

L-Pipecolic acid, N(O,S)-ethoxycarbonyl, (S)-(+)-3-methyl-2-butyl ester

Inchi: InChI=1S/C14H25NO4/c1-5-18-14(17)15-9-7-6-8-12(15)13(16)19-11(4)10(2)3/h10-12H,5
InchiKey: VFRHGYSFWMKXKS-PXYINDEMSA-N
Formula: C14H25NO4
SMILES: CCOC(=O)N1CCCCC1C(=O)OC(C)C(C)C
Mol. weight [g/mol]: 271.35

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.83		Crippen Method
logp	2.585		Crippen Method
mcvol	222.120	ml/mol	McGowan Method
rinpol	1755.10		NIST Webbook
rinpol	1755.10		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=R502186&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Legend

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

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