

Cinchonidine

Other names:	Cinchonan-9-ol, (8«alpha»,9R)- «alpha»-Quinidine (-)-Cinchonidine Cinchovatine L-Cinchonidine 2-Quinuclidinemethanol, «alpha»-4-quinolyl-5-vinyl- (8«alpha»,9R)-Cinchonan-9-ol (8S,9R)-Cinchonidine NSC 5364
Inchi:	InChI=1S/C19H22N2O/c1-2-13-12-21-10-8-14(13)11-18(21)19(22)16-7-9-20-17-6-4-3-5-
InchiKey:	KMPWYEUPVWOPIM-AREIWFIDSA-N
Formula:	C19H22N2O
SMILES:	<chem>C=CC1CN2CCC1CC2C(O)c1ccnc2ccccc12</chem>
Mol. weight [g/mol]:	294.39
CAS:	485-71-2

Physical Properties

Property code	Value	Unit	Source
chs	-10652.00	kJ/mol	NIST Webbook
log10ws	-4.76		Crippen Method
logp	3.165		Crippen Method
mcvol	235.160	ml/mol	McGowan Method
rinpole	2600.00		NIST Webbook
rinpole	2580.00		NIST Webbook
rinpole	2598.00		NIST Webbook

Sources

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=C485712&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

chs:	Standard solid enthalpy of combustion
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.cheméo.com/cid/68-707-2/Cinchonidine.pdf>

Generated by Cheméo on 2024-04-27 08:52:01.877654443 +0000 UTC m=+16497170.798231758.

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