

2,6-Pyridinedicarboxylic acid, pentyl undecyl ester

Inchi:	InChI=1S/C23H37NO4/c1-3-5-7-8-9-10-11-12-14-19-28-23(26)21-17-15-16-20(24-21)22
InchiKey:	ADYQWZDABDPFKV-UHFFFAOYSA-N
Formula:	C23H37NO4
SMILES:	CCCCCCCCCOC(=O)c1cccc(C(=O)OCCCC)n1
Mol. weight [g/mol]:	391.54

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.58		Crippen Method
logp	6.116		Crippen Method
mcvol	336.030	ml/mol	McGowan Method
rinsol	2848.00		NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U368289&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

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

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