7-acetyl-5,6-dimethyl-2,3-dihydro-(1H)-pyrrolizine

Inchi: InChI=1S/C11H15NO/c1-7-8(2)12-6-4-5-10(12)11(7)9(3)13/h4-6H2,1-3H3
InchiKey: BJTRWJSVQKOMJY-UHFFFAOYSA-N
Formula: C11H15NO
SMILES: CC(=O)c1c(C)c(C)n2c1CCC2
Mol. weight [g/mol]: 177.24

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.62		Crippen Method
logp	2.254		Crippen Method
mcvol	147.080	ml/mol	McGowan Method
rinpol	1778.00		NIST Webbook
rinpol	1778.00		NIST Webbook

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R231106&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

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

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