

# Demeton-S

<b>Other names:</b>	1-(2-diethoxyphosphorylsulfanylethylsulfanyl)ethane Demeton thiol Demetonthiol Diethyl S-(2-ethioethyl)thiophosphate Disulfoton O-analog Disulfoton oxon Ethylthionodemeton IS IS (pesticide) Isodemeton Isosystox Izosystox O,O-Diaethyl-S(2-aethylthio-aethyl)-monothiophosphat O,O-Diethyl S-(2-(ethylthio)ethyl) phosphorothiolate O,O-Diethyl S-(2-eththioethyl)phosphorothioate O,O-Diethyl S-[2-(Ethylthio)ethyl] phosphorothioate O,O-Diethyl S-[2-(ethylthio)ethyl] thiophosphate O,O-Diethyl S-ethyl-2-ethylmercaptophosphorothiolate O,O-Diethyl-S-(2-ethylthio-ethyl)-monothiofosfaat O,O-Dietil-S-(2-etiltio-etil)-monotiofosfato O,O-Dietyl-S-2-etylmerkaptioetylthiofosfat Phosphorothioic acid, O,O-diethyl S-[2-(ethylthio)ethyl] ester Po-systox Systox-thiol Thiol systox Thioldemeton Thiophosphate de O,O-diethyle et de S-(2-ethylthio-ethyle)
<b>Inchi:</b>	InChI=1S/C8H19O3PS2/c1-4-10-12(9,11-5-2)14-8-7-13-6-3/h4-8H2,1-3H3
<b>InchiKey:</b>	GRPRVIYRYGLIJU-UHFFFAOYSA-N
<b>Formula:</b>	C8H19O3PS2
<b>SMILES:</b>	CCOP(=O)(OCC)SCCSCC
<b>Mol. weight [g/mol]:</b>	258.34
<b>CAS:</b>	126-75-0

## Physical Properties

Property code	Value	Unit	Source
---------------	-------	------	--------

hvap	76.40	kJ/mol	NIST Webbook
log10ws	-2.11		Aqueous Solubility Prediction Method
logp	3.654		Crippen Method
mcvol	194.350	ml/mol	McGowan Method
rinpol	1706.00		NIST Webbook
rinpol	1738.00		NIST Webbook
rinpol	1706.00		NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	2.10887e+01
Coeff. B	-9.18962e+03
Temperature range (K), min.	441.79
Temperature range (K), max.	582.46

## Sources

<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>
<b>Aqueous Solubility Prediction Method:</b>	<a href="http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDataset002.xlsx">http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDataset002.xlsx</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=C126750&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C126750&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>

## Legend

<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<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>pvap:</b>	Vapor pressure
<b>rinpol:</b>	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/48-430-1/Demeton-S.pdf>

Generated by Cheméo on 2024-04-24 09:57:27.283641271 +0000 UTC m=+16241896.204218596.

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