Phenylphosphonic acid, butyl dodecyl ester

Inchi: InChI=1S/C22H39O3P/c1-3-5-7-8-9-10-11-12-13-17-21-25-26(23,24-20-6-4-2)22-18-15-1
InchiKey: XQDRHBJCMWPCLO-UHFFFAOYSA-N
Formula: C22H39O3P
SMILES: CCCCCCCCCCCCOP(=O)(OCCCC)c1ccccc1
Mol. weight [g/mol]: 382.52

Physical Properties

Property code	Value	Unit	Source
log10ws	-12.97		Crippen Method
logp	7.259		Crippen Method
mcvol	335.150	ml/mol	McGowan Method
rinpol	2697.00		NIST Webbook
rinpol	2697.00		NIST Webbook

Sources

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=U393242&Units=SI

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.