



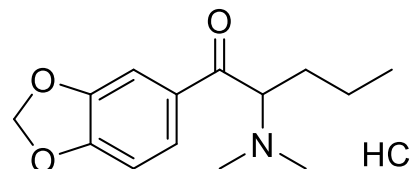
CERTIFIED REFERENCE MATERIAL CERTIFICATE OF ANALYSIS

NMIA D1094: N,N'-Dimethylpentylone hydrochloride

Report ID: D1094.2026.01

Chemical Formula: $C_{14}H_{19}NO_3 \cdot HCl$

Molecular Weight: 285.8 $gmol^{-1}$ (HCl), 249.3 $gmol^{-1}$ (base)



Certified value

Batch No.	CAS No.	Purity (mass fraction)
26-D-01	17763-13-2 (HCl) 803614-36-0 (base)	99.0 ± 1.3%

The uncertainty has been calculated according to ISO 33405 and is stated at the 95 % confidence limit ($k = 2$).

IUPAC name: 1-(benzo[d][1,3]dioxol-5-yl)-2-(dimethylamino)pentan-1-one hydrochloride.

Expiration of certification: The property values are valid till 12 May 2029, three years from the date of 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 expiry date/shelf life does not apply to sample bottles that have been opened. In such cases it is recommended that the end-user conduct their own in-house stability trials. The material will be re-tested periodically to ensure that the property values are still valid. In the event a product fails the stability trial, notification will be sent to impacted customers.

Description: White powder prepared by synthesis and certified for identity and purity by NMI Australia. Packaged in amber glass bottles with a septum and crimped aluminium cap.

Intended use: This certified reference material is suitable for use as a primary calibrator.

Instructions for use: Equilibrate the bottled material to room temperature before opening.

Recommended storage: When not in use this material should be stored at or below 25 °C in a closed container in a dry, dark area.

Metrological traceability: The certified purity value is traceable to the SI unit for mass (kg) through Australian national standards via balance calibration. In the mass balance all impurities are quantified as a mass fraction and subtracted from 100 %.

Stability: In the absence of long-term stability data, the measurement uncertainty at the 95 % coverage interval has been expanded to accommodate any potential change in the property value. The stability component has been estimated from stability trials conducted on similar materials by NMI Australia over the last ten years. The measurement uncertainty at the 95 % confidence interval also includes a stability component determined from accelerated stability trials conducted at 40 °C and 75 % humidity for 7 days.

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 ten randomly selected 1–2 mg sub samples 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.
30 July 2026

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

CIPM MRA notice: This certificate is consistent with the capabilities that are included in Appendix C of the MRA drawn up by the CIPM. Under the CIPM MRA, all participating institutes recognise the validity of each other's calibration and measurement certificates for the quantities, ranges and measurement uncertainties specified in the KCDB (for details see <http://www.bipm.org/kcdb/>). The "CIPM MRA Logo" and this statement attest only to the measurement(s) applied for determining the certified values on the certificate.

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

Characterisation Report:

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

$$\text{Purity} = (100 \% - I_{\text{ORG}}) \times (100 \% - I_{\text{VOL}} - I_{\text{NVR}}) \quad \text{Equation 1}$$

I_{ORG} = Organic impurities of related structure, I_{VOL} = volatile impurities, I_{NVR} = non-volatile residue.

Supporting evidence is provided by elemental microanalysis.

GC-FID:	Instrument:	Agilent 8890
	Column:	HP-1MS (30 m × 0.32 mm I.D. × 0.25 μm)
	Program:	60 °C (2 min), 15 °C/min to 240 °C (14.5 min), and 20 °C/min to 280 °C (14.5 min).
	Injector:	180 °C
	Detector Temp:	320 °C
	Carrier:	Helium (2 mL/min)
	Split ratio:	20/1
	Relative mass fraction of the main component as the free base:	
	Initial analysis:	Mean = 99.2 %, s = 0.04 % (10 sub samples in duplicate, May 2026)
Karl Fischer analysis:	Moisture content < 0.2 % mass fraction (May 2026)	
Thermogravimetric analysis:	Non-volatile residue < 0.2 % mass fraction (May 2026)	

Spectroscopic and other characterisation data

GC-MS:	Instrument:	Agilent 6890/5973
	Column:	HP-5MS, 30 m × 0.25 mm I.D. × 0.25 µm
	Program:	90 °C (1 min), 15 °C/min to 240 °C (5 min), 25 °C/min to 300 °C (4 min)
	Injector:	250 °C
	Split ratio:	20/1
	Transfer line temp:	280 °C
	Carrier:	Helium, 1.0 mL/min
	Scan range:	50–300 <i>m/z</i>
	The retention time of <i>N,N'</i> -dimethylpentylone free base is reported 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.	
	Free base (11.0 min):	149 (6), 101 (7), 100 (100), 71 (8), 65 (5), 58 (15), 56 (5), 44 (7), and 42 (8) <i>m/z</i>
ESI-MS:	Instrument:	Shimadzu LC-TQ-MS 8045
	Operation:	Direct infusion at 10 µL/min
	Ionisation mode:	Electrospray positive ion mode
	Interface voltage:	4.0 kV
	Peak:	250 [M+H] ⁺ , 251 [M+2H] ⁺ <i>m/z</i>
IR:	Instrument:	Bruker ALPHA II PLATINUM-ATR
	Range:	4000–400 cm ⁻¹ , neat
	Peaks:	2961, 2637, 1674, 1613, 1504, 1439, 1355, 1259, 1095, 1034, 932, 814, 784 cm ⁻¹
¹ H-NMR:	Instrument:	Bruker Avance III-500
	Field strength:	500 MHz
	Solvent:	D ₂ O (4.79 ppm)
	Spectral data:	δ 0.84 (3H, t, <i>J</i> = 7.3 Hz), 1.19 (2H, m), 2.05 (2H, m), 2.90 (3H, s), 3.02 (3H, s), 5.12 (1H, t, <i>J</i> = 5.5 Hz), 6.14 (2H, d, <i>J</i> = 2.6 Hz), 7.03 (1H, d, <i>J</i> = 8.3 Hz), 7.48 (1H, d, <i>J</i> = 1.8 Hz), 7.72 (1H, dd, <i>J</i> = 1.8, 8.3 Hz) ppm
	Acetone and 2-propanol were quantified at 0.01% and 0.02% mass fractions, respectively. Similarly, ¹ H NMR confirmed the presence of pentylone at ~0.4% mass fraction.	
¹³ C-NMR:	Instrument:	Bruker Avance III-500
	Field strength:	126 MHz
	Solvent:	D ₂ O
	Spectral data:	δ 12.9, 16.8, 30.8, 40.3, 43.6, 69.3, 102.7, 107.8, 108.6, 126.8, 128.3, 148.5, 153.9, 195.0 ppm
Melting point (DSC):		235–236 °C
Melting point (hot stage):		>222 °C (Decomp.)
Microanalysis:	Found:	C = 59.1%; H = 6.9%; N = 4.9%; (May 2026)
	Calculated:	C = 58.8%; H = 7.1%; N = 4.9%; (Calculated for C ₁₄ H ₁₉ NO ₃ ·HCl)