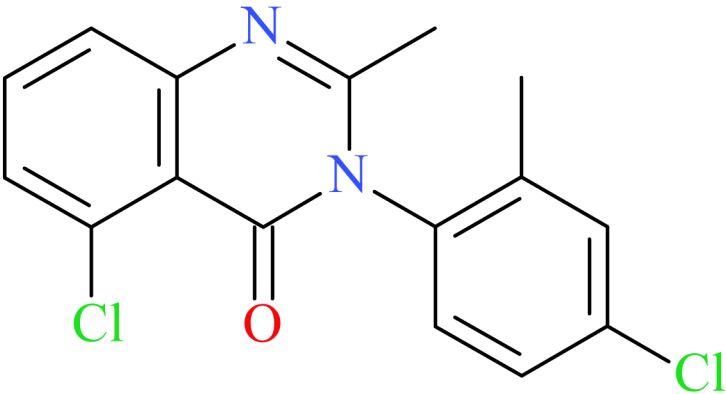




Dicloqualone



NPS SUBCLASS
Miscellaneous
REPORT DATE
June 17, 2025
SAMPLE RECEIVED
November 20, 2024
SAMPLE TYPE
Drug Material

Preferred Name	Dicloqualone
Synonyms	SL-164
Formal Name	5-chloro-3-(4-chloro-2-methyl-phenyl)-2-methyl-quinazolin-4-one
InChI Key	KUIHLOHNUGOCTO-UHFFFAOYSA-N
CAS Number	3476-88-8
Chemical Formula	C ₁₆ H ₁₂ Cl ₂ N ₂ O
Molecular Weight	319.2
Molecular Ion [M ⁺]	318
Exact Mass [M+H] ⁺	319.0399

Characterization & Intelligence

The following information was compiled in June 2025 and is subject to change as new research is conducted and as new information becomes available:

Description: Dicloqualone is categorized as a quinazolinone and bears structural similarity to other “quaaludes” such as methaqualone. Dicloqualone was first identified in California in early 2022 and has subsequently been identified across the United States.^{1,2,3} Dicloqualone was identified by our laboratory in November 2024 and confirmed after acquiring standard reference material.

Sample Source: Chicago Recovery Alliance (Chicago, IL)

Sample Appearance: White powder

Pharmacology: Dicloqualone has sedative properties and increases the activity of GABA receptors similar to that of methaqualone.^{4,5}

Toxicology: Dicloqualone has not been detected in toxicology cases to date at the CFSRE.

Drug Materials: Dicloqualone has been detected in one drug material to date at the CFSRE.

Demographics / Geographics: The drug material originated from Chicago, IL, and dicloqualone was found alongside opioids (e.g., fentanyl, heroin) and diphenhydramine.

Legal Status: Dicloqualone is not currently scheduled in the United States.

References:

- ▶ Cayman Chemical: [SL-164](#)
- ▶ ¹Lund et al. [Qua-alluding to the past: a case of methaqualone analog ingestion](#)
- ▶ ²Fels et al. [Step-by-step procedure to identify previously unknown compounds with LC-QTOF-MS...](#)
- ▶ ³Romanek et al. [Return of the quaaludes? Prolonged agitated delirium after intentional ingestion...](#)
- ▶ ⁴Saito et al. [Pharmacological studies on 2-methyl-3\(2'-methyl-4'-chlorophenyl\)-5-chloro-4\[H\]-quinazolinone](#)
- ▶ ⁵Dongwook et al. [Facile synthesis of aryl and benzyl substituted 2-methylquinazoline-4\(3H\)-one...](#)

About: In collaboration with medical examiner and coroner offices, crime laboratories, clinical partners, and other stakeholders, the Center for Forensic Science Research and Education (CFSRE) is documenting first confirmations of NPS through analysis of drug materials and/or toxicology samples. These reports are generated using comprehensive analytical techniques (e.g., GC-MS, LC-QTOF-MS, NMR) and include available information about the new substances identified at the time of reporting, as well as the analytical data generated during testing. Our new drug monographs are intended to assist with the rapid identification of NPS in forensic casework and related disciplines, and should not be used for confirmatory purposes alone.

Analytical Notes: All identifications were made based on evaluation of analytical data (GC-MS and LC-QTOF-MS) in comparison to analysis of acquired reference material.

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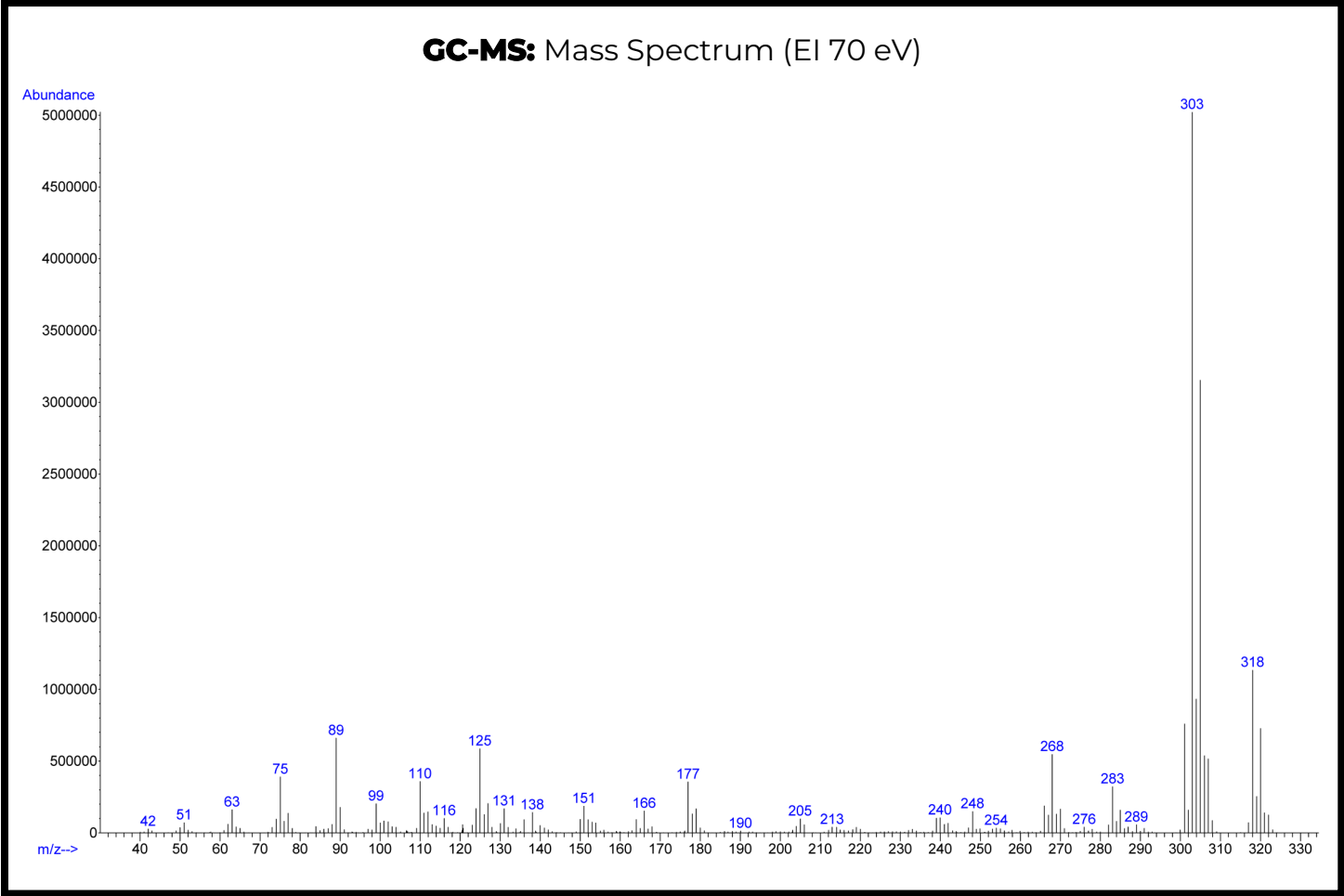
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Gas Chromatography Mass Spectrometry (GC-MS)

Laboratory: Center for Forensic Science Research and Education (CFSRE, Horsham PA, USA)
Sample Preparation: Acid-base extraction

Instrument: Agilent 5975 Series GC/MSD
Methods: [GC-MS Method Details](#) & [Monographs](#)



Liquid Chromatography Quadrupole Time-of-Flight Mass Spectrometry (LC-QTOF-MS)

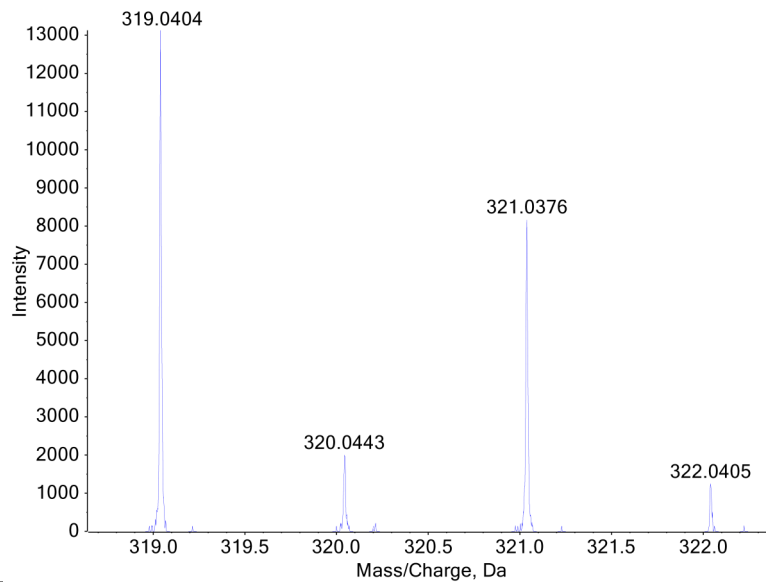
Laboratory: Center for Forensic Science Research and Education (CFSRE, Horsham, PA, USA)

Instrument: Sciex 5600+ LC-QTOF-MS

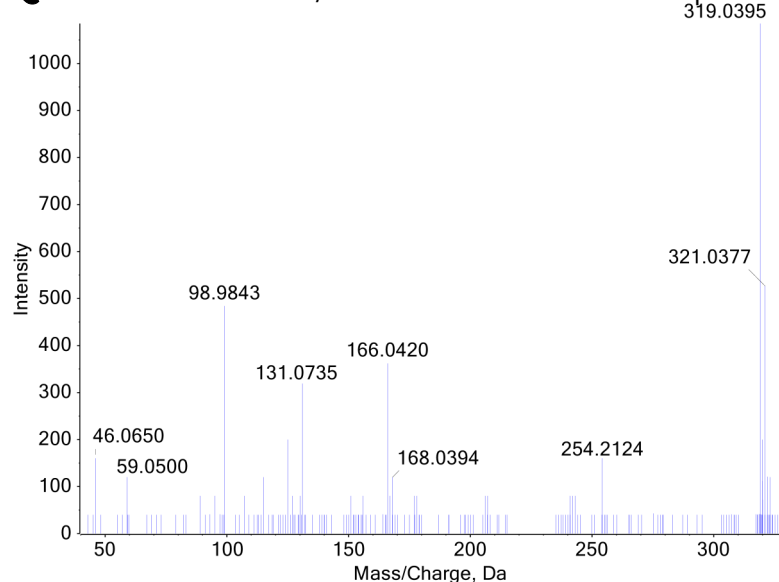
Methods: [LC-QTOF-MS Method Details](#) & [Monographs](#)

Sample Preparation: Dilution in mobile phase

LC-QTOF-MS: TOF-MS Precursor Ion Mass Spectrum



LC-QTOF-MS: TOF-MS/MS Product Ion Mass Spectrum



Confirmation Using Drug Standard: Reference material (Batch: 0587630-14) was purchased from Cayman Chemical (Ann Arbor, MI, USA). The analyte was confirmed to be dicloqualone based on retention time (sample: 8.56 min vs. standard: 8.54 min) and mass spectral data comparisons.