

Certificate of Analysis

Product Name: CP 809101 hydrochloride

Catalog No.: 3041

Batch No.: 2

CAS Number: 1215721-40-6

IUPAC Name: 2-[(3-Chlorophenyl)methoxy]-6-(1-piperazinyl)pyrazine hydrochloride

1. PHYSICAL AND CHEMICAL PROPERTIES

Batch Molecular Formula: C₁₅H₁₇N₄OCl.HCl

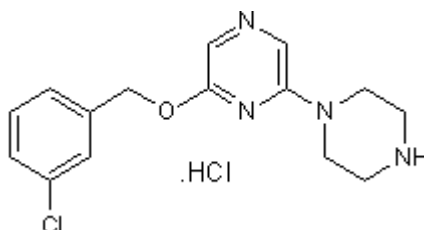
Batch Molecular Weight: 341.24

Physical Appearance: solid

Solubility: water to 20 mM
DMSO to 100 mM

Storage: Desiccate at RT

Batch Molecular Structure:



2. ANALYTICAL DATA

TLC: R_f = 0.47 (Dichloromethane:Ethanol:Ammonia soln. [90:9:1])

HPLC: Shows >99.4% purity

¹H NMR: Consistent with structure

Mass Spectrum: Consistent with structure

Microanalysis:

	Carbon	Hydrogen	Nitrogen
Theoretical	52.8	5.32	16.42
Found	52.84	5.27	16.33

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

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IUPAC Name: 2-[(3-Chlorophenyl)methoxy]-6-(1-piperazinyl)pyrazine hydrochloride

Description:

Potent and selective 5-HT_{2C} receptor agonist (pEC₅₀ values are 9.96, 7.19 and 6.81 for human 5-HT_{2C}, 5-HT_{2B} and 5-HT_{2A} receptors respectively). Displays antipsychotic activity; suppresses condition avoidance responding (CAR) and inhibits PCP and amphetamine-stimulated hyperactivity in rats following subcutaneous administration.

Physical and Chemical Properties:

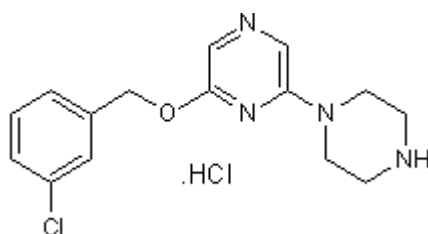
Batch Molecular Formula: C₁₅H₁₇N₄OCl.HCl

Batch Molecular Weight: 341.24

Physical Appearance: solid

Minimum Purity: >99%

Batch Molecular Structure:



References:

Siuciak et al (2007) CP-809,101, a selective 5-HT_{2C} agonist, shows activity in animal models of antipsychotic activity. *Neuropharmacology* **52** 279. PMID: 16949622.

Kalgutkar et al (2007) Genotoxicity of 2-(3-chlorobenzyloxy)-6-(piperazinyl)pyrazine, a novel 5-hydroxytryptamine_{2c} receptor agonist for the treatment of obesity: role of metabolic activation. *Drug Metab.Dispos.* **35** 848. PMID: 17344339.

Jensen et al (2013) Design, synthesis, and pharmacological characterization of *N*- and *O*-substituted 5,6,7,8-tetrahydro-4*H*-isoxazolo [4,5-*d*]azepin-3-ol analogues: novel 5-HT_{2A}/5-HT_{2C} receptor agonists with pro-cognitive properties. *J.Med.Chem* **56** 1211. PMID: 23301527.

Storage: Desiccate at RT

Solubility & Usage Info:

water to 20 mM

DMSO to 100 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.

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