

# Biuret, 1,5-bis(2-chloroethyl)-1-nitroso-

**Inchi:** InChI=1S/C6H10Cl2N4O3/c7-1-3-9-5(13)10-6(14)12(11-15)4-2-8/h1-4H2,(H2,9,10,13,14)  
**InchiKey:** KEUPEGWEPJSQDJ-UHFFFAOYSA-N  
**Formula:** C6H10Cl2N4O3  
**SMILES:** O=NN(CCCl)C(O)=NC(O)=NCCCl  
**Mol. weight [g/mol]:** 257.07  
**CAS:** 13857-12-0

## Physical Properties

Property code	Value	Unit	Source
hf	-458.91	kJ/mol	Joback Method
hvap	89.01	kJ/mol	Joback Method
log10ws	-1.35		Crippen Method
logp	1.275		Crippen Method
mcvol	164.510	ml/mol	McGowan Method
pc	2758.46	kPa	Joback Method
tb	824.86	K	Joback Method
tc	1024.11	K	Joback Method

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C13857120&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>

## Legend

**hf:** Enthalpy of formation at standard conditions  
**hvap:** Enthalpy of vaporization at standard conditions  
**log10ws:** 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>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

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