

# 1-Piperazineethanol

<b>Other names:</b>	(2-Hydroxyethyl)piperazine 1-(2-Hydroxyethyl)piperazine 1-(«beta»-Hydroxyethyl)piperazine 1-Hydroxy-2-(piperazin-1-yl)ethane 1-Piperazinethanol 2-(1-Piperazinyl)ethanol 2-Piperazinoethanol 2-piperazin-1-ylethanol 4-(2-Hydroxyethyl)piperazine Ethanol, 2-(1-piperazinyl)- Hydroxyethylpiperazine N-(2-hydroxyethyl)piperazine N-(Hydroxyethyl)piperazine N-(«beta»-Hydroxyethyl)piperazine NSC 26884 USAF DO-22 piperazine, N-(2-hydroxyethyl)-
<b>Inchi:</b>	InChI=1S/C6H14N2O/c9-6-5-8-3-1-7-2-4-8/h7,9H,1-6H2
<b>InchiKey:</b>	WFCSWCVEJLETKA-UHFFFAOYSA-N
<b>Formula:</b>	C6H14N2O
<b>SMILES:</b>	OCCN1CCNCC1
<b>Mol. weight [g/mol]:</b>	130.19
<b>CAS:</b>	103-76-4

## Physical Properties

Property code	Value	Unit	Source
chl	-4134.30 ± 1.10	kJ/mol	NIST Webbook
hfl	-227.60 ± 1.10	kJ/mol	NIST Webbook
hvap	78.80 ± 0.70	kJ/mol	NIST Webbook
log10ws	0.75		Crippen Method
logp	-1.116		Crippen Method
mcvol	110.370	ml/mol	McGowan Method
tb	519.20	K	NIST Webbook
tb	519.55	K	NIST Webbook

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpl	319.00	J/molxK	353.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	285.00	J/molxK	298.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	289.90	J/molxK	303.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	292.90	J/molxK	308.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	296.00	J/molxK	313.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	298.90	J/molxK	318.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	301.70	J/molxK	323.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	305.00	J/molxK	328.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	308.10	J/molxK	333.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K

cpl	311.00	J/mol×K	338.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	313.90	J/mol×K	343.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
cpl	316.50	J/mol×K	348.15	Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K
hvapt	77.30 ± 0.70	kJ/mol	325.50	NIST Webbook

## Sources

**Molar Heat Capacity (Cp) of Aqueous Cyclic Amine Solutions from (298.15 to 353.15) K**  
Crippen Method:

<https://www.doi.org/10.1021/je400178k>

**NIST Webbook:**

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

**Crippen Method:**

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C103764&Units=SI>

**Crippen Method:**

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

**Molar excess enthalpy (Hm E) for systems of aqueous piperazine derivatives:**

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

<https://www.doi.org/10.1016/j.jct.2015.06.006>

## Legend

<b>chl:</b>	Standard liquid enthalpy of combustion
<b>cpl:</b>	Liquid phase heat capacity
<b>hfl:</b>	Liquid phase enthalpy of formation 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>tb:</b>	Normal Boiling Point Temperature

Latest version available from:

<https://www.chemeo.com/cid/21-034-0/1-Piperazineethanol.pdf>

Generated by Cheméo on 2024-04-27 06:14:27.917266351 +0000 UTC m=+16487716.837843669.

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