

Certificate of Analysis

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Print Date: Jan 14th 2016

Product Name: 1-Deazaadenosine Catalog No.: 4488 Batch No.: 1

CAS Number: 14432-09-8

Batch Molecular Structure:

IUPAC Name: $3-\beta$ -D-Ribofuranosyl-3*H*-imidazo[4,5-*b*]pyridin-7-amine

1. PHYSICAL AND CHEMICAL PROPERTIES

Batch Molecular Formula: $C_{11}H_{14}N_4O_4$. $^{1/4}H_2O_4$

Batch Molecular Weight:270.75Physical Appearance:White solidSolubility:DMSO to 25 mMStorage:Store at +4°C

NH₂ N N N O N O HO

2. ANALYTICAL DATA

TLC: $R_f = 0.33$ (Ethyl acetate:Methanol [19:1])

HPLC: Shows 98.9% purity

¹H NMR: Consistent with structure

Mass Spectrum: Consistent with structure

Microanalysis: Carbon Hydrogen Nitrogen

Theoretical 48.8 5.4 20.69 Found 48.76 5.3 20.54



Product Information

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IUPAC Name: $3-\beta$ -D-Ribofuranosyl-3*H*-imidazo[4,5-*b*]pyridin-7-amine

Description:

Inhibitor of adenosine deaminase (K_i = 0.66 µM). Demonstrates antitumor activity in a range of leukemia cell lines.

Physical and Chemical Properties:

Batch Molecular Formula: C₁₁H₁₄N₄O₄. ½ H₂O

Batch Molecular Weight: 270.75 Physical Appearance: White solid

Minimum Purity: >98%

Batch Molecular Structure:

Storage: Store at +4°C

Solubility & Usage Info:

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:

Cristalli et al (1987) Improved synthesis and antitumor activity of 1-deazaadenosine. J.Med.Chem. 30 1686. PMID: 3625714.

Cristalli et al (1993) Adenosine deaminase inhibitors: Structure-activity relationships in 1-deazaadenosine and erythro-9-(2-hydroxy-3nonyl)adenine analogues. Drug Dev.Res. 28 253.