

1H-Indole, 1,2-dimethyl-

Other names:	Indole, 1,2-dimethyl- N-Methyl-2-methylindole 1,2-Dimethyl-1H-indole 1,2-Dimethylindole
Inchi:	InChI=1S/C10H11N/c1-8-7-9-5-3-4-6-10(9)11(8)2/h3-7H,1-2H3
InchiKey:	BJMUOUXGBFNLSN-UHFFFAOYSA-N
Formula:	C10H11N
SMILES:	Cc1cc2ccccc2n1C
Mol. weight [g/mol]:	145.20
CAS:	875-79-6

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.04		Crippen Method
logp	2.487		Crippen Method
mcvol	122.820	ml/mol	McGowan Method
rinpol	244.42		NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	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=C875796&Units=SI

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/79-612-5/1H-Indole-1-2-dimethyl.pdf>

Generated by Cheméo on 2025-12-05 09:30:18.933713328 +0000 UTC m=+4675216.463753992.

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