2H-quinolizine-1-methanol, octahydro-, (1s-cis)-

Other names: (+)-Epilupinine
1-Epilupinine
Epi-Lupinine

Inchi: InChI=1S/C10H19NO/c12-8-9-4-3-7-11-6-2-1-5-10(9)11/h9-10,12H,1-8H2/t9-,10?/m1/s1

InchiKey: HDVAWXXJVMJBAR-RGURZIINSA-N

Formula: C10H19NO

SMILES: OC[C@H]1CCCN2CCCCC12

Mol. weight [g/mol]: 169.27

Physical Properties

Property code	Value	Unit	Source
logp	1.243		Crippen Method
mcvol	145.890	ml/mol	McGowan Method
rinpol	1415.00		NIST Webbook
rinpol	1400.00		NIST Webbook
rinpol	1418.00		NIST Webbook
rinpol	1415.00		NIST Webbook
rinpol	1415.00		NIST Webbook
rinpol	1416.00		NIST Webbook
rinpol	1415.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C486715&Units=SI

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

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