2,6-Pyridinedicarboxylic acid, heptadecyl 3-methylbutyl ester

Inchi: InChI=1S/C29H49NO4/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-17-18-23-33-28(31)26-20-1

InchiKey: LHWFCUGADFVUPV-UHFFFAOYSA-N

Formula: C29H49NO4

SMILES: CCCCCCCCCCCCCCCCCOC(=O)c1cccc(C(=O)OCCC(C)C)n1

Mol. weight [g/mol]: 475.70

Physical Properties

Property code	Value	Unit	Source
log10ws	-9.85		Crippen Method
logp	8.313		Crippen Method
mcvol	420.570	ml/mol	McGowan Method
rinpol	3260.00		NIST Webbook
rinpol	3260.00		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=U368329&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.