2-pentyl-4-methyl-5-ethyloxazole

Inchi: InChI=1S/C11H19NO/c1-4-6-7-8-11-12-9(3)10(5-2)13-11/h4-8H2,1-3H3
InchiKey: CJVUTPBIHCDRQD-UHFFFAOYSA-N
Formula: C11H19NO
SMILES: CCCCCc1nc(C)c(CC)o1
Mol. weight [g/mol]: 181.27

Physical Properties

Property code	Value	Unit	Source
log10ws	-8.31		Crippen Method
logp	3.278		Crippen Method
mcvol	162.240	ml/mol	McGowan Method
rinpol	1263.00		NIST Webbook
rinpol	1263.00		NIST Webbook

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R161443&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.