

# Methanimidamide, N'-(3-hydroxyphenyl)-N,N-dimethyl-

<b>Other names:</b>	Formamidine, N'-(m-hydroxyphenyl)-N,N-dimethyl- m-Dimethylaminomethyleneiminophenol 3-Dimethylaminomethyleniminophenol N'-(3-hydroxyphenyl)-N,N-dimethylformamidine formamidine, N,N-dimethyl-N'-(3-hydroxyphenyl)
<b>Inchi:</b>	InChI=1S/C9H12N2O/c1-11(2)7-10-8-4-3-5-9(12)6-8/h3-7,12H,1-2H3
<b>InchiKey:</b>	RKXRPBBZGFLVTF-UHFFFAOYSA-N
<b>Formula:</b>	C9H12N2O
<b>SMILES:</b>	CN(C)C=Nc1cccc(O)c1
<b>Mol. weight [g/mol]:</b>	164.20
<b>CAS:</b>	25635-97-6

## Physical Properties

Property code	Value	Unit	Source
hf	-20.12	kJ/mol	Joback Method
hvap	56.28	kJ/mol	Joback Method
log10ws	-1.07		Crippen Method
logp	1.614		Crippen Method
mcvol	135.440	ml/mol	McGowan Method
pc	3423.86	kPa	Joback Method
tb	601.74	K	Joback Method
tc	837.54	K	Joback Method

## Sources

<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>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C25635976&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C25635976&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>

# Legend

<b>hf:</b>	Enthalpy of formation at standard conditions
<b>h<sub>vap</sub>:</b>	Enthalpy of vaporization at standard conditions
<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>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

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