2-Furancarboxamide, N-dodecyl-

Inchi: InChI=1S/C17H29NO2/c1-2-3-4-5-6-7-8-9-10-11-14-18-17(19)16-13-12-15-20-16/h12-13
InchiKey: VMKSSOXODLIIOE-UHFFFAOYSA-N
Formula: C17H29NO2
SMILES: CCCCCCCCCCCCNC(=O)c1ccco1
Mol. weight [g/mol]: 279.42

Physical Properties

Property code	Value	Unit	Source
log10ws	-10.15		Crippen Method
logp	4.930		Crippen Method
mcvol	248.350	ml/mol	McGowan Method
rinpol	2295.00		NIST Webbook
rinpol	2295.00		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=U407252&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

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.