

2-(P-n-nonylphenoxy)-3-tridecyl pyrazine

Inchi:	InChI=1S/C32H52N2O/c1-3-5-7-9-11-12-13-14-16-18-20-22-31-32(34-28-27-33-31)35-3
InchiKey:	JGTVXBDSQFGHLZ-UHFFFAOYSA-N
Formula:	C32H52N2O
SMILES:	CCCCCCCCCCCCc1nccnc1Oc1ccc(CCCCCCCCC)cc1
Mol. weight [g/mol]:	480.77
CAS:	2879-00-7

Physical Properties

Property code	Value	Unit	Source
log10ws	-11.78		Crippen Method
logp	10.415		Crippen Method
mcvol	440.050	ml/mol	McGowan Method

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=C2879007&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume

Latest version available from:

<https://www.chemeo.com/cid/86-205-9/2-P-n-nonylphenoxy-3-tridecyl-pyrazine.pdf>

Generated by Cheméo on 2024-04-24 19:00:33.101263874 +0000 UTC m=+16274482.021841186.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.