

Certificate of Analysis

Product Name: TC-G 24

Catalog No.: 4353

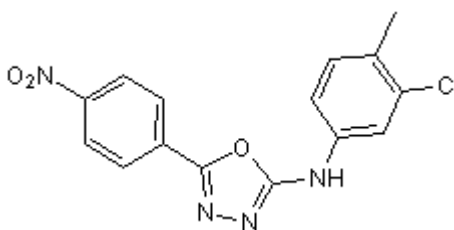
Batch No.: 1

CAS Number: 1257256-44-2

IUPAC Name: *N*-(3-Chloro-4-methylphenyl)-5-(4-nitrophenyl)-1,3,4-oxadiazol-2-amine

1. PHYSICAL AND CHEMICAL PROPERTIES

Batch Molecular Formula: C₁₅H₁₁ClN₄O₃·¼H₂O
Batch Molecular Weight: 335.23
Physical Appearance: Yellow solid
Solubility: DMSO to 10 mM with gentle warming
Storage: Store at +4°C
Batch Molecular Structure:



2. ANALYTICAL DATA

HPLC: Shows 98.8% purity
¹H NMR: Consistent with structure
Mass Spectrum: Consistent with structure

Microanalysis:

| | Carbon | Hydrogen | Nitrogen |
|-------------|--------|----------|----------|
| Theoretical | 53.74 | 3.46 | 16.71 |
| Found | 53.65 | 3.32 | 16.87 |

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

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Description:

Potent GSK-3 β inhibitor (IC₅₀ = 17 nM); displays selectivity for GSK-3 β over CDK2 (22% inhibition at 10 μ M). Increases liver glycogen reserves in rodents. Brain penetrant.

Physical and Chemical Properties:

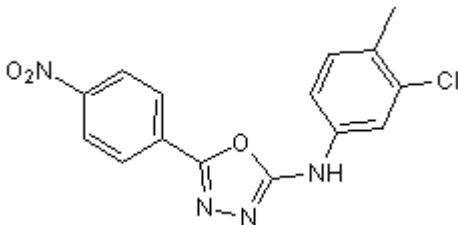
Batch Molecular Formula: C₁₅H₁₁ClN₄O₃·¼H₂O

Batch Molecular Weight: 335.23

Physical Appearance: Yellow solid

Minimum Purity: >98%

Batch Molecular Structure:



Storage: Store at +4°C

Solubility & Usage Info:

DMSO to 10 mM with gentle warming

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:

Khanfar et al (2010) Discovery of novel GSK-3 β inhibitors with potent in vitro and in vivo activities and excellent brain permeability using combined ligand- and structure-based virtual screening. *J.Med.Chem.* **53** 8534. PMID: 21082766.

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