

# 1,8-Diamino-3,6-dioxaoctane

<b>Other names:</b>	Ethanamine, 2,2'-[1,2-ethanediylbis(oxy)]bis- 3,6-Dioxaoctane-1,8-diamine 3,6-dioxaoctamethylenediamine 1,2-bis(2-aminoethoxy)ethane
<b>Inchi:</b>	InChI=1S/C6H16N2O2/c7-1-3-9-5-6-10-4-2-8/h1-8H2
<b>InchiKey:</b>	IWBOPFCKHIJFMS-UHFFFAOYSA-N
<b>Formula:</b>	C6H16N2O2
<b>SMILES:</b>	NCCOCCOCCN
<b>Mol. weight [g/mol]:</b>	148.20
<b>CAS:</b>	929-59-9

## Physical Properties

Property code	Value	Unit	Source
gf	-77.46	kJ/mol	Joback Method
hf	-364.03	kJ/mol	Joback Method
hfus	24.07	kJ/mol	Joback Method
hvap	58.80	kJ/mol	NIST Webbook
log10ws	0.62		Crippen Method
logp	-1.063		Crippen Method
mcvol	127.100	ml/mol	McGowan Method
pc	3403.91	kPa	Joback Method
tb	526.58	K	Joback Method
tc	716.93	K	Joback Method
tf	368.36	K	Joback Method
vc	0.466	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	311.78	J/molxK	526.58	Joback Method
cpg	323.37	J/molxK	558.30	Joback Method
cpg	334.53	J/molxK	590.03	Joback Method
cpg	345.26	J/molxK	621.75	Joback Method
cpg	355.55	J/molxK	653.48	Joback Method

cpg	365.39	J/mol×K	685.20	Joback Method
cpg	374.79	J/mol×K	716.93	Joback Method
hvapt	56.20	kJ/mol	323.00	NIST Webbook

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C929599&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C929599&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>
<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>

## 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>hvapt:</b>	Enthalpy of vaporization at a given temperature
<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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