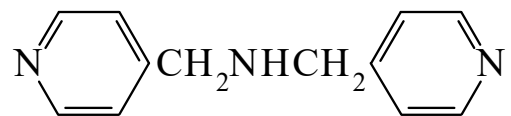




>>Entry Name: **Gapicomine**, INN

Synonym(s): *N*-(4-Pyridinylmethyl)-4-pyridinemethanamine, 9Cl. 4,4'-(Iminodimethylene)dipyridine, 8Cl. 1,1'-Di(4-pyridyl)dimethylamine. Bis(4-pyridylmethyl)amine



CAS Registry Number: 1539-39-5

Molecular Formula: C₁₂H₁₃N₃

Molecular Weight: 199.255

Accurate Mass: 199.110947

Percentage Composition: C 72.34%; H 6.58%; N 21.09%

Biological Use/Importance: Sedative, peripheral vasodilator

Boiling Point: Bp_{0.23} 164-166°

Calculated Log P Data: Log P -0.18 (uncertain value) (calc)

>>Derivative: Hydrochloride (1:3)

CAS Registry Number: 1539-38-4

Physical Description: Cryst. (H₂O)

Melting Point: Mp 228-230°

>>Derivative: Citrate salt

Synonym(s): Bicordin. CCB

CAS Registry Number: 24631-38-7

Melting Point: Mp 191°

References:

Biniecki, S. *et al.*, *Ann. Pharm. Fr.*, 1964, **22**, 685-687; *CA*, **63**, 2949a, (*synth, pharmacol*)

U.K. Pat., 1967, *Polfa*, 1 058 356; *CA*, **66**, 115607, (*synth, pharmacol, deriv*)

Kubikowski, P. *et al.*, *Diss. Pharm. Pharmacol.*, 1968, **20**, 17, (*pharmacol*)

Martindale, The Extra Pharmacopoeia, 28th/29th edn., *Pharmaceutical Press*, 1982, 9234