

# Piperidine, 1-nitroso-

<b>Other names:</b>	1-Nitrosopiperidine N-N-Pip N-Nitroso-piperidin N-Nitrosopiperidine NPIP Nitrosopiperidin No-Pip Pyridine, hexahydro-N-nitroso- Rcra waste number U179
<b>Inchi:</b>	InChI=1S/C5H10N2O/c8-6-7-4-2-1-3-5-7/h1-5H2
<b>InchiKey:</b>	UWSDONTXWQOZFN-UHFFFAOYSA-N
<b>Formula:</b>	C5H10N2O
<b>SMILES:</b>	O=NN1CCCCC1
<b>Mol. weight [g/mol]:</b>	114.15
<b>CAS:</b>	100-75-4

## Physical Properties

Property code	Value	Unit	Source
chl	-3366.00 ± 1.00	kJ/mol	NIST Webbook
hf	17.00 ± 2.00	kJ/mol	NIST Webbook
hfl	-31.00 ± 1.00	kJ/mol	NIST Webbook
log10ws	-0.17		Aqueous Solubility Prediction Method
logp	1.154		Crippen Method
mcvol	91.980	ml/mol	McGowan Method
rinpol	185.84		NIST Webbook
rinpol	1040.00		NIST Webbook
rinpol	184.19		NIST Webbook
rinpol	1040.00		NIST Webbook
rinpol	184.19		NIST Webbook

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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hvapt	47.70	kJ/mol	358.00	NIST Webbook
hvapt	47.70 ± 0.80	kJ/mol	333.00	NIST Webbook

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	382.20	K	2.70	NIST Webbook

## Sources

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

**Aqueous Solubility Prediction Method:** <http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa>

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

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C100754&Units=SI>

## Legend

<b>chl:</b>	Standard liquid enthalpy of combustion
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfl:</b>	Liquid phase enthalpy of formation 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>rinpol:</b>	Non-polar retention indices
<b>tbrp:</b>	Boiling point at reduced pressure

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