

2,6-Pyridinedicarboxylic acid, 2-octyl tetradecyl ester

Inchi: InChI=1S/C29H49NO4/c1-4-6-8-10-11-12-13-14-15-16-17-19-24-33-28(31)26-22-20-23-2
InchiKey: LNAMEXWCLVGLX-UHFFFAOYSA-N
Formula: C₂₉H₄₉NO₄
SMILES: CCCCCCCCCCCCCCOC(=O)c1cccc(C(=O)OC(C)CCCCC)n1
Mol. weight [g/mol]: 475.70

Physical Properties

Property code	Value	Unit	Source
log10ws	-10.21		Crippen Method
logp	8.455		Crippen Method
mcvol	420.570	ml/mol	McGowan Method
rinpol	3193.00		NIST Webbook

Sources

McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U368310&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

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