

# 2,4-Dicyanopyridine

Other names:	Pyridine-2,4-dicarbonitrile 2,4-Pyridinedicarbonitrile
Inchi:	InChI=1S/C7H3N3/c8-4-6-1-2-10-7(3-6)5-9/h1-3H
InchiKey:	HLAQMFURMNTLW-UHFFFAOYSA-N
Formula:	C7H3N3
SMILES:	N#Cc1ccnc(C#N)c1
Mol. weight [g/mol]:	129.12
CAS:	29181-50-8

## Physical Properties

Property code	Value	Unit	Source
log10ws	-1.94		Crippen Method
logp	0.825		Crippen Method
mcvol	98.470	ml/mol	McGowan Method

## Sources

Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
McGowan Method:	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C29181508&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C29181508&amp;Units=SI</a>

## Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume

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