

Pyridine, 2,4,6-triethyl

Inchi: InChI=1S/C11H17N/c1-4-9-7-10(5-2)12-11(6-3)8-9/h7-8H,4-6H2,1-3H3
InchiKey: XVMNXGCQCHNBEN-UHFFFAOYSA-N
Formula: C11H17N
SMILES: CCc1cc(CC)nc(CC)c1
Mol. weight [g/mol]: 163.26
CAS: 50286-76-5

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.65		Crippen Method
logp	2.769		Crippen Method
mcvol	152.070	ml/mol	McGowan Method
rinpol	1212.00		NIST Webbook
rinpol	1212.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
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=C50286765&Units=SI>

Legend

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.chemeo.com/cid/29-876-8/Pyridine-2-4-6-triethyl.pdf>

Generated by Cheméo on 2024-04-19 15:57:32.876585061 +0000 UTC m=+15831501.797162376.

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