2-Furancarboxamide, N-octadecyl-

Inchi: InChI=1S/C23H41NO2/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17-20-24-23(25)22-19-
InchiKey: RGMFNNIZXFLPPX-UHFFFAOYSA-N
Formula: C23H41NO2
SMILES: CCCCCCCCCCCCCCCCCCNC(=O)c1ccco1
Mol. weight [g/mol]: 363.58

Physical Properties

Property code	Value	Unit	Source
log10ws	-12.67		Crippen Method
logp	7.271		Crippen Method
mcvol	332.890	ml/mol	McGowan Method
rinpol	2929.00		NIST Webbook
rinpol	2929.00		NIST Webbook

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U407255&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.