

2,6-Pyridinedicarboxylic acid, dibutyl ester

Inchi: InChI=1S/C15H21NO4/c1-3-5-10-19-14(17)12-8-7-9-13(16-12)15(18)20-11-6-4-2/h7-9H,
InchiKey: SJGUKBIVRJZPJZ-UHFFFAOYSA-N
Formula: C15H21NO4
SMILES: CCCOC(=O)c1cccc(C(=O)OCCCC)n1
Mol. weight [g/mol]: 279.33

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.23		Crippen Method
logp	2.995		Crippen Method
mcvol	223.310	ml/mol	McGowan Method
rinpole	2068.00		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=U369100&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
rinpole: Non-polar retention indices

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