

1-Piperazineethanamine, 4-methyl-

Other names:	1-(2-Aminoethyl)-4-methylpiperazine 1-(2-aminoethyl)-4-methyl-piperazine 1-Methyl-4-(2-aminoethyl)piperazine 4-methylpiperazine-1-ethylamine Piperazine, 1-(2-aminoethyl)-4-methyl-
Inchi:	InChI=1S/C7H17N3/c1-9-4-6-10(3-2-8)7-5-9/h2-8H2,1H3
InchiKey:	GOWUDHPKGOIDIX-UHFFFAOYSA-N
Formula:	C7H17N3
SMILES:	CN1CCN(CCN)CC1
Mol. weight [g/mol]:	143.23
CAS:	934-98-5

Physical Properties

Property code	Value	Unit	Source
hvap	64.00 ± 0.20	kJ/mol	NIST Webbook
log10ws	0.78		Crippen Method
logp	-0.807		Crippen Method
mcvol	128.570	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hvapt	64.00	kJ/mol	298.15	Vapour pressure and enthalpy of vaporization of aliphatic poly-amines

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Vapour pressure and enthalpy of
vaporization of aliphatic poly-amines:
McGowan Method:

<https://www.doi.org/10.1016/j.jct.2009.09.003>

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

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C934985&Units=SI>

Legend

h_{vap}:	Enthalpy of vaporization at standard conditions
h_{vapt}:	Enthalpy of vaporization at a given temperature
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume

Latest version available from:

<https://www.chemeo.com/cid/81-121-7/1-Piperazineethanamine-4-methyl.pdf>

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