Pyridine, 4-amino-2,6-dichloro-

Inchi: InChI=1S/C5H4Cl2N2/c6-4-1-3(8)2-5(7)9-4/h1-2H,(H2,8,9)
InchiKey: WAEZOSSWRXDWAX-UHFFFAOYSA-N
Formula: C5H4Cl2N2
SMILES: Nc1cc(Cl)nc(Cl)c1
Mol. weight [g/mol]: 163.01

Physical Properties

Property code	Value	Unit	Source
hvap	64.11	kJ/mol	Cheméo Relay (1.0)
logp	1.971		Crippen Method
mcvol	101.990	ml/mol	McGowan Method
tb	538.51	K	Cheméo Relay (1.0)
tf	450.76	K	Cheméo Relay (1.0)

Sources

Cheméo Relay (1.0, beta):
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C2587022&Units=SI
McGowan Method: http://link.springer.com/article/10.1007/BF02311772
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):

Legend

hvap: Enthalpy of vaporization at standard conditions
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
tb: Normal Boiling Point Temperature
tf: Normal melting (fusion) point