

# N-Formylmorpholine

<b>Other names:</b>	4-Morpholinecarboxaldehyde 4-formylmorpholine 4-morpholinecarbaldehyde Morpholine, 4-formyl- N-Formylmorpholin N-formyltetrahydro-1,4-oxazine morpholine-4-carbaldehyde
<b>Inchi:</b>	InChI=1S/C5H9NO2/c7-5-6-1-3-8-4-2-6/h5H,1-4H2
<b>InchiKey:</b>	LCEDQNDDFOCWGG-UHFFFAOYSA-N
<b>Formula:</b>	C5H9NO2
<b>SMILES:</b>	O=CN1CCOCC1
<b>Mol. weight [g/mol]:</b>	115.13
<b>CAS:</b>	4394-85-8

## Physical Properties

Property code	Value	Unit	Source
log10ws	0.76		Crippen Method
logp	-0.525		Crippen Method
mcvol	87.870	ml/mol	McGowan Method
pc	5080.00	kPa	Critical Point and Vapor Pressure Measurements for 17 Compounds by a Low Residence Time Flow Method
tb	513.10	K	Isobaric (vapour + liquid) equilibria for N-formylmorpholine with ethylbenzene, n-butylbenzene, iso-propylbenzene and 1,2,4-trimethylbenzene at 101.33 kPa
tb	509.95	K	Cosolvent Selection for Benzene-Cyclohexane Separation in Extractive Distillation

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpl	188.00	J/mol×K	313.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	202.00	J/mol×K	353.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	201.00	J/mol×K	348.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	200.00	J/mol×K	343.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	197.00	J/mol×K	338.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K

cpl	186.00	J/mol×K	308.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	195.00	J/mol×K	333.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	193.00	J/mol×K	328.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	191.00	J/mol×K	323.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
cpl	185.00	J/mol×K	303.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K

cpl	189.00	J/mol×K	318.15	Molar Heat Capacity of Aqueous Sulfolane, 4-Formylmorpholine, 1-Methyl-2-pyrrolidinone, and Triethylene Glycol Dimethyl Ether Solutions from (303.15 to 353.15) K
dvisc	0.0027070	Paxs	343.15	Densities and Viscosities of N-Formylmorpholine + Benzene, + Toluene at Different Temperatures and Atmospheric Pressures
dvisc	0.0022740	Paxs	353.15	Densities and Viscosities of N-Formylmorpholine + Benzene, + Toluene at Different Temperatures and Atmospheric Pressures
dvisc	0.0076556	Paxs	298.15	Densities and Viscosities of N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different Temperatures and Atmospheric Pressure
dvisc	0.0065981	Paxs	303.15	Densities and Viscosities of N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different Temperatures and Atmospheric Pressure
dvisc	0.0050678	Paxs	313.15	Densities and Viscosities of N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different Temperatures and Atmospheric Pressure

dvisc	0.0040357	Paxs	323.15	Densities and Viscosities of N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different Temperatures and Atmospheric Pressure
dvisc	0.0033035	Paxs	333.15	Densities and Viscosities of N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different Temperatures and Atmospheric Pressure
dvisc	0.0027822	Paxs	343.15	Densities and Viscosities of N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different Temperatures and Atmospheric Pressure
dvisc	0.0023680	Paxs	353.15	Densities and Viscosities of N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different Temperatures and Atmospheric Pressure
dvisc	0.0078700	Paxs	298.15	Volumetric Properties and Viscosities for Aqueous NFM Solutions from 25 deg C to 70 deg C
dvisc	0.0067500	Paxs	303.15	Volumetric Properties and Viscosities for Aqueous NFM Solutions from 25 deg C to 70 deg C

dvisc	0.0051300	Paxs	313.15	Volumetric Properties and Viscosities for Aqueous NFM Solutions from 25 deg C to 70 deg C
dvisc	0.0040300	Paxs	323.15	Volumetric Properties and Viscosities for Aqueous NFM Solutions from 25 deg C to 70 deg C
dvisc	0.0032500	Paxs	333.15	Volumetric Properties and Viscosities for Aqueous NFM Solutions from 25 deg C to 70 deg C
dvisc	0.0026800	Paxs	343.15	Volumetric Properties and Viscosities for Aqueous NFM Solutions from 25 deg C to 70 deg C
dvisc	0.0079550	Paxs	298.15	Densities and Viscosities of N-Formylmorpholine + Benzene, + Toluene at Different Temperatures and Atmospheric Pressures
dvisc	0.0068770	Paxs	303.15	Densities and Viscosities of N-Formylmorpholine + Benzene, + Toluene at Different Temperatures and Atmospheric Pressures
dvisc	0.0052140	Paxs	313.15	Densities and Viscosities of N-Formylmorpholine + Benzene, + Toluene at Different Temperatures and Atmospheric Pressures

dvisc	0.0040800	Paxs	323.15	Densities and Viscosities of N-Formylmorpholine + Benzene, + Toluene at Different Temperatures and Atmospheric Pressures
dvisc	0.0032840	Paxs	333.15	Densities and Viscosities of N-Formylmorpholine + Benzene, + Toluene at Different Temperatures and Atmospheric Pressures
pvap	101.32	kPa	509.95	Cosolvent Selection for Benzene-Cyclohexane Separation in Extractive Distillation
rfi	1.48400		293.10	Liquid liquid equilibria of methylcyclohexane benzene N-formylmorpholine at several temperatures
rfi	1.48400		293.10	Measurement and Correlation of Liquid-Liquid Equilibria of Methylcyclohexane + Toluene + N-Formylmorpholine at (293, 303, 313, and 323) K
rfi	1.48440		293.15	Vapor Liquid Equilibrium Data for the Morpholine-4-carbaldehyde + n-Hexane or n-Heptane Binary Systems Using a Static-Synthetic Apparatus
rfi	1.48490		293.15	Experimental solubility for betulin and estrone in various solvents within the temperature range T = (293.2 to 328.2) K

rfi	1.48400	293.10	(Liquid + liquid) equilibria of three ternary systems: (heptane + benzene + N-formylmorpholine), (heptane + toluene + N-formylmorpholine), (heptane + xylene + N-formylmorpholine) from T = (298.15 to 353.15) K
rfi	1.47990	313.15	Refractive properties, speed of sound and FT-IR study of binary mixtures of N-formylmorpholine with some halobenzenes at 303.15, 308.15 and 313.15 K
rfi	1.48190	308.15	Refractive properties, speed of sound and FT-IR study of binary mixtures of N-formylmorpholine with some halobenzenes at 303.15, 308.15 and 313.15 K
rfi	1.48330	303.15	Refractive properties, speed of sound and FT-IR study of binary mixtures of N-formylmorpholine with some halobenzenes at 303.15, 308.15 and 313.15 K
rfi	1.48400	293.10	(Liquid + liquid) equilibria of {heptane + xylene + N-formylmorpholine} ternary system

## Sources

**McGowan Method:**

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

**Volumetric Properties and Viscosities for Aqueous NFM Solutions from 25 deg C to 70 deg C:**

<https://www.doi.org/10.1021/je0498208>

**(Liquid + liquid) equilibria for the ternary system of (N-formylmorpholine + Mesitylene + 4-(liquid + liquid) equilibria for ternary systems of  $\text{C}_6\text{H}_6$  + N-formylmorpholine + 4-(liquid + liquid) equilibria for ternary systems of  $\text{C}_6\text{H}_6$  + N-formylmorpholine + Ethane and Ethane + N-formylmorpholine + Solvent + liquid equilibria for ternary systems of N-formylmorpholine with alkanes: liquid equilibria of methylcyclohexane benzene N-formylmorpholine for the Binary Systems of N-Formylmorpholine with Cycloalkanes: liquid equilibria of three ternary systems: (heptane + benzene + N-formylmorpholine), (heptane + xylene + N-formylmorpholine) ternary experimental solubility of diosgenin in N-formylmorpholine various solvents between 303.15 K and 326.27 K in N-Formylmorpholine: Crippen Method:**

**Refractive properties, speed of sound and FT-IR study of binary mixtures of N-Formylmorpholine with some halobenzenes at 303.15, 308.15 and 313.15 K: Measurement and Correlation of Liquid-Liquid Equilibria of Densities and X-ray densities of Benzene + N-Formylmorpholine at Temperatures +313, 308.15 and 303.15 K and at different temperatures and Atmospheric Pressures: (N-Formylmorpholine) + Benzene equilibria for N-formylmorpholine with ethylbenzene, N-Formylmorpholine + Benzene and esterbenzene mixtures and (N-Formylmorpholine) + p-Xylene (303.15, 308.15, 313.15 K) N-Formylmorpholine (NFM) + p-Xylene, + o-Xylene, + m-Xylene at Different N-Formylmorpholine with Toluene and Benzene at 101.995 kPa or Benzene-Cyclohexane Separation in Extractive Distillation: for the Binary Systems of N-Formylmorpholine with Molar Heat Capacities of Aqueous Sulfolane, 4-Formylmorpholine, Cinnamyl Polidinone, and Triethylene Glycol Dimethyl Ether Critical Point and Vapor Pressure Measurements from 300 to 333.15 K: Measurements for 17 Compounds by a Vapour Residue Compressibility Method: Morpholine-4-carbaldehyde + n-Hexane or n-Heptane Binary Systems Using a Static-Synthetic Apparatus:**

- <https://www.doi.org/10.1016/j.fluid.2012.05.010>  
<https://www.doi.org/10.1016/j.jct.2011.01.011>  
<https://www.doi.org/10.1016/j.jct.2013.02.006>  
<https://www.doi.org/10.1021/je050172h>  
<https://www.doi.org/10.1016/j.jct.2011.10.022>  
<https://www.doi.org/10.1016/j.fluid.2007.03.013>  
<https://www.doi.org/10.1021/je020089j>  
<https://www.doi.org/10.1016/j.jct.2006.12.004>  
<https://www.doi.org/10.1016/j.jct.2007.05.001>  
<https://www.doi.org/10.1016/j.jct.2016.11.017>  
<https://www.doi.org/10.1021/je0255271>  
<http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
<https://www.doi.org/10.1016/j.fluid.2014.12.012>  
<http://webbook.nist.gov/cgi/cbook.cgi?ID=C4394858&Units=SI>  
<https://www.doi.org/10.1021/je700036v>  
<https://www.doi.org/10.1021/je060475r>  
<https://www.doi.org/10.1016/j.jct.2009.10.010>  
<https://www.doi.org/10.1016/j.jct.2012.04.016>  
<https://www.doi.org/10.1016/j.jct.2016.02.006>  
<https://www.doi.org/10.1021/je7002513>  
<https://www.doi.org/10.1021/je7005665>  
<https://www.doi.org/10.1021/je100081v>  
<https://www.doi.org/10.1021/je020208v>  
<https://www.doi.org/10.1021/je800999s>  
[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
<https://www.doi.org/10.1021/je060269j>  
<https://www.doi.org/10.1021/je4004417>

## Legend

cpl:	Liquid phase heat capacity
dvisc:	Dynamic viscosity
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rfi:	Refractive Index
tb:	Normal Boiling Point Temperature

Latest version available from:

<https://www.chemeo.com/cid/98-386-6/N-Formylmorpholine.pdf>

Generated by Cheméo on 2024-04-17 02:48:43.437900633 +0000 UTC m=+15611372.358477946.

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