

# Pyridine, 2-chloro-6-methyl-

<b>Other names:</b>	2-Picoline, 6-chloro- 2-Chloro-6-methylpyridine 2-Chloro-6-picoline 6-Chloro-2-picoline 2-Cl-6-(CH <sub>3</sub> )-pyridine 6-chloro-2-methylpyridine
<b>Inchi:</b>	InChI=1S/C6H6ClN/c1-5-3-2-4-6(7)8-5/h2-4H,1H3
<b>InchiKey:</b>	GXZDYRYYNXYPMQ-UHFFFAOYSA-N
<b>Formula:</b>	C <sub>6</sub> H <sub>6</sub> ClN
<b>SMILES:</b>	Cc1cccc(Cl)n1
<b>Mol. weight [g/mol]:</b>	127.57
<b>CAS:</b>	18368-63-3

## Physical Properties

Property code	Value	Unit	Source
affp	908.00	kJ/mol	NIST Webbook
basg	876.20	kJ/mol	NIST Webbook
log10ws	-2.40		Crippen Method
logp	2.043		Crippen Method
mcvol	93.860	ml/mol	McGowan Method
tb	456.70	K	NIST Webbook

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	339.20	K	1.30	NIST Webbook
tbrp	333.20	K	1.00	NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C18368633&Units=SI>

## Legend

**affp:** Proton affinity  
**basg:** Gas basicity  
**log10ws:** Log10 of Water solubility in mol/l  
**logp:** Octanol/Water partition coefficient  
**mcvol:** McGowan's characteristic volume  
**tb:** Normal Boiling Point Temperature  
**tbrp:** Boiling point at reduced pressure

Latest version available from:

<https://www.chemeo.com/cid/24-645-9/Pyridine-2-chloro-6-methyl.pdf>

Generated by Cheméo on 2024-04-26 17:34:24.598131365 +0000 UTC m=+16442113.518708687.

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