

# Dixanthogen

**Other names:** Formic acid, dithiobis[thio-, O,O-diethyl ester  
Thioperoxydicarbonic acid ( $[(HO)C(S)]_2S_2$ ), diethyl ester  
Antigal  
Auligen  
Aulin  
Aulinogen  
Bexide  
Bis[ethylxanthogen]  
Diethyl Dixanthogenate  
Diethyl xanthogenate  
Diethylxanthogen disulfide  
Dithiobis(thioformic acid) O,O-diethyl ester  
Dixan  
EXD  
Galasan  
Herbisan  
Herbisan 5  
K Preparation  
Lenisarin  
O,O-Diethyl Dithiobis[thioformate]  
Scabacidol  
Thioperoxydicarbonic acid, diethyl ester  
Xantoscabin  
Diethyl dithiobis(thiono formate)  
Bis(ethylxanthic)disulfide  
Diethylxanthic disulfide  
DEX  
Ethyl xanthogen disulfide  
O,O-Diethyl ester of dithiobis(thioformic acid)  
Preparation K  
Silfasan  
Sulfasan  
Xanthogen, bis(ethyl-  
Bisethylxanthogen disulfide  
Di-ethoxythiokarbonyl-disulfid  
Diethyl dixanthogen  
Ethylxanthic disulfide  
Skabilan  
Thioperoxydicarbonic acid ( $(((HO)C(S))_2S_2$ ), OC,OC'-diethyl ester  
NSC 402561

**Inchi:** InChI=1S/C6H10O2S4/c1-3-7-5(9)11-12-6(10)8-4-2/h3-4H2,1-2H3  
**InchiKey:** FVIGODVHAVLZOO-UHFFFAOYSA-N  
**Formula:** C6H10O2S4  
**SMILES:** CCOC(=S)SSC(=S)OCC  
**Mol. weight [g/mol]:** 242.40  
**CAS:** 502-55-6

## Physical Properties

Property code	Value	Unit	Source
gf	90.00	kJ/mol	Joback Method
hf	-54.87	kJ/mol	Joback Method
hfus	31.14	kJ/mol	Joback Method
hvap	60.86	kJ/mol	Joback Method
log10ws	-3.96		Crippen Method
logp	3.011		Crippen Method
mcvol	163.940	ml/mol	McGowan Method
pc	3857.88	kPa	Joback Method
tb	659.16	K	Joback Method
tc	916.10	K	Joback Method
tf	339.18	K	Joback Method
vc	0.588	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	348.82	J/molxK	659.16	Joback Method
cpg	358.50	J/molxK	701.98	Joback Method
cpg	367.46	J/molxK	744.81	Joback Method
cpg	375.75	J/molxK	787.63	Joback Method
cpg	383.44	J/molxK	830.46	Joback Method
cpg	390.56	J/molxK	873.28	Joback Method
cpg	397.18	J/molxK	916.10	Joback Method

# 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>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</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=C502556&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C502556&amp;Units=SI</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>cpg:</b>	Ideal gas heat capacity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvac:</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>mccol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/98-527-9/Dixanthogen.pdf>

Generated by Cheméo on 2024-05-02 03:45:37.557923882 +0000 UTC m=+16910786.478501191.

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