

# Pipecolic acid, N-propoxycarbonyl-, dodecyl ester

<b>Inchi:</b>	InChI=1S/C22H41NO4/c1-3-5-6-7-8-9-10-11-12-15-19-26-21(24)20-16-13-14-17-23(20)2
<b>InchiKey:</b>	FICWGBKONFFNEM-UHFFFAOYSA-N
<b>Formula:</b>	C22H41NO4
<b>SMILES:</b>	CCCCCCCCCCCCOC(=O)C1CCCN1C(=O)OCCC
<b>Mol. weight [g/mol]:</b>	383.57

## Physical Properties

Property code	Value	Unit	Source
log10ws	-6.31		Crippen Method
logp	5.852		Crippen Method
mcvol	334.840	ml/mol	McGowan Method
rinpole	2683.00		NIST Webbook
rinpole	2683.00		NIST Webbook

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U393003&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U393003&amp;Units=SI</a>

## Legend

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

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