

# **Certificate of Analysis**

BDG Synthesis certifies that this reference material meets or exceeds the specifications stated herein.

leil Beare

Neil Beare, PhD, Director 9 September 2016

Name: E-Rilpivirine-d<sub>4</sub>

CAS Number: none

**Structure:** 

**Molecular Weight:**  $C_{22}H_{14}D_4N_6 = 370.44$ 

Lot Number: BDG 16780

**Appearance:** White, crystalline solid

**Purity By HPLC:** 97.6 %

**Isotopic Purity:** Under  $0.5 \% d_0$ 

**Re-test Date:** 9 September 2021

**Storage and Handling:** Temperature: refrigerate for prolonged storage; may be handled and shipped at

ambient temperature.

Humidity: not believed to be hygroscopic; may be handled in normal laboratory

atmosphere.

Light: protect from strong sunlight.

Caution: only experienced laboratory personnel should handle the material.

Version 1 (Id929) 1/5

# **Identity and Purity**

### **Proton NMR Spectrum**

Identity: the signals are consistent with the proposed structure and in accord with literature where available. Isotopic Labelling: signals at the sites of deuteration are greatly diminished, compared with the spectrum of unlabelled material, indicating clean deuteration.

Residual Solvents: no residual solvents are observed.

Impurities: no significant impurities are evident in the spectrum.

### **Carbon-13 NMR Spectrum**

Identity: the signals are consistent with the proposed structure and in accord with literature where available. Isotopic Labelling: signals at the sites of deuteration have collapsed to small multiplets compared with the spectrum of unlabelled material, indicating clean deuteration.

### **High-resolution Mass Spectrum (TOF MS ES+)**

Found m/z 371.1920.  $C_{22}H_{15}D_4N_6$  [M+H]<sup>+</sup> requires m/z 371.1922. The deviation of 0.5 ppm is within normally accepted limits for the establishment of identity by HRMS. No signal for  $d_0$  material was seen (detection limit about 0.5 %).

#### **HPLC**

A somewhat broadened, slightly tailing peak is observed (97.6 %). A small peak for Z-Rilpivirine is observed (13.16 min). Note: in the absence of reference materials for preparing calibration curves, it is assumed that all peaks have the same detector response. Where possible, the conditions of analysis follow a pharmacopeial or literature method, or have been adapted from same.

#### **Elemental Analysis**

Found: C 71.19, H 3.81, D 2.18, N 22.93 % Requires: C 71.33, H 3.81, D 2.17, N 22.69 %

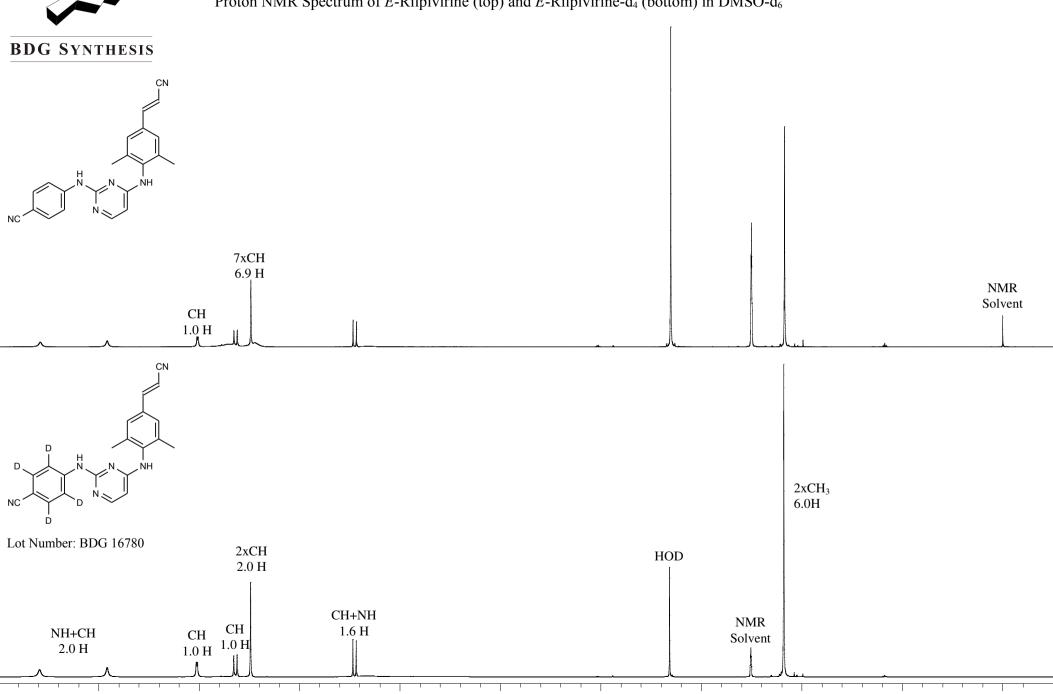
The elemental analyses fall within generally accepted limits for establishing the molecular formula given. The results may also be taken to imply the absence of significant quantities of water or inorganic salts (which have not been elsewhere tested for because of sample size limitations).

The available quantity of custom-synthesised material is always small, and this limits the extent and type of analytical data which can be obtained. This Certificate is presented in descriptive format for use by analytical chemists who are trained in the use of custom-synthesised materials. Custom materials often contain higher levels of residual solvents and/or water, and we urge you to use the corrected purity where needed rather than the raw HPLC purity. This compound is intended for use as an analytical reference material and it is not for human administration. Structures are shown with relative stereochemistry unless otherwise specified.

The re-test date is assigned from experience gained with the material in the laboratory and/or on storage. It is not possible to perform formal storage studies because of the small amount of material available.



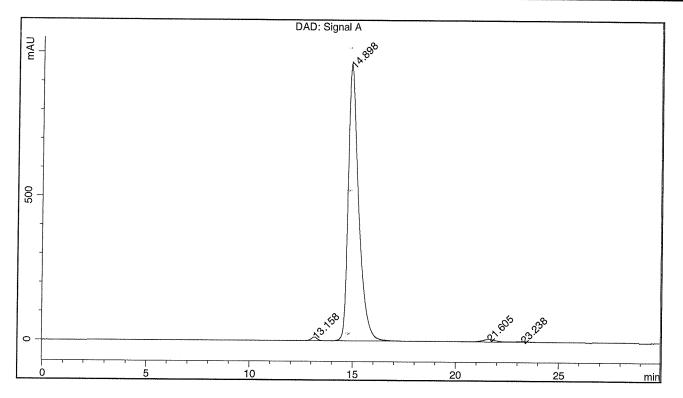
# Proton NMR Spectrum of E-Rilpivirine (top) and E-Rilpivirine-d<sub>4</sub> (bottom) in DMSO-d<sub>6</sub>



## BDG - Analysis of E-Rilpivirine

Column: Phenomenex Luna C18(2) 5um 250 x 4.6 mm
Guard: Phenomenex Security Guard C18 RP 4 x 3 mm
Mobile Phase: 50:50:0.05 Water: Methanol: Trifluoroacetic Acid
Column Temperature: 30 C . . . . Flow Rate: 1.0 mL/min . . . . Injection Volume: 10 uL
Sample Solvent: 1:1 Water: Methanol . . . . . Detection: UV 290 nm

Sample Name	BDG 16780	Instrument	AnalyticalLC01
Acquisition	09/09/2016, 16:04:08	Method (rev.)	LC10678c (10)
Sequence	BDG_09Sep2016c - Reprocessed	Vial Position	46
Operator	solvation010\cerityadmin	Injection	1 of 1



## **Area Percent Report**

Peak#	RT	Peak Height	Peak Area	Width	Area %
1	13.16 min	12.8454	292.4940	0.3457 min	0.845 %
2	14.90 min	956.9956	33801.0622	0.5161 min	97.622 %
3	21.60 min	8.7052	412.1879	0.7284 min	1.190 %
4	23.24 min	2.5499	118.5919	0.6626 min	0.343 %