

# 2-Pyridinamine, N-methyl-

<b>Other names:</b>	Pyridine, 2-(methylamino)- N-2-Pyridylmethylamine 2-(Methylamino)pyridine N-Methyl-2-pyridinamine Methylamino-2 pyridine
<b>Inchi:</b>	InChI=1S/C6H8N2/c1-7-6-4-2-3-5-8-6/h2-5H,1H3,(H,7,8)
<b>InchiKey:</b>	SVEUVITYHIHZQE-UHFFFAOYSA-N
<b>Formula:</b>	C6H8N2
<b>SMILES:</b>	CNc1ccccn1
<b>Mol. weight [g/mol]:</b>	108.14
<b>CAS:</b>	4597-87-9

## Physical Properties

Property code	Value	Unit	Source
ie	8.26 ± 0.05	eV	NIST Webbook
log10ws	-1.25		Crippen Method
logp	1.123		Crippen Method
mcvol	91.600	ml/mol	McGowan Method
tb	473.70	K	NIST Webbook
tb	475.15 ± 3.00	K	NIST Webbook

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hvapt	49.00	kJ/mol	315.50	NIST Webbook

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C4597879&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C4597879&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>

## Legend

<b>hvapt:</b>	Enthalpy of vaporization at a given temperature
<b>ie:</b>	Ionization energy
<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.cheméo.com/cid/65-433-9/2-Pyridinamine-N-methyl.pdf>

Generated by Cheméo on 2024-05-02 16:42:06.619163793 +0000 UTC m=+16957375.539741115.

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