

Oxamyl

Other names:	2-(Dimethylamino)-N-(((methylamino)carbonyl)oxy)-2-oxoethanimidothioic acid methyl ester 2-Dimethylamino-1-(methylthio)glyoxal O-methylcarbamoylmonoxime D 1410 DPX 1410 DPX 1410L Du Pont 1410 Ethanimidothioic acid, 2-(dimethylamino)-N-[[[(methylamino)carbonyl]oxy]-2-oxo-, methyl ester Insecticide-nematicide 1410 Methyl 1-(dimethylcarbamoyl)-N-(methylcarbamoyloxy)thioformimidate Methyl N',N'-dimethyl-N-((methylcarbamoyl)oxy)-1-thiooxamimidate N',N'-Dimethyl-N-((methylcarbamoyl)oxy)-1-thiooxamimidic acid methyl ester N',N'-dimethylcarbamoyl(methylthio)methylenamine N-methylcarbamate N,N-Dimethyl-«alpha»-methylcarbamoyloxyimino-«alpha»-(methylthio)acetamide N,N-Dimethyl-Â«alphaÂ»-methylcarbamoyloxyimino-Â«alphaÂ»-(methylthio)acetamide N,N-dimethyl-2-methylcarbamoyloxyimino-2-(methylthio)acetamide NSC 379588 Nematicide 1410 Oxamimidic acid, N',N'-dimethyl-N-((methylcarbamoyl)oxy)-1-methylthio- Oxamimidic acid, N',N'-dimethyl-N-[(methylcarbamoyl)oxy]-1-thio-, methyl ester Oxamyl (pesticide) S-Methyl 1-(dimethylcarbamoyl)-N-[(methylcarbamoyl)oxy]thioformimidate Thioxamyl Vydate Vydate L Vydate L insecticide/nematicide Vydate L oxamyl insecticide/nematocide
Inchi:	InChI=1S/C7H13N3O3S/c1-8-7(12)13-9-5(14-4)6(11)10(2)3/h1-4H3,(H,8,12)
InchiKey:	KZAUOCCYDRDERY-UHFFFAOYSA-N
Formula:	C7H13N3O3S
SMILES:	CNC(=O)ON=C(SC)C(=O)N(C)C
Mol. weight [g/mol]:	219.26
CAS:	23135-22-0

Physical Properties

Property code	Value	Unit	Source
hf	-309.89	kJ/mol	Joback Method

hvap	65.77			kJ/mol	Joback Method
log10ws	0.11				Aqueous Solubility Prediction Method
log10ws	0.11				Estimated Solubility Method
logp	0.107				Crippen Method
mcvol	160.490			ml/mol	McGowan Method
pc	2912.39			kPa	Joback Method
tb	697.67			K	Joback Method
tc	917.68			K	Joback Method
tf	373.01 ± 0.20			K	NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hfust	30.17	kJ/mol	372.20	NIST Webbook

Sources

Joback Method:

https://en.wikipedia.org/wiki/Joback_method

Aqueous Solubility Prediction Method:

<http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa>

Estimated Solubility Method:

http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C23135220&Units=SI>

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Legend

hf:	Enthalpy of formation at standard conditions
hfust:	Enthalpy of fusion at a given temperature
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature

tc: Critical Temperature
tf: Normal melting (fusion) point

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