

# (E)-N-Isopropylacetaldimine

**Inchi:** InChI=1S/C5H11N/c1-4-6-5(2)3/h4-5H,1-3H3  
**InchiKey:** MXOWOGWCSFJPNN-UHFFFAOYSA-N  
**Formula:** C5H11N  
**SMILES:** CC=NC(C)C  
**Mol. weight [g/mol]:** 85.15  
**CAS:** 27719-14-8

## Physical Properties

Property code	Value	Unit	Source
hf	-69.59	kJ/mol	Joback Method
hvap	29.65	kJ/mol	Joback Method
log10ws	-1.19		Crippen Method
logp	1.486		Crippen Method
mcvol	86.990	ml/mol	McGowan Method
pc	3103.64	kPa	Joback Method
tb	390.04	K	Joback Method
tc	583.65	K	Joback Method

## Sources

**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C27719148&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307I>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

## Legend

**hf:** Enthalpy of formation at standard conditions  
**hvap:** Enthalpy of vaporization at standard conditions  
**log10ws:** Log10 of Water solubility in mol/l

<b>logp:</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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