

# **Certificate of Analysis**

Print Date: Jan 13th 2016

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Product Name: DTG Catalog No.: 0841 Batch No.: 3

CAS Number: 97-39-2 EC Number: 202-577-6

IUPAC Name: 1,3-Di-(2-tolyl)guanidine

### 1. PHYSICAL AND CHEMICAL PROPERTIES

**Solubility:** ethanol to 25 mM DMSO to 25 mM

Storage: Store at RT

Batch Molecular Structure:

#### 2. ANALYTICAL DATA

**TLC:**  $R_f = 0.71$  (Dichloromethane:Methanol [9:1])

Melting Point: At 185°C

<sup>1</sup>H NMR: Consistent with structure

Microanalysis: Carbon Hydrogen Nitrogen

Theoretical 75.28 7.16 17.56 Found 75.1 7.19 17.6



## **Product Information**

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CAS Number: 97-39-2 EC Number: 202-577-6

**IUPAC Name:** 1,3-Di-(2-tolyl)guanidine

**Description:** 

Displays high and roughly equal affinity for both  $\sigma_1$  and  $\sigma_2$ 

receptors.

**Physical and Chemical Properties:** 

Batch Molecular Formula: C<sub>15</sub>H<sub>17</sub>N<sub>3</sub> Batch Molecular Weight: 239.32 Physical Appearance: White solid

**Batch Molecular Structure:** 

Storage: Store at RT

Solubility & Usage Info:

ethanol to 25 mM DMSO to 25 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:

Quirion et al (1992) A proposal for the classification of sigma binding sites. TiPS 13 85. PMID: 1315463.

He et al (1993) Synthesis and binding characteristics of potential SPECT imaging agents for σ1 and σ2 binding sites. J.Med.Chem. 36 566. PMID: 8496936.

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