National Measurement Institute

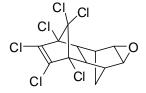


CERTIFIED REFERENCE MATERIAL CERTIFICATE OF ANALYSIS

NMIA P1747: Dieldrin

Report ID: P1747.2023.01

Chemical Formula: C₁₂H₈Cl₆O Molecular Weight: 380.9 g/mol



Certified value

Batch No.	CAS No.	Purity (mass fraction)
02-AV-01	60-57-1	99.9 ± 0.3%

The uncertainty has been calculated according to ISO Guide 35 and is stated at the 95% confidence limit (k = 2).

IUPAC name: (1R,2S,3S,7R,8S,9S,11R)-3,4,5,6,13,13-Hexachloro-10-oxapentacyclo[6.3.1.13,6.02,7.09,11]tridec-4-ene.

Expiration of certification: The property values are valid till 10 May 2033, ten 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 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.

Description: White crystalline powder sourced from an external supplier and certified for identity and purity by NMIA. Packaged in amber glass bottles with a septum and crimped aluminium cap or screw top 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.

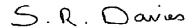
Recommended storage: When not in use this material should be stored at or below 4 °C in a closed container in a dry, dark area.

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: This material has demonstrated stability over a minimum period of ten years. The measurement uncertainty at the 95% confidence interval includes a stability component which has been estimated from annual stability trials. 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 2-2.5 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.



Dr Stephen R. Davies, Team Leader, Chemical Reference Materials, NMI. 6 June 2023

This report supersedes any issued prior to 06 June 2023

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:

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 and Karl Fischer. The purity value is calculated as per Equation 1.

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

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: HP5890

Column: BPX5 Capillary, 30 m \times 0.32 mm I.D. \times 0.25 μ m

Program: 180 °C (1 min), 15 °C/min to 300 °C

Injector: 250 °C

Detector Temp: 340 °C

Carrier: Helium

Split ratio: 30/1

Relative mass fraction of the main component:

Initial analysis: Mean = 99.9%, s = 0.01% (10 sub samples in duplicate, February 2002)

Re-analysis: Mean = 99.95%, s = 0.001% (5 sub samples in duplicate, February 2008)GC-FID:

Instrument: Agilent 6890N

Column: HP-1, 30 m x 0.32 mm I.D. x 0.25 μ m Program: 180 °C (1 min), 10 °C/min to 300 °C (3 min)

180 °C (1 min), 10 °C/min to 300 °C (5 min)

Injector: 230 °C

Detector Temp: 320 °C

Carrier: Helium

Split ratio: 20/1

Relative mass fraction of the main component:

Initial analysis: Mean = 99.94%, s = 0.003% (5 sub samples in duplicate, January 2014) Re-analysis: Mean = 99.95%, s = 0.001% (5 sub samples in duplicate, December 2018) Re-analysis: Mean = 99.95%, s = 0.001% (5 sub samples in duplicate, May 2023)

Karl Fischer analysis: Moisture content ≤ 0.2% mass fraction (February 2008 - May 2023)

Thermogravimetric analysis: Due to the volatility of the material, unable to estimate volatiles content

Non-volatile content < 0.2% mass fraction (August 2002)

Spectroscopic and other characterisation data

GC-MS: Instrument: HP5890/5971A

Column: ZB-5, 12 m x 0.25 mm l.D. x 0.20 μ m Program: 130 °C (1 min), 15 °C /min to 300 °C (3 min)

 $\begin{array}{lll} \mbox{Injector:} & 250 \ ^{\circ}\mbox{C} \\ \mbox{Split ratio:} & 40/1 \\ \mbox{Transfer line temp:} & 280 \ ^{\circ}\mbox{C} \\ \mbox{Carrier:} & \mbox{Helium} \\ \mbox{Scan range:} & 40-550 \ \mbox{\it m/z} \\ \end{array}$

The retention time of the parent compound 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.

Parent (8.7 min): 380 (M⁺, 11), 345 (11), 277 (32), 263 (39), 237 (19), 108 (19), 79 (100) m/z

The fragmentation pattern matches published reference spectra for dieldrin.

TLC: Conditions: Kieselgel 60F₂₅₄. Hexane/chloroform 50/50

Single spot observed, $R_f = 0.4$.

IR: Instrument: FT-IR, Biorad Merlin

Range: 4000-400 cm⁻¹, KBr pellet

Peaks: 1598, 1248, 1176, 1076, 1040, 1005, 845, 800, 702, 431 cm⁻¹

Microanalysis: Found: C = 38.0%, H = 2.0%, CI = 55.8% (February 2002)

Found: C = 38.0%, H = 2.1%, CI = not determined (November 2005) Calculated: C = 37.8%, H = 2.1%, CI = 55.8% (Calculated for $C_{12}H_8CI_6O$)