



BDG SYNTHESIS

Certificate of Analysis

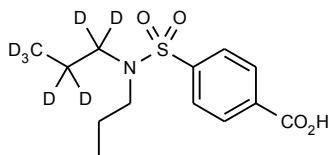
BDG Synthesis certifies that this reference material meets or exceeds the specifications stated herein.

Barry Dent

Barry R. Dent, PhD, Director
22 January 2011

Name: Probenecid-d₇
CAS Number: 57-66-9 (unlabelled)

Structure:



Molecular Weight: C₁₃H₁₂D₇NO₄S = 292.40

Lot Number: BDG 11165.2

Appearance: White powder

Purity By HPLC: 100.0 %

Isotopic Purity: Under 0.5 % d₀

Re-test Date: 22 January 2016

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. The material is susceptible to static.

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 absent, 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 (ESI+)

Found m/z 315.1361. $C_{13}H_{12}D_7NNaO_4S$ $[M+Na]^+$ requires m/z 315.1366. The deviation of 1.7 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 (100.0 %). 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 53.51, H 4.24, D 4.94, N 4.88 %
$C_{13}H_{12}D_7NO_4S$	Requires:	C 53.40, H 4.14, D 4.82, N 4.79 %

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).

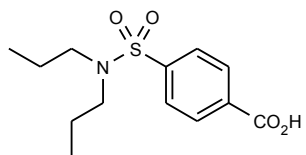
The available quantity of custom-synthesised material is always small, and this limits the extent and type of analytical data which can be obtained. This Certificate is presented in descriptive format for use by analytical chemists who are trained in the use of custom-synthesised materials. Custom materials 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 Probenecid (top) and Probenecid-d₇ (bottom) in DMSO-d₆

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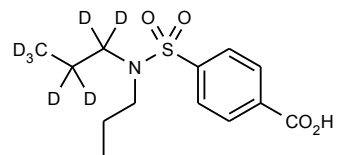
CO₂H

4H
4 x Ar-H

4H
2 x CH₂

4H
2 x CH₂

6H
2 x CH₃



Lot Number: BDG 11165.2

CO₂H

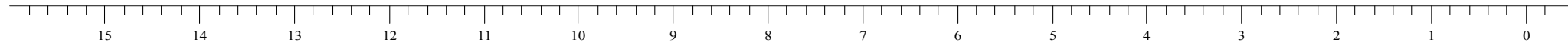
4H
4 x Ar-H

2H
CH₂

NMR
Solvent

2H
CH₂

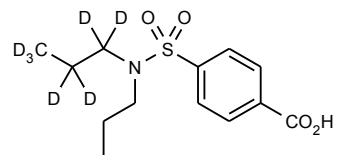
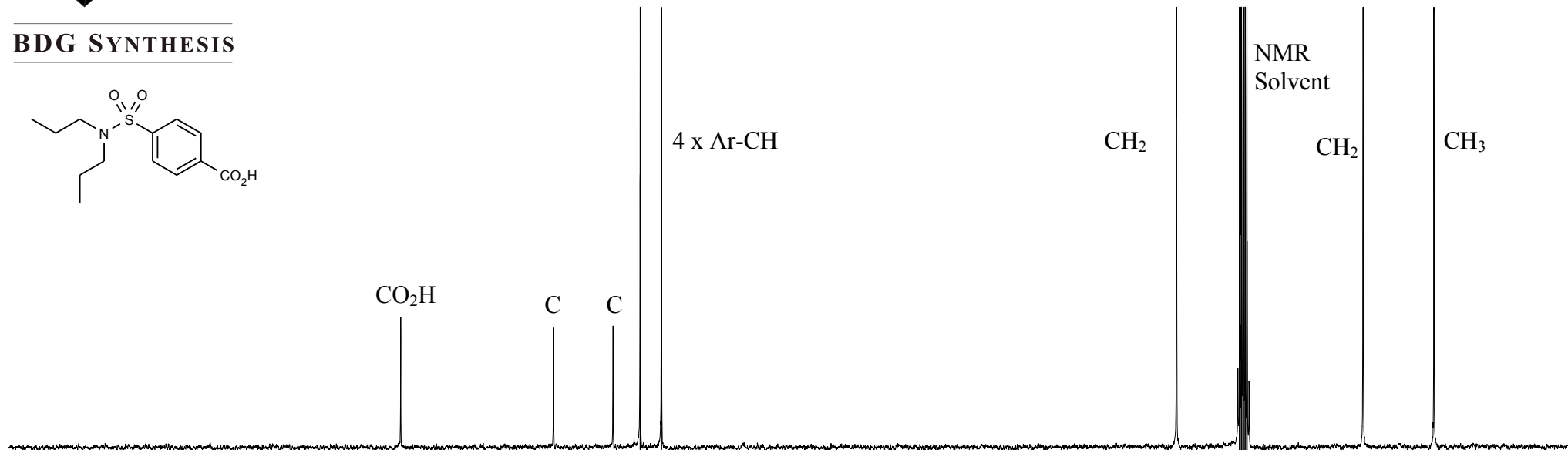
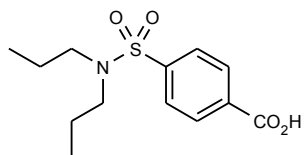
3H
CH₃



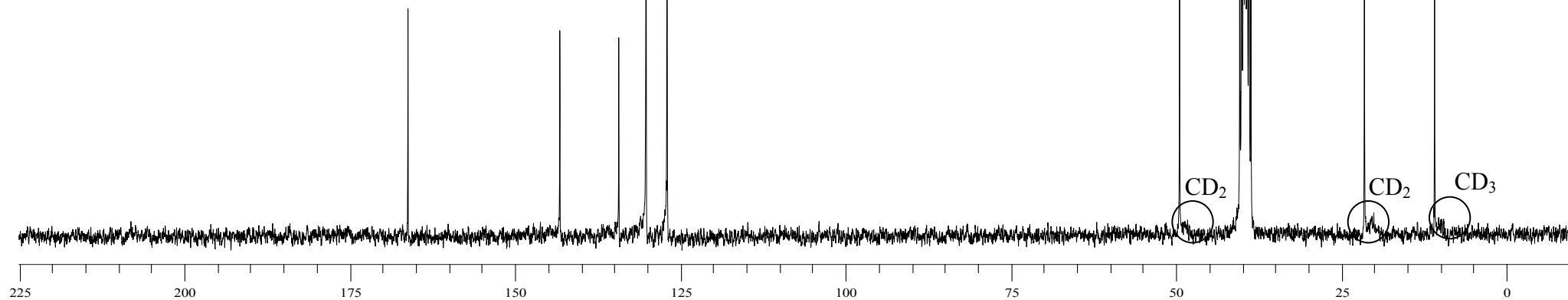


Carbon-13 NMR Spectrum of Probenecid (top) and Probenecid-d₇ (bottom) in DMSO-d₆

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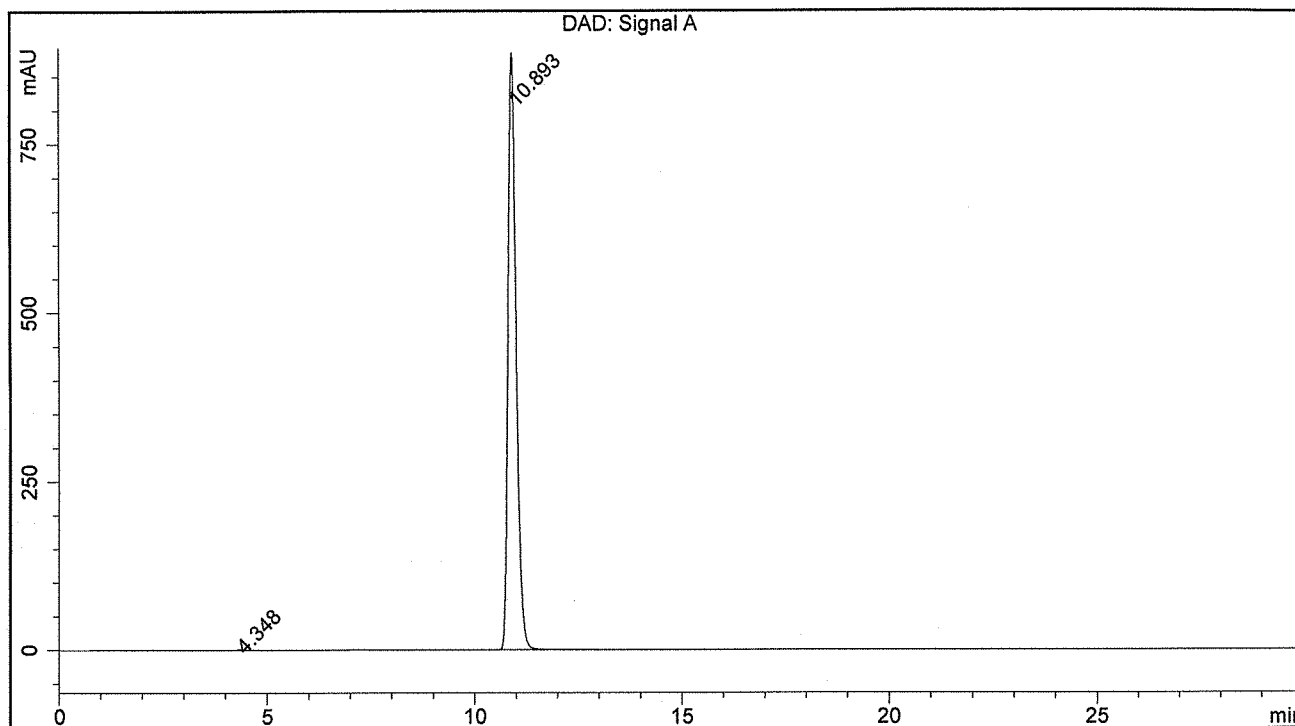
Lot Number: BDG 11165.2



BDG - Analysis of Probenecid-d7

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 50mM Potassium diHydrogen Phosphate pH=3.0 : Acetonitrile
 Flow Rate : 1.0 mL/min
 Sample Solvent : Initial Mobile Phase
 Injection Volume : 10 uL
 Column Temperature : 20C
 Detection : UV at 246 nm

Sample Name	BDG 11165.2	Instrument	AnalyticalLC01
Acquisition	22/01/2011, 16:05:01	Method (rev.)	LC10421a (5)
Sequence	BDG_22Jan2011g - Reprocessed	Vial Position	1
Operator	solvation010\cerityadmin	Injection	1 of 1



Area Percent Report

Peak#	RT	Peak Height	Peak Area	Width	Area %
1	4.35 min	0.8630	5.6840	0.1028 min	0.047 %
2	10.89 min	883.9055	12122.7627	0.2100 min	99.953 %