BDG SYNTHESIS

Certificate of Analysis

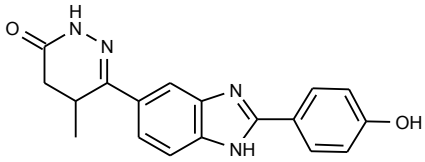
BDG Synthesis certifies that this reference material meets or exceeds the specifications stated herein.

Neil Beare, PhD, Director
25 January 2017

Name: ***O*-Desmethylpimobendan**

CAS Number: 108381-22-2

Structure:



Molecular Weight: $C_{18}H_{16}N_4O_2$ = 320.35

Lot Number: BDG 10621

Appearance: White, crystalline solid

Corrected Purity: 98.3 % (HPLC) - 5.5 % (water) = 92.8 %

Re-test Date: 25 January 2022

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.

• Analytical reference standards, metabolites, stable isotope labelled compounds • Custom synthesis • FTE contract research

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Identity and Purity

Proton NMR Spectrum

Identity: the signals are consistent with the proposed structure and in accord with literature where available.
Residual Solvents: no residual solvents are observed.
Impurities: traces of unidentified impurities are seen in the baseline.

Carbon-13 NMR Spectrum

Identity: the signals are consistent with the proposed structure and in accord with literature where available. Several peaks associated with the benzimidazole moiety have collapsed into the baseline.

High-resolution Mass Spectrum (ESI+)

Found *m/z* 321.1358. $C_{18}H_{17}N_4O_2$ $[M+H]^+$ requires *m/z* 321.1346. The deviation of 3.7 ppm is within normally accepted limits for the establishment of identity by HRMS.

HPLC

A sharp, symmetrical peak is observed (98.3 %). 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 63.85, H 5.36, N 16.45 %
$C_{18}H_{16}N_4O_2 \cdot 1.0H_2O$	Requires:	C 63.89, H 5.36, N 16.56 %
$C_{18}H_{16}N_4O_2$	Requires:	C 67.49, H 5.03, N 17.49 %

The elemental analyses fall somewhat outside those expected for anhydrous material; the presence of water is reasonably expected from the method of purification and/or the type of material, and the “best-fit” hydrated molecular formula is given.

Karl-Fischer Analysis

	Found:	H_2O 5.5 %
$C_{18}H_{16}N_4O_2 \cdot 1.0H_2O$	Requires:	H_2O 5.3 %

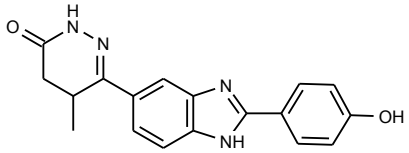
Of necessity, only a small sample could be used and only a single or duplicate analysis performed. We are unable to state what the errors in the reported water content are, but recommend that the result be used, as the best available, when determining corrected purity.

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.

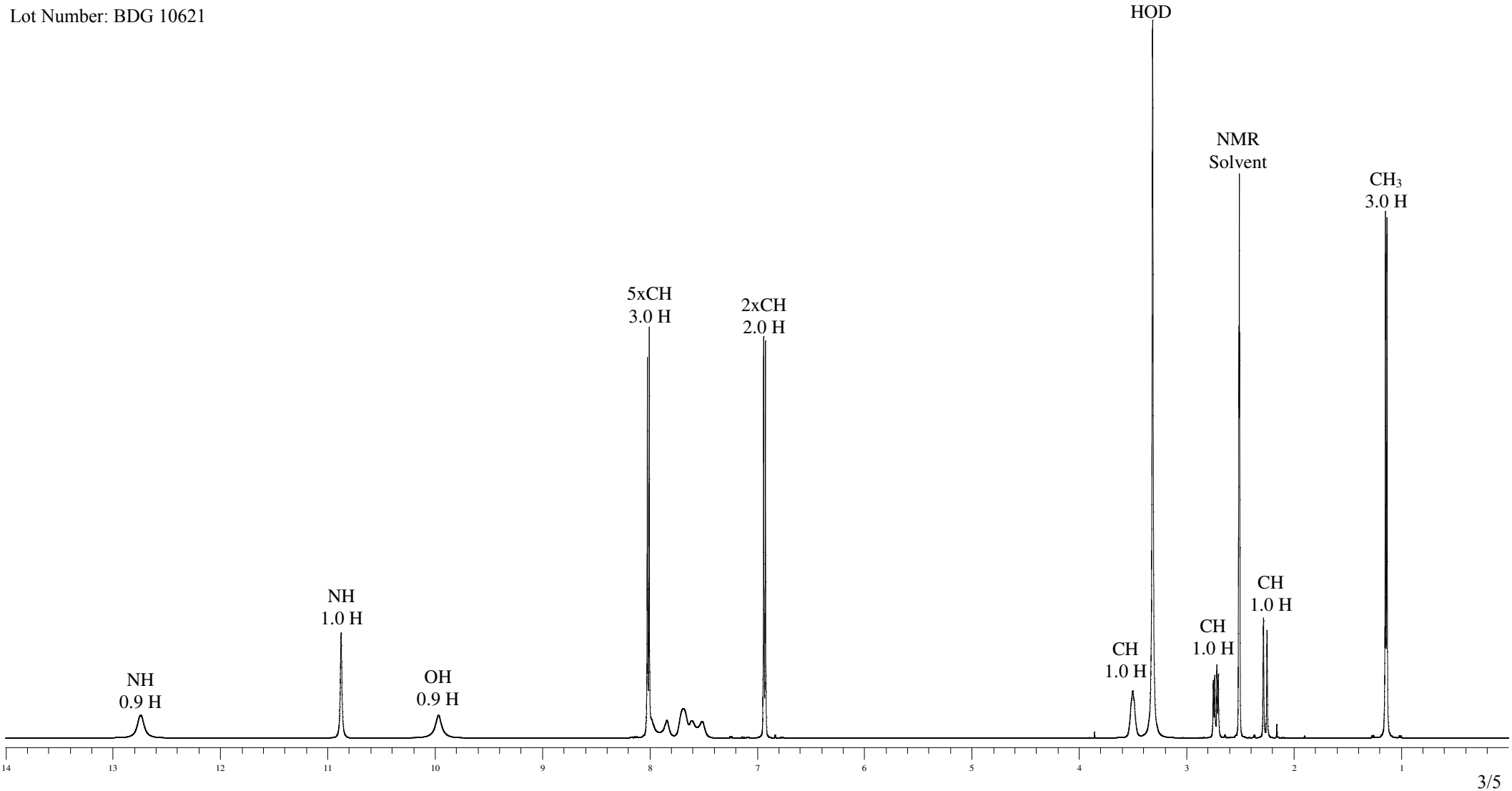
The re-test date is assigned from experience gained with the material in the laboratory and/or on storage. It is not possible to perform formal storage studies because of the small amount of material available.

BDG SYNTHESIS

Proton NMR Spectrum of *O*-Desmethylpimobendan in DMSO-d_6

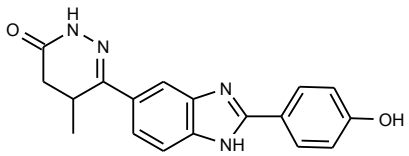


Lot Number: BDG 10621



BDG SYNTHESIS

Carbon-13 NMR Spectrum of *O*-Desmethylpimobendan in DMSO-d_6

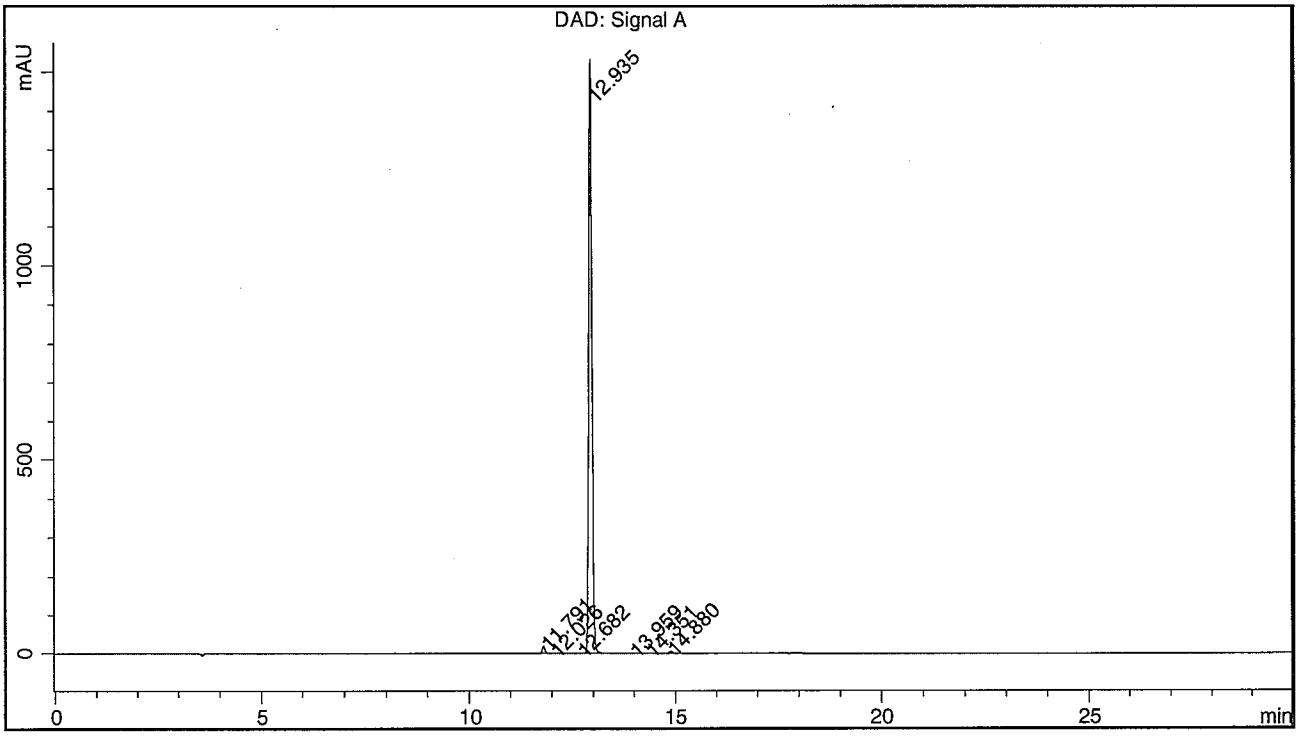


Lot Number: BDG 10621

BDG - Analysis of O-Desmethylpimobendan

Column : Phenomenex Luna C18(2) 5um 250 x 4.6 mm
Guard : Phenomenex Security Guard C18 RP 4 x 3 mm
Mobile Phase A : 90:10:0.1 Water : Acetonitrile : Trifluoroacetic Acid
Mobile Phase B : 50:50:0.1 Water : Acetonitrile : Trifluoroacetic Acid
Gradient (A:B) : T0=100:0, T24=0:100, T27=100:0, T30=100:0
Flow Rate : 1.0 mL/min . . . . . Sample Solvent : 1:1 Water : Acetonitrile
Column Temperature : 20 C . . . . . Injection Volume : 10 uL. . . . . Detection : UV 328 nm

Sample Name	BDG 10621	**Instrument**	AnalyticalLC01
Acquisition	25/01/2017, 13:50:39	**Method (rev.)**	LC10347b (12)
Sequence	BDG_25Jan2017b - Reprocessed	**Vial Position**	1
Operator	solvation010\cerityadmin	**Injection**	1 of 2



Area Percent Report

Peak#	RT	Peak Height	Peak Area	Width	Area %
1	11.79 min	18.1993	96.2687	0.0837 min	1.131 %
2	12.03 min	1.4203	8.3420	0.0885 min	0.098 %
3	12.68 min	0.3603	2.6494	0.1078 min	0.031 %
4	12.93 min	1535.2497	8361.0385	0.0855 min	98.259 %
5	13.96 min	0.5602	3.2469	0.0896 min	0.038 %
6	14.35 min	0.5249	4.9628	0.1343 min	0.058 %
7	14.88 min	5.0008	32.6401	0.1001 min	0.384 %