

Certificate of Analysis

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Product Name: BD 1008 dihydrobromide

Catalog No.: 0511

Batch No.: 2

CAS Number: 138356-09-9

IUPAC Name: *N*-[2-(3,4-Dichlorophenyl)ethyl]-*N*-methyl-2-(1-pyrrolidinyl)ethylamine dihydrobromide

1. PHYSICAL AND CHEMICAL PROPERTIES

Batch Molecular Formula: C₁₅H₂₂Cl₂N₂.2HBr

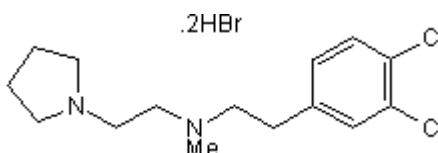
Batch Molecular Weight: 463.08

Physical Appearance: White solid

Solubility: water to 50 mM

Storage: Store at -20°C

Batch Molecular Structure:



2. ANALYTICAL DATA

HPLC: Shows 99.1% purity

¹H NMR: Consistent with structure

Mass Spectrum: Consistent with structure

Microanalysis:

	Carbon Hydrogen Nitrogen		
Theoretical	38.9	5.22	6.05
Found	38.7	5.14	6.16

Caution - Not Fully Tested • Research Use Only • Not For Human or Veterinary Use

Product Information

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IUPAC Name: *N*-[2-(3,4-Dichlorophenyl)ethyl]-*N*-methyl-2-(1-pyrrolidinyl)ethylamine dihydrobromide

Description:

Potent and selective σ -ligand (K_i against [3 H]-(+)-3-PPP = 0.34 nM).

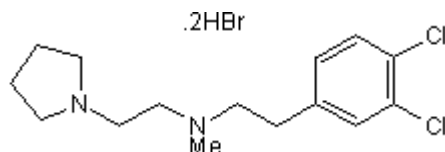
Physical and Chemical Properties:

Batch Molecular Formula: $C_{15}H_{22}Cl_2N_2 \cdot 2HBr$

Batch Molecular Weight: 463.08

Physical Appearance: White solid

Batch Molecular Structure:



Storage: Store at -20°C

Solubility & Usage Info:

water to 50 mM

Stability and Solubility Advice:

Some solutions can be difficult to obtain and can be encouraged by rapid stirring, sonication or gentle warming (in a 45-60°C water bath).

Information concerning product stability, particularly in solution, has rarely been reported and in most cases we can only offer a general guide. Our standard recommendations are:

SOLIDS: Provided storage is as stated on the product label and the vial is kept tightly sealed, the product can be stored for up to 6 months from date of receipt.

SOLUTIONS: We recommend that stock solutions, once prepared, are stored aliquoted in tightly sealed vials at -20°C or below and used within 1 month. Wherever possible solutions should be made up and used on the same day.

References:

De Costa et al (1992) Synthesis, characterization, and biological evaluation of a novel class of *N*-(arylethyl)-*N*-alkyl-2-(1-pyrrolidinyl)ethylamines - structural requirements and binding affinity at the σ receptor. *J.Med.Chem.* **35** 38. PMID: 1310114.

Monnet et al (1996) Differentiation of σ ligand-activated receptor subtypes that modulate NMDA-evoked [3 H]-noradrenaline release in rat hippocampal slices. *Br.J.Pharmacol.* **119** 65. PMID: 8872358.

Whittemore et al (1997) Antagonism of N-methyl-D-aspartate receptors by σ site ligands: potency, subtype-selectivity and mechanisms of inhibition. *J.Pharmacol.Exp.Ther.* **282** 326. PMID: 9223571.

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