Furan-2-carboxamide, N,N-diundecyl-

Inchi: InChI=1S/C27H49NO2/c1-3-5-7-9-11-13-15-17-19-23-28(27(29)26-22-21-25-30-26)24-20
InchiKey: YKPPDGQZLXOLPD-UHFFFAOYSA-N
Formula: C27H49NO2
SMILES: CCCCCCCCCCCN(CCCCCCCCCCC)C(=O)c1ccco1
Mol. weight [g/mol]: 419.68

Physical Properties

Property code	Value	Unit	Source
log10ws	-13.72		Crippen Method
logp	8.783		Crippen Method
mcvol	389.250	ml/mol	McGowan Method
rinpol	3084.00		NIST Webbook
rinpol	3084.00		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=U308212&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.