

Australian Government

Department of Industry, Science and Resources

# National Measurement Institute



# DEUTERATED INTERNAL STANDARD PRODUCT INFORMATION SHEET

## NMIA D584c: d<sub>4</sub>-19-Norandrosterone

Report ID: D584c.2020.03 (Ampouled 140731)

Chemical Formula: C18H24D4O2

Molecular Weight: 280.4 g/mol

## **Property value**

Batch No.	CAS No.	Mass per ampoule
13-S-09	361432-47-5	1003 ± 20 μg

**IUPAC name:**  $(3\alpha,5\alpha)$ -3-Hydroxy $(2,2,4,4-^2H_4)$ estran-17-one.

**Expiration of certification:** The property values are valid till 17 April 2025, i.e. five years from the date of re-certification provided the **unopened** material is handled and stored in accordance with the recommendations below. The material as issued in the unopened container and stored as recommended below should be suitable for use beyond this date, subject to confirmation of batch stability from the issuing body. The shelf life does not apply to ampoules that have been opened. In such cases it is recommended that the end-user conduct their own in-house stability trials.

**Description:** The compound is supplied as a dried aliquot in a sealed ampoule under an atmosphere of argon. The deuterated internal standard is intended for a single use to prepare a standard solution containing D584c. The material was prepared by synthesis, and certified for identity and purity by NMIA.

**Intended use:** The isotopic purity of this material is an estimate only. This material should be considered for use as an internal standard only.

**Instructions for use:** Open the ampoule and carefully rinse the interior at least three times with a suitable organic solvent (e.g. TBME). This will transfer approximately 1003  $\mu$ g of anhydrous 19-norandrosterone (d<sub>4</sub>, d<sub>3</sub>, d<sub>2</sub>, d<sub>1</sub> and d<sub>0</sub>). The mass of analyte in each ampoule is calculated from the assigned purity of the bulk and the concentration of bulk material in a stock solution used to prepare the ampoules.

**Recommended storage:** When not in use, this material should be stored at or below 4 °C in a closed container in a dry, dark area.

**Stability:** This material has demonstrated stability over a minimum period of three years. The long-term stability of the compound in solution has not been examined.

**Homogeneity assessment:** The homogeneity of the material was assessed using purity assay by GC-FID on seven randomly selected ampoules of the material. The material was judged to be sufficiently homogeneous at this level of sampling as the variation in analysis results between samples was not significantly different at a 95% confidence level from that observed on repeat analysis of the same sample.

**Safety:** Treat as a hazardous substance. Use appropriate work practices when handling to avoid skin or eye contact, ingestion or inhalation of dust. Refer to the provided safety data sheet.

S.R. Davies

Dr Stephen R. Davies, Team Leader, Chemical Reference Materials, NMI. 1 November 2022.

This report supersedes any issued prior to 1 November 2022.

NATA Accreditation No. 198 / Corporate Site No. 14214.

Legal notice: Terms and Conditions associated with the provision of this reference material can be found on the NMIA website.

### **Characterisation Report:**

GC-FID:	Instrument:	Agilent 7890		
	Column:	HP-1MS, 30 m × 0	HP-1MS, 30 m × 0.32 mm I.D. × 0.25 μm	
	Program:	200 °C (13 min), 2	200 °C (13 min), 20 °C/min to 300 °C (3 min)	
	Injector:	250 °C	Detector Temp: 320 °C	
	Carrier:	Helium	Split ratio: 20/1	
	Relative peak area of the main component:			
	Initial analysis:	Mean = 99.3%, s	= 0.01% (7 ampoules in duplicate, August 2014)	
	Re- analysis:	Mean = 99.5%, s	= 0.02% (5 ampoules in duplicate, May 2017)	
	Re- analysis:	Mean = 99.3%, s	= 0.05% (5 ampoules in duplicate, April 2020)	

#### The following analytical data was obtained on the bulk material subsequently used in the preparation of the ampoules.

The identity was confirmed by a range of spectroscopic techniques, NMR, IR and MS. The purity value was obtained by mass balance from a combination of traditional analytical techniques, including GC-FID, thermogravimetric analysis, Karl Fischer analysis and <sup>1</sup>H NMR spectroscopy. The purity value is calculated as per Equation 1.

Purity = 
$$(100 \% - I_{ORG}) \times (100 \% - I_{VOL} - I_{NVR})$$

Equation 1

lorg = Organic impurities of related structure, IvoL = volatile impurities, INVR = non-volatile residue.

Supporting evidence is provided by qualitative elemental microanalysis.

The main component of this material is  $d_4$ -19-norandrosterone.  $d_3$ -,  $d_2$ -,  $d_1$ - and  $d_0$ -19-Norandrosterone are also present. The stated chemical purity of the analyte represents the combined mass fractions of deuterated ( $d_4$ ,  $d_3$ ,  $d_2$  and  $d_1$ ) and  $d_0$ -19-norandrosterone in the material.

The isotopic purity of this material is an estimate only. This material should be considered for use as an internal standard only.

Isotopic Purity:	$d_4 \approx 90\% \ [ = d_4/(d_4))$	$d_4 \approx 90\% \ [ = d_4/(d_4 + d_3 + d_2 + d_1 + d_0) \times 100]$		
	$d_0 < 0.2\%$ [ = $d_0/(d_0)$	4 + d <sub>3</sub> +d <sub>2</sub> +d <sub>1</sub> +d <sub>0</sub> ) x 100]		
GC-FID:	Instrument: Column: Program: Injector: Carrier:	Agilent 6890 HP-1MS, 30 m × 0.32 mm 200 °C (13 min), 20 °C/mi 250 °C Helium	•	
	Relative peak area of I Initial analysis:	•	(10 sub samples in duplicate, October 2013)	
Thermogravimetric analysis:		Non volatile residue < 0.2% mass fraction (October 2013). The volatile content (e.g. organic solvents and/or water) could not be determined because of the inherent volatility of the material and/or degradation at elevated temperatures.		
Kaal <b>F</b> ischen andersier		Majatura content : 0.40/ m	and fraction (Ostahar 2012)	

Karl Fischer analysis: Moisture content < 0.1% mass fraction (October 2013)

## Spectroscopic and other characterisation data

GC-MS:	Parent compound:			
	Instrument:	Agilent 6890/5973		
	Column:	TG1-MS, 30 m x 0.25 mm l.D. x 0.25 μm		
	Program:	180 °C (1 min), 10 °C/min to 300 °C (3 min) 250 °C Transfer line temp: 280 °C		
	Injector: Carrier:	Helium, 1.0 mL/min Split ratio: 20/1		
	Bis-TMS derivative:			
	Instrument: Column:	Agilent 6890/5973		
	Program:	TG1-MS, 30 m x 0.25 mm l.D. x 0.25 μm 180 °C (1 min), 10 °C/min to 300 °C (3 min)		
	Injector:	250 °C Transfer line temp: 280 °C		
	Carrier:	Helium, 1.0 mL/min Split ratio: 20/1		
		•		
	The retention times of the parent compound and <i>bis</i> -TMS derivative are reported along with the major peaks in the mass spectra. The latter are reported as mass/charge ratios and (in brackets) as a percentage relative to the base peak.			
	Parent (9.2 min):	280 (M <sup>+</sup> , 100), 262 (22), 236 (36), 218 (19), 206 (51), 191 (29), 163 (20), 151 (27), 135 (18), 124 (20), 108 (21), 97 (26), 92 (31), 79 (31), 67 (30) <i>m/z</i>		
	<i>Bis</i> -TMS (9.9 min):	424 (M <sup>+</sup> , 64), 409 (100), 337 (12), 319 (24), 229 (14), 169 (33), 73 (100) <i>m</i> /z		
TLC:	Conditions:	Kieselgel 60F <sub>254</sub> . Chloroform/ethyl acetate (4/1) Single spot observed, $R_f = 0.6$ Visualisation with vanillin		
IR:	Instrument: Range: Peaks:	FT-IR, Biorad WIN FTS40 4000-400 cm <sup>-1</sup> , KBr pellet 3479, 1723, 1443, 1091 cm <sup>-1</sup>		
<sup>1</sup> H NMR:	Instrument: Field strength: Solvent: Spectral data:	Bruker Avance III-400 400 MHz CDCI₃ (7.26 ppm) δ 0.68-0.82 (2H, m), 0.86 (3H, s), 0.96-1.44 (9H, m), 1.49 (1H, m), 1.58-1.62 (1H, m), 1.69 (1H, dd, <i>J</i> = 3.1, 13.4 Hz), 1.74-1.80 (2H, m), 1.84-1.95 (2H, m), 2.06 (1H, dt, <i>J</i>		
		=19.2, 9.0 Hz), 2.43 (1H, ddd, <i>J</i> = 0.7, 8.7, 19.1 Hz), 4.07 (1H, d, <i>J</i> = 2.4 Hz) ppm		
		Ethyl acetate estimated at 0.6% mass fraction was observed in the <sup>1</sup> H NMR		
<sup>13</sup> C NMR:	Instrument: Field strength: Solvent: Spectral data:	Bruker DMX-500 126 MHz CDCl <sub>3</sub> (77.19) $\delta$ 14.0, 21.8, 23.7, 25.1, 30.0, 31.8, 32.4 (m), 33.6, 36.0, 39.9 (m), 41.0, 47.1, 48.1, 48.5, 50.9, 66.3, 221.6 ppm.		
Melting point:		164-166 °C		
Microanalysis:	Found: Calculated:	C = 76.6%; H+D = 10.3% (November, 2013) C = 77.1%; H+D = 10.0% (Calculated for C <sub>18</sub> H <sub>24</sub> D <sub>4</sub> O <sub>2</sub> )		