

# Parathion

## Other names:

AAT  
AATP  
AC 3422  
ACC 3422  
Alkron  
Alleron  
American Cyanamid 3422  
Aphamite  
Aralo  
B 404  
BAY E-605  
Bayer E-605  
Bladan  
Bladan F  
Compound 3422  
Corothion  
Corthion  
Corthione  
DNTP  
DPP  
Danthion  
Diethoxy, nitro-phenoxy phosphorothioate  
Diethyl 4-nitrophenyl phosphorothioate  
Diethyl 4-nitrophenyl phosphorothionate  
Diethyl p-Nitrophenyl phosphorothionate  
Diethyl p-nitrophenyl thionophosphate  
Diethyl p-nitrophenyl thiophosphate  
Diethyl para-nitrophenol thiophosphate  
Diethyl parathion  
Drexel parathion 8E  
E 605  
E 605 FORTE  
E 605F  
ENT 15,108  
ENT 15108  
Ecatox  
Ecatox 20  
Ekatox  
Ethlon  
Ethyl Parathion

Ethyle-parathion  
Etilon  
Etylparation  
Folidol  
Folidol E  
Folidol E & E 605  
Folidol E-605  
Folidol Oil  
Fosfermo  
Fosferno  
Fosfex  
Fosfive  
Fosova  
Fostern  
Fostox  
Galpar  
Gearphos  
Genithion  
Isotox  
Jacutin  
Kokotine  
Kolphos  
Kwell  
Kypthion  
Lethalaire G-54  
Lirothion  
Murfos  
Murphos  
NA 2783  
NCI-C00226  
NIUIF 100  
Niran  
Niran E-4  
Nitrostigmin  
Nitrostigmine  
Nitrostygmine  
Nourithion  
O,O-Diethyl O-(4-nitrophenyl) thiophosphate  
O,O-Diethyl O-(p-Nitrophenyl) phosphorothioate  
O,O-Diethyl O-(p-nitrophenyl) thionophosphate  
O,O-Diethyl O-(p-nitrophenyl) thiophosphate  
O,O-Diethyl-O-(4-nitrophenyl) phosphorothioate  
O,O-Diethyl-O-p-nitrofenylester kyseliny thiofosforecne

O,O-Dietil-O-(4-nitro-fenil)-monotiofosfato  
O,O-diaethyl-O-(4-nitro-phenyl)-monothiophosphat  
O,O-diethyl O-4-nitrophenyl phosphorothioate (parathion)  
O,O-diethyl-O-(4-nitro-fenil)monothiiofosfaat  
OMS 19  
Oleofos 20  
Oleoparaphene  
Oleoparathene  
Oleoparathion  
Orthophos  
PAC  
Pacol  
Panthion  
Paradust  
Paramar  
Paramar 50  
Paraphos  
Parathene  
Parathion A  
Parathion mixture  
Parathion-aethyl  
Parathion-ethyl  
Parawet  
Penncap E  
Pestox plus  
Pethion  
Phenol, p-nitro-, O-ester with O,O-diethylphosphorothioate  
Phoskil  
Phosphemol  
Phosphenol  
Phosphorothioic acid, O,O-diethyl O-(4-nitrophenyl) ester  
Phosphorothioic acid, O,O-diethyl O-(p-nitrophenyl) ester  
Phosphostigmine  
RB  
Rcra waste number P089  
Rhodiasol  
Rhodiatox  
Rhodiatrox  
Rodiatox  
S.N.P.  
SNP  
Selephos  
Sixty-three special E.C. insecticide

Soprathion  
 Stabilized ethyl parathion  
 Stathion  
 Strathion  
 Sulphos  
 Super Rodiatox  
 T-47  
 TOX 47  
 Thiofos  
 Thiomex  
 Thiophos  
 Thiophos 3422  
 Thiophos Parathion 4 E.C.  
 Thiophosphate de O,O-diethyle et de O-(4-nitrophenyle)  
 Tiofos  
 Vapophos  
 Vitrex  
 Vitrex HGI

diethoxy-(4-nitrophenoxy)-sulfanylidene phosphorane

**Inchi:** InChI=1S/C10H14NO5PS/c1-3-14-17(18,15-4-2)16-10-7-5-9(6-8-10)11(12)13/h5-8H,3-4H  
**InchiKey:** LCCNCVORNKJIRZ-UHFFFAOYSA-N  
**Formula:** C10H14NO5PS  
**SMILES:** CCOP(=S)(OCC)Oc1ccc([N+](=O)[O-])cc1  
**Mol. weight [g/mol]:** 291.26  
**CAS:** 56-38-2

## Physical Properties

Property code	Value	Unit	Source
log10ws	-4.29		Aqueous Solubility Prediction Method
log10ws	-4.66		Estimated Solubility Method
logp	3.271		Crippen Method
mcvol	199.840	ml/mol	McGowan Method
rinpol	1961.00		NIST Webbook
rinpol	1989.00		NIST Webbook
rinpol	2000.00		NIST Webbook
rinpol	1947.00		NIST Webbook
rinpol	1946.00		NIST Webbook
rinpol	1935.00		NIST Webbook

rinpol	2008.00		NIST Webbook
rinpol	1962.00		NIST Webbook
rinpol	1950.00		NIST Webbook
rinpol	1940.00		NIST Webbook
rinpol	1975.00		NIST Webbook
rinpol	1975.00		NIST Webbook
rinpol	1994.00		NIST Webbook
rinpol	1996.00		NIST Webbook
rinpol	1941.00		NIST Webbook
rinpol	1935.00		NIST Webbook
rinpol	1971.00		NIST Webbook
rinpol	1942.00		NIST Webbook
rinpol	1989.00		NIST Webbook
rinpol	1953.00		NIST Webbook
rinpol	1953.00		NIST Webbook
rinpol	1942.00		NIST Webbook
rinpol	1994.00		NIST Webbook
rinpol	1942.00		NIST Webbook
rinpol	1956.00		NIST Webbook
rinpol	1961.00		NIST Webbook
rinpol	1956.00		NIST Webbook
rinpol	1940.00		NIST Webbook
rinpol	2008.00		NIST Webbook
rinpol	1950.00		NIST Webbook
rinpol	1941.00		NIST Webbook
rinpol	1935.00		NIST Webbook
ripol	2967.00		NIST Webbook
ripol	2967.00		NIST Webbook
tf	279.25	K	Aqueous Solubility Prediction Method
tf	278.53 ± 0.20	K	NIST Webbook
tf	279.40 ± 0.10	K	NIST Webbook

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hfust	15.72	kJ/mol	278.10	NIST Webbook
hsubt	100.60	kJ/mol	308.00	NIST Webbook

# Sources

<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=C56382&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C56382&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</a>
<b>Aqueous Solubility Prediction Method:</b>	<a href="http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa">http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa</a>
<b>Estimated Solubility Method:</b>	<a href="http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt">http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt</a>

# Legend

<b>hfust:</b>	Enthalpy of fusion at a given temperature
<b>hsubt:</b>	Enthalpy of sublimation at a given temperature
<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>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	Polar retention indices
<b>tf:</b>	Normal melting (fusion) point

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