Australian Government
Department of Industry, Science and Resources

National Measurement Institute

NATA
WORLD RECOGNISED ACCREDITATION

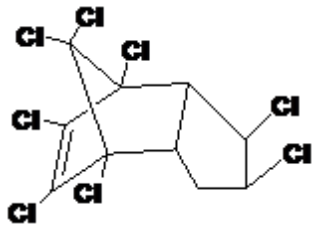
CERTIFIED REFERENCE MATERIAL CERTIFICATE OF ANALYSIS

NMIA P1624: cis-Chlordane

Report ID: P1624.2021.03

Chemical Formula: $C_{10}H_6Cl_8$

Molecular Weight: 409.8 g/mol



Certified value

Batch No.	CAS No.	Purity (mass fraction)
97-000663	5103-71-9	99.6 ± 0.3%

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

IUPAC name: (1S,2S,3R,4S,6S,7R)-1,3,4,7,8,9,10,10-Octachlorotricyclo[5.2.1.$0^{2,6}$]dec-8-ene.

Expiration of certification: The property values are valid till 22 January 2031, i.e. 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. The material will be re-tested on an annual basis to ensure that the property values are still valid. In the event a product fails the stability trial, notification will be sent to all impacted customers.

Description: White crystals 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 ºC 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 approach 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 seven 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.
6 September 2022

This report supersedes any issued prior to 6 September 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:

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 ^{1}H NMR spectroscopy. The purity value is calculated as per Equation 1.

Purity = (100 % - I_{ORG}) x (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:	Hewlett Packard 5890		
	Column:	BP-5 Capillary, 30 m x 0.32 mm I.D. x 0.25 μm		
	Program:	130 °C (1 min), 10 °C/min to 280 °C (3 min)		
	Injector:	250 °C	Detector Temp:	315°C
	Carrier:	Helium	Split ratio:	10/1
	Relative mass fraction of the main component:			
	Initial analysis:	Mean = 99.8%, s = 0.1% (10 subsamples in duplicate, May 1999)		
	Re-analysis:	Mean = 99.8%, s = 0.04% (5 subsamples in duplicate, May 2006)		
GC-FID:	Instrument:	Agilent 7890		
	Column:	HP-1MS, 30 m × 0.32 mm I.D. × 0.25 μm		
	Program:	130 °C (1 min), 10 °C/min to 300 °C (3 min)		
	Injector:	250 °C	Detector Temp:	320 °C
	Carrier:	Helium	Split ratio:	20/1
	Relative mass fraction of the main component:			
	Re-analysis:	Mean = 99.8%, s = 0.04% (5 subsamples in duplicate, May 2011)		
	Re-analysis:	Mean = 99.8%, s = 0.04% (5 subsamples in duplicate, March 2016)		
	Re-analysis:	Mean = 99.8%, s = 0.03% (5 subsamples in duplicate, January 2021)		

Karl Fischer analysis:	Moisture content ≤ 0.1% mass fraction (July 2006) Moisture content 0.22% mass fraction (May 2011) Moisture content 0.13% mass fraction (March 2016) Moisture content ≤ 0.1% mass fraction (January 2021)
Thermogravimetric analysis:	Non-volatile residue < 0.2% mass fraction (November 1999)

Spectroscopic and other characterisation data

GC-MS:	Instrument:	Varian Saturn 3400/2000 GC-MS Ion Trap
	Column:	J&W DB-17MS, 30 m x 0.25 mm I.D. x 0.17 μm
	Program:	130 °C (1 min), 15 °C /min to 300 °C (3 min)
	Injector:	250 °C
	Split ratio:	10/1
	Transfer line temp:	280 °C
	Carrier:	Helium
	Scan range:	50-550 *m/z*
	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 (8.0 min):	408 (M^+), 373 (100), 337, 301, 266, 237, 196, 169, 109, 65 *m/z*
		Matches published literature spectra for *cis*-chlordane
IR:	Instrument:	FT-IR, Biorad WIN FTS40
	Range:	4000-400 cm^{-1}, KBr pellet
	Peaks:	1602, 1269, 1166, 1020, 977, 901, 826, 751, 723, 699 cm^{-1}
		Matches published literature spectra for *cis*-chlordane
^{1}H NMR:	Instrument:	Bruker DMX-300
	Field strength:	300 MHz
	Solvent:	$CDCl_3$ (7.26 ppm)
	Spectral data:	δ 1.83 (1H, m), 2.40 (1H, m), 3.56 (1H, t), 3.72 (1H, dd), 3.97 (1H, m), 4.43 (1H, m) ppm
		The observed spectrum is consistent with a reference spectrum for *cis*-chlordane obtained at lower field strength.
^{13}C NMR:	Instrument:	Bruker Avance III-500
	Field strength:	126 MHz
	Solvent:	$CDCl_3$ (77.16 ppm)
	Spectral data:	δ 34.1, 52.0, 59.3, 59.6, 66.3, 80.2, 80.9, 105.0, 132.0, 132.6 ppm
Melting point:		104-106 °C
Microanalysis:	Found:	C = 29.4%, H = 1.3%, Cl = 69.0%
	Calculated:	C = 29.3%, H = 1.5%, Cl = 69.2% (Calculated for $C_{10}H_6Cl_8$)