

## BDG SYNTHESIS

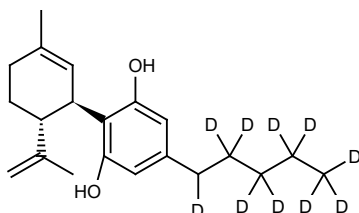
### Certificate of Analysis

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

*Neil Beare*

Neil Beare, PhD, Director  
1 November 2019

**Name:** Cannabidiol- $d_{10}$   
**CAS Number:** 13596-29-1 (unlabelled)  
**Structure:**



**Molecular Weight:**  $C_{21}H_{20}D_{10}O_2 = 324.52$   
**Lot Number:** BDG 11648  
**Appearance:** White, crystalline solid  
**Purity By HPLC:** 99.8 %  
**Isotopic Purity:** Under 0.5 %  $d_0$   
**Re-test Date:** 1 November 2024  
**Storage and Handling:** Temperature: Freeze (-20 °C) 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.

## 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: a trace (under 0.1 % w/w) of heptane is 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  $m/z$  325.2952.  $C_{21}H_{21}D_{10}O_2$   $[M+H]^+$  requires  $m/z$  325.2952. The deviation of 0.0 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 sharp, symmetrical peak is observed (99.8 %). 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 77.77, H 6.29, D 6.28 %
$C_{21}H_{20}D_{10}O_2$	Requires:	C 77.72, H 6.21, D 6.21 %

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).

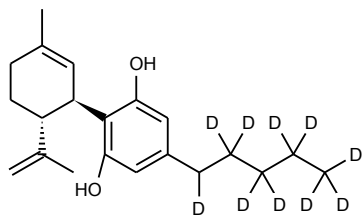
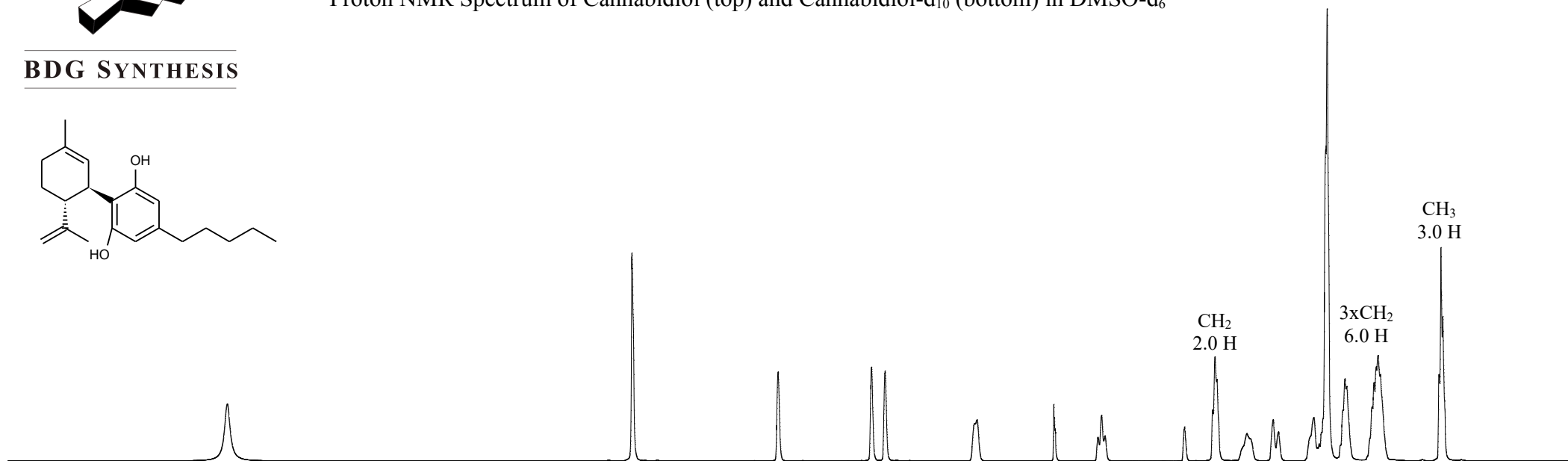
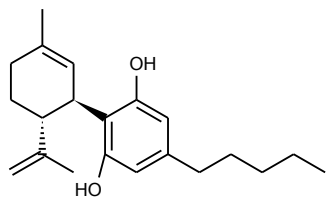
The available quantity of custom-synthesised material is always small, and this limits the extent and type of analytical data which can be obtained. This Certificate is presented in descriptive format for use by analytical chemists who are trained in the use of custom-synthesised materials. Custom materials 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.

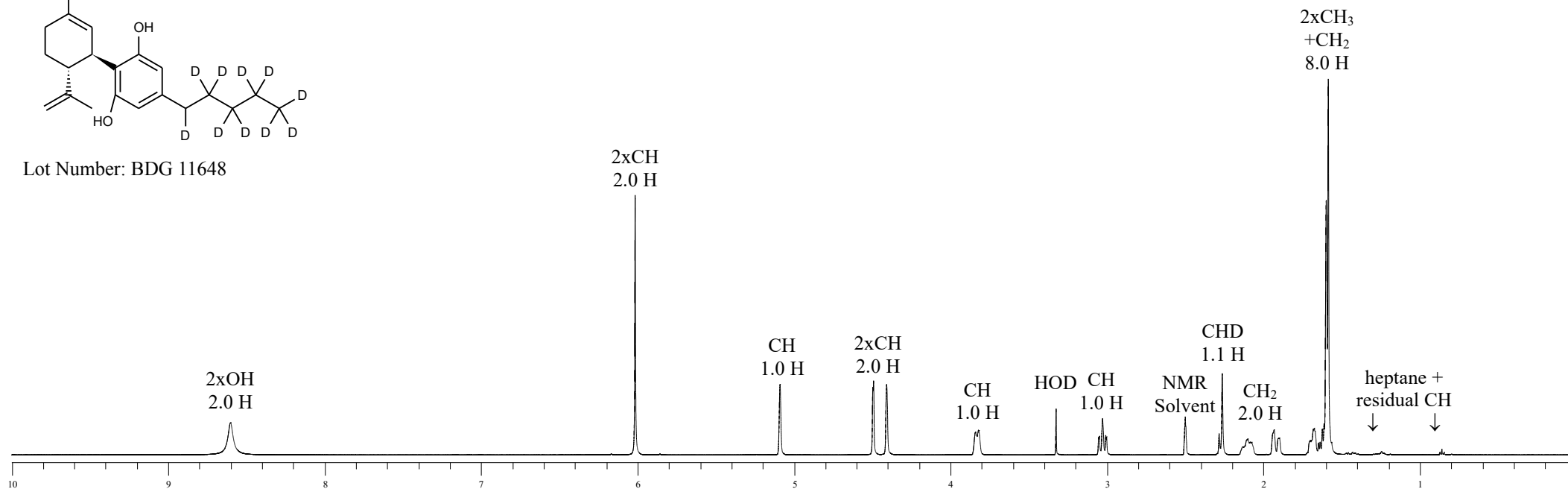


Proton NMR Spectrum of Cannabidiol (top) and Cannabidiol-d<sub>10</sub> (bottom) in DMSO-d<sub>6</sub>

**BDG SYNTHESIS**



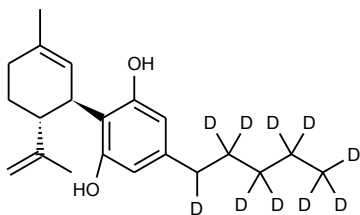
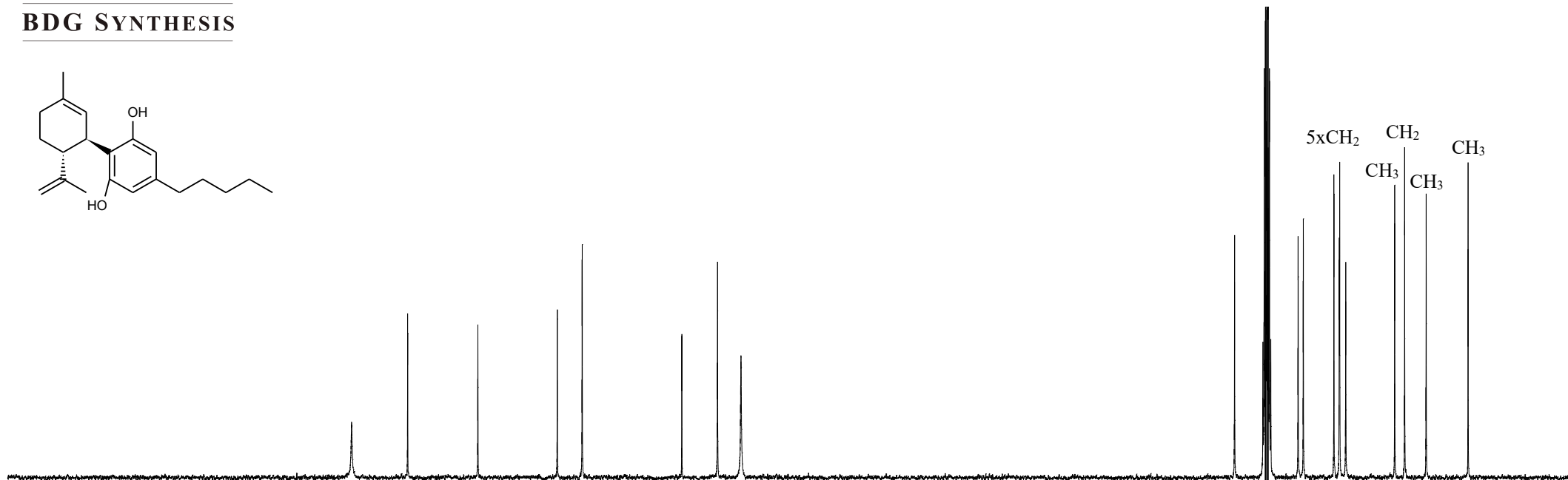
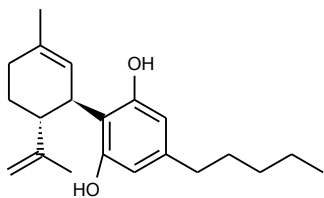
Lot Number: BDG 11648



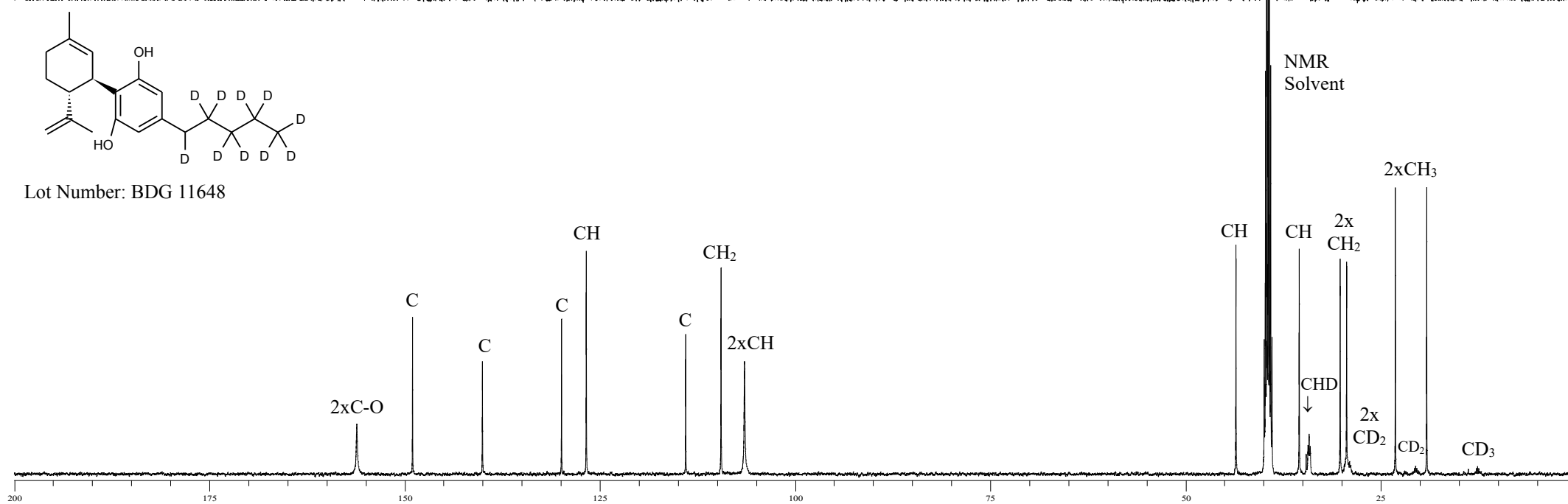


Carbon-13 NMR Spectrum of Cannabidiol (top) and Cannabidiol-d<sub>10</sub> (bottom) in DMSO-d<sub>6</sub>

**BDG SYNTHESIS**



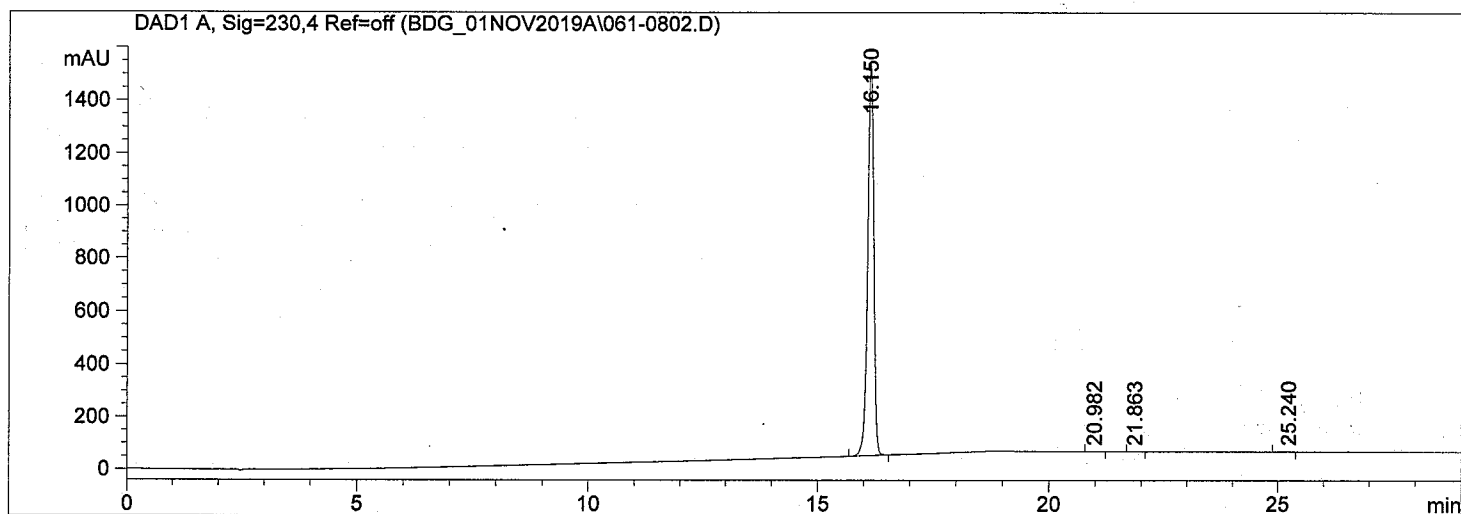
Lot Number: BDG 11648



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Acq. Operator   : Bruce Hamilton           Seq. Line :    8
Acq. Instrument : Instrument 1             Location  : Vial 61
Injection Date  : 11/1/2019 3:36:38 PM    Inj       :    2
                                           Inj Volume: 10 µl
Acq. Method     : C:\CHEM32\1\METHODS\2018\LC20102E.M
Last changed    : 11/1/2019 2:30:30 PM by Bruce Hamilton
Analysis Method : C:\CHEM32\1\METHODS\2018\LC20102E.M
Last changed    : 11/1/2019 5:49:00 PM by Bruce Hamilton
Method Info     : BDG - Analysis of Cannabidiol-d10
                  Column   : Phenomenex Luna C18(2) 5 µm 250 x 4.6 mm
                  Guard    : Phenomenex SecurityGuard C18 4 x 3 mm
                  Mobile Phase A : 30:70:0.05 Water : Methanol : Trifluoroacetic Acid
                  Mobile Phase C : 10:90:0.05 Water : Methanol : Trifluoroacetic Acid
                  Gradient (A:C) : T0=100:0, T15=0:100, T22=0:100, T23=100:0,
                  T27=100:0.
                  Sample Solvent : 30:70 Water : Methanol,   Detection : UV 230 nm,
                  Flow : 1 ml/min.,   Column Temperature : 35 C,   Injection : 10 µl.
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Area Percent Report  
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Sorted By           : Signal
Multiplier          : 1.0000
Dilution            : 1.0000
Use Multiplier & Dilution Factor with ISTDs

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Signal 1: DAD1 A, Sig=230,4 Ref=off

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	16.150	BB	0.1431	1.37101e4	1493.39783	99.8418
2	20.982	BB	0.1440	6.07138	5.90236e-1	0.0442
3	21.863	BB	0.1263	3.52594	3.59762e-1	0.0257
4	25.240	BB	0.1792	12.12168	8.56212e-1	0.0883

Totals :                                    1.37319e4  1495.20404

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\*\*\* End of Report \*\*\*