2,6-Pyridinedicarboxylic acid, isobutyl 2-methylpentyl ester

Inchi: InChI=1S/C17H25NO4/c1-5-7-13(4)11-22-17(20)15-9-6-8-14(18-15)16(19)21-10-12(2)3/h

InchiKey: NNPIMAXRPRVLCH-UHFFFAOYSA-N

Formula: C17H25NO4

SMILES: CCCC(C)COC(=O)c1cccc(C(=O)OCC(C)C)n1

Mol. weight [g/mol]: 307.38

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.59		Crippen Method
logp	3.487		Crippen Method
mcvol	251.490	ml/mol	McGowan Method
rinpol	2151.00		NIST Webbook
rinpol	2151.00		NIST Webbook

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U369082&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.