

# Phosphine

<b>Other names:</b>	Fosforowodor Gas-ex-B HYDROGEN PHOSPHIDE PH3 PHOSPHOROUS TRIHYDRIDE Phosphene Phosphorus hydride Phosphorus trihydride Phosphorwasserstoff Rcra waste number P096 Trihydrogen phosphide UN 2199
<b>Inchi:</b>	InChI=1S/H3P/h1H3
<b>InchiKey:</b>	XYFCBTPGUUZFHU-UHFFFAOYSA-N
<b>Formula:</b>	H3P
<b>SMILES:</b>	P
<b>Mol. weight [g/mol]:</b>	34.00
<b>CAS:</b>	7803-51-2

## Physical Properties

Property code	Value	Unit	Source
af	0.0380		KDB
affp	785.00	kJ/mol	NIST Webbook
basg	750.90	kJ/mol	NIST Webbook
dm	0.60	debye	KDB
gf	25.41	kJ/mol	KDB
hf	22.90	kJ/mol	KDB
ie	9.87 ± 0.01	eV	NIST Webbook
ie	9.96	eV	NIST Webbook
ie	10.00	eV	NIST Webbook
ie	9.96	eV	NIST Webbook
ie	9.96 ± 0.01	eV	NIST Webbook
ie	10.13 ± 0.02	eV	NIST Webbook
ie	9.87 ± 0.00	eV	NIST Webbook
ie	10.28	eV	NIST Webbook
ie	10.05 ± 0.05	eV	NIST Webbook
ie	9.87 ± 0.00	eV	NIST Webbook

ie	10.30 ± 0.10	eV	NIST Webbook
ie	9.98	eV	NIST Webbook
ie	10.59	eV	NIST Webbook
ie	10.59 ± 0.05	eV	NIST Webbook
ie	9.97 ± 0.02	eV	NIST Webbook
ie	9.87 ± 0.01	eV	NIST Webbook
ie	10.10	eV	NIST Webbook
log10ws	3.76		Crippen Method
logp	0.058		Crippen Method
mcvol	31.320	ml/mol	McGowan Method
pc	6540.00	kPa	KDB
tb	185.40	K	KDB
tc	324.50	K	KDB
tf	140.00	K	KDB
tt	139.41 ± 0.20	K	NIST Webbook

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hsubt	17.20	kJ/mol	134.50	NIST Webbook
hvapt	14.60 ± 0.10	kJ/mol	186.00	NIST Webbook
hvapt	14.60	kJ/mol	185.00	NIST Webbook
rho1	1529.00	kg/m <sup>3</sup>	298.00	KDB

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.40657e+01
Coeff. B	-1.66816e+03
Coeff. C	-8.84000e+00
Temperature range (K), min.	139.37
Temperature range (K), max.	324.75

# Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>KDB:</b>	<a href="https://www.cheric.org/files/research/kdb/mol/mol1921.mol">https://www.cheric.org/files/research/kdb/mol/mol1921.mol</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C7803512&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C7803512&amp;Units=SI</a>
<b>The Yaws Handbook of Vapor Pressure:</b>	<a href="https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure">https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>

# Legend

<b>af:</b>	Acentric Factor
<b>affp:</b>	Proton affinity
<b>basg:</b>	Gas basicity
<b>dm:</b>	Dipole Moment
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hsubt:</b>	Enthalpy of sublimation at a given temperature
<b>hvapt:</b>	Enthalpy of vaporization at a given temperature
<b>ie:</b>	Ionization energy
<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>pc:</b>	Critical Pressure
<b>pvap:</b>	Vapor pressure
<b>rho:</b>	Liquid Density
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>tt:</b>	Triple Point Temperature

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