1',3',5',6'-Tetramethylspiro[cyclopentane-1,2'-(2,3-

Inchi: InChI=1S/C15H22N2/c1-11-9-13-14(10-12(11)2)17(4)15(16(13)3)7-5-6-8-15/h9-10H,5-8H
InchiKey: WWHRITWAHDSGEL-UHFFFAOYSA-N
Formula: C15H22N2
SMILES: Cc1cc2c(cc1C)N(C)C1(CCCC1)N2C
Mol. weight [g/mol]: 230.35
CAS: 75751-20-1

Physical Properties

Property code	Value	Unit	Source
ie	6.35	eV	NIST Webbook
log10ws	-3.89		Crippen Method
logp	3.460		Crippen Method
mcvol	196.690	ml/mol	McGowan Method

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C75751201&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

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

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