

# Bis(2-chloroethyl) sulfone

<b>Other names:</b>	1-Chloro-2-(2-chloroethanesulfonyl)ethane 1-chloro-2-(2-chloroethylsulfonyl)ethane Bis(«beta»-chloroethyl) sulfone Bis(Â«betaÂ»-chloroethyl) sulfone Ethane, 1,1'-sulfonylbis(2-chloro- H Sulfone Mustard gas sulfone Mustard sulfone NSC 26284 Sulfone, bis(2-chloroethyl) Yperite sulfone
<b>Inchi:</b>	InChI=1S/C4H8Cl2O2S/c5-1-3-9(7,8)4-2-6/h1-4H2
<b>InchiKey:</b>	LUYAMNYBNTVQJG-UHFFFAOYSA-N
<b>Formula:</b>	C4H8Cl2O2S
<b>SMILES:</b>	O=S(=O)(CCCl)CCCl
<b>Mol. weight [g/mol]:</b>	191.08
<b>CAS:</b>	471-03-4

## Physical Properties

Property code	Value	Unit	Source
gf	-509.60	kJ/mol	Joback Method
hf	-610.72	kJ/mol	Joback Method
hfus	25.89	kJ/mol	Joback Method
hvap	51.90	kJ/mol	Joback Method
log10ws	-1.50		Aqueous Solubility Prediction Method
logp	0.879		Crippen Method
mccvol	119.790	ml/mol	McGowan Method
pc	4322.57	kPa	Joback Method
tb	413.56	K	Joback Method
tc	592.91	K	Joback Method
tf	233.24	K	Joback Method
vc	0.483	m3/kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	206.99	J/mol×K	413.56	Joback Method
cpg	215.40	J/mol×K	443.45	Joback Method
cpg	223.48	J/mol×K	473.34	Joback Method
cpg	231.24	J/mol×K	503.24	Joback Method
cpg	238.68	J/mol×K	533.13	Joback Method
cpg	245.79	J/mol×K	563.02	Joback Method
cpg	252.57	J/mol×K	592.91	Joback Method

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C471034&Units=SI>

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**Aqueous Solubility Prediction Method:** <http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa>

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

## 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>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>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

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