

# 2-Imidazolidinone, 1,3-dimethyl-

<b>Other names:</b>	1,3-Dimethyl-2-imidazolidone 1,3-Dimethylethyleneurea 1,3-Dimethylimidazolidin-2-one 1,3-Dimethylimidazolidinone 1,3-dimethyl-2-imidazolidinone Dimethylethyleneurea N,N'-Dimethyl-2-imidazolidinone N,N'-dimethylethyleneurea Rhonite 1
<b>Inchi:</b>	InChI=1S/C5H10N2O/c1-6-3-4-7(2)5(6)8/h3-4H2,1-2H3
<b>InchiKey:</b>	CYSGHNMQYZDMIA-UHFFFAOYSA-N
<b>Formula:</b>	C5H10N2O
<b>SMILES:</b>	CN1CCN(C)C1=O
<b>Mol. weight [g/mol]:</b>	114.15
<b>CAS:</b>	80-73-9

## Physical Properties

Property code	Value	Unit	Source
affp	918.40	kJ/mol	NIST Webbook
basg	886.00	kJ/mol	NIST Webbook
log10ws	0.30		Crippen Method
logp	-0.016		Crippen Method
mcvol	91.980	ml/mol	McGowan Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hvapt	54.30	kJ/mol	426.50	NIST Webbook
hvapt	48.50	kJ/mol	426.50	NIST Webbook

rho	1064.86	kg/m <sup>3</sup>	283.15	Excess molar volumes of binary mixtures of 1,3-dimethylimidazolidin-2-one with an alkan-1-ol at the temperatures 283.15 K, 298.15 K, and 313.15 K
rho	1051.59	kg/m <sup>3</sup>	298.15	Excess molar volumes of binary mixtures of 1,3-dimethylimidazolidin-2-one with an alkan-1-ol at the temperatures 283.15 K, 298.15 K, and 313.15 K
rho	1038.38	kg/m <sup>3</sup>	313.15	Excess molar volumes of binary mixtures of 1,3-dimethylimidazolidin-2-one with an alkan-1-ol at the temperatures 283.15 K, 298.15 K, and 313.15 K
rho	1069.30	kg/m <sup>3</sup>	278.15	D2O H2O solvent isotope effects on the volumetric properties of aqueous 1,3-dimethyl-2-imidazolidinone between (278.15 and 318.15) K
rho	1060.43	kg/m <sup>3</sup>	288.15	D2O H2O solvent isotope effects on the volumetric properties of aqueous 1,3-dimethyl-2-imidazolidinone between (278.15 and 318.15) K
rho	1051.60	kg/m <sup>3</sup>	298.15	D2O H2O solvent isotope effects on the volumetric properties of aqueous 1,3-dimethyl-2-imidazolidinone between (278.15 and 318.15) K
rho	1042.79	kg/m <sup>3</sup>	308.15	D2O H2O solvent isotope effects on the volumetric properties of aqueous 1,3-dimethyl-2-imidazolidinone between (278.15 and 318.15) K

rhoI	1038.38	kg/m3	313.15	D2O H2O solvent isotope effects on the volumetric properties of aqueous 1,3-dimethyl-2-imidazolidinone between (278.15 and 318.15) K
rhoI	1034.00	kg/m3	318.15	D2O H2O solvent isotope effects on the volumetric properties of aqueous 1,3-dimethyl-2-imidazolidinone between (278.15 and 318.15) K
rhoI	1058.00	kg/m3	293.15	Investigation of the Solubilities of Carbon Dioxide in Some Low Volatile Solvents and Their Thermodynamic Properties

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	380.20	K	2.30	NIST Webbook
tbrp	382.50 ± 0.50	K	2.40	NIST Webbook

## Sources

- (Liquid + liquid) equilibria of (heptane, or hexane, or cyclohexane + toluene + Investigation of the Solubilities of Binary Systems Dioxide in Some Low Volatile Solvents and Their Thermodynamic Properties: Solubility and thermodynamic properties of SO<sub>2</sub> in three low volatile liquids: NIST Webbook: <https://www.doi.org/10.1016/j.jct.2009.11.009>
- Crippen Method: <https://www.doi.org/10.1021/acs.jced.5b00893>
- Enthalpy-related interaction parameters in H/D isotopically distinguishable Molecules: Methods of tetramethylurea cyclic derivatives at 298.15 K: Excess molar volumes of binary mixtures of D<sub>2</sub>O and H<sub>2</sub>O: D<sub>2</sub>O H<sub>2</sub>O solvent isotope effects on the volumetric properties of aqueous 1,3-dimethyl-2-imidazolidinone between (278.15 and 318.15) K: [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)
- <https://www.doi.org/10.1016/j.jct.2016.05.004>
- <http://webbook.nist.gov/cgi/cbook.cgi?ID=C80739&Units=SI>
- <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
- <https://www.doi.org/10.1016/j.tca.2011.05.019>
- <http://link.springer.com/article/10.1007/BF02311772>
- <https://www.doi.org/10.1016/j.jct.2006.09.017>
- <https://www.doi.org/10.1016/j.jct.2009.06.019>

# Legend

<b>affp:</b>	Proton affinity
<b>basg:</b>	Gas basicity
<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>rho:</b>	Liquid Density
<b>tbrp:</b>	Boiling point at reduced pressure

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