2-Furancarboxamide, N-hexyl-

Inchi: InChI=1S/C11H17NO2/c1-2-3-4-5-8-12-11(13)10-7-6-9-14-10/h6-7,9H,2-5,8H2,1H3,(H,1
InchiKey: QEALKLTYTGLOIV-UHFFFAOYSA-N
Formula: C11H17NO2
SMILES: CCCCCCNC(=O)c1ccco1
Mol. weight [g/mol]: 195.26

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.64		Crippen Method
logp	2.590		Crippen Method
mcvol	163.810	ml/mol	McGowan Method
rinpol	1664.00		NIST Webbook
rinpol	1664.00		NIST Webbook

Sources

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=U407245&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

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