

Triethyl phosphite

Other names:	Phosphorous acid, triethyl ester Triethoxyphosphine (C ₂ H ₅ O) ₃ P Ethyl phosphite, (EtO) ₃ P UN 2323 Phosphorus ether
Inchi:	InChI=1S/C6H15O3P/c1-4-7-10(8-5-2)9-6-3/h4-6H2,1-3H3
InchiKey:	BDZBKCUKTQZUTL-UHFFFAOYSA-N
Formula:	C ₆ H ₁₅ O ₃ P
SMILES:	CCOP(OCC)OCC
Mol. weight [g/mol]:	166.16
CAS:	122-52-1

Physical Properties

Property code	Value	Unit	Source
hvap	53.00	kJ/mol	NIST Webbook
ie	8.99	eV	NIST Webbook
ie	8.40	eV	NIST Webbook
ie	8.80	eV	NIST Webbook
ie	8.92	eV	NIST Webbook
ie	8.30	eV	NIST Webbook
ie	8.20	eV	NIST Webbook
ie	8.40	eV	NIST Webbook
ie	9.15	eV	NIST Webbook
ie	8.99	eV	NIST Webbook
ie	8.90	eV	NIST Webbook
log10ws	1.58		Crippen Method
logp	2.323		Crippen Method
mcvol	133.470	ml/mol	McGowan Method
rmpol	922.00		NIST Webbook
rmpol	922.00		NIST Webbook
rmpol	922.00		NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.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=C122521&Units=SI

Legend

h_{vap}:	Enthalpy of vaporization at standard conditions
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

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