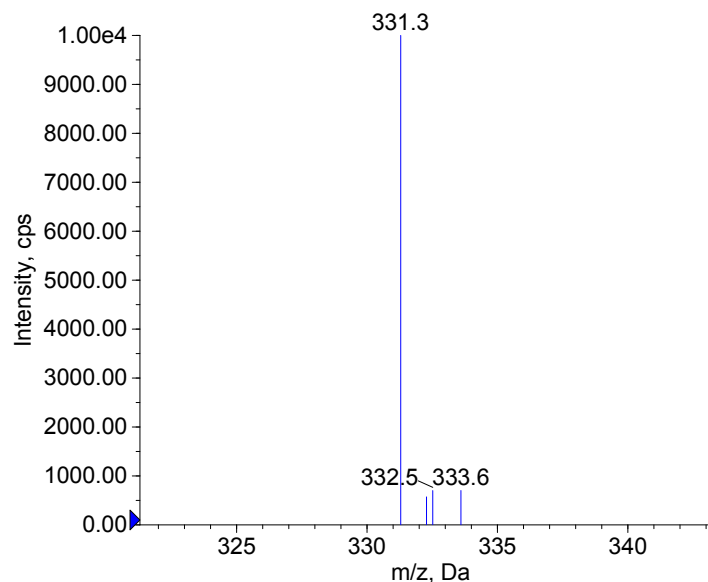


Compound Name: CP 47,497 C8 homologue
Synonyms: cannabicyclohexanol
Formula: C22H36O2
CAS Number: 70434-92-3
Molecular weight(Da): 332.2715
Compound Class: Synthetic Cannabinoids
Compound ID:
User Value: 0.0000
Keyword:
Internal Standard: No
Comment1: 332.271530

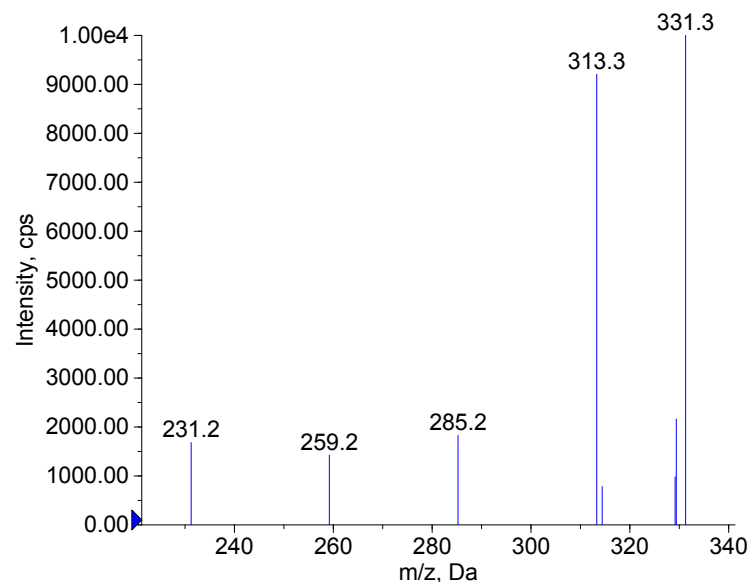
Instrument Model: 3200 Q TRAP
Ion Source: Turbo Spray
Polarity: Negative
Scan Type: EPI
1st Precursor m/z: 331.4000
Collision Energy1(V): -20.0000
Charge State1: -1
CAD Gas Type:
CAD Gas Value: Level 4
Retention Time(min): 1.20
Resolution: unit
Comment2:
Comment3:

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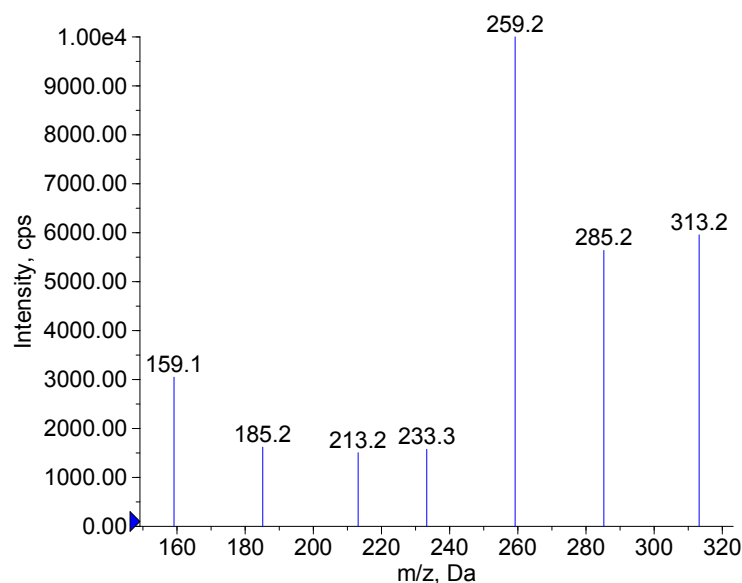
■ - CP 47,497 C8 homologue Prec(331.4) CE(-20.0) RT(1.2)



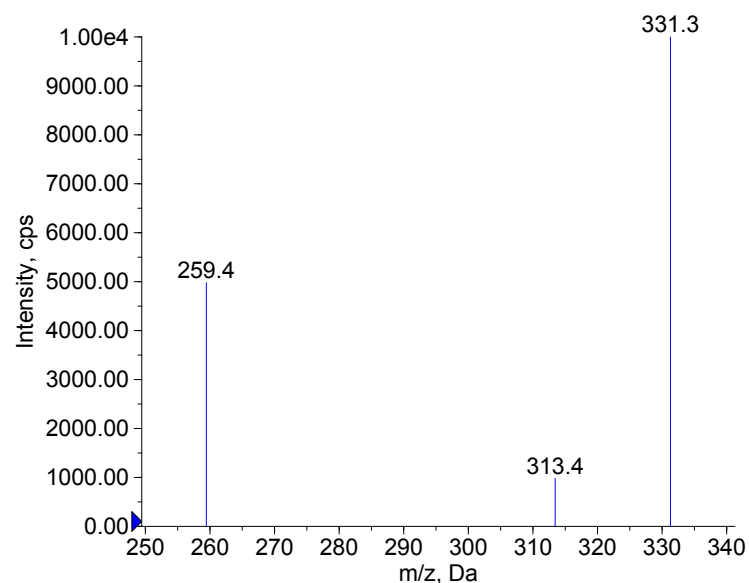
■ - CP 47,497 C8 homologue Prec(331.4) CE(-35.0) RT(1.2)



■ - CP 47,497 C8 homologue Prec(331.4) CE(-50.0) RT(1.3)



■ - CP 47,497 C8 homologue Prec(331.4) CE(-99.0) RT(1.3)



Comment1: monoisotopic mass
"CE(99.0)": CE spread 35 +/- 15 eV

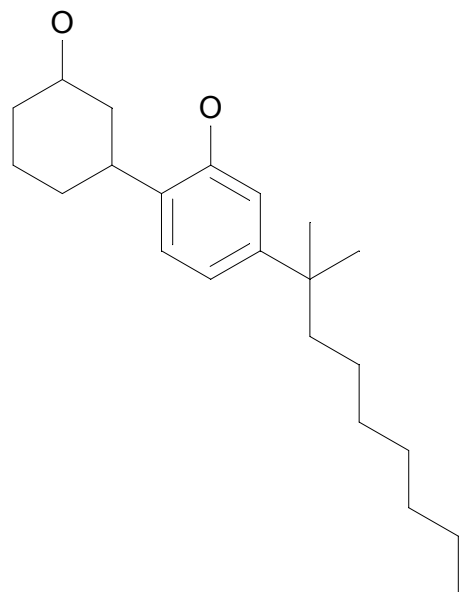
C22H36O2

Chemist Name:

Project Name:

Workbook Name:

Workbook Page:



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Comment1: monoisotopic mass
"CE(99.0)": CE spread 35 +/- 15 eV