

Our Formula. Your Success.

Compound Data Sheet					
Name	Androst-4-ene-3,17-dione				
Synonym(s)	17-Ketotestosterone, 4-Androstene-3,17-dione				
Epichem Item #	EPL-DV2 Batch 1				
CAS#	63-05-8				
Molecular Formula	$C_{19}H_{26}O_2$				
Molecular Weight	286.42 g/mol				
Appearance	White solid				
Melting Point	170.8-172.9°C				
Purity*	98.1% at 242nm by HPLC (assuming all components detected with the same response factor).				
	No impurities detected by ¹ H NMR analysis.				
	LC/MS: $[M+H]^+ = 287.2$				
	IR vmax cm-1 1659(s), 1732(s)				
Date of Manufacture	3 April 2019				
Storage Requirements	Protect from heat, light and moisture.				
Special Precautions	This compound is for laboratory use only. Its toxicological properties may not				
	have been fully established. It should be handled only by suitably qualified				
	personnel.				
Date of Shipment	TBA				
	This certificate is valid for one year from the date of shipment provided the				
	substance is stored under the recommended conditions.				
Retest Date	TBA (Proper Storage and Handling Required)				

Product Reviewed By: Product Released By:

James Rixson, PhD
Head of Production

Carol Worth, PhD
Quality Manager

Release Date: 11 February 2022

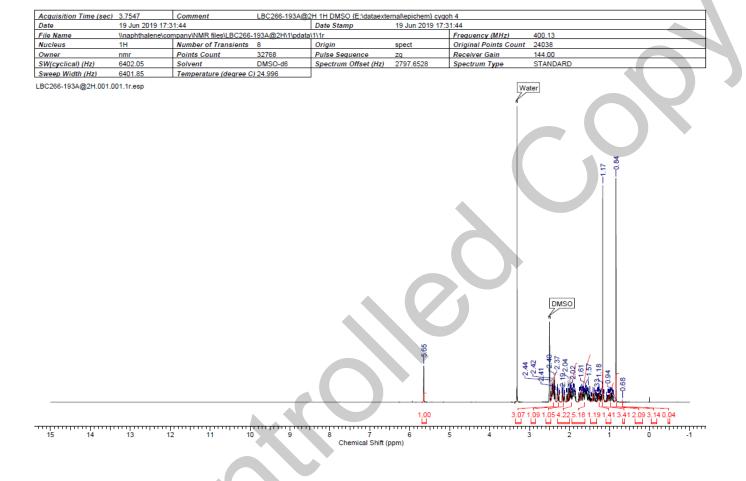
I. Identity

The identity of this product was established using the following analyses:

Ia. ¹HNMR Spectrum

Conditions: 400 MHz, DMSO-d₆

¹HNMR spectrum consistent with chemical structure.



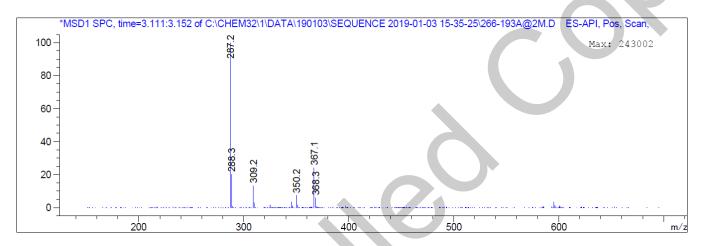
Ib. Mass Spectrum

The mass spectrum of this material was analysed by Liquid Chromatography Mass Spectroscopy (LCMS) using in-house EM005.WI08.

Method: ACN/water gradient (+ 0.1% formic acid).

ZORBAX SB-C8, 4.6 x 30 mm, 3.5 micron.

Retention		Mol. Weight
Time (MS)	MS Area	or Ion
3.129	2223576	367.10 I
		309.20 I
		288.25 I
		287.20 I

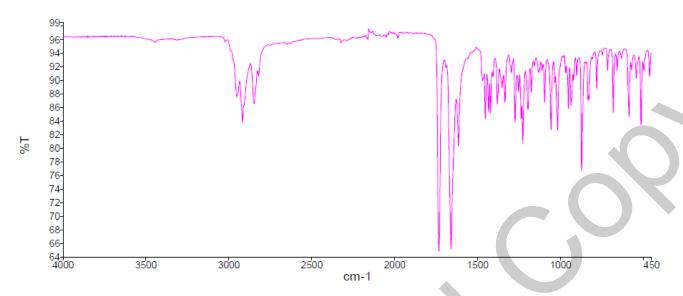


Theoretical value: 287.2 [M+H]⁺.

The signal of the Mass Spectrum is consistent with the theoretical value and its interpretation is consistent with the structural formula.

Ic. IR Spectrum

The infra-red spectrum of this material was analysed by Fourier-Transform Infrared Spectroscopy (FTIR) using in-house EM005.WI09.



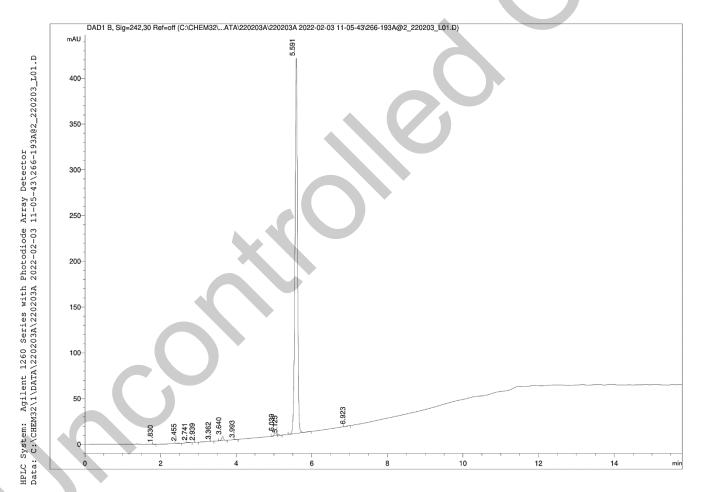
The interpretation of the signals of the Fourier-Transform Infrared Spectrum is consistent with the structural formula.

II. Purity

The purity of this material was analysed by high performance liquid chromatography (HPLC) using inhouse EM005.WI07.

HPLC Conditions:

Column	Conditions				Detector	Injector
Agilent Zorbax	25°C					Auto
SB-C18 4.6 x 250mm	Time (min)	% Line A (Water + 0.1% (v/v) TFA)	% Line B (Acetonitrile + 0.1% (v/v) TFA)	Flow rate (mL/min)	242nm	1.0 μL 1.2 mg/mL in
	0.00	50	50	1.5		100% acetonitrile
5 micron	9.00	5	95	1.5		(NO MODIFIERS)
	14.00	5	95	1.5		
	15.00	50	50	1.5		
	20.70	50	50	1.5		



Area Percent Report - Sorted by Signal

Peak Number	Retention Time (rounded)	Area	Area % (rounded)
1	1.83	0.05	0.00
2	2.45	1.70	0.09
3	2.74	1.86	0.10
4	2.94	0.13	0.01
5	3.36	0.13	0.01
6	3.64	15.74	0.82
7	3.99	0.15	0.01
8	5.04	13.77	0.72
9	5.12	2.73	0.14
10	5.59	1878.24	98.07
11	6.92	0.71	0.04
Totals			100 (rounded)

For the calculation the system peaks were ignored. The content of the analyte was determined as a ratio of the peak area of the analyte and the cumulative areas of the purities, added up to 100%.

Results:

Average 98.1% (average of duplicate runs)