

# Ethanamine, N-ethyl-N-nitroso-

<b>Other names:</b>	Diethylamine, N-nitroso- Diethylnitrosamine DENA N-Nitroso-N,N-Diethylamine N-Nitrosodiethylamine N,N-Diethylnitrosoamine Nitrosodiethylamine Diaethylnitrosamin Diethylnitrosoamine DEN Ethylamine, N-nitrosodi- N-Ethyl-N-nitrosoethanamine N,N-Diethylnitrosamine NDEA Diethylnitrosamide N-Nitroso-diaethylamine Rcra waste number U174 N-Ethyl-N-nitroso-ethylamine DEN (mutagen) NSC 132
<b>Inchi:</b>	InChI=1S/C4H10N2O/c1-3-6(4-2)5-7/h3-4H2,1-2H3
<b>InchiKey:</b>	WBNQDOYYEUMPFS-UHFFFAOYSA-N
<b>Formula:</b>	C4H10N2O
<b>SMILES:</b>	CCN(CC)N=O
<b>Mol. weight [g/mol]:</b>	102.14
<b>CAS:</b>	55-18-5

## Physical Properties

Property code	Value	Unit	Source
hf	-226.55	kJ/mol	Joback Method
hvap	35.64	kJ/mol	Joback Method
ie	8.76	eV	NIST Webbook
log10ws	-1.25		Crippen Method
logp	1.010		Crippen Method
mcvol	88.750	ml/mol	McGowan Method
pc	3801.00	kPa	Joback Method
rinpol	144.23		NIST Webbook

rinpol	139.17		NIST Webbook
rinpol	876.00		NIST Webbook
rinpol	881.00		NIST Webbook
rinpol	876.00		NIST Webbook
rinpol	139.17		NIST Webbook
tb	450.10	K	NIST Webbook
tc	533.59	K	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C55185&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C55185&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>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>

## Legend

<b>hf:</b>	Enthalpy of formation at standard conditions
<b>h<sub>vap</sub>:</b>	Enthalpy of vaporization at standard conditions
<b>ie:</b>	Ionization energy
<b>log<sub>10</sub>ws:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</b>	Octanol/Water partition coefficient
<b>m<sub>cvol</sub>:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

Latest version available from:

<https://www.chemeo.com/cid/36-262-1/Ethanamine-N-ethyl-N-nitroso.pdf>

Generated by Cheméo on 2024-04-23 06:29:23.413465453 +0000 UTC m=+16143012.334042765.

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