

Phensuximide

Other names:	2,5-Pyrrolidinedione, 1-methyl-3-phenyl-Succinimide, N-methyl-2-phenyl-Epimid Lifene Milontin Milonton Mirontin N-Methyl-«alpha»-phenylsuccinimide N-Methyl-2-phenylsuccinimide Phensuximid PM 334 Succitimal 1-Methyl-3-Phenylsuccinimide 1-Methyl-3-phenylpyrrolidin-2,5-dione Phenylsuximide (. +/- .)-N-Methyl-2-phenylsuccinimide
Inchi:	InChI=1S/C11H11NO2/c1-12-10(13)7-9(11(12)14)8-5-3-2-4-6-8/h2-6,9H,7H2,1H3
InchiKey:	WLWFNJKHKGIJNW-UHFFFAOYSA-N
Formula:	C11H11NO2
SMILES:	CN1C(=O)CC(c2ccccc2)C1=O
Mol. weight [g/mol]:	189.21
CAS:	86-34-0

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.51		Crippen Method
logp	1.159		Crippen Method
mcpvol	144.350	ml/mol	McGowan Method
rinpol	1634.00		NIST Webbook
rinpol	1613.00		NIST Webbook
rinpol	1603.00		NIST Webbook
rinpol	1620.00		NIST Webbook
rinpol	1601.00		NIST Webbook
rinpol	1607.00		NIST Webbook
rinpol	1620.00		NIST Webbook
rinpol	1620.00		NIST Webbook
rinpol	1611.00		NIST Webbook

rinpol	284.01	NIST Webbook
rinpol	1634.00	NIST Webbook
rinpol	1607.00	NIST Webbook
rinpol	1620.00	NIST Webbook
rinpol	284.01	NIST Webbook
rinpol	1618.00	NIST Webbook
rinpol	1601.00	NIST Webbook
rinpol	1607.00	NIST Webbook
rinpol	1607.00	NIST Webbook
rinpol	1618.00	NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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=C86340&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

Latest version available from:

<https://www.chemeo.com/cid/91-027-1/Phensuximide.pdf>

Generated by Cheméo on 2024-05-03 17:04:52.044179453 +0000 UTC m=+17045140.964756763.

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