

2-Pyrimidinamine, 4,6-dimethyl-

Other names:	2-amino-4,6-dimethylpyrimidine 4,6-Dimethyl-pyrimidin-2-ylamine 4,6-dimethyl-2-aminopyrimidine 4,6-dimethyl-2-pyrimidinamine Acetylacetoneguanidine Pyrimidine, 2-amino-4,6-dimethyl-
Inchi:	InChI=1S/C6H9N3/c1-4-3-5(2)9-6(7)8-4/h3H,1-2H3,(H2,7,8,9)
InchiKey:	IDQNBVFPZMCDDN-UHFFFAOYSA-N
Formula:	C6H9N3
SMILES:	Cc1cc(C)nc(N)n1
Mol. weight [g/mol]:	123.16
CAS:	767-15-7

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.59		Crippen Method
logp	0.676		Crippen Method
mcvol	101.580	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
rhos	1240.00	kg/m3	298.15	Energetics of aminomethylpyrimidines: An examination of the aromaticity of nitrogen heteromonocyclic derivatives

Sources

McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C767157&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Energetics of aminomethylpyrimidines: <https://www.doi.org/10.1016/j.jct.2013.03.010>
An examination of the aromaticity of nitrogen heteromonocyclic derivatives:

Legend

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

Latest version available from:

<https://www.chemeo.com/cid/44-064-2/2-Pyrimidinamine-4-6-dimethyl.pdf>

Generated by Cheméo on 2025-12-24 09:03:20.370626428 +0000 UTC m=+6315197.900667092.

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