

# n-Hexyl-2,2'-iminodiethanol

<b>Other names:</b>	2,2'-(n-Hexylamino)diethanol 2,2'-(Hexylamino)diethanol
<b>Inchi:</b>	InChI=1S/C10H23NO2/c1-2-3-4-5-6-11(7-9-12)8-10-13/h12-13H,2-10H2,1H3
<b>InchiKey:</b>	GIACMWUKBLHAAG-UHFFFAOYSA-N
<b>Formula:</b>	C10H23NO2
<b>SMILES:</b>	CCCCCN(CCO)CCO
<b>Mol. weight [g/mol]:</b>	189.30
<b>CAS:</b>	6752-33-6

## Physical Properties

Property code	Value	Unit	Source
gf	-129.54	kJ/mol	Joback Method
hf	-486.66	kJ/mol	Joback Method
hfus	32.85	kJ/mol	Joback Method
hvap	73.25	kJ/mol	Joback Method
log10ws	-1.11		Crippen Method
logp	0.853		Crippen Method
mcvol	173.480	ml/mol	McGowan Method
pc	2510.03	kPa	Joback Method
tb	625.00	K	Joback Method
tc	782.17	K	Joback Method
tf	356.57	K	Joback Method
vc	0.651	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	479.06	J/molxK	625.00	Joback Method
cpg	491.16	J/molxK	651.19	Joback Method
cpg	502.75	J/molxK	677.39	Joback Method
cpg	513.84	J/molxK	703.58	Joback Method
cpg	524.47	J/molxK	729.78	Joback Method
cpg	534.63	J/molxK	755.97	Joback Method
cpg	544.36	J/molxK	782.17	Joback Method

# Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	427.50 ± 0.50	K	0.50	NIST Webbook

## Sources

<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=C6752336&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C6752336&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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>

## 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>mvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tbrp:</b>	Boiling point at reduced pressure
<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/48-010-7/n-Hexyl-2-2-iminodiethanol.pdf>

Generated by Cheméo on 2024-04-19 22:02:50.877127553 +0000 UTC m=+15853419.797704925.

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