Webbook
rinpol	953.00		NIST Webbook
rinpol	953.00		NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hvapt	38.10	kJ/mol	404.50	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=C762049&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

h_{vap}:	Enthalpy of vaporization at standard conditions
h_{vapt}:	Enthalpy of vaporization at a given temperature
ie:	Ionization energy
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/10-408-7/Diethyl-phosphite.pdf>

Generated by Cheméo on 2024-04-19 02:10:16.75827552 +0000 UTC m=+15781865.678852837.

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