

Pyrrolidine, 1-acetyl-

Other names:	N-Acetylpyrrolidine 1-Acetylpyrrolidine N-Acetypyrrolidine Pyrrolidine, N-acetyl Acetylpyrrolidine
Inchi:	InChI=1S/C6H11NO/c1-6(8)7-4-2-3-5-7/h2-5H2,1H3
InchiKey:	LNWWQYYLZVZXKS-UHFFFAOYSA-N
Formula:	C6H11NO
SMILES:	CC(=O)N1CCCC1
Mol. weight [g/mol]:	113.16
CAS:	4030-18-6

Physical Properties

Property code	Value	Unit	Source
affp	925.40	kJ/mol	NIST Webbook
basg	894.40	kJ/mol	NIST Webbook
log10ws	-0.57		Crippen Method
logp	0.629		Crippen Method
mcvol	96.090	ml/mol	McGowan Method
rinpol	1128.00		NIST Webbook
rinpol	1128.00		NIST Webbook
ripol	1766.00		NIST Webbook
ripol	1766.00		NIST Webbook

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	379.00 ± 1.00	K	2.00	NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C4030186&Units=SI>
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>

Legend

affp: Proton affinity
basg: Gas basicity
log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
rinpol: Non-polar retention indices
ripol: Polar retention indices
tbrp: Boiling point at reduced pressure

Latest version available from:

<https://www.chemeo.com/cid/45-707-7/Pyrrolidine-1-acetyl.pdf>

Generated by Cheméo on 2024-04-24 10:10:07.352487946 +0000 UTC m=+16242656.273065258.

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