

**Fleming  
Lee Shue**

*Sent via electronic mail (michael.maccabe@dec.ny.gov)*

June 13, 2023

Michael D. MacCabe, P.E.  
Senior Environmental Engineer  
Division of Environmental Remediation  
NYS Department of Environmental Conservation  
625 Broadway, 12th Floor  
Albany, NY 12233-7016

**Re: Annual Groundwater Monitoring Report  
388 Bridge Street Site - Brooklyn, New York  
BCP Site #C224134**

Dear Mr. MacCabe:

Fleming-Lee Shue Environmental Engineering and Geology, D.P.C. (FLS) presents this Annual Groundwater Monitoring Report for the 388 Bridge Street property (Site) located at 388 Bridge Street in Brooklyn, NY. The Site is currently in Site Management and this groundwater monitoring event was completed in accordance with the New York State Department of Environmental Conservation (NYSDEC) approved Site Management Plan (SMP) dated December 2013 and subsequent SMP Modifications dated October 2016 and July 2019.

## **Background**

Results from subsurface investigations performed by FLS from 2008 to 2010 detected tetrachloroethylene (PCE) in both soil and groundwater. In August 2009, the Site was accepted into the NYSDEC Brownfield Cleanup Program (BCP). Remedial activities were conducted in accordance with the NYSDEC-approved Remedial Action Work Plan dated April 2012. As a part of the approved remedy, on-site structures were demolished and the entire footprint of the site was excavated to a depth of up to 25 feet below surface grade (bsg) in preparation for redevelopment. The BCP Volunteer achieved a Track 2 remedy at the Site. After completion of the remedial work, residual contamination remained on-Site. Therefore, institutional and engineering controls were incorporated into the Site remedy to control exposure to the remaining contamination. This

included the installation of a sub-slab depressurization system (SSDS) within the neighboring 80 Willoughby Property (former St. Joseph's High School) and a soil vapor extraction system (SVE) within the new on-Site building in June 2013. The SVE system was installed to remove Volatile Organic Compounds (VOCs) from soil gas beneath the building slab. The system operated from 2013 through 2016 and included six extraction points (SVE-1, SVE-2, SVE-3, SVE-4, SVE-5 and SVE-6).

In 2016, after monitoring of PCE concentrations and prior approval of NYSDEC, the 2013 SVE system was downsized to limit extraction to where the bulk of the PCE mass remains (SVE-2). Each of the vapor extraction points, except for one location (SVE-2), were converted into groundwater monitoring wells (SVE-MW-1, SVE-MW-3, SVE-MW 4, SVE-MW-5 and SVE-MW-6) to monitor natural attenuation of VOCs. In July 2016 and with the prior approval of NYSDEC (dated July 29, 2016), SVE-MW-3 and SVE-MW-6 were abandoned because they did not extend into the groundwater table and were therefore not usable as groundwater monitoring wells. Off-Site monitoring wells, MW-3 and MW-7, were destroyed during construction activities.

At the time of this report, extraction well SVE-2 continues to operate as a part of the SVE system. Once remediation has been completed, extraction well SVE-2 will be converted to a groundwater monitoring well and serve as the downgradient well. Figure 2 presents the well locations and results from the last four rounds of groundwater sampling.

In its July 2019 Semi-Annual Groundwater Monitoring Report, FLS recommended a reduction in the frequency of groundwater monitoring from semi-annual to annual based on stable concentrations of PCE and TCE on-Site. NYSDEC approved this reduction in an email, dated July 18, 2019. The SMP was subsequently updated to reflect this amendment.

## **Geology and Hydrogeology**

### *Geology*

Regionally, Brooklyn at the western end of Long Island is underlain by sedimentary layers that strike northeast and are inclined gently to the southeast. These layers appear at or near the surface in the vicinity of Long Island Sound, where differential erosion has left relatively tough sands and clays at elevations of more than 60 feet above sea level.

Resting on top of these sands and clays and forming the highest elevation is a belt of glacially deposited debris composed of an unsorted, unstratified mixture of boulders, sand, silt, and clay. This debris was deposited in the interval between 75,000 and 17,000 years ago when the area was covered by a massive sheet of glacial ice. In the vicinity of New York, the ice was moving in a generally southerly direction, bringing with it a huge load of detached bedrock, sediment, and soil that it had scoured from more northerly regions

The geology observed at the Site can be characterized as two major strata groups. The upper stratum is a fill material layer that ranges in thickness from 3 to 17 feet. Beneath the fill layer is a natural glacial till deposit consisting of brown to red-brown, fine- to coarse-grained sand with trace cobbles and boulders.

### *Hydrology*

Historically, groundwater has been encountered at approximately 43 to 45 feet below ground surface. Local groundwater flow during this sampling event was to the north-northeast. This localized groundwater flow direction may be influenced by the subway tunnels located north and southwest of the Site and pumping operations at a Metropolitan Transportation Authority (MTA) de-watering station located within 1.5 miles northeast of the Site. Figure 3 presents a groundwater contour map.

### **Groundwater Monitoring Program**

The groundwater monitoring program for the Site began in March 2016. The groundwater monitoring program was implemented to monitor natural attenuation of VOCs in groundwater following the downsizing of the SVE system. The SVE system, installed in 2013, was downsized and modified in 2016 to target the area where the bulk of the contaminant mass remains, primarily in the area of well SVE-2. Selected soil vapor extraction wells were converted to monitoring wells and included in the groundwater monitoring program. The objectives of the groundwater monitoring program include the following:

- Provide a current round of groundwater analytical data from the monitoring wells;
- Evaluate the existing and time-based groundwater conditions at the Site; and
- Evaluate the time-based trends of VOCs.

The groundwater monitoring program involves the following activities:

- Measurement of groundwater field parameters including depth to water, pH, dissolved oxygen (DO), total dissolved solids (TDS), conductivity, oxidation-reduction potential (ORP), turbidity, salinity, and temperature to determine groundwater conditions;
- Collection of groundwater samples for VOCs to evaluate chlorinated VOC concentration trends and monitor natural attenuation;
- Collection of groundwater samples for geochemical parameters including nitrate, nitrite, sulfate, iron (II), total organic carbon, and dissolved organic carbon to evaluate evidence supporting natural attenuation.

## Groundwater Sampling Procedures

On March 28, 2023, groundwater samples were collected from the three (3) on-Site monitoring wells (SVE-MW-1, SVE-MW 4, and SVE-MW-5). Prior to sampling, FLS screened the well headspace for VOCs with a MiniRAE 3000 Photoionization detector (PID) and collected depth to water measurements. Based on depth to water measurements, groundwater flow is estimated to be to the north-northeast (Figure 3). Groundwater samples were collected using the low-flow sampling method (EPA Low-Flow Groundwater Sampling Procedures, April 1996). Each monitoring well was purged prior to sampling using a peristaltic pump until groundwater parameters (temperature, pH, DO, conductivity, ORP, TDS, and turbidity) stabilized, or three well volumes were purged. Water-quality measurements were monitored using a Horiba U-52 multi-parameter water-quality meter. The monitoring well purge logs are included in Appendix A.

After the stabilization of the groundwater parameters, samples were collected via dedicated pump tubing directly into laboratory-supplied containers. After sample collection each container was labeled, placed on ice in an insulated cooler and transported under chain-of-custody protocol to SGS/Accutest Laboratories of Dayton, New Jersey, a New York Environmental Laboratory Approval Program Certified Laboratory. The groundwater samples were analyzed for Target Compounds List VOCs by EPA Method 8260C and several geochemical parameters.

## Summary of Analytical Results

The groundwater analytical results, from this event and the past ten sampling events, were compared to the NYSDEC Division of Water Technical and Operational Guidance Series 1.1.1 Ambient Water Quality Standards and Guidance Values (TOGS) and are summarized in Table 1. The laboratory data report for this most recent event is provided in Appendix B.

Overall, the groundwater analytical results were non-detect for all VOCs with the exception of chlorinated solvent PCE and its associated daughter products. Analytical results indicate that PCE remain at similar concentrations to previous events with the exception of SVE-MW-5 where PCE was non-detect for the first time since monitoring began in March 2016. PCE only exceeded the TOGS standard of 5 µg/L at SVE-MW-4 (15 µg/L) this event and was the only detection above its standard in all three samples. PCE was detected in monitoring well SVE-MW-1 just under the TOGS standard at 4.9 µg/L. Concentrations of trichloroethylene (TCE) and cis-1, 2-dichloroethylene continue to remain below the TOGS standards (5 µg/L) in all three monitoring wells. Chloroform, which has historically been observed on Site, was detected in all monitoring wells, but remained well below the TOGS standard of 7 µg/L.

An overview of the trends of PCE and TCE concentrations from March 2016 to March 2023 is presented in the attached graphs. Analytical results from this event show that PCE concentrations generally remained at similar low-level concentrations with decreases observed in SVE-MW-4 (28.8 µg/L to 15 µg/L) and in SVE-MW-5 (21.4 µg/L to non-detect). Overall PCE, concentrations remain well below historical maximum concentrations of 72 µg/L in SVE-MW-4 and 39.3 µg/L in SVE-MW-5 in September 2018. TCE concentrations have remained below the 5 µg/L TOGS standard in all wells since September 2017, and continue to display an asymptotic trend.

## **Conclusions and Recommendations**

Overall, groundwater results were non-detect for all VOCs on Site with the exception of chlorinated solvent PCE and its associated daughter products. That said, concentrations of chlorinated VOCs have largely remained at similar or lower concentrations within all three monitoring wells. PCE decreased in SVE-MW-5 to non-detect for the first time since groundwater monitoring began in March 2016. The only exceedance from this sampling event was of PCE in SVE-MW-4 at a concentration of 15 µg/L, but overall concentrations remain well below historical maximums from September 2018 and have been at or approaching asymptotic levels since groundwater monitoring began in 2016. TCE, cis-1, 2 DCE and Chloroform concentrations remain below their respective TOGS standards in all monitoring wells.

FLS recommends continuing groundwater monitoring on an annual basis in order to monitor natural attenuation trends on Site and assess groundwater quality.

Please contact us with any comments or questions.

Sincerely,

Fleming-Lee Shue Environmental Engineering and Geology, D.P.C.

Arnold F. Flemming

Arnold Fleming  
President

cc: Roger Fortune Stahl Realty  
Jennifer Coghlan, Esq. Sive, Paget & Riesel  
Joel Kane Fleming-Lee-Shue, D.P.C.

enc: Table 1 Groundwater Sampling Analytical Results

Figure 1 Site Location Map

- Figure 2      Site Plan and Groundwater Sampling Results  
Figure 3      Groundwater Contour Map
- Graphs        PCE and TCE Concentration Trends
- Appendix A    Monitoring Well Purge Logs  
Appendix B    Laboratory Analytical Data Report

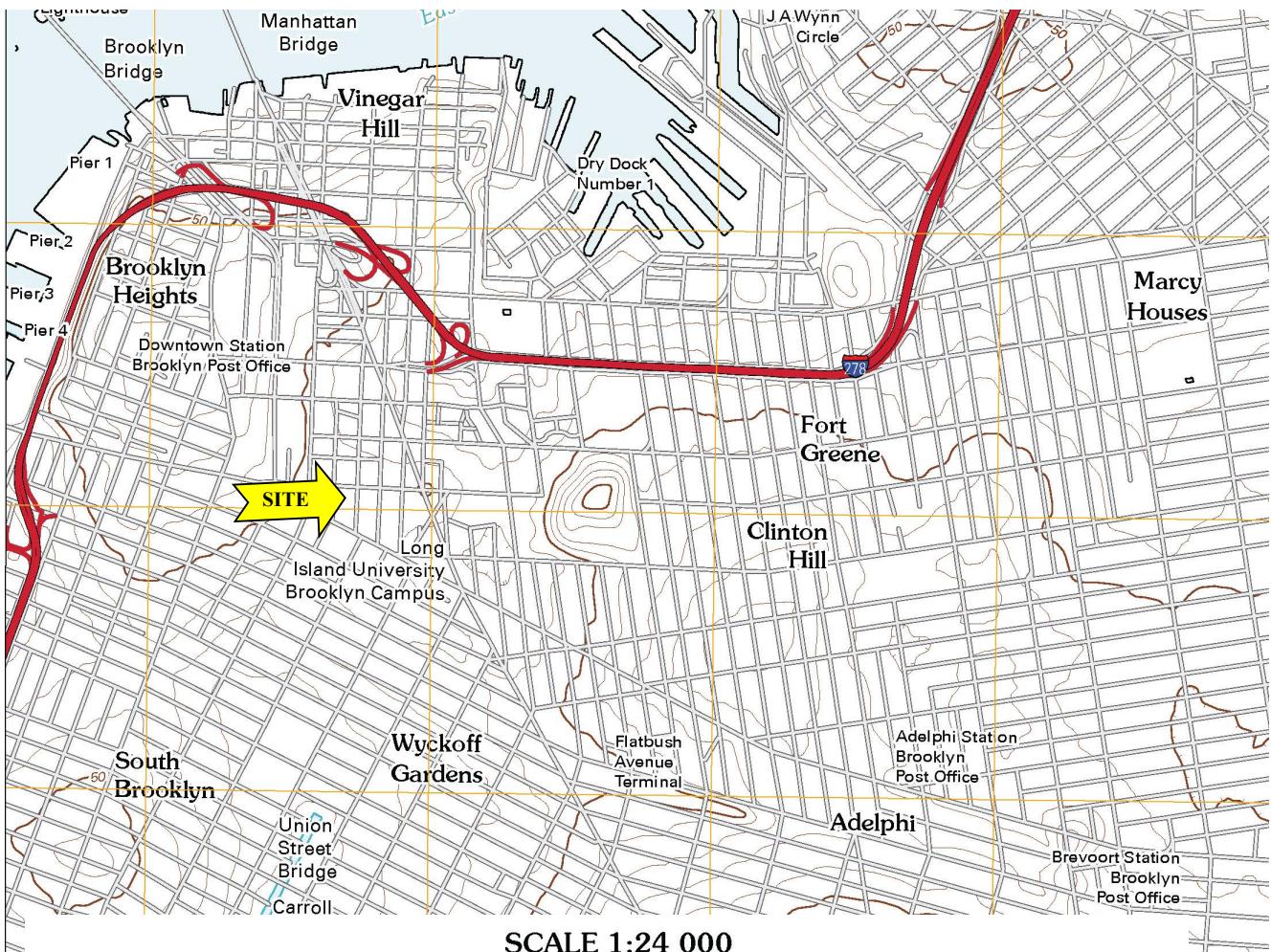
# Tables



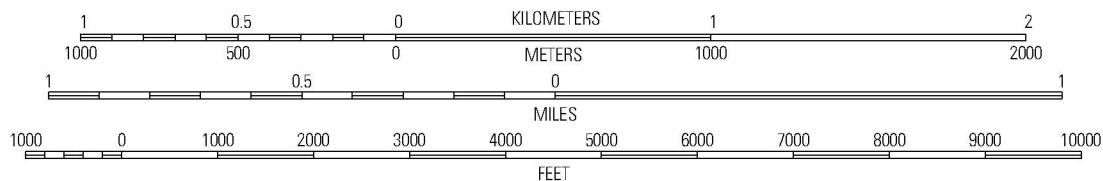


# Figures





SCALE 1:24 000



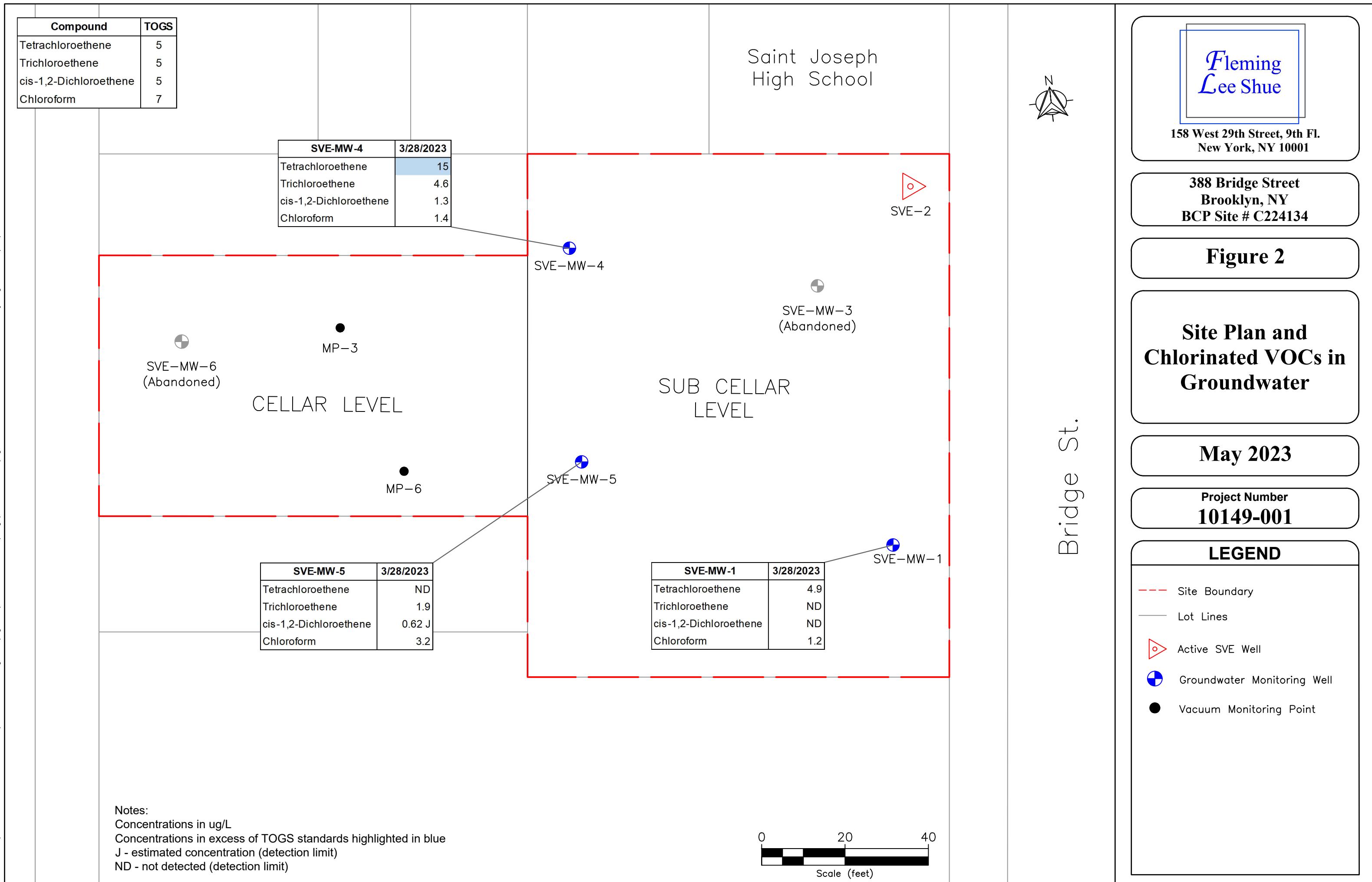
CONTOUR INTERVAL 10 FEET

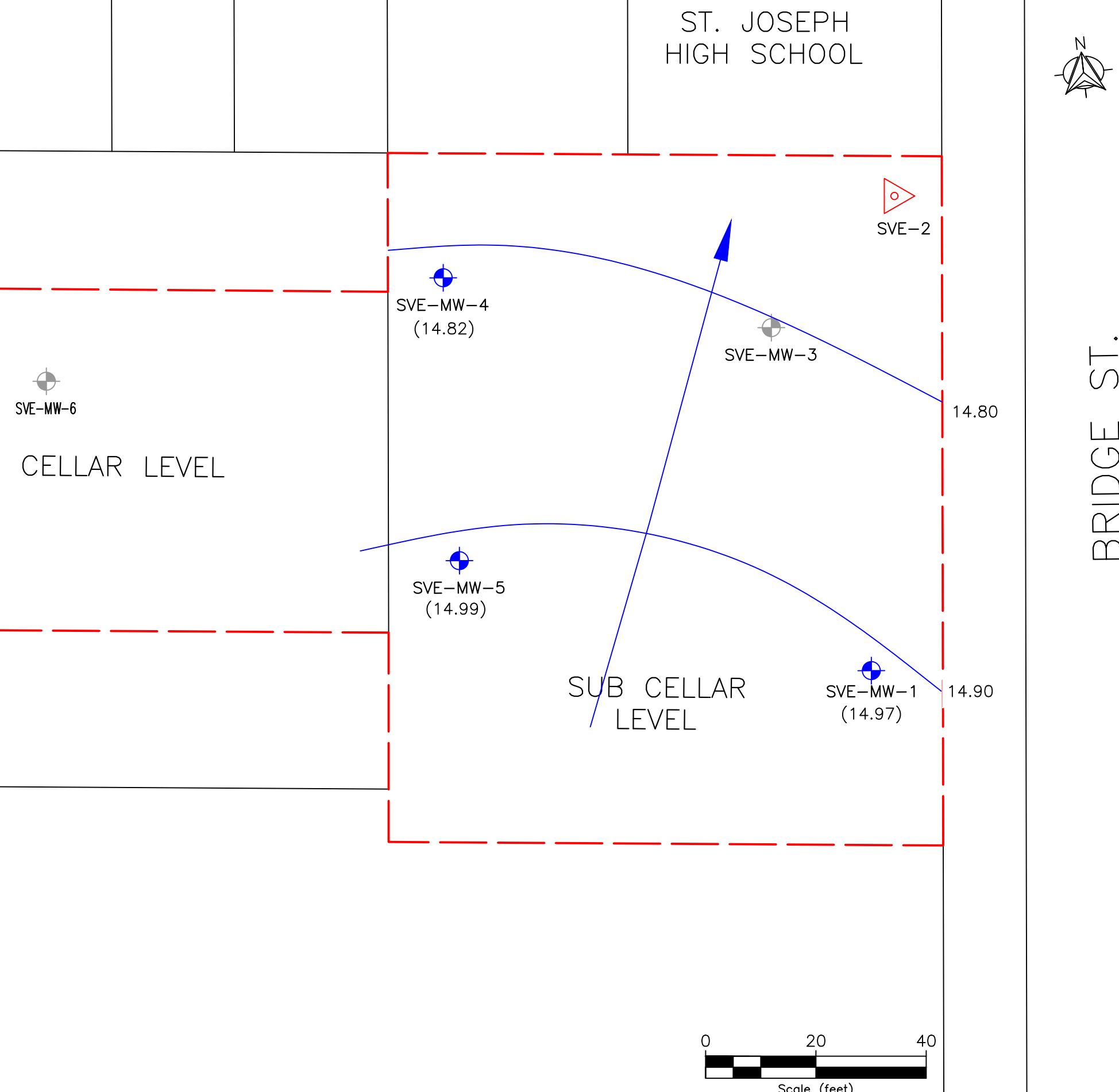
Site: Brooklyn Quadrangle, New York 7.5 Minute series USGS Topographic Map (79287)  
Obtained from United States Geological Survey topography compiled 2010

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## FIGURE 1: SITE LOCATION MAP

SITE: 388 Bridge Street  
Brooklyn, New York





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Lee Shue

158 West 29th Street, 9th Fl.  
New York, NY 10001

388 Bridge Street  
Brooklyn, NY  
BCP Site # C224134

**Figure 3**

**Groundwater  
Elevation  
Contour**

**May 2023**

**Project Number**  
**10149-001**

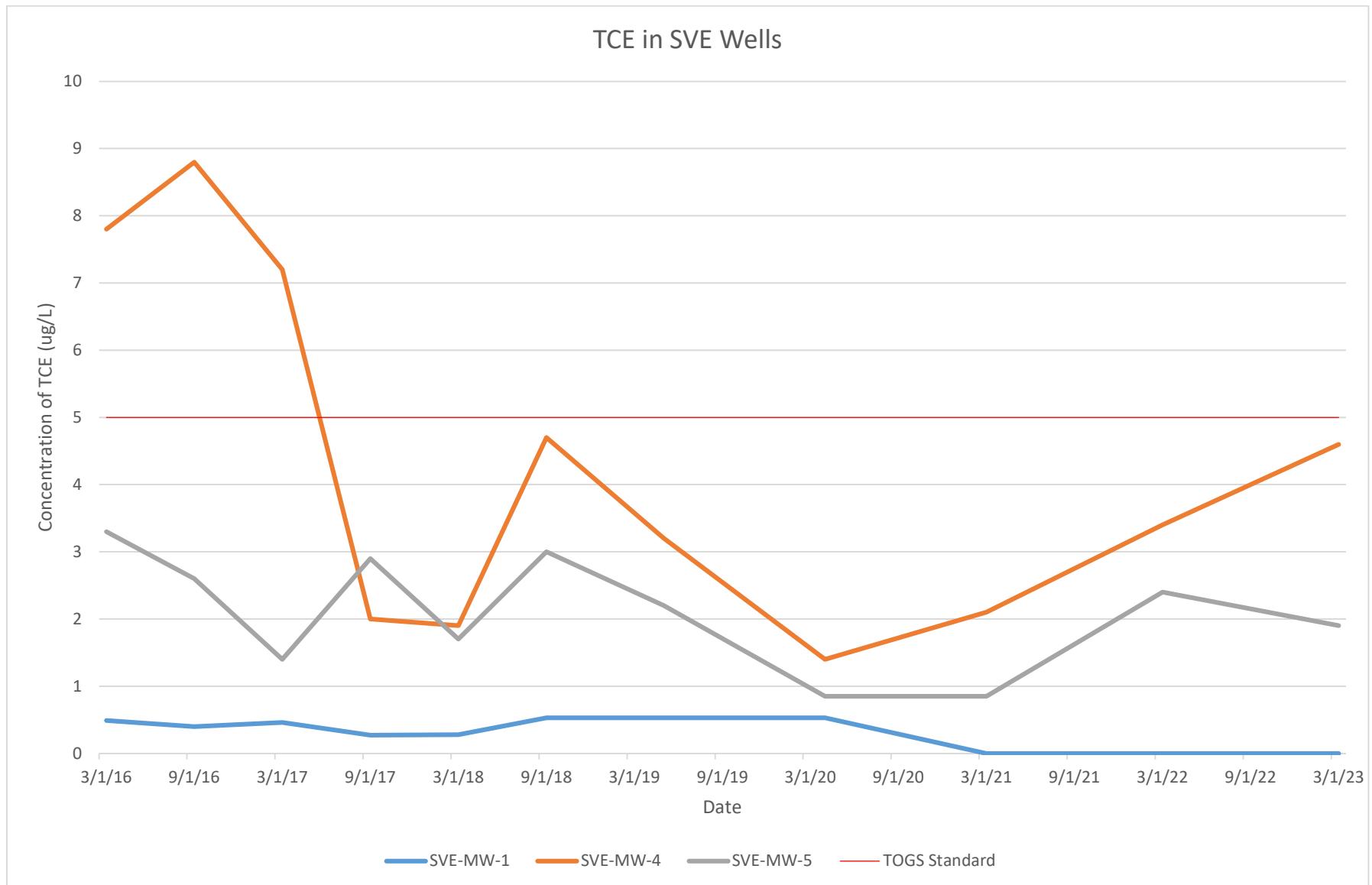
**LEGEND**

- SITE BOUNDARY
- SVE WELL
- MONITORING WELL
- ABANDONED MONITORING WELL
- GROUNDWATER ELEVATION
- GROUNDWATER FLOW DIRECTION

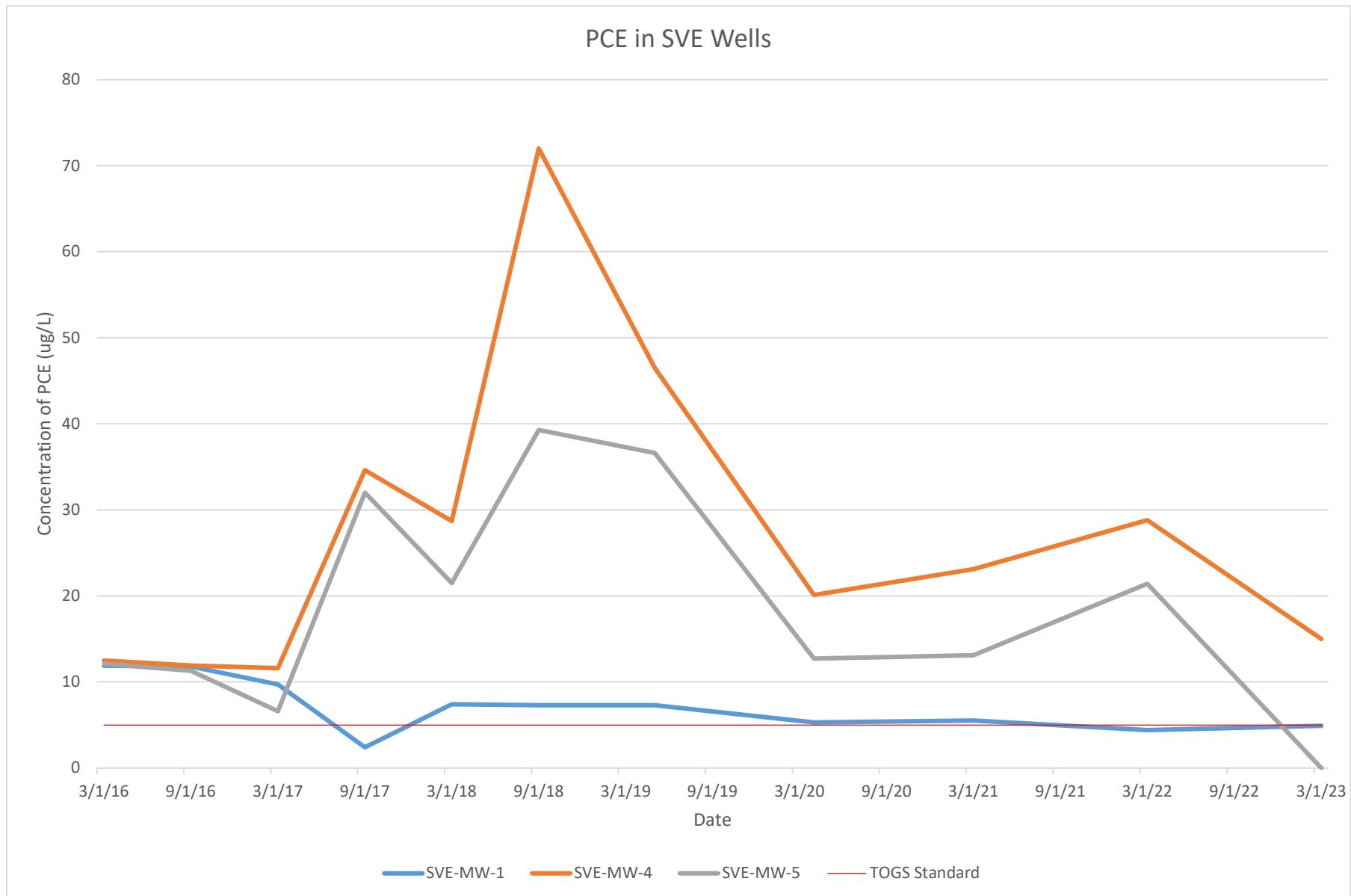
# Graphs



Graphs – Contaminant Concentration Trends  
Annual Groundwater Report  
388 Bridge Street, Brooklyn, NY



Graphs – Contaminant Concentration Trends  
Annual Groundwater Report  
388 Bridge Street, Brooklyn, NY



# Appendix A

## Monitoring Well Purge Logs









# Appendix B

## Laboratory Analytical Data Report



The results set forth herein are provided by SGS North America Inc.

**e-Hardcopy 2.0**  
*Automated Report*

## Technical Report for

**Fleming-Lee Shue, Inc.**

**388 Bridge Street, Brooklyn, NY**

**10188**

**SGS Job Number: JD62888**

**Sampling Date: 03/28/23**



### Report to:

**Fleming-Lee Shue, Inc.  
158 West 29th Street 9th Floor  
New York, NY 10001  
joel@flemingleeshue.com**

**ATTN: Steve Panter**

**Total number of pages in report: 628**



Test results contained within this data package meet the requirements  
of the National Environmental Laboratory Accreditation Program  
and/or state specific certification programs as applicable.

**David Chastain  
General Manager**

**Client Service contact: Tammy McCloskey 732-329-0200**

Certifications: NJ(12129), NY(10983), CA, CT, FL, IL, IN, KS, KY, LA, MA, MD, ME, MN, NC,  
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Test results relate only to samples analyzed.

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## Sample Summary

Fleming-Lee Shue, Inc.

Job No: JD62888

388 Bridge Street, Brooklyn, NY  
Project No: 10188

Sample Number	Collected Date	Time By	Matrix Received	Code Type	Client Sample ID
---------------	----------------	---------	-----------------	-----------	------------------

This report contains results reported as ND = Not detected. The following applies:  
Organics ND = Not detected above the MDL

JD62888-1 03/28/23 11:55 BH 03/29/23 AQ Ground Water SVE-1

JD62888-1F 03/28/23 11:55 BH 03/29/23 AQ Groundwater Filtered SVE-1

JD62888-2 03/28/23 11:50 BH 03/29/23 AQ Ground Water SVE-4

JD62888-2F 03/28/23 11:50 BH 03/29/23 AQ Groundwater Filtered SVE-4

JD62888-3 03/28/23 10:05 BH 03/29/23 AQ Ground Water SVE-5

JD62888-3F 03/28/23 10:05 BH 03/29/23 AQ Groundwater Filtered SVE-5

JD62888-4 03/28/23 11:30 BH 03/29/23 AQ Field Blank Water FB-1

JD62888-4F 03/28/23 11:30 BH 03/29/23 AQ Field Blank Filtered FB-1

JD62888-5 03/28/23 11:55 03/29/23 AQ Trip Blank Water TRIP BLANK

## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** Fleming-Lee Shue, Inc.

**Job No:** JD62888

**Site:** 388 Bridge Street, Brooklyn, NY

**Report Date** 4/7/2023 3:38:20 PM

On 03/29/2023, 3 sample(s), 1 Trip Blank(s), and 2 Field Blank(s) were received at SGS North America Inc. (SGS) at a temperature of 2.1 °C. The samples were intact and properly preserved, unless noted below. An SGS Job Number of JD62888 was assigned to the project. The lab sample ID, client sample ID, and date of sample collection are detailed in the report's Results Summary.

Specified quality control criteria were achieved for this job except as noted below. For more information, please refer to the analytical results and QC summary pages.

### MS Volatiles By Method SW846 8260D

**Matrix:** AQ

**Batch ID:** V1R100

- All samples were analyzed within the recommended method holding time.
- Sample(s) JD62953-1MS, JD62953-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- Matrix Spike/Matrix Spike Duplicate recovery(s) of Toluene are outside control limits. Outside control limits due to high level in sample relative to spike amount.
- JD62888-4 for Vinyl chloride: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-3 for Vinyl chloride: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-1 for Acetone: Associated CCV outside of control limits high, sample was ND.
- JD62888-1 for Bromomethane: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-1 for Vinyl chloride: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-4 for Bromomethane: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-5 for Acetone: Associated CCV outside of control limits high, sample was ND.
- JD62888-5 for Bromomethane: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-5 for Vinyl chloride: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-2 for Acetone: Associated CCV outside of control limits high, sample was ND.
- JD62888-2 for Bromomethane: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-2 for Vinyl chloride: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-3 for Acetone: Associated CCV outside of control limits high, sample was ND.
- JD62888-3 for Bromomethane: Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.
- JD62888-4 for Acetone: Associated CCV outside of control limits high, sample was ND.

## General Chemistry By Method EPA 300/SW846 9056A

**Matrix:** AQ

**Batch ID:** GP45907

- All samples were prepared within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JD62888-3DUP, JD62888-3MS were used as the QC samples for the Sulfate analysis.

## General Chemistry By Method EPA 353.2/LACHAT

**Matrix:** AQ

**Batch ID:** GP45893

- All samples were prepared within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JD62498-5DUP, JD62498-5MS were used as the QC samples for the Nitrogen, Nitrate + Nitrite analysis.
- The matrix spike (MS) recovery(s) of Nitrogen, Nitrate + Nitrite are outside control limits. Spike recovery indicates possible matrix interference.

## General Chemistry By Method EPA353.2/SM4500NO2B

**Matrix:** AQ

**Batch ID:** R202695

- The data for EPA353.2/SM4500NO2B meets quality control requirements.
- JD62888-1 for Nitrogen, Nitrate: Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

**Matrix:** AQ

**Batch ID:** R202696

- The data for EPA353.2/SM4500NO2B meets quality control requirements.
- JD62888-2 for Nitrogen, Nitrate: Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

**Matrix:** AQ

**Batch ID:** R202697

- The data for EPA353.2/SM4500NO2B meets quality control requirements.
- JD62888-3 for Nitrogen, Nitrate: Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

**Matrix:** AQ

**Batch ID:** R202698

- The data for EPA353.2/SM4500NO2B meets quality control requirements.
- JD62888-4 for Nitrogen, Nitrate: Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

## General Chemistry By Method SM3500FE B-11

**Matrix:** AQ

**Batch ID:** GN39908

- All method blanks for this batch meet method specific criteria.
- Sample(s) JD62888-1MS, JD62888-1MSD were used as the QC samples for the Iron, Ferrous analysis.
- JD62888-4 for Iron, Ferrous: Field analysis required. Received out of hold time and analyzed by request.
- JD62888-3 for Iron, Ferrous: Field analysis required. Received out of hold time and analyzed by request.
- JD62888-2 for Iron, Ferrous: Field analysis required. Received out of hold time and analyzed by request.
- JD62888-1 for Iron, Ferrous: Field analysis required. Received out of hold time and analyzed by request.

## General Chemistry By Method SM4500NO2 B-11

**Matrix:** AQ

**Batch ID:** GN39776

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JD62888-1MS, JD62888-1MSD were used as the QC samples for the Nitrogen, Nitrite analysis.

**General Chemistry By Method SM5310 B-11****Matrix:** AQ**Batch ID:** GP45881

- All samples were prepared within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JD62888-2FMS, JD62888-2FMSD were used as the QC samples for the Dissolved Organic Carbon analysis.

**General Chemistry By Method SM5310 B-11/14****Matrix:** AQ**Batch ID:** GP45861

- All samples were prepared within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JD62808-1MS, JD62808-1MSD were used as the QC samples for the Total Organic Carbon analysis.

SGS certifies that data reported for samples received, listed on the associated custody chain or analytical task order, were produced to specifications meeting SGS's Quality System precision, accuracy and completeness objectives except as noted.

Estimated non-standard method measurement uncertainty data is available on request, based on quality control bias and implicit for standard methods. Acceptable uncertainty requires tested parameter quality control data to meet method criteria.

SGS is not responsible for data quality assumptions if partial reports are used and recommends that this report be used in its entirety. This report is authorized by SGS indicated via signature on the report cover.

## Summary of Hits

Page 1 of 2

Job Number: JD62888

Account: Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Collected: 03/28/23

3

Lab Sample ID Analyte	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
--------------------------	------------------	-----------------	----	-----	-------	--------

JD62888-1 SVE-1

Chloroform	1.2	1.0	0.50	ug/l	SW846 8260D
Tetrachloroethene	4.9	1.0	0.56	ug/l	SW846 8260D
Nitrogen, Nitrate <sup>a</sup>	9.2	0.41		mg/l	EPA353.2/SM4500NO2B
Nitrogen, Nitrate + Nitrite	9.2	0.40		mg/l	EPA 353.2/LACHAT
Sulfate	93.5	2.0		mg/l	EPA 300/SW846 9056A
Total Organic Carbon	1.0	1.0		mg/l	SM5310 B-11/14

JD62888-1F SVE-1

Dissolved Organic Carbon	1.0	1.0		mg/l	SM5310 B-11
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JD62888-2 SVE-4

Chloroform	1.4	1.0	0.50	ug/l	SW846 8260D
cis-1,2-Dichloroethene	1.3	1.0	0.51	ug/l	SW846 8260D
Tetrachloroethene	15.0	1.0	0.56	ug/l	SW846 8260D
Trichloroethene	4.6	1.0	0.53	ug/l	SW846 8260D
Nitrogen, Nitrate <sup>a</sup>	5.5	0.21		mg/l	EPA353.2/SM4500NO2B
Nitrogen, Nitrate + Nitrite	5.5	0.20		mg/l	EPA 353.2/LACHAT
Sulfate	79.9	2.0		mg/l	EPA 300/SW846 9056A

JD62888-2F SVE-4

No hits reported in this sample.

JD62888-3 SVE-5

Chloroform	3.2	1.0	0.50	ug/l	SW846 8260D
cis-1,2-Dichloroethene	0.62 J	1.0	0.51	ug/l	SW846 8260D
Trichloroethene	1.9	1.0	0.53	ug/l	SW846 8260D
Nitrogen, Nitrate <sup>a</sup>	7.5	0.31		mg/l	EPA353.2/SM4500NO2B
Nitrogen, Nitrate + Nitrite	7.5	0.30		mg/l	EPA 353.2/LACHAT
Sulfate	96.8	2.0		mg/l	EPA 300/SW846 9056A
Total Organic Carbon	1.1	1.0		mg/l	SM5310 B-11/14

JD62888-3F SVE-5

No hits reported in this sample.

JD62888-4 FB-1

No hits reported in this sample.

**Summary of Hits**

Job Number: JD62888  
Account: Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY  
Collected: 03/28/23

C3

Lab Sample ID	Client Sample ID	Result/ Analyte	Qual	RL	MDL	Units	Method
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JD62888-4F FB-1

No hits reported in this sample.

JD62888-5 TRIP BLANK

No hits reported in this sample.

(a) Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

**Sample Results**

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**Report of Analysis**

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SGS North America Inc.

## Report of Analysis

Page 1 of 2

**Client Sample ID:** SVE-1  
**Lab Sample ID:** JD62888-1  
**Matrix:** AQ - Ground Water  
**Method:** SW846 8260D  
**Project:** 388 Bridge Street, Brooklyn, NY

**Date Sampled:** 03/28/23  
**Date Received:** 03/29/23  
**Percent Solids:** n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1R03027.D	1	04/04/23 15:45	NW	n/a	n/a	V1R100
Run #2							

**Purge Volume**  
Run #1 5.0 ml  
Run #2

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone <sup>a</sup>	ND	10	3.1	ug/l	
71-43-2	Benzene	ND	0.50	0.43	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.45	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>b</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	2.7	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.46	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	1.2	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
110-82-7	Cyclohexane	ND	5.0	0.78	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	0.53	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.56	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.51	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.54	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
123-91-1	1,4-Dioxane	ND	130	19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
76-13-1	Freon 113	ND	5.0	0.58	ug/l	

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

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## Report of Analysis

Page 2 of 2

Client Sample ID:	SVE-1	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-1	Date Received:	03/29/23
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	SW846 8260D		
Project:	388 Bridge Street, Brooklyn, NY		

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
591-78-6	2-Hexanone	ND	5.0	2.0	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
79-20-9	Methyl Acetate	ND	5.0	0.80	ug/l	
108-87-2	Methylcyclohexane	ND	5.0	0.60	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.9	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
100-42-5	Styrene	ND	1.0	0.49	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	4.9	1.0	0.56	ug/l	
108-88-3	Toluene	ND	1.0	0.49	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.40	ug/l	
75-01-4	Vinyl chloride <sup>b</sup>	ND	1.0	0.52	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
1330-20-7	Xylene (total)	ND	1.0	0.59	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	109%		80-120%
17060-07-0	1,2-Dichloroethane-D4	108%		80-120%
2037-26-5	Toluene-D8	94%		80-120%
460-00-4	4-Bromofluorobenzene	102%		82-114%

- (a) Associated CCV outside of control limits high, sample was ND.  
 (b) Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

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**Report of Analysis**

Page 1 of 1

Client Sample ID:	SVE-1	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-1	Date Received:	03/29/23
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Iron, Ferrous <sup>a</sup>	< 0.20	0.20	mg/l	1	04/02/23 13:49	MK	SM3500FE B-11
Nitrogen, Nitrate <sup>b</sup>	9.2	0.41	mg/l	1	03/30/23 15:29	MM	EPA353.2/SM4500NO2B
Nitrogen, Nitrate + Nitrite	9.2	0.40	mg/l	4	03/30/23 15:29	MM	EPA 353.2/LACHAT
Nitrogen, Nitrite	< 0.010	0.010	mg/l	1	03/30/23 08:36	MP	SM4500NO2 B-11
Sulfate	93.5	2.0	mg/l	1	03/30/23 22:25	SS	EPA 300/SW846 9056A
Total Organic Carbon	1.0	1.0	mg/l	1	03/30/23 15:53	MB	SM5310 B-11/14

(a) Field analysis required. Received out of hold time and analyzed by request.

(b) Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

Client Sample ID:	SVE-1	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-1F	Date Received:	03/29/23
Matrix:	AQ - Groundwater Filtered	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Dissolved Organic Carbon	1.0	1.0	mg/l	1	03/31/23 21:17	MB	SM5310 B-11

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RL = Reporting Limit

SGS North America Inc.

## Report of Analysis

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4.3  
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**Client Sample ID:** SVE-4  
**Lab Sample ID:** JD62888-2  
**Matrix:** AQ - Ground Water  
**Method:** SW846 8260D  
**Project:** 388 Bridge Street, Brooklyn, NY

**Date Sampled:** 03/28/23  
**Date Received:** 03/29/23  
**Percent Solids:** n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1R03034.D	1	04/04/23 18:45	NW	n/a	n/a	V1R100
Run #2							

Purge Volume  
Run #1 5.0 ml  
Run #2

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone <sup>a</sup>	ND	10	3.1	ug/l	
71-43-2	Benzene	ND	0.50	0.43	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.45	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>b</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	2.7	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.46	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	1.4	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
110-82-7	Cyclohexane	ND	5.0	0.78	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	0.53	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.56	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
156-59-2	cis-1,2-Dichloroethene	1.3	1.0	0.51	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.54	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
123-91-1	1,4-Dioxane	ND	130	19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
76-13-1	Freon 113	ND	5.0	0.58	ug/l	

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

Page 2 of 2

Client Sample ID:	SVE-4	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-2	Date Received:	03/29/23
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	SW846 8260D		
Project:	388 Bridge Street, Brooklyn, NY		

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
591-78-6	2-Hexanone	ND	5.0	2.0	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
79-20-9	Methyl Acetate	ND	5.0	0.80	ug/l	
108-87-2	Methylcyclohexane	ND	5.0	0.60	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.9	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
100-42-5	Styrene	ND	1.0	0.49	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	15.0	1.0	0.56	ug/l	
108-88-3	Toluene	ND	1.0	0.49	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	4.6	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.40	ug/l	
75-01-4	Vinyl chloride <sup>b</sup>	ND	1.0	0.52	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
1330-20-7	Xylene (total)	ND	1.0	0.59	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	111%		80-120%
17060-07-0	1,2-Dichloroethane-D4	108%		80-120%
2037-26-5	Toluene-D8	94%		80-120%
460-00-4	4-Bromofluorobenzene	102%		82-114%

- (a) Associated CCV outside of control limits high, sample was ND.  
 (b) Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

**Report of Analysis**

Page 1 of 1

Client Sample ID:	SVE-4	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-2	Date Received:	03/29/23
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Iron, Ferrous <sup>a</sup>	< 0.20	0.20	mg/l	1	04/02/23 13:49	MK	SM3500FE B-11
Nitrogen, Nitrate <sup>b</sup>	5.5	0.21	mg/l	1	03/30/23 15:30	MM	EPA353.2/SM4500NO2B
Nitrogen, Nitrate + Nitrite	5.5	0.20	mg/l	2	03/30/23 15:30	MM	EPA 353.2/LACHAT
Nitrogen, Nitrite	< 0.010	0.010	mg/l	1	03/30/23 08:36	MP	SM4500NO2 B-11
Sulfate	79.9	2.0	mg/l	1	03/30/23 22:39	SS	EPA 300/SW846 9056A
Total Organic Carbon	< 1.0	1.0	mg/l	1	03/30/23 16:04	MB	SM5310 B-11/14

(a) Field analysis required. Received out of hold time and analyzed by request.

(b) Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

Client Sample ID:	SVE-4	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-2F	Date Received:	03/29/23
Matrix:	AQ - Groundwater Filtered	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Dissolved Organic Carbon	< 1.0	1.0	mg/l	1	03/31/23 19:59	MB	SM5310 B-11

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RL = Reporting Limit

SGS North America Inc.

## Report of Analysis

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4.5  
4

**Client Sample ID:** SVE-5  
**Lab Sample ID:** JD62888-3  
**Matrix:** AQ - Ground Water  
**Method:** SW846 8260D  
**Project:** 388 Bridge Street, Brooklyn, NY

**Date Sampled:** 03/28/23  
**Date Received:** 03/29/23  
**Percent Solids:** n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1R03035.D	1	04/04/23 19:10	NW	n/a	n/a	V1R100
Run #2							

Purge Volume  
Run #1 5.0 ml  
Run #2

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone <sup>a</sup>	ND	10	3.1	ug/l	
71-43-2	Benzene	ND	0.50	0.43	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.45	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>b</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	2.7	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.46	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	3.2	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
110-82-7	Cyclohexane	ND	5.0	0.78	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	0.53	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.56	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
156-59-2	cis-1,2-Dichloroethene	0.62	1.0	0.51	ug/l	J
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.54	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
123-91-1	1,4-Dioxane	ND	130	19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
76-13-1	Freon 113	ND	5.0	0.58	ug/l	

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Report of Analysis

Page 2 of 2

Client Sample ID:	SVE-5	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-3	Date Received:	03/29/23
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	SW846 8260D		
Project:	388 Bridge Street, Brooklyn, NY		

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
591-78-6	2-Hexanone	ND	5.0	2.0	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
79-20-9	Methyl Acetate	ND	5.0	0.80	ug/l	
108-87-2	Methylcyclohexane	ND	5.0	0.60	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.9	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
100-42-5	Styrene	ND	1.0	0.49	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.56	ug/l	
108-88-3	Toluene	ND	1.0	0.49	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	1.9	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.40	ug/l	
75-01-4	Vinyl chloride <sup>b</sup>	ND	1.0	0.52	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
1330-20-7	Xylene (total)	ND	1.0	0.59	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	110%		80-120%
17060-07-0	1,2-Dichloroethane-D4	108%		80-120%
2037-26-5	Toluene-D8	94%		80-120%
460-00-4	4-Bromofluorobenzene	102%		82-114%

- (a) Associated CCV outside of control limits high, sample was ND.  
 (b) Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

**Report of Analysis**

Page 1 of 1

Client Sample ID:	SVE-5	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-3	Date Received:	03/29/23
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Iron, Ferrous <sup>a</sup>	< 0.20	0.20	mg/l	1	04/02/23 13:49	MK	SM3500FE B-11
Nitrogen, Nitrate <sup>b</sup>	7.5	0.31	mg/l	1	03/30/23 15:31	MM	EPA353.2/SM4500NO2B
Nitrogen, Nitrate + Nitrite	7.5	0.30	mg/l	3	03/30/23 15:31	MM	EPA 353.2/LACHAT
Nitrogen, Nitrite	< 0.010	0.010	mg/l	1	03/30/23 08:36	MP	SM4500NO2 B-11
Sulfate	96.8	2.0	mg/l	1	03/30/23 21:07	SS	EPA 300/SW846 9056A
Total Organic Carbon	1.1	1.0	mg/l	1	03/30/23 16:15	MB	SM5310 B-11/14

(a) Field analysis required. Received out of hold time and analyzed by request.

(b) Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

Client Sample ID:	SVE-5	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-3F	Date Received:	03/29/23
Matrix:	AQ - Groundwater Filtered	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Dissolved Organic Carbon	< 1.0	1.0	mg/l	1	03/31/23 21:28	MB	SM5310 B-11

---

RL = Reporting Limit

SGS North America Inc.

## Report of Analysis

Page 1 of 2

Client Sample ID:	FB-1	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-4	Date Received:	03/29/23
Matrix:	AQ - Field Blank Water	Percent Solids:	n/a
Method:	SW846 8260D		
Project:	388 Bridge Street, Brooklyn, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1R03031.D	1	04/04/23 17:28	NW	n/a	n/a	V1R100
Run #2							

Purge Volume
Run #1    5.0 ml
Run #2

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone <sup>a</sup>	ND	10	3.1	ug/l	
71-43-2	Benzene	ND	0.50	0.43	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.45	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>b</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	2.7	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.46	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
110-82-7	Cyclohexane	ND	5.0	0.78	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	0.53	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.56	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.51	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.54	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
123-91-1	1,4-Dioxane	ND	130	19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
76-13-1	Freon 113	ND	5.0	0.58	ug/l	

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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## Report of Analysis

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Client Sample ID:	FB-1	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-4	Date Received:	03/29/23
Matrix:	AQ - Field Blank Water	Percent Solids:	n/a
Method:	SW846 8260D		
Project:	388 Bridge Street, Brooklyn, NY		

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
591-78-6	2-Hexanone	ND	5.0	2.0	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
79-20-9	Methyl Acetate	ND	5.0	0.80	ug/l	
108-87-2	Methylcyclohexane	ND	5.0	0.60	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.9	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
100-42-5	Styrene	ND	1.0	0.49	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.56	ug/l	
108-88-3	Toluene	ND	1.0	0.49	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.40	ug/l	
75-01-4	Vinyl chloride <sup>b</sup>	ND	1.0	0.52	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
1330-20-7	Xylene (total)	ND	1.0	0.59	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	109%		80-120%
17060-07-0	1,2-Dichloroethane-D4	106%		80-120%
2037-26-5	Toluene-D8	95%		80-120%
460-00-4	4-Bromofluorobenzene	103%		82-114%

- (a) Associated CCV outside of control limits high, sample was ND.  
 (b) Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.

ND = Not detected      MDL = Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

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**Report of Analysis**

Page 1 of 1

Client Sample ID:	FB-1	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-4	Date Received:	03/29/23
Matrix:	AQ - Field Blank Water	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Iron, Ferrous <sup>a</sup>	< 0.20	0.20	mg/l	1	04/02/23 13:49	MK	SM3500FE B-11
Nitrogen, Nitrate <sup>b</sup>	< 0.11	0.11	mg/l	1	03/30/23 14:57	MM	EPA353.2/SM4500NO2B
Nitrogen, Nitrate + Nitrite	< 0.10	0.10	mg/l	1	03/30/23 14:57	MM	EPA 353.2/LACHAT
Nitrogen, Nitrite	< 0.010	0.010	mg/l	1	03/30/23 08:36	MP	SM4500NO2 B-11
Sulfate	< 2.0	2.0	mg/l	1	03/30/23 23:30	SS	EPA 300/SW846 9056A
Total Organic Carbon	< 1.0	1.0	mg/l	1	03/30/23 14:35	MB	SM5310 B-11/14

(a) Field analysis required. Received out of hold time and analyzed by request.

(b) Calculated as: (Nitrogen, Nitrate + Nitrite) - (Nitrogen, Nitrite)

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

Client Sample ID:	FB-1	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-4F	Date Received:	03/29/23
Matrix:	AQ - Field Blank Filtered	Percent Solids:	n/a
Project:	388 Bridge Street, Brooklyn, NY		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Dissolved Organic Carbon	< 1.0	1.0	mg/l	1	03/31/23 20:35	MB	SM5310 B-11

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RL = Reporting Limit

SGS North America Inc.

## Report of Analysis

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Client Sample ID:	TRIP BLANK	Date Sampled:	03/28/23
Lab Sample ID:	JD62888-5	Date Received:	03/29/23
Matrix:	AQ - Trip Blank Water	Percent Solids:	n/a
Method:	SW846 8260D		
Project:	388 Bridge Street, Brooklyn, NY		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	1R03032.D	1	04/04/23 17:53	NW	n/a	n/a	V1R100
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

## VOA TCL List (SOM0 1.1)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone <sup>a</sup>	ND	10	3.1	ug/l	
71-43-2	Benzene	ND	0.50	0.43	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.45	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>b</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	2.7	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.46	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
110-82-7	Cyclohexane	ND	5.0	0.78	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	0.53	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.56	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.51	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.54	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
123-91-1	1,4-Dioxane	ND	130	19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
76-13-1	Freon 113	ND	5.0	0.58	ug/l	

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

**Report of Analysis**

Page 2 of 2

<b>Client Sample ID:</b>	TRIP BLANK	<b>Date Sampled:</b>	03/28/23
<b>Lab Sample ID:</b>	JD62888-5	<b>Date Received:</b>	03/29/23
<b>Matrix:</b>	AQ - Trip Blank Water	<b>Percent Solids:</b>	n/a
<b>Method:</b>	SW846 8260D		
<b>Project:</b>	388 Bridge Street, Brooklyn, NY		

**VOA TCL List (SOM0 1.1)**

CAS No.	Compound	Result	RL	MDL	Units	Q
591-78-6	2-Hexanone	ND	5.0	2.0	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
79-20-9	Methyl Acetate	ND	5.0	0.80	ug/l	
108-87-2	Methylcyclohexane	ND	5.0	0.60	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.9	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
100-42-5	Styrene	ND	1.0	0.49	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.56	ug/l	
108-88-3	Toluene	ND	1.0	0.49	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.40	ug/l	
75-01-4	Vinyl chloride <sup>b</sup>	ND	1.0	0.52	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
1330-20-7	Xylene (total)	ND	1.0	0.59	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	110%		80-120%
17060-07-0	1,2-Dichloroethane-D4	109%		80-120%
2037-26-5	Toluene-D8	95%		80-120%
460-00-4	4-Bromofluorobenzene	105%		82-114%

- (a) Associated CCV outside of control limits high, sample was ND.  
 (b) Associated CCV outside of control limits low. A sensitivity check was analyzed to demonstrate system suitability to detect affected analyte. Sample was ND.

ND = Not detected      MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

## Misc. Forms

### Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody
- Sample Tracking Chronicle
- Internal Chain of Custody

GS

GW  
FB  
TB

## CHAIN OF CUSTODY

SGS North America Inc. - Dayton  
 2235 Route 130, Dayton, NJ 08810  
 TEL. 732-329-0200  
[www.sgs.com/ehsusa](http://www.sgs.com/ehsusa)

TM-03223-72

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Client / Reporting Information		Project Information										FED-EX Tracking #		Bottle Order Control #						
Company Name: <b>Fleming Lee Shue</b>		Project Name: <b>388 Bridge St. LBN Sampling</b>										SGS Quote #		SGS Job # <b>JD 62888</b>						
Street Address: <b>159 W. 24th St. FL9</b>		Street: <b>388 Bridge St.</b>		Billing Information (if different from Report to)																
City <b>NYC</b>	State <b>NY</b>	Zip <b>10001</b>	City <b>Bronxton</b>	State <b>NY</b>	Company Name															
Project Contact <b>J. Moore</b>		E-mail <b>jmoore@flemingleeshue.com</b>		Project # <b>10100</b>		Street Address														
Phone # <b>212-675-3525</b>				Client Purchase Order #		City		State		Zip										
Sampler(s) Name(s) <b>B. Hesse</b>		Phone #		Project Manager <b>J. McChesky</b>		Attention:														
SGS Sample #	Field ID / Point of Collection	Collection				Number of Bottles								pH Check (Lab Use Only)						
		MEOH/HD1 Vial #	Date	Time	Sampled by	Grab (G) Comp (C)	Source Chlorinated (Y/N)	Matrix	# of bottles	HCl	NaOH	H <sub>2</sub> SO <sub>4</sub>	None	Dil Water	MEOH	ENCORE				
1	SVE-1	3/29/23	11:55	BH	G	N	GW	9	4		1	4			X	X	X	X	X	C29
2	SVE-4		11:50	BH	G	N	GW	9	4		1	4			X	X	X	X	X	19/25
3	SVE-5		10:05	BH	G	N	GW	9	4		1	4			X	X	X	X	X	1739
4	FBI		11:30	BH	G	N	FB	8	3		1	4			X	X	X	X	X	
5	Trip Blank	3/29/23	8:44	-	-	-	-	2	2						X					
Turn Around Time (Business Days)												Deliverable				Comments / Special Instructions				
Approved By (SGS PM) / Date:				Deliverable								Comments / Special Instructions								
<input checked="" type="checkbox"/> 10 Business Days <input type="checkbox"/> 5 Business Days <input type="checkbox"/> 3 Business Days* <input type="checkbox"/> 2 Business Days* <input type="checkbox"/> 1 Business Day* <input type="checkbox"/> Other _____				<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input checked="" type="checkbox"/> NJ Reduced (Level 3) <input type="checkbox"/> Full Tier I (Level 4) <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NJ DKP								<input type="checkbox"/> NYASP Category A <input checked="" type="checkbox"/> NYASP Category B <input type="checkbox"/> MA MCP Criteria _____ <input type="checkbox"/> CT RCP Criteria _____ <input type="checkbox"/> State Forms _____ <input type="checkbox"/> EDD Format _____								
All data available via SGS Engage				Commercial "A" = Results only; Commercial "B" = Results + QC Summary Commercial "C" = Results + QC Summary + Partial Raw data								<b>SGS COURIER</b> Initial Assessment <b>A/C3B</b> Label Verification _____ <a href="http://www.sgs.com/en/terms-and-conditions">http://www.sgs.com/en/terms-and-conditions</a>								

Sample Custody must be documented below each time samples change possession, including courier delivery.

Relinquished by:	Date / Time:	Received By:	Relinquished By:	Date / Time:	Received By:
1 <i>John C</i>	3/29/23 13:10	1	2		
3 <i>John Luberman</i>	Date / Time:	Received By: 3	4 <i>John Luberman</i>	Date / Time: 3/29/23 13:10	Received By: 4
5 <i>John Luberman</i>	Date / Time:	Received By: 5	Custody Seal #		Therm ID: On Ice

EHSA-QAC-0023-05 Rev.Date:8/5/22

JD62888: Chain of Custody

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SGS

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JD62888

## SGS Sample Receipt Summary

Job Number: JD62888 Client: FLEMING-LEE SHUE, INC. Project: 388 BRIDGE STREET, BROOKLYN, NY  
 Date / Time Received: 3/29/2023 6:00:00 PM Delivery Method: IMPULSE Airbill #'s:

Cooler Temps (Raw Measured) °C: Cooler 1: (2.3);

Cooler Temps (Corrected) °C: Cooler 1: (2.1);

<u>Cooler Security</u>	<u>Y or N</u>	<u>Y or N</u>	<u>Sample Integrity - Documentation</u>	<u>Y or N</u>
1. Custody Seals Present:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	3. COC Present:	<input checked="" type="checkbox"/>
2. Custody Seals Intact:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	4. Smpl Dates/Time OK	<input checked="" type="checkbox"/>
<u>Cooler Temperature</u>	<u>Y or N</u>		<u>Sample Integrity - Condition</u>	<u>Y or N</u>
1. Temp criteria achieved:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	1. Sample recv'd within HT:	<input checked="" type="checkbox"/>
2. Cooler temp verification:	IR Gun		2. All containers accounted for:	<input checked="" type="checkbox"/>
3. Cooler media:	Ice (Bag)		3. Condition of sample:	Intact
4. No. Coolers:	1			
<u>Quality Control Preservation</u>	<u>Y or N</u>	<u>N/A</u>	<u>Sample Integrity - Instructions</u>	<u>Y or N</u> <u>N/A</u>
1. Trip Blank present / cooler:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	1. Analysis requested is clear:	<input checked="" type="checkbox"/>
2. Trip Blank listed on COC:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	2. Bottles received for unspecified tests	<input type="checkbox"/> <input checked="" type="checkbox"/>
3. Samples preserved properly:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	3. Sufficient volume recv'd for analysis:	<input checked="" type="checkbox"/>
4. VOCs headspace free:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	4. Compositing instructions clear:	<input type="checkbox"/> <input type="checkbox"/> <input checked="" type="checkbox"/>
			5. Filtering instructions clear:	<input type="checkbox"/> <input type="checkbox"/> <input checked="" type="checkbox"/>

Test Strip Lot #s:	pH 1-12: 231619	pH 12+: 203117A	Other: (Specify)
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Comments

SM089-03  
Rev. Date 12/7/17

JD62888: Chain of Custody

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5.1

SGS North America Inc.

## Internal Sample Tracking Chronicle

Fleming-Lee Shue, Inc.

Job No: JD62888

388 Bridge Street, Brooklyn, NY  
Project No: 10188

Sample Number	Method	Analyzed	By	Prepped	By	Test Codes
JD62888-1	Collected: 28-MAR-23 11:55 By: BH	Received: 29-MAR-23 By: JP	SVE-1			
JD62888-1	SM4500NO2 B-11	30-MAR-23 08:36	MP			NO2
JD62888-1	EPA353.2/SM4500NO2	30-MAR-23 15:29	MM			NO30
JD62888-1	EPA 353.2/LACHAT	30-MAR-23 15:29	MM	30-MAR-23	MM	NO32
JD62888-1	SM5310 B-11/14	30-MAR-23 15:53	MB	30-MAR-23	MB	TOC
JD62888-1	EPA 300/SW846 9056A	30-MAR-23 22:25	SS	30-MAR-23	SS	SO4
JD62888-1	SM3500FE B-11	02-APR-23 13:49	MK			FE2
JD62888-1	SW846 8260D	04-APR-23 15:45	NW			V8260TCL11
JD62888-2	Collected: 28-MAR-23 11:50 By: BH	Received: 29-MAR-23 By: JP	SVE-4			
JD62888-2	SM4500NO2 B-11	30-MAR-23 08:36	MP			NO2
JD62888-2	EPA353.2/SM4500NO2	30-MAR-23 15:30	MM			NO30
JD62888-2	EPA 353.2/LACHAT	30-MAR-23 15:30	MM	30-MAR-23	MM	NO32
JD62888-2	SM5310 B-11/14	30-MAR-23 16:04	MB	30-MAR-23	MB	TOC
JD62888-2	EPA 300/SW846 9056A	30-MAR-23 22:39	SS	30-MAR-23	SS	SO4
JD62888-2	SM3500FE B-11	02-APR-23 13:49	MK			FE2
JD62888-2	SW846 8260D	04-APR-23 18:45	NW			V8260TCL11
JD62888-3	Collected: 28-MAR-23 10:05 By: BH	Received: 29-MAR-23 By: JP	SVE-5			
JD62888-3	SM4500NO2 B-11	30-MAR-23 08:36	MP			NO2
JD62888-3	EPA353.2/SM4500NO2	30-MAR-23 15:31	MM			NO30
JD62888-3	EPA 353.2/LACHAT	30-MAR-23 15:31	MM	30-MAR-23	MM	NO32
JD62888-3	SM5310 B-11/14	30-MAR-23 16:15	MB	30-MAR-23	MB	TOC
JD62888-3	EPA 300/SW846 9056A	30-MAR-23 21:07	SS	30-MAR-23	SS	SO4
JD62888-3	SM3500FE B-11	02-APR-23 13:49	MK			FE2
JD62888-3	SW846 8260D	04-APR-23 19:10	NW			V8260TCL11
JD62888-4	Collected: 28-MAR-23 11:30 By: BH	Received: 29-MAR-23 By: JP	FB-1			
JD62888-4	SM4500NO2 B-11	30-MAR-23 08:36	MP			NO2
JD62888-4	SM5310 B-11/14	30-MAR-23 14:35	MB	30-MAR-23	MB	TOC
JD62888-4	EPA353.2/SM4500NO2	30-MAR-23 14:57	MM			NO30
JD62888-4	EPA 353.2/LACHAT	30-MAR-23 14:57	MM	30-MAR-23	MM	NO32

SGS North America Inc.

## Internal Sample Tracking Chronicle

Fleming-Lee Shue, Inc.

Job No: JD62888

388 Bridge Street, Brooklyn, NY  
Project No: 10188

Sample Number	Method	Analyzed	By	Prepped	By	Test Codes
JD62888-4	EPA 300/SW846 9056A30-MAR-23 23:30	SS		30-MAR-23 SS		SO4
JD62888-4	SM3500FE B-11	02-APR-23 13:49	MK			FE2
JD62888-4	SW846 8260D	04-APR-23 17:28	NW			V8260TCL11
JD62888-5 Collected: 28-MAR-23 11:55 By:				Received: 29-MAR-23 By: JP		
TRIP BLANK						
JD62888-5	SW846 8260D	04-APR-23 17:53	NW			V8260TCL11
JD62888-1F Collected: 28-MAR-23 11:55 By: BH				Received: 29-MAR-23 By: JP		
SVE-1						
JD62888-1F	SM5310 B-11	31-MAR-23 21:17	MB	30-MAR-23 MB	DOC	
JD62888-2F Collected: 28-MAR-23 11:50 By: BH				Received: 29-MAR-23 By: JP		
SVE-4						
JD62888-2F	SM5310 B-11	31-MAR-23 19:59	MB	30-MAR-23 MB	DOC	
JD62888-3F Collected: 28-MAR-23 10:05 By: BH				Received: 29-MAR-23 By: JP		
SVE-5						
JD62888-3F	SM5310 B-11	31-MAR-23 21:28	MB	30-MAR-23 MB	DOC	
JD62888-4F Collected: 28-MAR-23 11:30 By: BH				Received: 29-MAR-23 By: JP		
FB-1						
JD62888-4F	SM5310 B-11	31-MAR-23 20:35	MB	30-MAR-23 MB	DOC	

# SGS Internal Chain of Custody

Page 1 of 4

Job Number: JD62888  
 Account: FLSNYNY Fleming-Lee Shue, Inc.  
 Project: 388 Bridge Street, Brooklyn, NY  
 Received: 03/29/23

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JD62888-1.1	Secured Storage	Dave Hunkel	03/31/23 10:48	Retrieve from Storage
JD62888-1.1	Dave Hunkel	Secured Staging Area	03/31/23 10:48	Return to Storage
JD62888-1.1	Secured Staging Area	Marcin Kotowski	03/31/23 17:24	Retrieve from Storage
JD62888-1.1	Marcin Kotowski	Secured Storage	04/02/23 16:07	Return to Storage
JD62888-1.2	Secured Storage	Todd Shoemaker	03/30/23 10:24	Retrieve from Storage
JD62888-1.2	Todd Shoemaker	Secured Staging Area	03/30/23 10:24	Return to Storage
JD62888-1.2	Secured Staging Area	Mira Michael	03/30/23 10:35	Retrieve from Storage
JD62888-1.2	Mira Michael	Secured Storage	03/30/23 16:56	Return to Storage
JD62888-1.3	Secured Storage	Dave Hunkel	03/30/23 13:16	Retrieve from Storage
JD62888-1.3	Dave Hunkel	Secured Staging Area	03/30/23 13:16	Return to Storage
JD62888-1.3	Secured Staging Area	Sarah Sarantopoulos	03/30/23 15:04	Retrieve from Storage
JD62888-1.3	Sarah Sarantopoulos	Secured Storage	04/03/23 13:24	Return to Storage
JD62888-1.4	Secured Storage	Dave Hunkel	03/30/23 07:05	Retrieve from Storage
JD62888-1.4	Dave Hunkel	Secured Staging Area	03/30/23 07:06	Return to Storage
JD62888-1.4	Secured Staging Area	Mahendra Patel	03/30/23 08:01	Retrieve from Storage
JD62888-1.4	Mahendra Patel	Secured Storage	03/30/23 17:27	Return to Storage
JD62888-1.5	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-1.5	Secured Storage	Dave Hunkel	03/30/23 10:20	Retrieve from Storage
JD62888-1.5	Dave Hunkel	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-1.7	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-1.8	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-1.8	Nicholas Weigand	GCMS1K	04/04/23 15:07	Load on Instrument
Analyst chain of custody update error.				
JD62888-1.8	GCMS1K	Nicholas Weigand	04/05/23 16:13	Unload from Instrument
JD62888-1.8	Nicholas Weigand	Secured Storage	04/05/23 16:13	Return to Storage
JD62888-1.9	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-1F.6	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-1F.6	Secured Storage	Dave Hunkel	03/30/23 10:20	Retrieve from Storage
JD62888-1F.6	Dave Hunkel	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-2.1	Secured Storage	Dave Hunkel	03/31/23 10:48	Retrieve from Storage
JD62888-2.1	Dave Hunkel	Secured Staging Area	03/31/23 10:48	Return to Storage
JD62888-2.1	Secured Staging Area	Marcin Kotowski	03/31/23 17:24	Retrieve from Storage
JD62888-2.1	Marcin Kotowski	Secured Storage	04/02/23 16:07	Return to Storage
JD62888-2.2	Secured Storage	Todd Shoemaker	03/30/23 10:24	Retrieve from Storage

# SGS Internal Chain of Custody

Page 2 of 4

**Job Number:** JD62888  
**Account:** FLSNYNY Fleming-Lee Shue, Inc.  
**Project:** 388 Bridge Street, Brooklyn, NY  
**Received:** 03/29/23

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JD62888-2.2	Todd Shoemaker	Secured Staging Area	03/30/23 10:24	Return to Storage
JD62888-2.2	Secured Staging Area	Mira Michael	03/30/23 10:35	Retrieve from Storage
JD62888-2.2	Mira Michael	Secured Storage	03/30/23 16:56	Return to Storage
JD62888-2.3	Secured Storage	Dave Hunkele	03/30/23 13:16	Retrieve from Storage
JD62888-2.3	Dave Hunkele	Secured Staging Area	03/30/23 13:16	Return to Storage
JD62888-2.3	Secured Staging Area	Sarah Sarantopoulos	03/30/23 15:04	Retrieve from Storage
JD62888-2.3	Sarah Sarantopoulos	Secured Storage	04/03/23 13:24	Return to Storage
JD62888-2.4	Secured Storage	Dave Hunkele	03/30/23 07:05	Retrieve from Storage
JD62888-2.4	Dave Hunkele	Secured Staging Area	03/30/23 07:06	Return to Storage
JD62888-2.4	Secured Staging Area	Mahendra Patel	03/30/23 08:01	Retrieve from Storage
JD62888-2.4	Mahendra Patel	Secured Storage	03/30/23 17:27	Return to Storage
JD62888-2.5	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-2.5	Secured Storage	Dave Hunkele	03/30/23 10:20	Retrieve from Storage
JD62888-2.5	Dave Hunkele	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-2.7	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-2.8	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-2.8	Nicholas Weigand	GCMS1K	04/04/23 15:07	Load on Instrument
Analyst chain of custody update error.				
JD62888-2.8	GCMS1K	Nicholas Weigand	04/05/23 16:13	Unload from Instrument
JD62888-2.8	Nicholas Weigand	Secured Storage	04/05/23 16:13	Return to Storage
JD62888-2.9	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-2F.6	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-2F.6	Secured Storage	Dave Hunkele	03/30/23 10:20	Retrieve from Storage
JD62888-2F.6	Dave Hunkele	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-3.1	Secured Storage	Dave Hunkele	03/31/23 10:48	Retrieve from Storage
JD62888-3.1	Dave Hunkele	Secured Staging Area	03/31/23 10:48	Return to Storage
JD62888-3.1	Secured Staging Area	Marcin Kotowski	03/31/23 17:24	Retrieve from Storage
JD62888-3.1	Marcin Kotowski	Secured Storage	04/02/23 16:07	Return to Storage
JD62888-3.2	Secured Storage	Todd Shoemaker	03/30/23 10:24	Retrieve from Storage
JD62888-3.2	Todd Shoemaker	Secured Staging Area	03/30/23 10:24	Return to Storage
JD62888-3.2	Secured Staging Area	Mira Michael	03/30/23 10:35	Retrieve from Storage
JD62888-3.2	Mira Michael	Secured Storage	03/30/23 16:56	Return to Storage
JD62888-3.3	Secured Storage	Dave Hunkele	03/30/23 13:16	Retrieve from Storage
JD62888-3.3	Dave Hunkele	Secured Staging Area	03/30/23 13:16	Return to Storage

# SGS Internal Chain of Custody

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Job Number: JD62888  
 Account: FLSNYNY Fleming-Lee Shue, Inc.  
 Project: 388 Bridge Street, Brooklyn, NY  
 Received: 03/29/23

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JD62888-3.3	Secured Staging Area	Sarah Sarantopoulos	03/30/23 15:04	Retrieve from Storage
JD62888-3.3	Sarah Sarantopoulos	Secured Storage	04/03/23 13:24	Return to Storage
JD62888-3.4	Secured Storage	Dave Hunkel	03/30/23 07:05	Retrieve from Storage
JD62888-3.4	Dave Hunkel	Secured Staging Area	03/30/23 07:06	Return to Storage
JD62888-3.4	Secured Staging Area	Mahendra Patel	03/30/23 08:01	Retrieve from Storage
JD62888-3.4	Mahendra Patel	Secured Storage	03/30/23 17:27	Return to Storage
JD62888-3.5	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-3.5	Secured Storage	Dave Hunkel	03/30/23 10:20	Retrieve from Storage
JD62888-3.5	Dave Hunkel	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-3.7	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-3.8	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-3.8	Nicholas Weigand	GCMS1K	04/04/23 15:07	Load on Instrument
Analyst chain of custody update error.				
JD62888-3.8	GCMS1K	Nicholas Weigand	04/05/23 16:13	Unload from Instrument
JD62888-3.8	Nicholas Weigand	Secured Storage	04/05/23 16:13	Return to Storage
JD62888-3.9	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-3F.6	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-3F.6	Secured Storage	Dave Hunkel	03/30/23 10:20	Retrieve from Storage
JD62888-3F.6	Dave Hunkel	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-4.1	Secured Storage	Dave Hunkel	03/31/23 10:48	Retrieve from Storage
JD62888-4.1	Dave Hunkel	Secured Staging Area	03/31/23 10:48	Return to Storage
JD62888-4.1	Secured Staging Area	Marcin Kotowski	03/31/23 17:24	Retrieve from Storage
JD62888-4.1	Marcin Kotowski	Secured Storage	04/02/23 16:07	Return to Storage
JD62888-4.2	Secured Storage	Todd Shoemaker	03/30/23 10:24	Retrieve from Storage
JD62888-4.2	Todd Shoemaker	Secured Staging Area	03/30/23 10:24	Return to Storage
JD62888-4.2	Secured Staging Area	Mira Michael	03/30/23 10:35	Retrieve from Storage
JD62888-4.2	Mira Michael	Secured Storage	03/30/23 16:56	Return to Storage
JD62888-4.3	Secured Storage	Dave Hunkel	03/30/23 13:16	Retrieve from Storage
JD62888-4.3	Dave Hunkel	Secured Staging Area	03/30/23 13:16	Return to Storage
JD62888-4.3	Secured Staging Area	Sarah Sarantopoulos	03/30/23 15:04	Retrieve from Storage
JD62888-4.3	Sarah Sarantopoulos	Secured Storage	04/03/23 13:24	Return to Storage
JD62888-4.4	Secured Storage	Dave Hunkel	03/30/23 07:05	Retrieve from Storage
JD62888-4.4	Dave Hunkel	Secured Staging Area	03/30/23 07:06	Return to Storage
JD62888-4.4	Secured Staging Area	Mahendra Patel	03/30/23 08:01	Retrieve from Storage

# SGS Internal Chain of Custody

Page 4 of 4

Job Number: JD62888  
Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY  
Received: 03/29/23

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JD62888-4.4	Mahendra Patel	Secured Storage	03/30/23 17:27	Return to Storage
JD62888-4.5	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-4.5	Secured Storage	Dave Hunkele	03/30/23 10:20	Retrieve from Storage
JD62888-4.5	Dave Hunkele	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-4.7	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-4.7	Nicholas Weigand	GCMS1K	04/04/23 15:07	Load on Instrument
Analyst chain of custody update error.				
JD62888-4.7	GCMS1K	Nicholas Weigand	04/05/23 16:13	Unload from Instrument
JD62888-4.7	Nicholas Weigand	Secured Storage	04/05/23 16:13	Return to Storage
JD62888-4.8	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-4F.6	Joshua Reitan	Secured Storage	03/29/23 21:30	Return to Storage
JD62888-4F.6	Secured Storage	Dave Hunkele	03/30/23 10:20	Retrieve from Storage
JD62888-4F.6	Dave Hunkele	Secured Staging Area	03/30/23 10:20	Return to Storage
JD62888-5.1	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage
JD62888-5.1	Nicholas Weigand	GCMS1K	04/04/23 15:07	Load on Instrument
Analyst chain of custody update error.				
JD62888-5.1	GCMS1K	Nicholas Weigand	04/05/23 16:13	Unload from Instrument
JD62888-5.1	Nicholas Weigand	Secured Storage	04/05/23 16:13	Return to Storage
JD62888-5.2	Joshua Reitan	Secured Storage	03/29/23 20:08	Return to Storage

**MS Volatiles****QC Data Summaries**

**Includes the following where applicable:**

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries
- Instrument Performance Checks (BFB)
- Internal Standard Area Summaries
- Surrogate Recovery Summaries
- Initial and Continuing Calibration Summaries
- Run Sequence Reports



**Method Blank Summary**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V1R100-MB	1R03017.D	1	04/04/23	NW	n/a	n/a	V1R100

The QC reported here applies to the following samples:

Method: SW846 8260D

JD62888-1, JD62888-2, JD62888-3, JD62888-4, JD62888-5

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	3.1	ug/l	
71-43-2	Benzene	ND	0.50	0.43	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.45	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	2.7	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.46	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
110-82-7	Cyclohexane	ND	5.0	0.78	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	0.53	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.56	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.51	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.54	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
123-91-1	1,4-Dioxane	ND	130	19	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
76-13-1	Freon 113	ND	5.0	0.58	ug/l	
591-78-6	2-Hexanone	ND	5.0	2.0	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
79-20-9	Methyl Acetate	ND	5.0	0.80	ug/l	
108-87-2	Methylcyclohexane	ND	5.0	0.60	ug/l	

## Method Blank Summary

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Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V1R100-MB	1R03017.D	1	04/04/23	NW	n/a	n/a	V1R100

The QC reported here applies to the following samples:

Method: SW846 8260D

JD62888-1, JD62888-2, JD62888-3, JD62888-4, JD62888-5

CAS No.	Compound	Result	RL	MDL	Units	Q
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
108-10-1	4-Methyl-2-pentanone(MIBK)	ND	5.0	1.9	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
100-42-5	Styrene	ND	1.0	0.49	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.56	ug/l	
108-88-3	Toluene	ND	1.0	0.49	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.40	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.52	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
1330-20-7	Xylene (total)	ND	1.0	0.59	ug/l	

CAS No. Surrogate Recoveries Limits

1868-53-7	Dibromofluoromethane	109%	80-120%
17060-07-0	1,2-Dichloroethane-D4	107%	80-120%
2037-26-5	Toluene-D8	95%	80-120%
460-00-4	4-Bromofluorobenzene	103%	82-114%

CAS No.	Tentatively Identified Compounds	R.T.	Est. Conc.	Units	Q
	system artifact	1.45	11	ug/l	J
	Total TIC, Volatile		0	ug/l	

**Blank Spike Summary**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V1R100-BS	1R03015.D	1	04/04/23	NW	n/a	n/a	V1R100

The QC reported here applies to the following samples:

Method: SW846 8260D

JD62888-1, JD62888-2, JD62888-3, JD62888-4, JD62888-5

CAS No.	Compound	Spike ug/l	BSP ug/l	BSP %	Limits
67-64-1	Acetone	200	240	120	51-151
71-43-2	Benzene	50	40.0	80	80-115
74-97-5	Bromochloromethane	50	47.3	95	78-124
75-27-4	Bromodichloromethane	50	45.1	90	77-128
75-25-2	Bromoform	50	48.2	96	67-141
74-83-9	Bromomethane	50	26.5	53	53-152
78-93-3	2-Butanone (MEK)	200	209	105	61-150
75-15-0	Carbon disulfide	50	39.5	79	59-140
56-23-5	Carbon tetrachloride	50	48.5	97	75-127
108-90-7	Chlorobenzene	50	43.8	88	80-115
75-00-3	Chloroethane	50	36.8	74	54-147
67-66-3	Chloroform	50	45.6	91	75-116
74-87-3	Chloromethane	50	38.1	76	46-144
110-82-7	Cyclohexane	50	49.6	99	62-127
96-12-8	1,2-Dibromo-3-chloropropane	50	51.0	102	64-134
124-48-1	Dibromochloromethane	50	48.8	98	76-132
106-93-4	1,2-Dibromoethane	50	48.5	97	75-130
95-50-1	1,2-Dichlorobenzene	50	47.3	95	74-125
541-73-1	1,3-Dichlorobenzene	50	43.9	88	73-124
106-46-7	1,4-Dichlorobenzene	50	43.1	86	80-114
75-71-8	Dichlorodifluoromethane	50	45.1	90	42-152
75-34-3	1,1-Dichloroethane	50	43.5	87	72-124
107-06-2	1,2-Dichloroethane	50	43.9	88	73-117
75-35-4	1,1-Dichloroethene	50	43.2	86	70-124
156-59-2	cis-1,2-Dichloroethene	50	45.5	91	71-119
156-60-5	trans-1,2-Dichloroethene	50	44.8	90	71-123
78-87-5	1,2-Dichloropropane	50	39.2	78	75-120
10061-01-5	cis-1,3-Dichloropropene	50	44.0	88	77-124
10061-02-6	trans-1,3-Dichloropropene	50	46.1	92	75-132
123-91-1	1,4-Dioxane	1250	1170	94	57-147
100-41-4	Ethylbenzene	50	41.8	84	77-124
76-13-1	Freon 113	50	41.6	83	58-149
591-78-6	2-Hexanone	200	174	87	58-136
98-82-8	Isopropylbenzene	50	46.0	92	60-136
79-20-9	Methyl Acetate	50	47.9	96	57-139
108-87-2	Methylcyclohexane	50	39.0	78	73-130

\* = Outside of Control Limits.

**Blank Spike Summary**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V1R100-BS	1R03015.D	1	04/04/23	NW	n/a	n/a	V1R100

The QC reported here applies to the following samples:

Method: SW846 8260D

JD62888-1, JD62888-2, JD62888-3, JD62888-4, JD62888-5

CAS No.	Compound	Spike ug/l	BSP ug/l	BSP %	Limits
1634-04-4	Methyl Tert Butyl Ether	50	47.3	95	72-127
108-10-1	4-Methyl-2-pentanone(MIBK)	200	168	84	63-135
75-09-2	Methylene chloride	50	46.5	93	69-122
100-42-5	Styrene	50	44.4	89	78-126
79-34-5	1,1,2,2-Tetrachloroethane	50	43.0	86	66-125
127-18-4	Tetrachloroethene	50	40.7	81	73-119
108-88-3	Toluene	50	41.1	82	76-126
87-61-6	1,2,3-Trichlorobenzene	50	46.1	92	62-132
120-82-1	1,2,4-Trichlorobenzene	50	47.2	94	67-132
71-55-6	1,1,1-Trichloroethane	50	45.2	90	77-136
79-00-5	1,1,2-Trichloroethane	50	42.6	85	75-123
79-01-6	Trichloroethene	50	43.2	86	80-118
75-69-4	Trichlorofluoromethane	50	51.6	103	56-154
75-01-4	Vinyl chloride	50	31.9	64	52-146
	m,p-Xylene	100	86.3	86	77-125
95-47-6	o-Xylene	50	42.1	84	76-126
1330-20-7	Xylene (total)	150	128	85	77-125

CAS No.	Surrogate Recoveries	BSP	Limits
1868-53-7	Dibromofluoromethane	106%	76-120%
17060-07-0	1,2-Dichloroethane-D4	103%	64-135%
2037-26-5	Toluene-D8	93%	76-117%
460-00-4	4-Bromofluorobenzene	98%	72-122%

\* = Outside of Control Limits.

**Matrix Spike/Matrix Spike Duplicate Summary**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JD62953-1MS	1R03028.D	4	04/04/23	NW	n/a	n/a	V1R100
JD62953-1MSD	1R03029.D	4	04/04/23	NW	n/a	n/a	V1R100
JD62953-1 <sup>a</sup>	1R03033.D	4	04/04/23	NW	n/a	n/a	V1R100

The QC reported here applies to the following samples:

Method: SW846 8260D

JD62888-1, JD62888-2, JD62888-3, JD62888-4, JD62888-5

CAS No.	Compound	JD62953-1		Spike	MS	MS	Spike	MSD	MSD	RPD	Limits Rec/RPD
		ug/l	Q	ug/l	ug/l	%	ug/l	ug/l	%		
67-64-1	Acetone	ND		800	717	90	800	677	85	6	22-134/19
71-43-2	Benzene	19.1		200	184	82	200	180	80	2	49-137/12
74-97-5	Bromochloromethane	ND		200	207	104	200	203	102	2	78-122/12
75-27-4	Bromodichloromethane	ND		200	187	94	200	182	91	3	76-121/12
75-25-2	Bromoform	ND		200	206	103	200	214	107	4	70-133/13
74-83-9	Bromomethane	ND		200	72.1	36	200	85.7	43	17	27-164/38
78-93-3	2-Butanone (MEK)	ND		800	762	95	800	733	92	4	52-137/17
75-15-0	Carbon disulfide	ND		200	180	90	200	176	88	2	54-136/16
56-23-5	Carbon tetrachloride	ND		200	210	105	200	211	106	0	70-132/13
108-90-7	Chlorobenzene	ND		200	187	94	200	185	93	1	68-123/12
75-00-3	Chloroethane	ND		200	152	76	200	149	75	2	48-152/17
67-66-3	Chloroform	ND		200	196	98	200	190	95	3	68-120/13
74-87-3	Chloromethane	ND		200	150	75	200	141	71	6	35-156/18
110-82-7	Cyclohexane	8.0	J	200	195	94	200	187	90	4	53-146/14
96-12-8	1,2-Dibromo-3-chloropropane	ND		200	217	109	200	219	110	1	63-134/16
124-48-1	Dibromochloromethane	ND		200	208	104	200	208	104	0	75-122/12
106-93-4	1,2-Dibromoethane	ND		200	207	104	200	204	102	1	63-134/12
95-50-1	1,2-Dichlorobenzene	ND		200	207	104	200	201	101	3	74-119/12
541-73-1	1,3-Dichlorobenzene	ND		200	194	97	200	189	95	3	75-117/12
106-46-7	1,4-Dichlorobenzene	ND		200	184	92	200	181	91	2	72-117/12
75-71-8	Dichlorodifluoromethane	ND		200	210	105	200	220	110	5	34-163/16
75-34-3	1,1-Dichloroethane	ND		200	189	95	200	183	92	3	68-129/13
107-06-2	1,2-Dichloroethane	ND		200	186	93	200	181	91	3	66-120/13
75-35-4	1,1-Dichloroethene	ND		200	192	96	200	190	95	1	59-133/15
156-59-2	cis-1,2-Dichloroethene	ND		200	196	98	200	190	95	3	52-140/12
156-60-5	trans-1,2-Dichloroethene	ND		200	195	98	200	193	97	1	70-125/13
78-87-5	1,2-Dichloropropane	ND		200	165	83	200	161	81	2	73-124/12
10061-01-5	cis-1,3-Dichloropropene	ND		200	182	91	200	179	90	2	75-125/13
10061-02-6	trans-1,3-Dichloropropene	ND		200	197	99	200	190	95	4	75-122/12
123-91-1	1,4-Dioxane	ND		5000	5360	107	5000	4490	90	18	57-145/40
100-41-4	Ethylbenzene	202		200	348	73	200	342	70	2	37-144/12
76-13-1	Freon 113	ND		200	190	95	200	192	96	1	61-142/14
591-78-6	2-Hexanone	ND		800	734	92	800	715	89	3	56-132/16
98-82-8	Isopropylbenzene	12.9		200	211	99	200	207	97	2	71-126/13
79-20-9	Methyl Acetate	ND		200	205	103	200	200	100	2	51-139/18
108-87-2	Methylcyclohexane	4.8	J	200	171	83	200	172	84	1	59-137/16

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

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Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JD62953-1MS	1R03028.D	4	04/04/23	NW	n/a	n/a	V1R100
JD62953-1MSD	1R03029.D	4	04/04/23	NW	n/a	n/a	V1R100
JD62953-1 <sup>a</sup>	1R03033.D	4	04/04/23	NW	n/a	n/a	V1R100

The QC reported here applies to the following samples:

Method: SW846 8260D

JD62888-1, JD62888-2, JD62888-3, JD62888-4, JD62888-5

CAS No.	Compound	JD62953-1		Spike	MS	MS	Spike	MSD	MSD	RPD	Limits Rec/RPD
		ug/l	Q	ug/l	ug/l	%	ug/l	ug/l	%		
1634-04-4	Methyl Tert Butyl Ether	ND		200	200	100	200	199	100	1	66-124/12
108-10-1	4-Methyl-2-pentanone(MIBK)	ND		800	701	88	800	680	85	3	65-135/14
75-09-2	Methylene chloride	ND		200	202	101	200	198	99	2	66-125/14
100-42-5	Styrene	ND		200	204	102	200	204	102	0	71-133/12
79-34-5	1,1,2,2-Tetrachloroethane	ND		200	181	91	200	172	86	5	68-127/14
127-18-4	Tetrachloroethene	ND		200	176	88	200	181	91	3	58-132/13
108-88-3	Toluene	1030	E	200	1030	0* <sup>b</sup>	200	1020	-5* <sup>b</sup>	1	46-139/12
87-61-6	1,2,3-Trichlorobenzene	ND		200	205	103	200	198	99	3	57-136/17
120-82-1	1,2,4-Trichlorobenzene	ND		200	206	103	200	205	103	0	61-137/16
71-55-6	1,1,1-Trichloroethane	ND		200	196	98	200	190	95	3	67-132/13
79-00-5	1,1,2-Trichloroethane	ND		200	179	90	200	175	88	2	75-120/12
79-01-6	Trichloroethene	ND		200	184	92	200	182	91	1	56-136/12
75-69-4	Trichlorofluoromethane	ND		200	211	106	200	220	110	4	61-145/16
75-01-4	Vinyl chloride	ND		200	134	67	200	125	63	7	41-156/16
	m,p-Xylene	381		400	706	81	400	695	79	2	32-151/12
95-47-6	o-Xylene	315		200	451	68	200	443	64	2	50-139/12
1330-20-7	Xylene (total)	696		600	1160	77	600	1140	74	2	38-147/12

CAS No. Surrogate Recoveries MS MSD JD62953-1 Limits

1868-53-7	Dibromofluoromethane	106%	105%	106%	80-120%
17060-07-0	1,2-Dichloroethane-D4	99%	99%	104%	80-120%
2037-26-5	Toluene-D8	94%	92%	93%	80-120%
460-00-4	4-Bromofluorobenzene	100%	99%	101%	82-114%

(a) Dilution required due to high concentration of target compound.

(b) Outside control limits due to high level in sample relative to spike amount.

\* = Outside of Control Limits.

6.3.1  
6

**Instrument Performance Check (BFB)**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample:	V1R91-BFB	Injection Date:	03/28/23
Lab File ID:	1R02813.D	Injection Time:	21:55
Instrument ID:	GCMS1R		

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	15.0 - 40.0% of mass 95	53808	22.7	Pass
75	30.0 - 60.0% of mass 95	112563	47.5	Pass
95	Base peak, 100% relative abundance	237077	100.0	Pass
96	5.0 - 9.0% of mass 95	16559	6.98	Pass
173	Less than 2.0% of mass 174	1620	0.68	(0.67) <sup>a</sup> Pass
174	50.0 - 150.0% of mass 95	241536	101.9	Pass
175	5.0 - 9.0% of mass 174	17164	7.24	(7.11) <sup>a</sup> Pass
176	95.0 - 101.0% of mass 174	233536	98.5	(96.7) <sup>a</sup> Pass
177	5.0 - 9.0% of mass 176	14108	5.95	(6.04) <sup>b</sup> Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V1R91-IC0091	1R02815.D	03/28/23	22:46	00:51	Initial cal 0.2
V1R91-IC0091	1R02817.D	03/28/23	23:19	01:24	Initial cal 0.5
V1R91-IC0091	1R02819.D	03/28/23	23:52	01:57	Initial cal 1
V1R91-IC0091	1R02821.D	03/29/23	00:25	02:30	Initial cal 2
V1R91-IC0091	1R02823.D	03/29/23	00:58	03:03	Initial cal 4
V1R91-IC0091	1R02825.D	03/29/23	01:31	03:36	Initial cal 8
V1R91-IC0091	1R02827.D	03/29/23	02:04	04:09	Initial cal 20
V1R91-ICC0091	1R02829.D	03/29/23	02:37	04:42	Initial cal 50
V1R91-IC0091	1R02831.D	03/29/23	03:10	05:15	Initial cal 100
V1R91-IC0091	1R02833.D	03/29/23	03:42	05:47	Initial cal 200
V1R91-ICV0091	1R02839.D	03/29/23	05:20	07:25	Initial cal verification 50
V1R91-ICV0091	1R02841.D	03/29/23	05:53	07:58	Initial cal verification 50

**Instrument Performance Check (BFB)**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample:	V1R91-BFB2	Injection Date:	03/29/23
Lab File ID:	1R02846.D	Injection Time:	16:59
Instrument ID:	GCMS1R		

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	15.0 - 40.0% of mass 95	51715	24.3	Pass
75	30.0 - 60.0% of mass 95	103344	48.6	Pass
95	Base peak, 100% relative abundance	212736	100.0	Pass
96	5.0 - 9.0% of mass 95	14480	6.81	Pass
173	Less than 2.0% of mass 174	1103	0.52	(0.53) <sup>a</sup> Pass
174	50.0 - 150.0% of mass 95	206571	97.1	Pass
175	5.0 - 9.0% of mass 174	15660	7.36	(7.58) <sup>a</sup> Pass
176	95.0 - 101.0% of mass 174	203776	95.8	(98.6) <sup>a</sup> Pass
177	5.0 - 9.0% of mass 176	13543	6.37	(6.65) <sup>b</sup> Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V1R91-ICV0091	1R02847.D	03/29/23	17:25	00:26	Initial cal verification 50

**Instrument Performance Check (BFB)**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample:	V1R100-BFB	Injection Date:	04/04/23
Lab File ID:	1R03013.D	Injection Time:	09:38
Instrument ID:	GCMS1R		

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	15.0 - 40.0% of mass 95	62083	19.1	Pass
75	30.0 - 60.0% of mass 95	143333	44.1	Pass
95	Base peak, 100% relative abundance	325376	100.0	Pass
96	5.0 - 9.0% of mass 95	21747	6.68	Pass
173	Less than 2.0% of mass 174	0	0.00	(0.00) <sup>a</sup> Pass
174	50.0 - 150.0% of mass 95	293013	90.1	Pass
175	5.0 - 9.0% of mass 174	21152	6.50	(7.22) <sup>a</sup> Pass
176	95.0 - 101.0% of mass 174	283755	87.2	(96.8) <sup>a</sup> Pass
177	5.0 - 9.0% of mass 176	19014	5.84	(6.70) <sup>b</sup> Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V1R100-CC91	1R03013.D	04/04/23	09:38	00:00	Continuing cal 20
V1R100-BS	1R03015.D	04/04/23	10:34	00:56	Blank Spike
V1R100-MB	1R03017.D	04/04/23	11:25	01:47	Method Blank
ZZZZZZ	1R03018R.D	04/04/23	11:54	02:16	(unrelated sample)
ZZZZZZ	1R03018.D	04/04/23	11:54	02:16	(unrelated sample)
ZZZZZZ	1R03019R.D	04/04/23	12:20	02:42	(unrelated sample)
ZZZZZZ	1R03019.D	04/04/23	12:20	02:42	(unrelated sample)
ZZZZZZ	1R03020.D	04/04/23	12:46	03:08	(unrelated sample)
ZZZZZZ	1R03021.D	04/04/23	13:11	03:33	(unrelated sample)
ZZZZZZ	1R03021R.D	04/04/23	13:11	03:33	(unrelated sample)
ZZZZZZ	1R03022.D	04/04/23	13:37	03:59	(unrelated sample)
ZZZZZZ	1R03023.D	04/04/23	14:03	04:25	(unrelated sample)
ZZZZZZ	1R03024.D	04/04/23	14:28	04:50	(unrelated sample)
ZZZZZZ	1R03025.D	04/04/23	14:54	05:16	(unrelated sample)
ZZZZZZ	1R03026.D	04/04/23	15:19	05:41	(unrelated sample)
JD62888-1	1R03027.D	04/04/23	15:45	06:07	SVE-1
JD62953-1MS	1R03028.D	04/04/23	16:11	06:33	Matrix Spike
JD62953-1MSD	1R03029.D	04/04/23	16:36	06:58	Matrix Spike Duplicate
JD62888-4	1R03031.D	04/04/23	17:28	07:50	FB-1
JD62888-5	1R03032.D	04/04/23	17:53	08:15	TRIP BLANK
JD62953-1	1R03033.D	04/04/23	18:19	08:41	(used for QC only; not part of job JD62888)
JD62888-2	1R03034.D	04/04/23	18:45	09:07	SVE-4
JD62888-3	1R03035.D	04/04/23	19:10	09:32	SVE-5
ZZZZZZ	1R03036.D	04/04/23	19:36	09:58	(unrelated sample)

# Instrument Performance Check (BFB)

Page 2 of 2

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Sample:	V1R100-BFB	Injection Date:	04/04/23
Lab File ID:	1R03013.D	Injection Time:	09:38
Instrument ID:	GCMS1R		

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
ZZZZZZZ	1R03037.D	04/04/23	20:01	10:23	(unrelated sample)
ZZZZZZZ	1R03038.D	04/04/23	20:27	10:49	(unrelated sample)
ZZZZZZZ	1R03039.D	04/04/23	20:53	11:15	(unrelated sample)
ZZZZZZZ	1R03040.D	04/04/23	21:18	11:40	(unrelated sample)

6.4.3  
6

# Internal Standard Area Summary

Page 1 of 2

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Check Std:	V1R100-CC91	Injection Date:	04/04/23
Lab File ID:	1R03013.D	Injection Time:	09:38
Instrument ID:	GCMS1R	Method:	SW846 8260D

	IS 1 AREA	IS 2 AREA	IS 3 AREA	IS 4 AREA	IS 5 AREA	
Check Std	649271	3.48	688798	5.02	1060871	5.64
Upper Limit <sup>a</sup>	1298542	3.98	1377596	5.52	2121742	6.14
Lower Limit <sup>b</sup>	324636	2.98	344399	4.52	530436	5.14

Lab Sample ID	IS 1 AREA	IS 2 AREA	IS 3 AREA	IS 4 AREA	IS 5 AREA	
V1R100-BS	614235	3.49	730438	5.02	1084886	5.65
V1R100-MB	612313	3.48	592616	5.02	922312	5.65
ZZZZZZ	605234	3.49	579738	5.02	907666	5.65
ZZZZZZ	605234	3.49	579738	5.02	907666	5.65
ZZZZZZ	600910	3.49	628676	5.02	968829	5.65
ZZZZZZ	600910	3.49	628676	5.02	968829	5.65
ZZZZZZ	624089	3.48	620072	5.02	961381	5.65
ZZZZZZ	595693	3.49	578679	5.03	915793	5.65
ZZZZZZ	595693	3.49	578679	5.03	915793	5.65
ZZZZZZ	614750	3.48	619030	5.02	953587	5.65
ZZZZZZ	598892	3.48	572375	5.02	901955	5.65
ZZZZZZ	572882	3.48	574792	5.03	904035	5.65
ZZZZZZ	578240	3.49	565295	5.03	883663	5.66
ZZZZZZ	521373	3.48	554367	5.02	870868	5.65
JD62888-1	590128	3.49	553552	5.03	865233	5.65
JD62953-1MS	507065	3.49	639724	5.02	974157	5.65
JD62953-1MSD	584603	3.49	697532	5.03	1038223	5.66
JD62888-4	631950	3.48	583702	5.03	908888	5.66
JD62888-5	592479	3.49	568084	5.03	880717	5.66
JD62953-1	534542	3.49	558173	5.03	856573	5.66
JD62888-2	575874	3.49	548691	5.03	858272	5.66
JD62888-3	525305	3.49	542116	5.03	848679	5.66
ZZZZZZ	586826	3.49	546029	5.03	843157	5.66
ZZZZZZ	529461	3.49	537394	5.03	832517	5.66
ZZZZZZ	583637	3.49	536623	5.03	839343	5.66
ZZZZZZ	531119	3.49	522955	5.03	814601	5.66
ZZZZZZ	543596	3.50	531147	5.03	824443	5.66

IS 1 = Tert Butyl Alcohol-D9

IS 2 = Pentafluorobenzene

IS 3 = 1,4-Difluorobenzene

IS 4 = Chlorobenzene-D5

IS 5 = 1,4-Dichlorobenzene-d4

(a) Upper Limit = + 100% of check standard area; Retention time + 0.5 minutes.

6.5.1  
6

## Internal Standard Area Summary

Page 2 of 2

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Check Std:	V1R100-CC91	Injection Date:	04/04/23
Lab File ID:	1R03013.D	Injection Time:	09:38
Instrument ID:	GCMS1R	Method:	SW846 8260D

Lab Sample ID	IS 1 AREA	IS 2 RT	IS 3 AREA	IS 4 RT	IS 5 AREA	RT
------------------	--------------	------------	--------------	------------	--------------	----

(b) Lower Limit = -50% of check standard area; Retention time -0.5 minutes.

6.5.1  
6

# Surrogate Recovery Summary

Page 1 of 1

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Method: SW846 8260D

Matrix: AQ

Samples and QC shown here apply to the above method

Lab Sample ID	Lab File ID	S1	S2	S3	S4
JD62888-1	1R03027.D	109	108	94	102
JD62888-2	1R03034.D	111	108	94	102
JD62888-3	1R03035.D	110	108	94	102
JD62888-4	1R03031.D	109	106	95	103
JD62888-5	1R03032.D	110	109	95	105
JD62953-1MS	1R03028.D	106	99	94	100
JD62953-1MSD	1R03029.D	105	99	92	99
V1R100-BS	1R03015.D	106	103	93	98
V1R100-MB	1R03017.D	109	107	95	103

Surrogate Compounds	Recovery Limits
------------------------	--------------------

S1 = Dibromofluoromethane	80-120%
S2 = 1,2-Dichloroethane-D4	80-120%
S3 = Toluene-D8	80-120%
S4 = 4-Bromofluorobenzene	82-114%

6.6.1  
6

**Initial Calibration Summary****Page 1 of 5**Job Number: **JD62888**Sample: **V1R91-ICC0091**Account: **FLSNYNY Fleming-Lee Shue, Inc.**Lab FileID: **1R02829.D**Project: **388 Bridge Street, Brooklyn, NY**

## Response Factor Report GCMSR

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)

Title : SW846 8260D, Rx1624Sil MS 60m x 0.25mm x 1.4um

Last Update : Thu Mar 30 09:13:25 2023

Response via : Initial Calibration

## Calibration Files

1	=1R02819.D	4	=1R02823.D	100	=1R02831.D	50	=1R02829.D
20	=1R02827.D	200	=1R02833.D	8	=1R02825.D	0.5	=1R02817.D
2	=1R02821.D	0.2	=1R02815.D	=	=		

## Compound

	1	4	100	50	20	200	8	0.5	2	0.2	Avg	%RSD
--	---	---	-----	----	----	-----	---	-----	---	-----	-----	------

-----

1) I	tert butyl alcohol-d9	-----ISTD-----									
2)	ethanol	0.087 0.098 0.080 0.084 0.090 0.058 0.100 0.096 0.119									
3)	tertiary butyl alcohol	0.090 18.29									
4)	1,4-dioxane	1.399 1.291 1.238 1.146 1.119 1.115 1.244 1.338 1.399									
5)	I	tertafluorobenzene -----ISTD-----									
6)	chlorodifluoromethane	0.437 0.352 0.355 0.378 0.266 0.407									
7)	dichlorodifluoromethane	0.476 0.381 17.86									
8)	chloromethane	0.355 0.380 0.342 0.328 0.335 0.246 0.378 0.408 0.413 0.341									
9)	vinyl chloride	0.353 13.58									
10)	1,3-butadiene	0.853 0.649 0.480 0.514 0.560 0.349 0.627									
11)	bromomethane	0.740 0.597 26.35									
12)	chloroethane	0.853 0.649 0.480 0.514 0.560 0.349 0.627									
13)	trichlorofluoromethane	0.740 0.597 26.35									
14)	ethyl ether	0.359 0.290 14.59									
15)	acrolein	0.359 0.290 14.59									
16)	freon 113	0.359 0.290 14.59									
17)	1,1-dichloroethene	0.359 0.290 14.59									
18)	acetone	0.359 0.290 14.59									
19)	acetonitrile	0.359 0.290 14.59									
20)	iodomethane	0.359 0.290 14.59									
		----- Quadratic regression ----- Coefficient = 0.9995									
		Response Ratio = 0.00583 + 0.58085 *A + -0.05787 *A^2									

6.7.1  
6

**Initial Calibration Summary**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NYSample: V1R91-ICC0091  
Lab FileID: 1R02829.D

21)	carbon disulfide	0.956 0.822 0.786 0.729 0.751 0.744 0.832 0.856 0.997	0.830	11.30
22)	methylene chloride	0.286 0.333 0.303 0.294 0.288 0.286 0.331 0.281 0.373	0.308	10.03
23)	methyl acetate	0.102 0.093 0.090 0.080 0.085 0.088 0.090	0.109	0.092 10.04
24)	methyl tert butyl ether	0.992 0.861 0.956 0.828 0.786 0.888 0.831 0.883 1.037 0.742	0.880	10.46
25)	trans-1,2-dichloroethene	0.346 0.292 0.332 0.292 0.281 0.304 0.304 0.311 0.363	0.314	8.66
26)	hexane *This compound fails initial calibration criteria.*	0.294 0.276 0.245 0.240 0.233 0.228 0.266 0.306 0.307	0.266	11.80
27)	di-isopropyl ether	1.332 1.287 1.458 1.200 1.119 1.146 1.236 1.243 1.387 1.083	1.249	9.58
28)	2-butanone	0.082 0.071 0.093 0.072 0.078 0.069 0.080 0.086 0.092	0.080	11.07
29)	1,1-dichloroethane	0.596 0.577 0.693 0.551 0.530 0.539 0.579 0.596 0.660 0.571	0.589	8.77
30)	chloroprene	0.583 0.471 0.621 0.490 0.431 0.516 0.476 0.457 0.556 0.477	0.508	11.92
31)	acrylonitrile	0.212 0.239 0.239 0.231 0.226 0.209 0.242 0.200 0.254	0.228	7.80
32)	vinyl acetate	0.092 0.074 0.086 0.078 0.067 0.079 0.079	0.089	0.080 10.50
33)	ethyl tert-butyl ether	1.028 0.965 1.110 0.969 0.898 1.009 0.973 0.806 1.138 0.970	0.987	9.68
34)	ethyl acetate	0.126 0.121 0.128 0.107 0.106 0.097 0.116	0.100	0.113 10.62
35)	2,2-dichloropropane	0.456 0.426 0.573 0.383 0.367 0.441 0.394 0.395 0.447	0.431	14.25
36)	cis-1,2-dichloroethene	0.377 0.345 0.431 0.321 0.310 0.350 0.327 0.311 0.362 0.374	0.351	10.63
37)	propionitrile	0.113 0.099 0.120 0.098 0.097 0.090 0.098 0.093 0.121 0.106	0.104	10.60
38)	methyl acrylate	0.090 0.073 0.107 0.084 0.076 0.079 0.084	0.085	0.085 12.19
39)	methacrylonitrile	0.246 0.216 0.264 0.219 0.193 0.202 0.209	0.244	0.224 10.95
40)	bromochloromethane	0.167 0.169 0.211 0.159 0.140 0.174 0.149 0.132 0.181 0.182	0.166	13.79
41)	tetrahydrofuran	0.327 0.288 0.292 0.260 0.252 0.213 0.278	0.325	0.279 13.61
42)	chloroform	0.606 0.523 0.638 0.532 0.489 0.549 0.526 0.684 0.642	0.577	11.72
43)	tert-Butyl Formate	0.298 0.312 0.365 0.331 0.278 0.333 0.301 0.268 0.340	0.314	9.98
44)	dibromofluoromethane (s)	0.428 0.442 0.453 0.439 0.434 0.431 0.432 0.422 0.433 0.433	0.435	1.94
45)	1,1,1-trichloroethane	0.541 0.466 0.510 0.455 0.440 0.503 0.464 0.508 0.584	0.497	9.27
46)	cyclohexane	0.467 0.434 0.487 0.450 0.428 0.435 0.427 0.517 0.507	0.461	7.56
47)	isobutyl alcohol	0.044 0.039 0.040 0.042 0.035 0.043	0.051	0.042 11.58
48)	1,1-dichloropropene	0.439 0.408 0.439 0.400 0.388 0.393 0.423 0.395 0.464	0.416	6.30
49)	carbon tetrachloride	0.424 0.397 0.451 0.405 0.379 0.483 0.398 0.362 0.430	0.414	8.95
50)	tert-amyl alcohol			

6.7.1  
6

**Initial Calibration Summary**

**Job Number:** JD62888  
**Account:** FLSNYNY Fleming-Lee Shue, Inc.  
**Project:** 388 Bridge Street, Brooklyn, NY

**Sample:** V1R91-ICC0091  
**Lab FileID:** 1R02829.D

51)	isopropyl acetate	0.050 0.052 0.053 0.047 0.050 0.046 0.050	0.051	0.050	4.87
		0.117 0.099 0.125 0.104 0.100 0.104 0.108	0.133	0.111	11.16
52)	I 1,4-difluorobenzene	-----ISTD-----			
53)	1,2-dichloroethane-d4 (s)	0.276 0.281 0.333 0.293 0.287 0.342 0.277 0.271 0.278 0.267 0.291	0.291	8.88	
54)	n-butyl alcohol	0.017 0.017 0.022 0.019 0.018 0.020 0.018	0.020	0.019	8.96
55)	2,2,4-trimethylpentane	0.772 0.663 0.758 0.666 0.614 0.681 0.674 0.628 0.754 0.726 0.694	0.694	8.01	
56)	benzene	0.977 0.971 1.168 0.943 0.861 1.003 0.898 0.908 1.118 1.010 0.986	0.986	9.74	
57)	tert-amyl methyl ether	0.700 0.703 0.844 0.692 0.626 0.755 0.671 0.640 0.759 0.619 0.701	0.701	9.94	
58)	heptane	0.172 0.163 0.156 0.149 0.136 0.147 0.154 0.139 0.194	0.157	11.41	
59)	1,2-dichloroethane	0.304 0.293 0.345 0.304 0.261 0.340 0.288 0.340 0.337 0.330 0.314	0.314	8.99	
60)	ethyl acrylate	0.533 0.485 0.599 0.502 0.438 0.493 0.483 0.495 0.542	0.508	8.97	
61)	trichloroethene	0.265 0.228 0.302 0.238 0.223 0.279 0.235 0.190 0.270 0.260 0.249	0.249	13.06	
62)	2-chloroethyl vinyl ether	0.205 0.211 0.283 0.234 0.193 0.231 0.203 0.177 0.231 0.169 0.214	0.214	15.40	
63)	methyl methacrylate	0.088 0.076 0.111 0.077 0.070 0.091 0.077	0.082	0.084	15.18
64)	methylcyclohexane	0.418 0.411 0.459 0.372 0.333 0.405 0.375 0.319 0.408 0.440 0.394	0.394	11.25	
65)	1,2-dichloropropane	0.302 0.270 0.330 0.263 0.235 0.258 0.256 0.233 0.296	0.271	11.87	
66)	dibromomethane	0.172 0.168 0.212 0.160 0.138 0.166 0.148 0.186 0.159	0.168	12.87	
67)	bromodichloromethane	0.324 0.281 0.324 0.298 0.255 0.327 0.271 0.301 0.311	0.299	8.60	
68)	2-nitropropane	0.153 0.159 0.138 0.119 0.137 0.140	0.141	9.75	
69)	epichlorohydrin	0.045 0.051 0.061 0.052 0.045 0.052 0.050	0.051	0.051	9.79
70)	cis-1,3-dichloropropene	0.375 0.356 0.464 0.385 0.340 0.414 0.363 0.311 0.402 0.438 0.385	0.385	12.00	
71)	4-methyl-2-pentanone	0.171 0.174 0.236 0.207 0.172 0.185 0.177 0.160 0.208 0.184 0.187	0.187	12.26	
72)	isoamyl alcohol	0.015 0.021 0.020 0.016 0.018 0.016	0.016	0.017	13.93
73)	I chlorobenzene-d5	-----ISTD-----			
74)	toluene-d8 (s)	1.328 1.339 1.484 1.533 1.394 1.347 1.329 1.335 1.354 1.364 1.381	1.381	5.15	
75)	toluene	0.667 0.620 0.703 0.675 0.604 0.649 0.607 0.492 0.698 0.761 0.648	0.648	11.31	
76)	ethyl methacrylate	0.339 0.350 0.413 0.361 0.345 0.344 0.342 0.302 0.385	0.353	8.79	
77)	trans-1,3-dichloropropene	0.342 0.340 0.410 0.341 0.329 0.366 0.337 0.326 0.382 0.293 0.347	0.347	9.35	
78)	1,1,2-trichloroethane	0.203 0.202 0.198 0.189 0.181 0.188 0.205 0.186 0.257 0.215 0.203	0.203	10.81	
79)	tetrachloroethene	0.295 0.248 0.220 0.215 0.223 0.256 0.251 0.274 0.299 0.270 0.255	0.255	11.77	
80)	2-hexanone				

6.7.1  
6

**Initial Calibration Summary**

Job Number: JD62888

Sample: V1R91-ICC0091

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Lab FileID: 1R02829.D

	0.199	0.203	0.210	0.207	0.198	0.184	0.200	0.173	0.237	0.234	0.205	9.63
81)	1,3-dichloropropane											
	0.369	0.352	0.356	0.346	0.338	0.348	0.369	0.297	0.428	0.415	0.362	10.39
82)	butyl acetate											
	0.266	0.272	0.278	0.258	0.250	0.234	0.261	0.286	0.306		0.268	7.79
83)	dibromochloromethane											
	0.274	0.237	0.282	0.248	0.231	0.303	0.231	0.205	0.286		0.255	12.67
84)	1,2-dibromoethane											
	0.267	0.237	0.233	0.232	0.227	0.254	0.229	0.220	0.282		0.242	8.60
85)	n-butyl ether											
	1.213	1.134	1.209	1.147	1.103	1.053	1.116	0.996	1.277	1.230	1.148	7.51
86)	chlorobenzene											
	0.725	0.673	0.795	0.679	0.653	0.756	0.667	0.650	0.840	0.723	0.716	8.97
87)	1,1,1,2-tetrachloroethane											
	0.243	0.230	0.313	0.272	0.230	0.302	0.229	0.201	0.257		0.253	14.53
88)	ethylbenzene											
	1.169	1.124	1.344	1.254	1.103	1.205	1.109	1.120	1.231	1.244	1.190	6.66
89)	m,p-xylene											
	0.477	0.425	0.529	0.515	0.425	0.513	0.433	0.411	0.483	0.409	0.462	10.05
90)	o-xylene											
	0.887	0.814	1.083	1.058	0.836	0.990	0.817	0.700	0.929	1.068	0.918	14.12
91)	styrene											
	0.691	0.689	0.927	0.981	0.715	0.863	0.671	0.621	0.779		0.771	16.27
92)	butyl acrylate											
	0.691	0.601	0.665	0.608	0.582	0.574	0.605	0.477	0.696	0.519	0.602	11.67
93)	n-amyl acetate											
	0.213	0.204	0.244	0.281	0.224	0.203	0.209		0.266		0.230	12.90
94)	isopropylbenzene											
	1.040	0.982	1.312	1.295	0.990	1.174	0.988	0.847	1.070		1.078	14.35
95)	bromoform											
	0.176	0.164	0.266	0.236	0.173	0.245	0.185	0.167	0.195	0.220	0.203	17.99
96)	cis-1,4-dichloro-2-butene											
	0.137	0.118	0.194	0.191	0.129	0.162	0.128		0.148		0.151	19.17
97)	I 1,4-dichlorobenzene-d -----ISTD-----											
98)	4-bromofluorobenzene (s)											
	0.834	0.855	0.801	0.686	0.769	0.778	0.853	0.867	0.868	0.871	0.818	7.30
99)	1,1,2,2-tetrachloroethane											
	0.897	0.720	0.672	0.548	0.700	0.652	0.863	0.746	0.966		0.752	17.65
100)	trans-1,4-dichloro-2-butene											
	0.256	0.258	0.238	0.216	0.247	0.228	0.267	0.261	0.302		0.253	9.84
101)	1,2,3-trichloropropane											
	0.237	0.244	0.221	0.169	0.203	0.211	0.226		0.271		0.223	13.56
102)	bromobenzene											
	0.761	0.574	0.599	0.475	0.505	0.619	0.611	0.618	0.686	0.462	0.591	15.69
103)	n-propylbenzene											
	2.824	2.609	2.531	2.119	2.340	2.397	2.640	2.579	2.963		2.556	9.89
104)	2-chlorotoluene											
	0.601	0.552	0.543	0.422	0.484	0.544	0.547	0.404	0.684	0.505	0.529	15.47
105)	4-chlorotoluene											
	1.551	1.525	1.563	1.432	1.389	1.531	1.509	1.637	1.752	1.813	1.570	8.38
106)	1,3,5-trimethylbenzene											
	1.842	1.764	1.850	1.621	1.594	1.806	1.834	1.540	1.980	1.764	1.760	7.71
107)	tert-butylbenzene											
	1.691	1.525	1.593	1.258	1.360	1.612	1.549	1.429	1.811		1.536	11.00
108)	1,2,4-trimethylbenzene											
	1.972	1.736	1.957	1.802	1.633	1.898	1.851	1.688	1.927	1.442	1.791	9.38
109)	sec-butylbenzene											
	2.277	2.066	2.222	1.972	1.917	2.143	2.161	2.051	2.397	1.724	2.093	9.21
110)	p-isopropyltoluene											

**Initial Calibration Summary**

**Job Number:** JD62888  
**Account:** FLSNYNY Fleming-Lee Shue, Inc.  
**Project:** 388 Bridge Street, Brooklyn, NY

**Sample:** V1R91-ICC0091  
**Lab FileID:** 1R02829.D

111)	1,2,3-trimethylbenzene	1.939 1.822 2.105 1.947 1.654 2.144 1.866 1.619 2.108 1.970 1.917 9.47
		1.994 1.762 1.980 1.918 1.699 1.997 1.846 1.804 2.176 1.909 7.68
112)	1,3-dichlorobenzene	1.105 1.055 1.232 1.130 0.952 1.331 1.073 1.091 1.301 1.100 1.137 10.30
113)	1,4-dichlorobenzene	1.110 1.124 1.194 1.117 0.997 1.290 1.050 1.302 1.193 1.238 1.161 8.59
114)	1,2-dichlorobenzene	1.097 1.005 1.116 1.040 0.930 1.174 1.078 0.875 1.189 0.931 1.043 10.24
115)	benzyl chloride	1.470 1.354 1.397 1.346 1.295 1.439 1.343 1.318 1.649 1.401 7.75
116)	n-butylbenzene	0.877 0.868 0.827 0.828 0.818 0.844 0.850 0.927 0.981 0.800 0.862 6.38
117)	hexachloroethane	0.143 0.168 0.246 0.170 0.179 0.204 0.166 0.182 18.27
118)	1,2-dibromo-3-chloropropane	0.204 0.199 0.195 0.151 0.184 0.230 0.227 0.204 0.239 0.204 13.25
119)	1,3,5-trichlorobenzene	0.692 0.684 0.632 0.458 0.564 0.767 0.688 0.779 0.658 16.10
120)	1,2,4-trichlorobenzene	0.577 0.557 0.521 0.406 0.478 0.643 0.555 0.692 0.554 16.18
121)	hexachlorobutadiene	0.234 0.202 0.162 0.138 0.185 0.231 0.215 0.242 0.201 18.33
122)	naphthalene	2.041 2.020 2.073 1.479 1.743 2.289 2.080 2.456 2.022 14.94
123)	1,2,3-trichlorobenzene	0.654 0.549 0.529 0.376 0.474 0.619 0.548 0.636 0.548 16.80
124)	2-methylnaphthalene	0.939 0.902 0.745 0.613 0.788 1.025 0.925 0.993 1.025 0.884 15.94

(#) = Out of Range   ### Number of calibration levels exceeded format   ###

M1R0091.M              Thu Mar 30 09:19:01 2023

6.7.1  
6

**Initial Calibration Verification**

Job Number: JD62888

Sample: V1R91-ICV0091

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Lab FileID: 1R02839.D

**Evaluate Continuing Calibration Report**

Data File : C:\MSDCHEM\1\VR0091\1R02839.D Vial: 14  
 Acq On : 29 Mar 2023 5:20 am Operator: PrashanS  
 Sample : ICV0091-50 Inst : GCMSR  
 Misc : MS67262,V1R0091,5,,,,1 Multiplr: 1.00  
 MS Integration Params: rteint.p

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)  
 Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 Last Update : Thu Mar 30 09:13:25 2023  
 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)R.T.	
1 I	tert butyl alcohol-d9	1.000	1.000	0.0	108	0.00	3.49
2	ethanol	0.090	0.085	5.6	109	0.00	2.69
3	tertiary butyl alcohol	1.254	1.119	10.8	106	0.00	3.56
4	1,4-dioxane	0.093	0.088	5.4	114	0.00	6.11
5 I	pentafluorobenzene	1.000	1.000	0.0	101	0.00	5.03
6	chlorodifluoromethane			-----NA-----			
7	dichlorodifluoromethane	0.353	0.261	26.1	81	0.00	1.64
8	chloromethane	50.000	36.870	26.3	79	0.00	1.80
9	vinyl chloride	0.576	0.420	27.1	84	0.00	1.90
10	1,3-butadiene	0.430	0.447	-4.0	114	0.00	1.93
11	bromomethane	0.322	0.221	31.4#	72	0.00	2.18
12	chloroethane	0.290	0.223	23.1	84	0.00	2.29
13	trichlorofluoromethane	0.441	0.358	18.8	88	0.00	2.54
14	ethyl ether	0.182	0.178	2.2	106	0.00	2.80
15	acrolein	0.105	0.097	7.6	100	0.00	2.94
16	freon 113	0.249	0.230	7.6	102	0.00	3.04
17	1,1-dichloroethene	0.283	0.259	8.5	103	0.00	3.03
18	acetone	0.068	0.062	8.8	116	0.00	3.07
19	acetonitrile	0.092	0.073	20.7	88	0.00	3.33
20	iodomethane	50.000	44.221	11.6	84	0.00	3.18
21	carbon disulfide	0.830	0.787	5.2	109	0.00	3.25
22	methylene chloride	0.308	0.307	0.3	106	0.00	3.50
23	methyl acetate	0.092	0.086	6.5	109	0.00	3.36
24	methyl tert butyl ether	0.880	0.872	0.9	107	0.00	3.73
25	trans-1,2-dichloroethene	0.314	0.306	2.5	107	0.00	3.74
26	hexane	0.266	0.173	35.0#	73	0.00	3.98
27	di-isopropyl ether	1.249	1.262	-1.0	107	0.00	4.14
28	2-butanone	0.080	0.081	-1.3	113	0.00	4.59
29	1,1-dichloroethane	0.589	0.577	2.0	106	0.00	4.13
30	chloroprene	0.508	0.502	1.2	104	0.00	4.19
31	acrylonitrile	0.228	0.227	0.4	100	0.00	3.70
32	vinyl acetate	0.080	0.067	16.2	87	0.00	4.12
33	ethyl tert-butyl ether	0.987	0.967	2.0	101	0.00	4.45

**Initial Calibration Verification**

Job Number: JD62888

Sample: V1R91-ICV0091

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Lab FileID: 1R02839.D

34	ethyl acetate	0.113	0.115	-1.8	109	0.00	4.62
35	2,2-dichloropropane	0.431	0.400	7.2	106	0.00	4.62
36	cis-1,2-dichloroethene	0.351	0.345	1.7	109	0.00	4.61
37	propionitrile	0.104	0.102	1.9	105	0.00	4.65
38	methyl acrylate	0.085	0.090	-5.9	109	0.00	4.67
39	methacrylonitrile	0.224	0.218	2.7	101	0.00	4.77
40	bromochloromethane	0.166	0.155	6.6	99	0.00	4.81
41	tetrahydrofuran	0.279	0.261	6.5	102	0.00	4.82
42	chloroform	0.577	0.539	6.6	103	0.00	4.89
43	tert-Butyl Formate	0.314	0.239	23.9	73	0.00	4.91
44 S	dibromofluoromethane (s)	0.435	0.453	-4.1	105	0.00	5.02
45	1,1,1-trichloroethane	0.497	0.473	4.8	106	0.00	5.04
46	cyclohexane	0.461	0.387	16.1	87	0.00	5.10
47	isobutyl alcohol	0.042	0.041	2.4	102	0.00	5.21
48	1,1-dichloropropene	0.416	0.414	0.5	105	0.00	5.16
49	carbon tetrachloride	0.414	0.414	0.0	104	0.00	5.16
50	tert-amyl alcohol	0.050	0.048	4.0	106	0.00	5.31
51	isopropyl acetate	0.111	0.108	2.7	106	0.00	5.33
52 I	1,4-difluorobenzene	1.000	1.000	0.0	105	0.00	5.65
53 S	1,2-dichloroethane-d4 (s)	0.291	0.297	-2.1	106	0.00	5.30
54	n-butyl alcohol	0.019	0.019	0.0	106	0.00	5.74
55	2,2,4-trimethylpentane	0.694	0.689	0.7	109	0.00	5.41
56	benzene	0.986	0.939	4.8	105	0.00	5.33
57	tert-amyl methyl ether	0.701	0.663	5.4	101	0.00	5.41
58	heptane	0.157	0.142	9.6	101	0.00	5.54
59	1,2-dichloroethane	0.314	0.289	8.0	100	0.00	5.36
60	ethyl acrylate	0.508	0.500	1.6	105	0.00	5.89
61	trichloroethene	0.249	0.239	4.0	106	0.00	5.85
62	2-chloroethyl vinyl ether	0.214	0.226	-5.6	102	0.00	6.47
63	methyl methacrylate	0.084	0.103	-22.6	141	0.00	6.08
64	methylcyclohexane	0.394	0.406	-3.0	115	0.00	6.05
65	1,2-dichloropropene	0.271	0.270	0.4	108	0.00	6.06
66	dibromomethane	0.168	0.165	1.8	109	0.00	6.13
67	bromodichloromethane	0.299	0.278	7.0	98	0.00	6.27
68	2-nitropropane	0.141	0.133	5.7	101	0.00	6.43
69	epichlorohydrin	0.051	0.049	3.9	100	0.00	6.51
70	cis-1,3-dichloropropene	0.385	0.370	3.9	101	0.00	6.59
71	4-methyl-2-pentanone	0.187	0.204	-9.1	104	0.00	6.69
72	isoamyl alcohol	0.017	0.020	-17.6	107	0.00	6.72
73 I	chlorobenzene-d5	1.000	1.000	0.0	100	0.00	7.82
74 S	toluene-d8 (s)	1.381	1.530	-10.8	100	0.00	6.80
75	toluene	0.648	0.678	-4.6	101	0.00	6.85
76	ethyl methacrylate	0.353	0.327	7.4	91	0.00	7.03
77	trans-1,3-dichloropropene	0.347	0.348	-0.3	103	0.00	7.01
78	1,1,2-trichloroethane	0.203	0.193	4.9	103	0.00	7.15
79	tetrachloroethene	0.255	0.305	-19.6	142	0.00	7.22
80	2-hexanone	0.205	0.209	-2.0	101	0.00	7.29
81	1,3-dichloropropane	0.362	0.354	2.2	103	0.00	7.27
82	butyl acetate	0.268	0.279	-4.1	108	0.00	7.35
83	dibromochloromethane	0.255	0.251	1.6	102	0.00	7.42
84	1,2-dibromoethane	0.242	0.238	1.7	103	0.00	7.51
85	n-butyl ether	1.148	1.216	-5.9	106	0.00	7.87
86	chlorobenzene	0.716	0.676	5.6	100	0.00	7.83
87	1,1,1,2-tetrachloroethane	0.253	0.285	-12.6	105	0.00	7.89
88	ethylbenzene	1.190	1.309	-10.0	105	0.00	7.89
89	m,p-xylene	0.462	0.527	-14.1	103	0.00	7.97
90	o-xylene	0.918	1.074	-17.0	102	0.00	8.22
91	styrene	0.771	0.856	-11.0	88	0.00	8.24

6.7.2  
6

**Initial Calibration Verification**

Job Number: JD62888

Sample: V1R91-ICV0091

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Lab FileID: 1R02839.D

92	butyl acrylate	0.602	0.634	-5.3	105	0.00	8.16
93	n-amyl acetate	0.230	0.273	-18.7	98	0.00	8.30
94	isopropylbenzene	1.078	1.303	-20.9	101	0.00	8.45
95	bromoform	0.203	0.232	-14.3	99	0.00	8.36
96	cis-1,4-dichloro-2-butene	0.151	0.176	-16.6	92	0.00	8.48
97 I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	100	0.00	9.26
98 S	4-bromofluorobenzene (s)	0.818	0.644	21.3	94	0.00	8.56
99	1,1,2,2-tetrachloroethane	0.752	0.597	20.6	108	0.00	8.63
100	trans-1,4-dichloro-2-bute	0.253	0.212	16.2	98	0.00	8.66
101	1,2,3-trichloropropane	0.223	0.177	20.6	104	0.00	8.67
102	bromobenzene	0.591	0.443	25.0	93	0.00	8.66
103	n-propylbenzene	2.556	2.096	18.0	99	0.00	8.70
104	2-chlorotoluene	0.529	0.424	19.8	100	0.00	8.77
105	4-chlorotoluene	1.570	1.429	9.0	99	0.00	8.84
106	1,3,5-trimethylbenzene	1.760	1.606	8.7	99	0.00	8.81
107	tert-butylbenzene	1.536	1.247	18.8	99	0.00	9.00
108	1,2,4-trimethylbenzene	1.791	1.805	-0.8	100	0.00	9.03
109	sec-butylbenzene	2.093	1.897	9.4	96	0.00	9.13
110	p-isopropyltoluene	1.917	1.909	0.4	98	0.00	9.22
111	1,2,3-trimethylbenzene			-----NA-----			
112	1,3-dichlorobenzene	1.137	1.103	3.0	97	0.00	9.22
113	1,4-dichlorobenzene	1.161	1.102	5.1	98	0.00	9.28
114	1,2-dichlorobenzene	1.043	0.975	6.5	93	0.00	9.49
115	benzyl chloride	1.401	1.138	18.8	84	0.00	9.34
116	n-butylbenzene	0.862	0.772	10.4	93	0.00	9.45
117	hexachloroethane	0.182	0.184	-1.1	108	0.00	9.65
118	1,2-dibromo-3-chloropropa	0.204	0.148	27.5	98	0.00	9.93
119	1,3,5-trichlorobenzene	0.658	0.493	25.1	107	0.00	10.03
120	1,2,4-trichlorobenzene	0.554	0.416	24.9	102	0.00	10.38
121	hexachlorobutadiene	0.201	0.148	26.4	107	0.00	10.44
122	naphthalene	2.022	1.496	26.0	101	0.00	10.53
123	1,2,3-trichlorobenzene	0.548	0.400	27.0	106	0.00	10.65
124	2-methylnaphthalene	0.884	0.667	24.5	108	0.00	11.13

( # ) = Out of Range  
1R02829.D M1R0091.MSPCC's out = 0 CCC's out = 0  
Thu Mar 30 09:18:42 20236.7.2  
6

**Initial Calibration Verification**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NYSample: V1R91-ICV0091  
Lab FileID: 1R02841.D**Evaluate Continuing Calibration Report**

Data File : C:\MSDCHEM\1\VR0091\1R02841.D Vial: 15  
 Acq On : 29 Mar 2023 5:53 am Operator: PrashanS  
 Sample : ICV0091-50 Inst : GCMSR  
 Misc : MS67262,V1R0091,5,,,,1 Multiplr: 1.00  
 MS Integration Params: rteint.p

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)  
 Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 Last Update : Thu Mar 30 09:07:12 2023  
 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1 I	tert butyl alcohol-d9	1.000	1.000	0.0	113	0.00	3.49
2	ethanol		-----NA-----				
3	tertiary butyl alcohol		-----NA-----				
4	1,4-dioxane		-----NA-----				
5 I	pentafluorobenzene	1.000	1.000	0.0	94	0.00	5.03
6	chlorodifluoromethane	0.381	0.432	-13.4	115	0.00	1.65
7	dichlorodifluoromethane		-----NA-----				
8	chloromethane		-----True-----	Calc.	% Drift	-----	
9	vinyl chloride		AvgRF	CCRF	% Dev	-----	
10	1,3-butadiene			-----NA-----			
11	bromomethane			-----NA-----			
12	chloroethane			-----NA-----			
13	trichlorofluoromethane			-----NA-----			
14	ethyl ether			-----NA-----			
15	acrolein			-----NA-----			
16	freon 113			-----NA-----			
17	1,1-dichloroethene			-----NA-----			
18	acetone			-----NA-----			
19	acetonitrile	0.092	0.099	-7.6	111	0.00	3.33
20	iodomethane		-----True-----	Calc.	% Drift	-----	
21	carbon disulfide		AvgRF	CCRF	% Dev	-----	
22	methylene chloride			-----NA-----			
23	methyl acetate			-----NA-----			
24	methyl tert butyl ether			-----NA-----			
25	trans-1,2-dichloroethene			-----NA-----			
26	hexane			-----NA-----			
27	di-isopropyl ether			-----NA-----			
28	2-butanone			-----NA-----			
29	1,1-dichloroethane			-----NA-----			
30	chloroprene			-----NA-----			
31	acrylonitrile	0.228	0.239	-4.8	98	0.00	3.70
32	vinyl acetate			-----NA-----			
33	ethyl tert-butyl ether			-----NA-----			

**Initial Calibration Verification**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NYSample: V1R91-ICV0091  
Lab FileID: 1R02841.D

34	ethyl acetate		-----	-NA-----					
35	2,2-dichloropropane		-----	-NA-----					
36	cis-1,2-dichloroethene		-----	-NA-----					
37	propionitrile		-----	-NA-----					
38	methyl acrylate		-----	-NA-----					
39	methacrylonitrile		-----	-NA-----					
40	bromochloromethane		-----	-NA-----					
41	tetrahydrofuran		-----	-NA-----					
42	chloroform		-----	-NA-----					
43	tert-Butyl Formate		-----	-NA-----					
44 S	dibromofluoromethane (s)	0.435	0.442	-1.6	95	0.00	5.02		
45	1,1,1-trichloroethane		-----	-NA-----					
46	cyclohexane		-----	-NA-----					
47	isobutyl alcohol		-----	-NA-----					
48	1,1-dichloropropene		-----	-NA-----					
49	carbon tetrachloride		-----	-NA-----					
50	tert-amyl alcohol		-----	-NA-----					
51	isopropyl acetate		-----	-NA-----					
52 I	1,4-difluorobenzene	1.000	1.000	0.0	98	0.00	5.66		
53 S	1,2-dichloroethane-d4 (s)	0.291	0.283	2.7	95	0.00	5.30		
54	n-butyl alcohol		-----	-NA-----					
55	2,2,4-trimethylpentane		-----	-NA-----					
56	benzene		-----	-NA-----					
57	tert-amyl methyl ether		-----	-NA-----					
58	heptane		-----	-NA-----					
59	1,2-dichloroethane		-----	-NA-----					
60	ethyl acrylate		-----	-NA-----					
61	trichloroethene		-----	-NA-----					
62	2-chloroethyl vinyl ether		-----	-NA-----					
63	methyl methacrylate		-----	-NA-----					
64	methylcyclohexane		-----	-NA-----					
65	1,2-dichloropropene		-----	-NA-----					
66	dibromomethane		-----	-NA-----					
67	bromodichloromethane		-----	-NA-----					
68	2-nitropropane		-----	-NA-----					
69	epichlorohydrin		-----	-NA-----					
70	cis-1,3-dichloropropene		-----	-NA-----					
71	4-methyl-2-pentanone		-----	-NA-----					
72	isoamyl alcohol		-----	-NA-----					
73 I	chlorobenzene-d5	1.000	1.000	0.0	88	0.00	7.82		
74 S	toluene-d8 (s)	1.381	1.334	3.4	77	0.00	6.80		
75	toluene		-----	-NA-----					
76	ethyl methacrylate		-----	-NA-----					
77	trans-1,3-dichloropropene		-----	-NA-----					
78	1,1,2-trichloroethane		-----	-NA-----					
79	tetrachloroethene	0.255	0.198	22.4	81	0.00	7.22		
80	2-hexanone		-----	-NA-----					
81	1,3-dichloropropane		-----	-NA-----					
82	butyl acetate		-----	-NA-----					
83	dibromochloromethane		-----	-NA-----					
84	1,2-dibromoethane		-----	-NA-----					
85	n-butyl ether		-----	-NA-----					
86	chlorobenzene		-----	-NA-----					
87	1,1,1,2-tetrachloroethane		-----	-NA-----					
88	ethylbenzene		-----	-NA-----					
89	m,p-xylene		-----	-NA-----					
90	o-xylene		-----	-NA-----					
91	styrene		-----	-NA-----					

6.7.3  
6

**Initial Calibration Verification**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NYSample: V1R91-ICV0091  
Lab FileID: 1R02841.D

92	butyl acrylate		-----	-NA-----					
93	n-amyl acetate		-----	-NA-----					
94	isopropylbenzene		-----	-NA-----					
95	bromoform		-----	-NA-----					
96	cis-1,4-dichloro-2-butene		-----	-NA-----					
97 I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	59	0.00	9.26		
98 S	4-bromofluorobenzene (s)	0.818	0.865	-5.7	74	0.00	8.56		
99	1,1,2,2-tetrachloroethane		-----	-NA-----					
100	trans-1,4-dichloro-2-bute		-----	-NA-----					
101	1,2,3-trichloropropane		-----	-NA-----					
102	bromobenzene		-----	-NA-----					
103	n-propylbenzene		-----	-NA-----					
104	2-chlorotoluene		-----	-NA-----					
105	4-chlorotoluene		-----	-NA-----					
106	1,3,5-trimethylbenzene		-----	-NA-----					
107	tert-butylbenzene		-----	-NA-----					
108	1,2,4-trimethylbenzene		-----	-NA-----					
109	sec-butylbenzene		-----	-NA-----					
110	p-isopropyltoluene		-----	-NA-----					
111	1,2,3-trimethylbenzene	1.909	1.967	-3.0	60	0.00	9.28		
112	1,3-dichlorobenzene		-----	-NA-----					
113	1,4-dichlorobenzene		-----	-NA-----					
114	1,2-dichlorobenzene		-----	-NA-----					
115	benzyl chloride		-----	-NA-----					
116	n-butylbenzene		-----	-NA-----					
117	hexachloroethane		-----	-NA-----					
118	1,2-dibromo-3-chloropropa		-----	-NA-----					
119	1,3,5-trichlorobenzene		-----	-NA-----					
120	1,2,4-trichlorobenzene		-----	-NA-----					
121	hexachlorobutadiene		-----	-NA-----					
122	naphthalene		-----	-NA-----					
123	1,2,3-trichlorobenzene		-----	-NA-----					
124	2-methylnaphthalene		-----	-NA-----					

(#= Out of Range  
1R02829.D M1R0091.MSPCC's out = 0 CCC's out = 0  
Thu Mar 30 09:11:41 2023

**Initial Calibration Verification**

Job Number: JD62888

Sample: V1R91-ICV0091

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Lab FileID: 1R02847.D

**Evaluate Continuing Calibration Report**

Data File : C:\MSDCHEM\1\VR0091\1R02847.D Vial: 19  
 Acq On : 29 Mar 2023 5:25 pm Operator: PrashanS  
 Sample : icv0091-50 Inst : GCMSR  
 Misc : MS67262,V1R0091,5,,,,1 Multiplr: 1.00  
 MS Integration Params: rteint.p

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)  
 Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 Last Update : Thu Mar 30 09:13:25 2023  
 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1 I	tert butyl alcohol-d9	1.000	1.000	0.0	89	-0.01	3.48
2	ethanol		-----NA-----				
3	tertiary butyl alcohol		-----NA-----				
4	1,4-dioxane		-----NA-----				
5 I	pentafluorobenzene	1.000	1.000	0.0	81	-0.01	5.02
6	chlorodifluoromethane		-----NA-----				
7	dichlorodifluoromethane	0.353	0.348	1.4	86	0.00	1.63
8	chloromethane	50.000	58.022	-16.0	95	0.00	1.80
9	vinyl chloride	0.576	0.583	-1.2	93	0.00	1.89
10	1,3-butadiene		-----NA-----				
11	bromomethane	0.322	0.263	18.3	68	0.00	2.18
12	chloroethane	0.290	0.249	14.1	75	-0.01	2.28
13	trichlorofluoromethane	0.441	0.416	5.7	82	0.00	2.53
14	ethyl ether		-----NA-----				
15	acrolein		-----NA-----				
16	freon 113		-----NA-----				
17	1,1-dichloroethene		-----NA-----				
18	acetone		-----NA-----				
19	acetonitrile		-----NA-----				
20	iodomethane		-----NA-----				
21	carbon disulfide		-----NA-----				
22	methylene chloride		-----NA-----				
23	methyl acetate		-----NA-----				
24	methyl tert butyl ether		-----NA-----				
25	trans-1,2-dichloroethene		-----NA-----				
26	hexane		-----NA-----				
27	di-isopropyl ether		-----NA-----				
28	2-butanone		-----NA-----				
29	1,1-dichloroethane		-----NA-----				
30	chloroprene		-----NA-----				
31	acrylonitrile		-----NA-----				
32	vinyl acetate		-----NA-----				
33	ethyl tert-butyl ether		-----NA-----				

**Initial Calibration Verification**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NYSample: V1R91-ICV0091  
Lab FileID: 1R02847.D

34	ethyl acetate		-----	-NA-----						
35	2,2-dichloropropane		-----	-NA-----						
36	cis-1,2-dichloroethene		-----	-NA-----						
37	propionitrile		-----	-NA-----						
38	methyl acrylate		-----	-NA-----						
39	methacrylonitrile		-----	-NA-----						
40	bromochloromethane		-----	-NA-----						
41	tetrahydrofuran		-----	-NA-----						
42	chloroform		-----	-NA-----						
43	tert-Butyl Formate		-----	-NA-----						
44 S	dibromofluoromethane (s)	0.435	0.434	0.2	80	-0.01			5.01	
45	1,1,1-trichloroethane		-----	-NA-----						
46	cyclohexane		-----	-NA-----						
47	isobutyl alcohol		-----	-NA-----						
48	1,1-dichloropropene		-----	-NA-----						
49	carbon tetrachloride		-----	-NA-----						
50	tert-amyl alcohol		-----	-NA-----						
51	isopropyl acetate		-----	-NA-----						
52 I	1,4-difluorobenzene	1.000	1.000	0.0	83	-0.01		5.64		
53 S	1,2-dichloroethane-d4 (s)	0.291	0.270	7.2	77	-0.01		5.29		
54	n-butyl alcohol		-----	-NA-----						
55	2,2,4-trimethylpentane		-----	-NA-----						
56	benzene		-----	-NA-----						
57	tert-amyl methyl ether		-----	-NA-----						
58	heptane		-----	-NA-----						
59	1,2-dichloroethane		-----	-NA-----						
60	ethyl acrylate		-----	-NA-----						
61	trichloroethene		-----	-NA-----						
62	2-chloroethyl vinyl ether		-----	-NA-----						
63	methyl methacrylate		-----	-NA-----						
64	methylcyclohexane		-----	-NA-----						
65	1,2-dichloropropene		-----	-NA-----						
66	dibromomethane		-----	-NA-----						
67	bromodichloromethane		-----	-NA-----						
68	2-nitropropane		-----	-NA-----						
69	epichlorohydrin		-----	-NA-----						
70	cis-1,3-dichloropropene		-----	-NA-----						
71	4-methyl-2-pentanone		-----	-NA-----						
72	isoamyl alcohol		-----	-NA-----						
73 I	chlorobenzene-d5	1.000	1.000	0.0	73	0.00		7.81		
74 S	toluene-d8 (s)	1.381	1.336	3.3	64	0.00		6.79		
75	toluene		-----	-NA-----						
76	ethyl methacrylate		-----	-NA-----						
77	trans-1,3-dichloropropene		-----	-NA-----						
78	1,1,2-trichloroethane		-----	-NA-----						
79	tetrachloroethene		-----	-NA-----						
80	2-hexanone		-----	-NA-----						
81	1,3-dichloropropane		-----	-NA-----						
82	butyl acetate		-----	-NA-----						
83	dibromochloromethane		-----	-NA-----						
84	1,2-dibromoethane		-----	-NA-----						
85	n-butyl ether		-----	-NA-----						
86	chlorobenzene		-----	-NA-----						
87	1,1,1,2-tetrachloroethane		-----	-NA-----						
88	ethylbenzene		-----	-NA-----						
89	m,p-xylene		-----	-NA-----						
90	o-xylene		-----	-NA-----						
91	styrene		-----	-NA-----						

6.7.4  
6

**Initial Calibration Verification**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NYSample: V1R91-ICV0091  
Lab FileID: 1R02847.D

92	butyl acrylate		-----NA-----
93	n-amyl acetate		-----NA-----
94	isopropylbenzene		-----NA-----
95	bromoform		-----NA-----
96	cis-1,4-dichloro-2-butene		-----NA-----
97 I	1,4-dichlorobenzene-d4	1.000	1.000 0.0 50 0.00 9.26
98 S	4-bromofluorobenzene (s)	0.818	0.856 -4.6 63 0.00 8.56
99	1,1,2,2-tetrachloroethane		-----NA-----
100	trans-1,4-dichloro-2-bute		-----NA-----
101	1,2,3-trichloropropane		-----NA-----
102	bromobenzene		-----NA-----
103	n-propylbenzene		-----NA-----
104	2-chlorotoluene		-----NA-----
105	4-chlorotoluene		-----NA-----
106	1,3,5-trimethylbenzene		-----NA-----
107	tert-butylbenzene		-----NA-----
108	1,2,4-trimethylbenzene		-----NA-----
109	sec-butylbenzene		-----NA-----
110	p-isopropyltoluene		-----NA-----
111	1,2,3-trimethylbenzene		-----NA-----
112	1,3-dichlorobenzene		-----NA-----
113	1,4-dichlorobenzene		-----NA-----
114	1,2-dichlorobenzene		-----NA-----
115	benzyl chloride		-----NA-----
116	n-butylbenzene		-----NA-----
117	hexachloroethane		-----NA-----
118	1,2-dibromo-3-chloropropane		-----NA-----
119	1,3,5-trichlorobenzene		-----NA-----
120	1,2,4-trichlorobenzene		-----NA-----
121	hexachlorobutadiene		-----NA-----
122	naphthalene		-----NA-----
123	1,2,3-trichlorobenzene		-----NA-----
124	2-methylnaphthalene		-----NA-----

(#= Out of Range  
1R02829.D M1R0091.MSPCC's out = 0 CCC's out = 0  
Thu Mar 30 09:23:24 2023

## Continuing Calibration Summary

Job Number: JD62888

Sample: V1R100-CC91

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Lab FileID: 1R03013.D

## Evaluate Continuing Calibration Report

Data File : C:\msdchem\1\data\Me...2023\VR0100\1R03013.D Vial: 2  
 Acq On : 4 Apr 2023 9:38 am Operator: nickw  
 Sample : cc91-20 Inst : GCMSR  
 Misc : MS67909,V1R0100,5,,,,1 Multiplr: 1.00  
 MS Integration Params: rteint.p

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)  
 Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 Last Update : Thu Mar 30 09:13:25 2023  
 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1	I tert butyl alcohol-d9	1.000	1.000	0.0	130	-0.01	3.48
2	ethanol	0.090	0.096	-6.7	139	-0.01	2.68
3	tertiary butyl alcohol	1.254	1.174	6.4	137	-0.01	3.56
4	1,4-dioxane	0.093	0.086	7.5	137	-0.01	6.11
5	I pentafluorobenzene	1.000	1.000	0.0	112	-0.01	5.02
6	chlorodifluoromethane	0.381	0.313	17.8	93	0.00	1.65
7	dichlorodifluoromethane	0.353	0.367	-4.0	123	0.00	1.63
8	chloromethane	20.000	16.338	18.3	95	0.00	1.79
9	vinyl chloride	0.576	0.409	29.0#	84	0.00	1.89
10	1,3-butadiene	0.430	0.261	39.3#	70	0.00	1.93
11	bromomethane	0.322	0.242	24.8#	92	0.00	2.18
12	chloroethane	0.290	0.233	19.7	96	0.00	2.29
13	trichlorofluoromethane	0.441	0.482	-9.3	131	0.00	2.53
14	ethyl ether	0.182	0.180	1.1	119	-0.01	2.79
15	acrolein	0.105	0.110	-4.8	125	0.00	2.93
16	freon 113	0.249	0.233	6.4	121	-0.01	3.03
17	1,1-dichloroethene	0.283	0.283	0.0	126	-0.01	3.03
18	acetone	0.068	0.087	-27.9#	146	0.00	3.06
19	acetonitrile	0.092	0.103	-12.0	129	-0.01	3.32
20	iodomethane	20.000	12.534	37.3#	68	0.00	3.17
21	carbon disulfide	0.830	0.740	10.8	110	-0.01	3.24
22	methylene chloride	0.308	0.320	-3.9	124	-0.01	3.48
23	methyl acetate	0.092	0.095	-3.3	125	0.00	3.35
24	methyl tert butyl ether	0.880	0.916	-4.1	131	-0.01	3.72
25	trans-1,2-dichloroethene	0.314	0.307	2.2	122	0.00	3.73
26	hexane	0.266	0.254	4.5	122	-0.01	3.97
27	di-isopropyl ether	1.249	1.261	-1.0	126	-0.01	4.13
28	2-butanone	0.080	0.093	-16.2	134	-0.01	4.58
29	1,1-dichloroethane	0.589	0.570	3.2	121	-0.01	4.12
30	chloroprene	0.508	0.508	0.0	132	-0.01	4.18
31	acrylonitrile	0.228	0.241	-5.7	119	-0.01	3.69
32	vinyl acetate	0.080	0.084	-5.0	141	-0.01	4.11
33	ethyl tert-butyl ether	0.987	1.065	-7.9	133	-0.01	4.44

## Continuing Calibration Summary

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NYSample: V1R100-CC91  
Lab FileID: 1R03013.D

34	ethyl acetate	0.113	0.106	6.2	113	-0.01	4.60
35	2,2-dichloropropane	0.431	0.451	-4.6	138	-0.01	4.61
36	cis-1,2-dichloroethene	0.351	0.361	-2.8	131	-0.01	4.60
37	propionitrile	0.104	0.107	-2.9	124	-0.01	4.64
38	methyl acrylate	0.085	0.088	-3.5	128	-0.01	4.65
39	methacrylonitrile	0.224	0.229	-2.2	133	-0.01	4.77
40	bromochloromethane	0.166	0.177	-6.6	141	0.00	4.81
41	tetrahydrofuran	0.279	0.250	10.4	111	-0.01	4.81
42	chloroform	0.577	0.577	0.0	132	-0.01	4.88
43	tert-Butyl Formate	0.314	0.345	-9.9	139	-0.01	4.90
44 S	dibromofluoromethane (s)	0.435	0.471	-8.3	122	0.00	5.01
45	1,1,1-trichloroethane	0.497	0.484	2.6	123	-0.01	5.02
46	cyclohexane	0.461	0.469	-1.7	123	-0.01	5.08
47	isobutyl alcohol	0.042	0.048	-14.3	127	-0.01	5.20
48	1,1-dichloropropene	0.416	0.422	-1.4	122	-0.01	5.15
49	carbon tetrachloride	0.414	0.426	-2.9	126	-0.01	5.15
50	tert-amyl alcohol	0.050	0.058	-16.0	128	-0.02	5.30
51	isopropyl acetate	0.111	0.115	-3.6	128	-0.01	5.32
52 I	1,4-difluorobenzene	1.000	1.000	0.0	121	-0.01	5.64
53 S	1,2-dichloroethane-d4 (s)	0.291	0.302	-3.8	127	-0.01	5.29
54	n-butyl alcohol	0.019	0.019	0.0	130	-0.02	5.73
55	2,2,4-trimethylpentane	0.694	0.589	15.1	116	-0.01	5.40
56	benzene	0.986	0.842	14.6	118	-0.01	5.32
57	tert-amyl methyl ether	0.701	0.660	5.8	128	-0.01	5.41
58	heptane	0.157	0.132	15.9	117	-0.01	5.54
59	1,2-dichloroethane	0.314	0.293	6.7	136	-0.01	5.35
60	ethyl acrylate	0.508	0.450	11.4	124	0.00	5.89
61	trichloroethene	0.249	0.229	8.0	125	-0.01	5.84
62	2-chloroethyl vinyl ether	0.214	0.190	11.2	119	-0.01	6.46
63	methyl methacrylate	0.084	0.080	4.8	137	-0.02	6.07
64	methylcyclohexane	0.394	0.334	15.2	121	0.00	6.04
65	1,2-dichloropropene	0.271	0.231	14.8	119	-0.01	6.05
66	dibromomethane	0.168	0.150	10.7	132	-0.01	6.13
67	bromodichloromethane	0.299	0.279	6.7	132	-0.01	6.25
68	2-nitropropane	0.141	0.126	10.6	127	0.00	6.42
69	epichlorohydrin	0.051	0.053	-3.9	141	-0.01	6.50
70	cis-1,3-dichloropropene	0.385	0.356	7.5	127	-0.01	6.59
71	4-methyl-2-pentanone	0.187	0.170	9.1	119	-0.01	6.68
72	isoamyl alcohol	0.017	0.017	0.0	126	-0.01	6.72
73 I	chlorobenzene-d5	1.000	1.000	0.0	124	0.00	7.81
74 S	toluene-d8 (s)	1.381	1.298	6.0	116	0.00	6.79
75	toluene	0.648	0.565	12.8	116	-0.01	6.84
76	ethyl methacrylate	0.353	0.342	3.1	123	-0.01	7.02
77	trans-1,3-dichloropropene	0.347	0.342	1.4	129	0.00	7.00
78	1,1,2-trichloroethane	0.203	0.186	8.4	127	0.00	7.15
79	tetrachloroethene	0.255	0.227	11.0	126	0.00	7.21
80	2-hexanone	0.205	0.193	5.9	121	0.00	7.28
81	1,3-dichloropropane	0.362	0.341	5.8	125	0.00	7.26
82	butyl acetate	0.268	0.247	7.8	122	-0.01	7.35
83	dibromochloromethane	0.255	0.252	1.2	136	0.00	7.41
84	1,2-dibromoethane	0.242	0.254	-5.0	139	0.00	7.51
85	n-butyl ether	1.148	0.979	14.7	110	0.00	7.86
86	chlorobenzene	0.716	0.659	8.0	125	0.00	7.83
87	1,1,1,2-tetrachloroethane	0.253	0.230	9.1	124	0.00	7.88
88	ethylbenzene	1.190	1.067	10.3	120	0.00	7.88
89	m,p-xylene	0.462	0.417	9.7	122	0.00	7.97
90	o-xylene	0.918	0.836	8.9	124	0.00	8.22
91	styrene	0.771	0.710	7.9	123	0.00	8.23

6.7.5  
6

## Continuing Calibration Summary

Job Number: JD62888

Sample: V1R100-CC91

Account: FLSNYNY Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Lab FileID: 1R03013.D

92	butyl acrylate	0.602	0.568	5.6	121	0.00	8.16
93	n-amyl acetate	0.230	0.203	11.7	112	0.00	8.29
94	isopropylbenzene	1.078	1.039	3.6	130	0.00	8.44
95	bromoform	0.203	0.183	9.9	131	0.00	8.36
96	cis-1,4-dichloro-2-butene	0.151	0.145	4.0	140	0.00	8.48
97 I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	116	0.00	9.26
98 S	4-bromofluorobenzene (s)	0.818	0.833	-1.8	125	0.00	8.56
99	1,1,2,2-tetrachloroethane	0.752	0.716	4.8	118	0.00	8.63
100	trans-1,4-dichloro-2-bute	0.253	0.238	5.9	112	0.00	8.65
101	1,2,3-trichloropropane	0.223	0.232	-4.0	132	0.00	8.67
102	bromobenzene	0.591	0.582	1.5	133	0.00	8.66
103	n-propylbenzene	2.556	2.446	4.3	121	0.00	8.70
104	2-chlorotoluene	0.529	0.545	-3.0	130	0.00	8.77
105	4-chlorotoluene	1.570	1.481	5.7	123	0.00	8.84
106	1,3,5-trimethylbenzene	1.760	1.683	4.4	122	0.00	8.80
107	tert-butylbenzene	1.536	1.490	3.0	127	0.00	9.00
108	1,2,4-trimethylbenzene	1.791	1.750	2.3	124	0.00	9.03
109	sec-butylbenzene	2.093	2.076	0.8	125	0.00	9.13
110	p-isopropyltoluene	1.917	1.812	5.5	127	0.00	9.21
111	1,2,3-trimethylbenzene	1.909	1.767	7.4	120	0.00	9.28
112	1,3-dichlorobenzene	1.137	1.058	6.9	129	0.00	9.22
113	1,4-dichlorobenzene	1.161	1.039	10.5	121	0.00	9.27
114	1,2-dichlorobenzene	1.043	1.040	0.3	129	0.00	9.48
115	benzyl chloride	1.401	1.509	-7.7	135	0.00	9.33
116	n-butylbenzene	0.862	0.793	8.0	112	0.00	9.45
117	hexachloroethane	0.182	0.186	-2.2	121	0.00	9.65
118	1,2-dibromo-3-chloropropa	0.204	0.228	-11.8	143	0.00	9.93
119	1,3,5-trichlorobenzene	0.658	0.647	1.7	133	0.00	10.02
120	1,2,4-trichlorobenzene	0.554	0.550	0.7	133	0.00	10.38
121	hexachlorobutadiene	0.201	0.190	5.5	119	0.00	10.44
122	naphthalene	2.022	2.068	-2.3	137	0.00	10.53
123	1,2,3-trichlorobenzene	0.548	0.542	1.1	132	0.00	10.65
124	2-methylnaphthalene	0.884	0.925	-4.6	136	0.00	11.13

( # ) = Out of Range  
1R02827.D M1R0091.MSPCC's out = 0 CCC's out = 0  
Thu Apr 06 16:37:46 20236.7.5  
6

**Run Sequence Report**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Run ID: V1R100		Method: SW846 8260D		Instrument ID: GCMS1R
Lab Sample ID	Lab File ID	Date/Time Analyzed	Prep QC Batch	Client Sample ID
V1R100-BFB	1R03013.D	04/04/23 09:38	n/a	BFB Tune
V1R100-CC91	1R03013.D	04/04/23 09:38	n/a	Continuing cal 20
V1R100-BS	1R03015.D	04/04/23 10:34	n/a	Blank Spike
V1R100-MB	1R03017.D	04/04/23 11:25	n/a	Method Blank
ZZZZZZ	1R03018.D	04/04/23 11:54	n/a	(unrelated sample)
ZZZZZZ	1R03018R.D	04/04/23 11:54	n/a	(unrelated sample)
ZZZZZZ	1R03019.D	04/04/23 12:20	n/a	(unrelated sample)
ZZZZZZ	1R03019R.D	04/04/23 12:20	n/a	(unrelated sample)
ZZZZZZ	1R03020.D	04/04/23 12:46	n/a	(unrelated sample)
ZZZZZZ	1R03021.D	04/04/23 13:11	n/a	(unrelated sample)
ZZZZZZ	1R03021R.D	04/04/23 13:11	n/a	(unrelated sample)
ZZZZZZ	1R03022.D	04/04/23 13:37	n/a	(unrelated sample)
ZZZZZZ	1R03023.D	04/04/23 14:03	n/a	(unrelated sample)
ZZZZZZ	1R03024.D	04/04/23 14:28	n/a	(unrelated sample)
ZZZZZZ	1R03025.D	04/04/23 14:54	n/a	(unrelated sample)
ZZZZZZ	1R03026.D	04/04/23 15:19	n/a	(unrelated sample)
JD62888-1	1R03027.D	04/04/23 15:45	n/a	SVE-1
JD62953-1MS	1R03028.D	04/04/23 16:11	n/a	Matrix Spike
JD62953-1MSD	1R03029.D	04/04/23 16:36	n/a	Matrix Spike Duplicate
JD62888-4	1R03031.D	04/04/23 17:28	n/a	FB-1
JD62888-5	1R03032.D	04/04/23 17:53	n/a	TRIP BLANK
JD62953-1	1R03033.D	04/04/23 18:19	n/a	(used for QC only; not part of job JD62888)
JD62888-2	1R03034.D	04/04/23 18:45	n/a	SVE-4
JD62888-3	1R03035.D	04/04/23 19:10	n/a	SVE-5
ZZZZZZ	1R03036.D	04/04/23 19:36	n/a	(unrelated sample)
ZZZZZZ	1R03037.D	04/04/23 20:01	n/a	(unrelated sample)
ZZZZZZ	1R03038.D	04/04/23 20:27	n/a	(unrelated sample)
ZZZZZZ	1R03039.D	04/04/23 20:53	n/a	(unrelated sample)
ZZZZZZ	1R03040.D	04/04/23 21:18	n/a	(unrelated sample)

**Run Sequence Report**

Job Number: JD62888

Account: FLSNYNY Fleming-Lee Shue, Inc.

Project: 388 Bridge Street, Brooklyn, NY

Run ID: V1R91		Method: SW846 8260D		Instrument ID: GCMS1R
Lab Sample ID	Lab File ID	Date/Time Analyzed	Prep QC Batch	Client Sample ID
V1R91-BFB	1R02813.D	03/28/23 21:55	n/a	BFB Tune
V1R91-IC0091	1R02815.D	03/28/23 22:46	n/a	Initial cal 0.2
V1R91-IC0091	1R02817.D	03/28/23 23:19	n/a	Initial cal 0.5
V1R91-IC0091	1R02819.D	03/28/23 23:52	n/a	Initial cal 1
V1R91-IC0091	1R02821.D	03/29/23 00:25	n/a	Initial cal 2
V1R91-IC0091	1R02823.D	03/29/23 00:58	n/a	Initial cal 4
V1R91-IC0091	1R02825.D	03/29/23 01:31	n/a	Initial cal 8
V1R91-IC0091	1R02827.D	03/29/23 02:04	n/a	Initial cal 20
V1R91-ICC0091	1R02829.D	03/29/23 02:37	n/a	Initial cal 50
V1R91-IC0091	1R02831.D	03/29/23 03:10	n/a	Initial cal 100
V1R91-IC0091	1R02833.D	03/29/23 03:42	n/a	Initial cal 200
V1R91-ICV0091	1R02839.D	03/29/23 05:20	n/a	Initial cal verification 50
V1R91-ICV0091	1R02841.D	03/29/23 05:53	n/a	Initial cal verification 50
V1R91-BFB2	1R02846.D	03/29/23 16:59	n/a	BFB Tune
V1R91-ICV0091	1R02847.D	03/29/23 17:25	n/a	Initial cal verification 50

**MS Volatiles****Raw Data**

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03027.D  
 Acq On : 4 Apr 2023 3:45 pm  
 Operator : nickw  
 Sample : jd62888-1  
 Misc : MS67955,V1R0100,5,,,1  
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Apr 06 17:03:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.491	65	590128	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	553552	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.651	114	865233	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	839012	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	372781	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) dibromofluoromethane (s)	5.018	113	263310	54.73	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	109.46%		
53) 1,2-dichloroethane-d4 (s)	5.298	65	270684	53.81	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	107.62%		
74) toluene-d8 (s)	6.800	98	1094401	47.24	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	94.48%		
98) 4-bromofluorobenzene (s)	8.565	174	311494	51.06	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	102.12%		

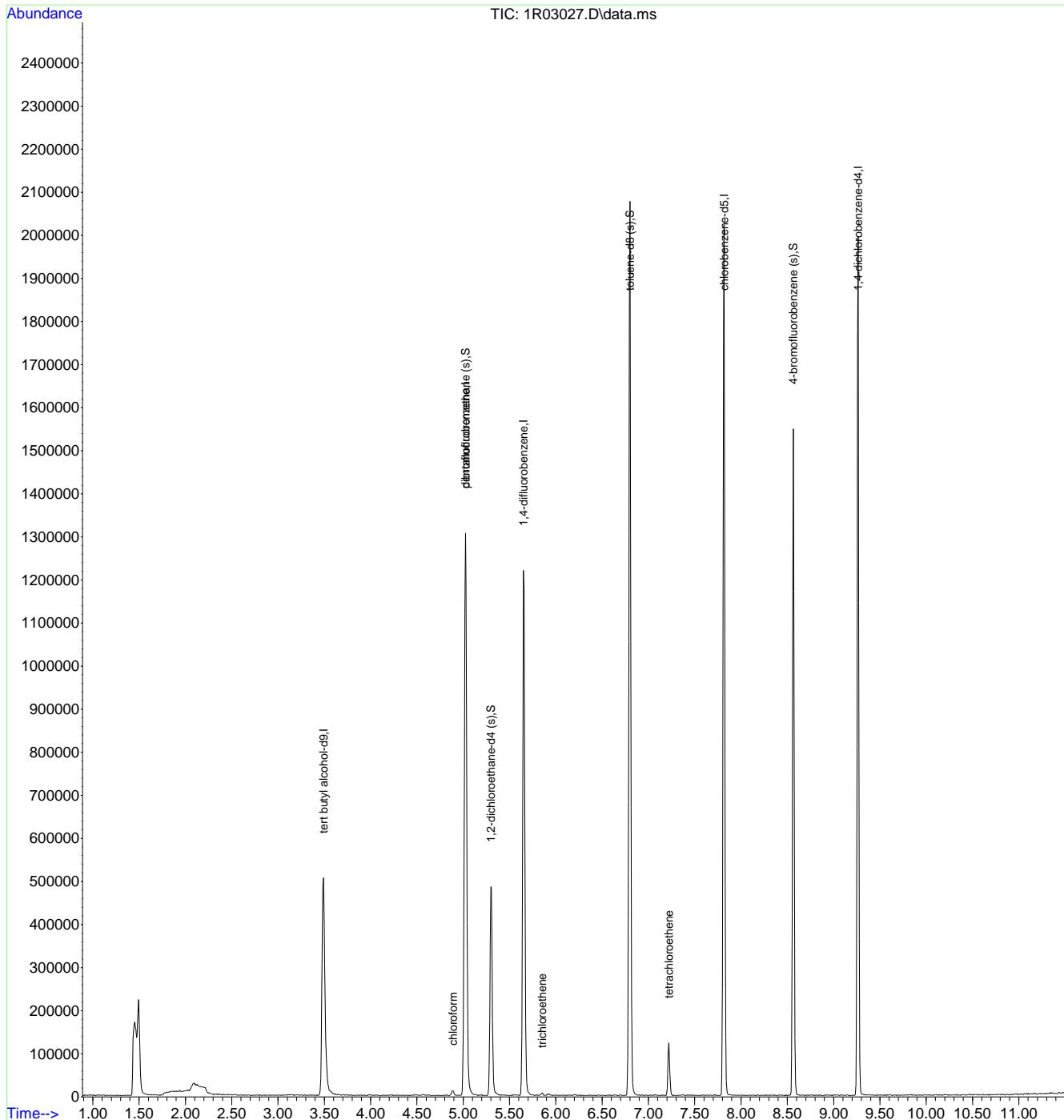
Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
42) chloroform	4.890	83	7565	1.19	ug/L	79
61) trichloroethene	5.851	95	1574	0.37	ug/L	# 63
79) tetrachloroethene	7.220	164	20820	4.86	ug/L	96

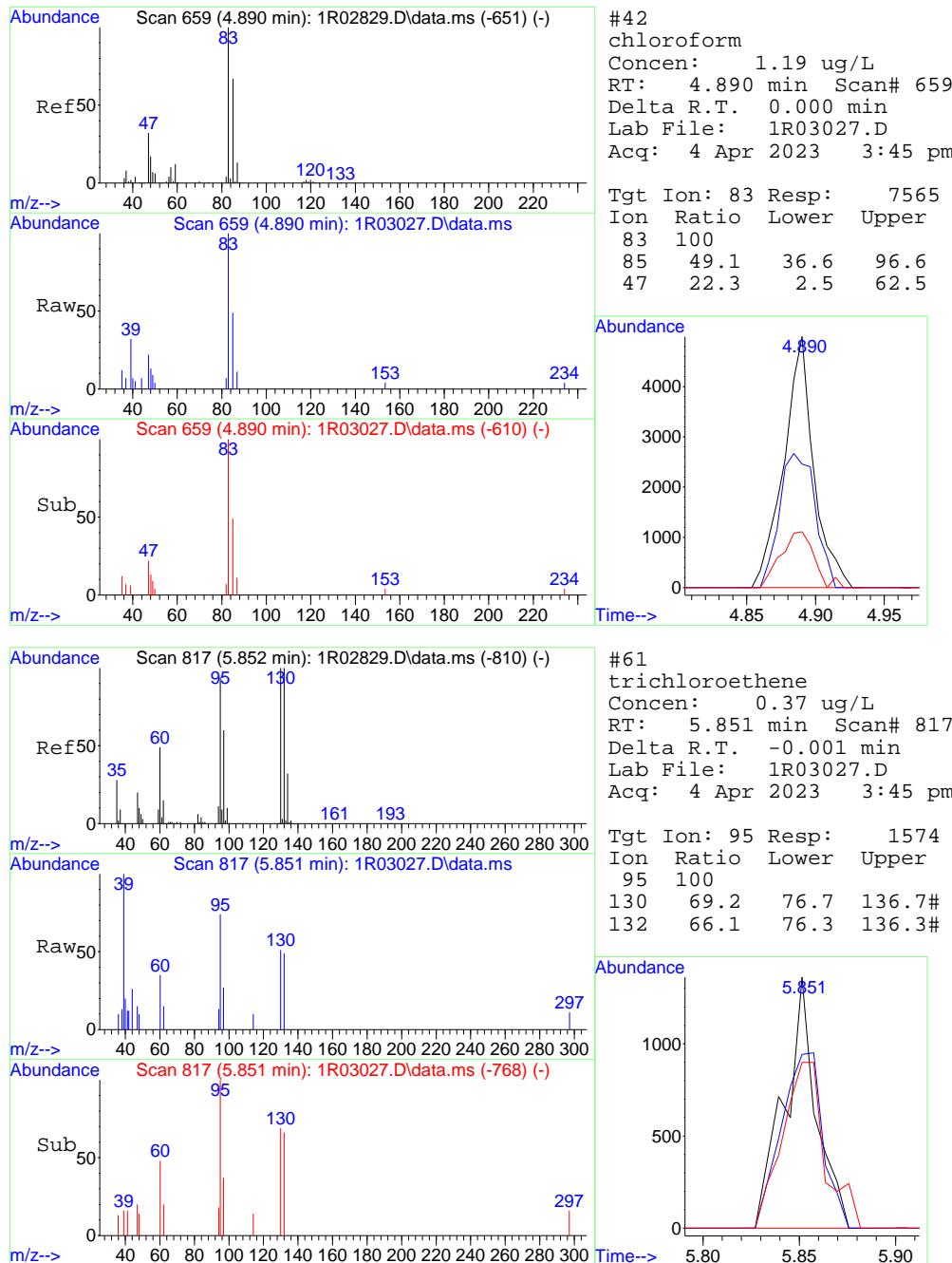
(#) = qualifier out of range (m) = manual integration (+) = signals summed

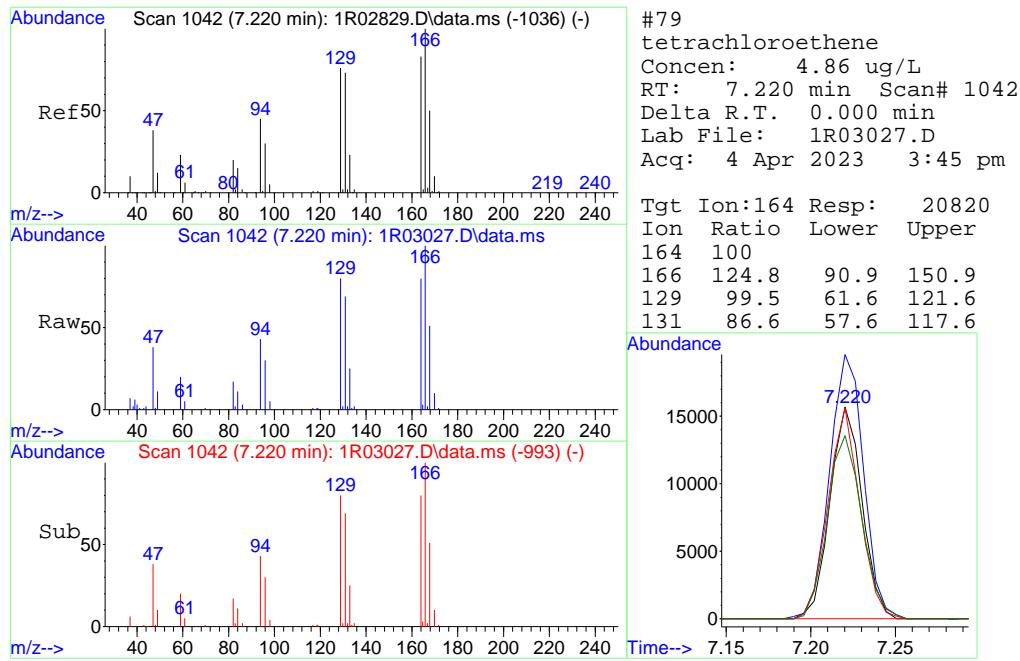
## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03027.D  
 Acq On : 4 Apr 2023 3:45 pm  
 Operator : nickw  
 Sample : jd62888-1  
 Misc : MS67955,V1R0100,5,,,1  
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Apr 06 17:03:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration







## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03034.D  
 Acq On : 4 Apr 2023 6:45 pm  
 Operator : nickw  
 Sample : jd62888-2  
 Misc : MS67891,V1R0100,5,,,1  
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Apr 06 17:21:31 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.491	65	575874	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	548691	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	858272	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	830614	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	379122	50.00	ug/L	0.00

System Monitoring Compounds						
44) dibromofluoromethane (s)	5.018	113	265101	55.59	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 111.18%	
53) 1,2-dichloroethane-d4 (s)	5.298	65	270071	54.12	ug/L	0.00
Spiked Amount	50.000	Range	81 - 124	Recovery	= 108.24%	
74) toluene-d8 (s)	6.800	98	1081414	47.15	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 94.30%	
98) 4-bromofluorobenzene (s)	8.565	174	315983	50.93	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 101.86%	

Target Compounds					Qvalue
36) cis-1,2-dichloroethene	4.610	96	5165	1.34	ug/L # 58
42) chloroform	4.890	83	8640	1.37	ug/L 84
61) trichloroethene	5.851	95	19678	4.61	ug/L 95
79) tetrachloroethene	7.220	164	63729	15.04	ug/L 93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\

Data File : 1R03034.D

Acq On : 4 Apr 2023 6:45 pm

Operator : nickw

Sample : jd62888-2

Misc : MS67891,V1R0100,5,,,1

ALS Vial : 23 Sample Multiplier: 1

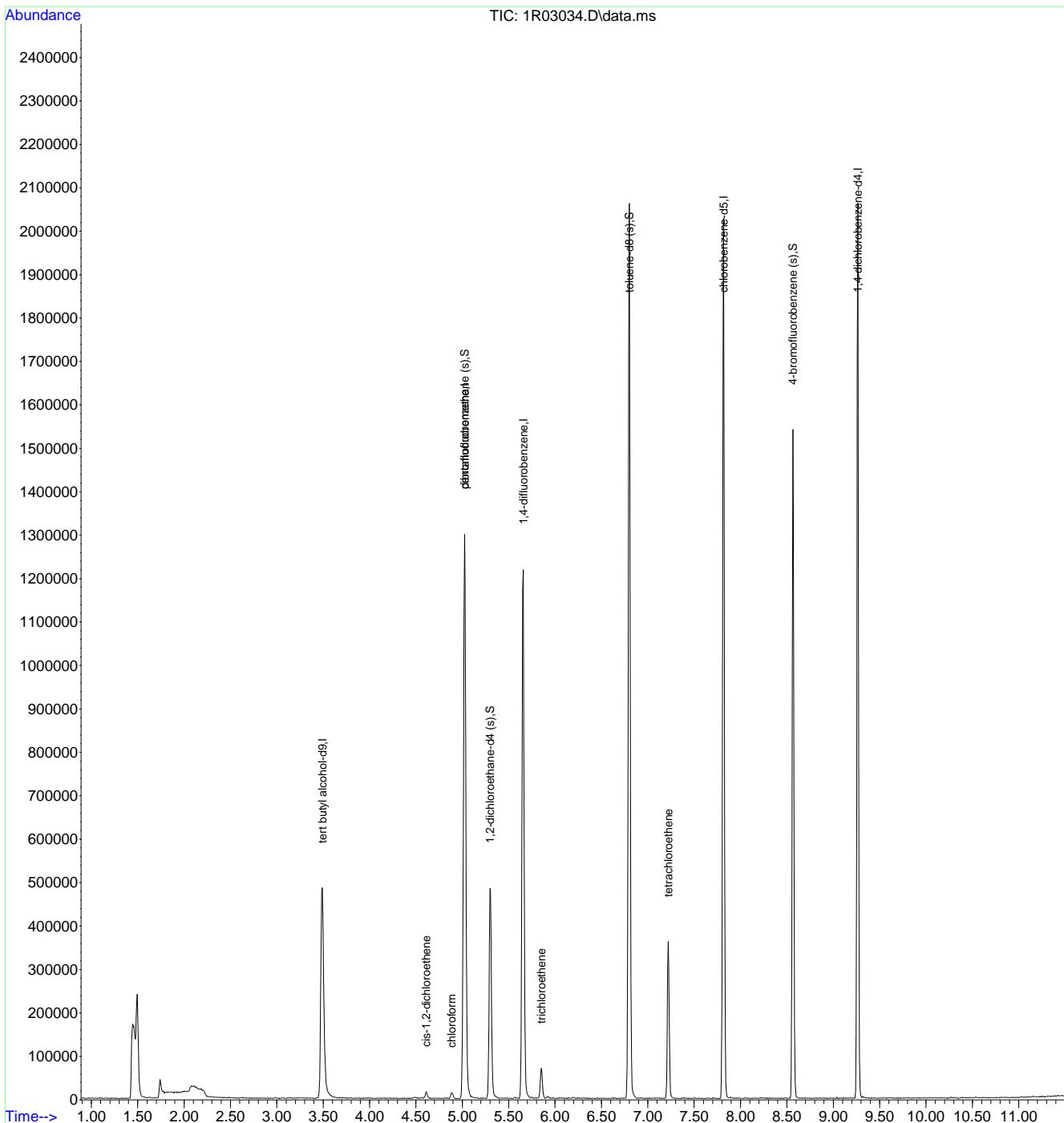
Quant Time: Apr 06 17:21:31 2023

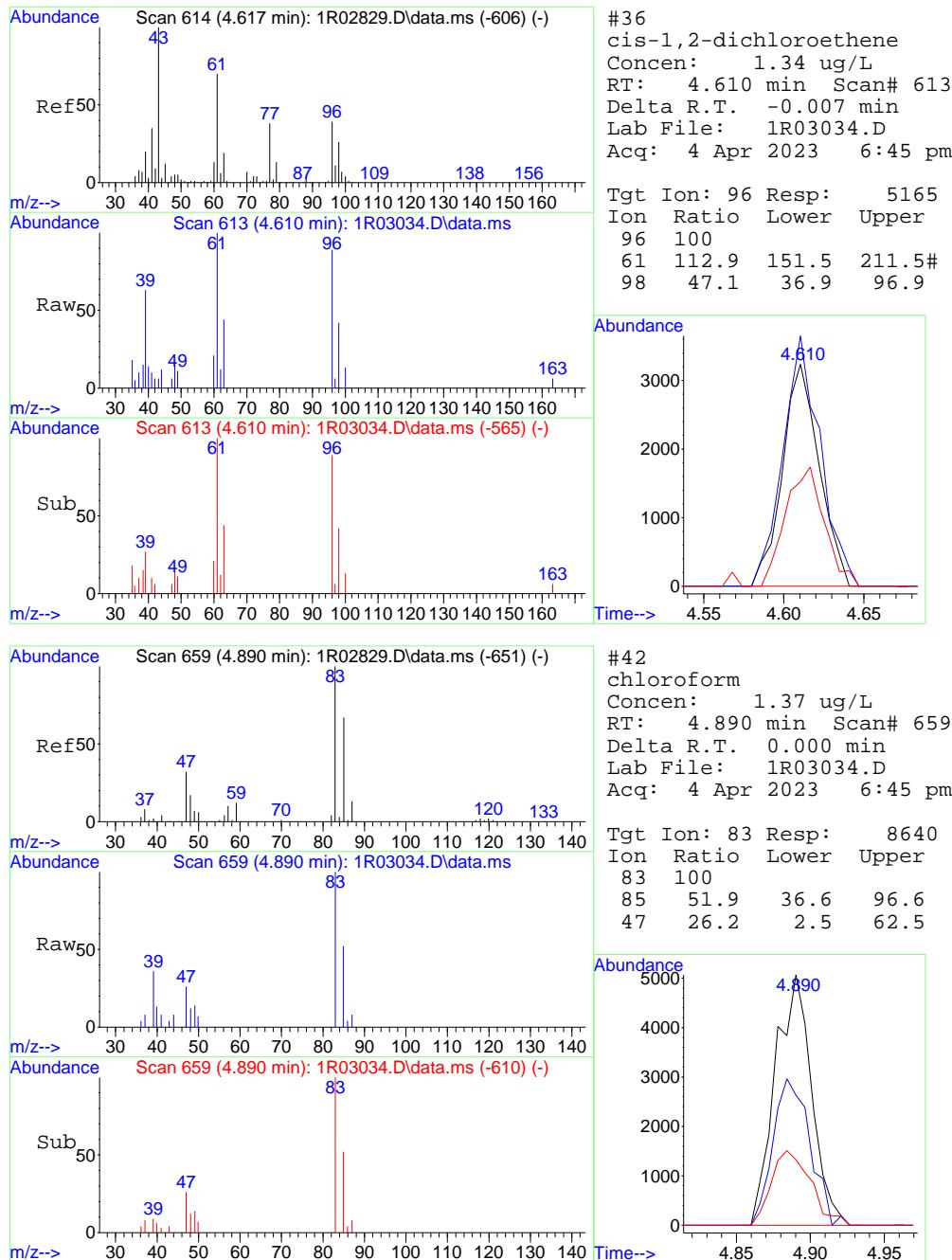
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M

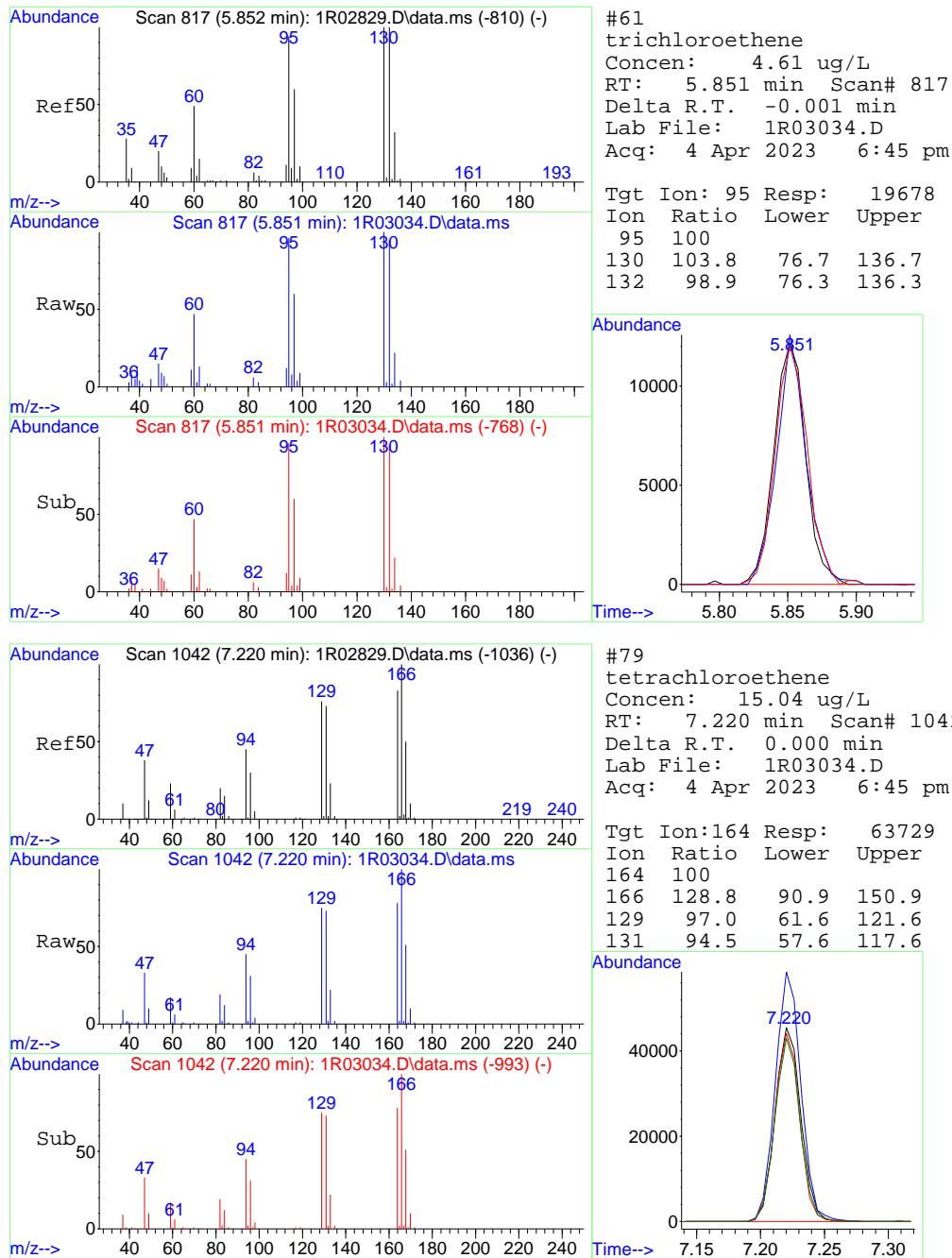
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um

QLast Update : Thu Mar 30 09:13:25 2023

Response via : Initial Calibration







## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03035.D  
 Acq On : 4 Apr 2023 7:10 pm  
 Operator : nickw  
 Sample : jd62888-3  
 Misc : MS67891,V1R0100,5,,,1  
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Apr 06 17:22:44 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.491	65	525305	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	542116	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	848679	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	827269	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	369643	50.00	ug/L	0.00

System Monitoring Compounds						
44) dibromofluoromethane (s)	5.024	113	258665	54.89	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 109.78%	
53) 1,2-dichloroethane-d4 (s)	5.304	65	266042	53.92	ug/L	0.00
Spiked Amount	50.000	Range	81 - 124	Recovery	= 107.84%	
74) toluene-d8 (s)	6.801	98	1072924	46.97	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 93.94%	
98) 4-bromofluorobenzene (s)	8.565	174	308412	50.99	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 101.98%	

Target Compounds					Qvalue
36) cis-1,2-dichloroethene	4.617	96	2349	0.62	ug/L # 45
42) chloroform	4.890	83	20271	3.24	ug/L 91
61) trichloroethene	5.851	95	8052	1.91	ug/L # 71

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\

Data File : 1R03035.D

Acq On : 4 Apr 2023 7:10 pm

Operator : nickw

Sample : jd62888-3

Misc : MS67891,V1R0100,5,,,1

ALS Vial : 24 Sample Multiplier: 1

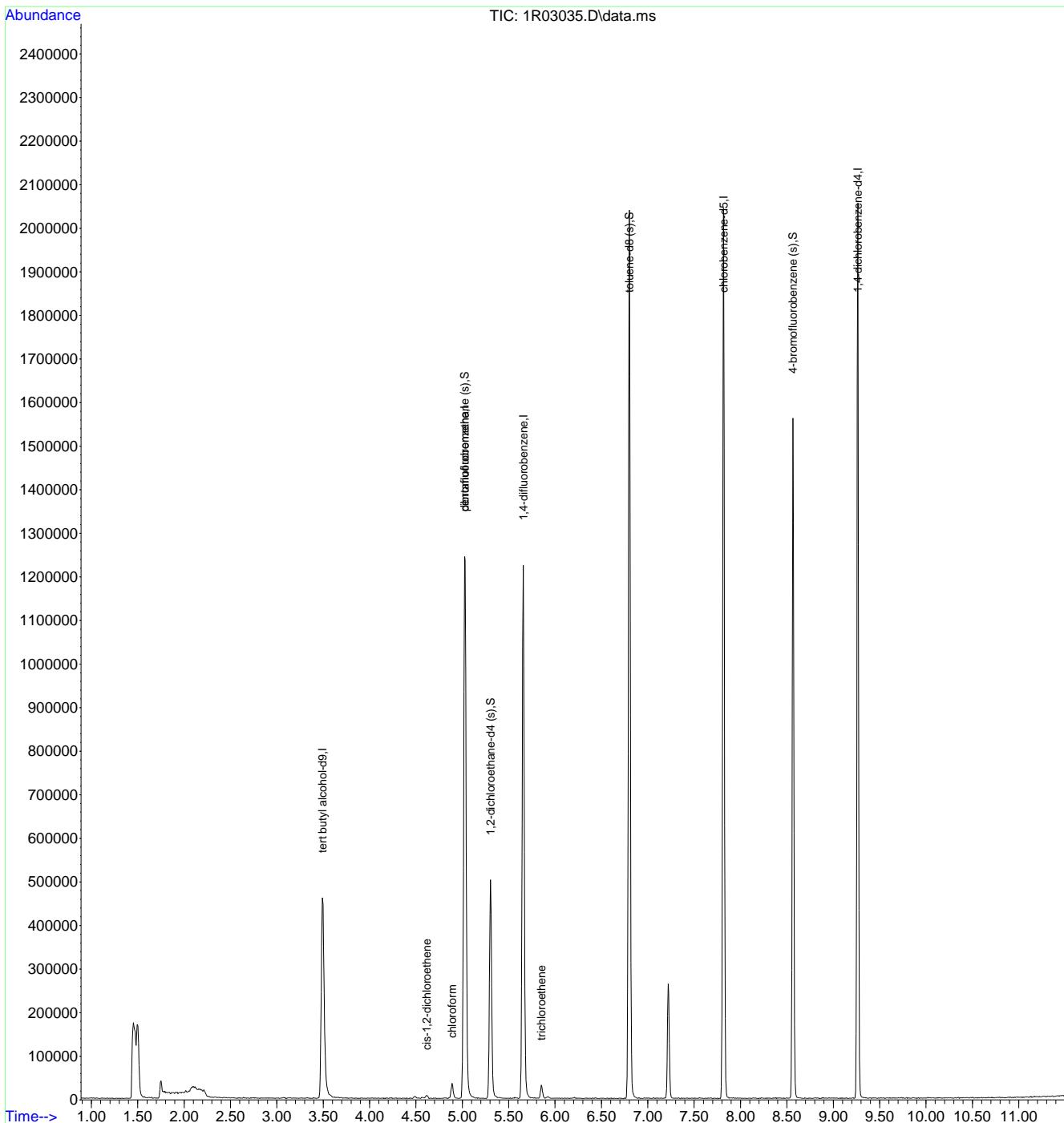
Quant Time: Apr 06 17:22:44 2023

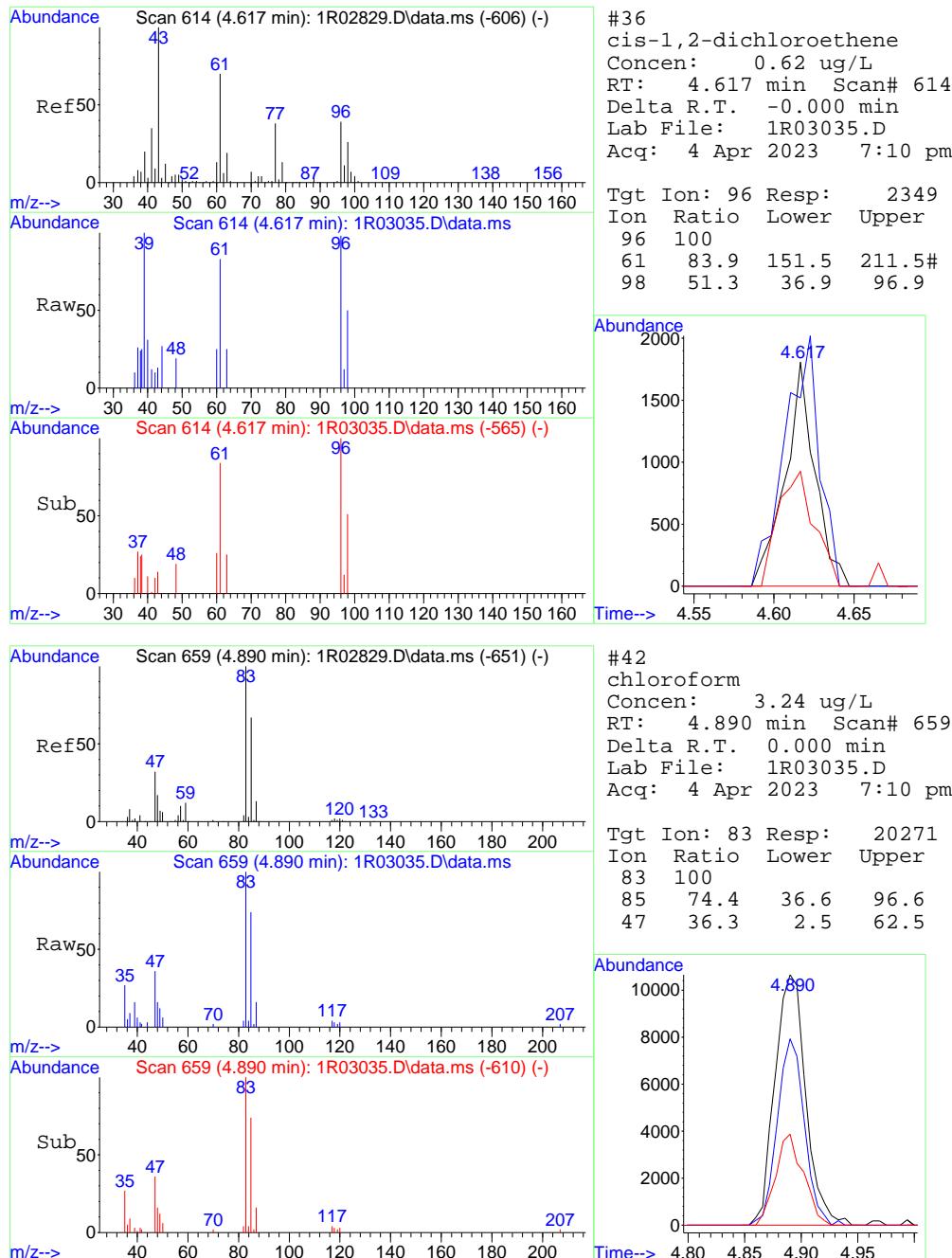
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M

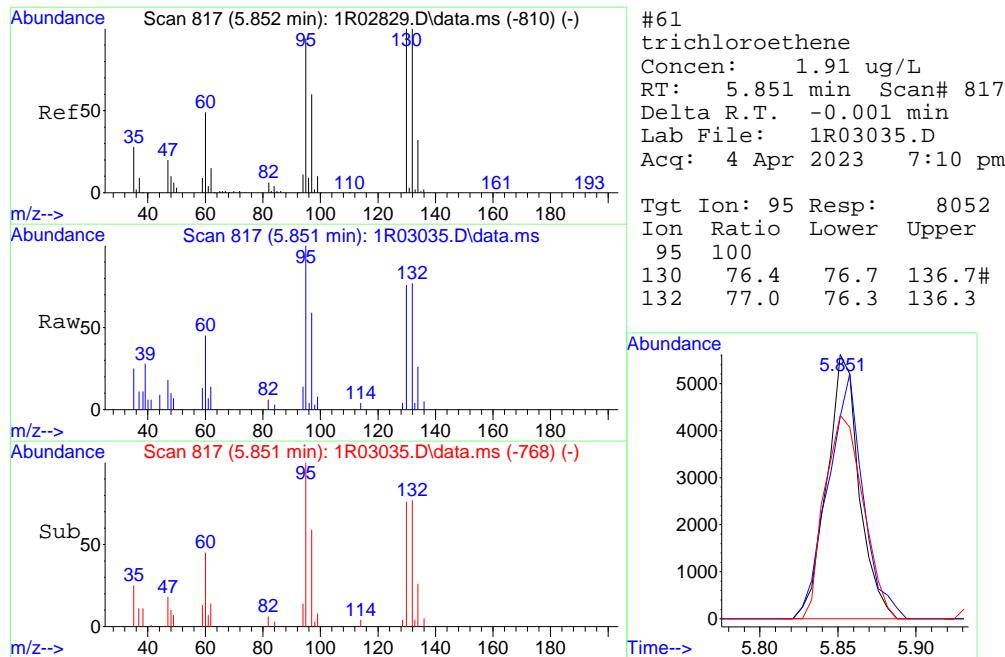
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um

QLast Update : Thu Mar 30 09:13:25 2023

Response via : Initial Calibration







## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03031.d

Acq On : 4 Apr 2023 5:28 pm  
 Operator : nickw  
 Sample : jd62888-4  
 Misc : MS67891,V1R0100,5,,,1  
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Apr 06 17:07:24 2023

Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M

Quant Title : SW846 8260D, Rx1624Sil MS 60m x 0.25mm x 1.4um

QLast Update : Wed Mar 29 12:39:36 2023

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.485	65	631950	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	583702	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	908888	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	871234	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	394162	50.00	ug/L	0.00

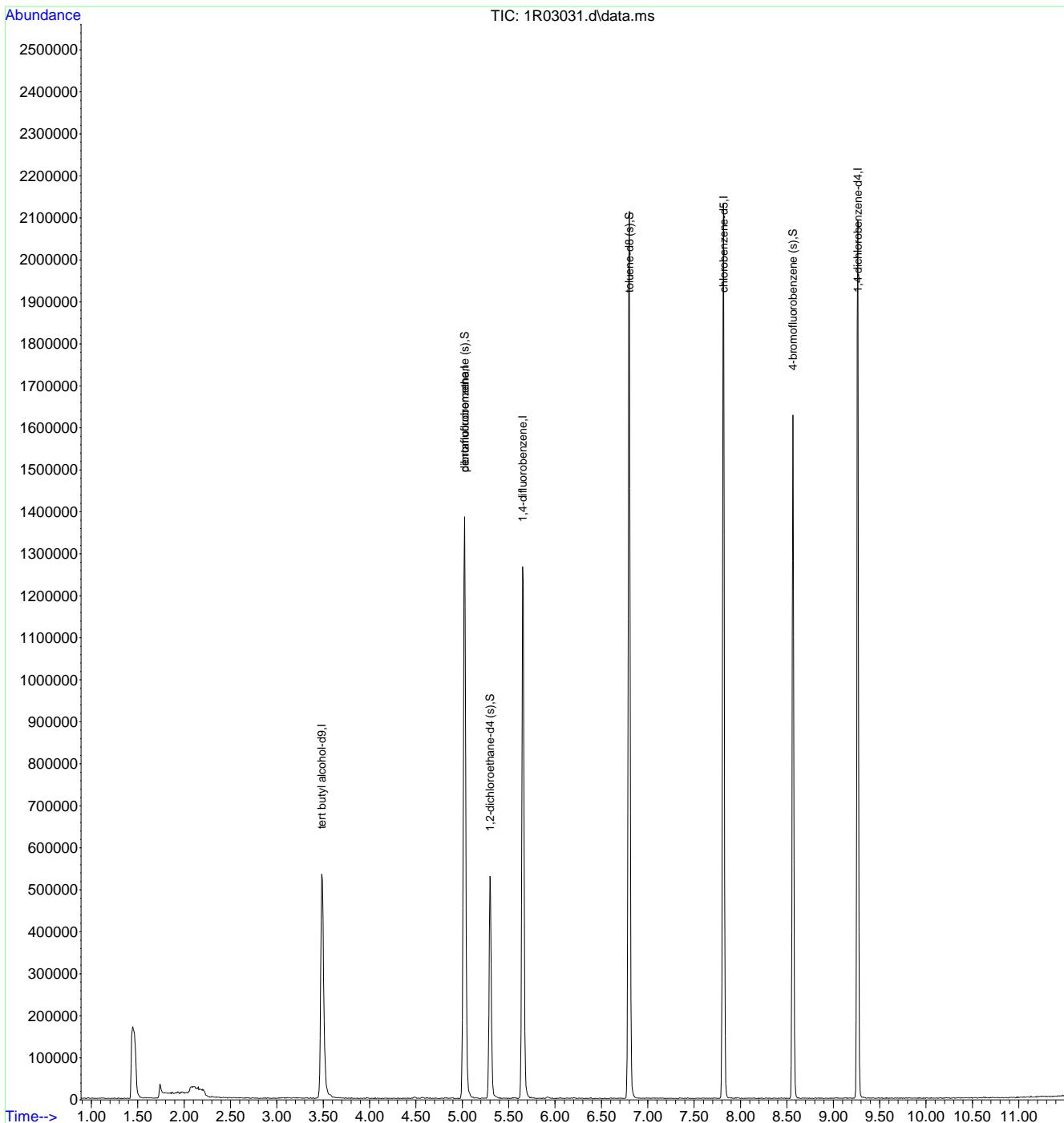
System Monitoring Compounds						
44) dibromofluoromethane (s)	5.018	113	275584	54.32	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 108.64%	
53) 1,2-dichloroethane-d4 (s)	5.298	65	281218	53.22	ug/L	0.00
Spiked Amount	50.000	Range	81 - 124	Recovery	= 106.44%	
74) toluene-d8 (s)	6.800	98	1143456	47.53	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 95.06%	
98) 4-bromofluorobenzene (s)	8.565	174	331187	51.35	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 102.70%	

Target Compounds	Qvalue
(#) = qualifier out of range (m) = manual integration (+) = signals summed	

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03031.d  
 Acq On : 4 Apr 2023 5:28 pm  
 Operator : nickw  
 Sample : jd62888-4  
 Misc : MS67891,V1R0100,5,,,1  
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Apr 06 17:07:24 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03032.D  
 Acq On : 4 Apr 2023 5:53 pm  
 Operator : nickw  
 Sample : jd62888-5  
 Misc : MS67891,V1R0100,5,,,1  
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Apr 06 17:08:20 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rx1624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.491	65	592479	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	568084	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	880717	50.00	ug/L	0.00
73) chlorobenzene-d5	7.817	117	851076	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	380661	50.00	ug/L	0.00

System Monitoring Compounds						
44) dibromofluoromethane (s)	5.018	113	271677	55.02	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	110.04%
53) 1,2-dichloroethane-d4 (s)	5.304	65	278371	54.36	ug/L	0.00
Spiked Amount	50.000	Range	81 - 124	Recovery	=	108.72%
74) toluene-d8 (s)	6.801	98	1121187	47.71	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	95.42%
98) 4-bromofluorobenzene (s)	8.565	174	326570	52.43	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	104.86%

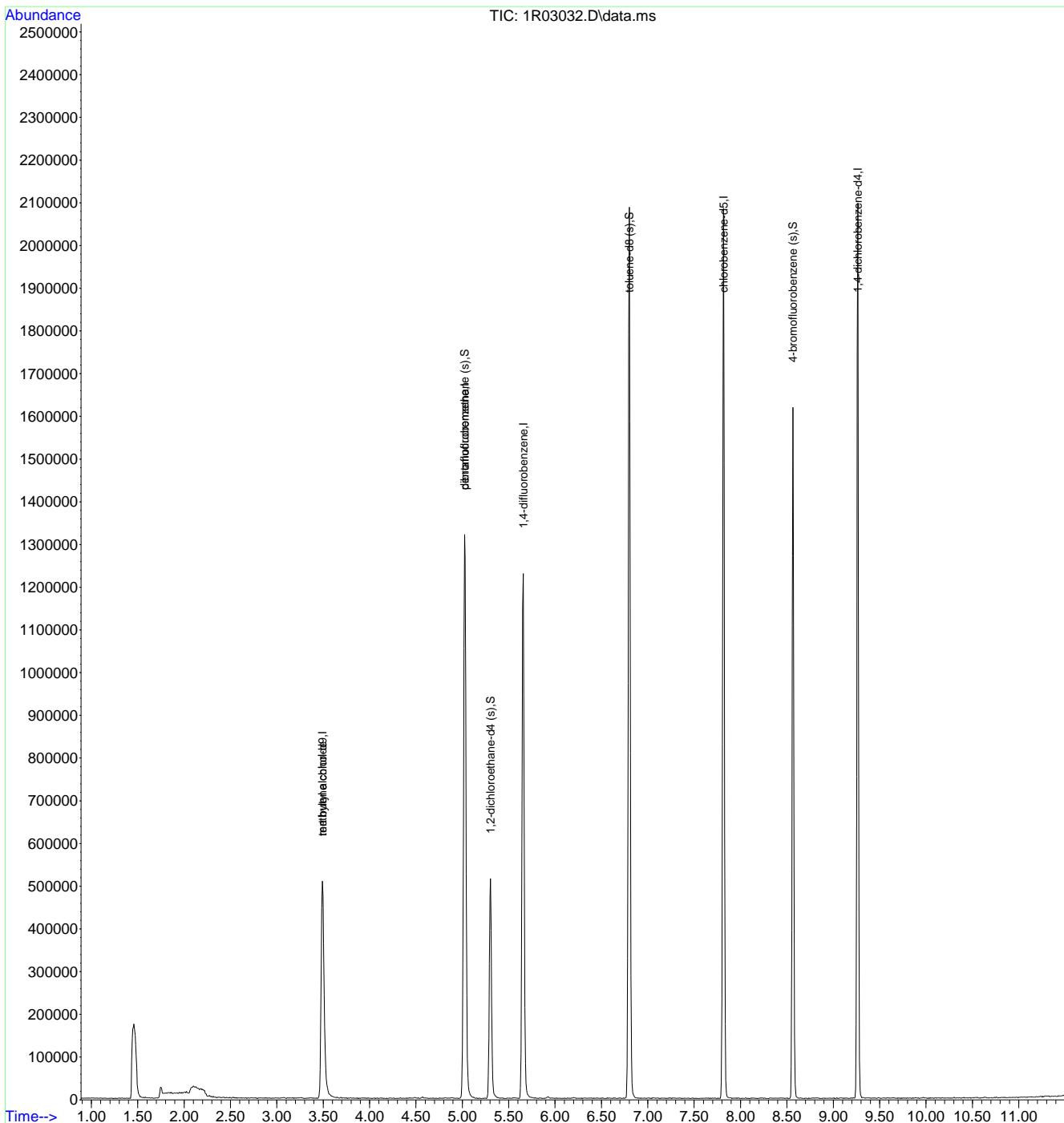
Target Compounds					Qvalue
22) methylene chloride	3.497	84	1653	0.47	ug/L

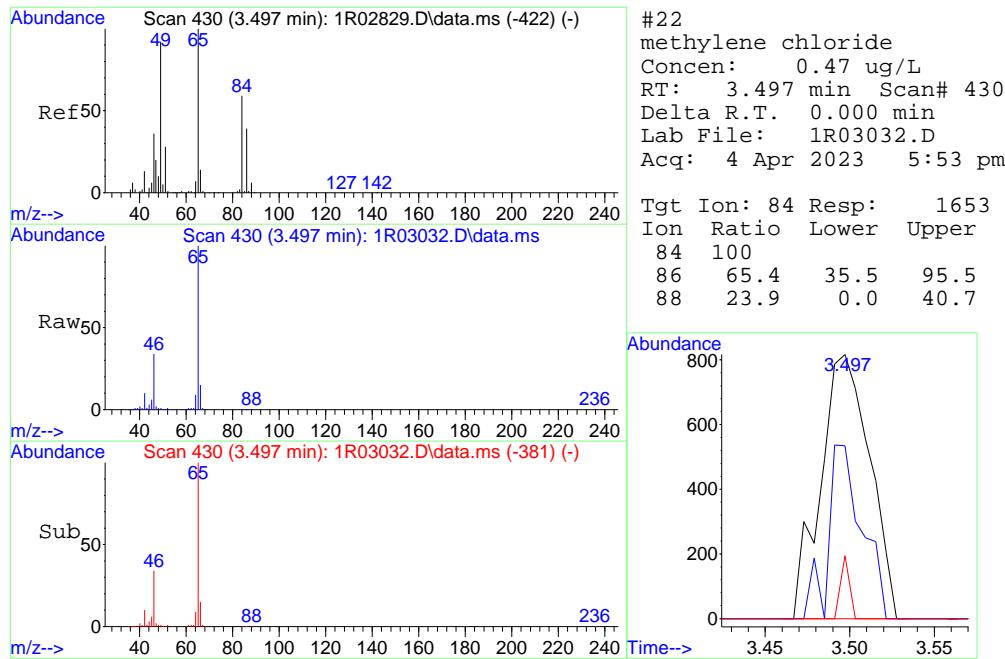
(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03032.D  
 Acq On : 4 Apr 2023 5:53 pm  
 Operator : nickw  
 Sample : jd62888-5  
 Misc : MS67891,V1R0100,5,,,1  
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Apr 06 17:08:20 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration





## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03017.D  
 Acq On : 4 Apr 2023 11:25 am  
 Operator : nickw  
 Sample : mb  
 Misc : MS67936,V1R0100,5,,,1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Apr 06 16:43:57 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.485	65	612313	500.00	ug/L	0.00
5) pentafluorobenzene	5.024	168	592616	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.651	114	922312	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	902372	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	401555	50.00	ug/L	0.00

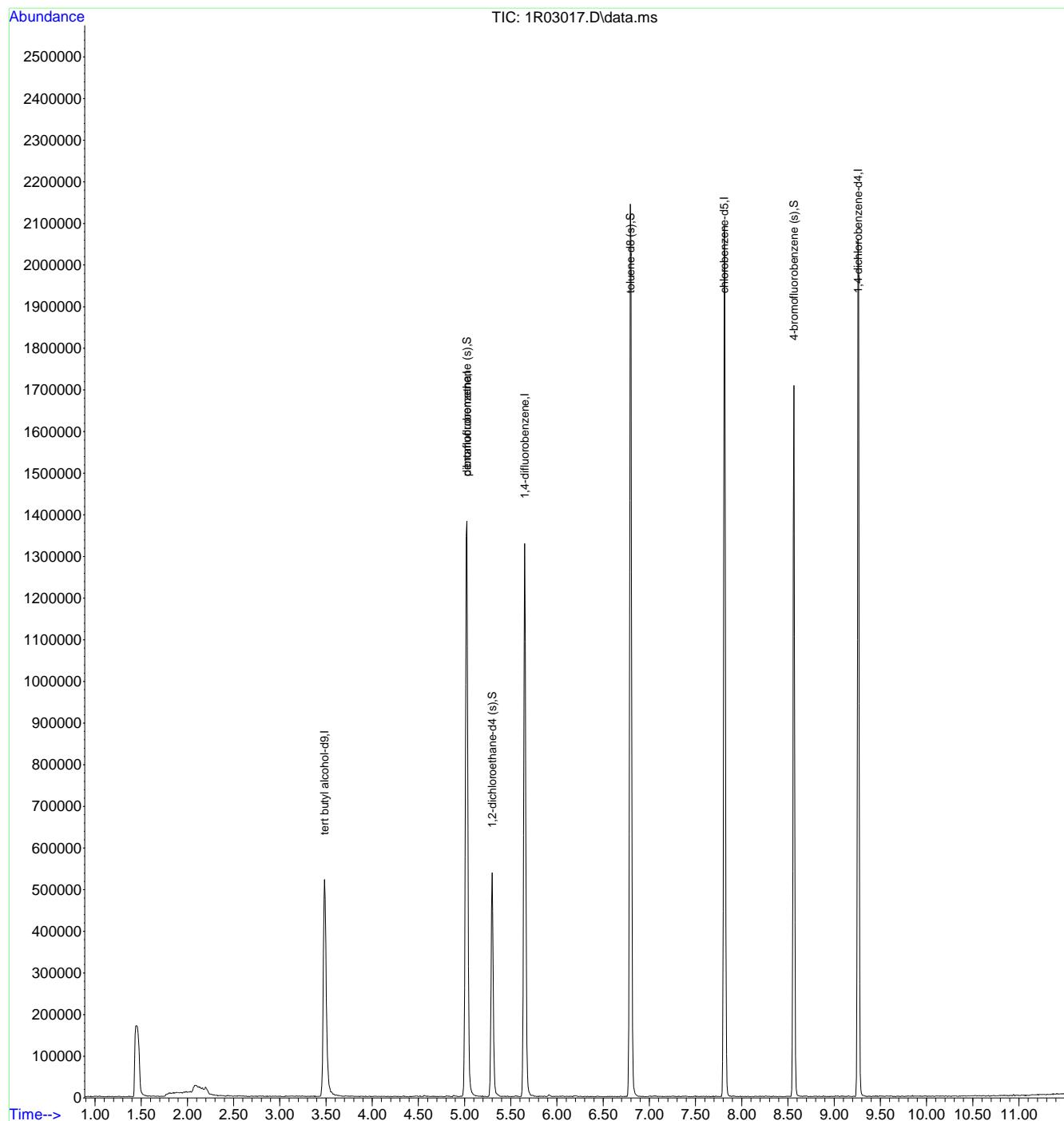
System Monitoring Compounds						
44) dibromofluoromethane (s)	5.018	113	281309	54.61	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 109.22%	
53) 1,2-dichloroethane-d4 (s)	5.298	65	287938	53.70	ug/L	0.00
Spiked Amount	50.000	Range	81 - 124	Recovery	= 107.40%	
74) toluene-d8 (s)	6.794	98	1180225	47.36	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 94.72%	
98) 4-bromofluorobenzene (s)	8.565	174	339546	51.67	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	= 103.34%	

Target Compounds	Qvalue
(#) = qualifier out of range (m) = manual integration (+) = signals summed	

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03017.D  
 Acq On : 4 Apr 2023 11:25 am  
 Operator : nickw  
 Sample : mb  
 Misc : MS67936,V1R0100,5,,,1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Apr 06 16:43:57 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03015.D  
 Acq On : 4 Apr 2023 10:34 am  
 Operator : nickw  
 Sample : bs  
 Misc : MS67909,V1R0100,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Apr 04 11:08:20 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.491	65	614235	500.00	ug/L	0.00
5) pentafluorobenzene	5.024	168	730438	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.651	114	1084886	50.00	ug/L	0.00
73) chlorobenzene-d5	7.810	117	1086387	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.258	152	571074	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) dibromofluoromethane (s)	5.012	113	337209	53.11	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	106.22%		
53) 1,2-dichloroethane-d4 (s)	5.298	65	326110	51.70	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	103.40%		
74) toluene-d8 (s)	6.795	98	1389050	46.30	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	92.60%		
98) 4-bromofluorobenzene (s)	8.565	174	458953	49.11	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	98.22%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) ethanol	2.688	45	574417	5184.18	ug/L	98
3) tertiary butyl alcohol	3.564	59	350002	227.15	ug/L	96
4) 1,4-dioxane	6.113	88	134188	1168.37	ug/L	95
6) chlorodifluoromethane	1.654	51	187011	33.56	ug/L	95
7) dichlorodifluoromethane	1.636	85	232241	45.08	ug/L	99
8) chloromethane	1.800	50	302762	38.07	ug/L	97
9) vinyl chloride	1.897	62	268130	31.88	ug/L	98
10) 1,3-butadiene	1.934	54	163161	25.96	ug/L	95
11) bromomethane	2.183	94	124864	26.52	ug/L	97
12) chloroethane	2.287	64	155898	36.80	ug/L	100
13) trichlorofluoromethane	2.530	101	332795	51.62	ug/L	97
14) ethyl ether	2.798	74	121370	45.69	ug/L	92
15) acrolein	2.938	56	75252	49.22	ug/L	98
16) freon 113	3.035	151	151025	41.55	ug/L	95
17) 1,1-dichloroethene	3.035	96	178484	43.22	ug/L	98
18) acetone	3.065	58	239222	239.72	ug/L	98
19) acetonitrile	3.327	41	641665	477.30	ug/L	98
20) iodomethane	3.181	142	110271	38.83	ug/L	97
21) carbon disulfide	3.248	76	478882	39.48	ug/L	99
22) methylene chloride	3.491	84	209619	46.55	ug/L	96
23) methyl acetate	3.357	74	64521	47.94	ug/L #	63
24) methyl tert butyl ether	3.722	73	608741	47.33	ug/L	98
25) trans-1,2-dichloroethene	3.734	96	205228	44.76	ug/L	97
26) hexane	3.978	56	161844	41.64	ug/L	98
27) di-isopropyl ether	4.136	45	814994	44.66	ug/L	98
28) 2-butanone	4.586	72	245257	209.09	ug/L	94
29) 1,1-dichloroethane	4.124	63	374203	43.48	ug/L	97
30) chloroprene	4.185	53	332575	44.84	ug/L	97
31) acrylonitrile	3.698	53	163585	49.10	ug/L	97
32) vinyl acetate	4.112	86	57072	48.60	ug/L #	79
33) ethyl tert-butyl ether	4.446	59	699862	48.55	ug/L	98
34) ethyl acetate	4.617	45	75998	46.20	ug/L #	69
35) 2,2-dichloropropane	4.617	77	302428	48.01	ug/L	89
36) cis-1,2-dichloroethene	4.604	96	233468	45.55	ug/L	89
37) propionitrile	4.647	54	692922	457.71	ug/L	99
38) methyl acrylate	4.659	85	57221	46.23	ug/L #	81
39) methacrylonitrile	4.775	67	144686	44.22	ug/L	94
40) bromochloromethane	4.805	128	114929	47.32	ug/L	88
41) tetrahydrofuran	4.817	42	160131	39.24	ug/L	97
42) chloroform	4.884	83	383734	45.56	ug/L	96
43) tert-Butyl Formate	4.903	59	239381	52.17	ug/L	95

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03015.D  
 Acq On : 4 Apr 2023 10:34 am  
 Operator : nickw  
 Sample : bs  
 Misc : MS67909,V1R0100,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Apr 04 11:08:20 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1,1-trichloroethane	5.030	97	328154	45.20	ug/L	92
46) cyclohexane	5.091	84	333997	49.56	ug/L	84
47) isobutyl alcohol	5.207	43	310889	504.80	ug/L	99
48) 1,1-dichloropropene	5.158	75	275986	45.37	ug/L	93
49) carbon tetrachloride	5.158	117	293329	48.47	ug/L	98
50) tert-amyl alcohol	5.304	73	181596	248.57	ug/L	97
51) isopropyl acetate	5.328	87	77308	47.55	ug/L #	88
54) n-butyl alcohol	5.742	56	960798	2359.13	ug/L	98
55) 2,2,4-trimethylpentane	5.407	57	586116	38.94	ug/L	98
56) benzene	5.322	78	856388	40.04	ug/L	99
57) tert-amyl methyl ether	5.407	73	656560	43.17	ug/L	99
58) heptane	5.541	57	134395	39.51	ug/L	89
59) 1,2-dichloroethane	5.359	62	299073	43.90	ug/L	93
60) ethyl acrylate	5.894	55	452569	41.07	ug/L	99
61) trichloroethene	5.845	95	233229	43.22	ug/L	98
62) 2-chloroethyl vinyl ether	6.466	63	960090	207.09	ug/L	96
63) methyl methacrylate	6.077	100	83901	45.99	ug/L #	91
64) methylcyclohexane	6.040	83	333325	38.99	ug/L	97
65) 1,2-dichloropropane	6.058	63	231044	39.22	ug/L	97
66) dibromomethane	6.131	93	153965	42.29	ug/L	97
67) bromodichloromethane	6.259	83	292900	45.13	ug/L	96
68) 2-nitropropane	6.423	41	131642	43.03	ug/L	90
69) epichlorohydrin	6.509	57	263735	238.55	ug/L	99
70) cis-1,3-dichloropropene	6.594	75	367837	44.05	ug/L	96
71) 4-methyl-2-pentanone	6.691	58	681179	167.56	ug/L	95
72) isoamyl alcohol	6.722	70	336472	888.55	ug/L	96
75) toluene	6.849	92	578719	41.12	ug/L	99
76) ethyl methacrylate	7.026	69	340526	44.36	ug/L	96
77) trans-1,3-dichloropropene	7.001	75	347252	46.12	ug/L	92
78) 1,1,2-trichloroethane	7.147	83	187356	42.57	ug/L	94
79) tetrachloroethene	7.214	164	225538	40.69	ug/L	96
80) 2-hexanone	7.281	58	771889	173.64	ug/L	98
81) 1,3-dichloropropane	7.263	76	345649	43.97	ug/L	93
82) butyl acetate	7.354	56	240041	41.22	ug/L	93
83) dibromochloromethane	7.421	129	270670	48.83	ug/L	98
84) 1,2-dibromoethane	7.506	107	255320	48.48	ug/L	97
85) n-butyl ether	7.859	57	1002667	40.20	ug/L	99
86) chlorobenzene	7.835	112	680818	43.75	ug/L	95
87) 1,1,1,2-tetrachloroethane	7.883	131	252330	45.87	ug/L	96
88) ethylbenzene	7.883	91	1080915	41.80	ug/L	99
89) m,p-xylene	7.969	106	866206	86.29	ug/L	95
90) o-xylene	8.218	91	839250	42.07	ug/L	96
91) styrene	8.230	104	743473	44.40	ug/L	97
92) butyl acrylate	8.157	55	571326	43.69	ug/L	98
93) n-amyl acetate	8.297	70	204443	40.83	ug/L	98
94) isopropylbenzene	8.443	105	1075960	45.96	ug/L	98
95) bromoform	8.358	173	212540	48.24	ug/L	97
96) cis-1,4-dichloro-2-butene	8.480	88	158082	48.18	ug/L	98
99) 1,1,2,2-tetrachloroethane	8.632	83	369188	43.01	ug/L	99
100) trans-1,4-dichloro-2-b...	8.656	53	123155	42.67	ug/L	78
101) 1,2,3-trichloropropane	8.674	110	117956	46.36	ug/L	90
102) bromobenzene	8.662	156	302094	44.75	ug/L #	81
103) n-propylbenzene	8.705	91	1228754	42.09	ug/L	97
104) 2-chlorotoluene	8.772	126	275114	45.56	ug/L	89
105) 4-chlorotoluene	8.839	91	738410	41.17	ug/L	98
106) 1,3,5-trimethylbenzene	8.808	105	863129	42.95	ug/L	96
107) tert-butylbenzene	9.003	119	757758	43.18	ug/L	95
108) 1,2,4-trimethylbenzene	9.033	105	917157	44.84	ug/L	98

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03015.D  
 Acq On : 4 Apr 2023 10:34 am  
 Operator : nickw  
 Sample : bs  
 Misc : MS67909,V1R0100,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Apr 04 11:08:20 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
109) sec-butylbenzene	9.131	105	1055167	44.14	ug/L	97
110) p-isopropyltoluene	9.216	119	959144	43.80	ug/L	97
111) 1,2,3-trimethylbenzene	9.283	105	922368	42.31	ug/L	96
112) 1,3-dichlorobenzene	9.216	146	569831	43.88	ug/L	99
113) 1,4-dichlorobenzene	9.277	146	571446	43.08	ug/L	97
114) 1,2-dichlorobenzene	9.490	146	563896	47.32	ug/L	96
115) benzyl chloride	9.337	91	792701	49.53	ug/L	94
116) n-butylbenzene	9.453	92	412790	41.93	ug/L	95
117) hexachloroethane	9.648	201	109204	52.41	ug/L	93
118) 1,2-dibromo-3-chloropr...	9.928	157	118714	51.04	ug/L	96
119) 1,3,5-trichlorobenzene	10.031	180	335856	44.69	ug/L	98
120) 1,2,4-trichlorobenzene	10.378	180	298375	47.17	ug/L	92
121) hexachlorobutadiene	10.445	225	103057	44.85	ug/L	97
122) naphthalene	10.530	128	1119638	48.47	ug/L	98
123) 1,2,3-trichlorobenzene	10.645	180	288473	46.09	ug/L	93
124) 2-methylnaphthalene	11.132	142	252014	24.96	ug/L	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

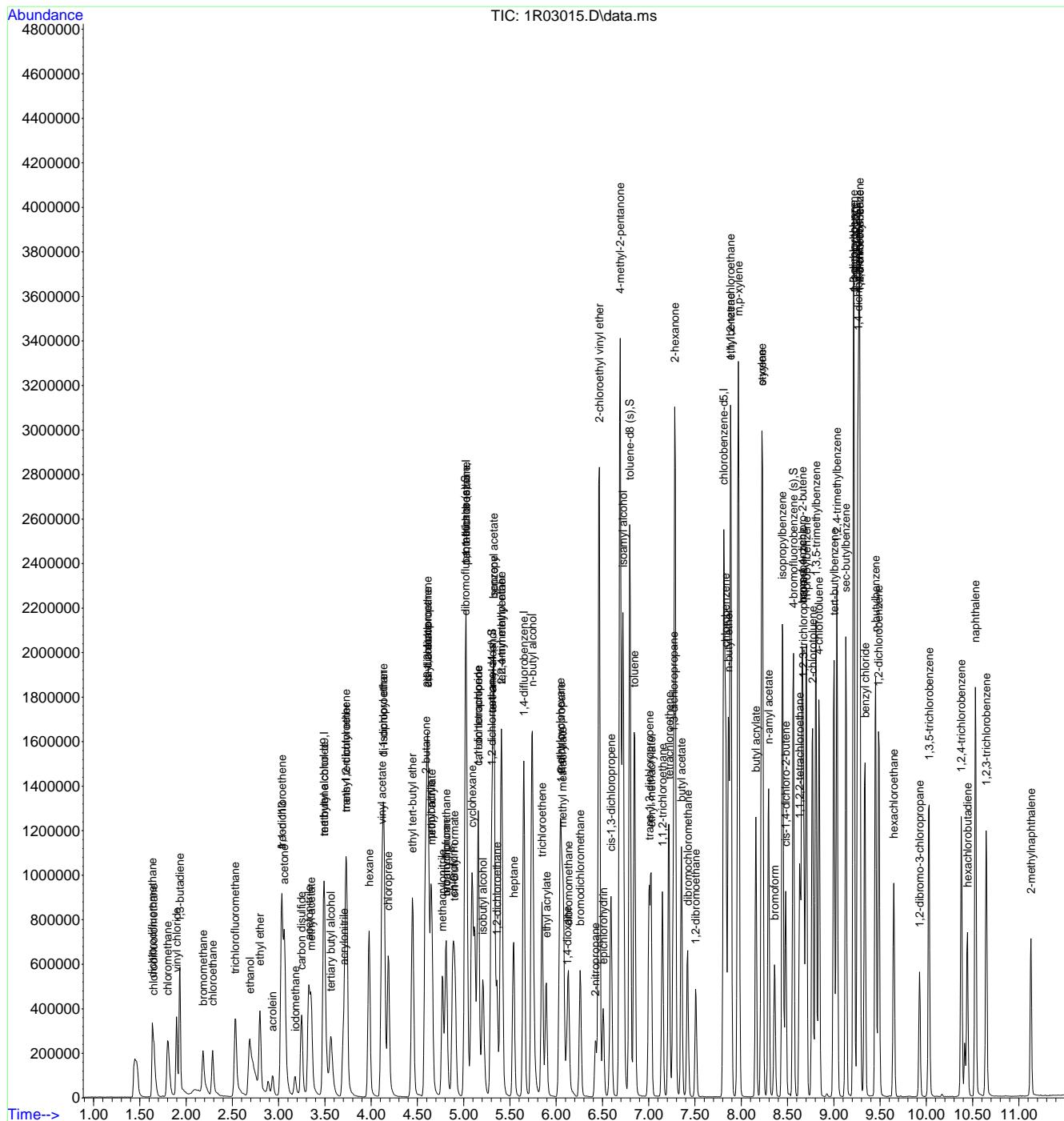
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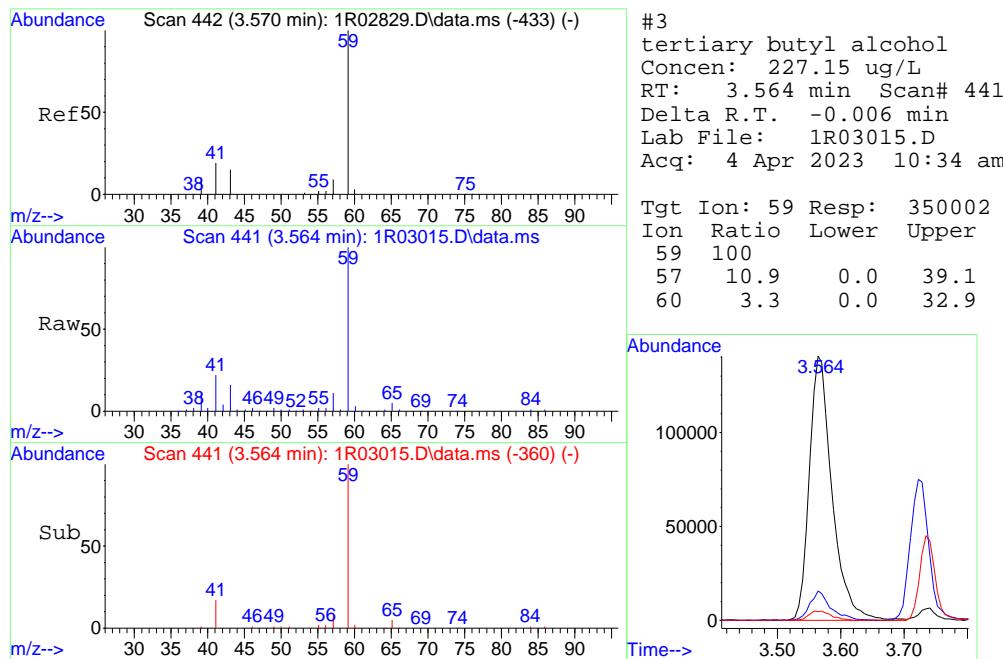
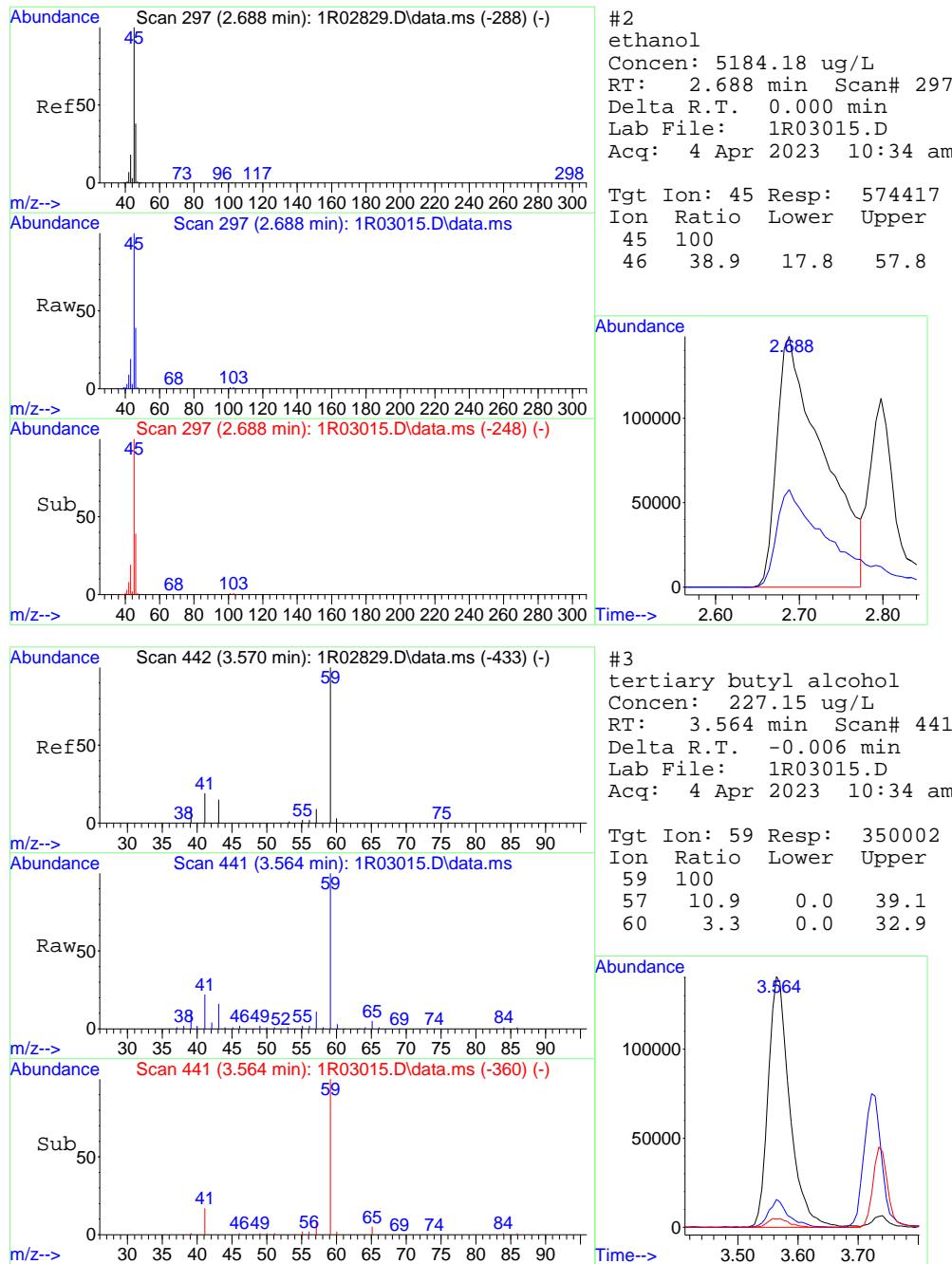
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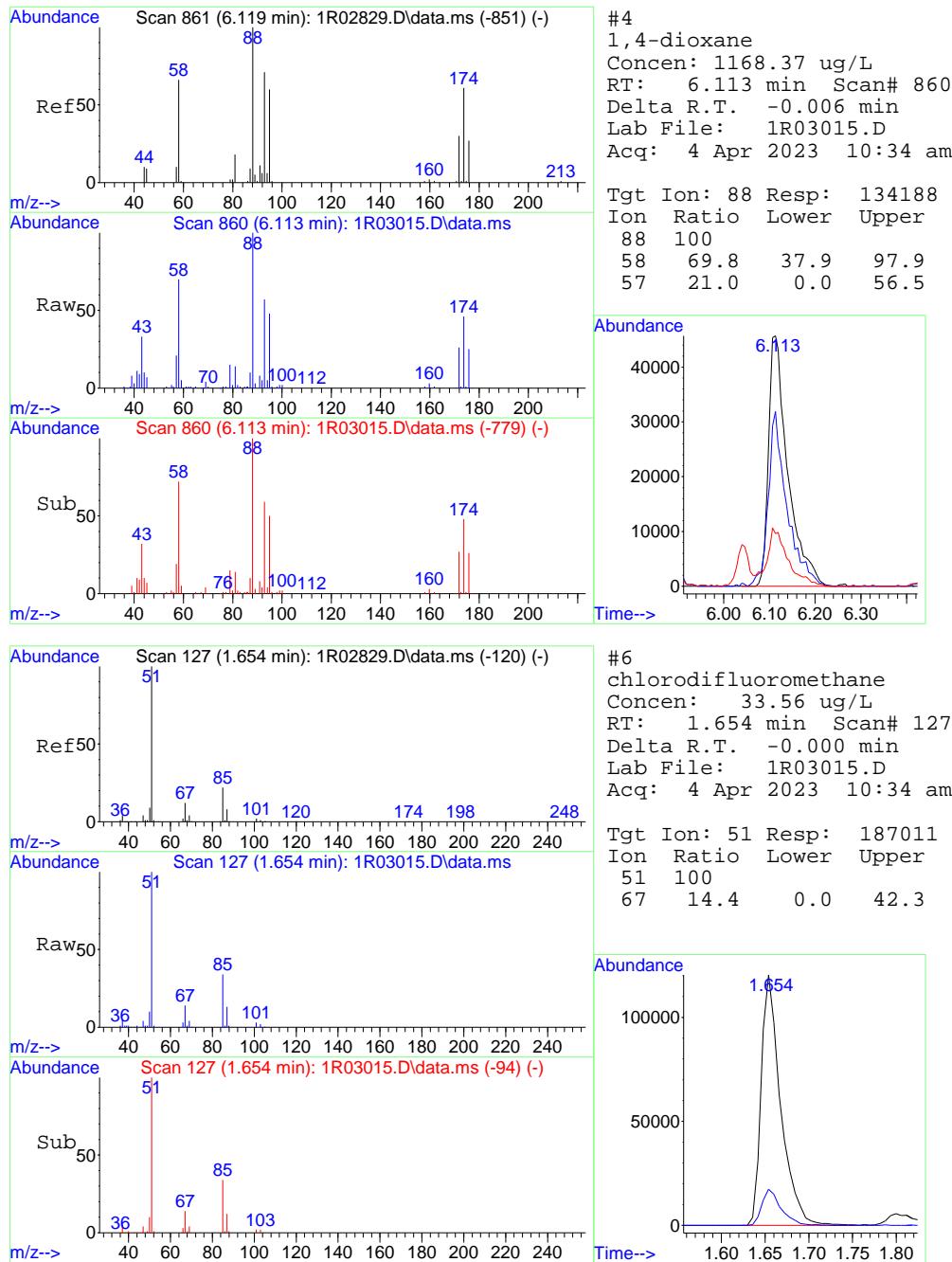
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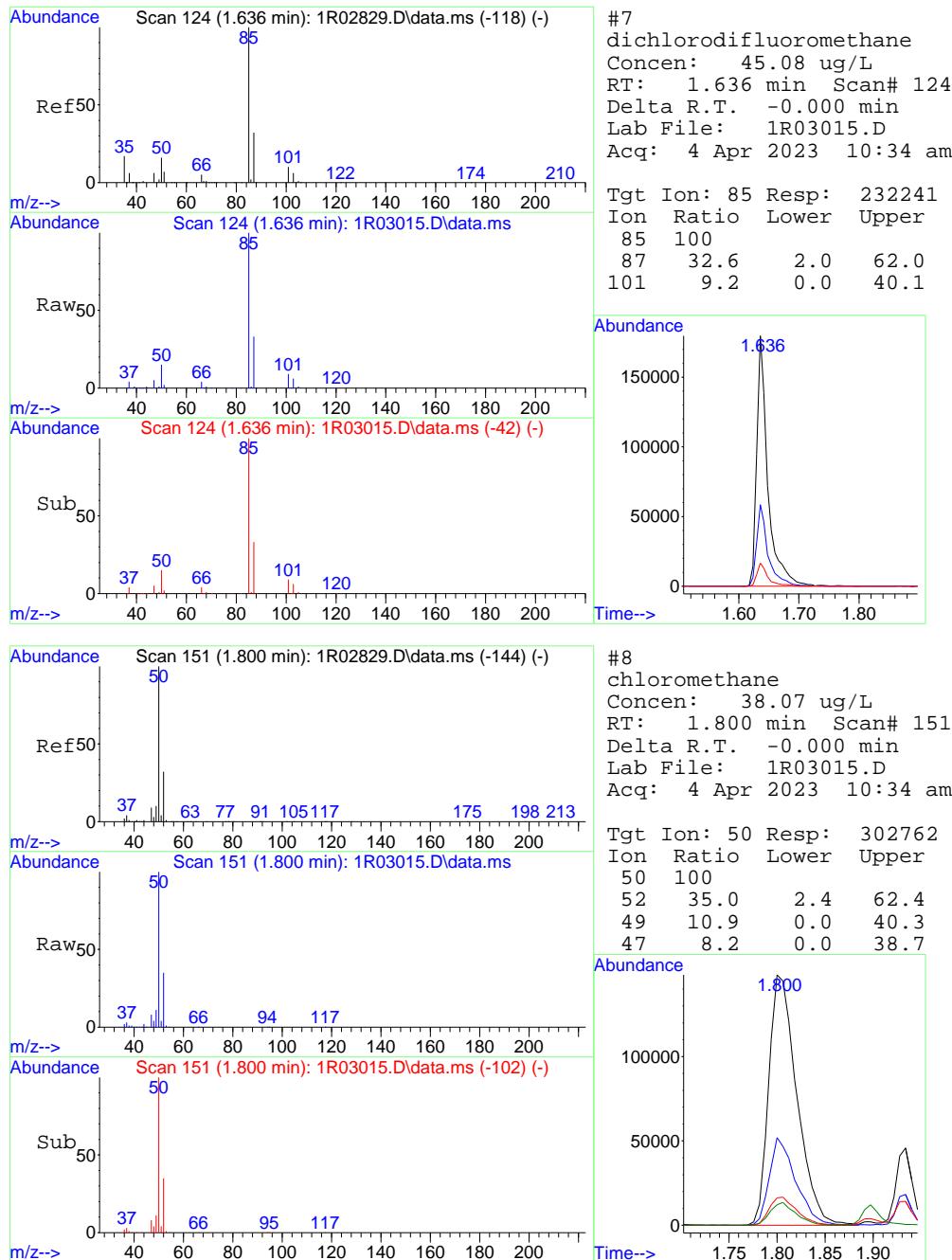
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Data File : 1R03015.D  
Acq On : 4 Apr 2023 10:34 am  
Operator : nickw  
Sample : bs  
Misc : MS67909,V1R0100,5,,,1  
ALS Vial : 4 Sample Multiplier: 1

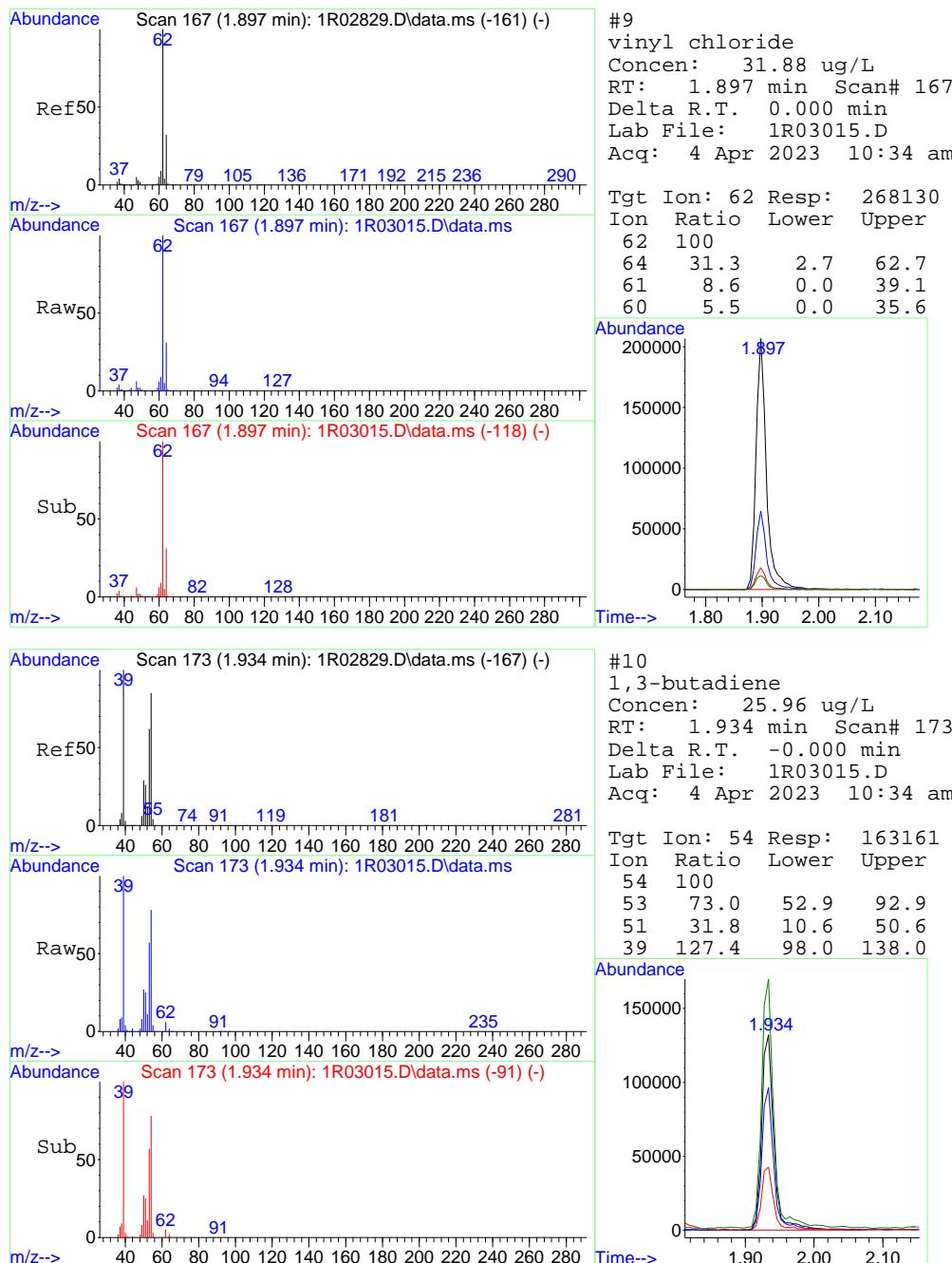
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QLast Update : Wed Mar 29 12:39:36 2023  
Response via : Initial Calibration

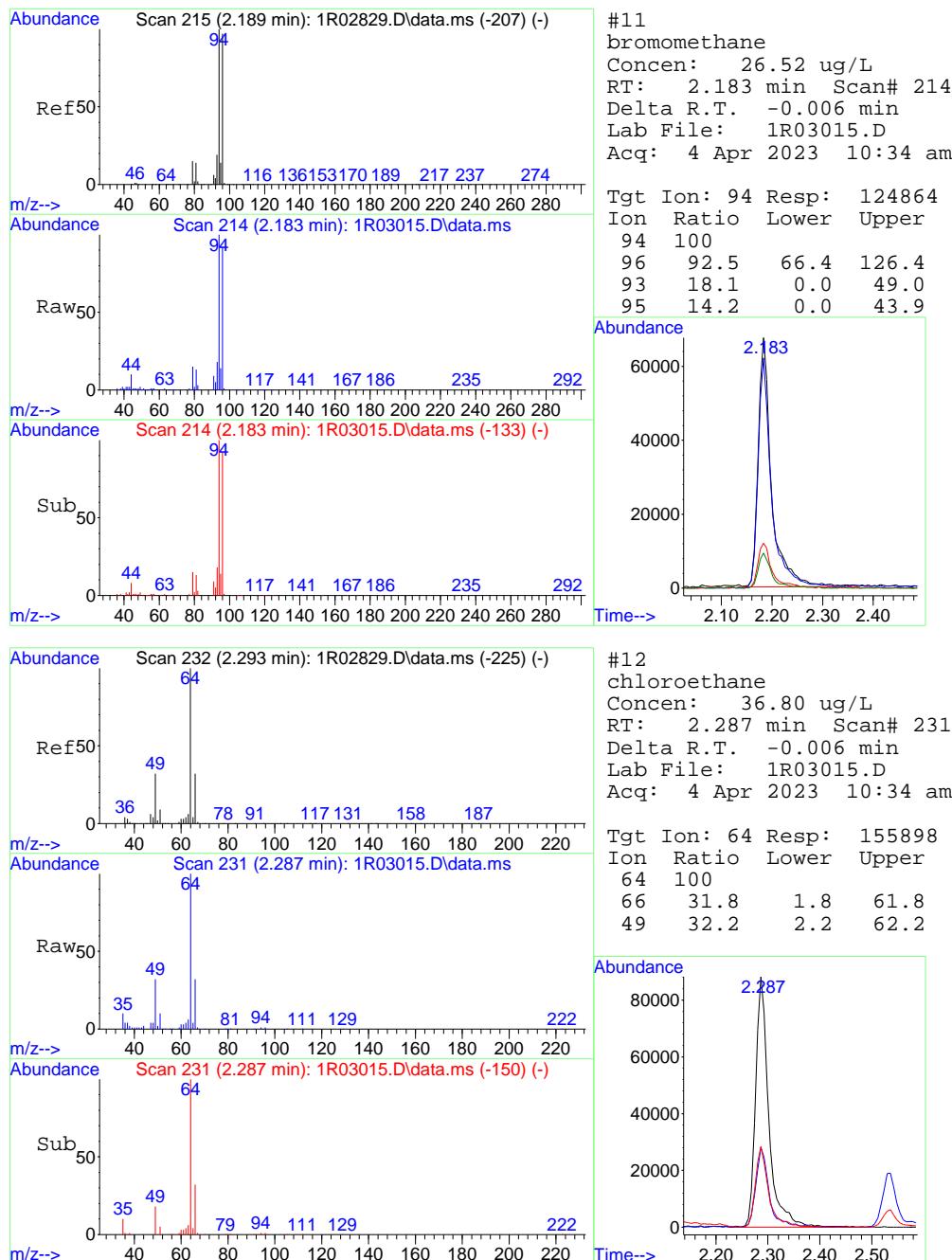


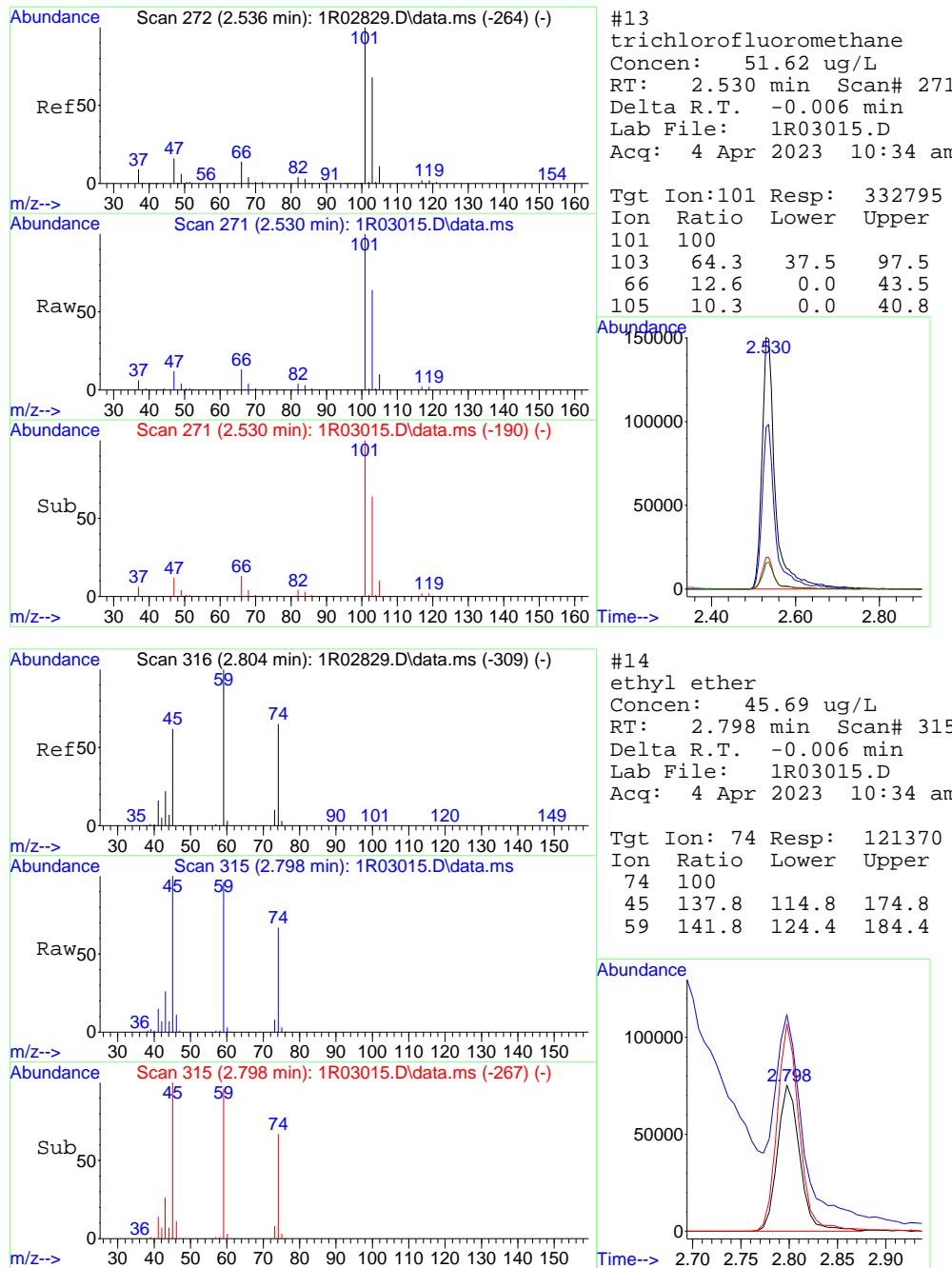


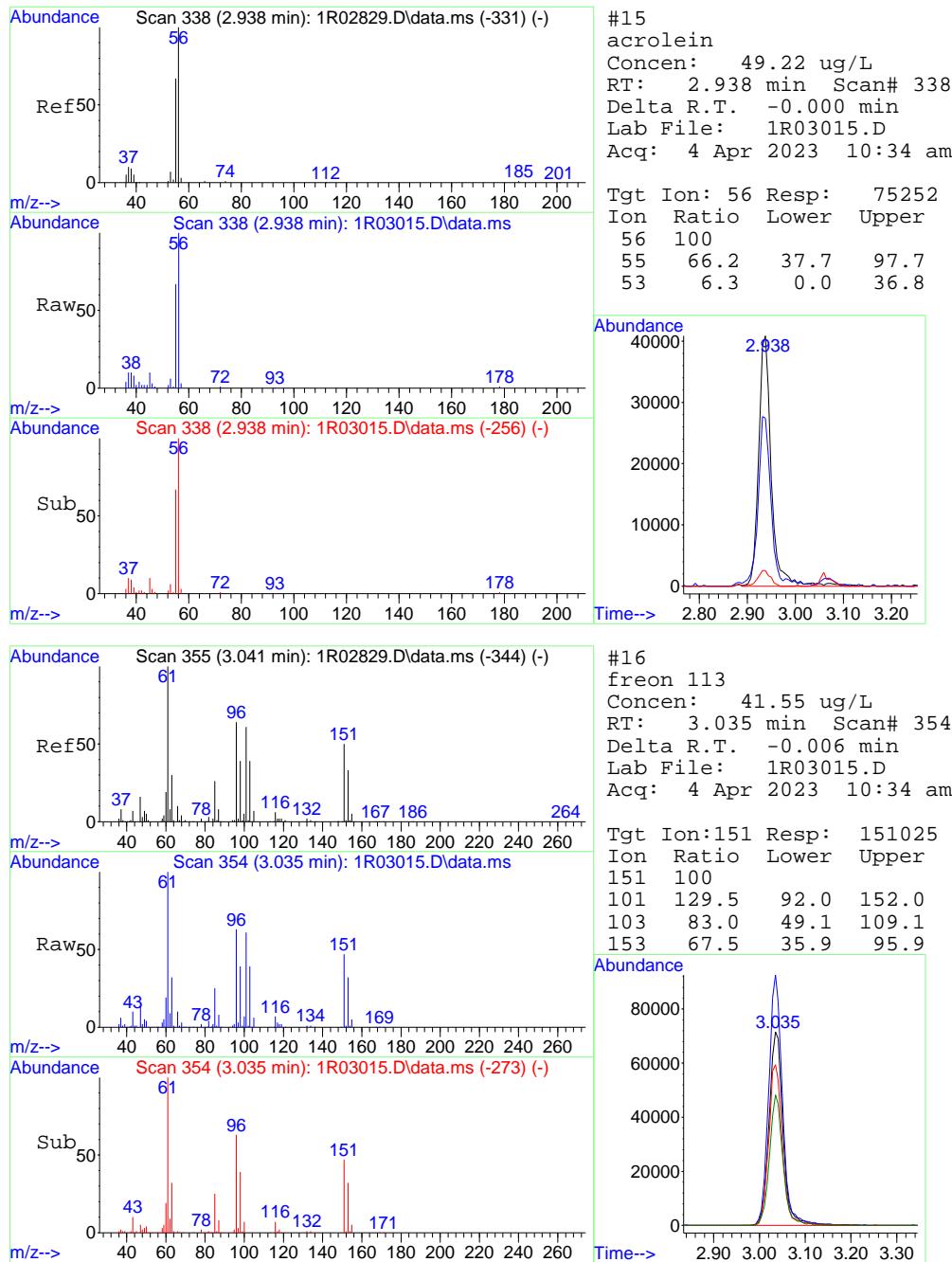


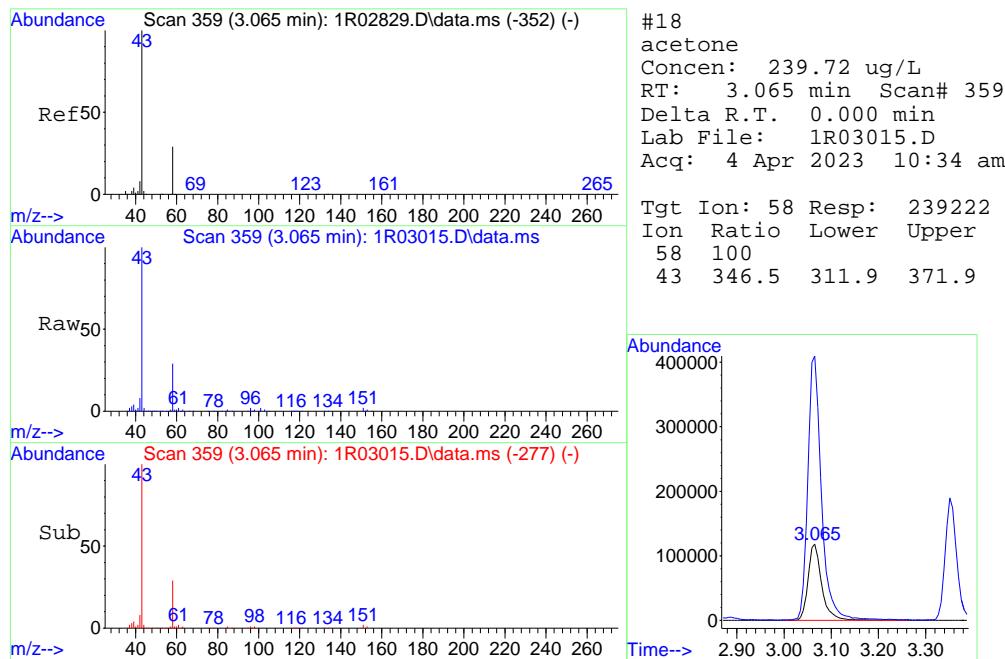
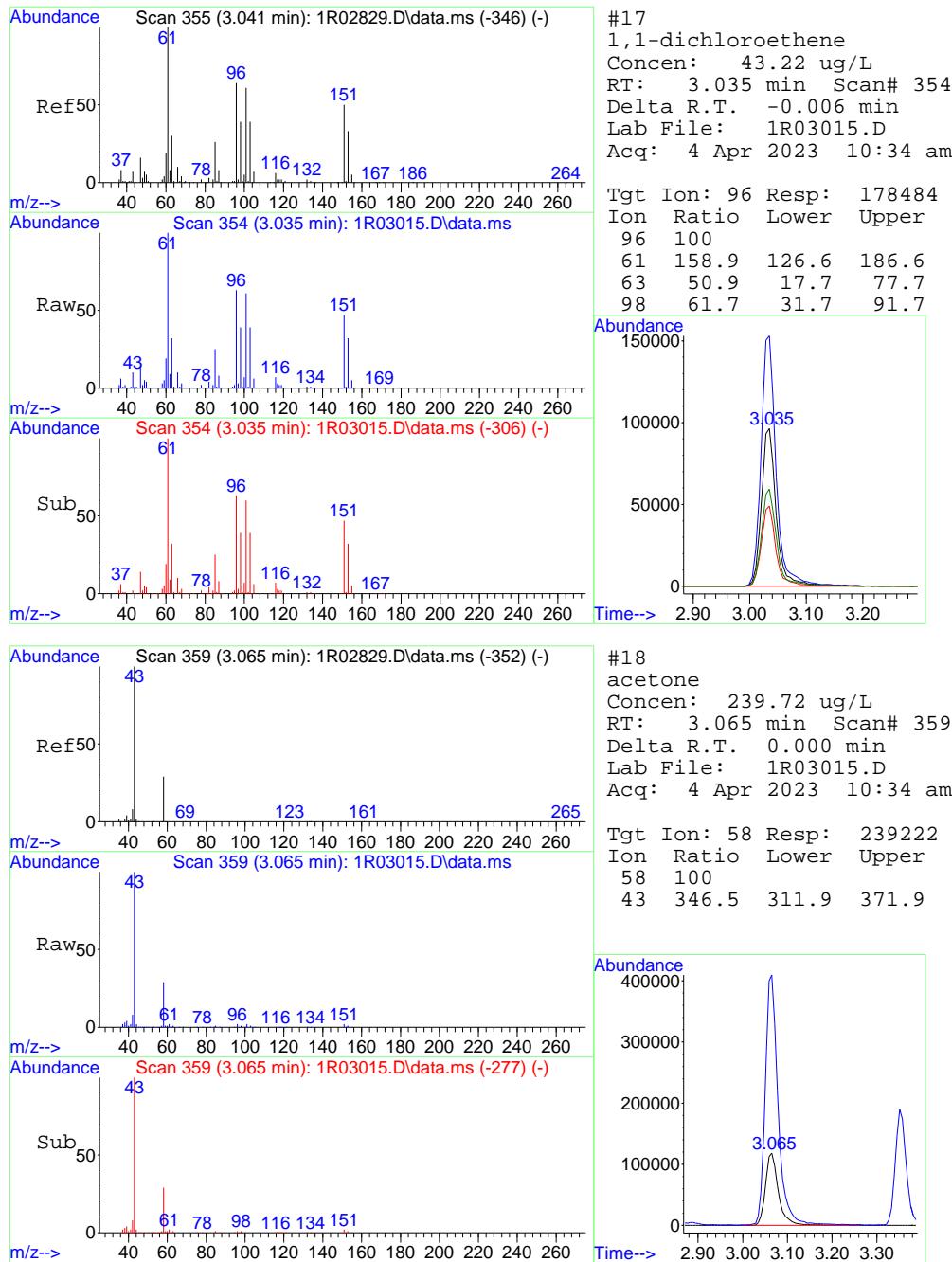


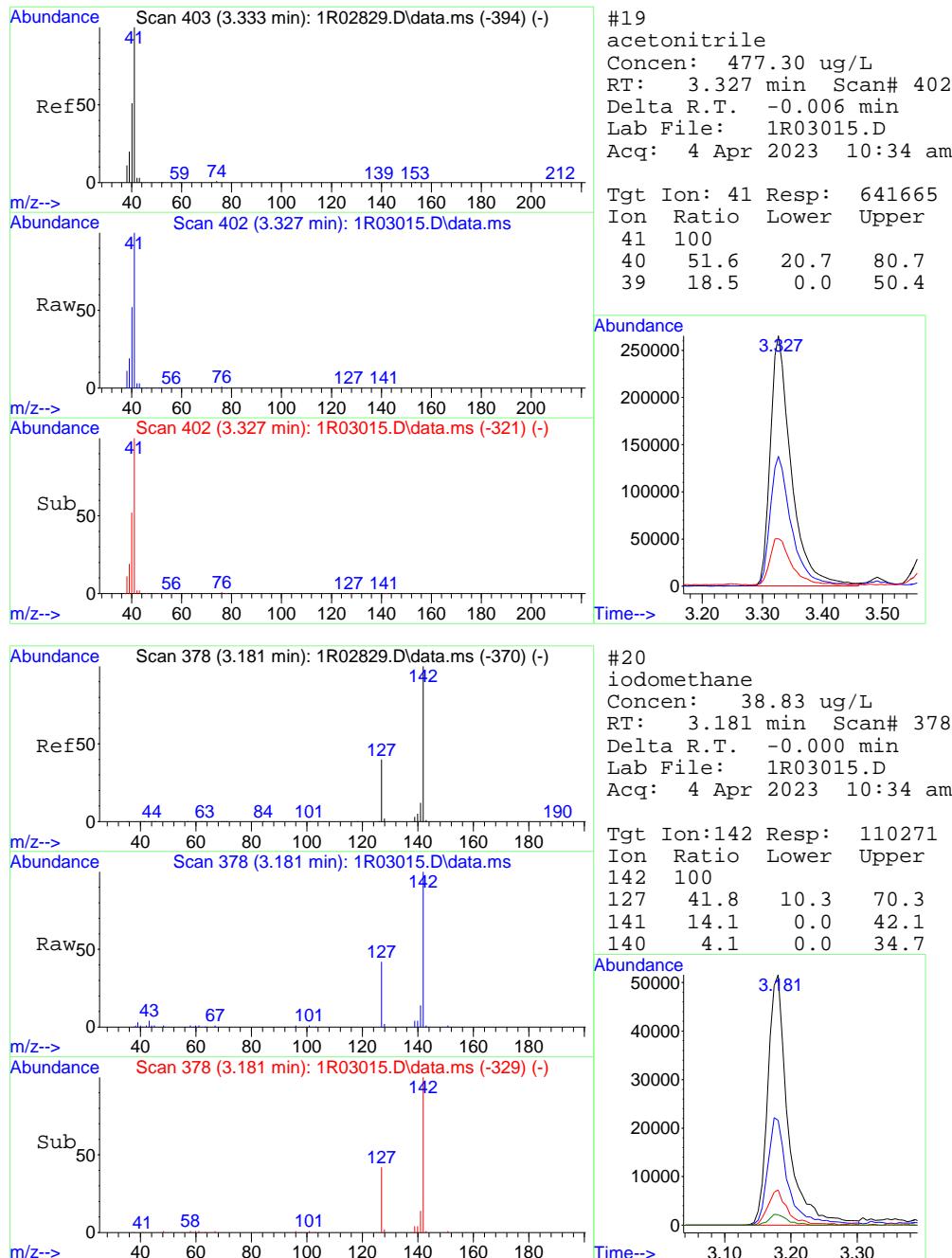


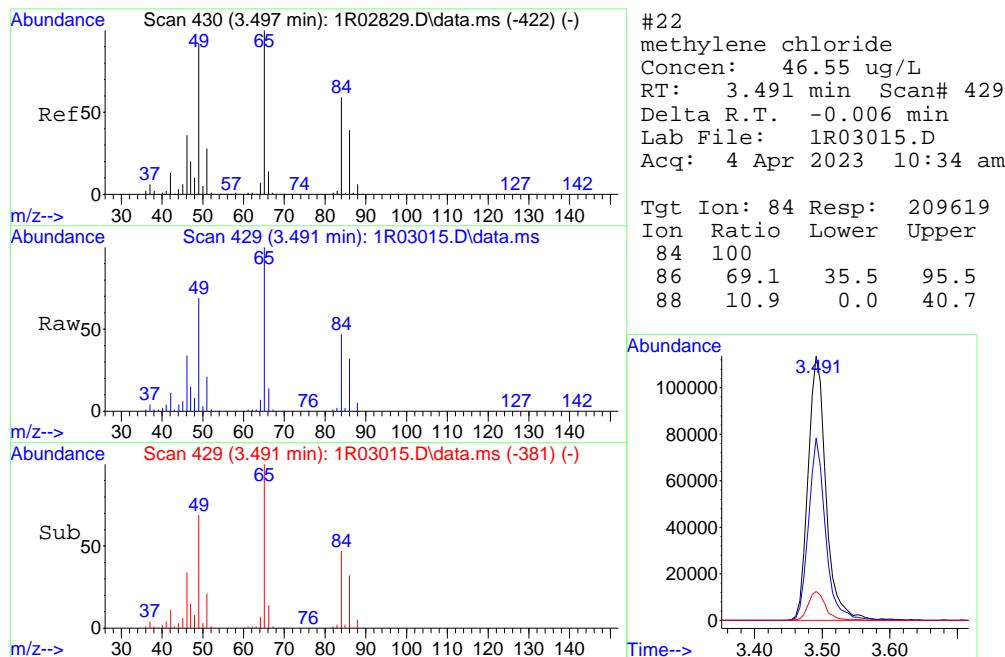
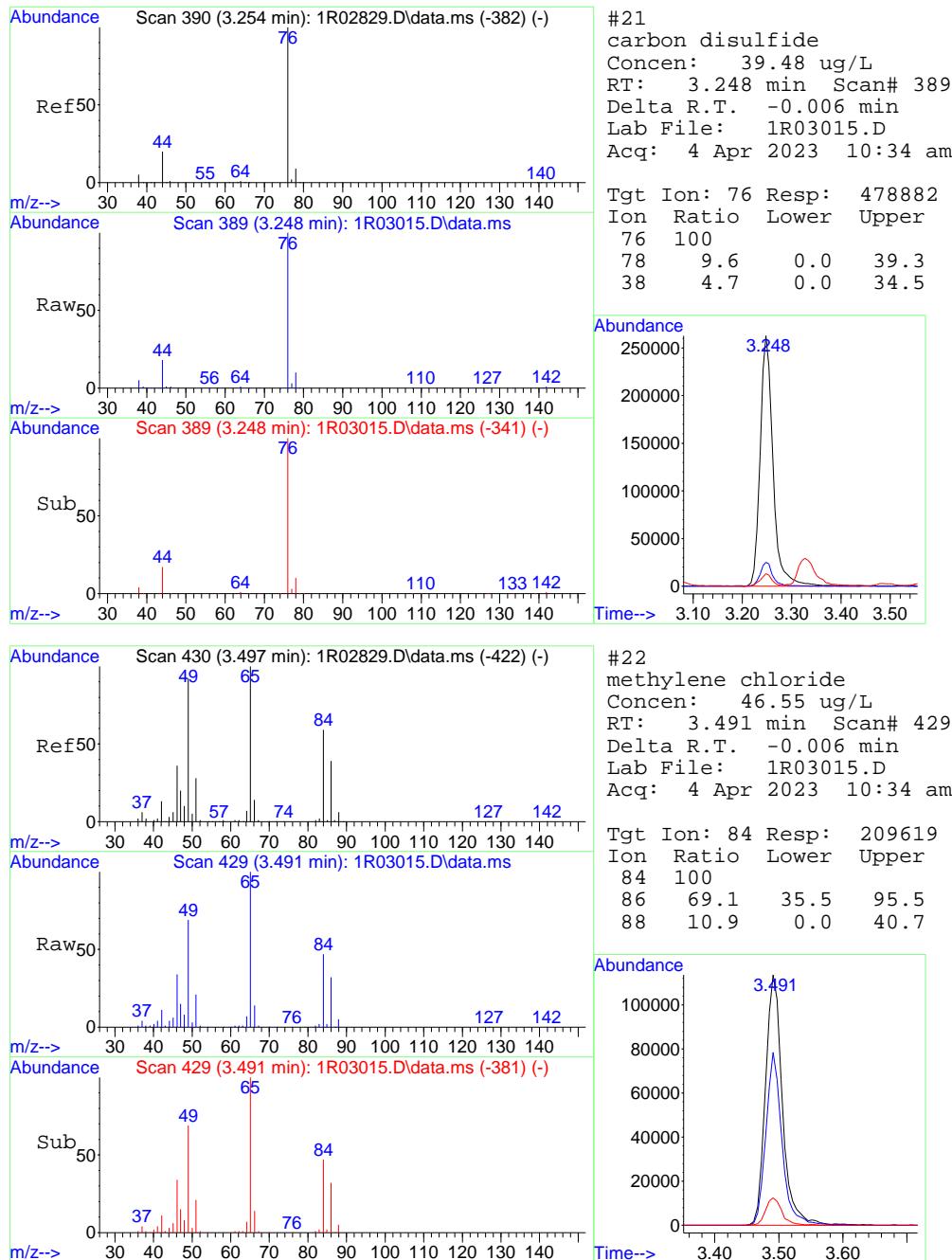


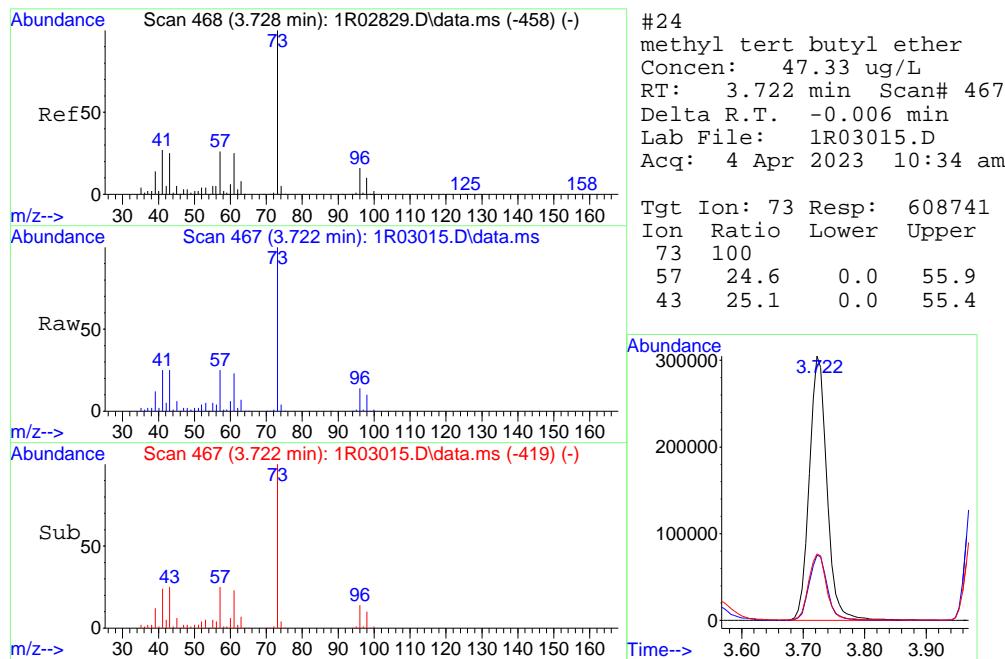
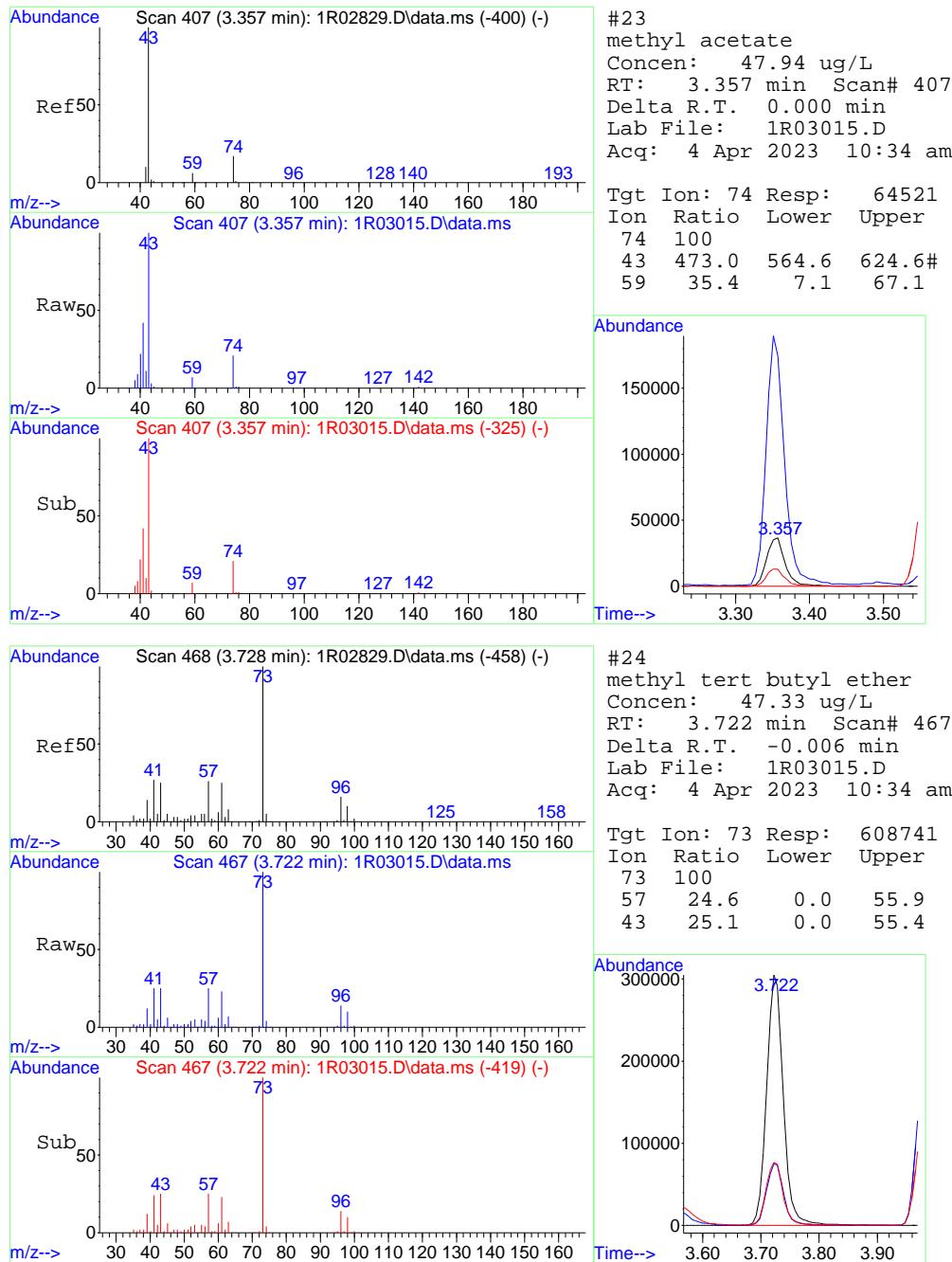


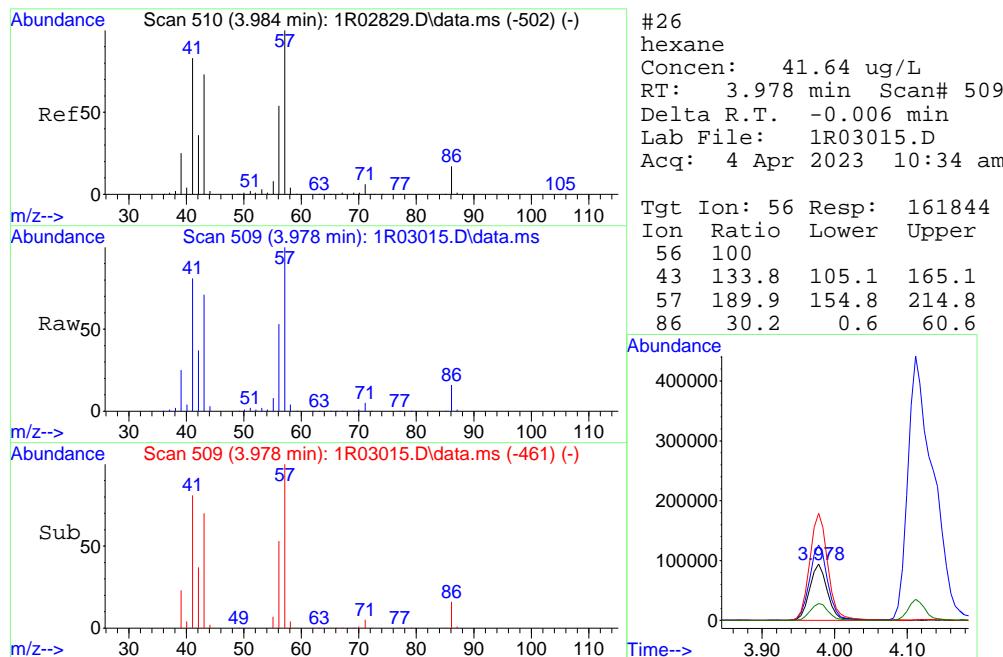
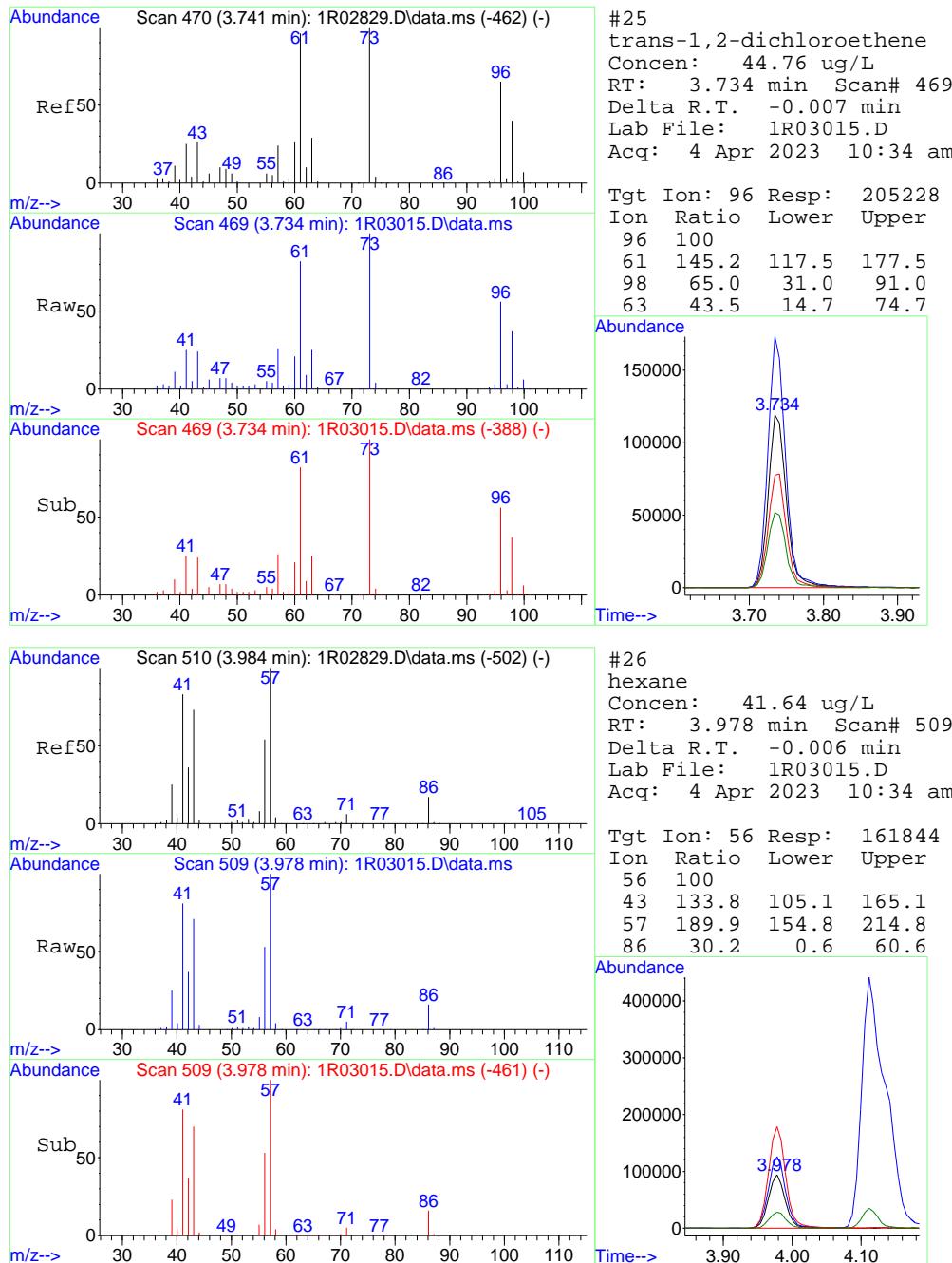


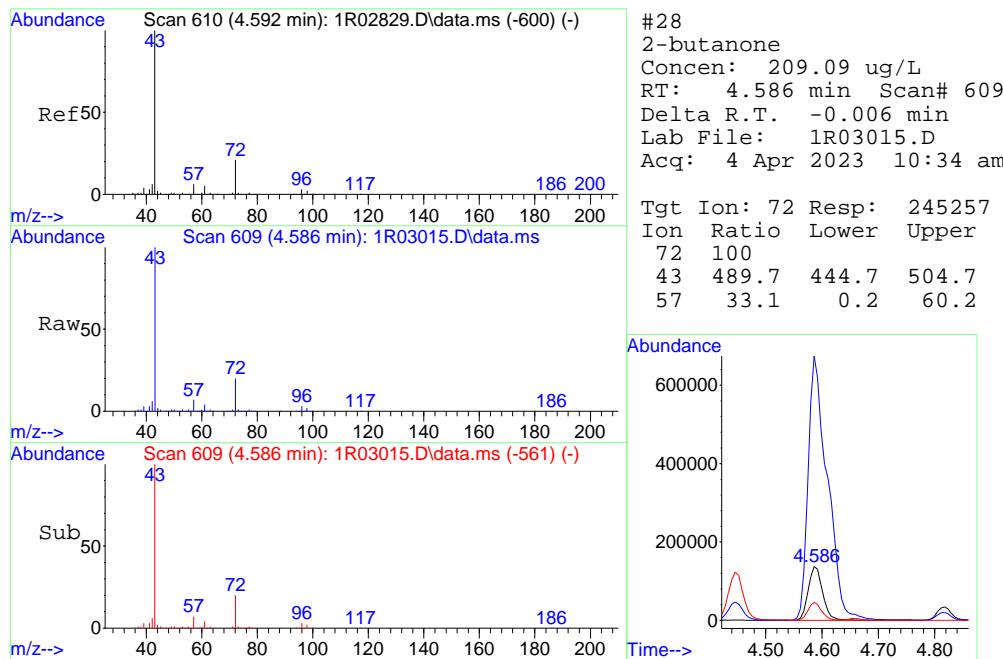
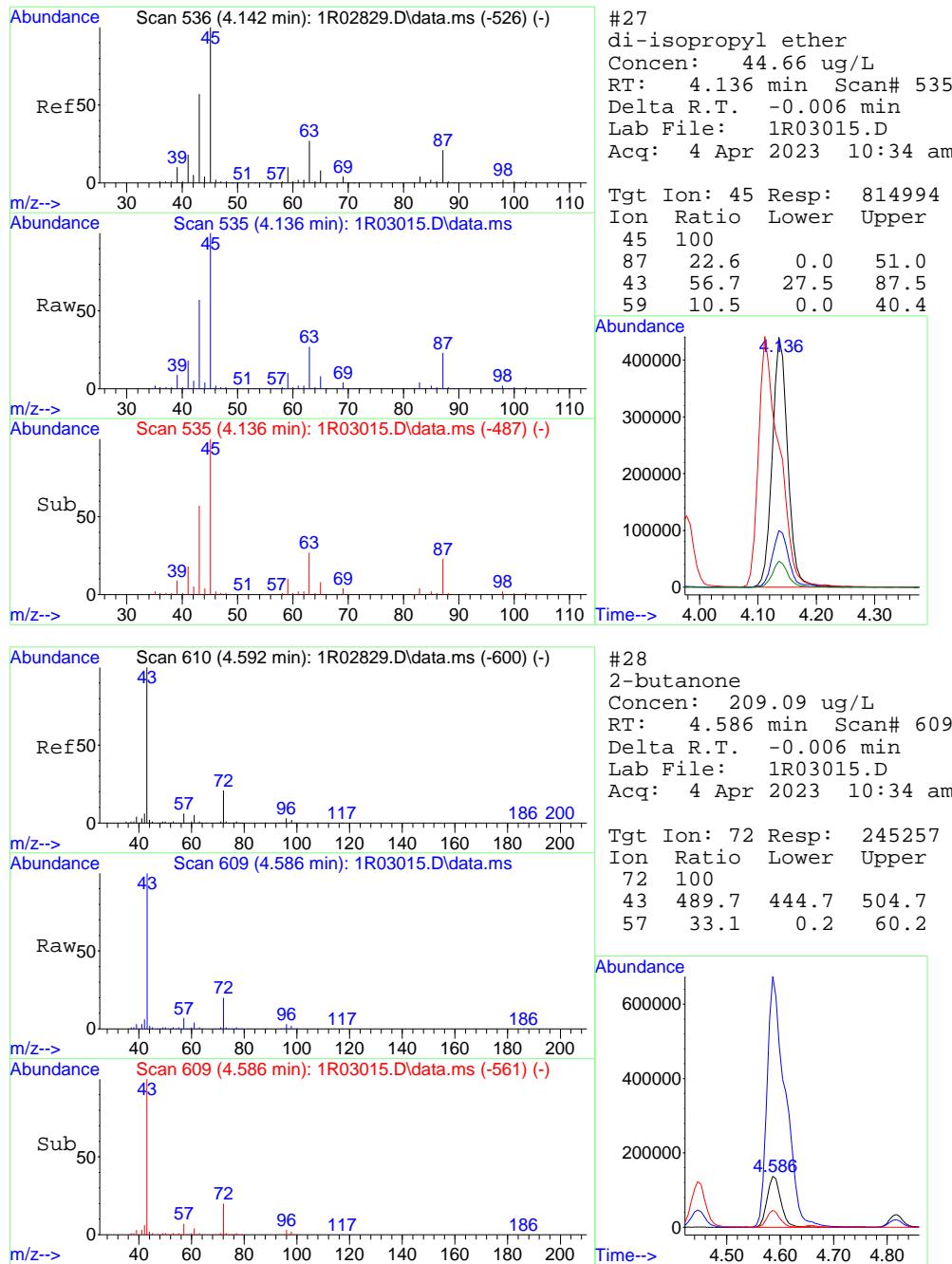


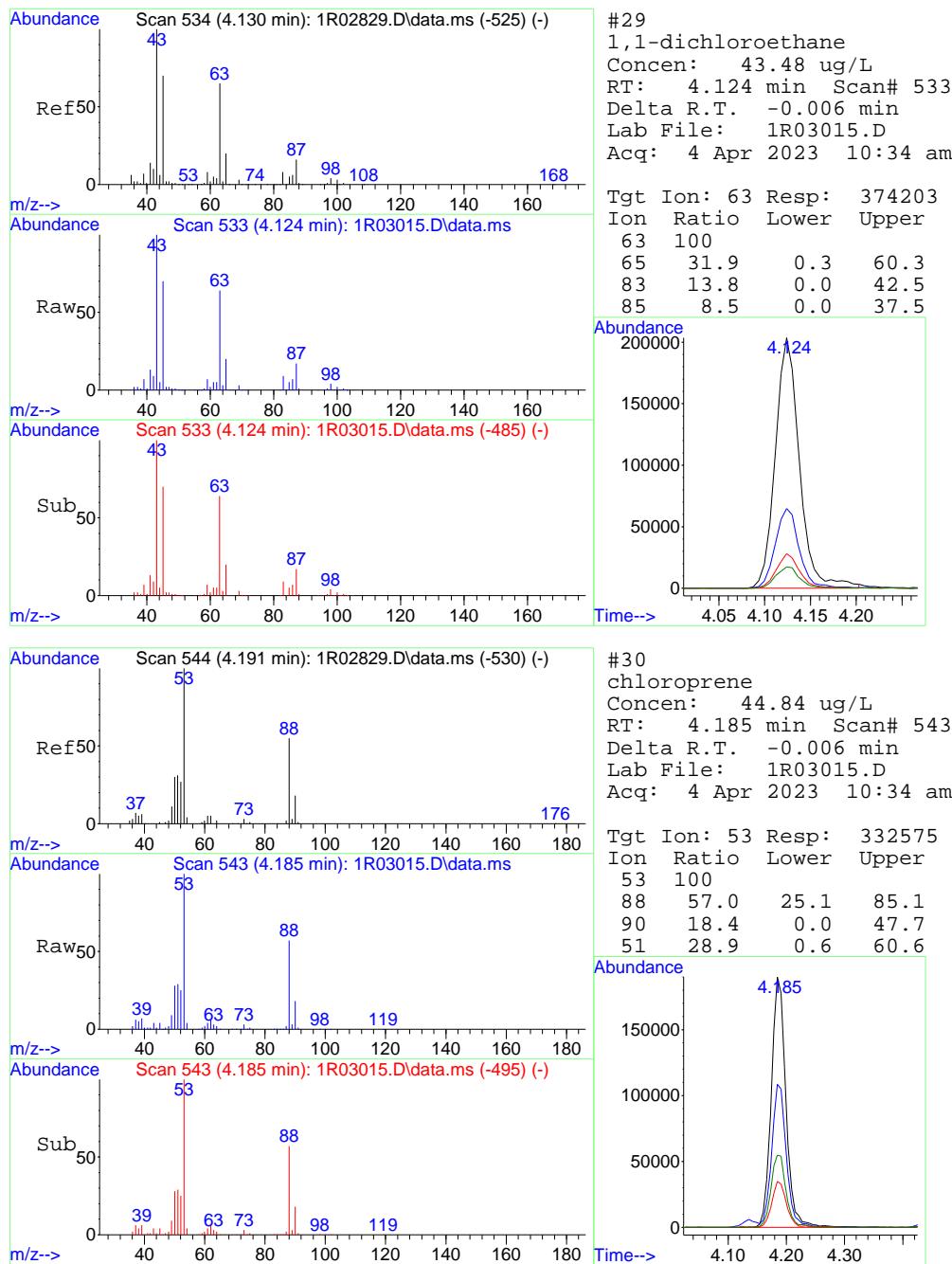


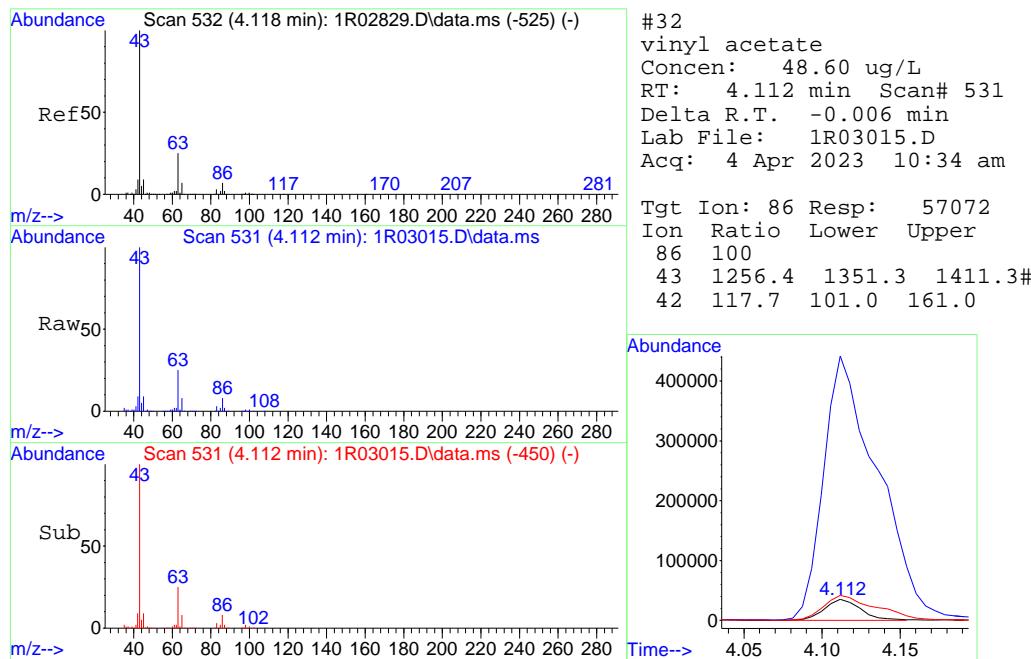
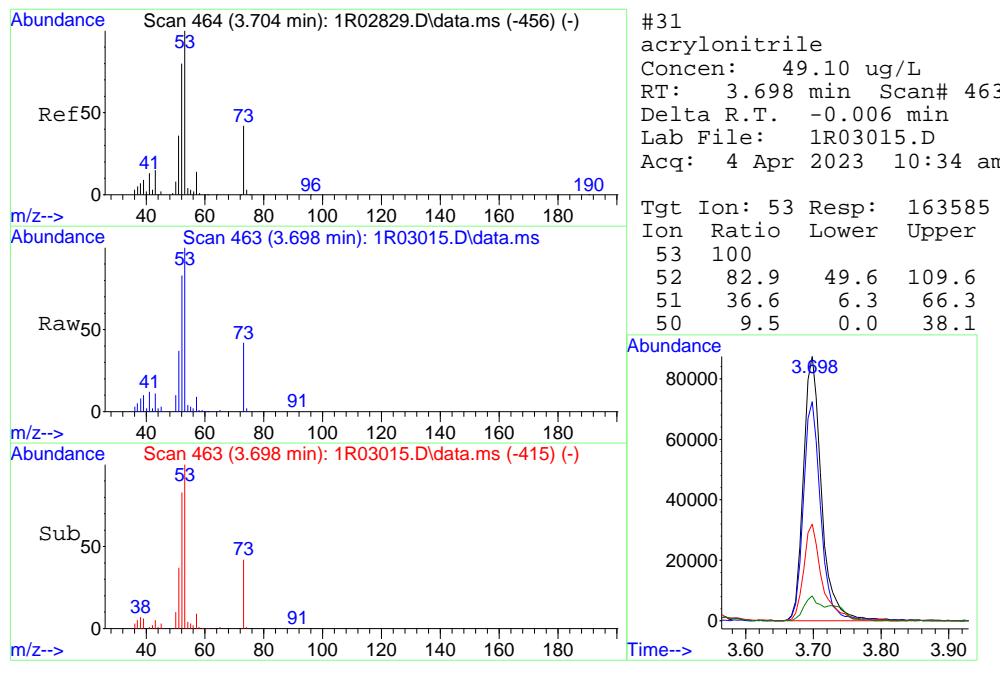


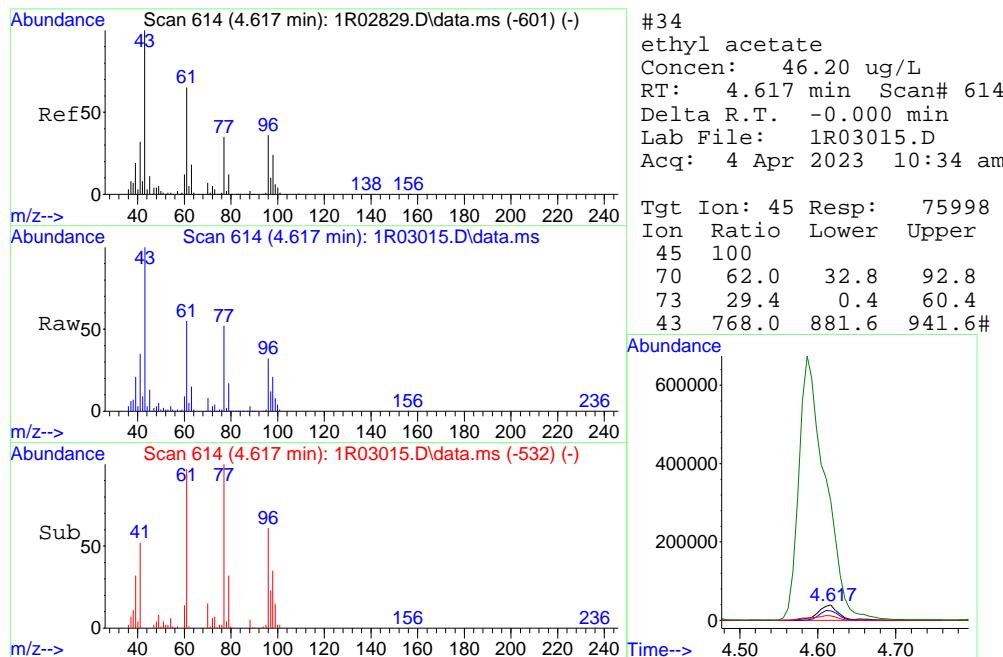
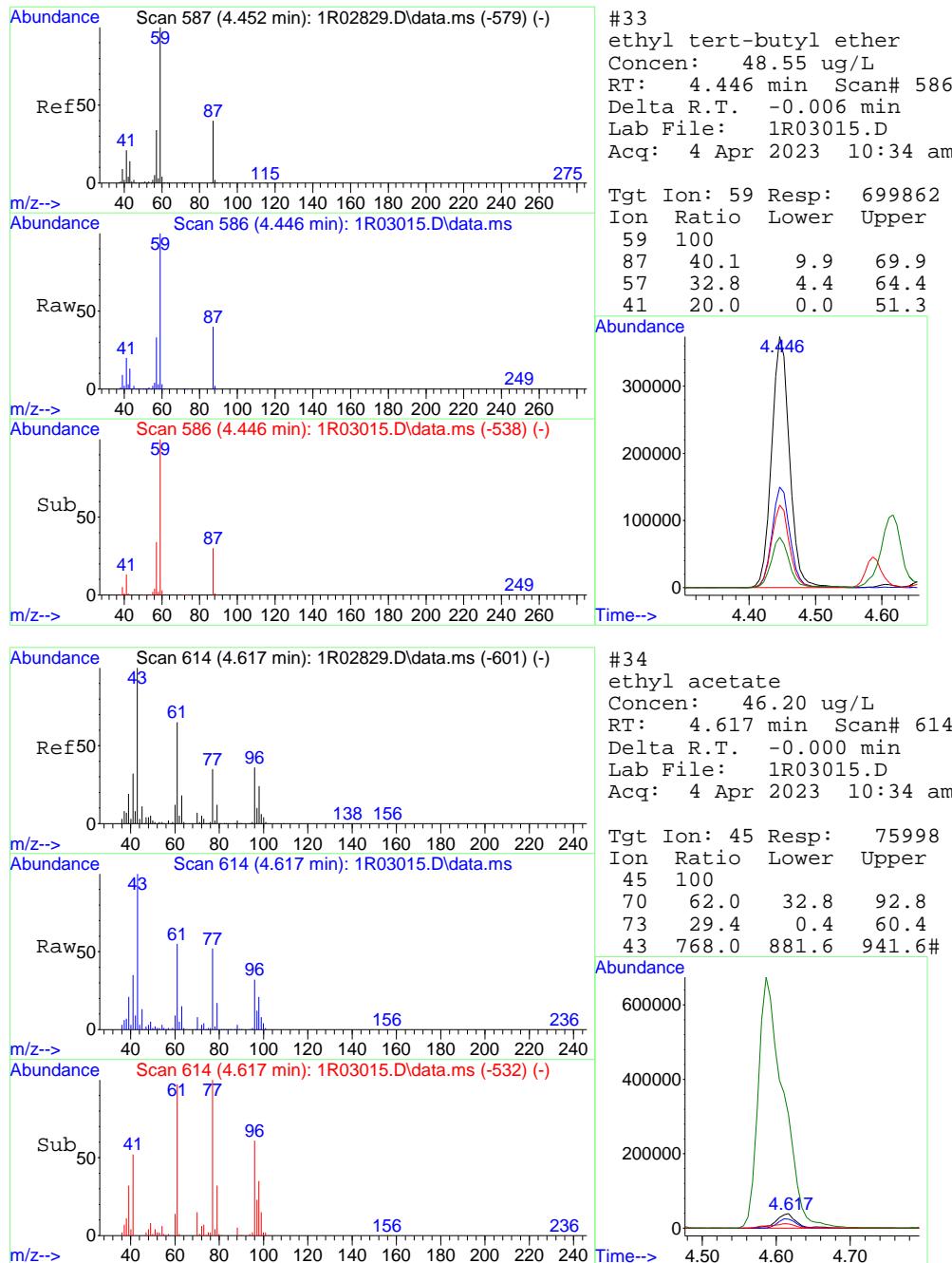


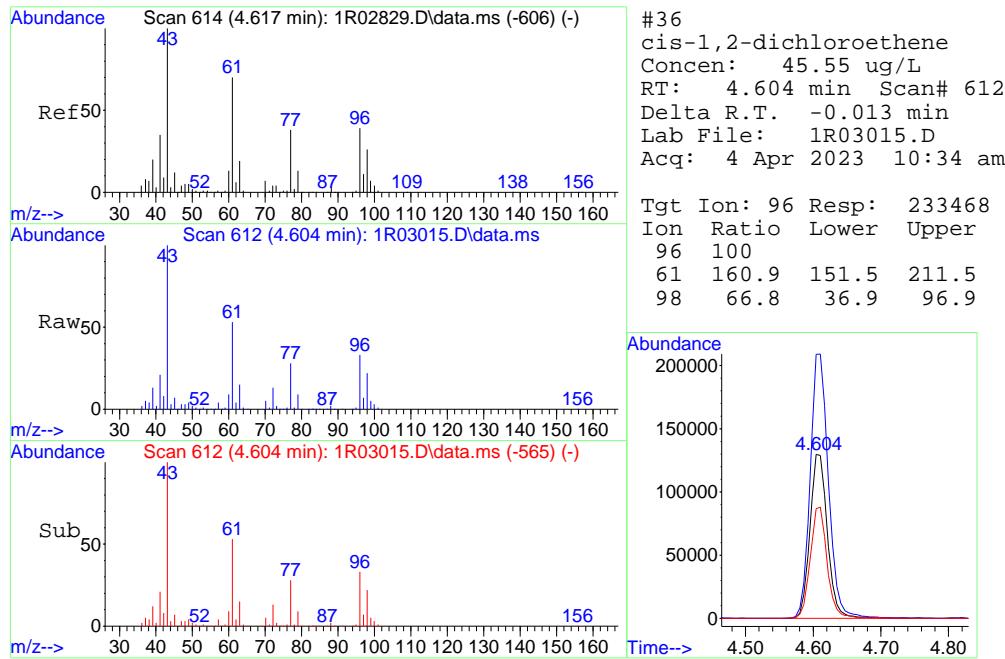
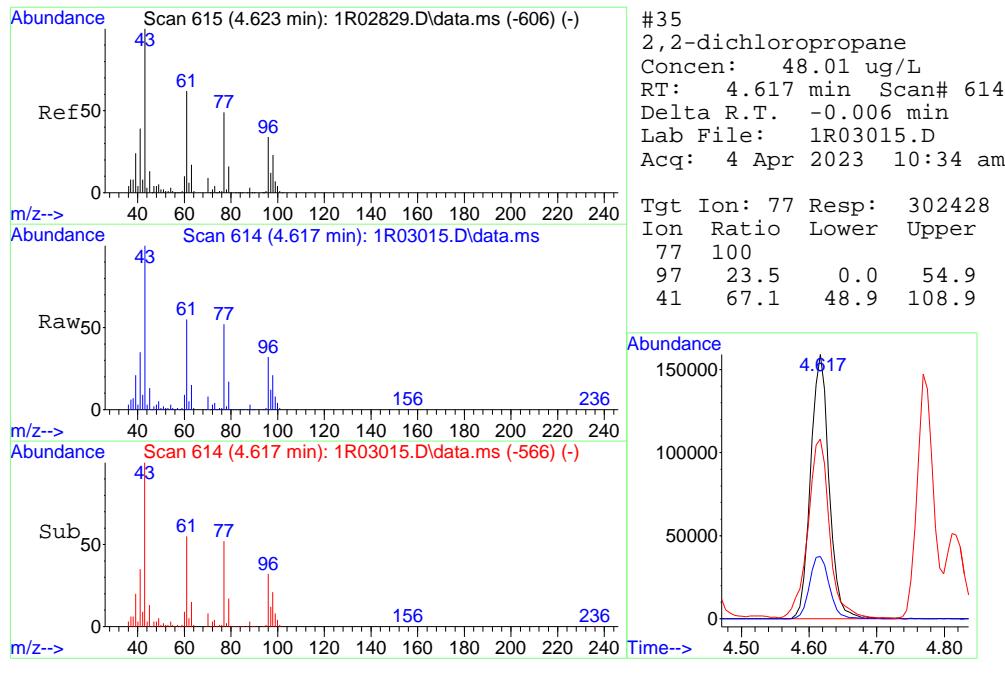


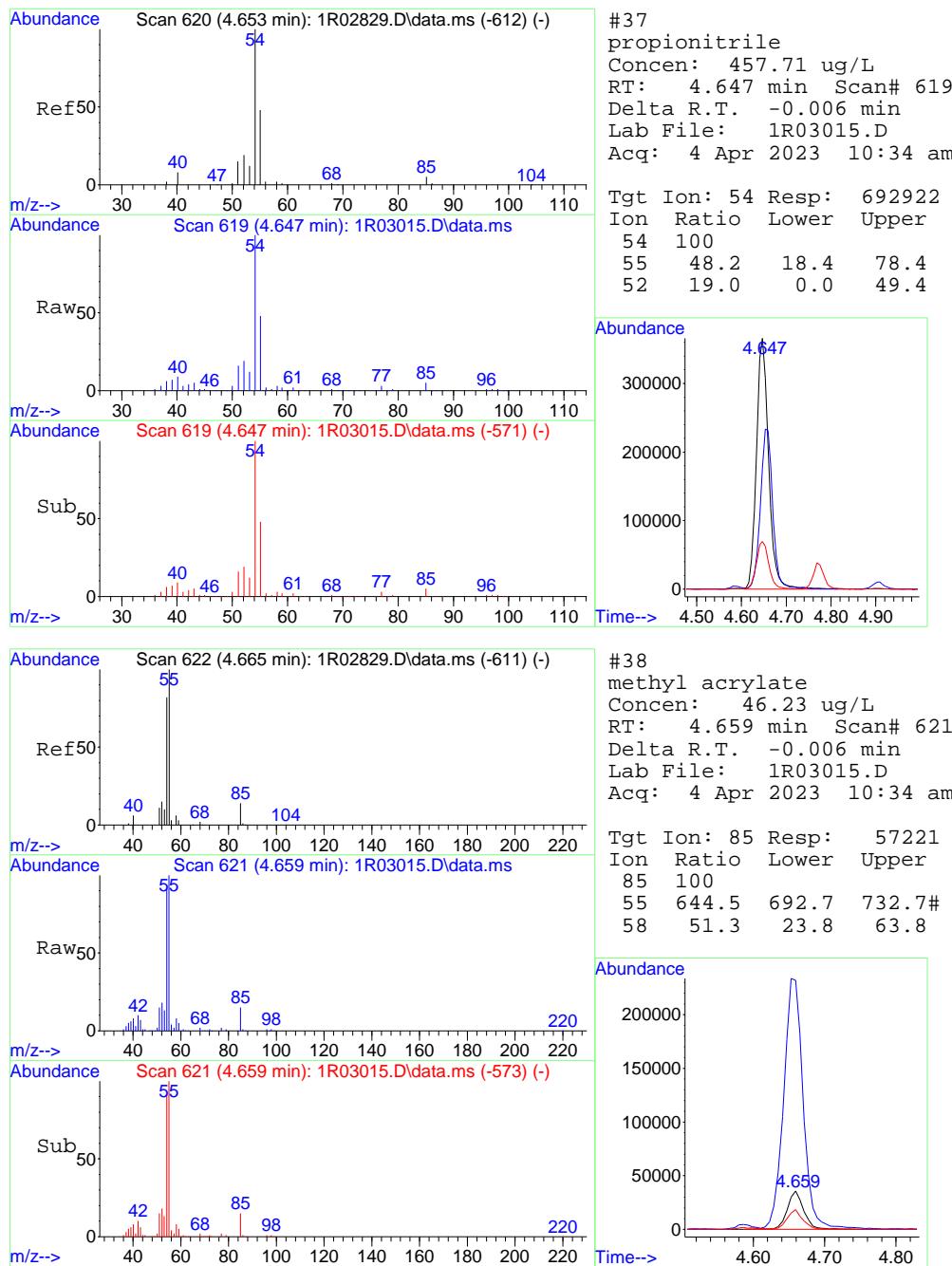


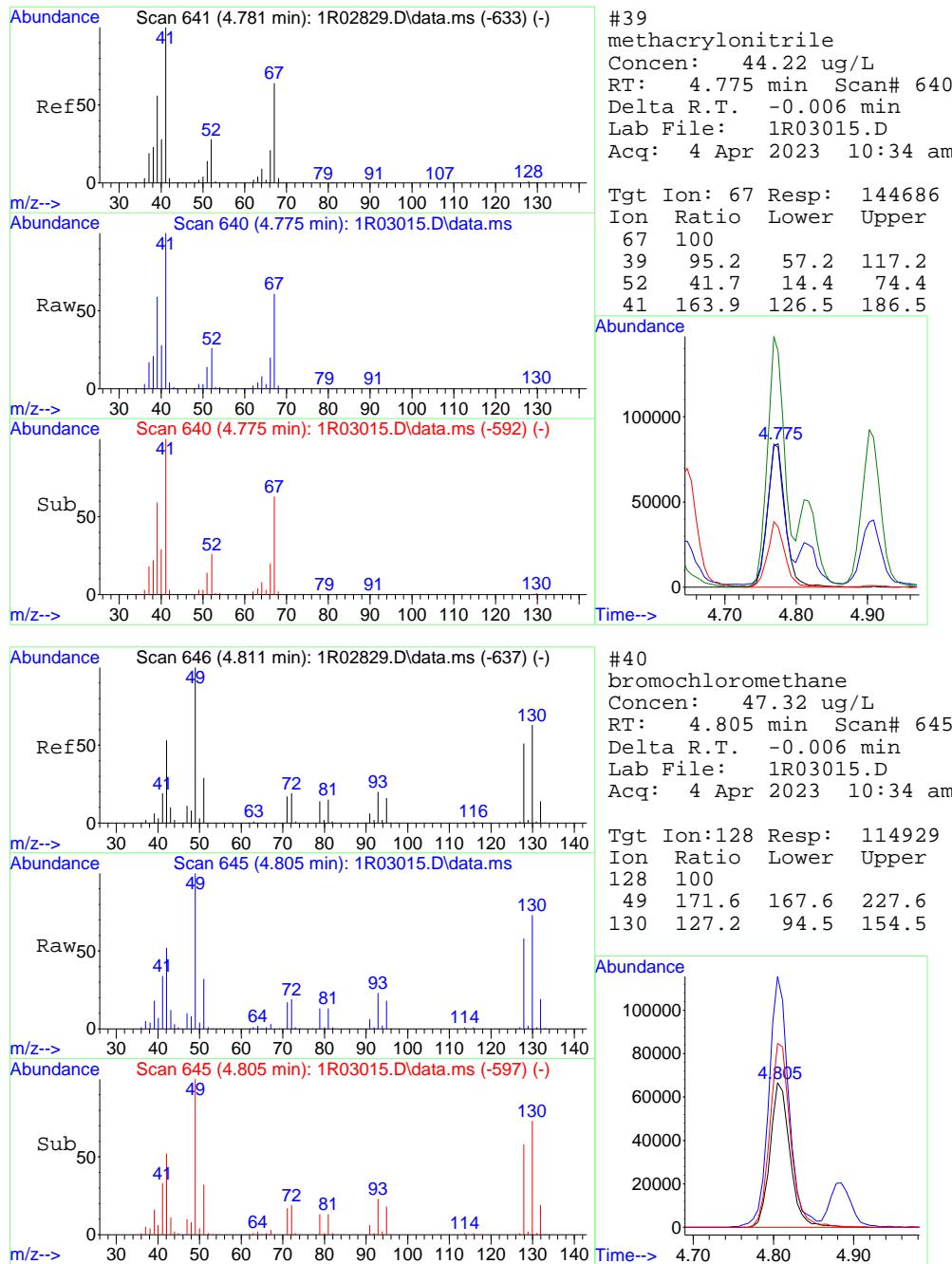


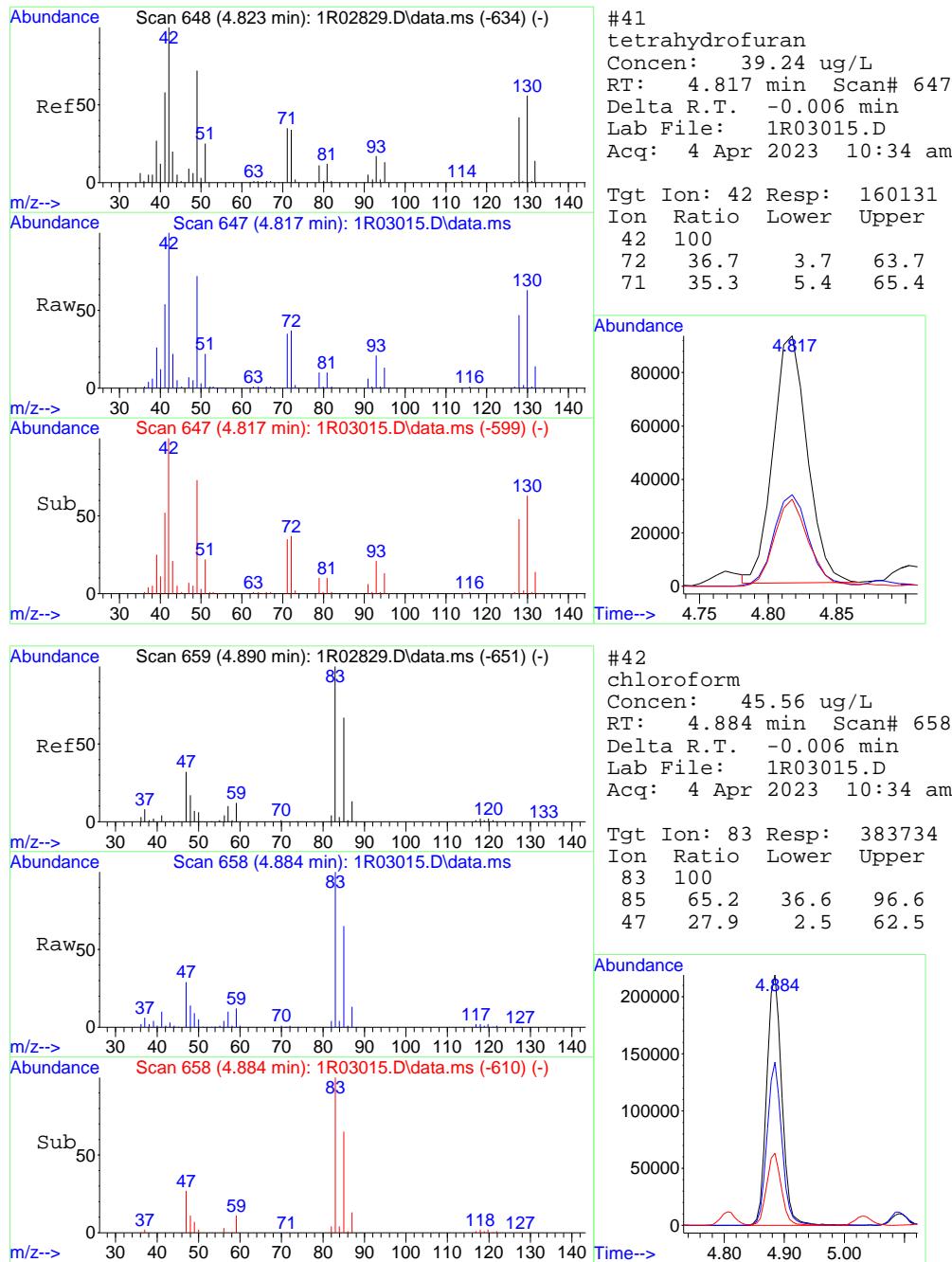


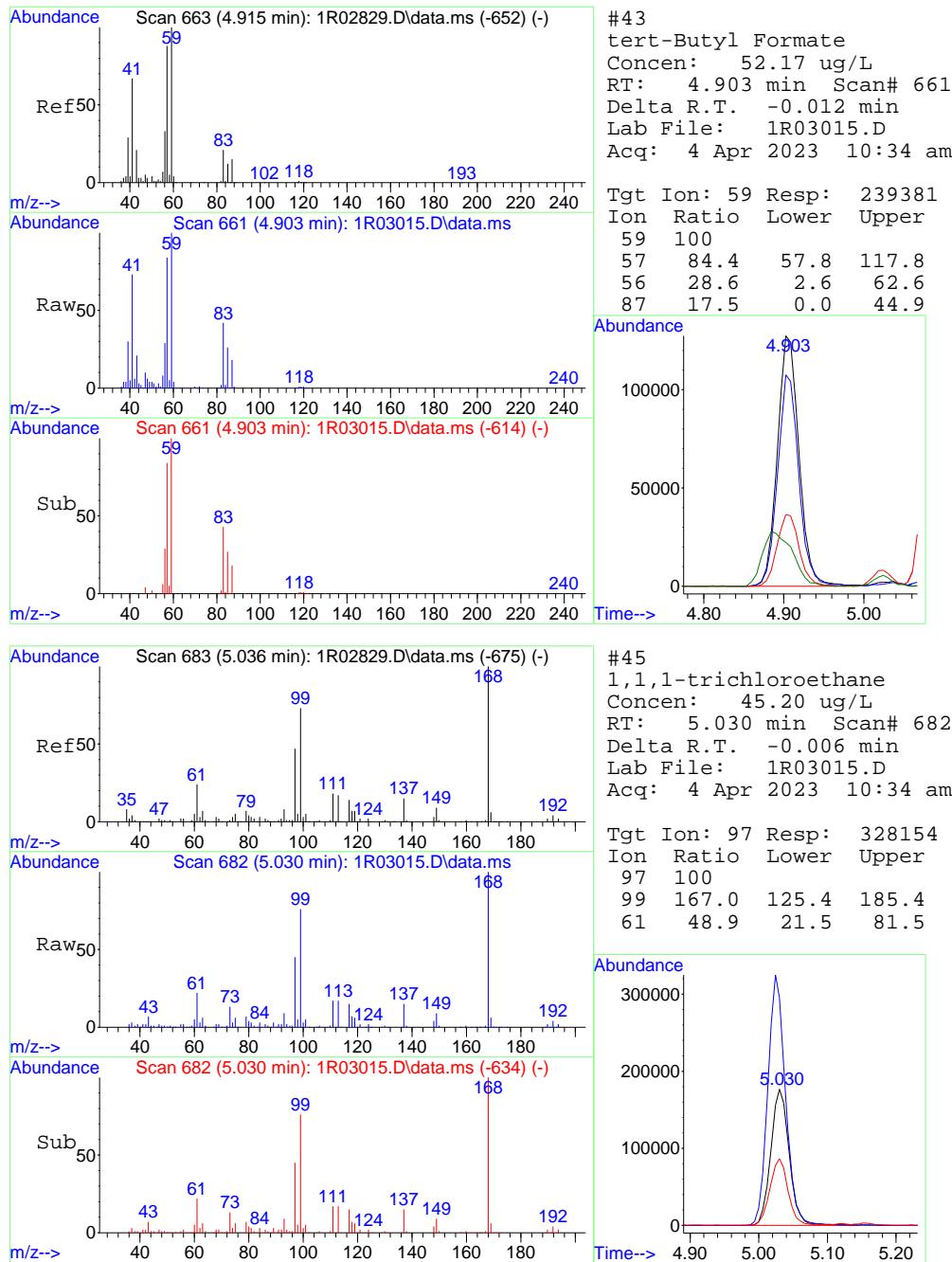


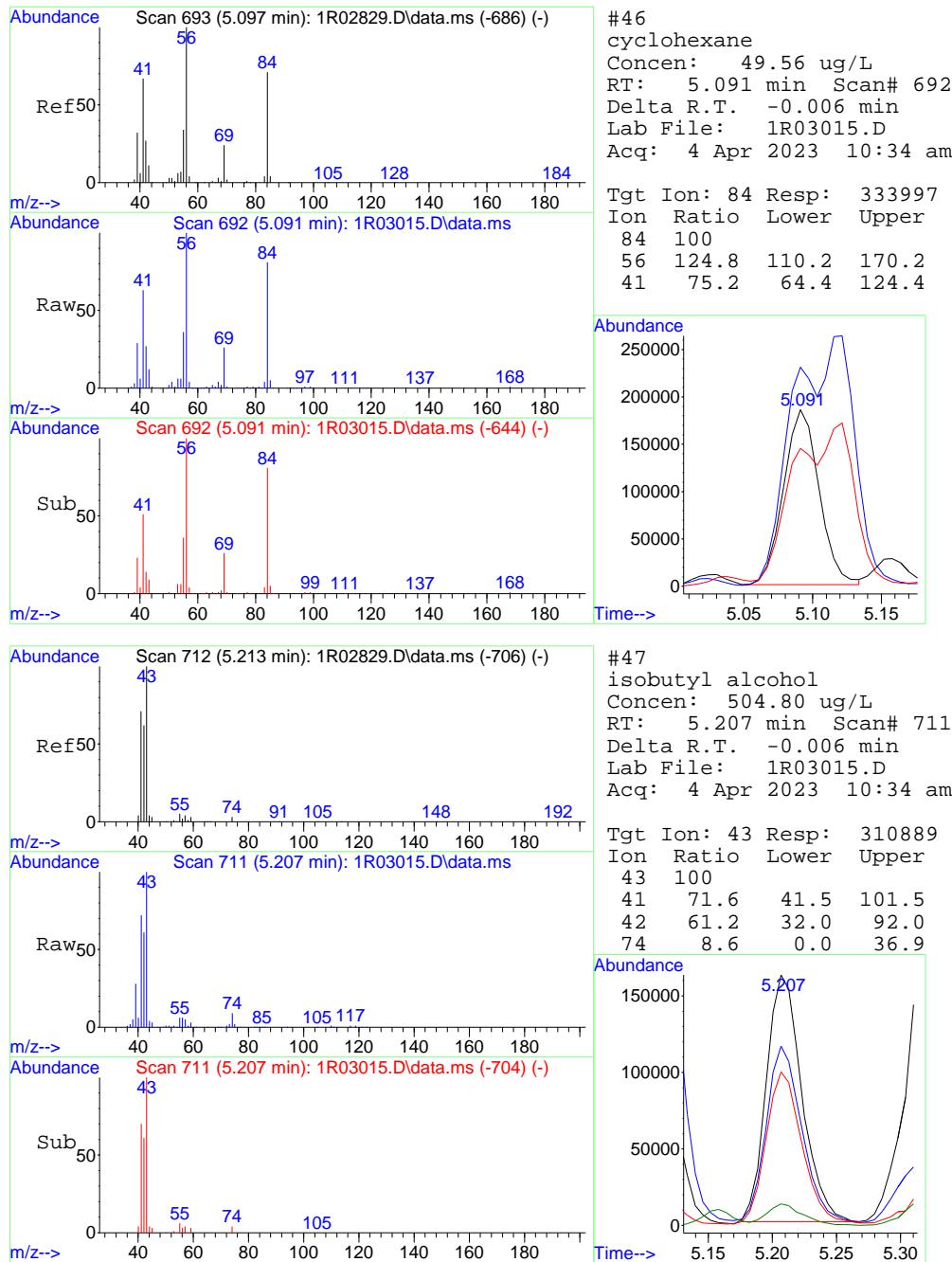


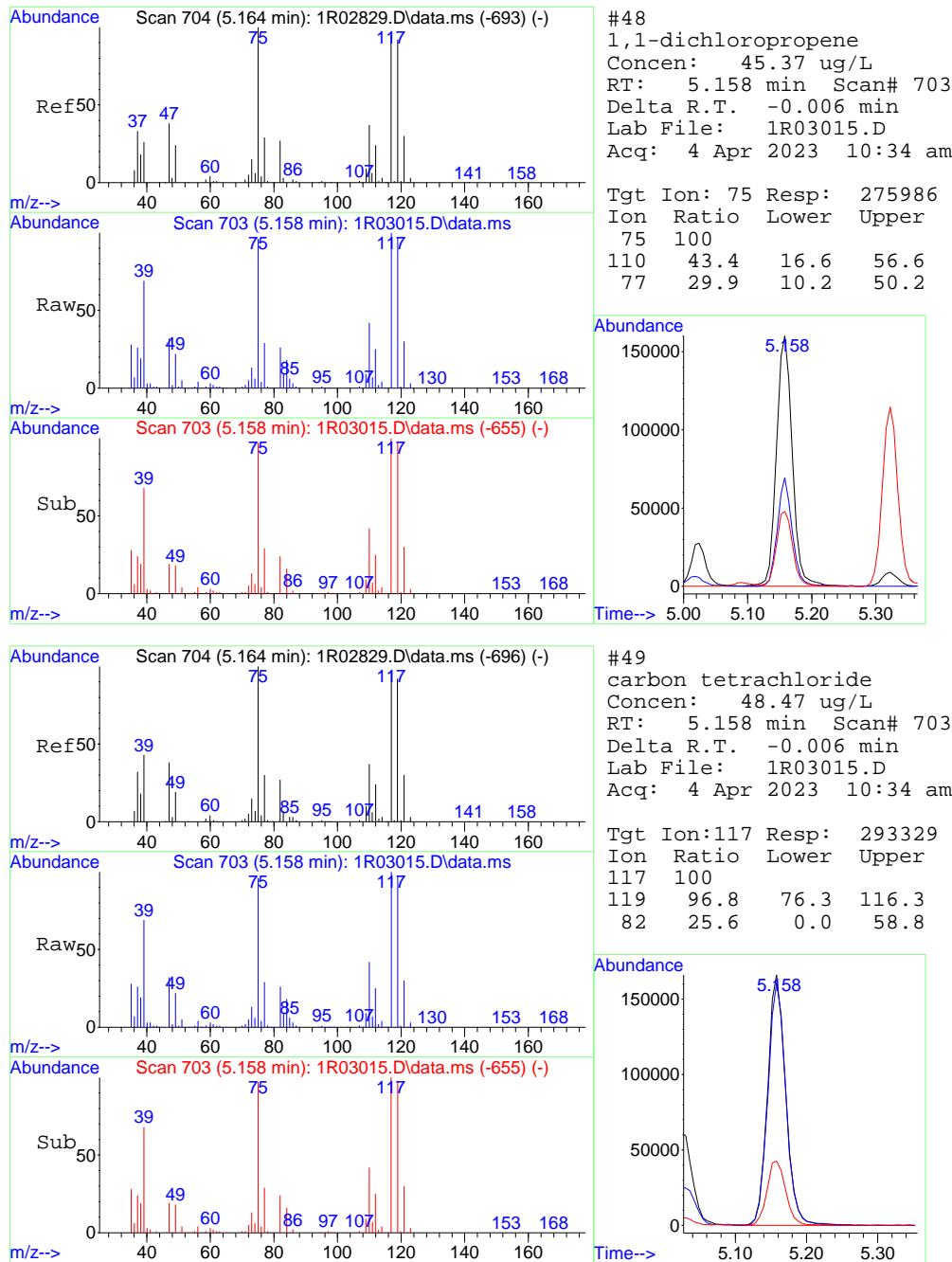


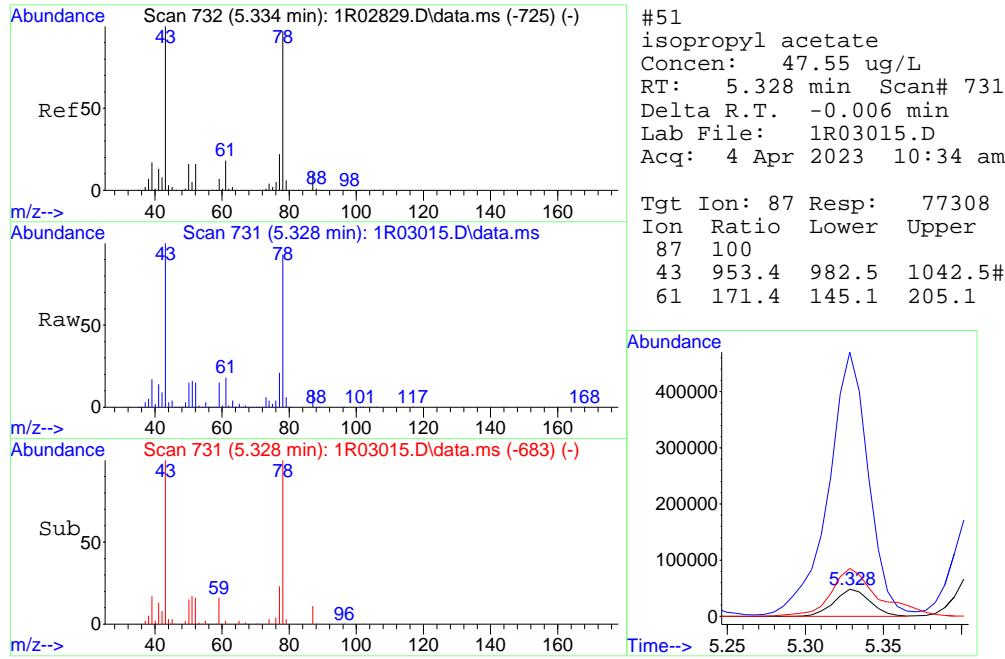
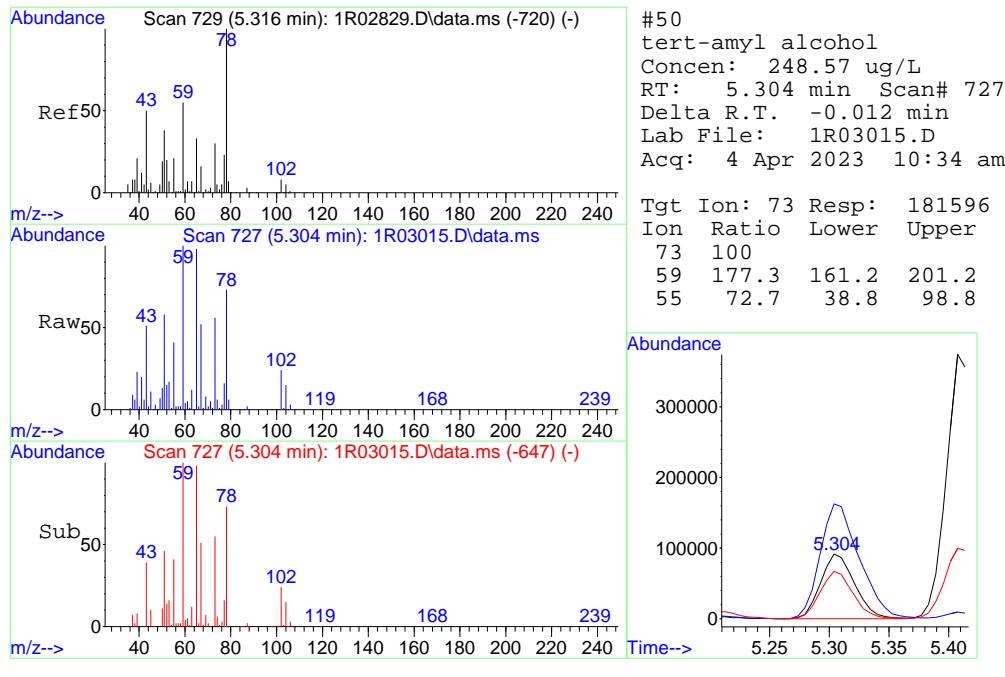


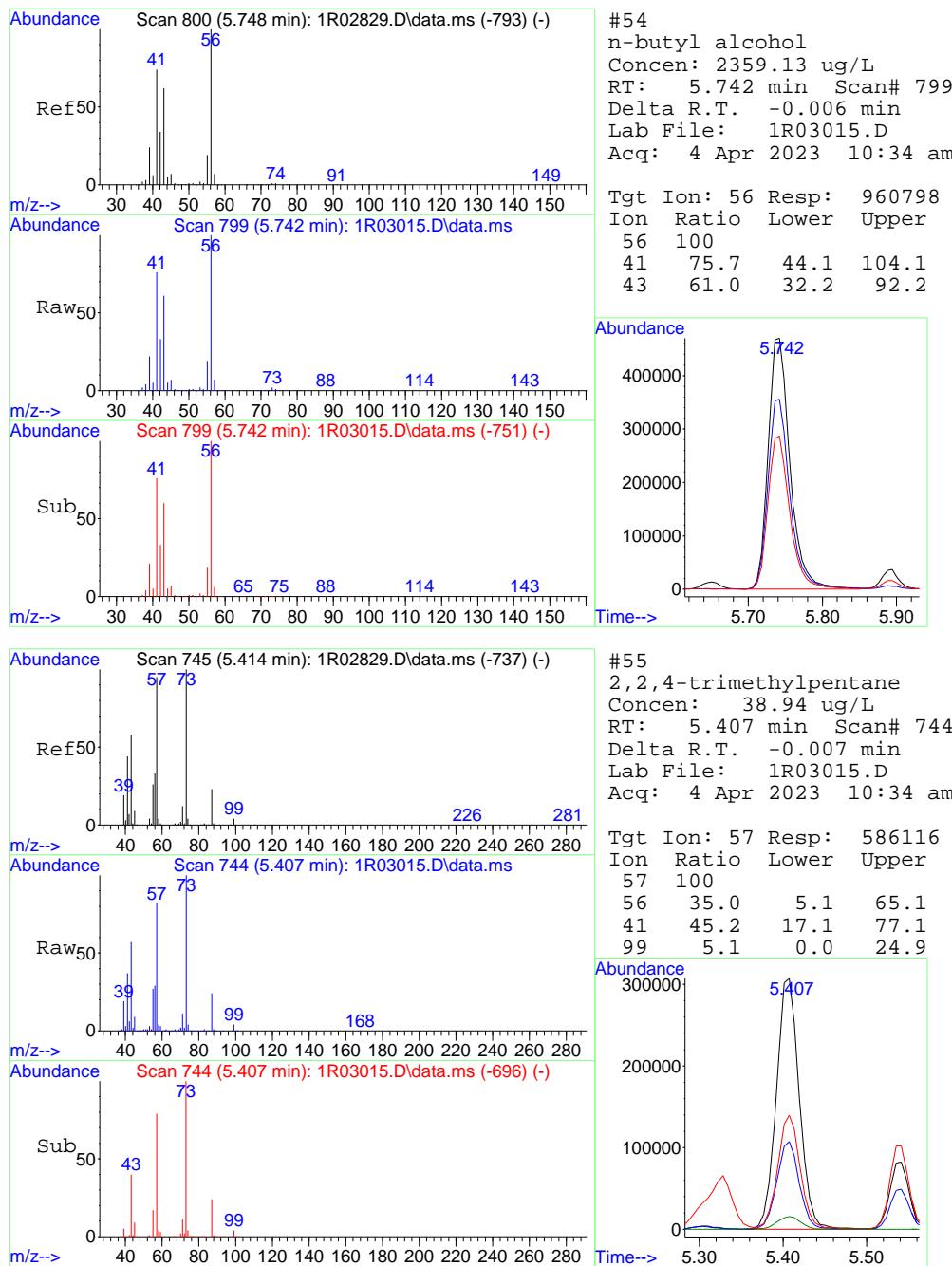


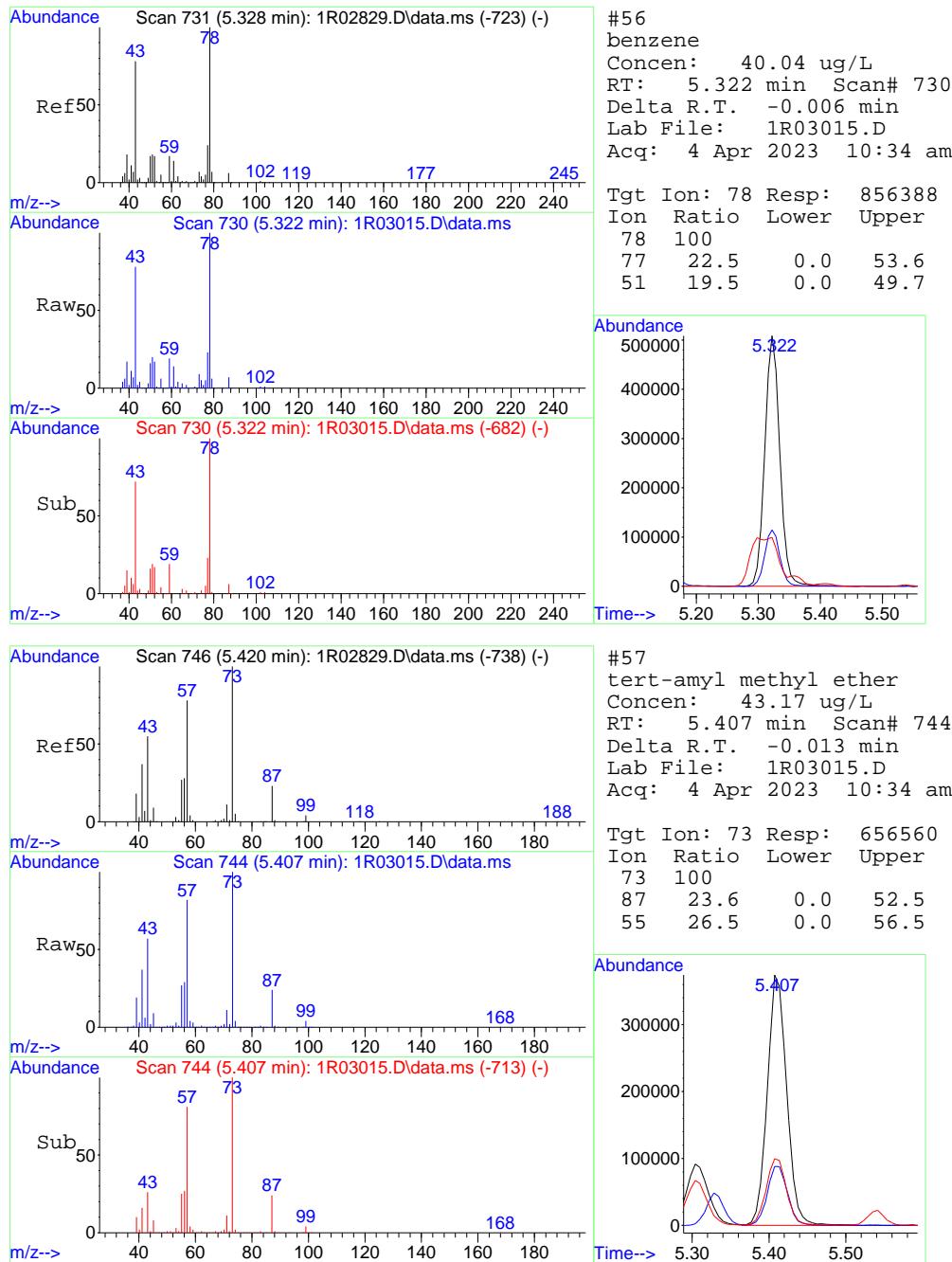


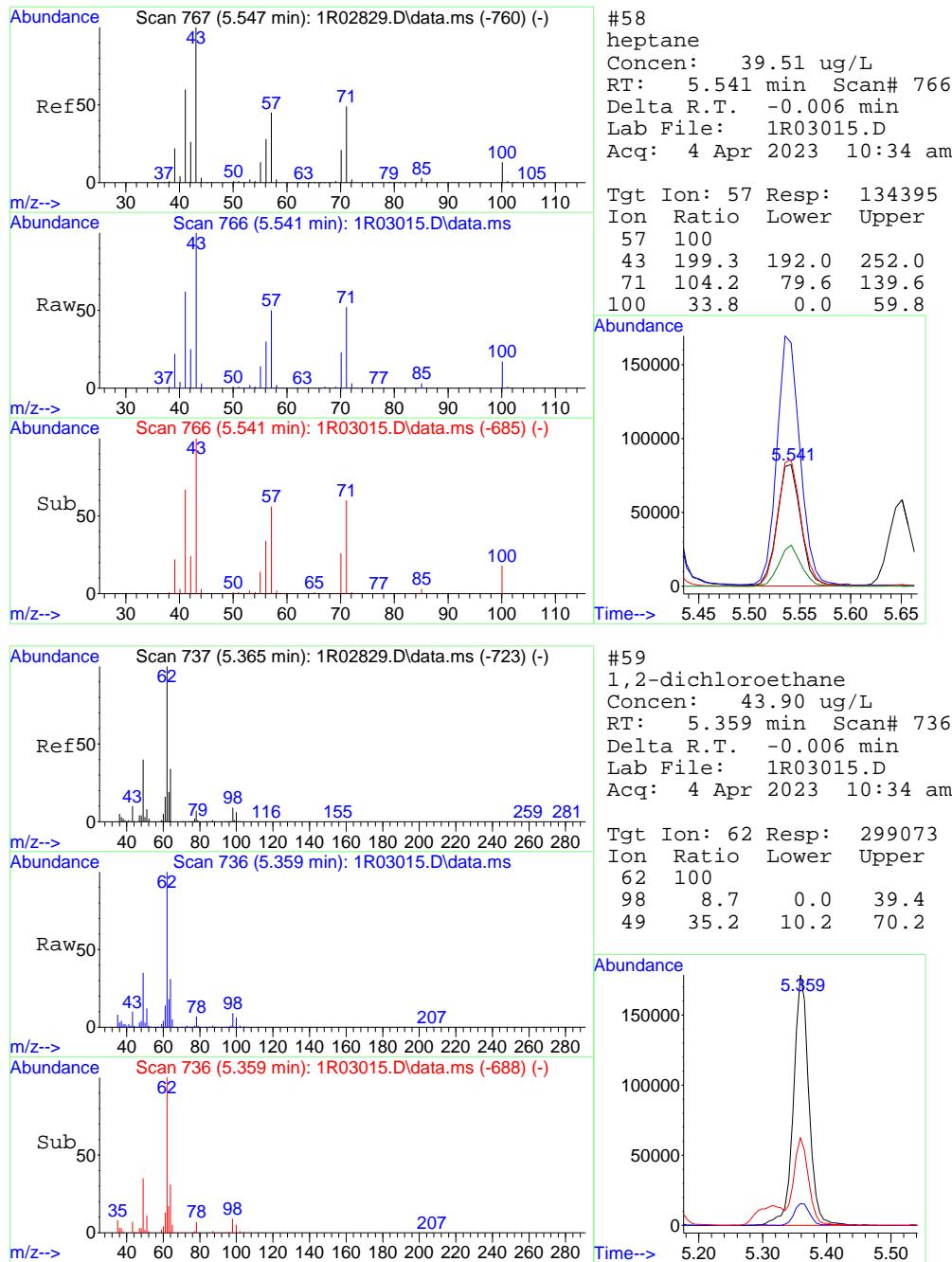


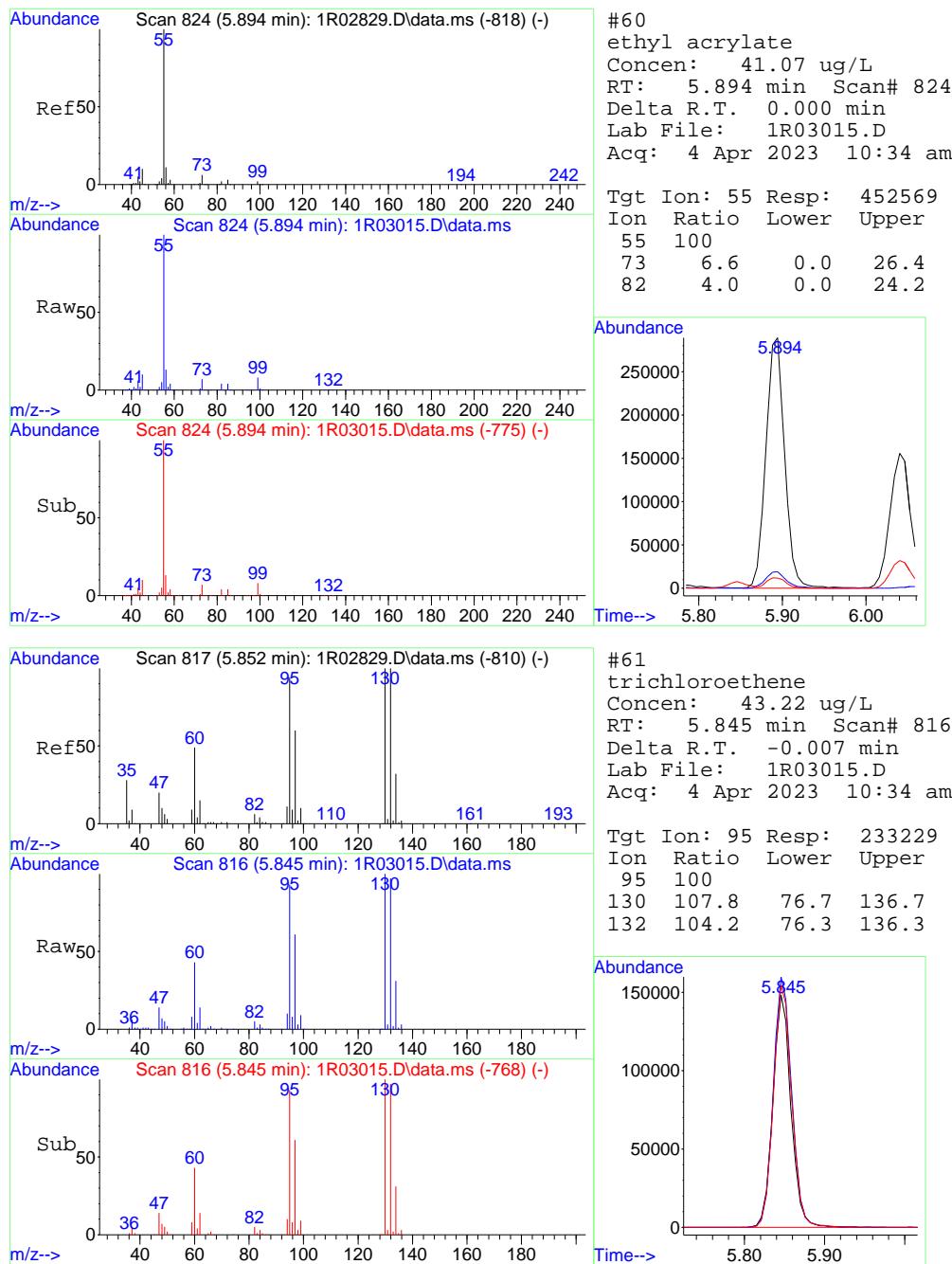


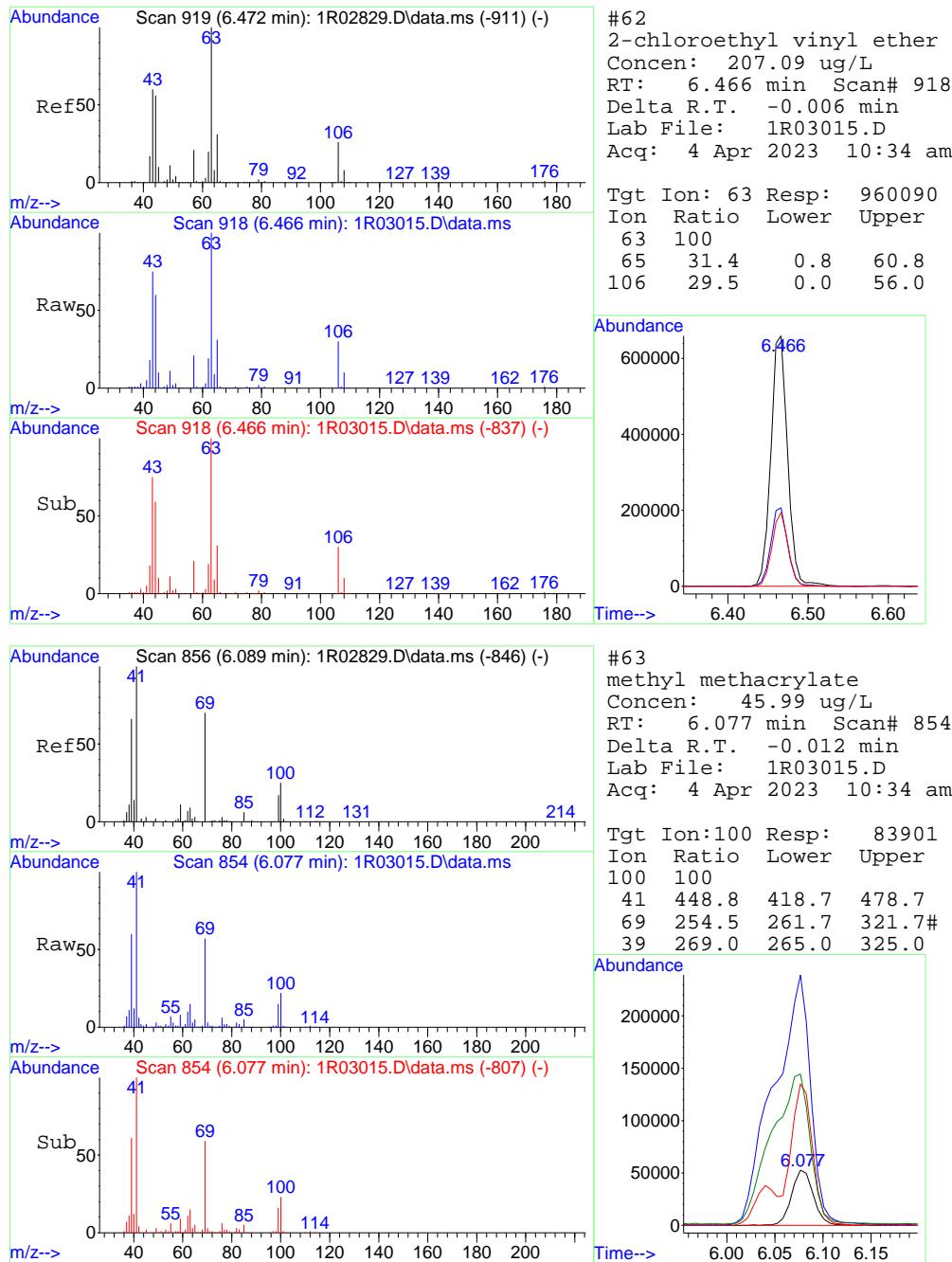


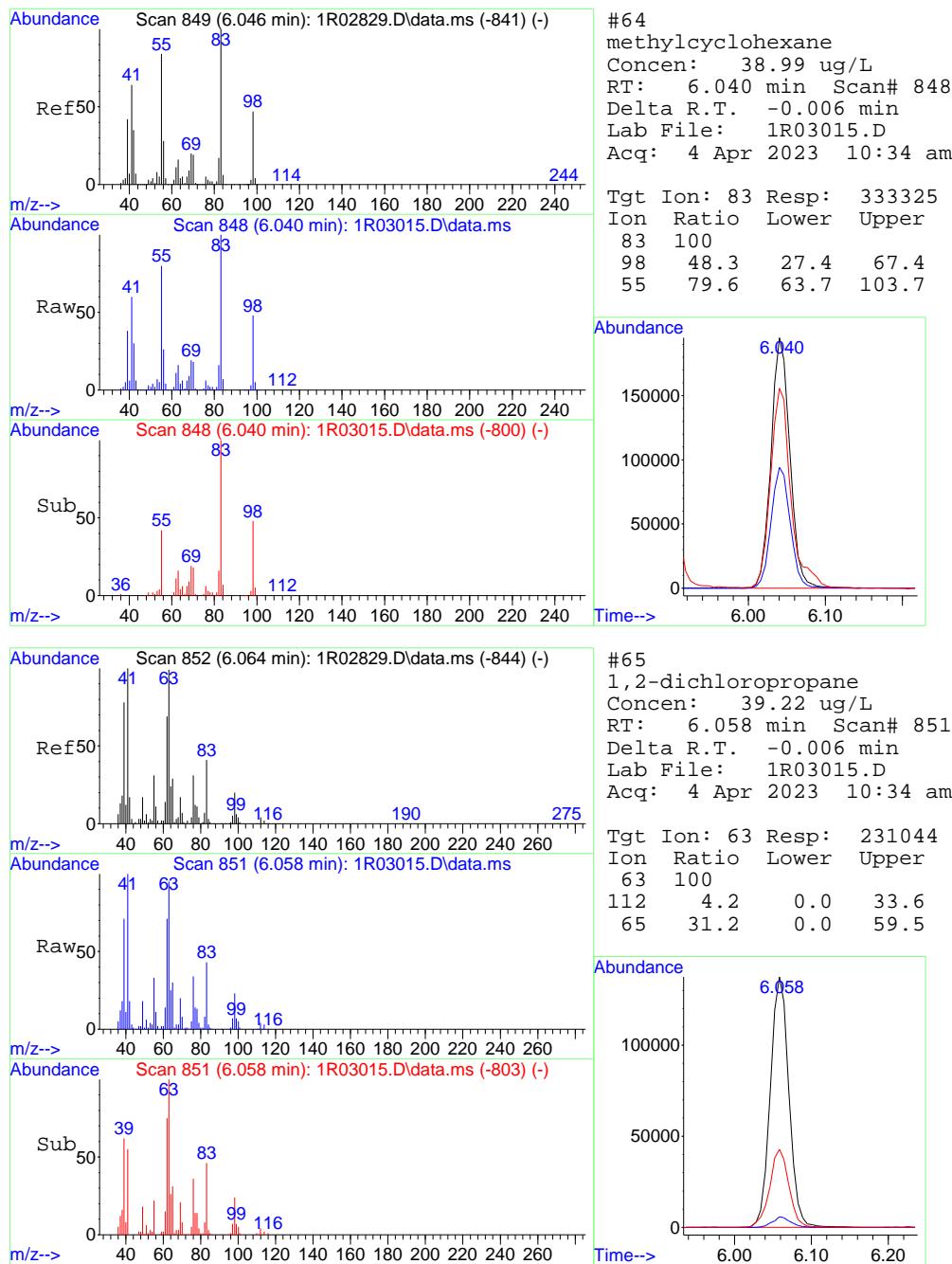


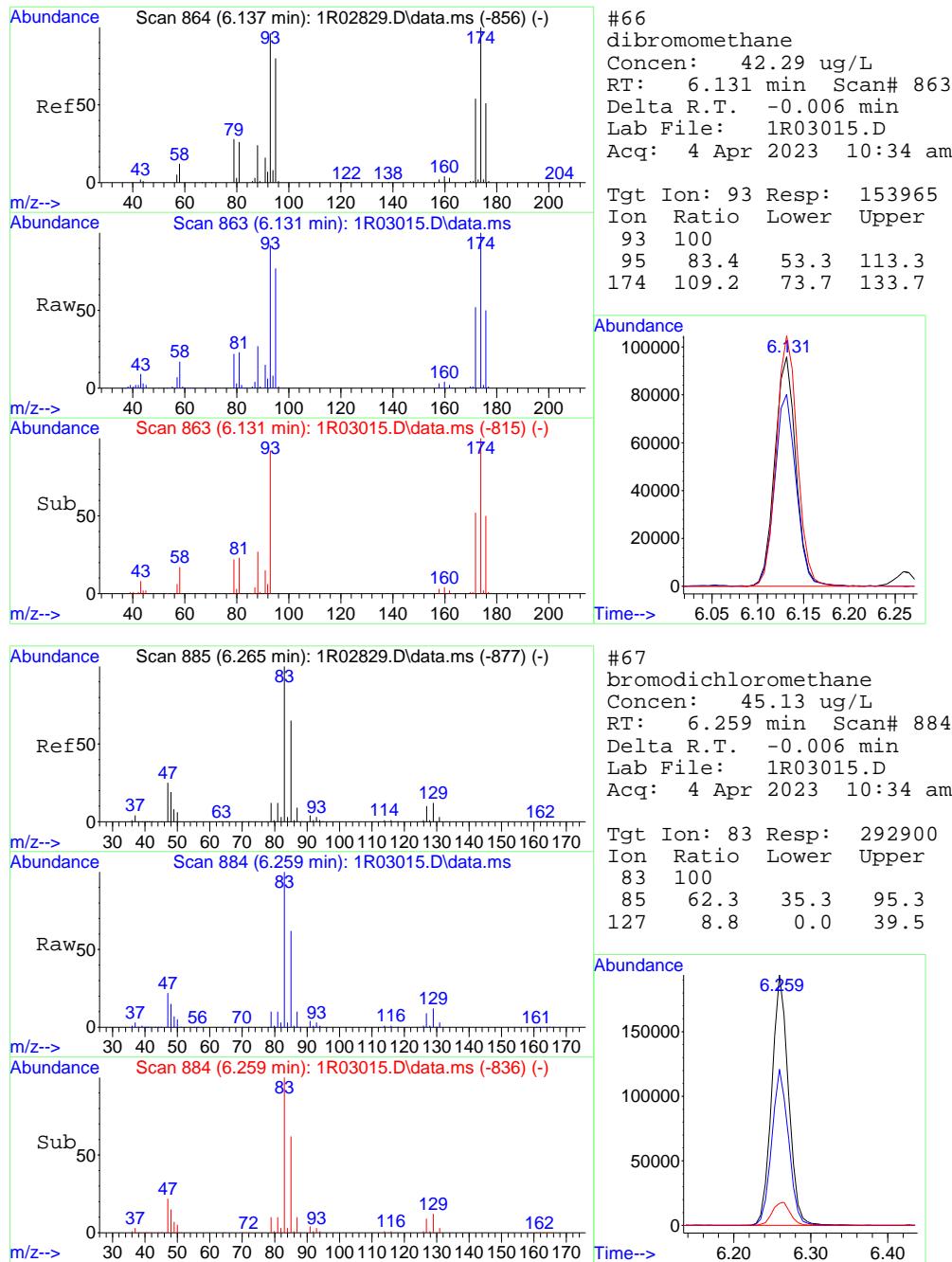


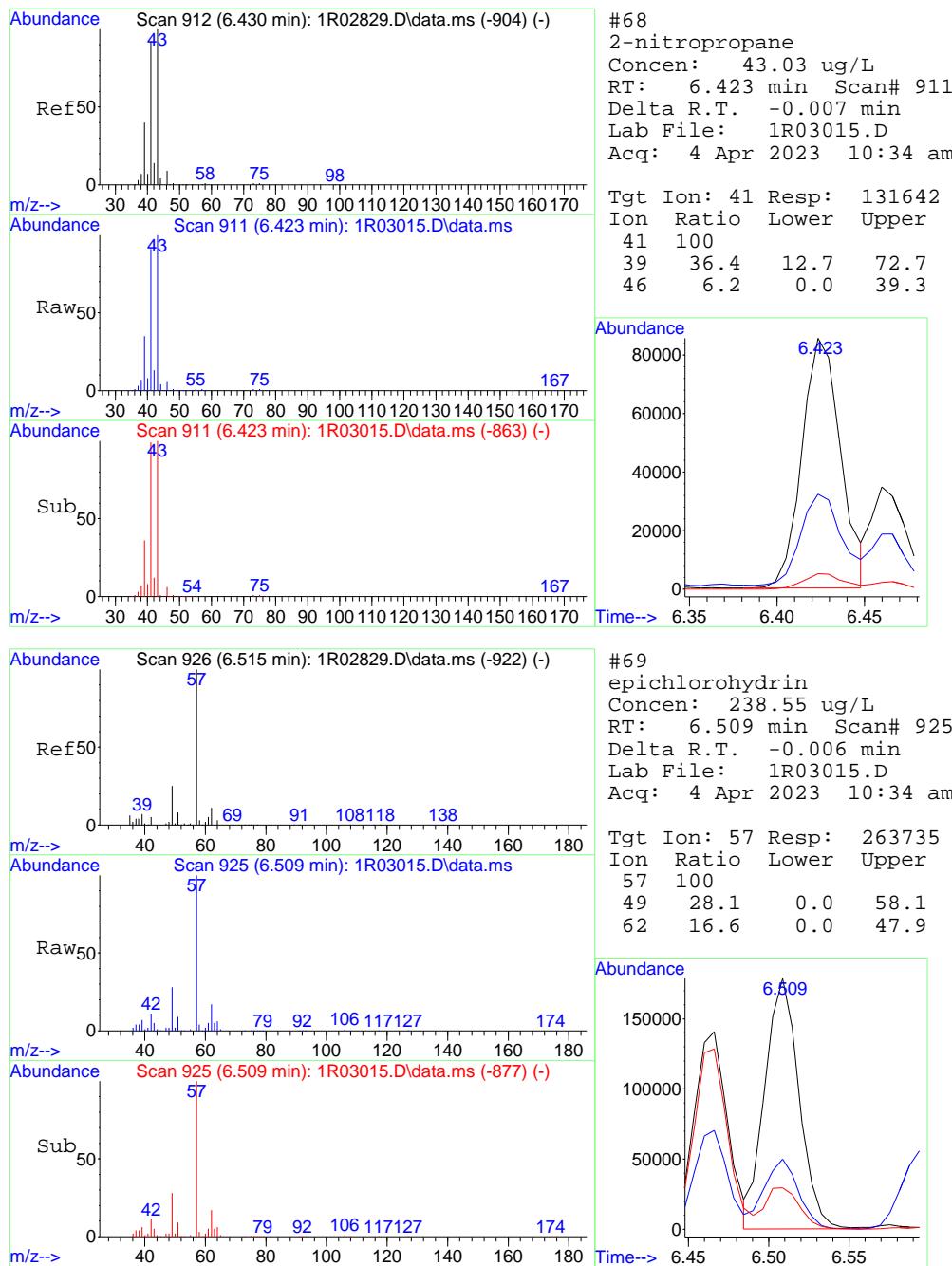


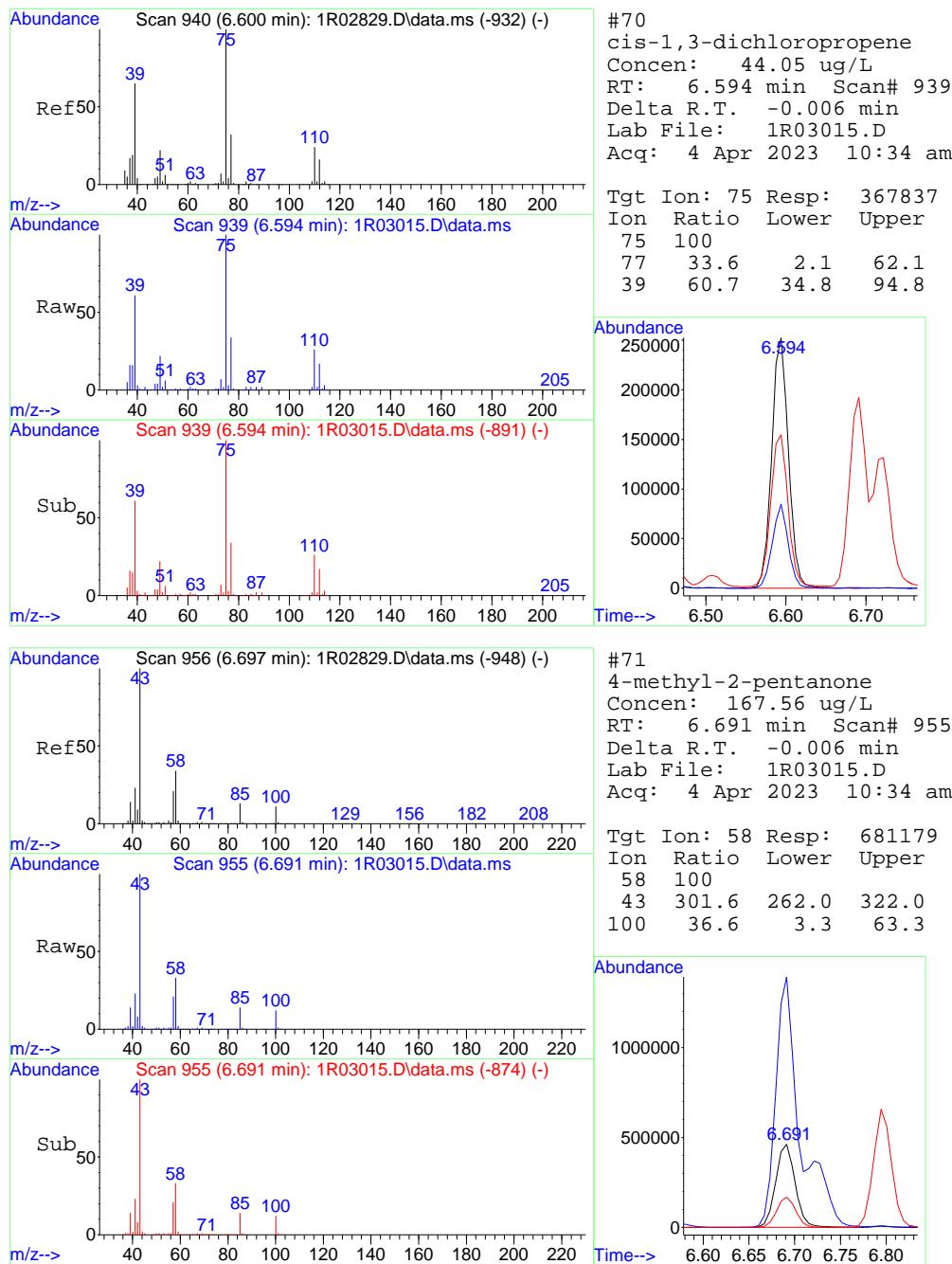


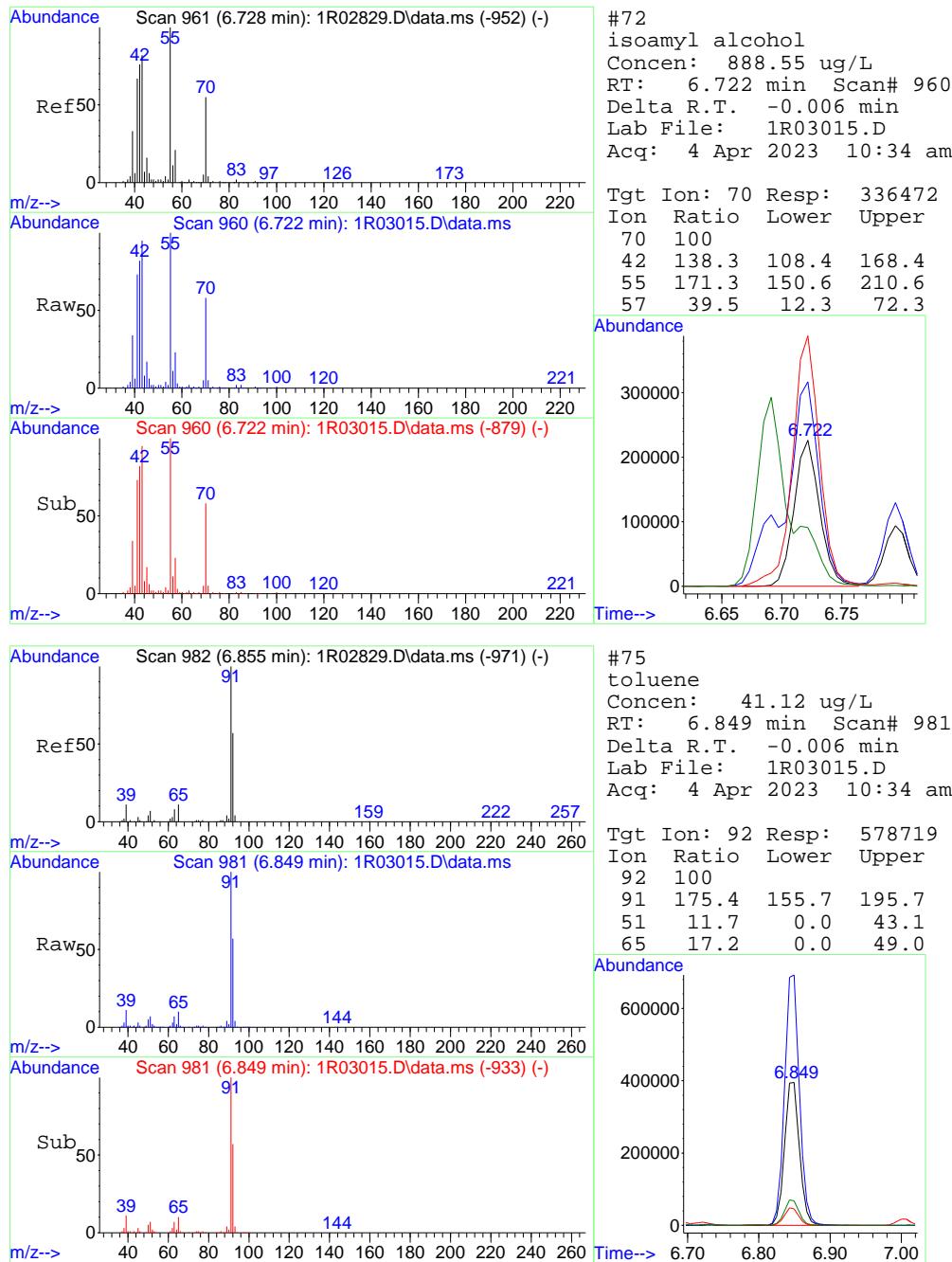


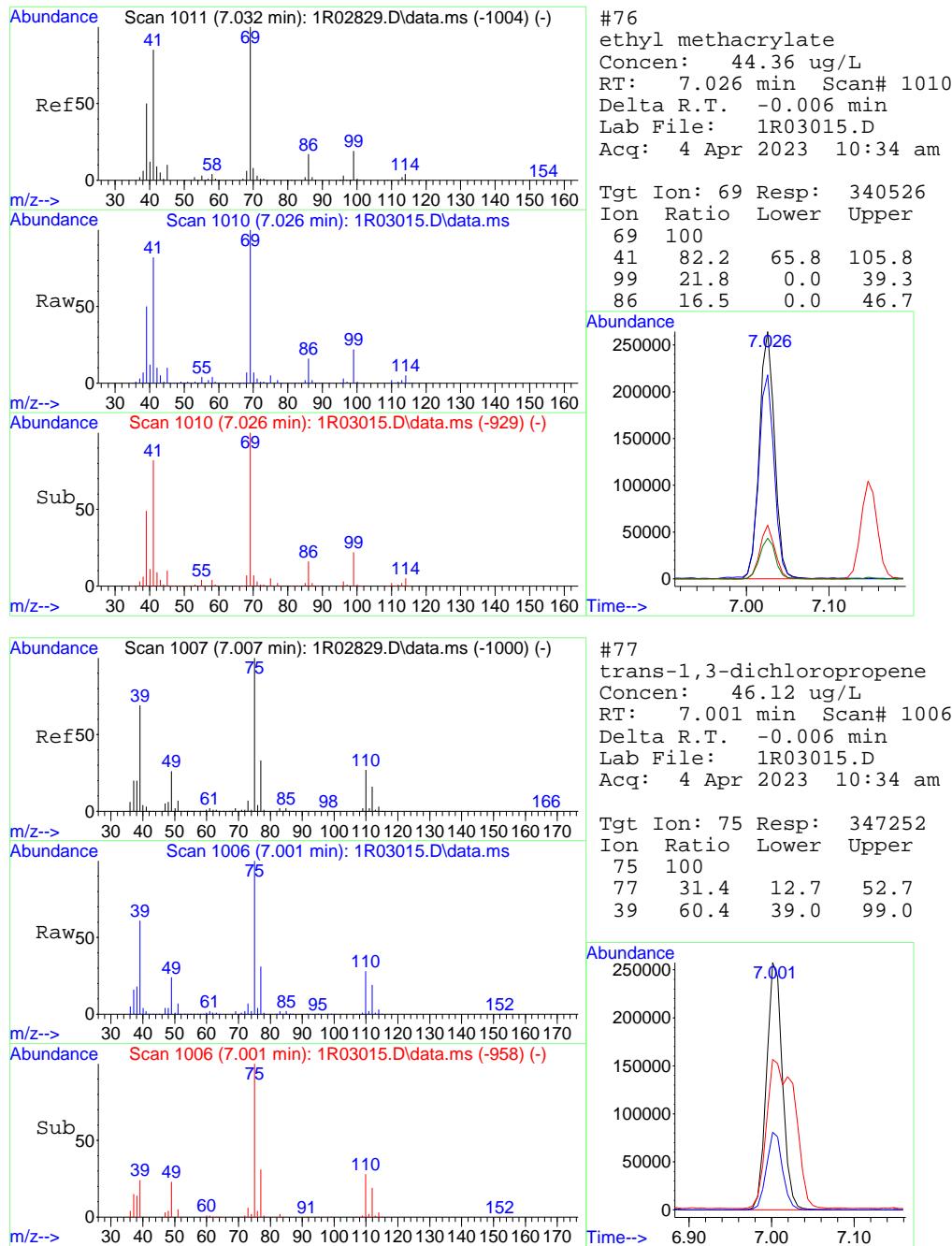


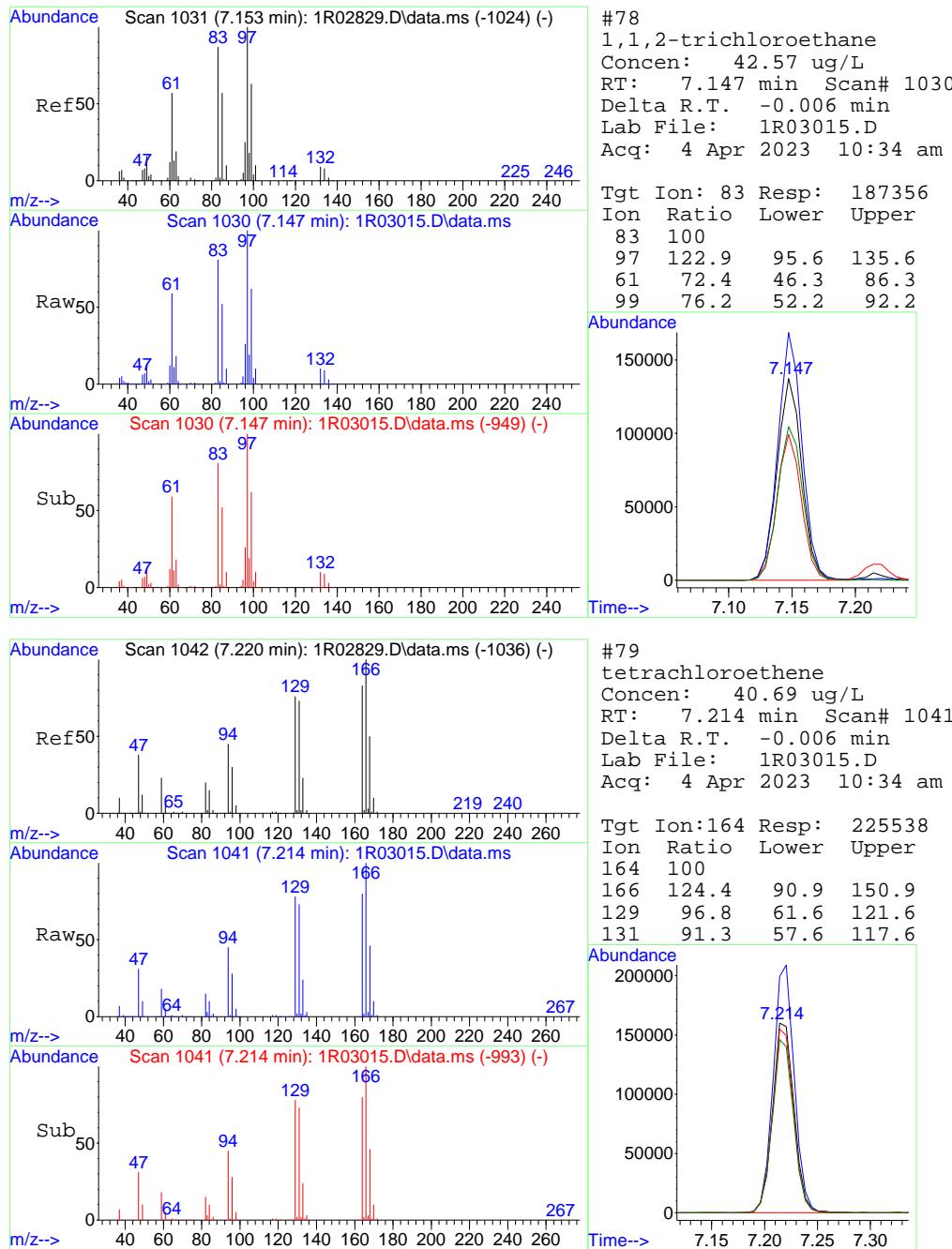


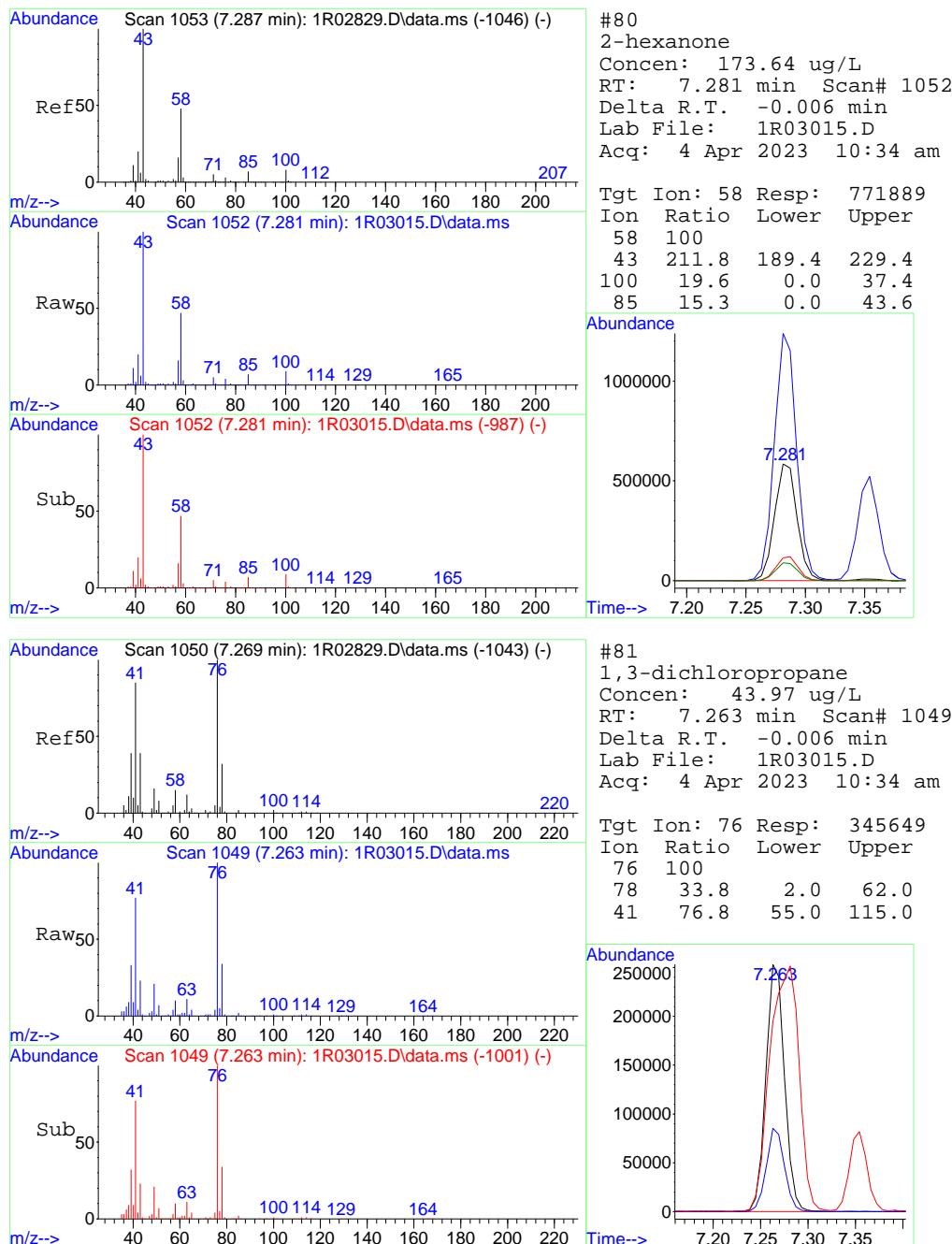


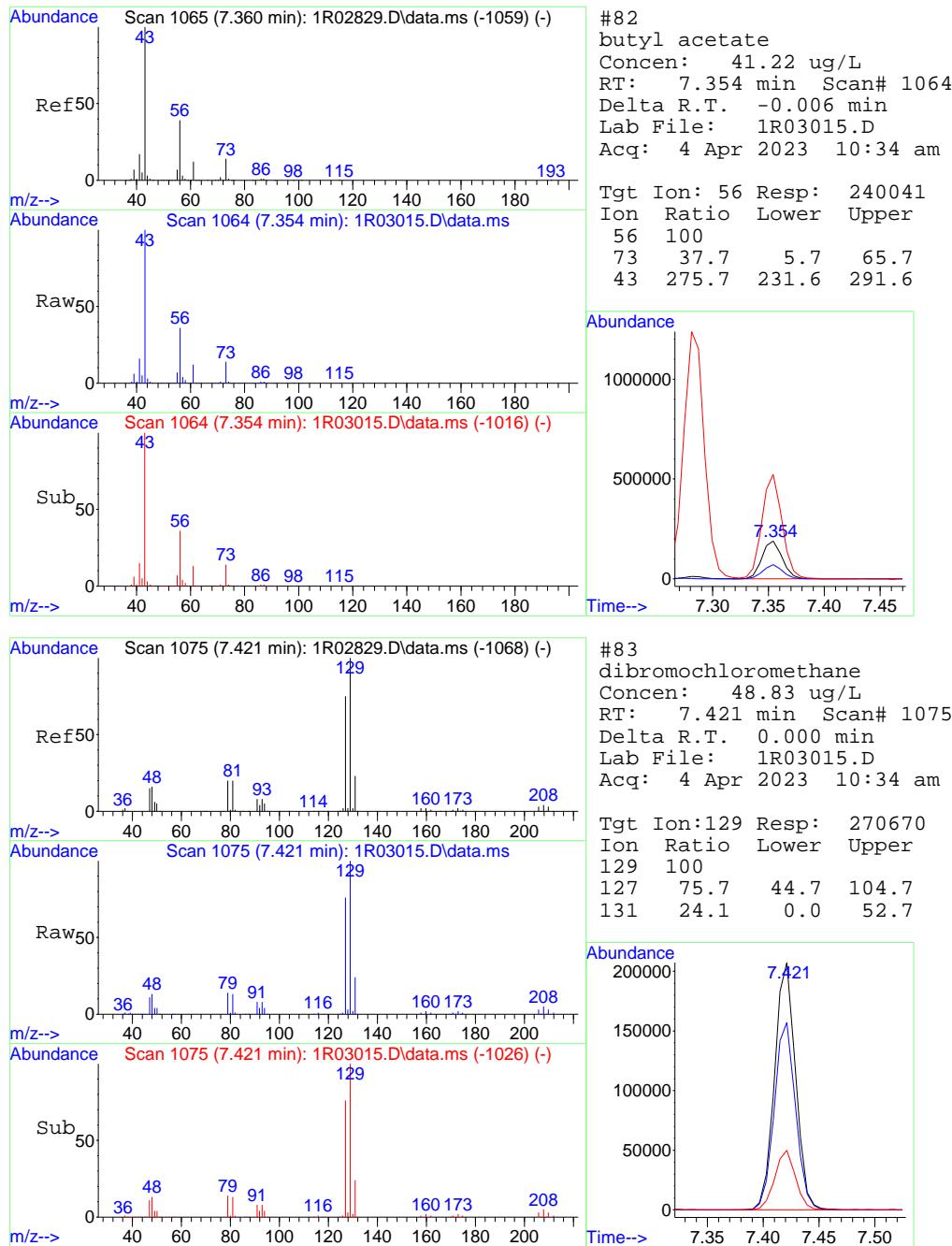


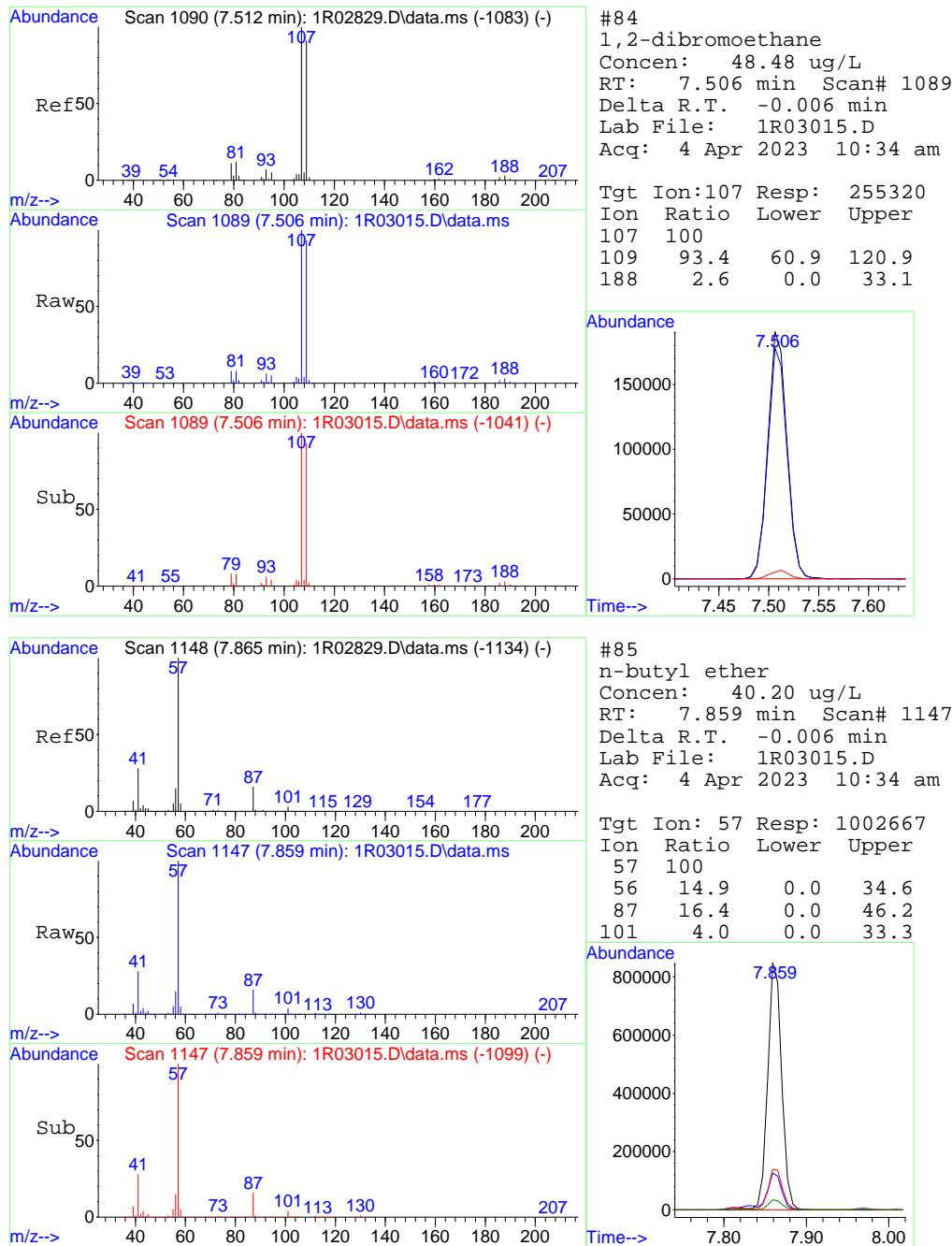


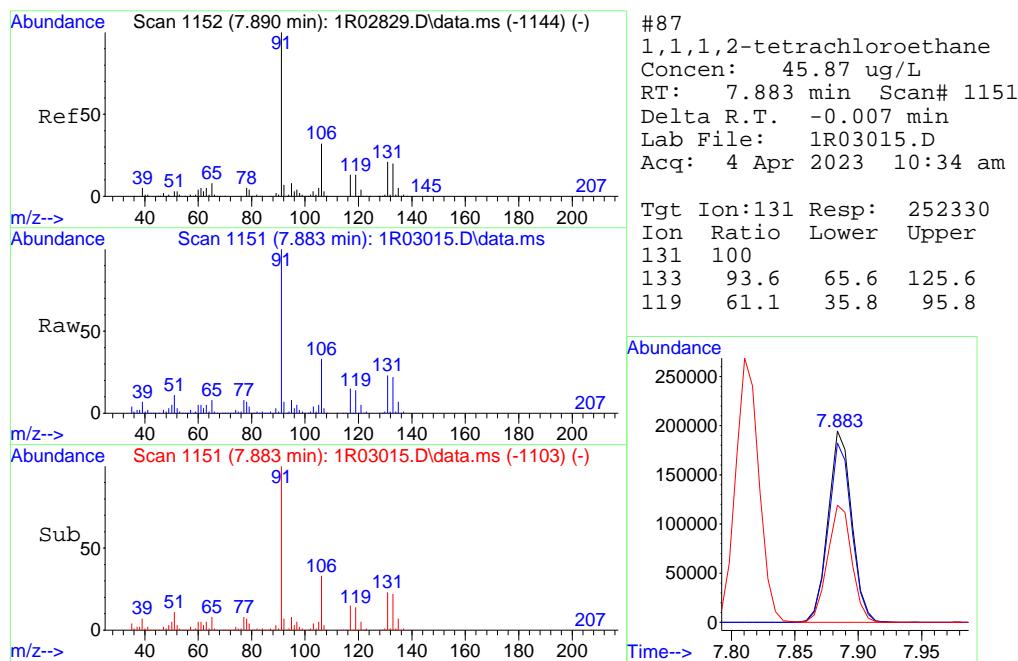
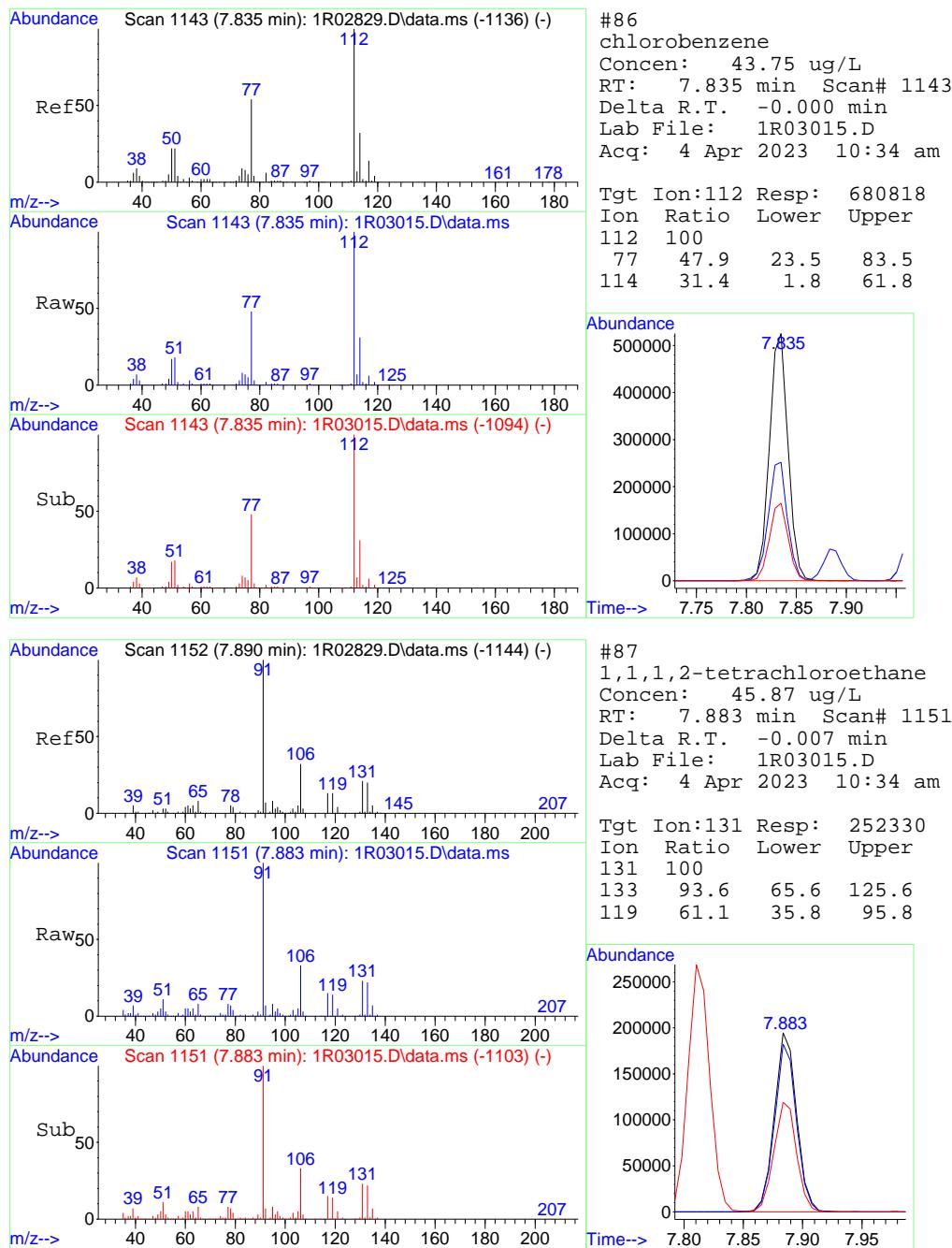


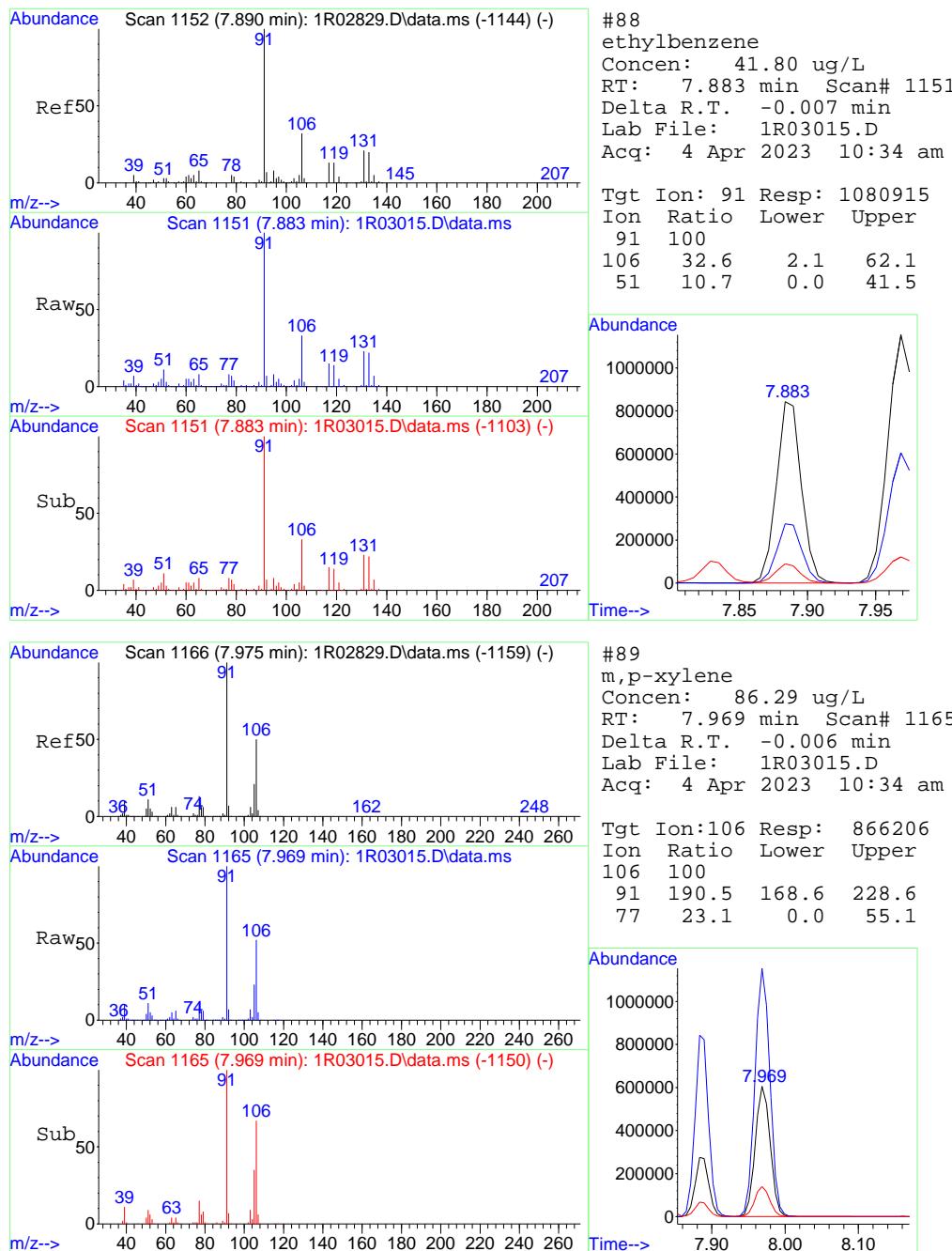


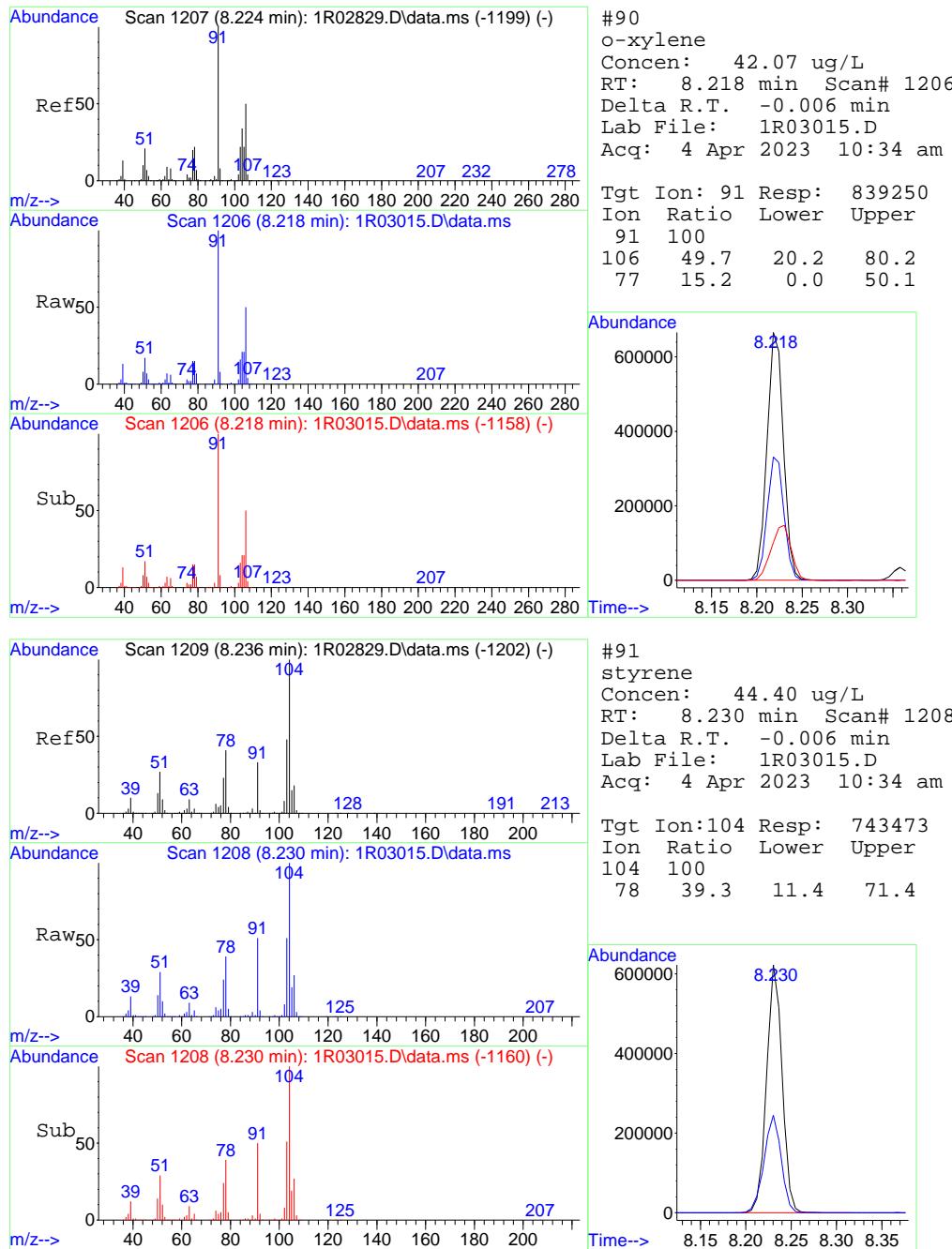


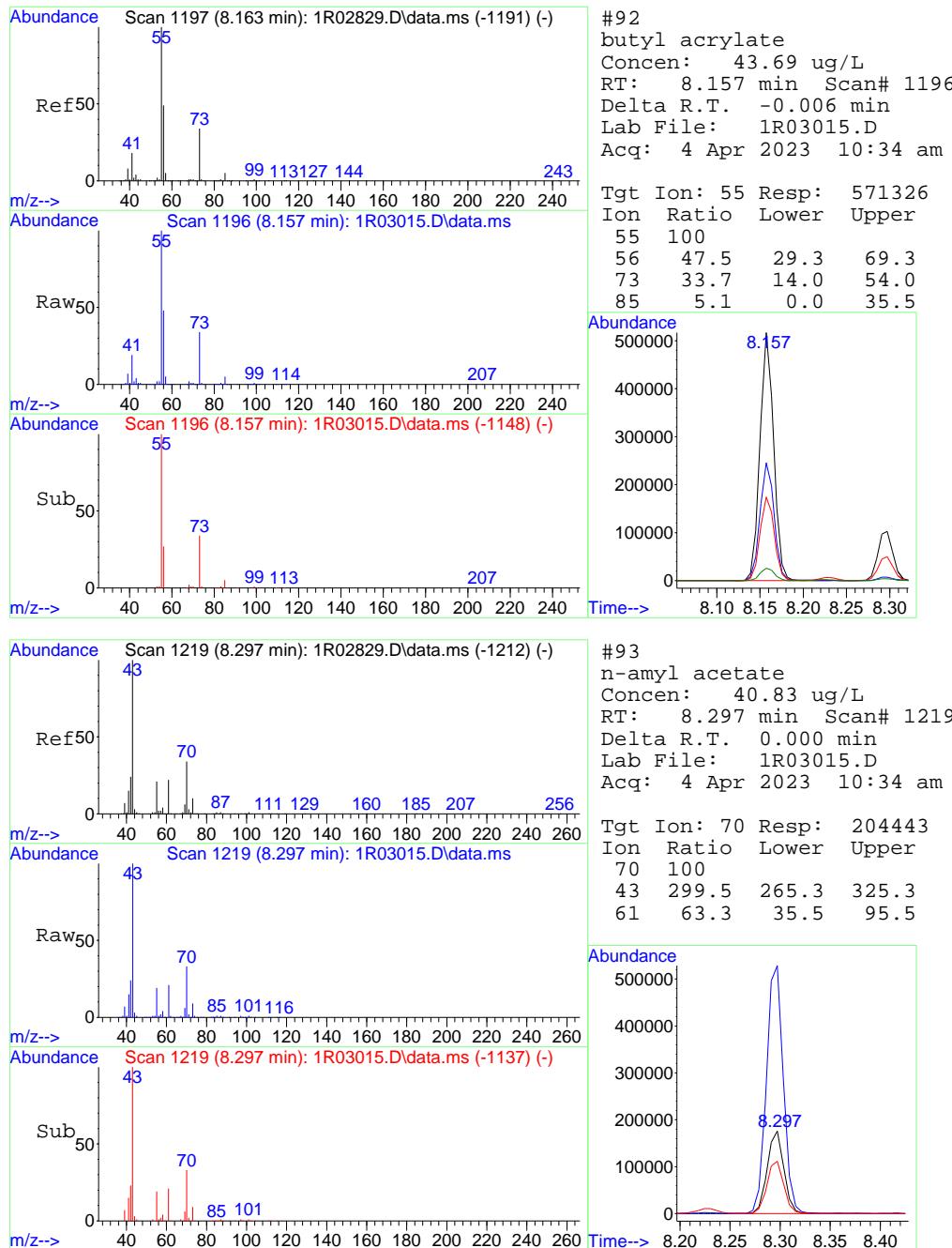


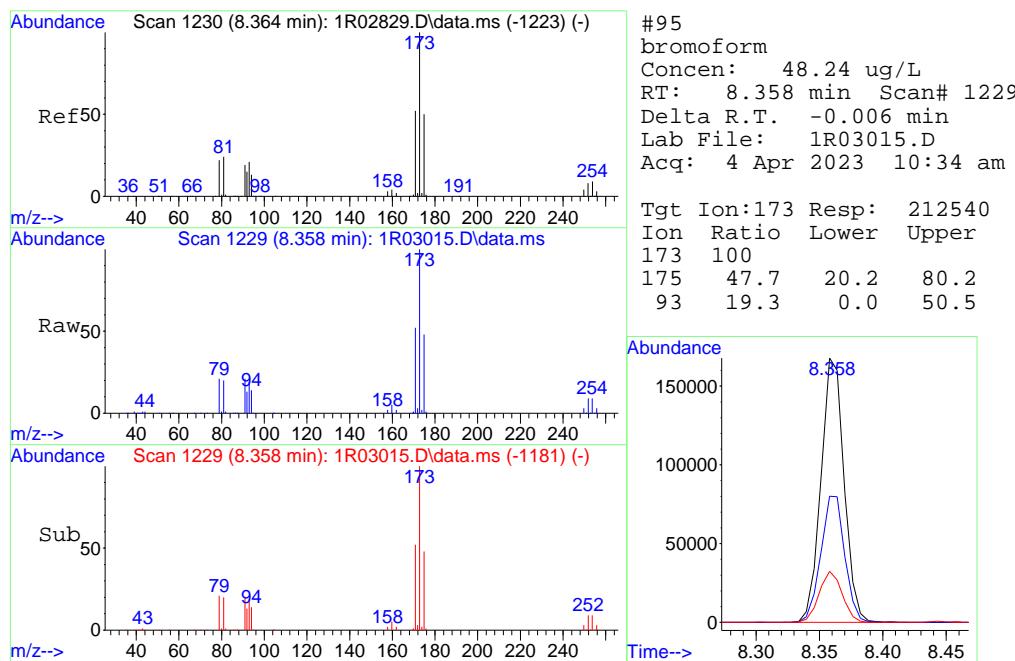
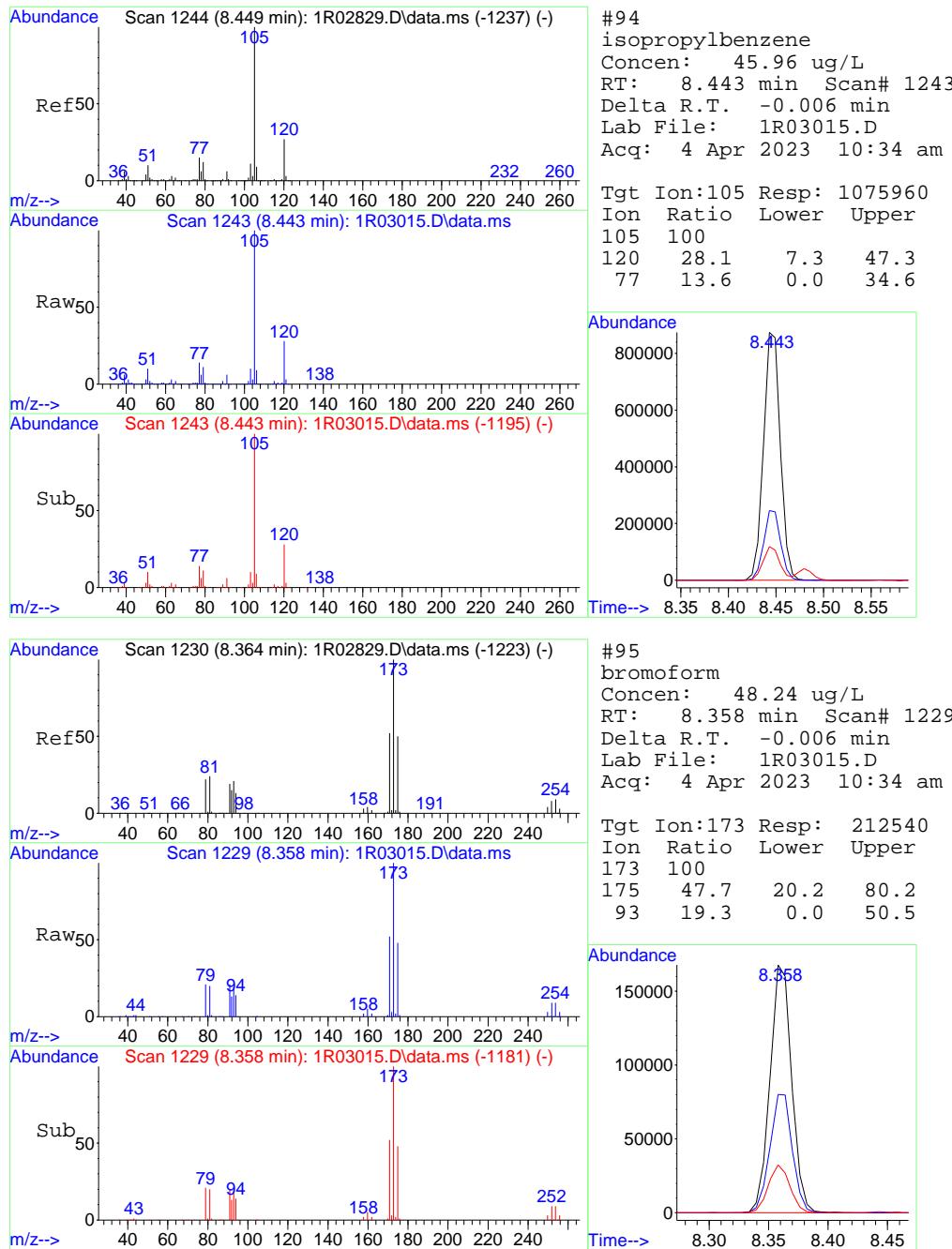


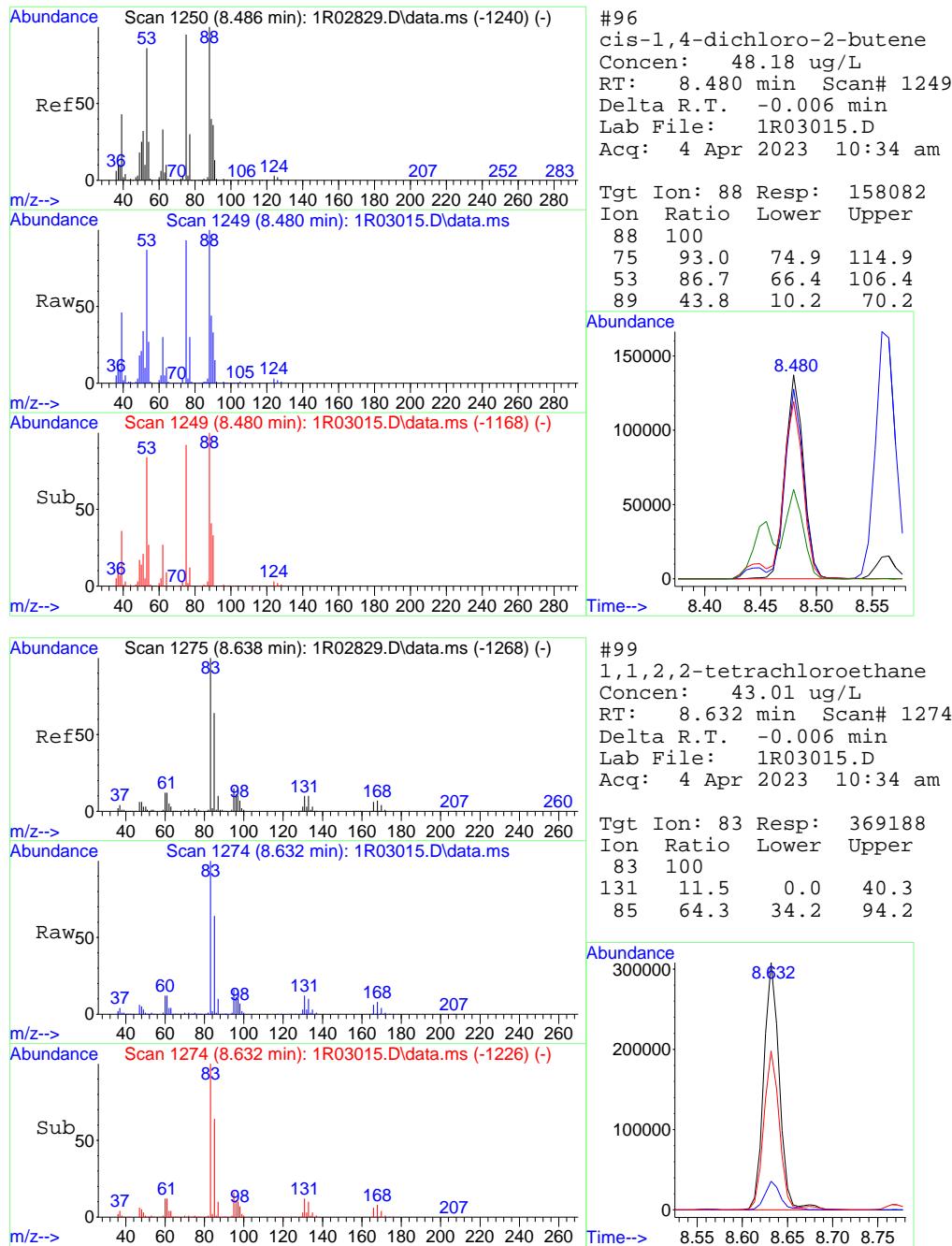


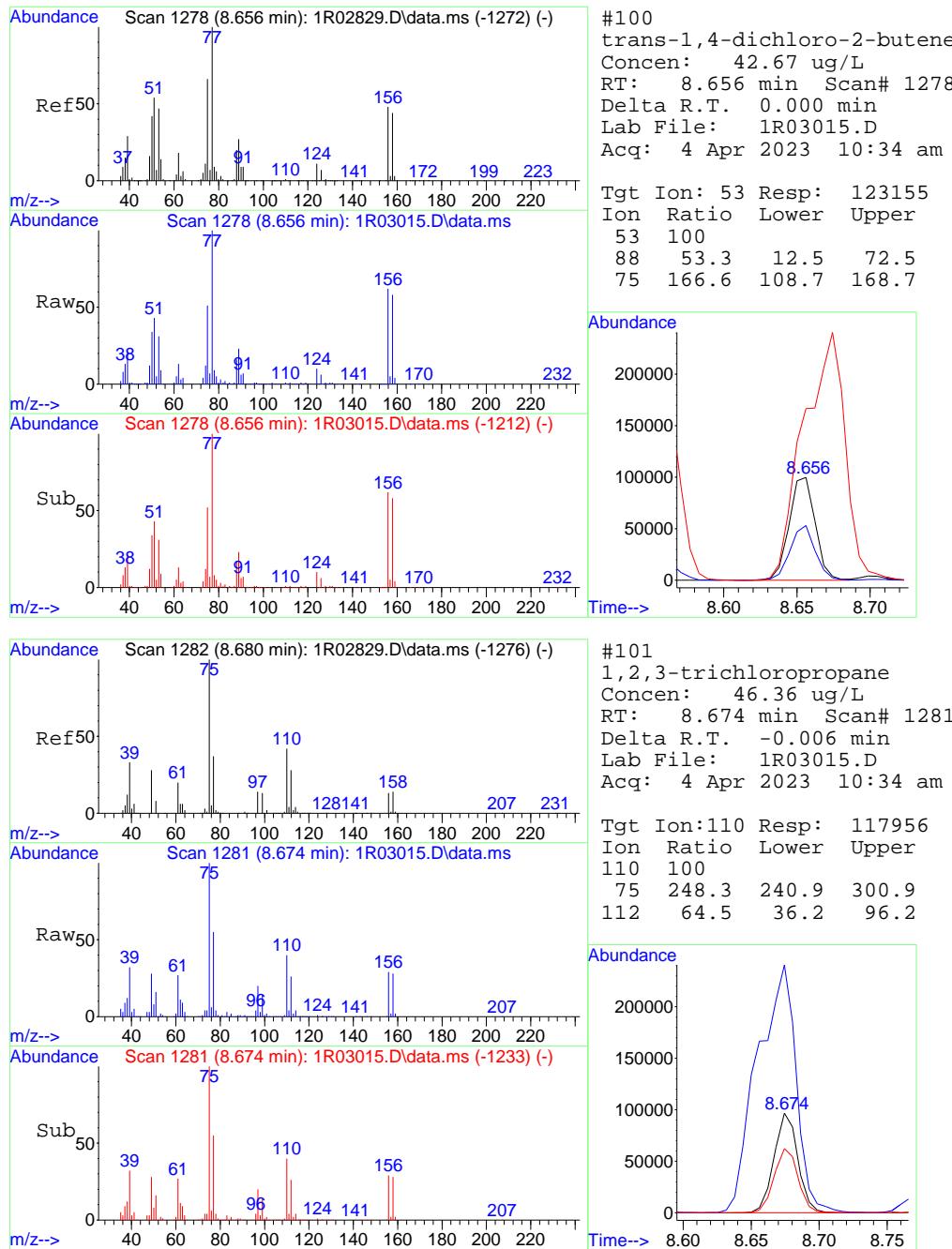


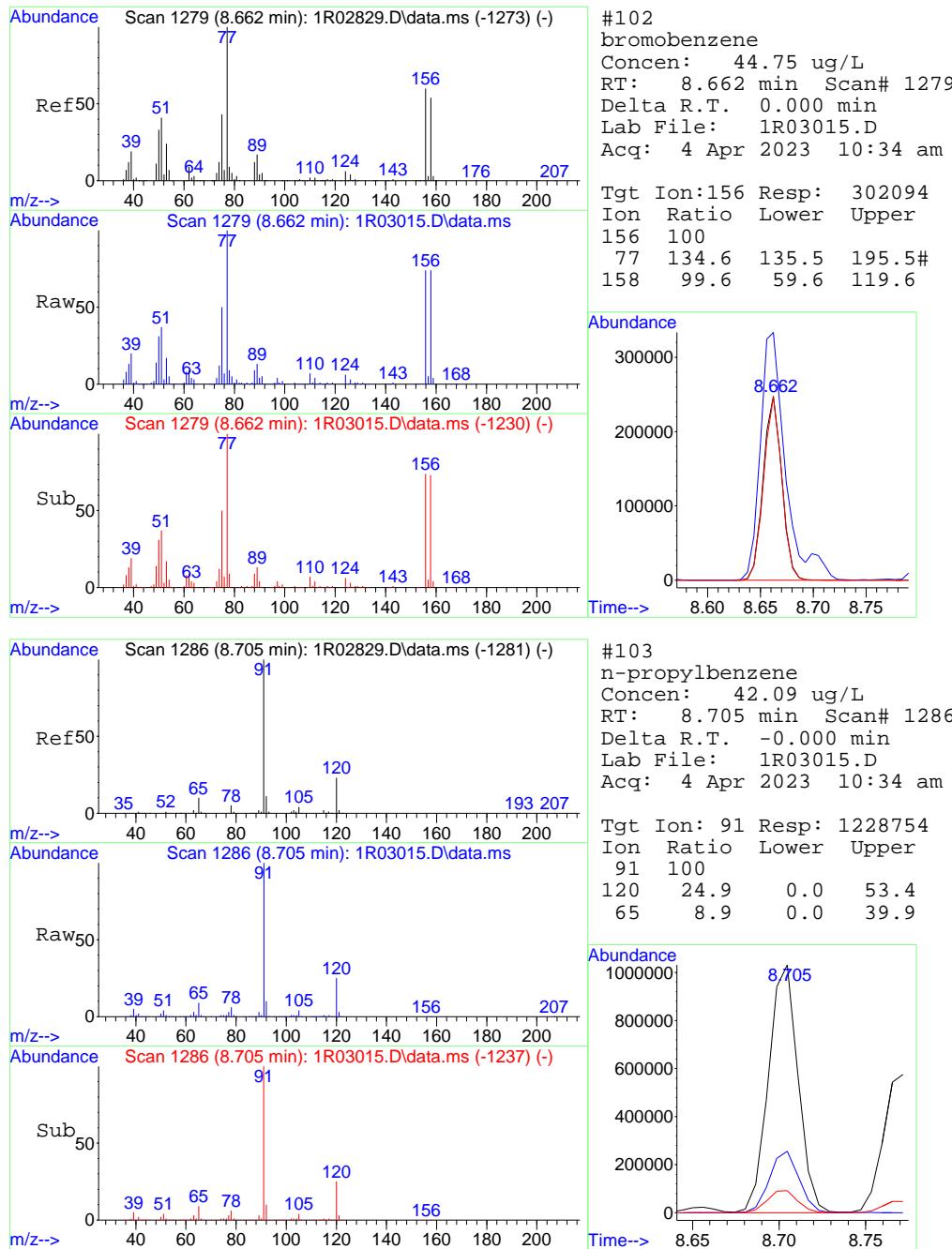


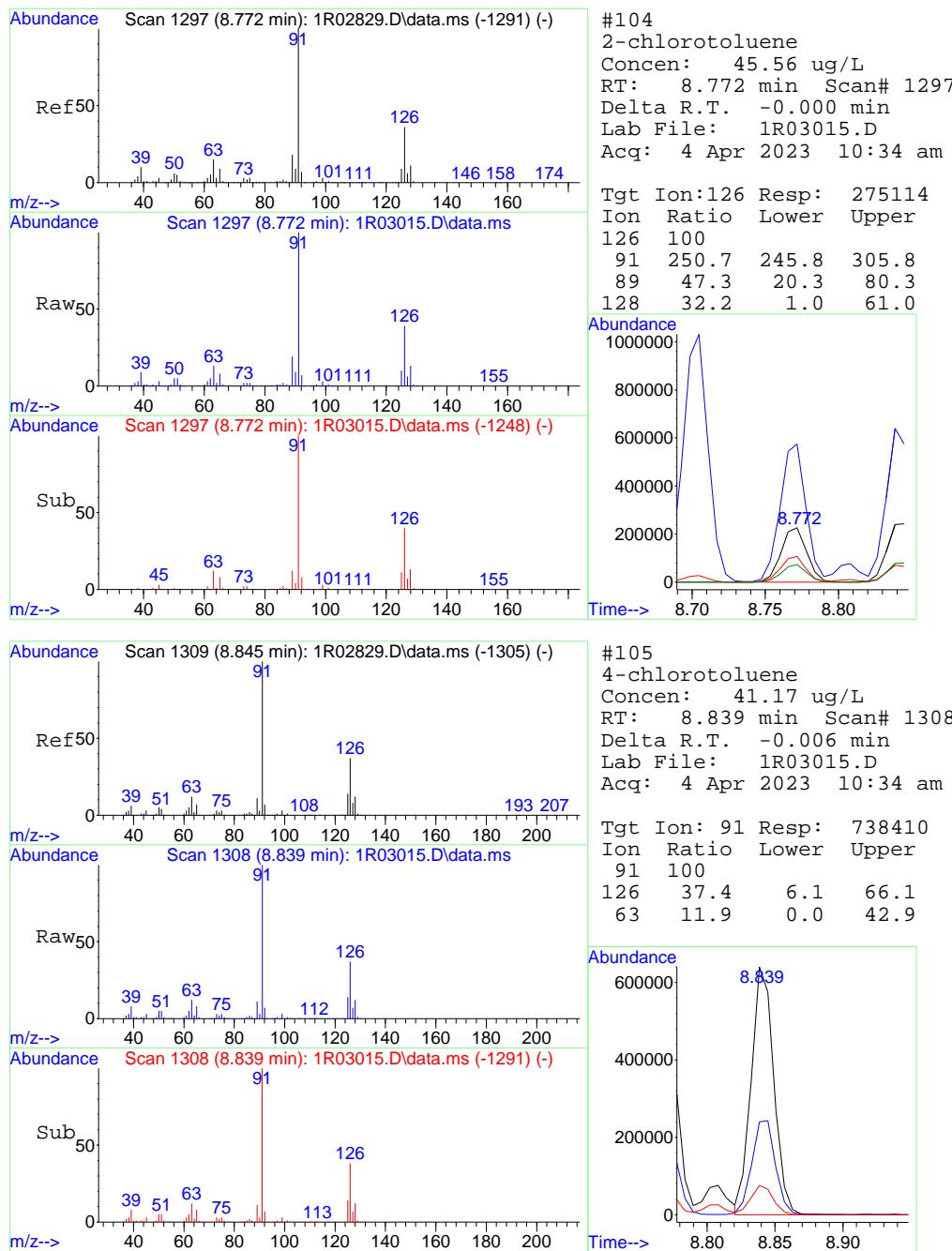


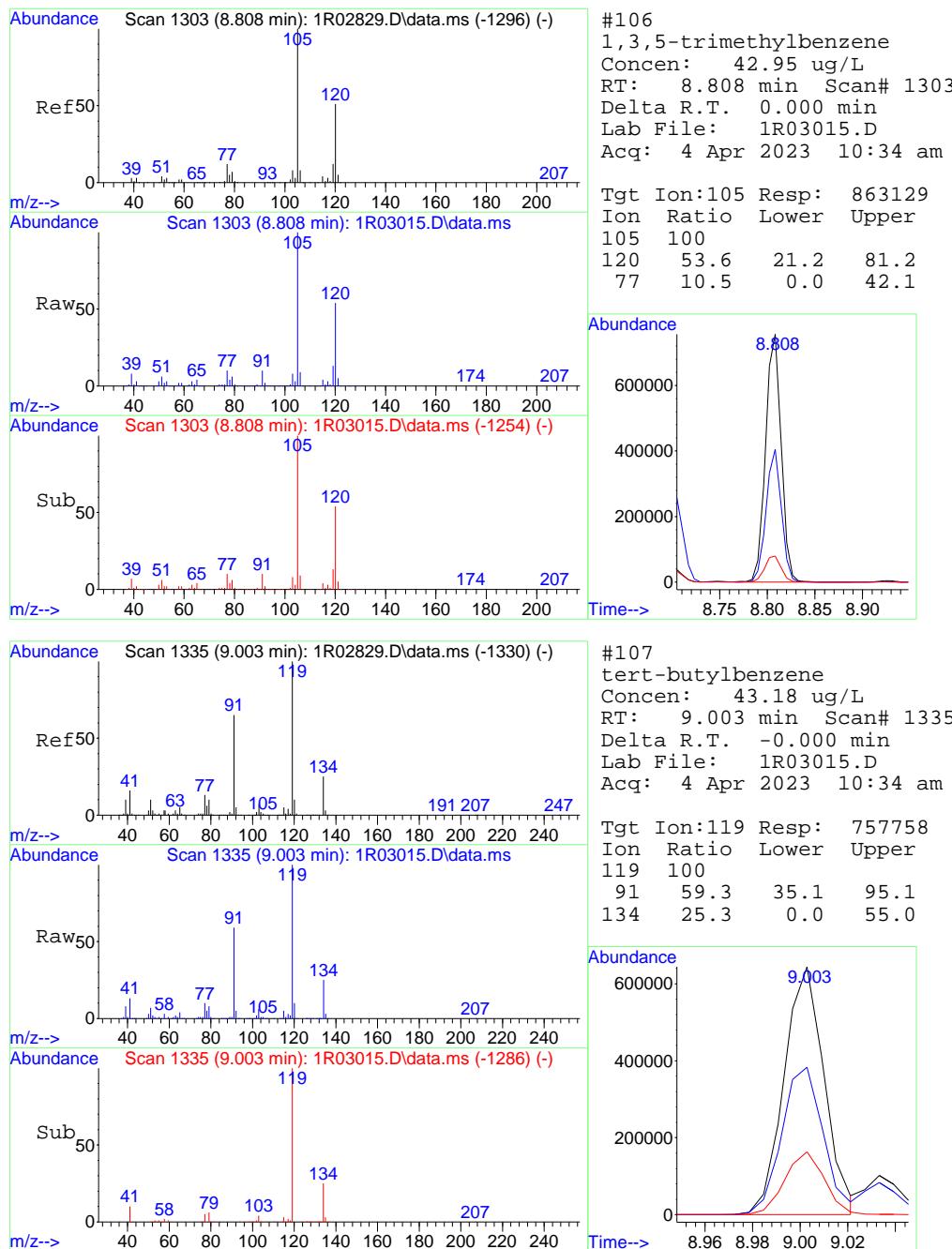


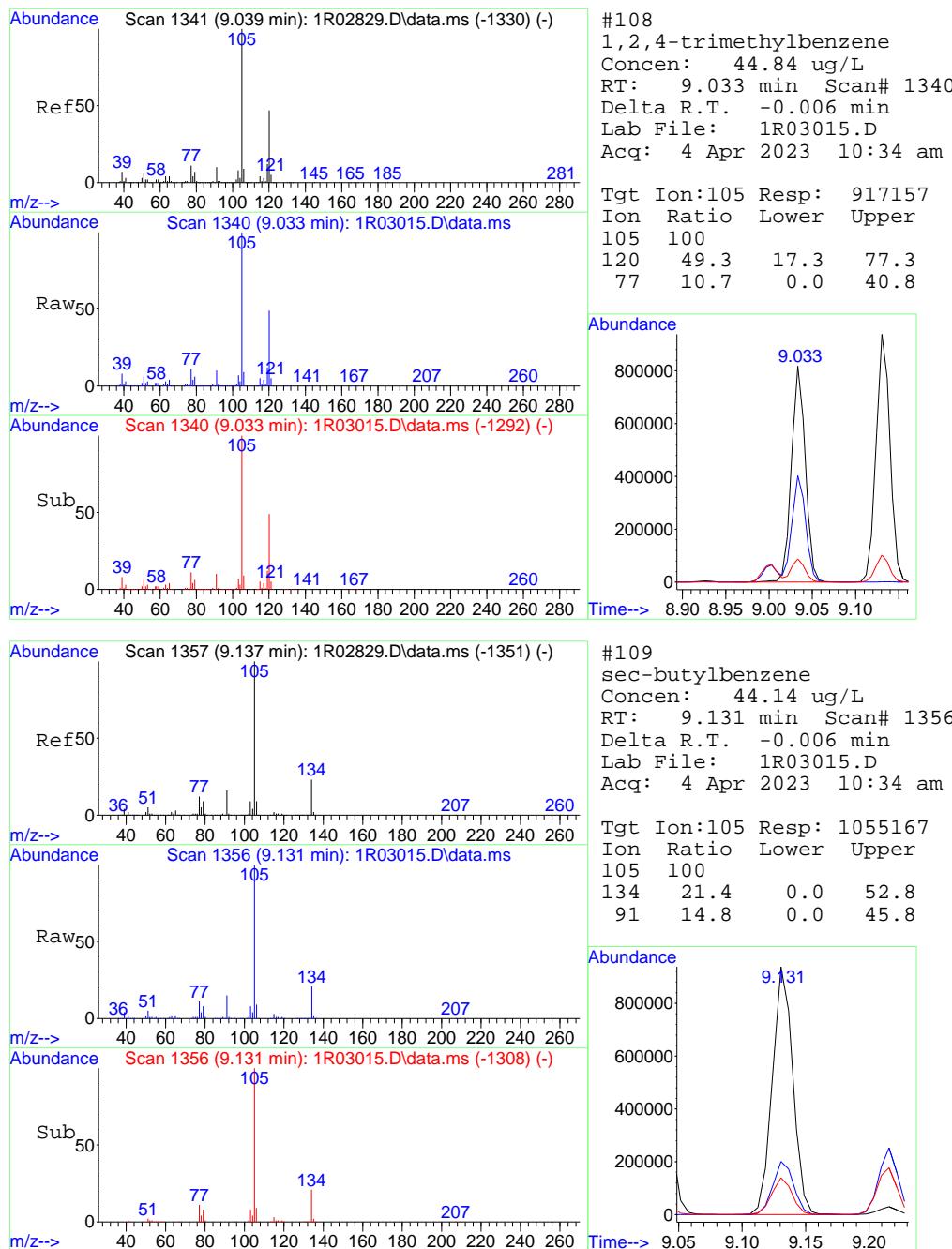


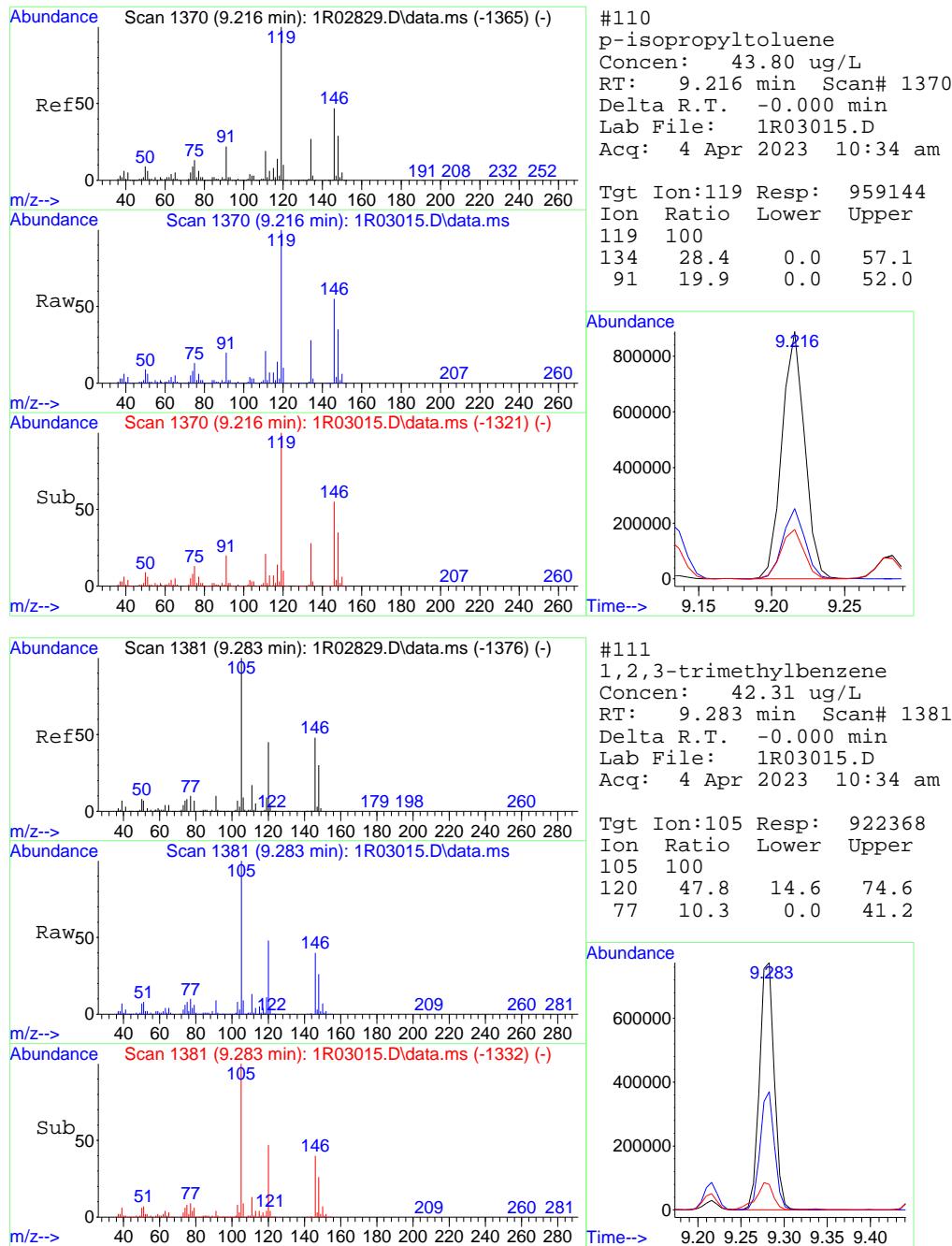


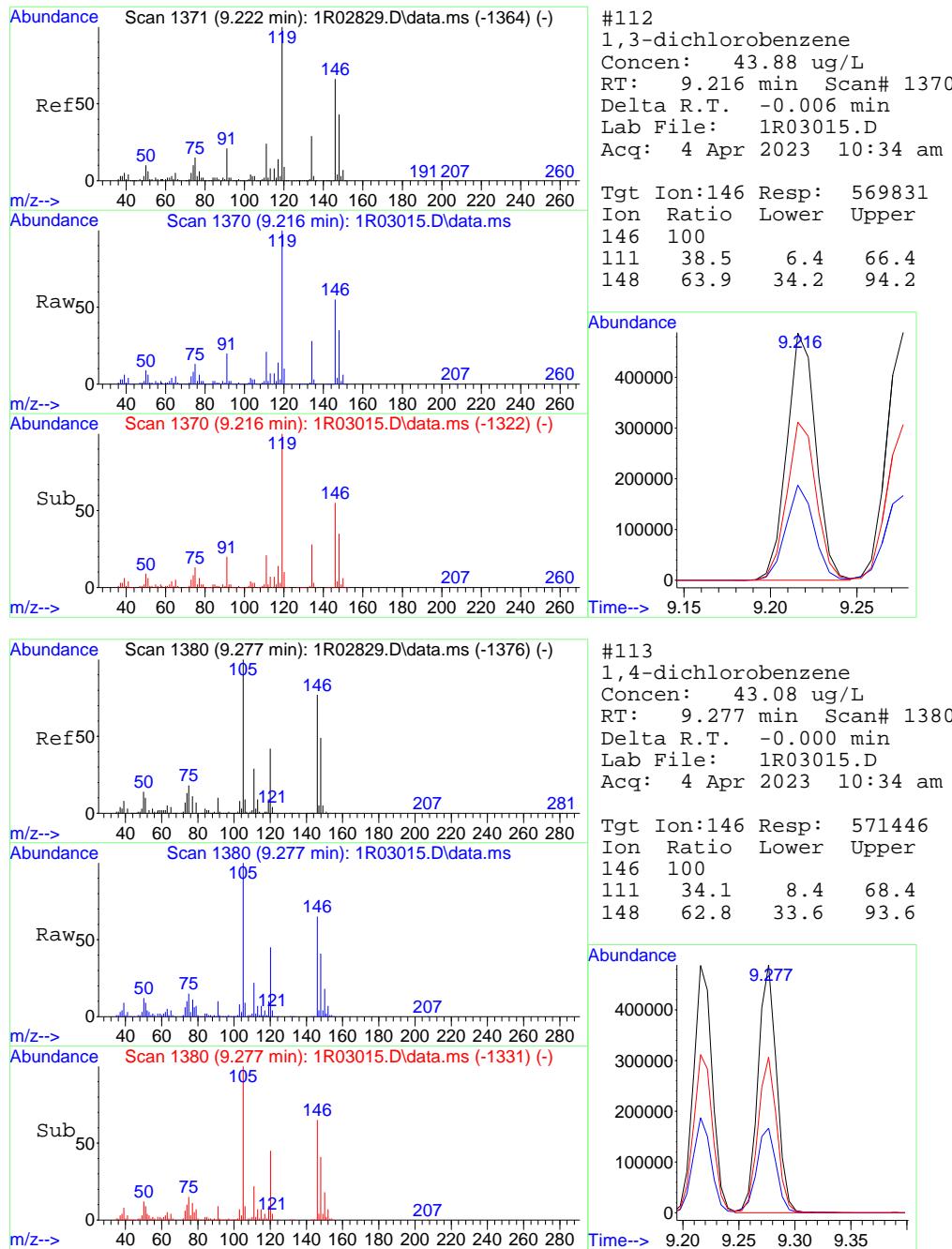


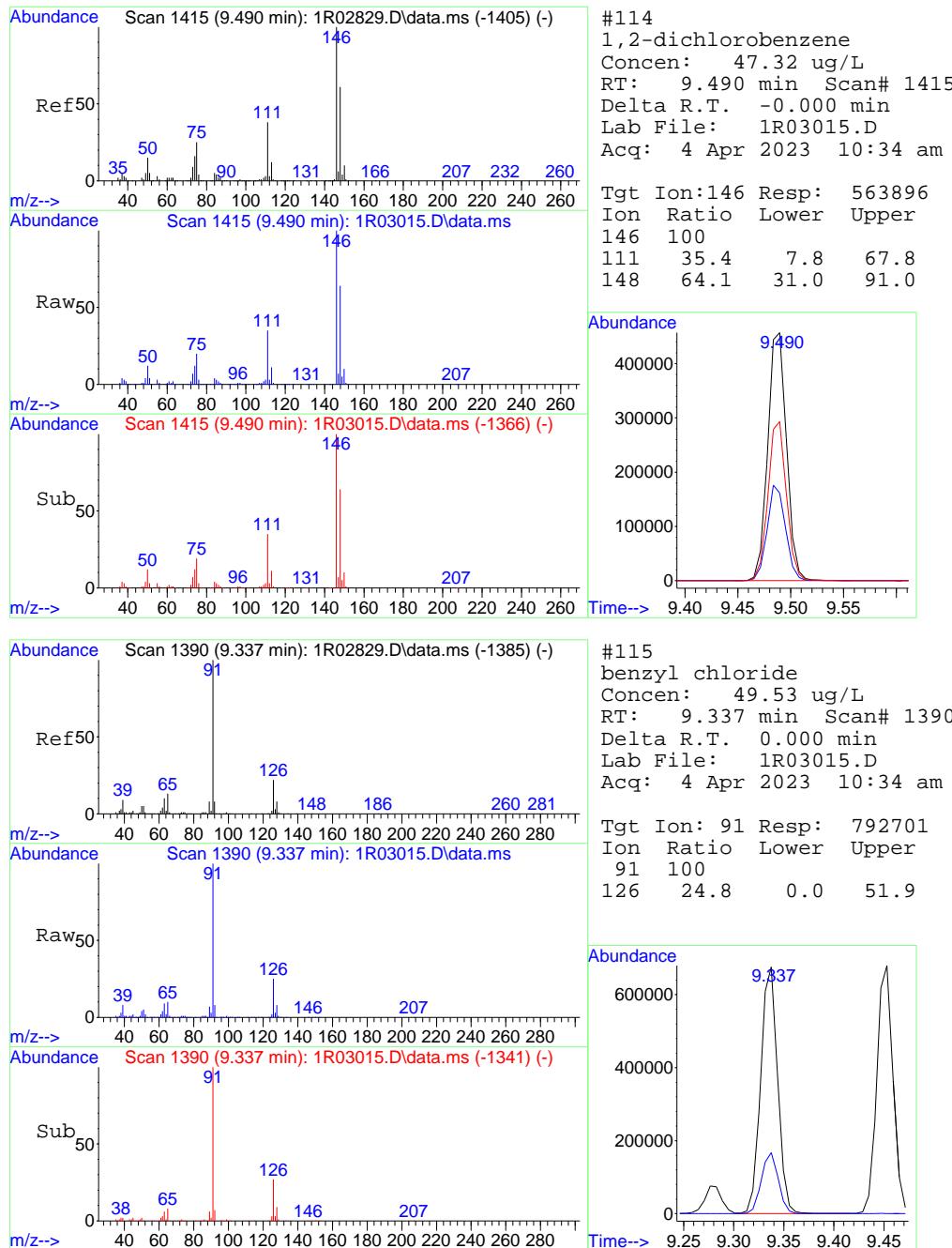


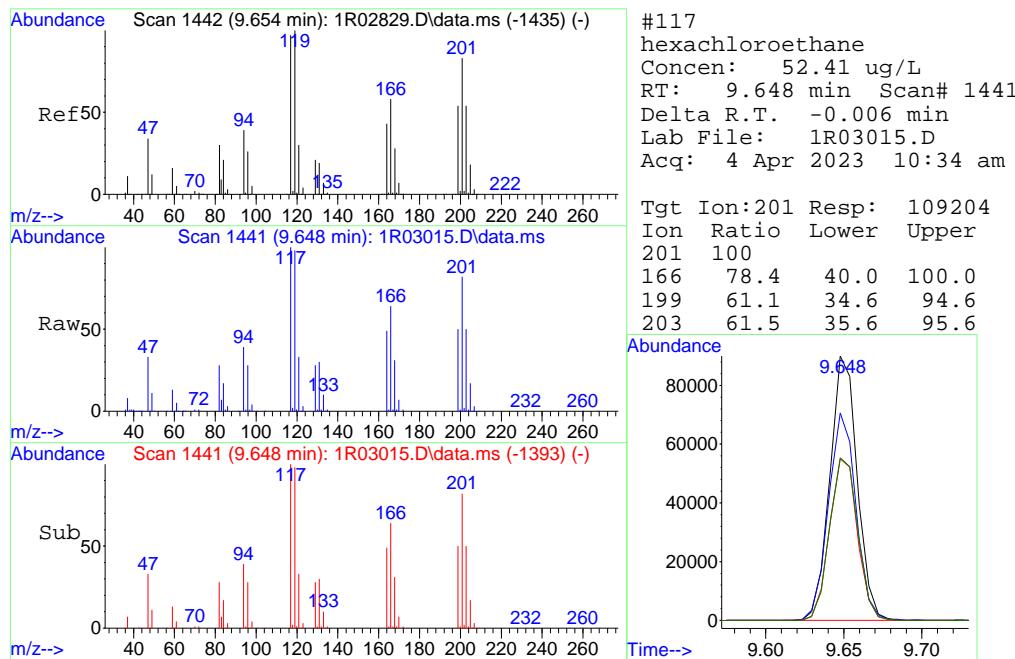
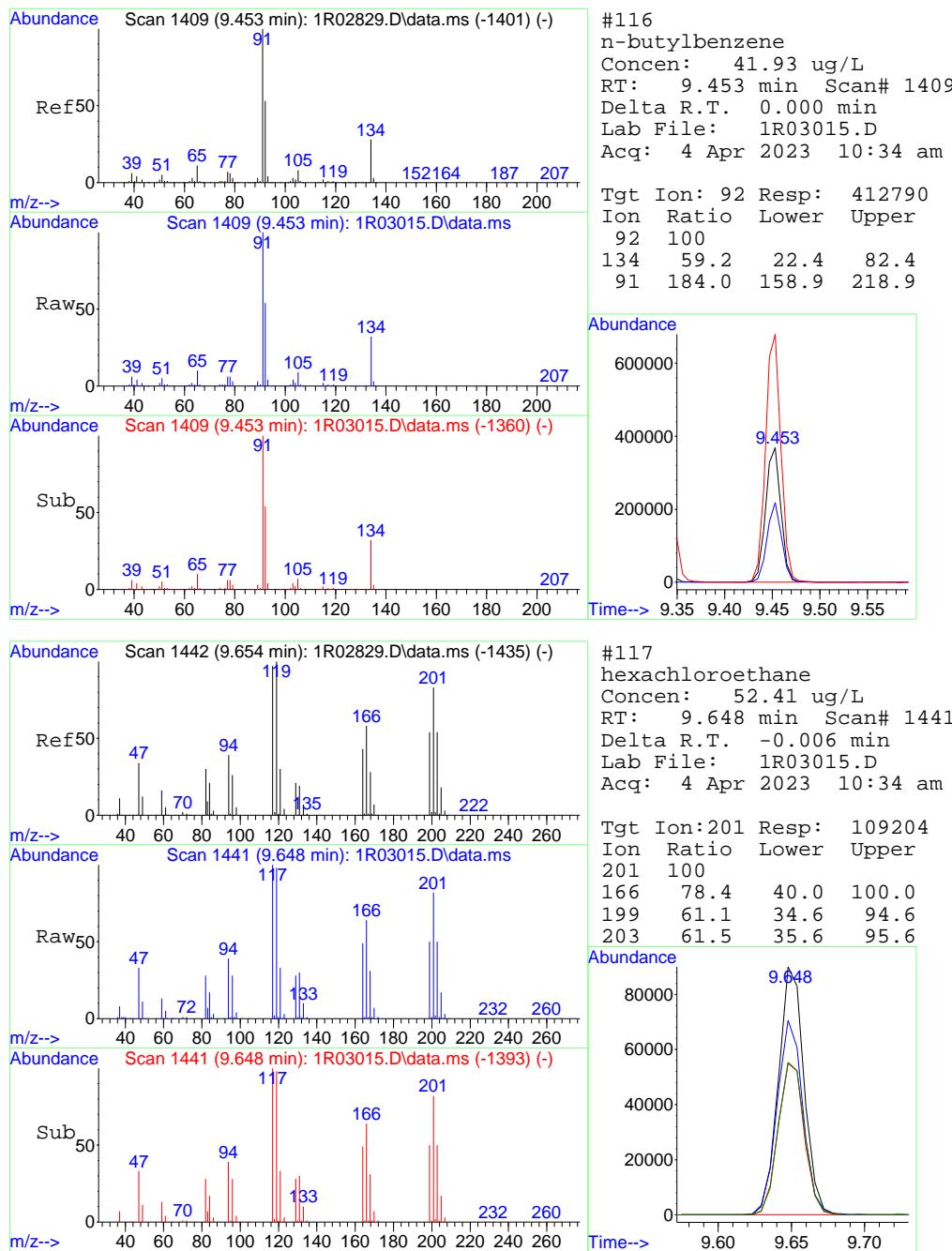


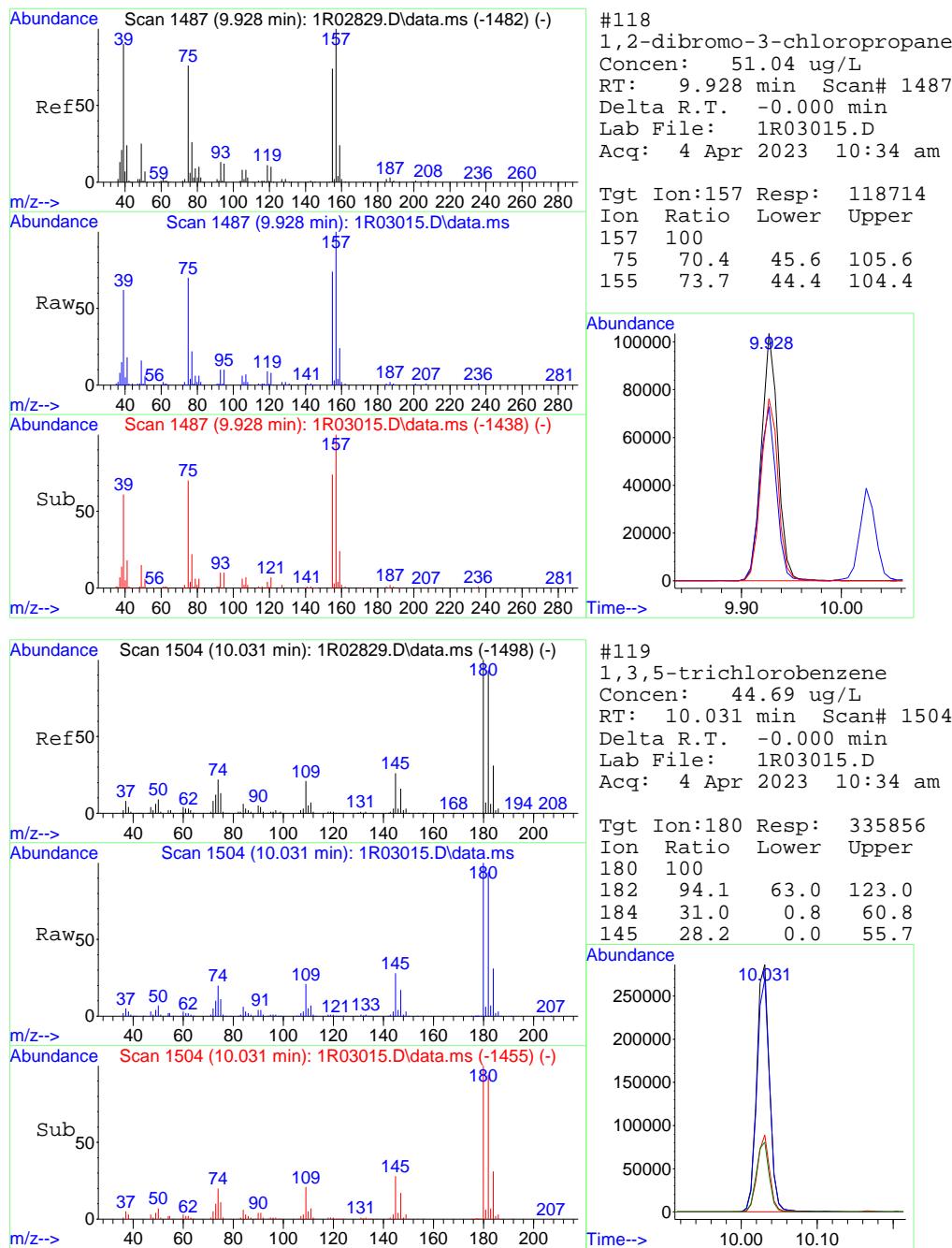


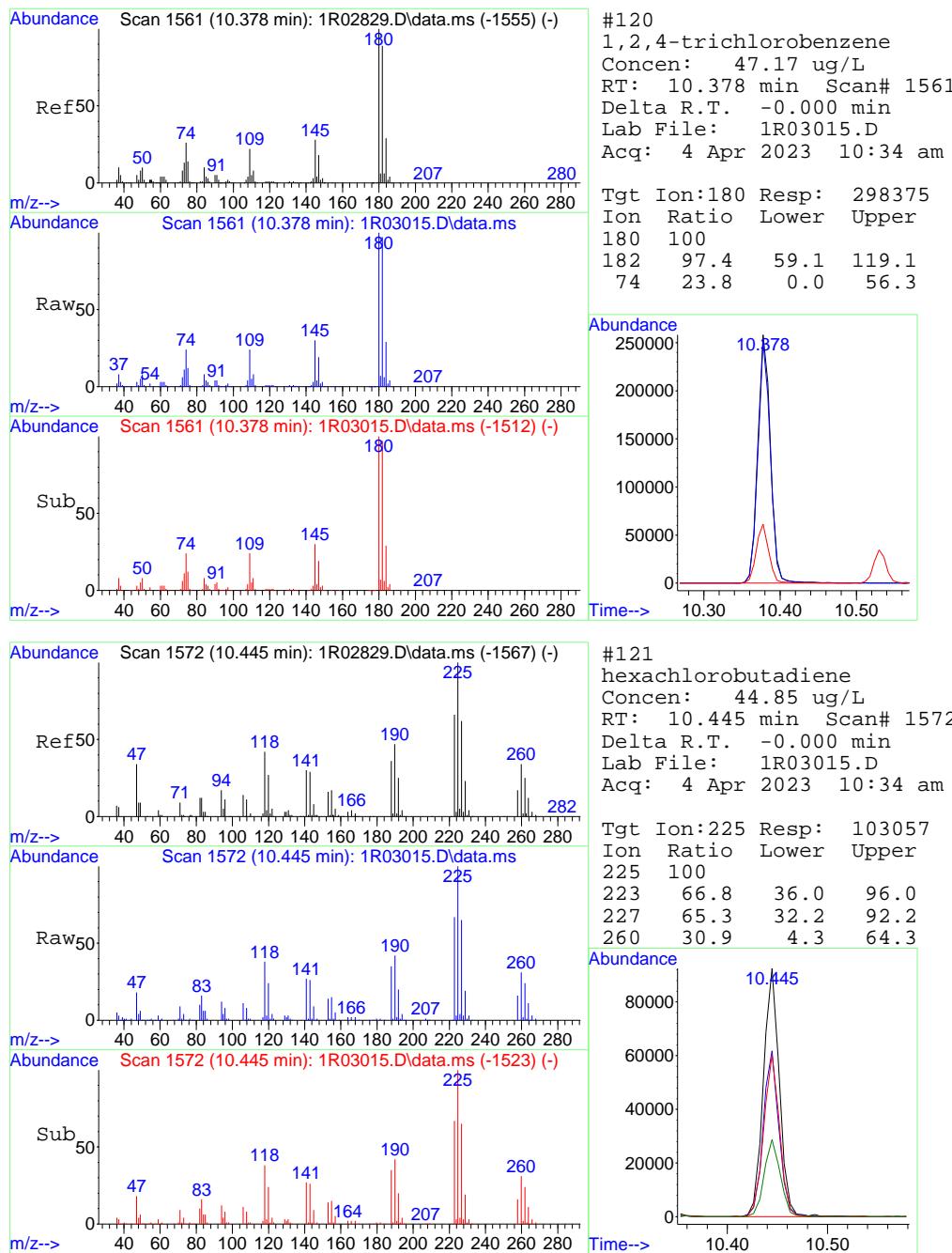


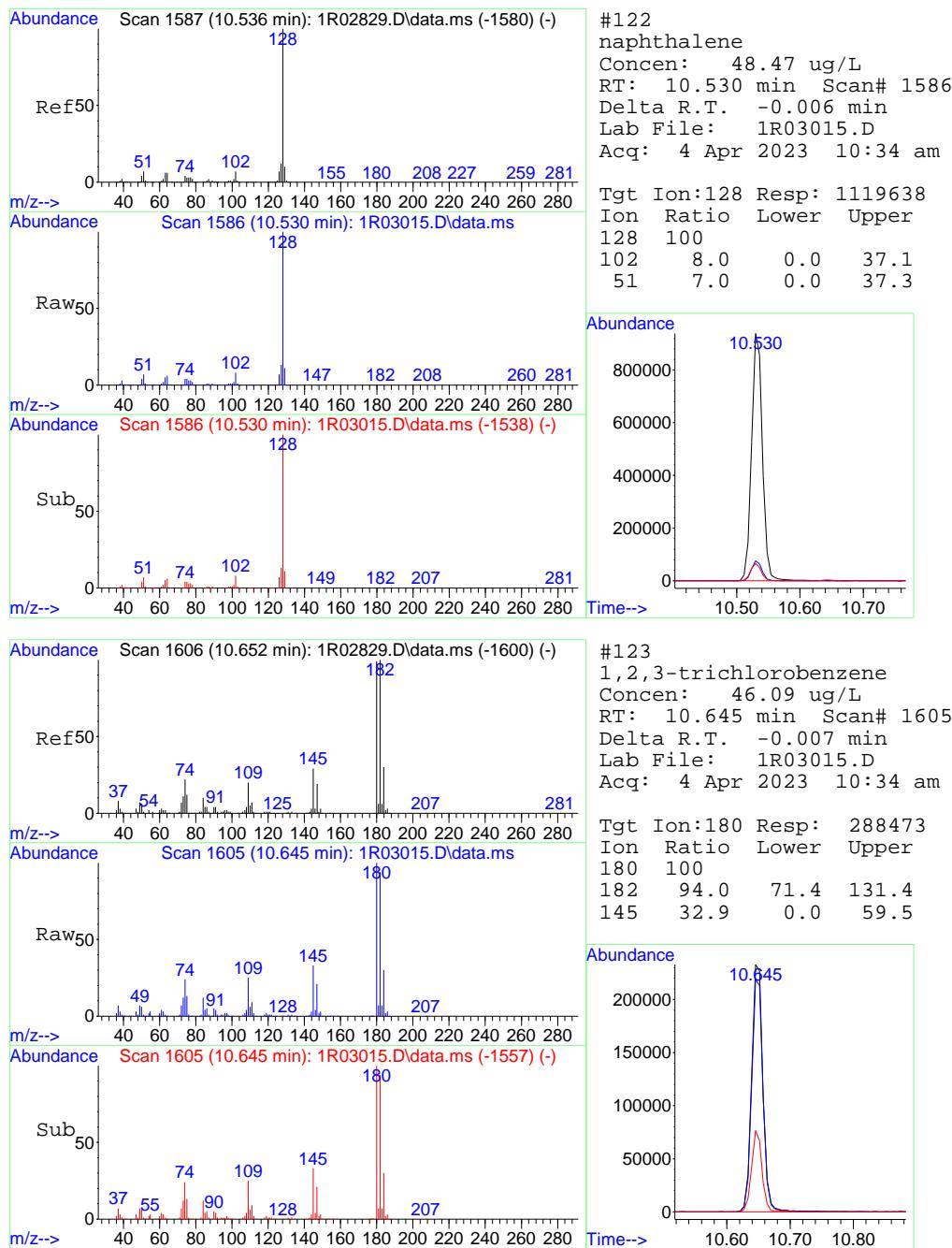


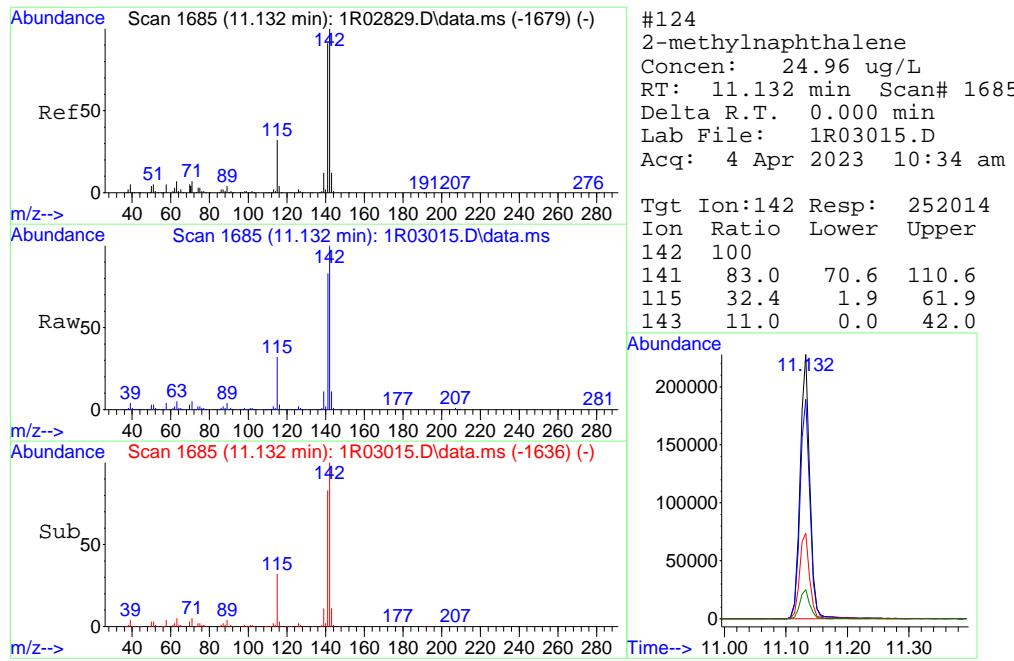












## Quantitation Report (Not Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03028.D  
 Acq On : 4 Apr 2023 4:11 pm  
 Operator : nickw  
 Sample : jd62953-1ms  
 Misc : MS67936,V1R0100,5,,,4  
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Apr 05 09:49:58 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rx1624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.491	65	507065	500.00	ug/L	0.00
5) pentafluorobenzene	5.024	168	639724	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.651	114	974157	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	955113	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	501250	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) dibromofluoromethane (s)	5.018	113	294874	53.03	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 106.06%			
53) 1,2-dichloroethane-d4 (s)	5.298	65	279744	49.39	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery = 98.78%			
74) toluene-d8 (s)	6.800	98	1245931	47.24	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 94.48%			
98) 4-bromofluorobenzene (s)	8.565	174	411533	50.17	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 100.34%			

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) ethanol	2.688	45	537287	5873.94	ug/L	98
3) tertiary butyl alcohol	3.564	59	338524	266.13	ug/L	98
4) 1,4-dioxane	6.113	88	127156	1341.14	ug/L	91
6) chlorodifluoromethane	1.654	51	229817	47.08	ug/L	96
7) dichlorodifluoromethane	1.635	85	237176	52.57	ug/L	99
8) chloromethane	1.800	50	261845	37.54	ug/L	98
9) vinyl chloride	1.897	62	246350	33.45	ug/L	96
10) 1,3-butadiene	1.927	54	201622	36.63	ug/L	95
11) bromomethane	2.183	94	74271	18.01	ug/L	98
12) chloroethane	2.286	64	140801	37.95	ug/L	99
13) trichlorofluoromethane	2.530	101	298148	52.80	ug/L	94
14) ethyl ether	2.797	74	113847	48.94	ug/L	94
15) acrolein	2.937	56	62620	46.77	ug/L	88
16) freon 113	3.035	151	150815	47.38	ug/L	98
17) 1,1-dichloroethene	3.035	96	173965	48.10	ug/L	97
18) acetone	3.065	58	156748	179.34	ug/L	99
19) acetonitrile	3.327	41	622512	528.71	ug/L	97
20) iodomethane	3.181	142	124300	47.94	ug/L	98
21) carbon disulfide	3.248	76	477001	44.90	ug/L	99
22) methylene chloride	3.491	84	199173	50.50	ug/L	97
23) methyl acetate	3.357	74	60551	51.37	ug/L #	71
24) methyl tert butyl ether	3.728	73	563599	50.03	ug/L	98
25) trans-1,2-dichloroethene	3.740	96	195405	48.66	ug/L	94
26) hexane	3.978	56	159926	46.98	ug/L	95
27) di-isopropyl ether	4.142	45	772008	48.30	ug/L	96
28) 2-butanone	4.592	72	195758	190.55	ug/L	94
29) 1,1-dichloroethane	4.124	63	356068	47.24	ug/L	98
30) chloroprene	4.191	53	323229	49.76	ug/L	99
31) acrylonitrile	3.698	53	135464	46.42	ug/L	99
32) vinyl acetate	4.111	86	54411	52.91	ug/L #	84
33) ethyl tert-butyl ether	4.452	59	656018	51.97	ug/L	98
34) ethyl acetate	4.616	45	72451	50.29	ug/L #	76
35) 2,2-dichloropropane	4.616	77	286511	51.93	ug/L	96
36) cis-1,2-dichloroethene	4.610	96	219660	48.93	ug/L	90
37) propionitrile	4.647	54	656643	495.25	ug/L	91
38) methyl acrylate	4.659	85	54553	50.32	ug/L	98
39) methacrylonitrile	4.775	67	138167	48.21	ug/L	91
40) bromochloromethane	4.811	128	109973	51.70	ug/L #	83
41) tetrahydrofuran	4.817	42	164054	45.91	ug/L	96
42) chloroform	4.884	83	361776	49.04	ug/L	94
43) tert-Butyl Formate	4.908	59	191559	47.67	ug/L	95

## Quantitation Report (Not Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\

Data File : 1R03028.D

Acq On : 4 Apr 2023 4:11 pm

Operator : nickw

Sample : jd62953-1ms

Misc : MS67936,V1R0100,5,,,4

ALS Vial : 17 Sample Multiplier: 1

Quant Time: Apr 05 09:49:58 2023

Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M

Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um

QLast Update : Thu Mar 30 09:13:25 2023

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1,1-trichloroethane	5.030	97	311462	48.99	ug/L	86
46) cyclohexane	5.091	84	287048	48.64	ug/L	88
47) isobutyl alcohol	5.207	43	303681	563.02	ug/L	98
48) 1,1-dichloropropene	5.158	75	269305	50.55	ug/L	97
49) carbon tetrachloride	5.158	117	277795	52.41	ug/L	97
50) tert-amyl alcohol	5.310	73	183011	286.03	ug/L	94
51) isopropyl acetate	5.334	87	72828	51.14	ug/L #	72
54) n-butyl alcohol	5.742	56	922127	2521.54	ug/L	99
55) 2,2,4-trimethylpentane	5.407	57	587864	43.50	ug/L	98
56) benzene	5.322	78	884752	46.07	ug/L	99
57) tert-amyl methyl ether	5.413	73	626840	45.90	ug/L	95
58) heptane	5.541	57	130013	42.57	ug/L	94
59) 1,2-dichloroethane	5.365	62	284395	46.49	ug/L	95
60) ethyl acrylate	5.894	55	428816	43.34	ug/L	99
61) trichloroethene	5.851	95	222336	45.88	ug/L	98
62) 2-chloroethyl vinyl ether	6.466	63	28495	6.84	ug/L	98
63) methyl methacrylate	6.083	100	78248	47.77	ug/L #	74
64) methylcyclohexane	6.040	83	328010	42.73	ug/L	98
65) 1,2-dichloropropane	6.064	63	218300	41.27	ug/L	98
66) dibromomethane	6.131	93	142926	43.72	ug/L	99
67) bromodichloromethane	6.265	83	272149	46.70	ug/L	97
68) 2-nitropropane	6.429	41	131196	47.76	ug/L	95
69) epichlorohydrin	6.508	57	242063	243.83	ug/L	98
70) cis-1,3-dichloropropene	6.594	75	341417	45.53	ug/L	96
71) 4-methyl-2-pentanone	6.691	58	639812	175.28	ug/L	95
72) isoamyl alcohol	6.721	70	334473	983.67	ug/L	98
75) toluene	6.849	92	3172100	256.39	ug/L	99
76) ethyl methacrylate	7.025	69	317829	47.09	ug/L	99
77) trans-1,3-dichloropropene	7.007	75	325232	49.13	ug/L	92
78) 1,1,2-trichloroethane	7.153	83	173043	44.73	ug/L	94
79) tetrachloroethene	7.220	164	214223	43.96	ug/L	97
80) 2-hexanone	7.287	58	717171	183.51	ug/L	97
81) 1,3-dichloropropane	7.269	76	319392	46.22	ug/L	98
82) butyl acetate	7.354	56	228657	44.66	ug/L	95
83) dibromochloromethane	7.421	129	253478	52.01	ug/L	97
84) 1,2-dibromoethane	7.512	107	239277	51.68	ug/L	97
85) n-butyl ether	7.865	57	967181	44.11	ug/L	100
86) chlorobenzene	7.835	112	641181	46.87	ug/L	95
87) 1,1,1,2-tetrachloroethane	7.889	131	235129	48.62	ug/L	98
88) ethylbenzene	7.889	91	1977730	86.99	ug/L	98
89) m,p-xylene	7.975	106	1557826	176.51	ug/L	94
90) o-xylene	8.224	91	1976417	112.69	ug/L	95
91) styrene	8.230	104	751059	51.02	ug/L	93
92) butyl acrylate	8.163	55	538792	46.86	ug/L	99
93) n-amyl acetate	8.297	70	191145	43.42	ug/L	92
94) isopropylbenzene	8.449	105	1084920	52.71	ug/L	97
95) bromoform	8.364	173	199793	51.58	ug/L	96
96) cis-1,4-dichloro-2-butene	8.486	88	144341	50.04	ug/L	97
99) 1,1,2,2-tetrachloroethane	8.638	83	341553	45.33	ug/L	98
100) trans-1,4-dichloro-2-b...	8.656	53	112529	44.42	ug/L	89
101) 1,2,3-trichloropropane	8.680	110	109080	48.84	ug/L	87
102) bromobenzene	8.662	156	283333	47.82	ug/L	84
103) n-propylbenzene	8.705	91	1251865	48.86	ug/L	97
104) 2-chlorotoluene	8.771	126	265486	50.09	ug/L	88
105) 4-chlorotoluene	8.844	91	712631	45.27	ug/L	95
106) 1,3,5-trimethylbenzene	8.808	105	886259	50.24	ug/L	98
107) tert-butylbenzene	9.003	119	730318	47.41	ug/L	97
108) 1,2,4-trimethylbenzene	9.033	105	1535392	85.53	ug/L	98

## Quantitation Report (Not Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03028.D

Acq On : 4 Apr 2023 4:11 pm  
 Operator : nickw  
 Sample : jd62953-1ms  
 Misc : MS67936,V1R0100,5,,,4  
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Apr 05 09:49:58 2023

Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M

Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um

QLast Update : Thu Mar 30 09:13:25 2023

Response via : Initial Calibration

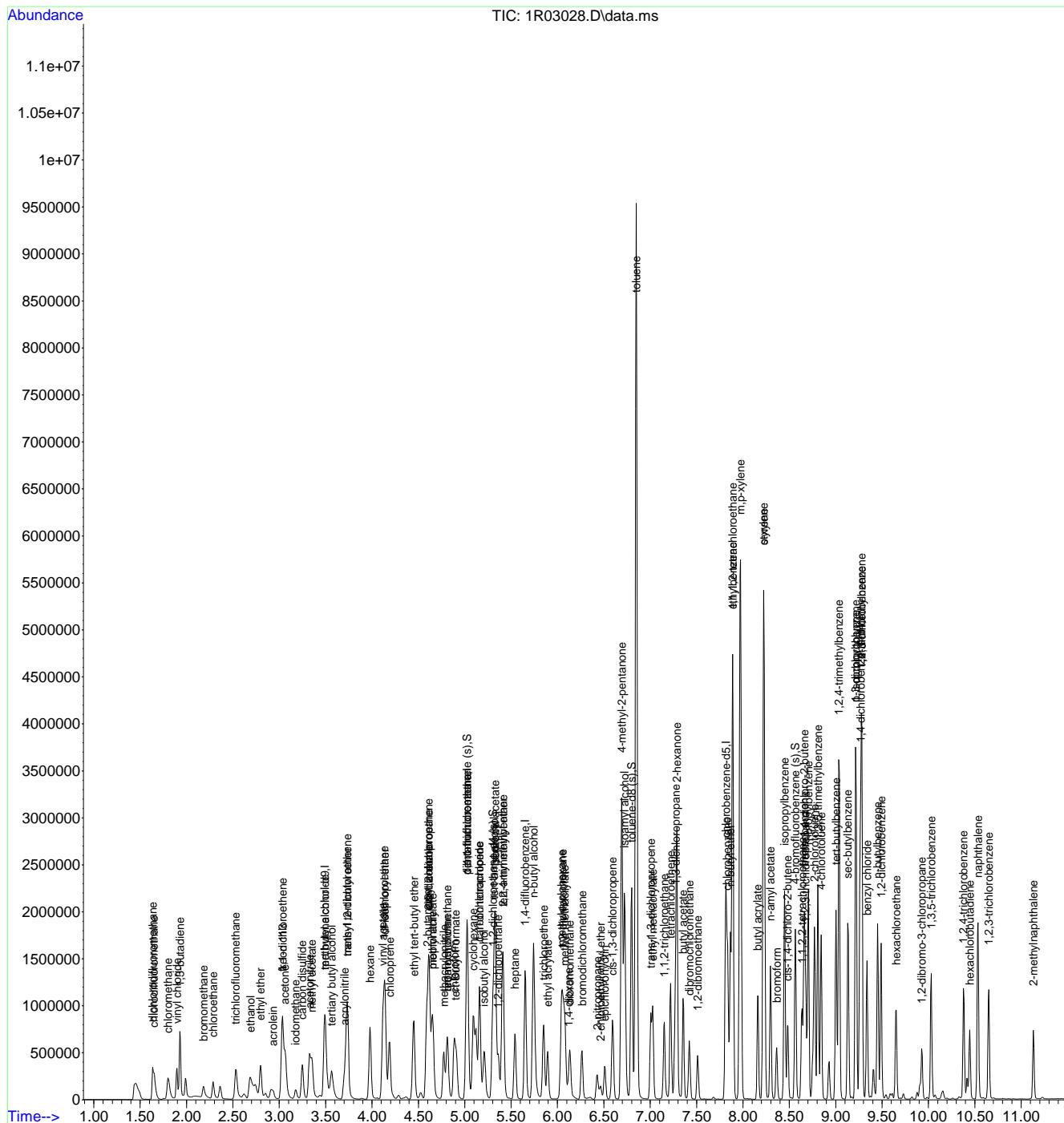
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
109) sec-butylbenzene	9.130	105	1018214	48.53	ug/L	96
110) p-isopropyltoluene	9.216	119	921100	47.92	ug/L	99
111) 1,2,3-trimethylbenzene	9.282	105	1033591	54.02	ug/L	96
112) 1,3-dichlorobenzene	9.222	146	552128	48.44	ug/L	99
113) 1,4-dichlorobenzene	9.276	146	536429	46.07	ug/L	98
114) 1,2-dichlorobenzene	9.489	146	541409	51.76	ug/L	98
115) benzyl chloride	9.337	91	726698	51.73	ug/L	95
116) n-butylbenzene	9.453	92	395014	45.71	ug/L	98
117) hexachloroethane	9.654	201	107246	58.63	ug/L	96
118) 1,2-dibromo-3-chloropr...	9.927	157	110777	54.26	ug/L	96
119) 1,3,5-trichlorobenzene	10.031	180	322479	48.88	ug/L	97
120) 1,2,4-trichlorobenzene	10.378	180	286358	51.57	ug/L	94
121) hexachlorobutadiene	10.444	225	100988	50.07	ug/L	96
122) naphthalene	10.536	128	1171779	57.79	ug/L	98
123) 1,2,3-trichlorobenzene	10.651	180	281951	51.32	ug/L	94
124) 2-methylnaphthalene	11.132	142	253134	28.57	ug/L	98

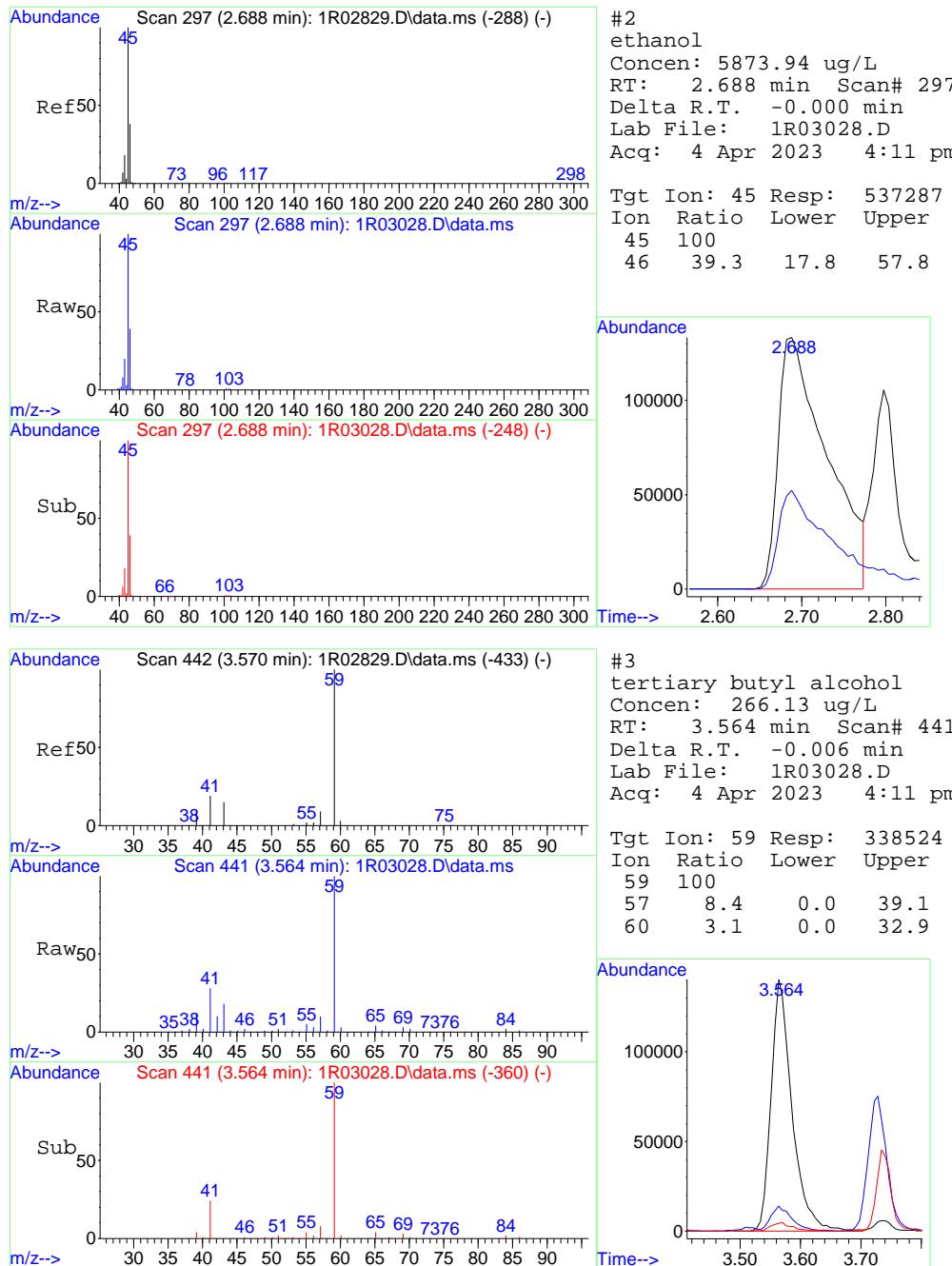
(#) = qualifier out of range (m) = manual integration (+) = signals summed

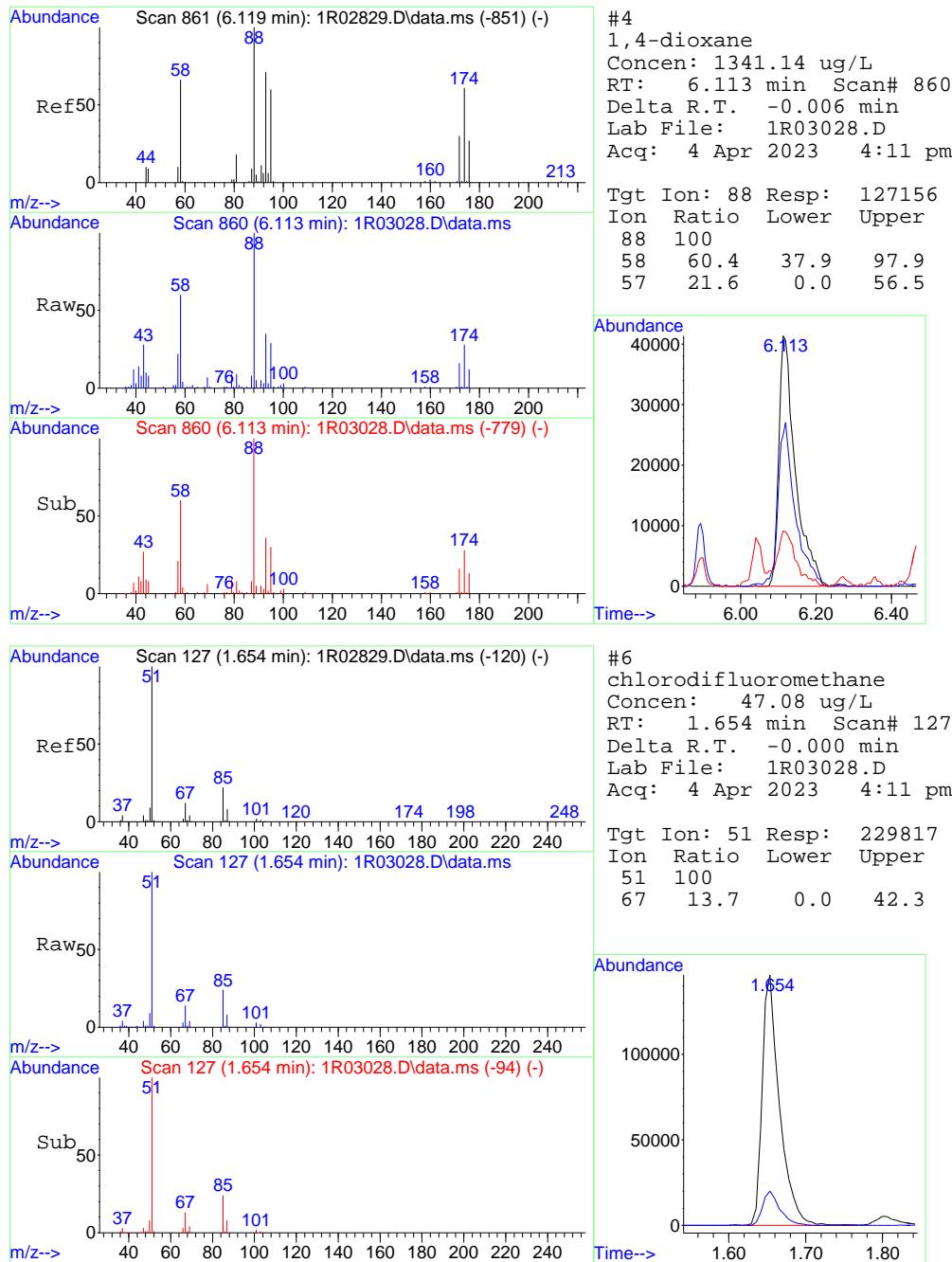
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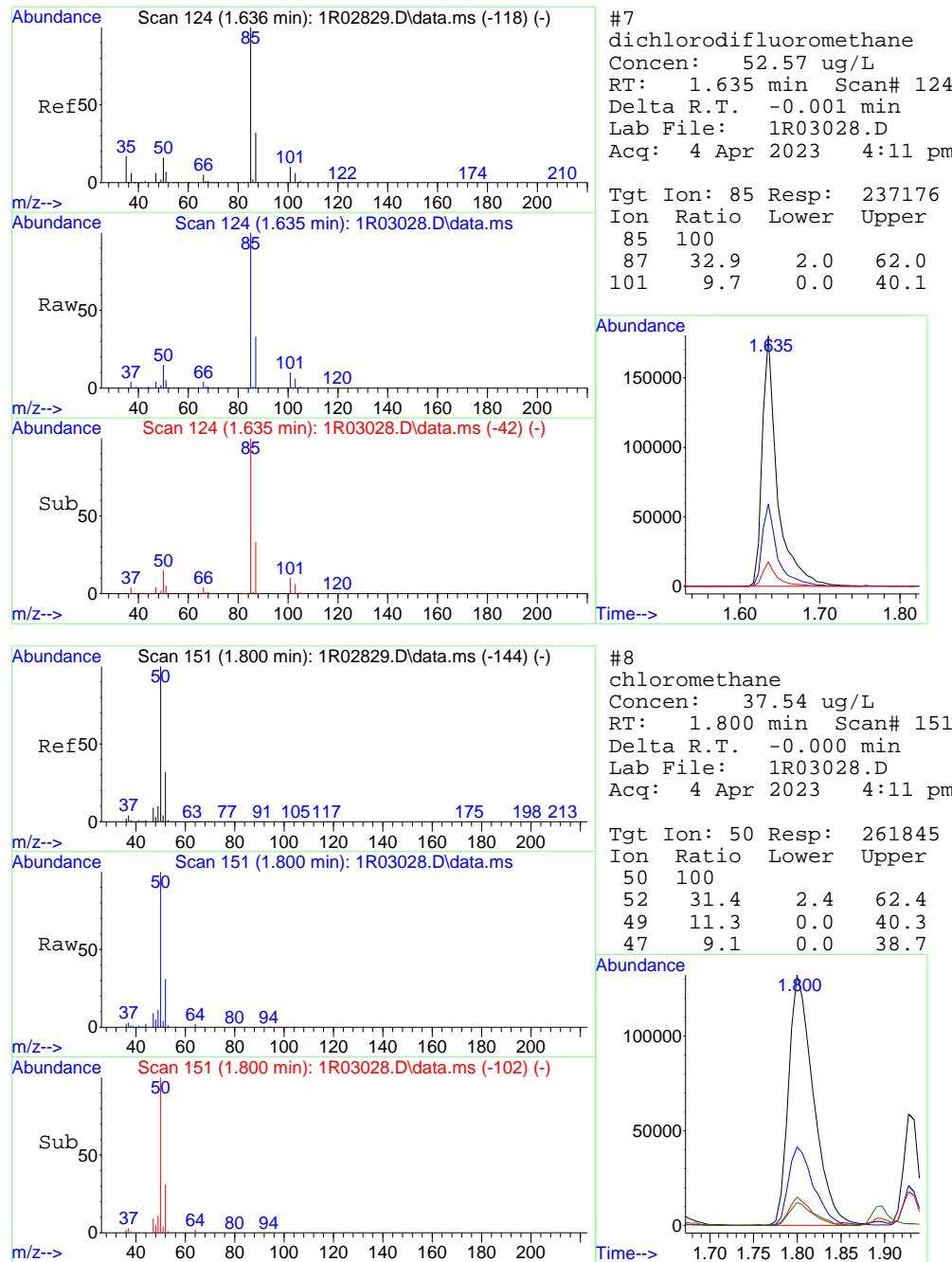
Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03028.D  
 Acq On : 4 Apr 2023 4:11 pm  
 Operator : nickw  
 Sample : jd62953-1ms  
 Misc : MS67936,V1R0100,5,,,4  
 ALS Vial : 17 Sample Multiplier: 1

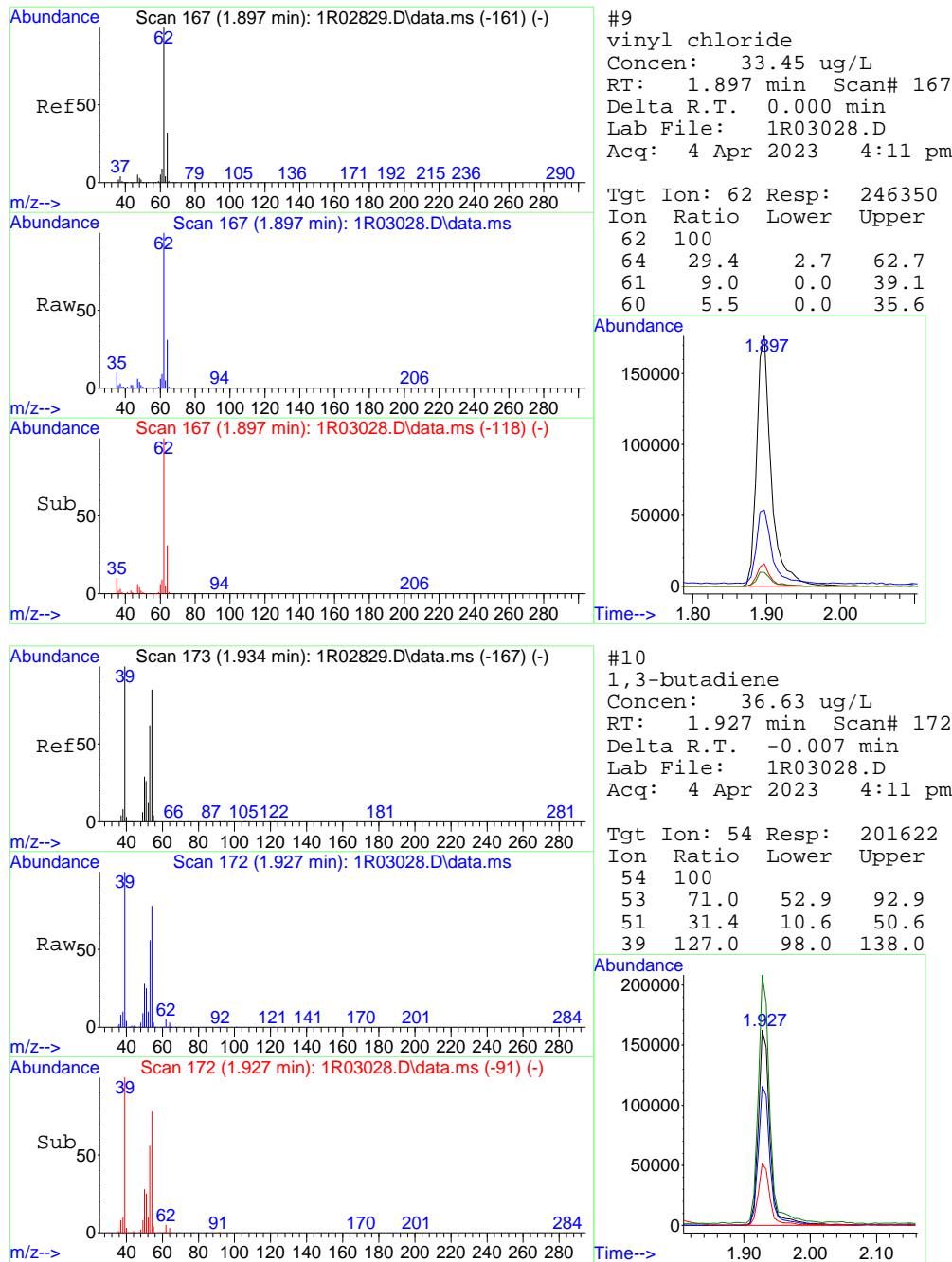
Quant Time: Apr 05 09:49:58 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

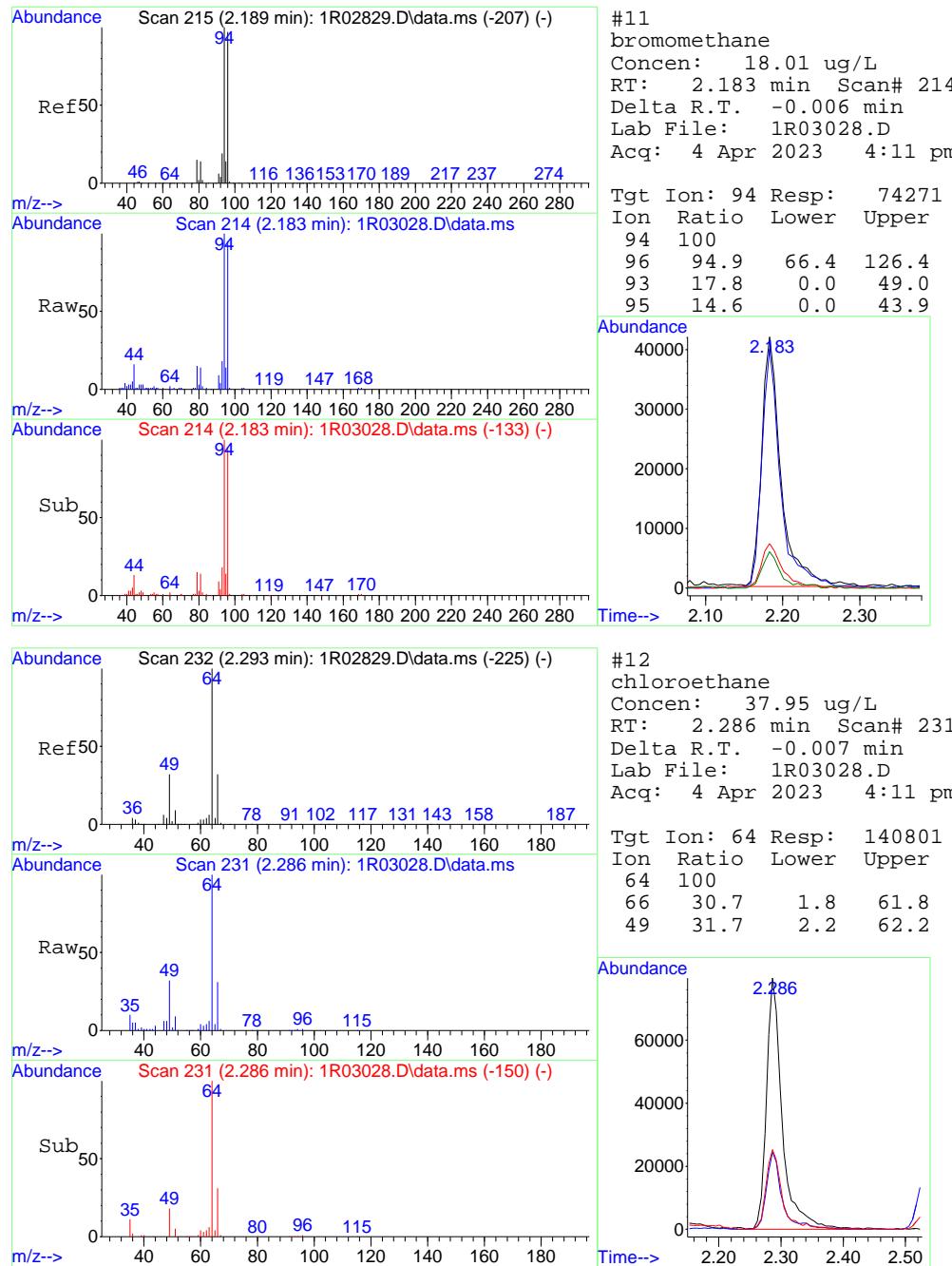


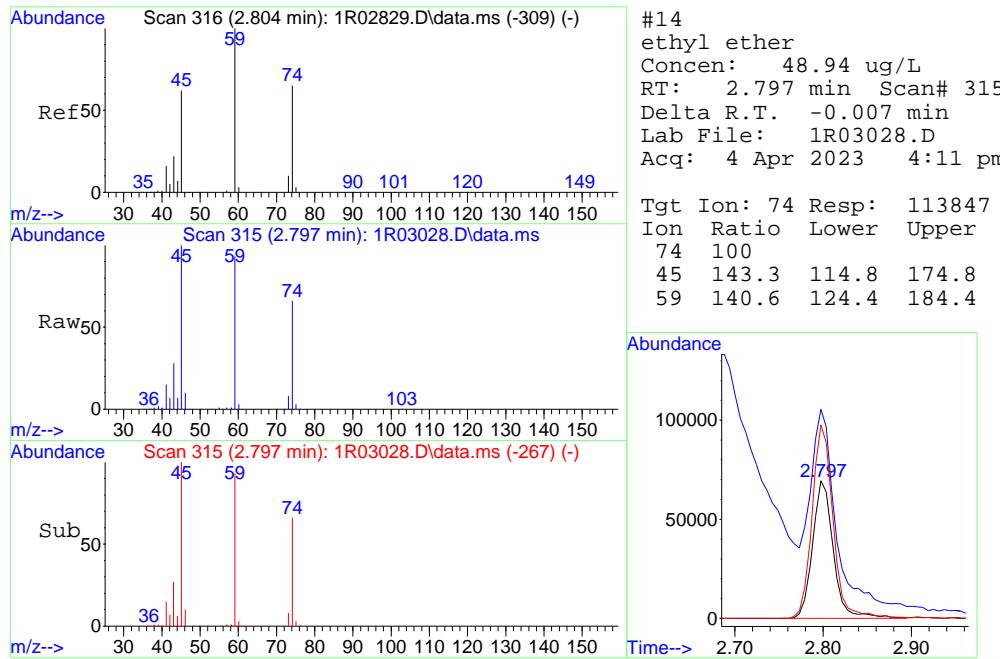
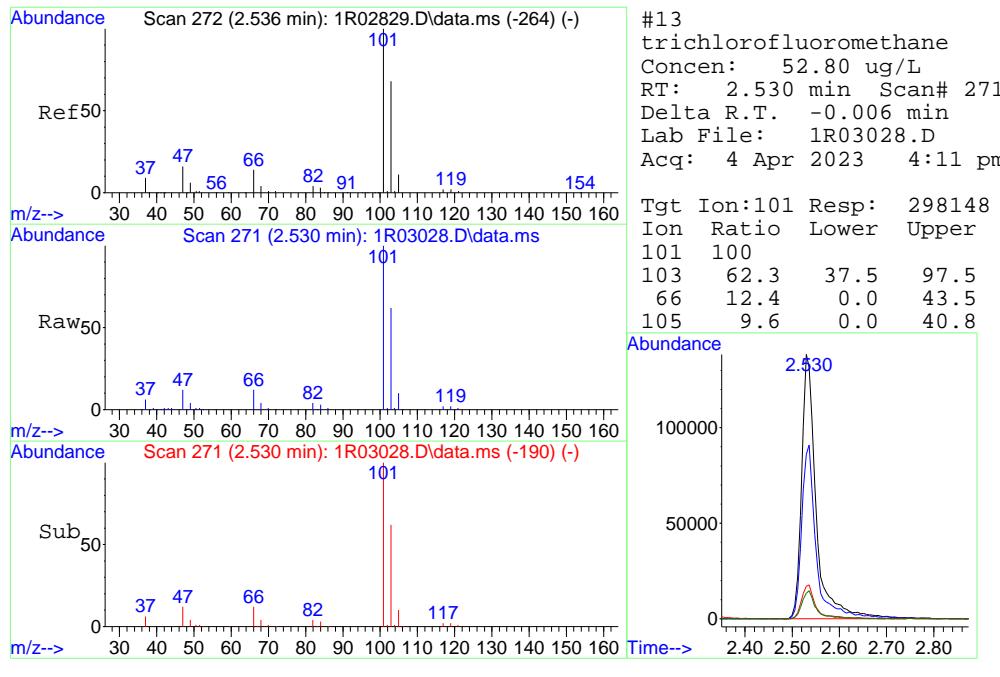


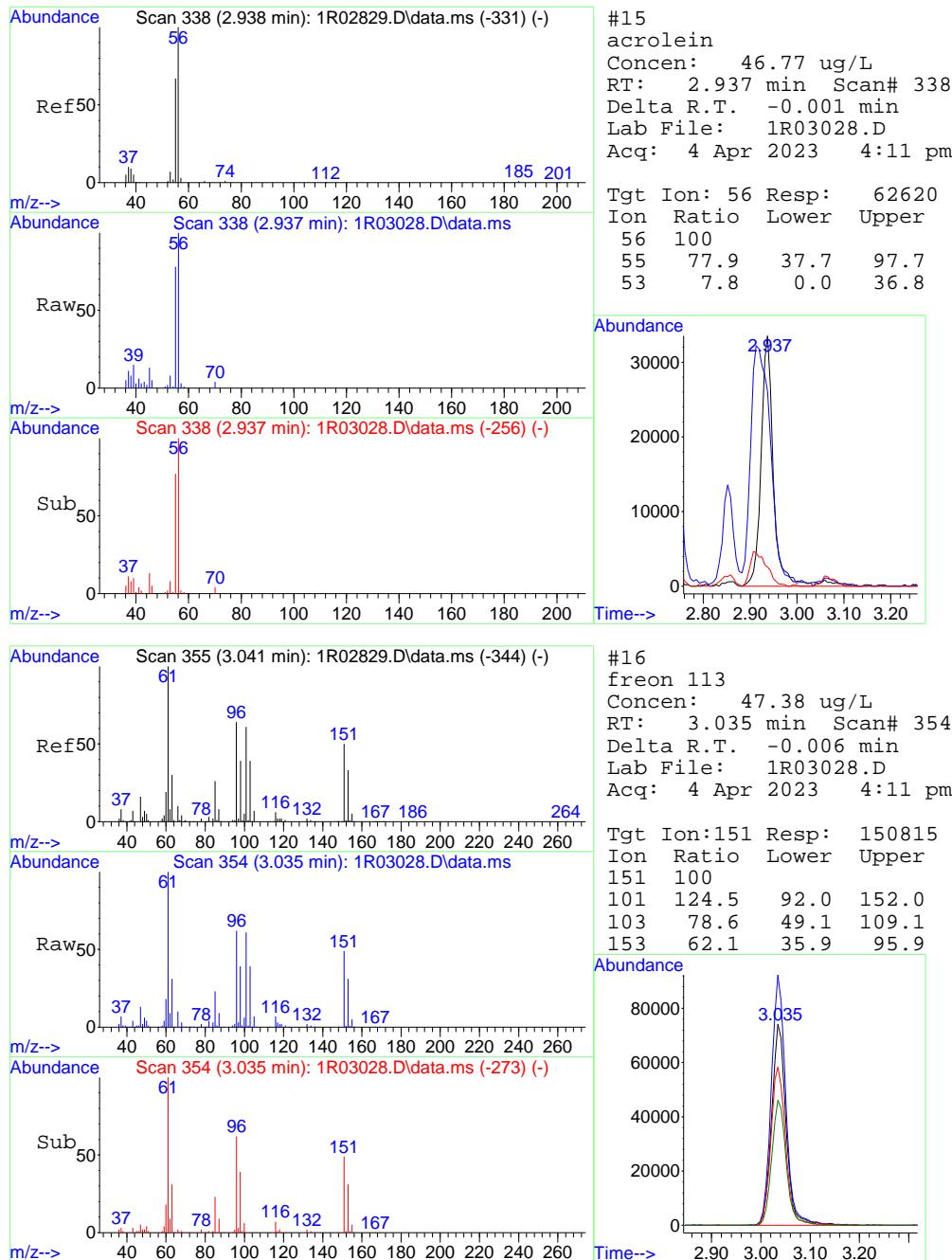


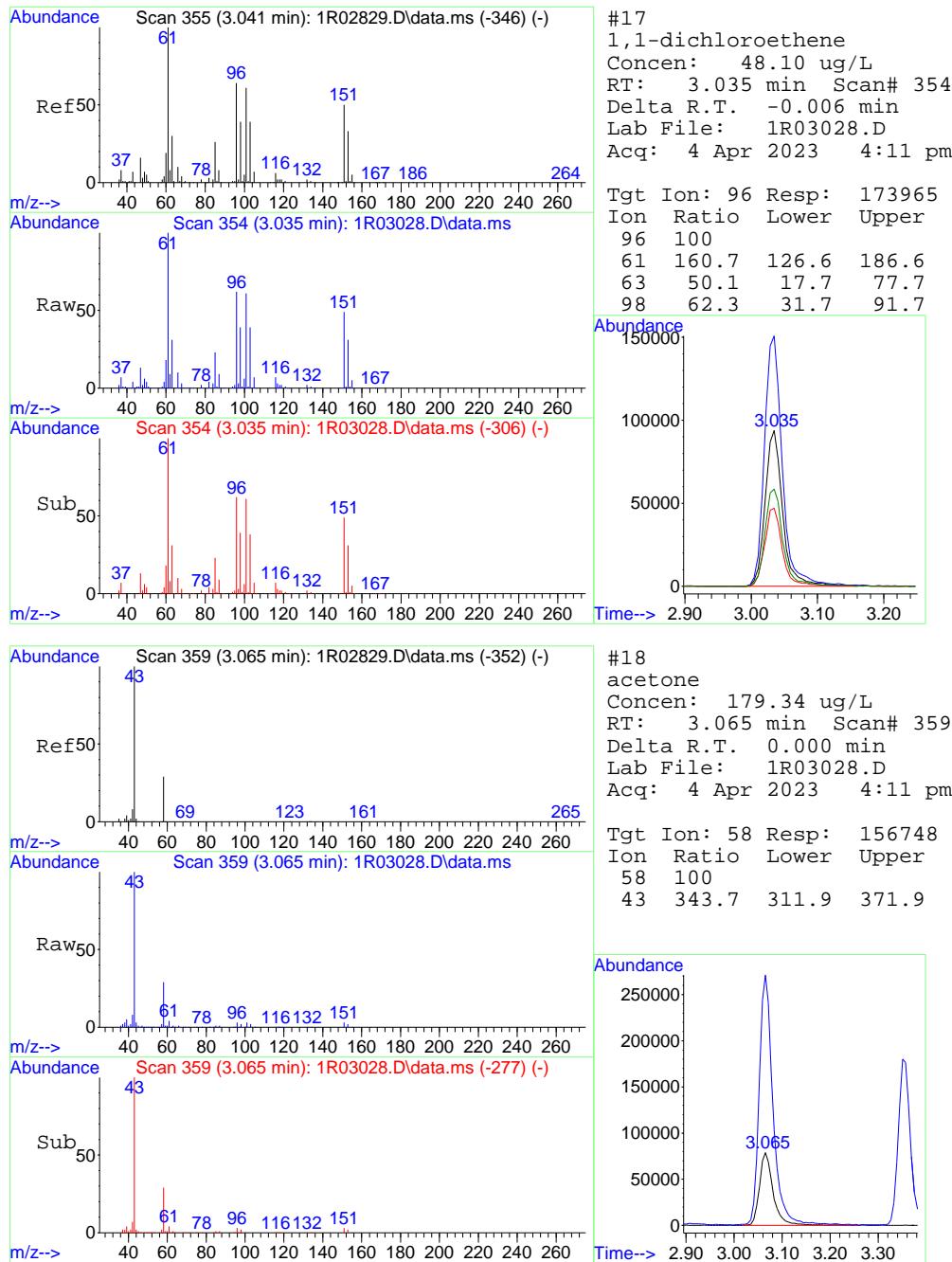


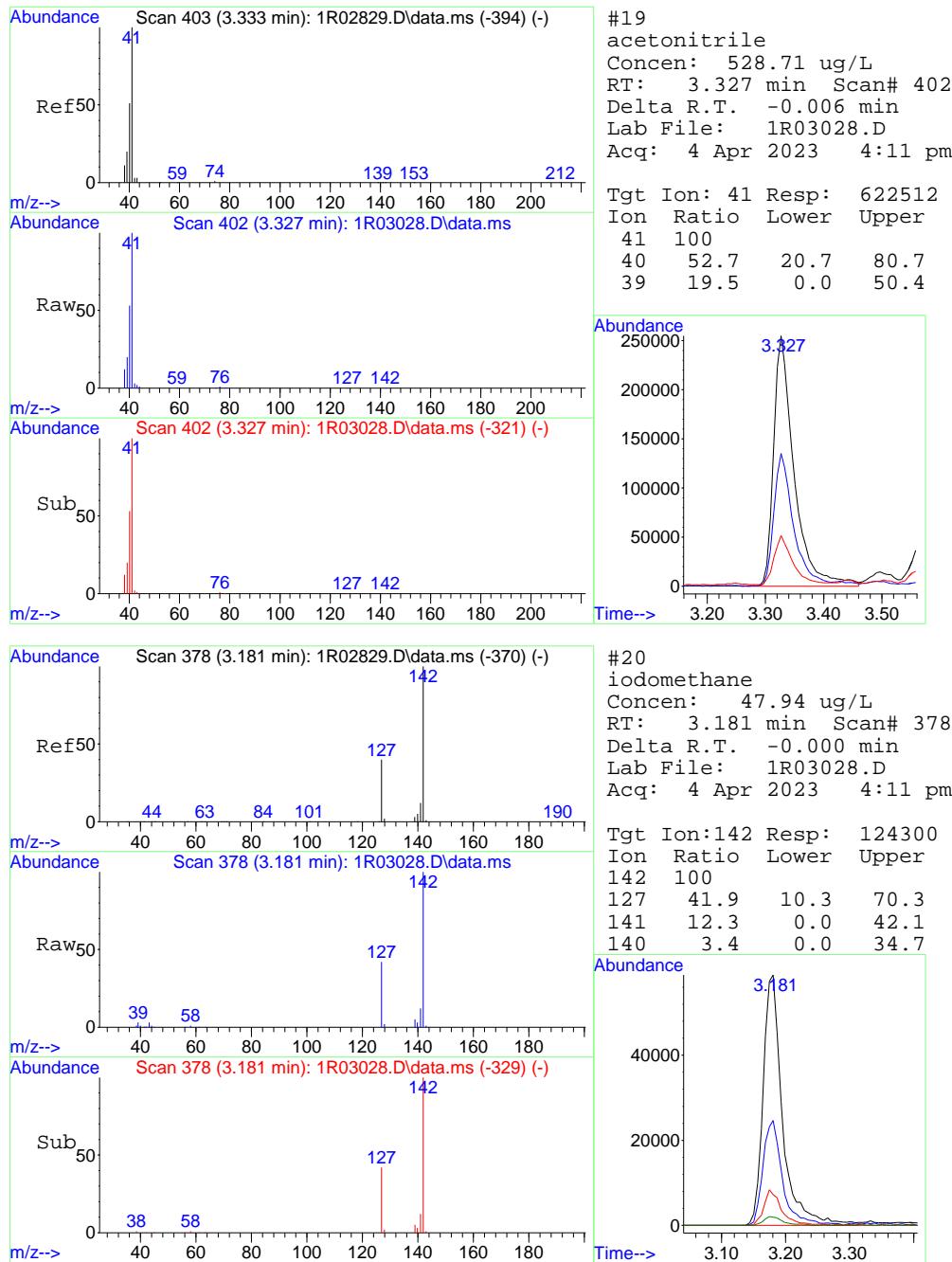


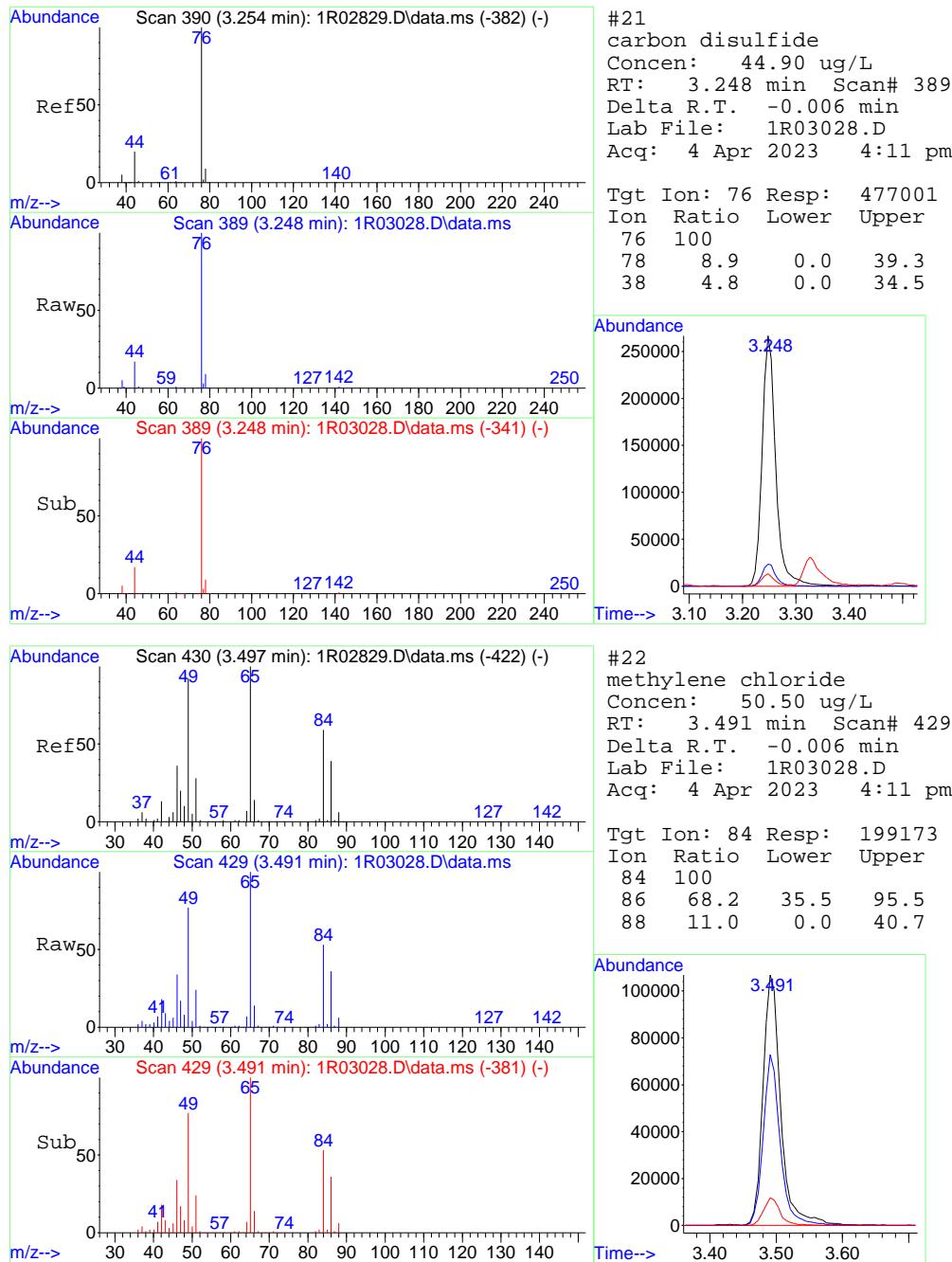


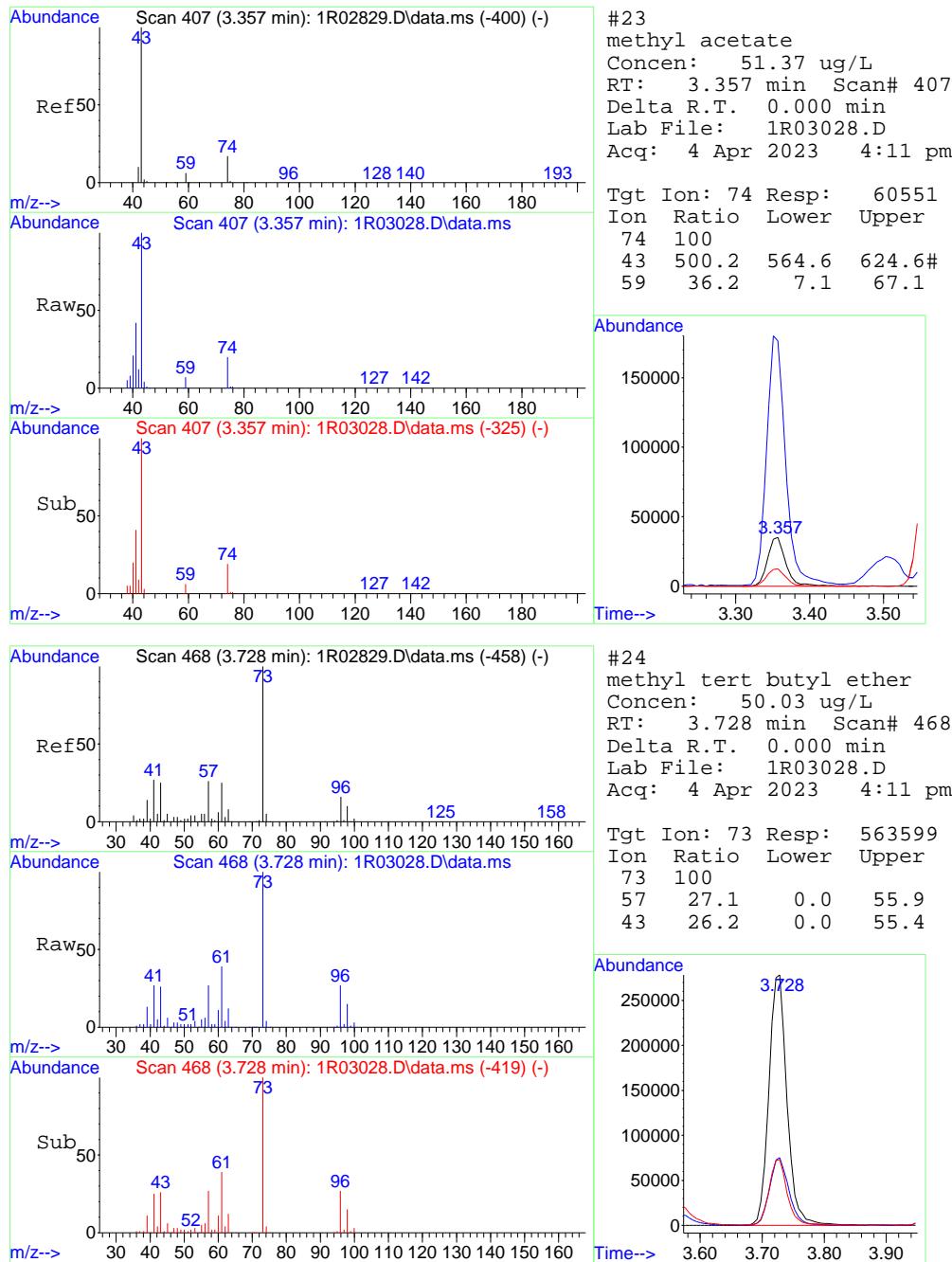


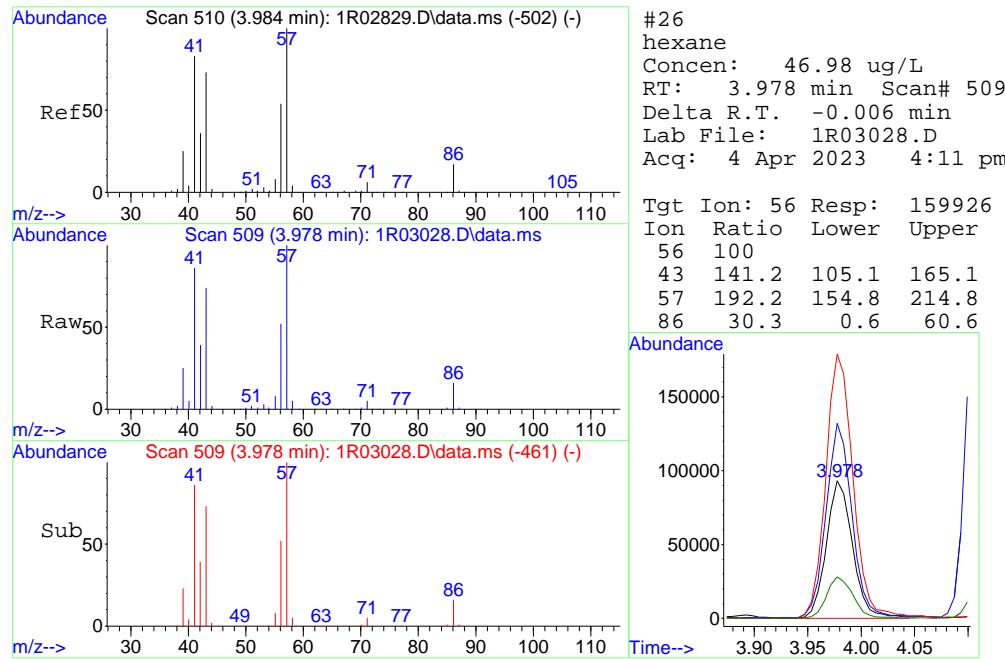
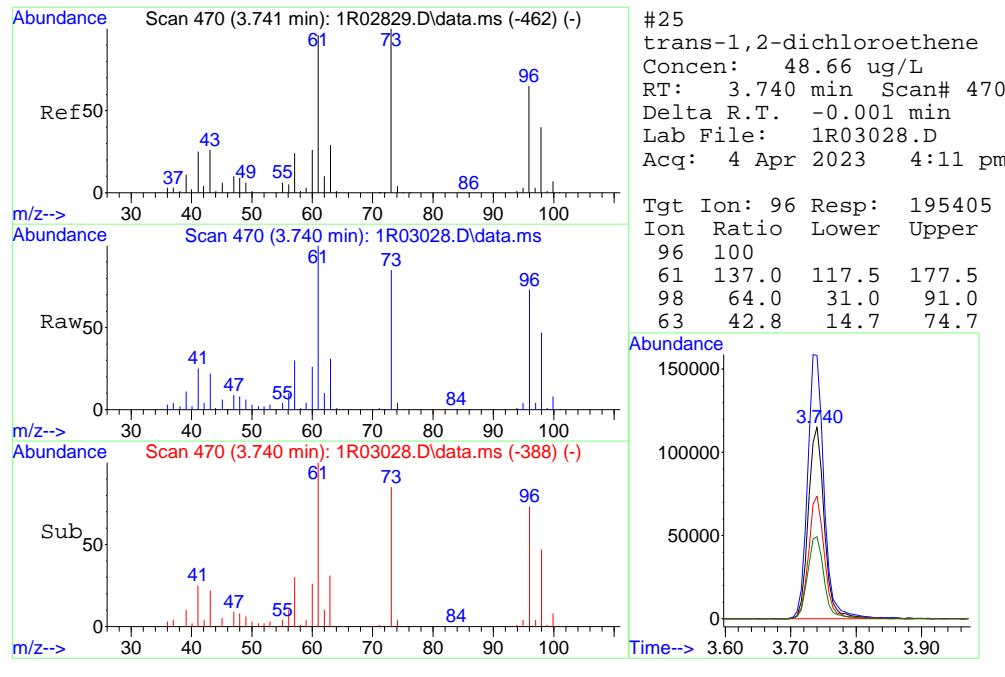


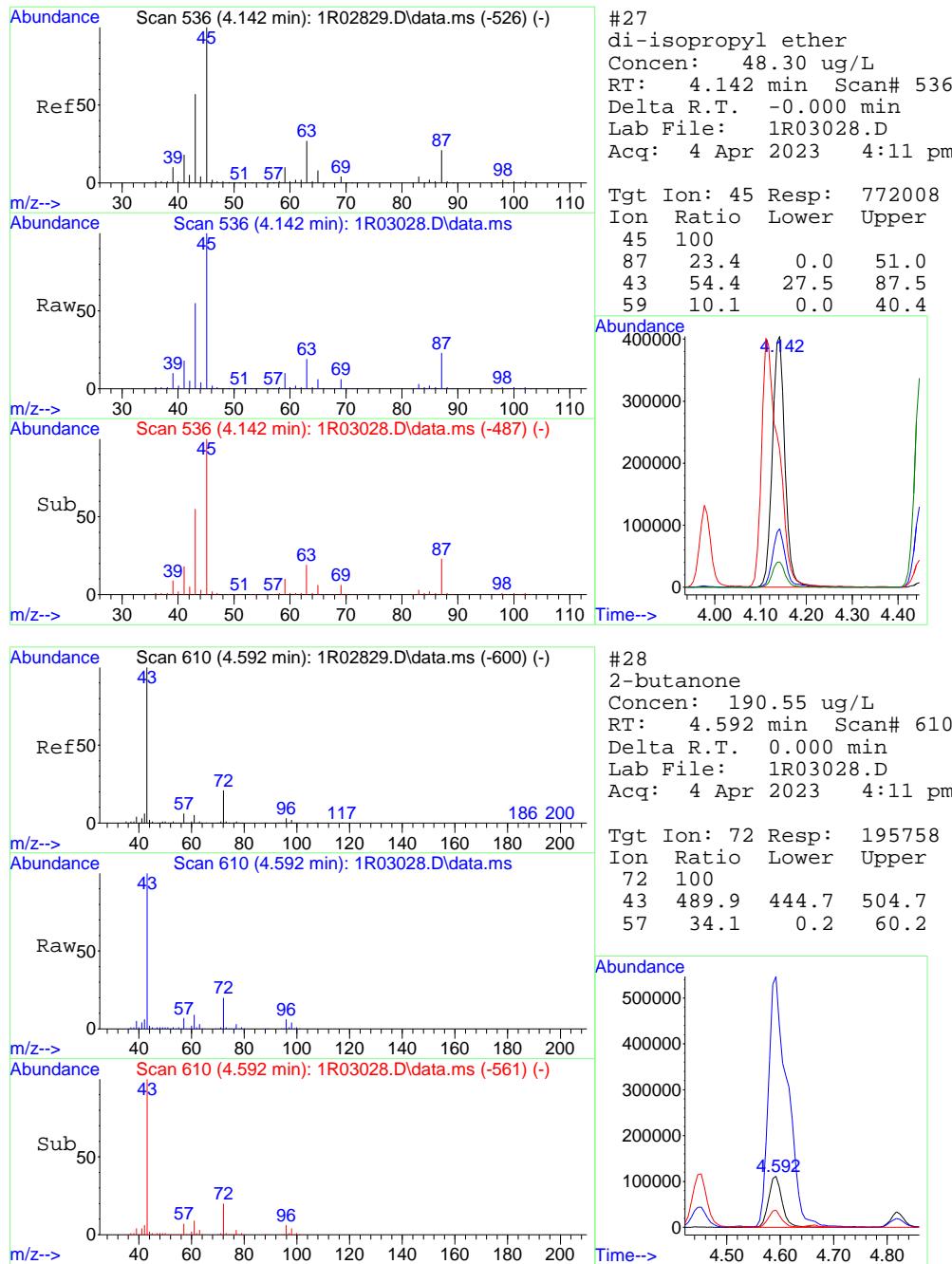


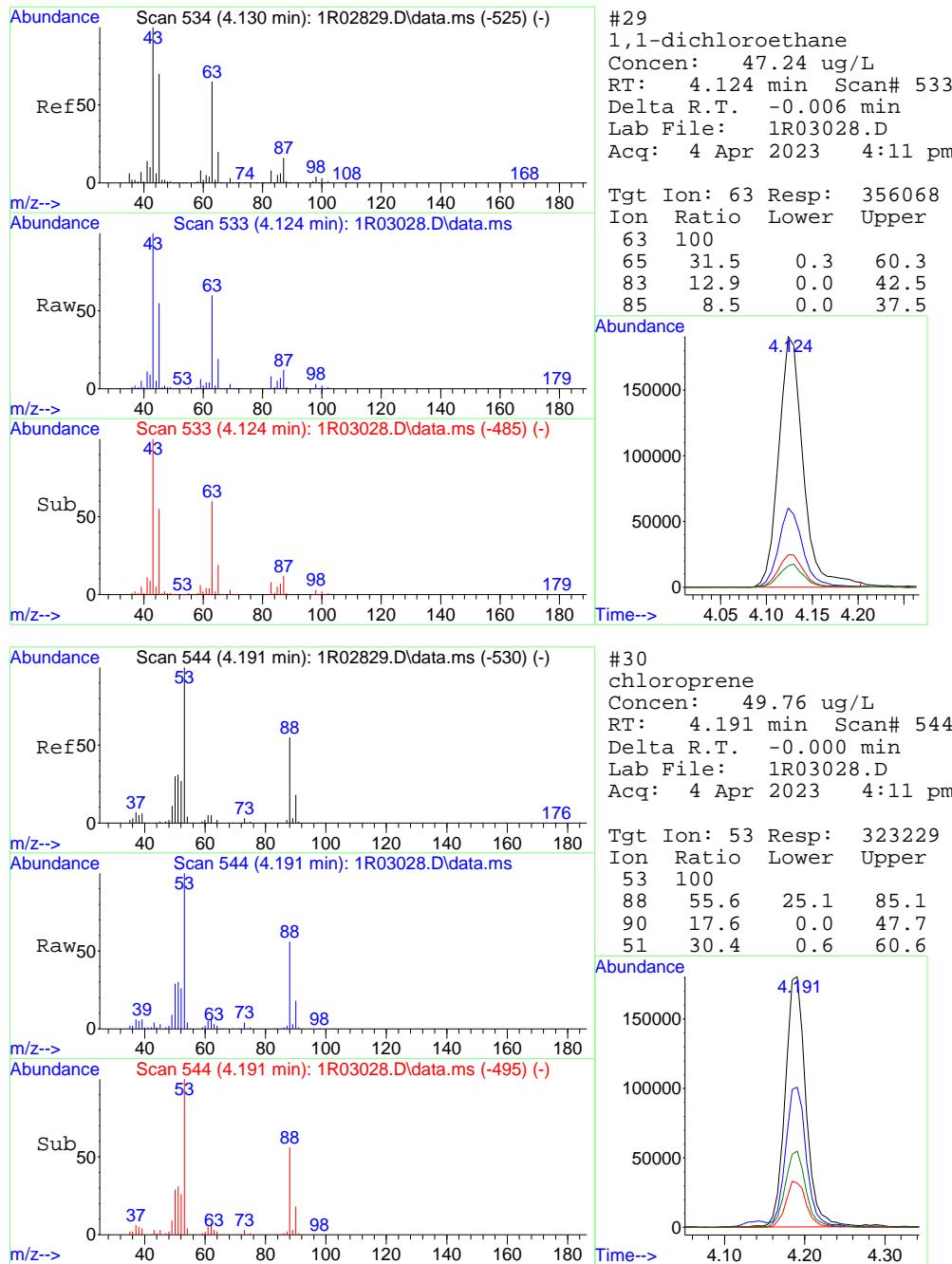


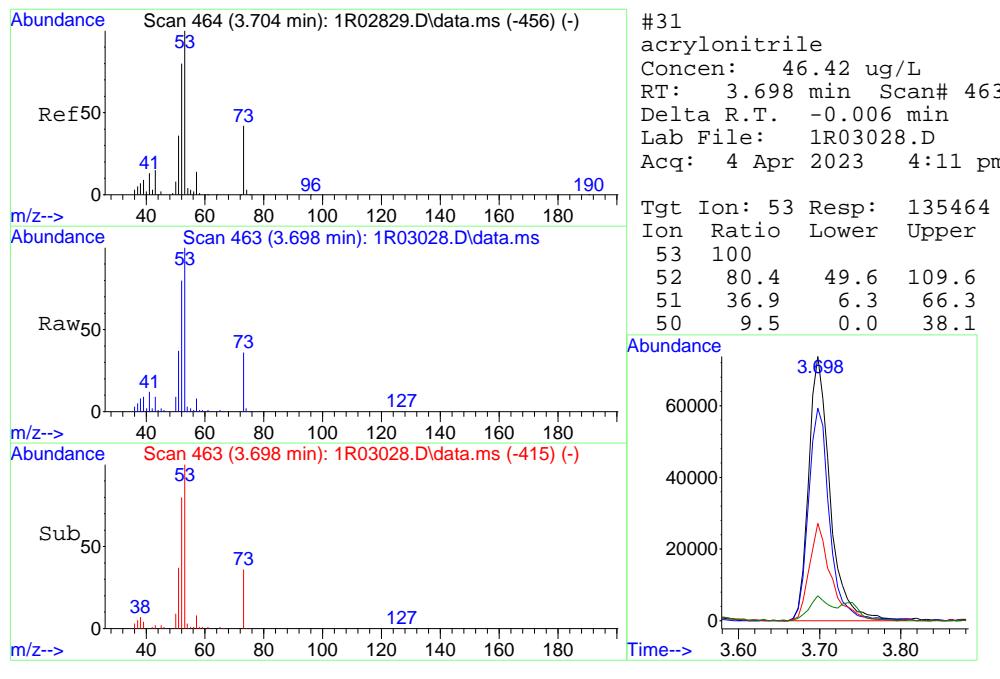






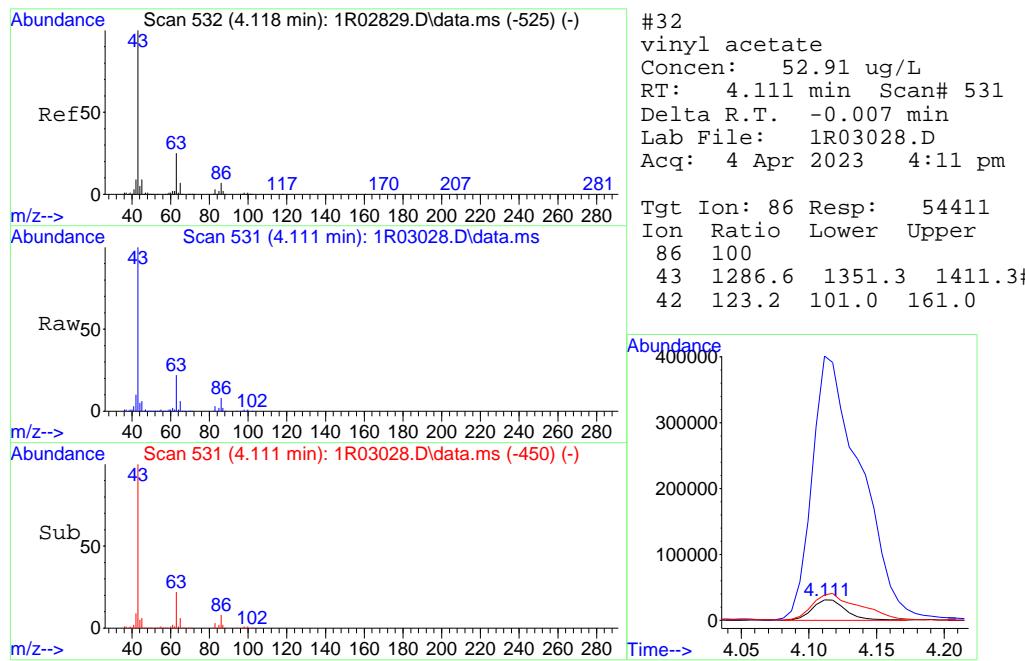






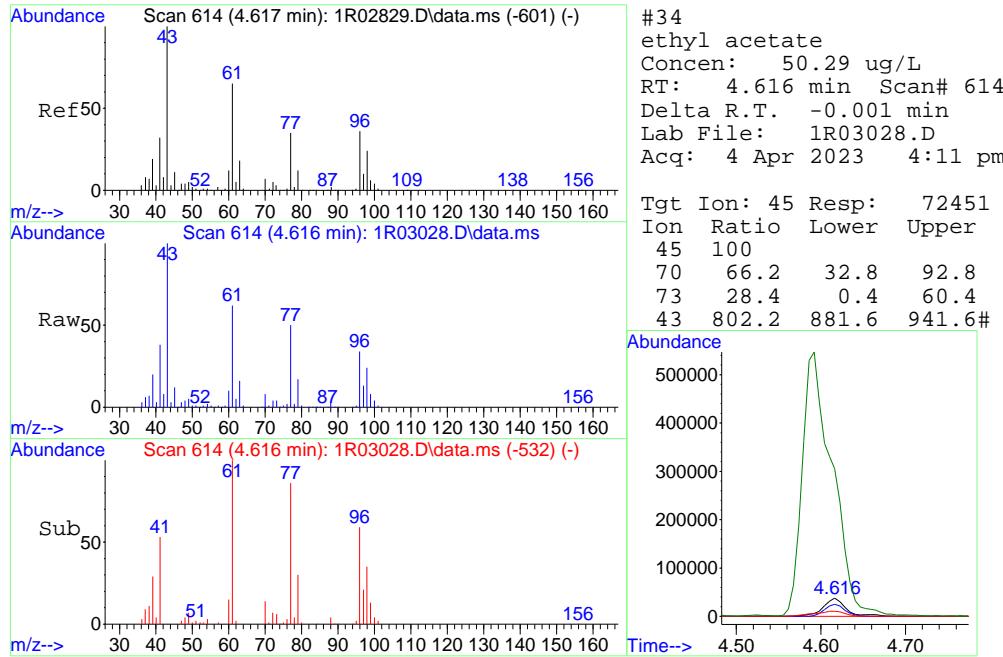
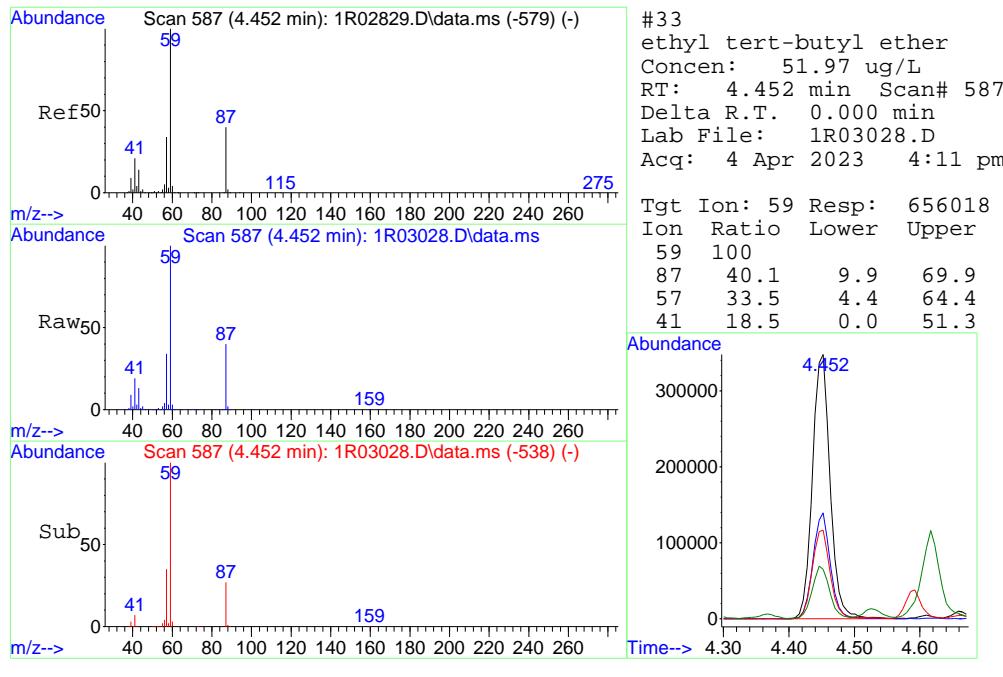
#31  
acrylonitrile  
Concen: 46.42 ug/L  
RT: 3.698 min Scan# 463  
Delta R.T. -0.006 min  
Lab File: 1R03028.D  
Acq: 4 Apr 2023 4:11 pm

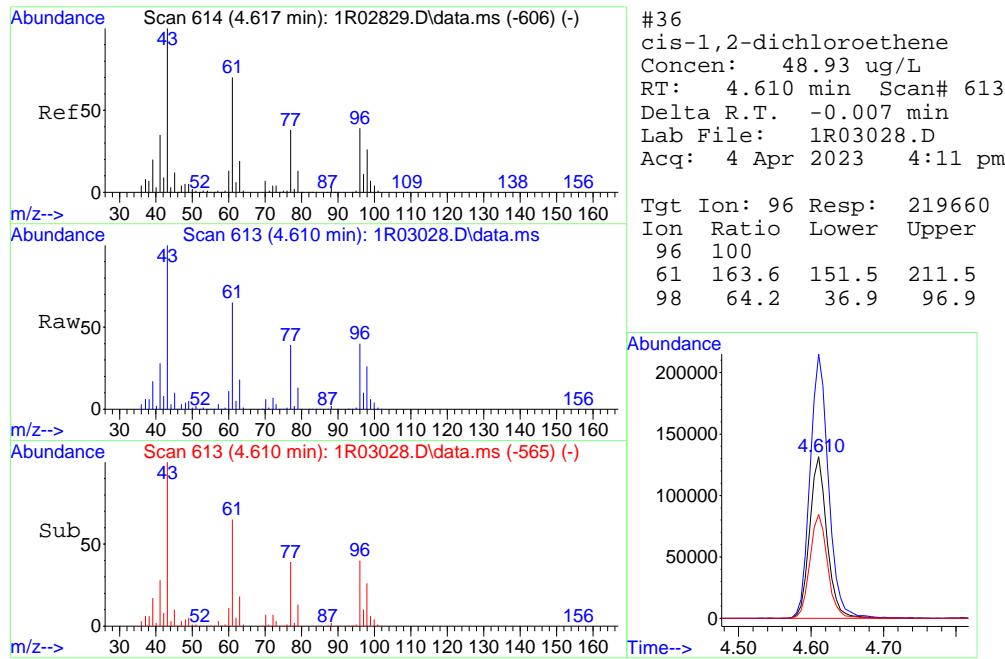
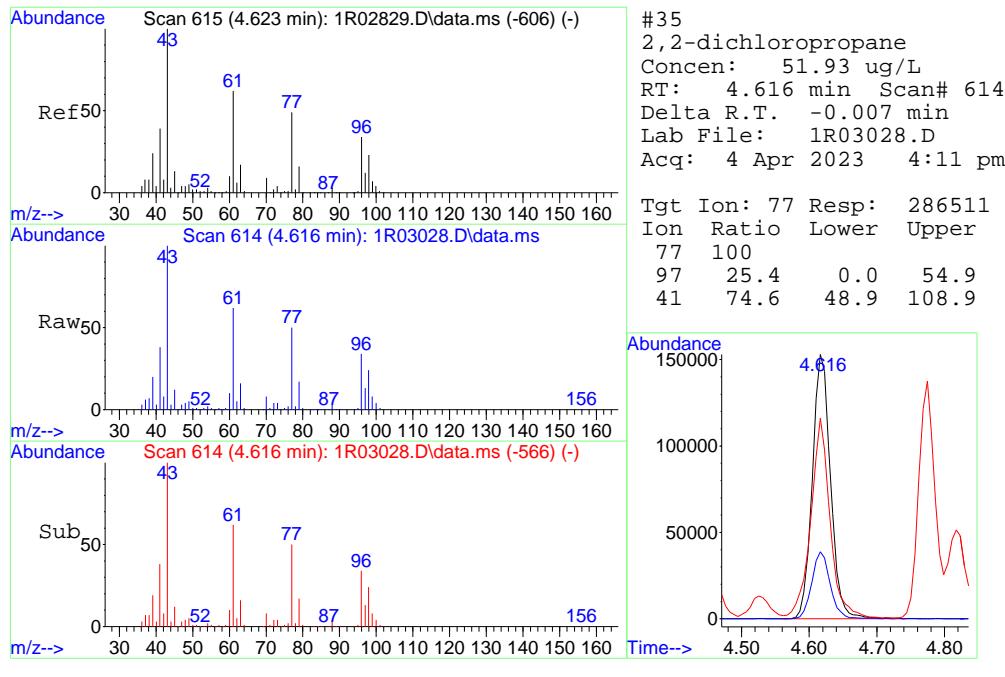
Tgt Ion: 53 Resp: 135464  
Ion Ratio Lower Upper  
53 100  
52 80.4 49.6 109.6  
51 36.9 6.3 66.3  
50 9.5 0.0 38.1

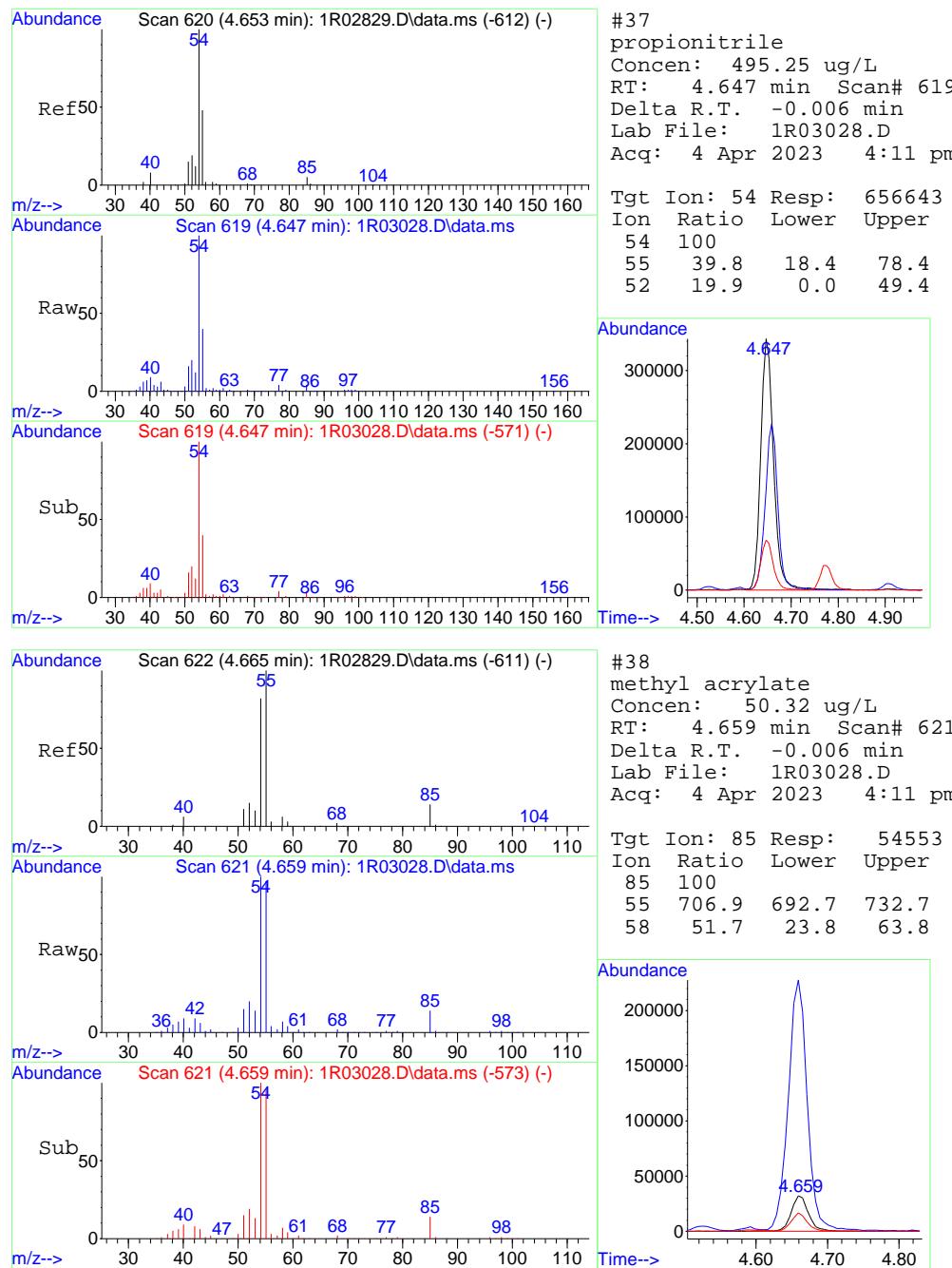


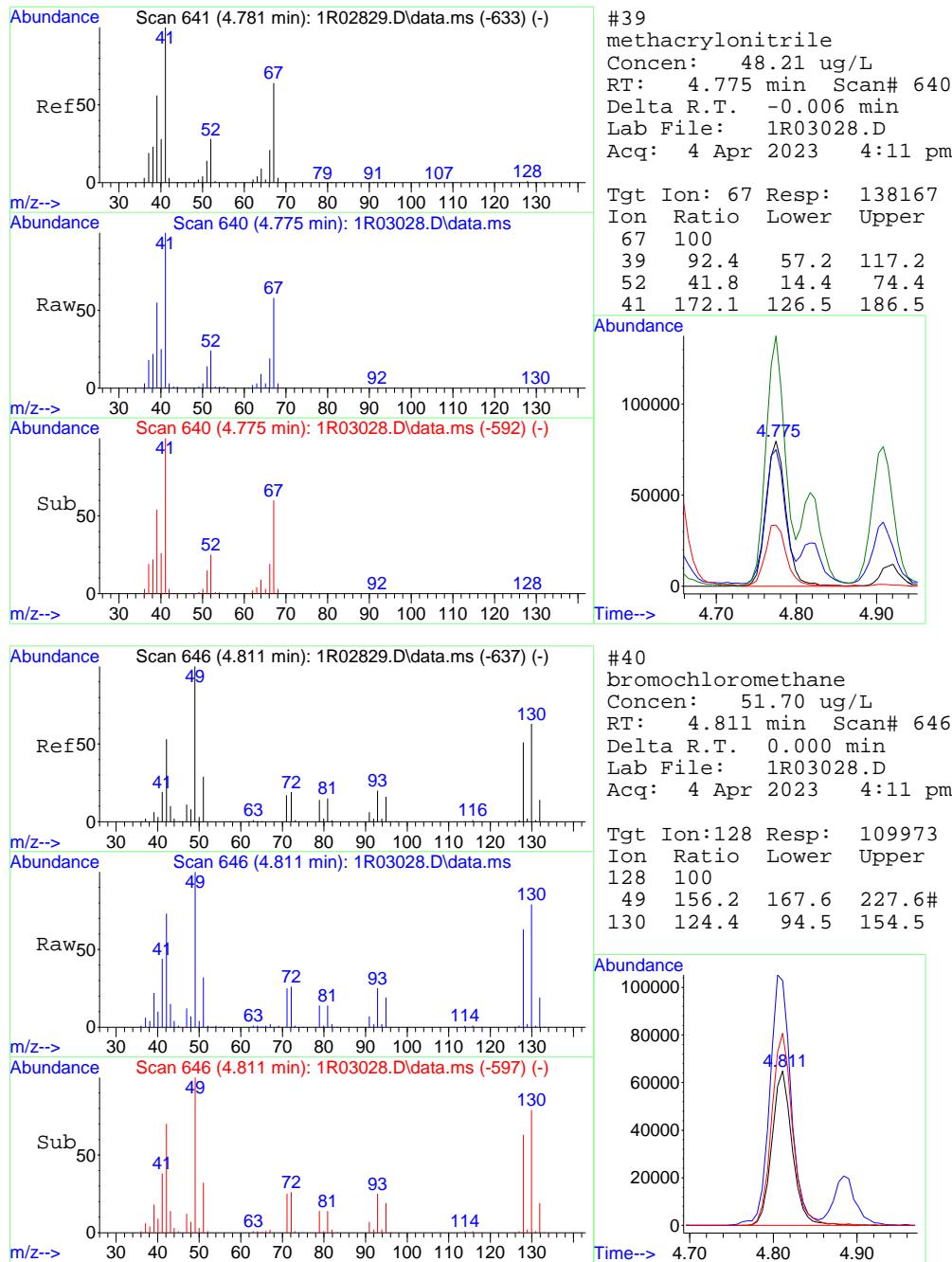
#32  
vinyl acetate  
Concen: 52.91 ug/L  
RT: 4.111 min Scan# 531  
Delta R.T. -0.007 min  
Lab File: 1R03028.D  
Acq: 4 Apr 2023 4:11 pm

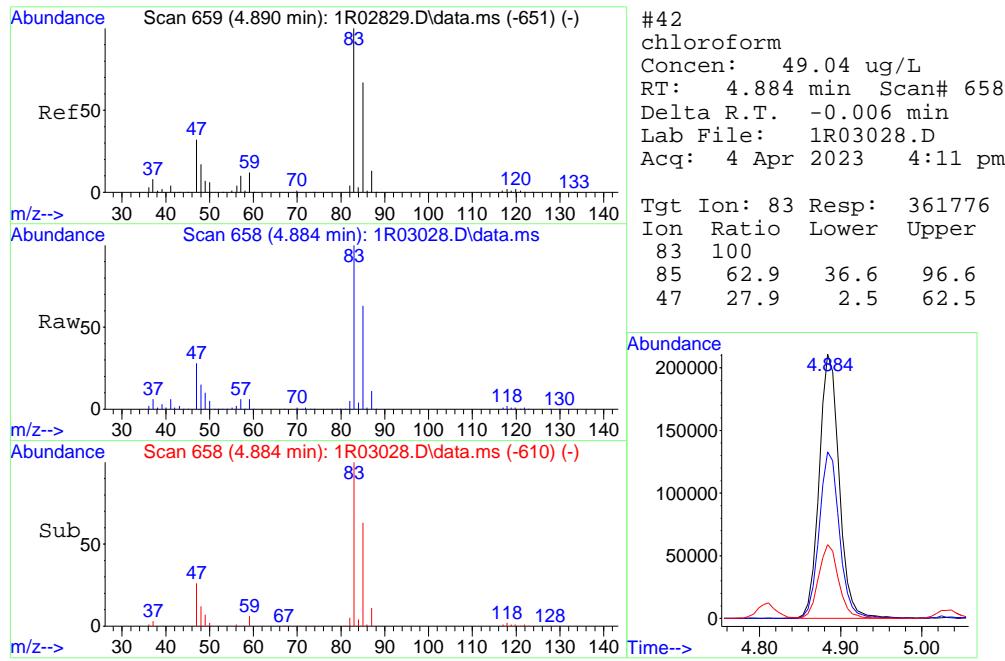
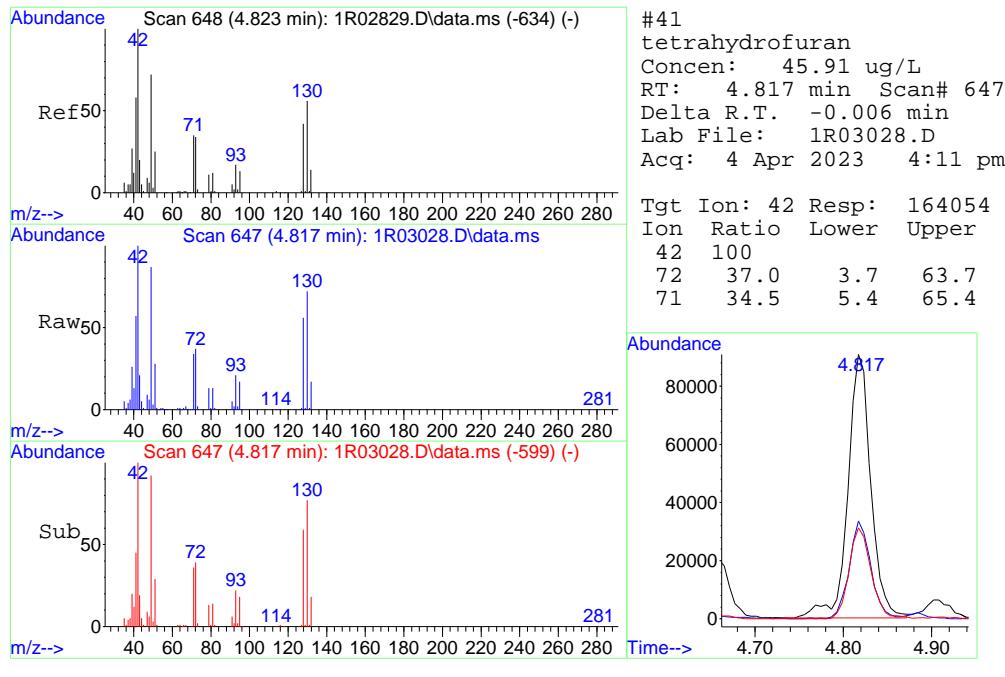
Tgt Ion: 86 Resp: 54411  
Ion Ratio Lower Upper  
86 100  
43 1286.6 1351.3 1411.3#  
42 123.2 101.0 161.0



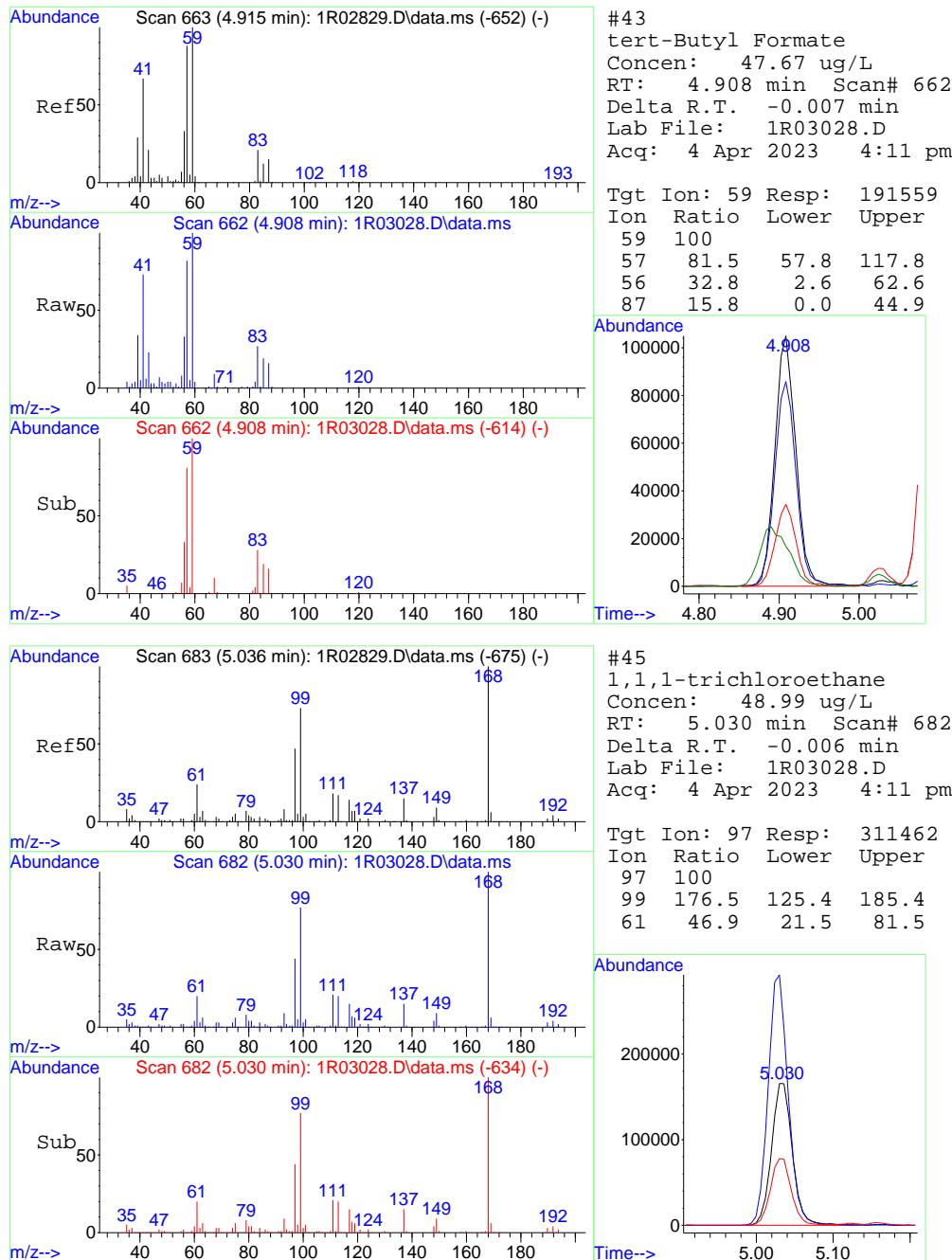


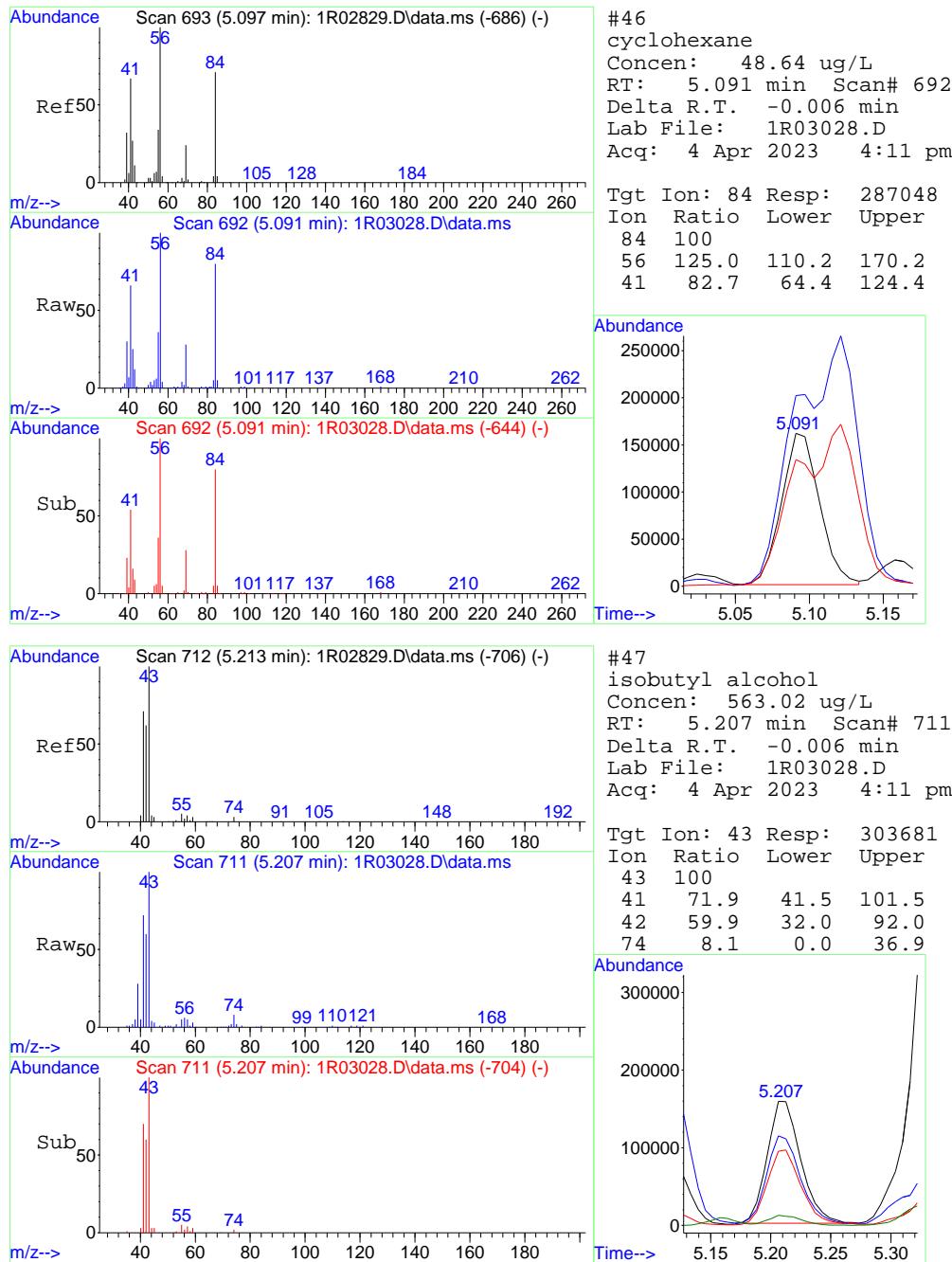


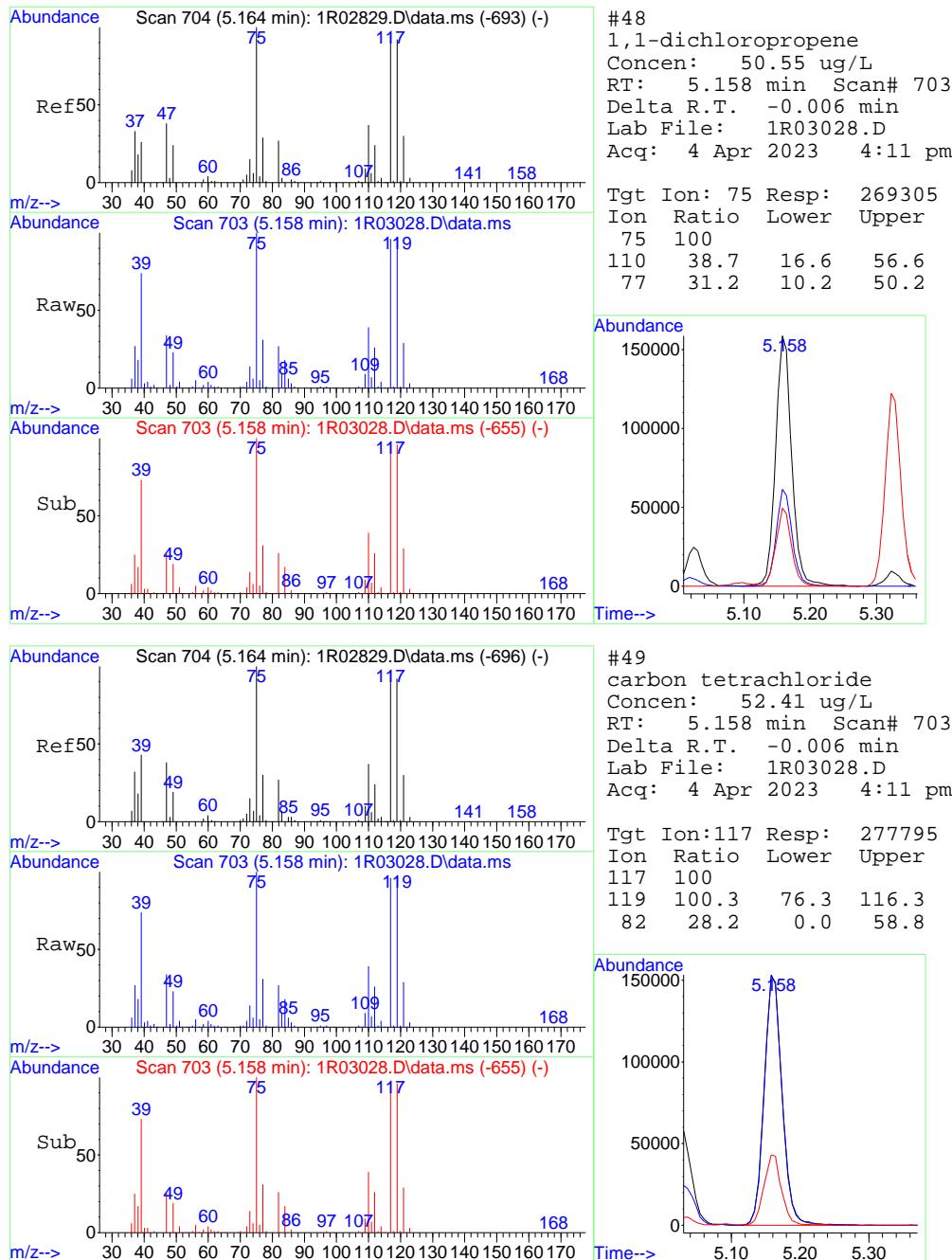


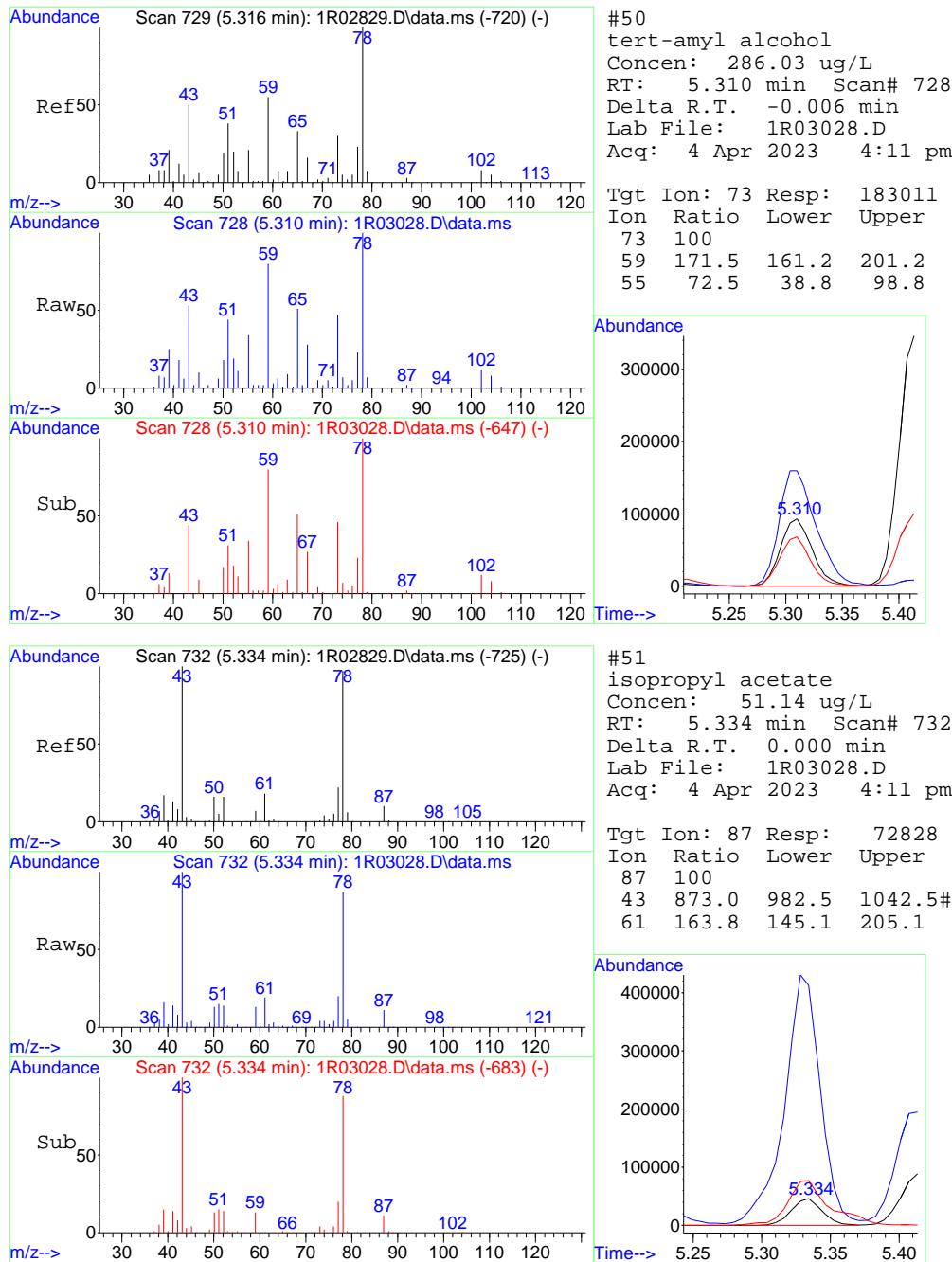


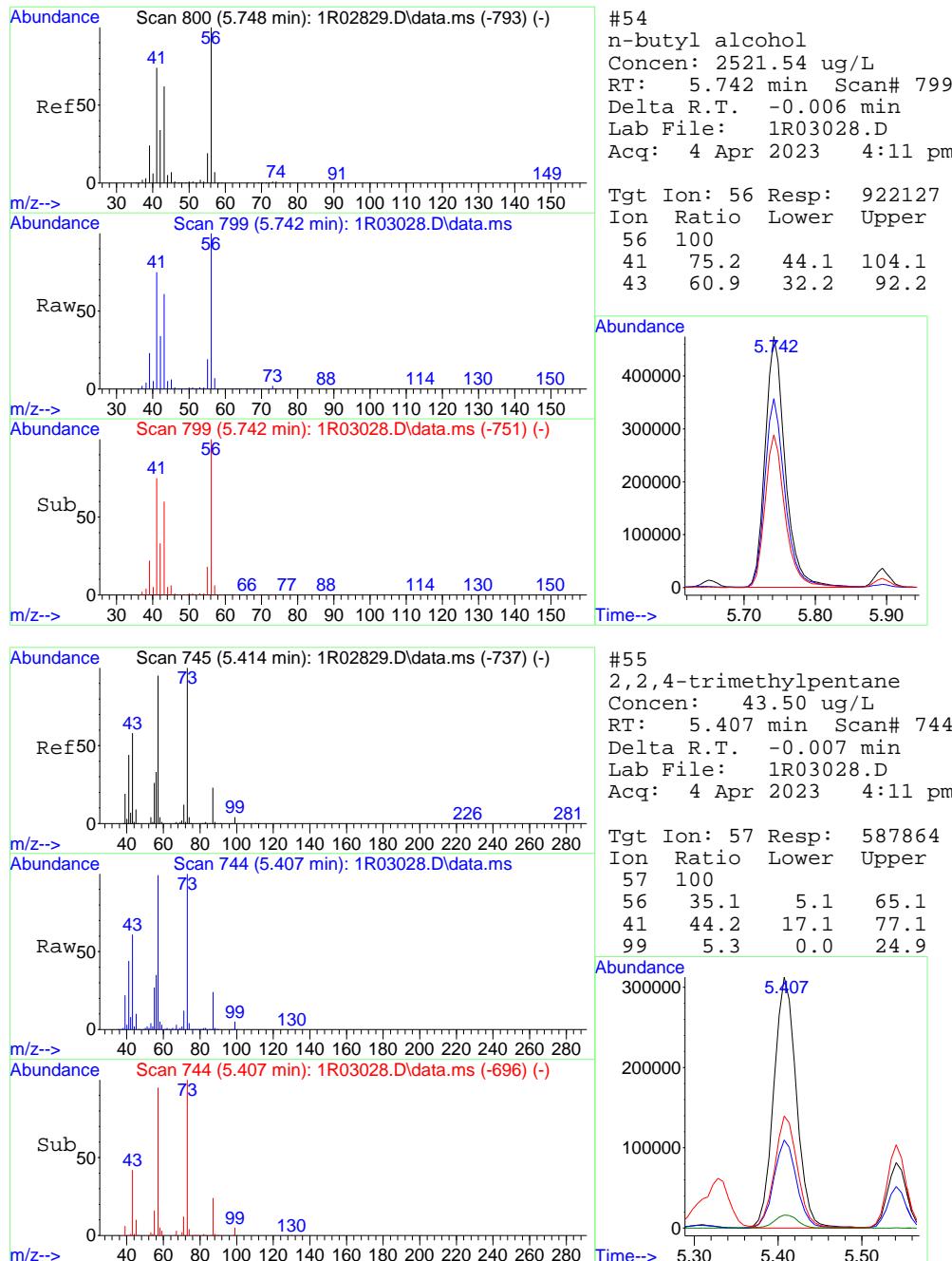
7.4.1  
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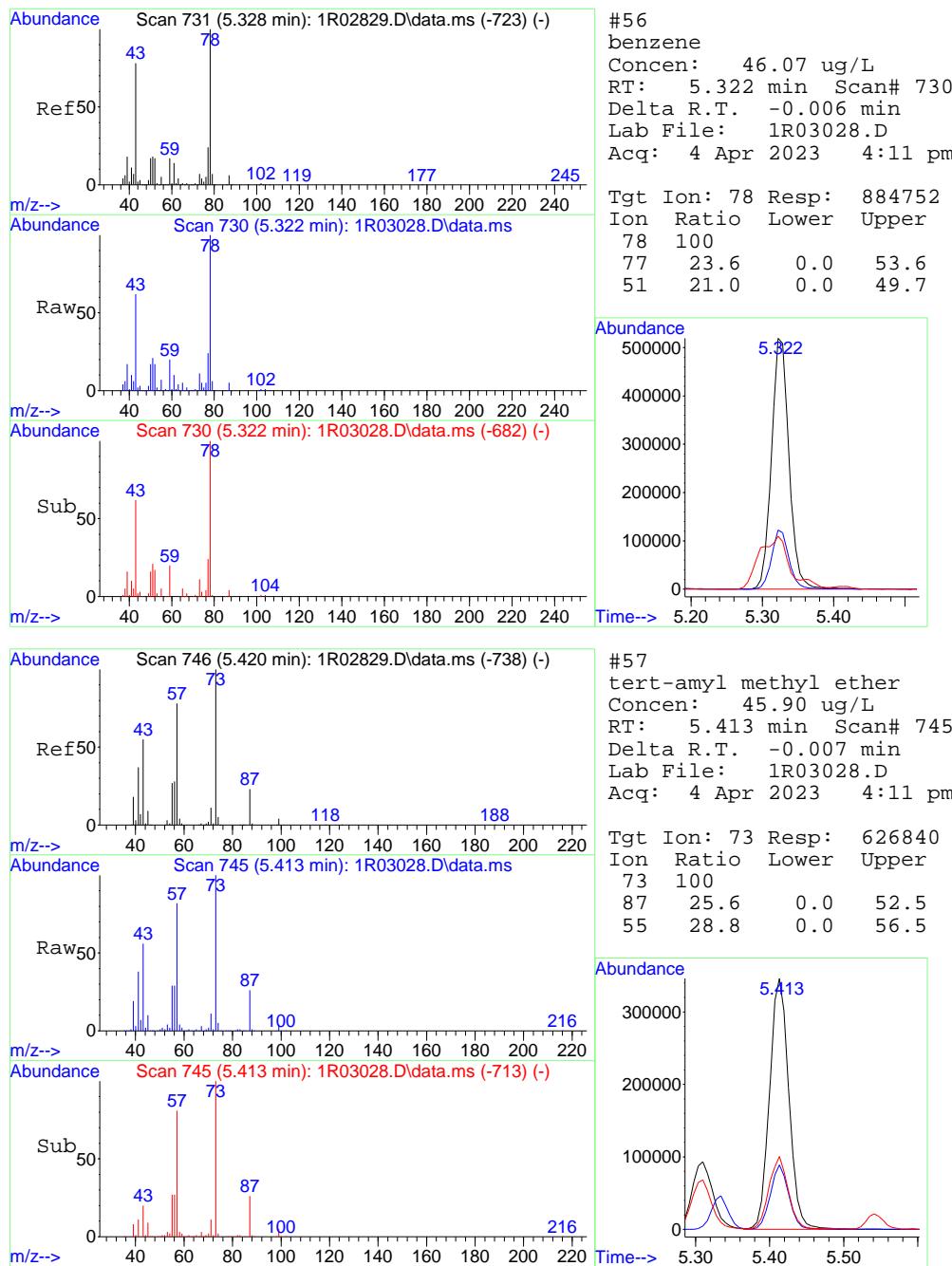


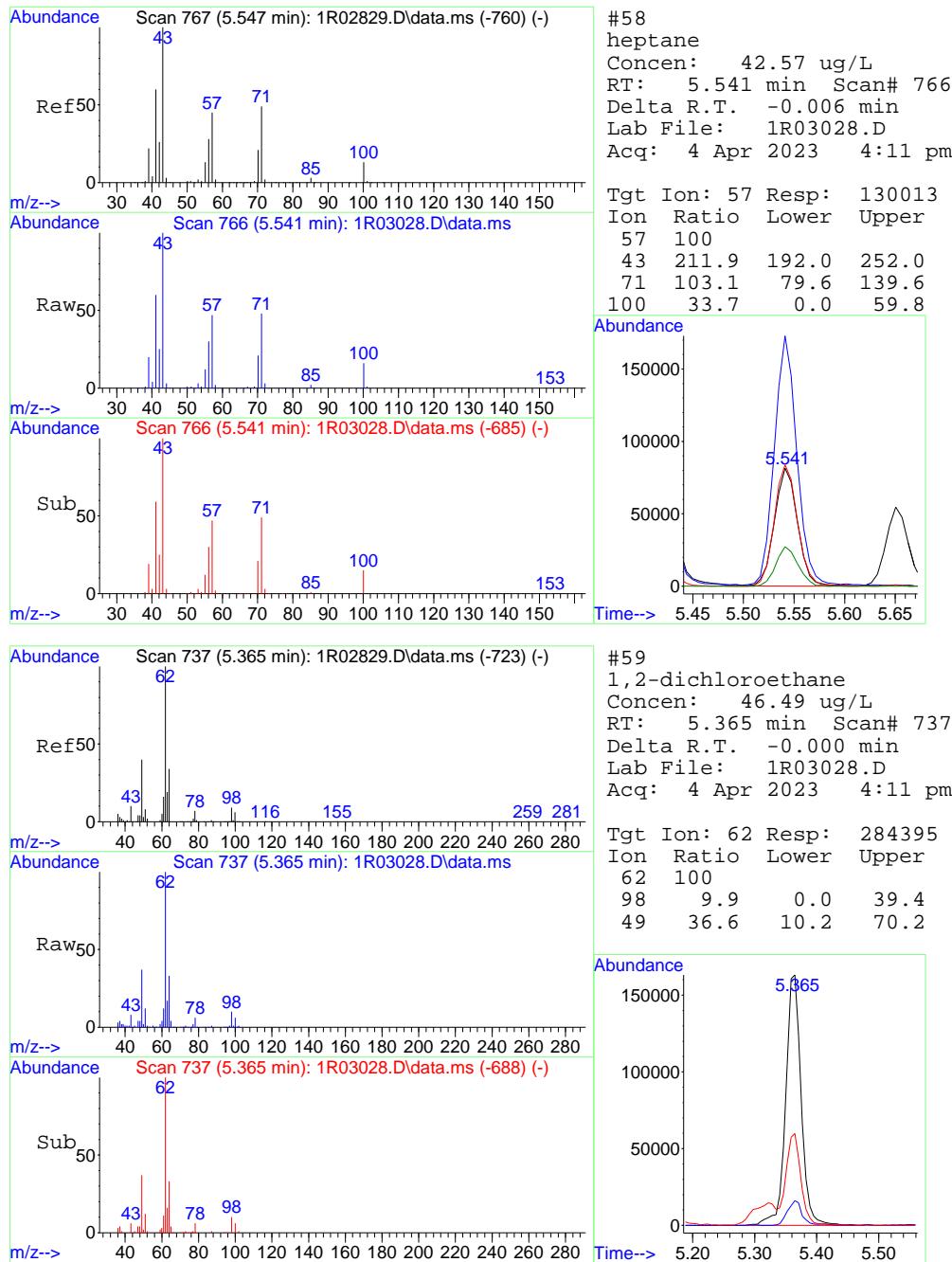


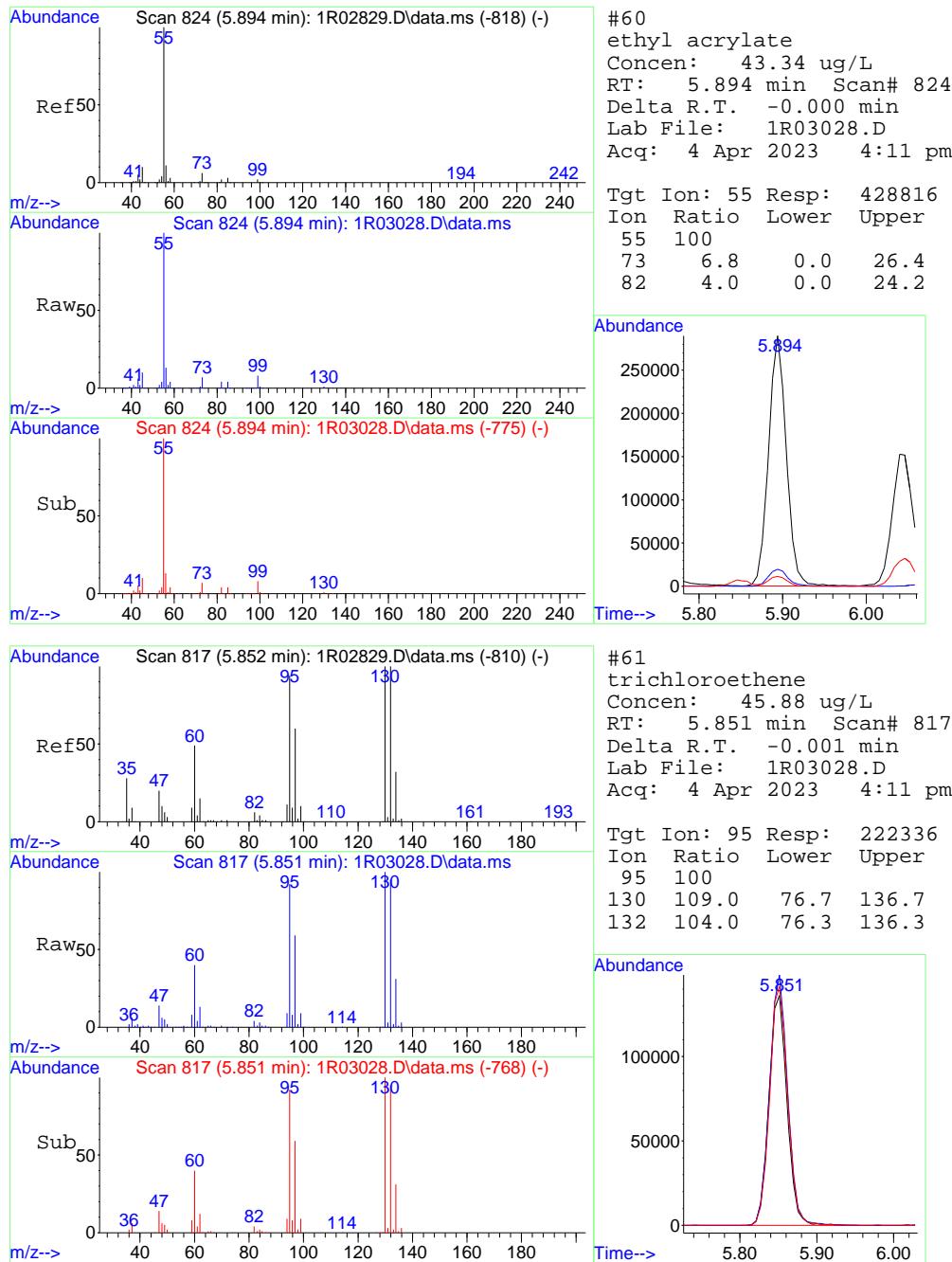


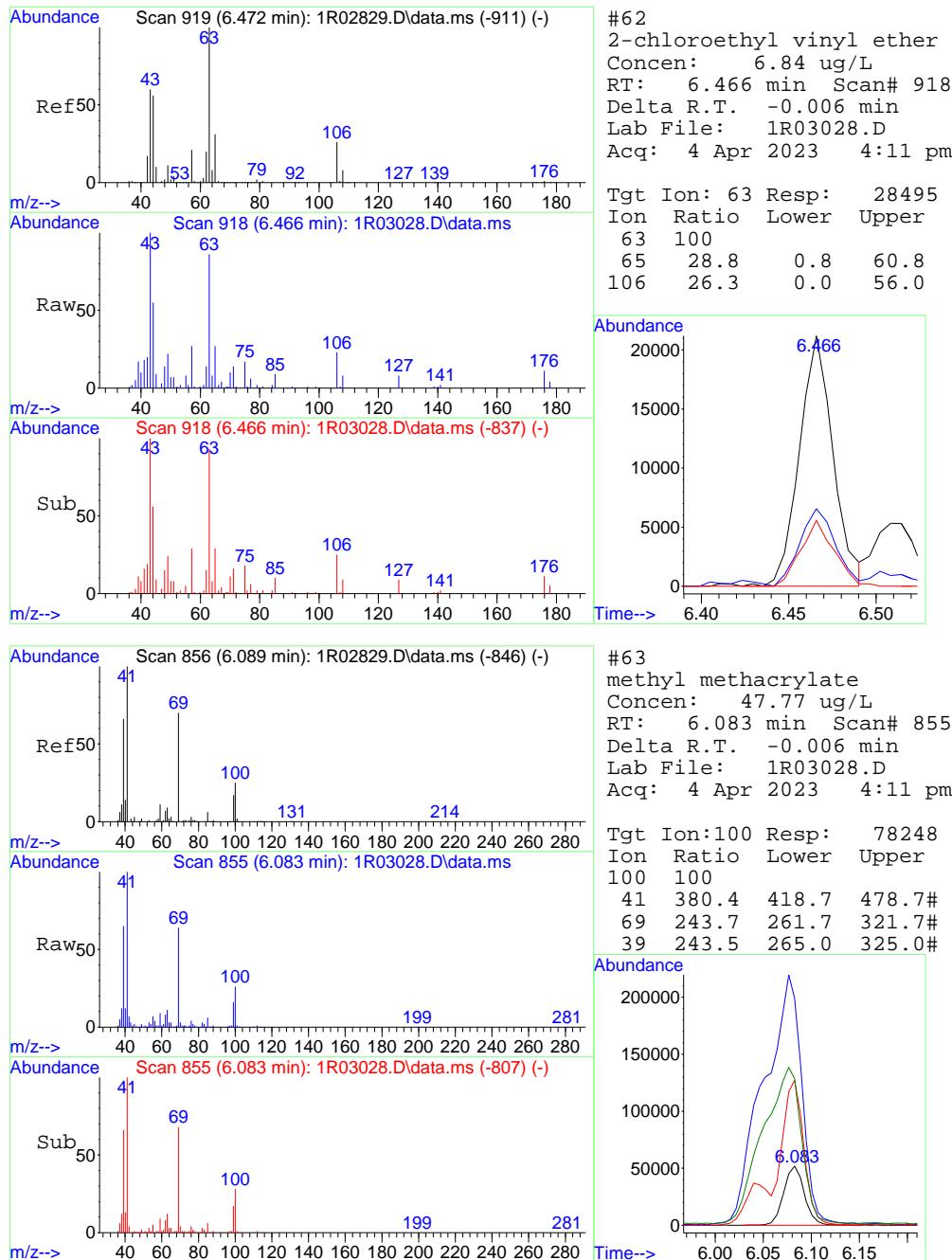


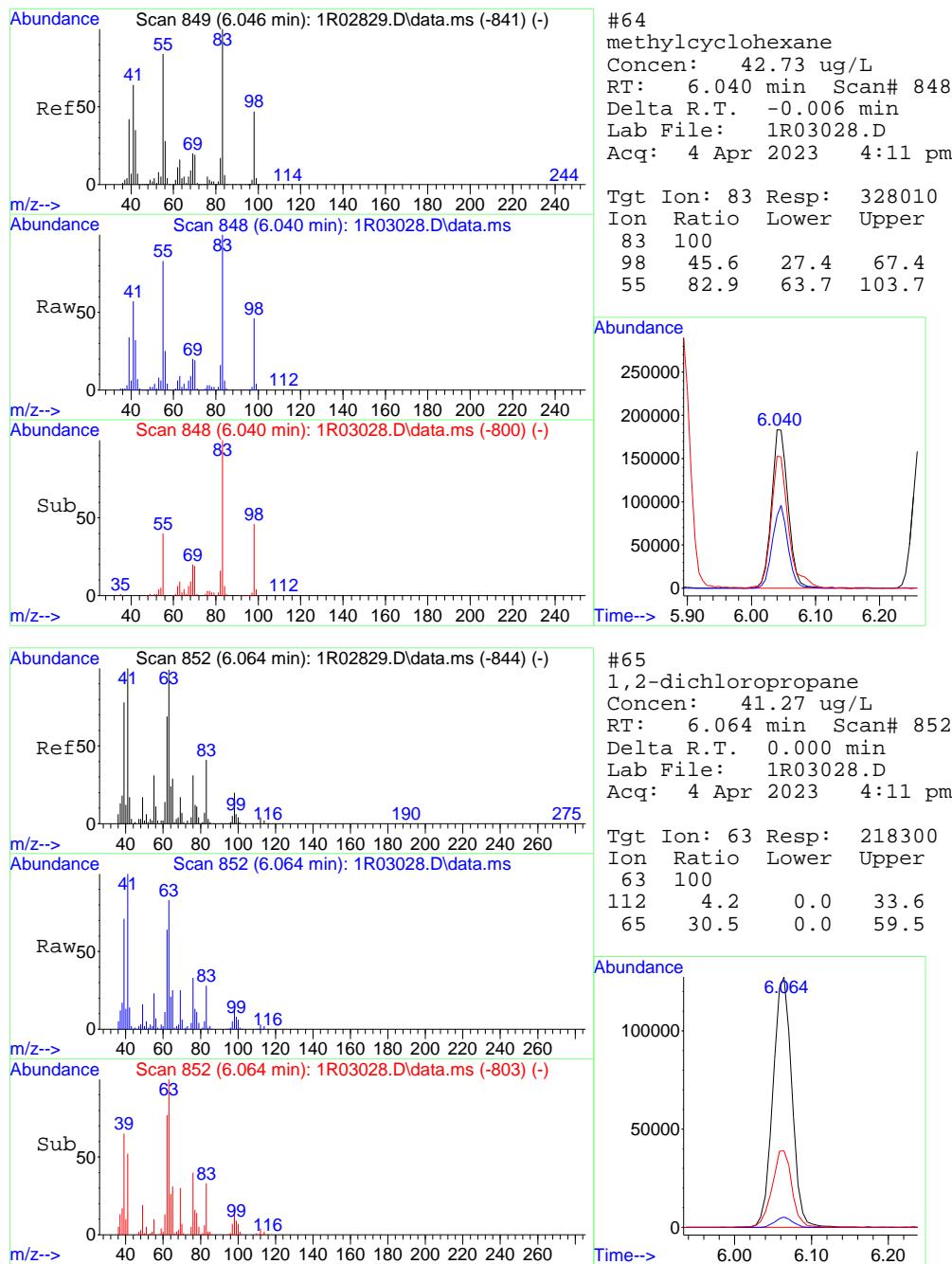


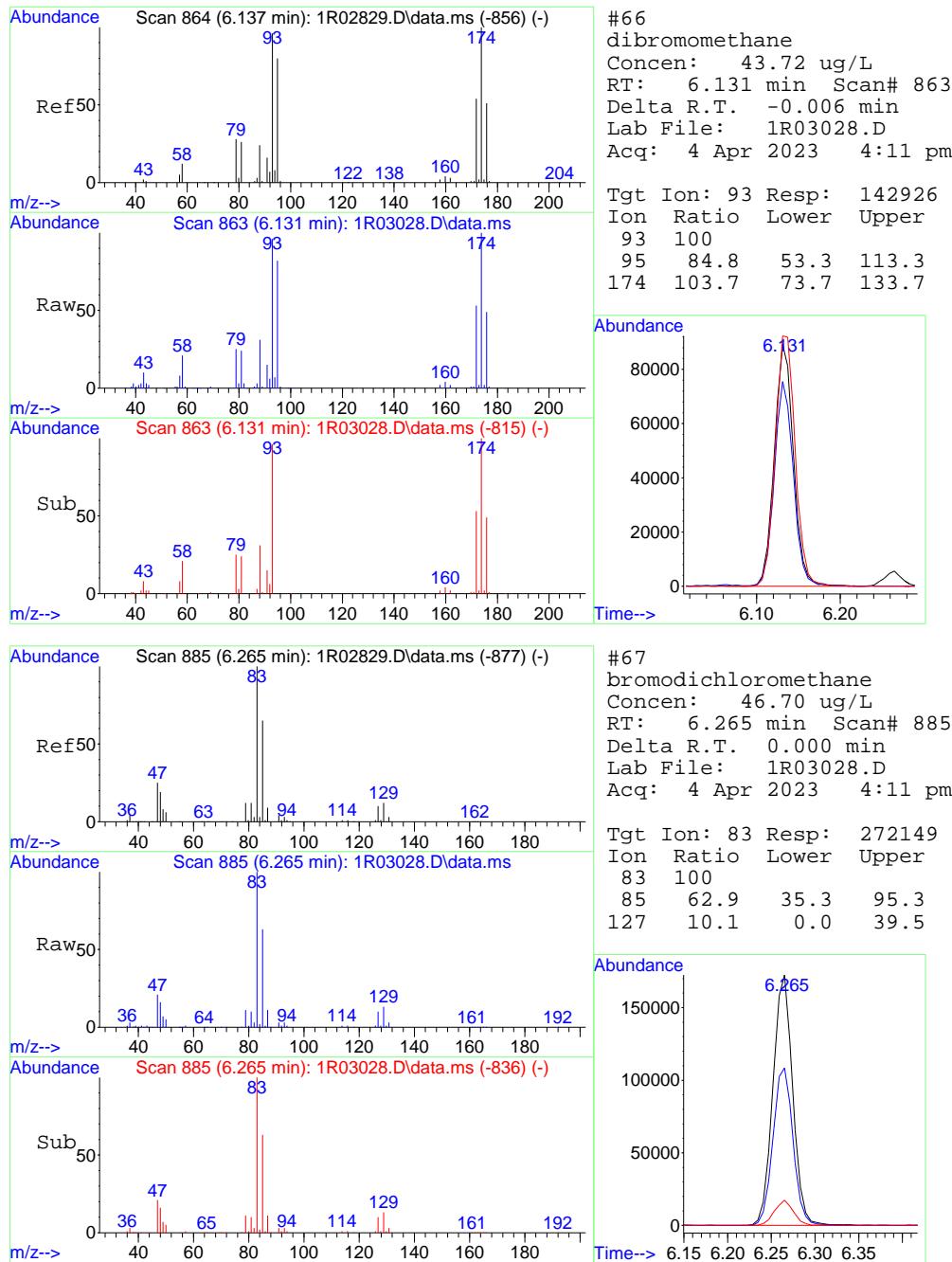


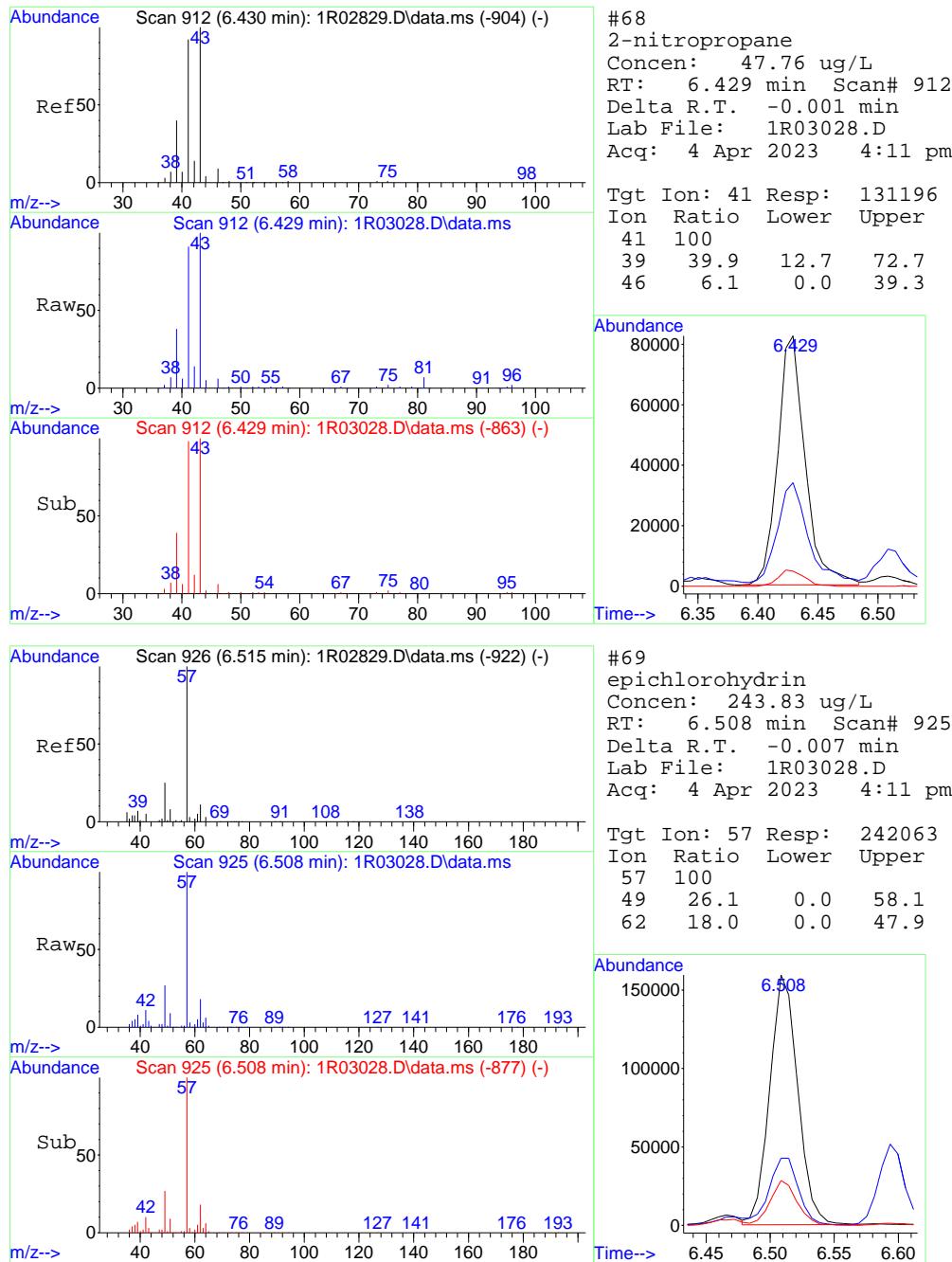


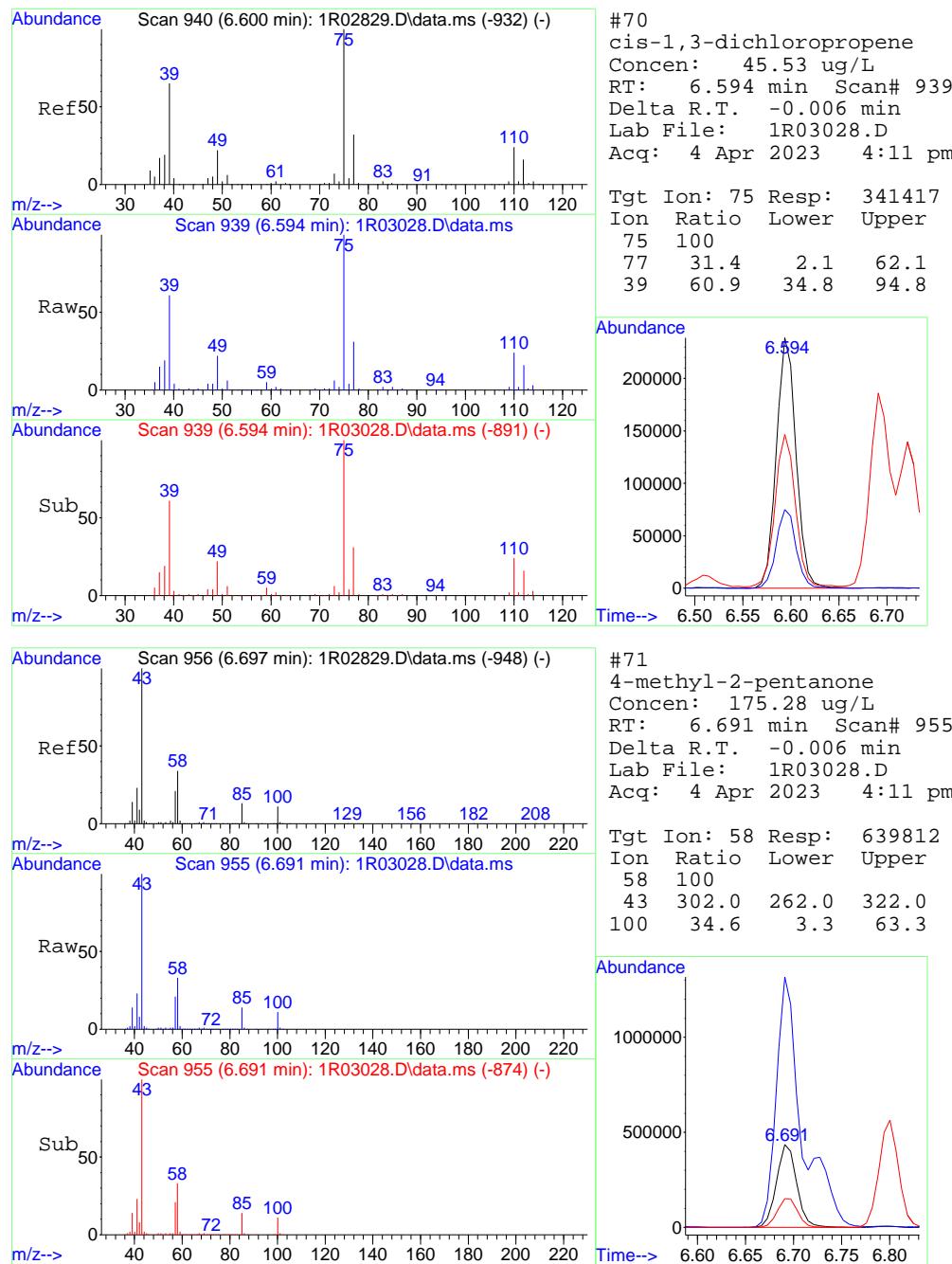


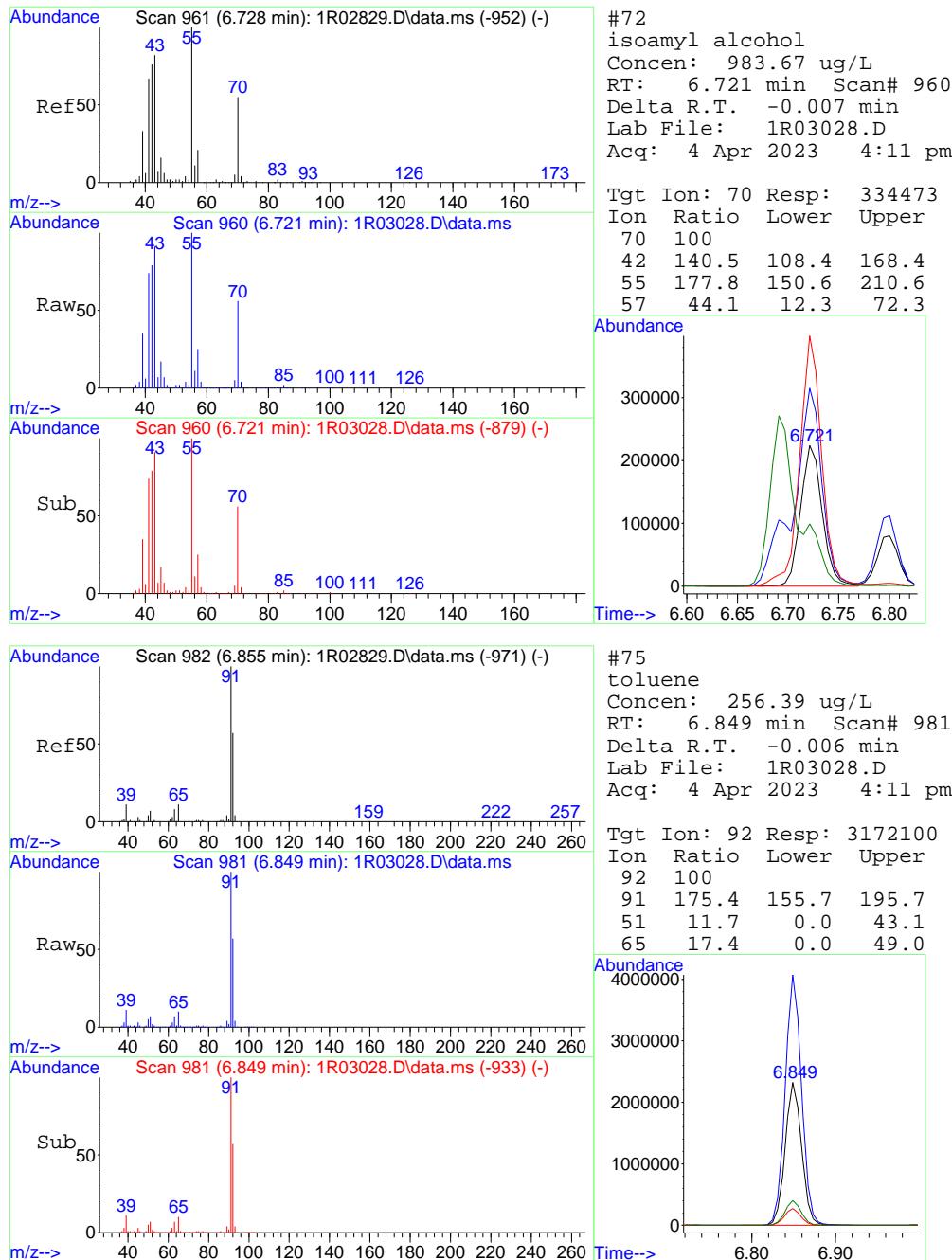


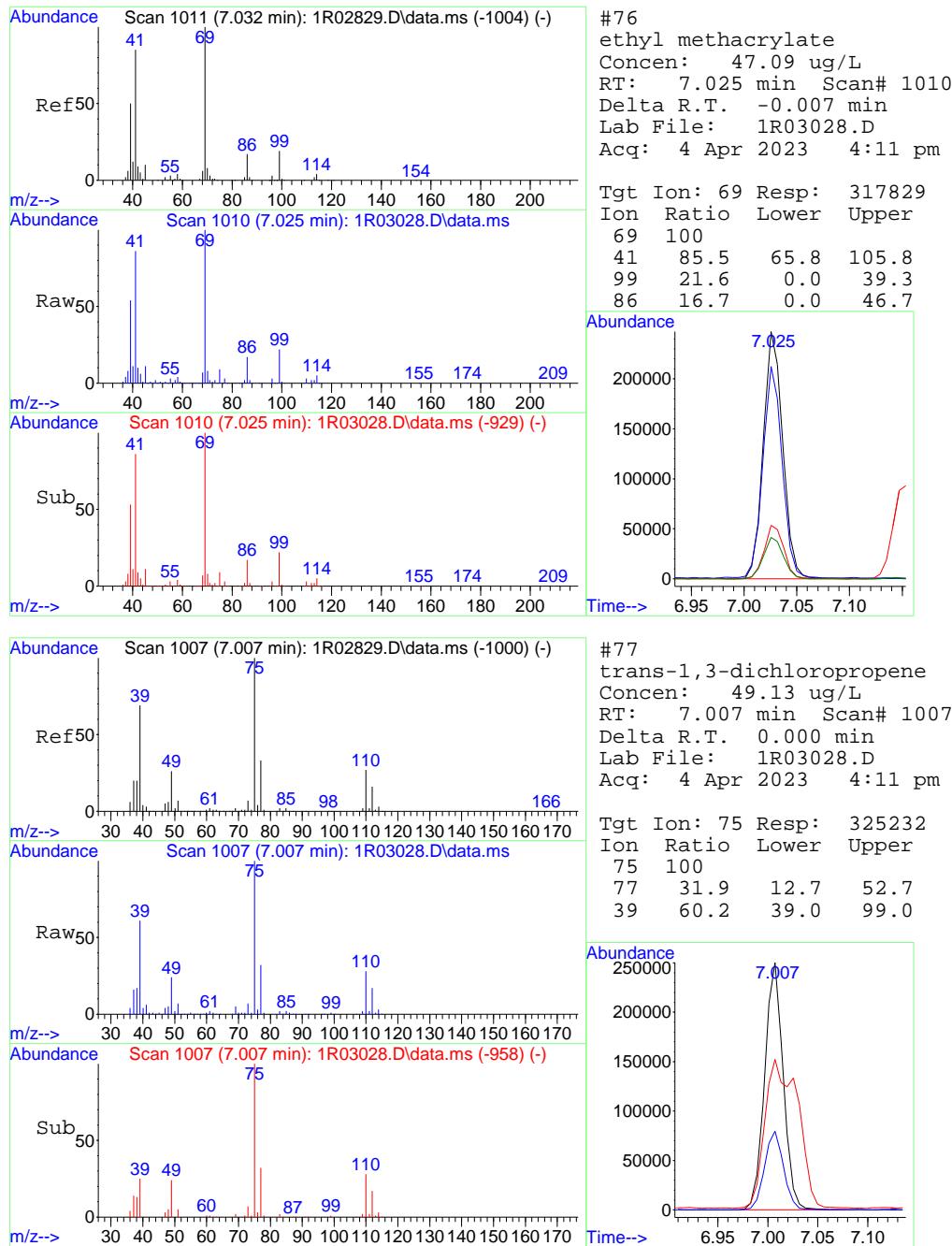


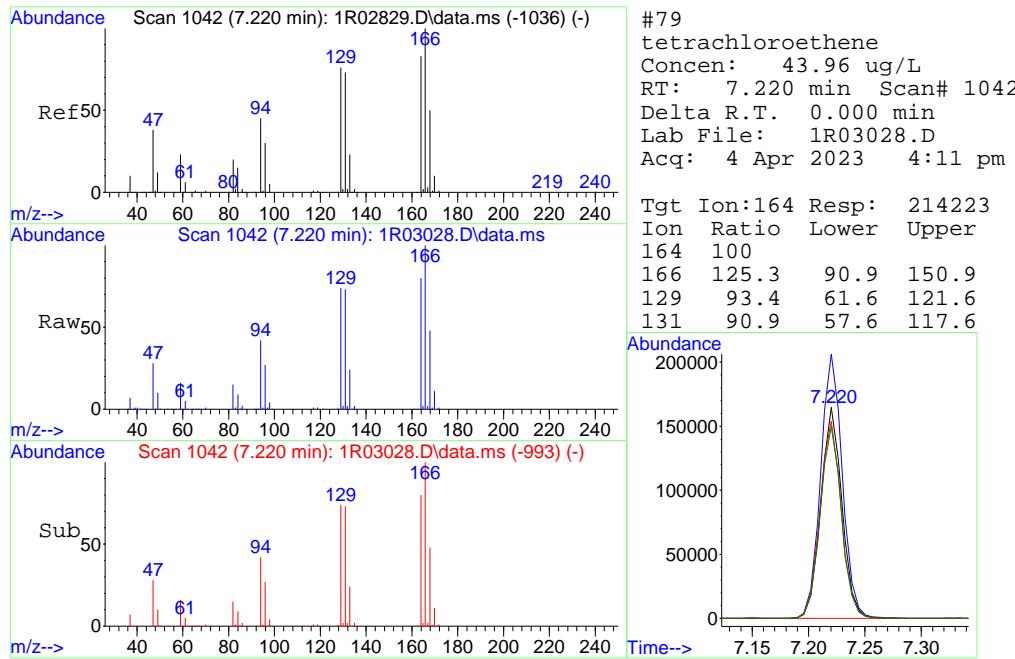
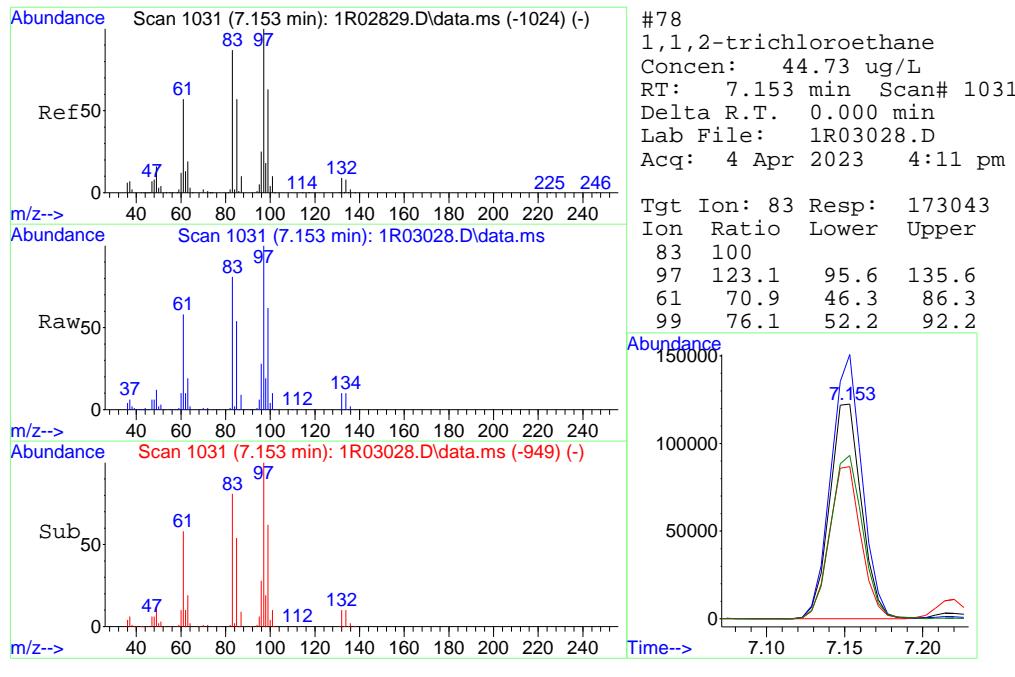


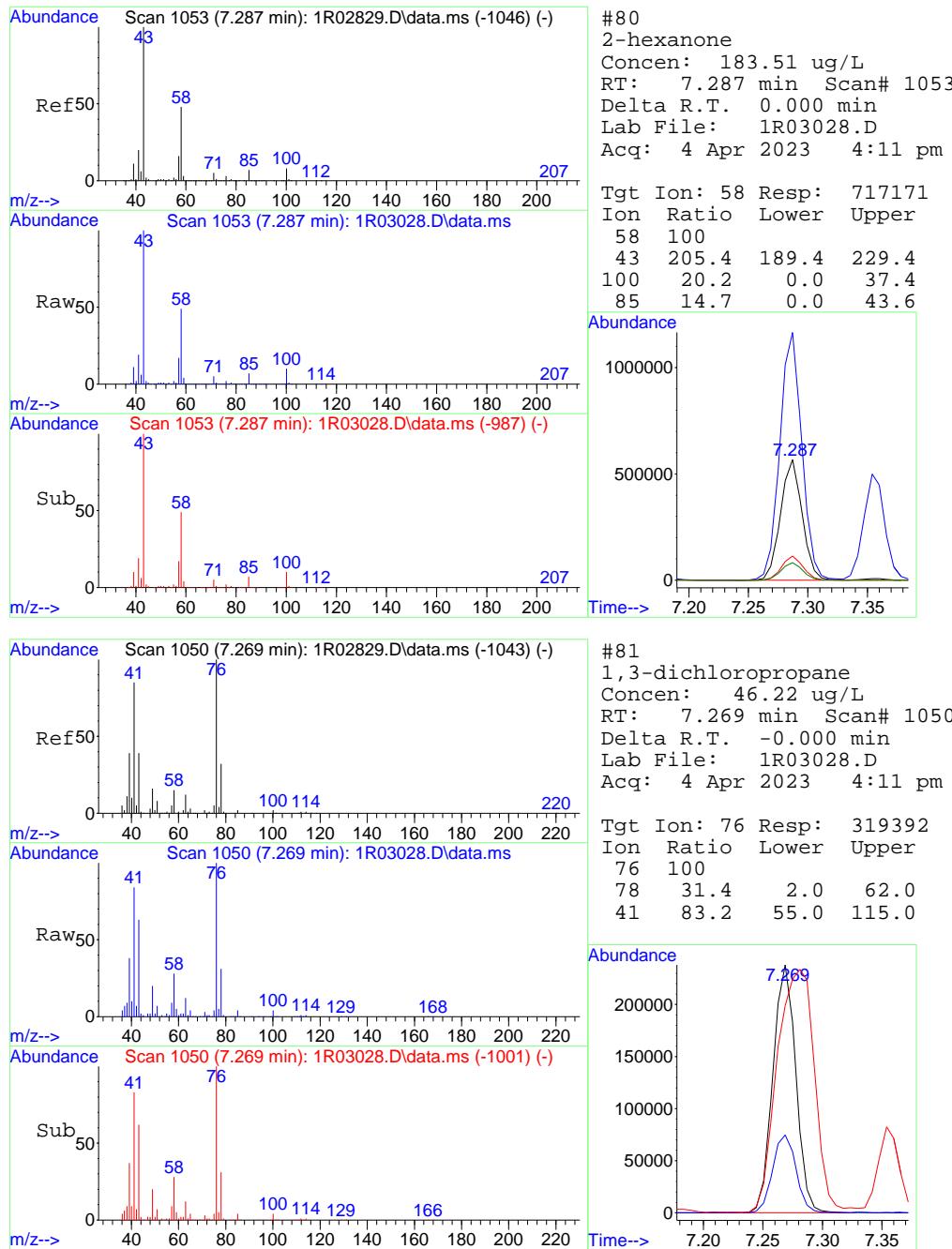


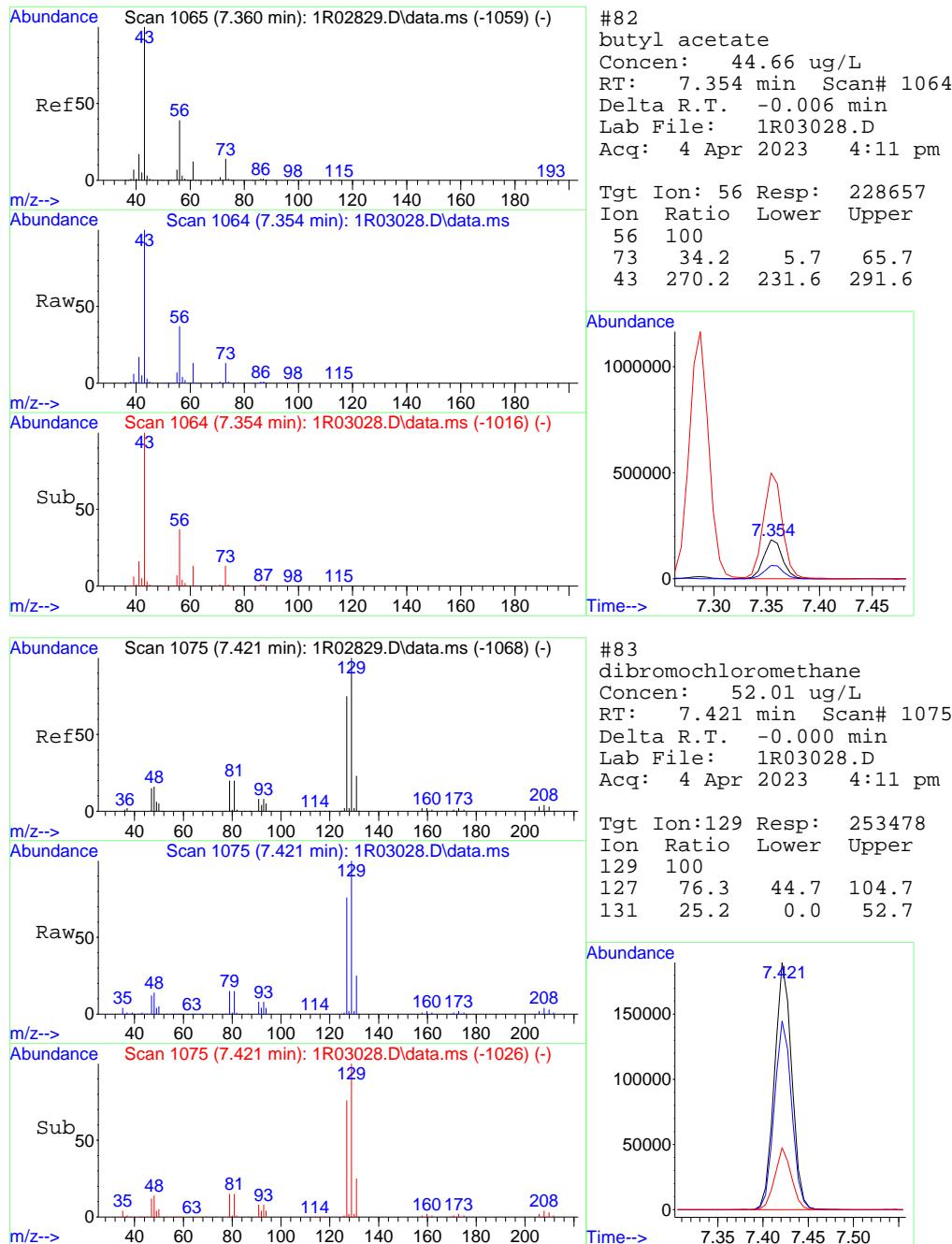


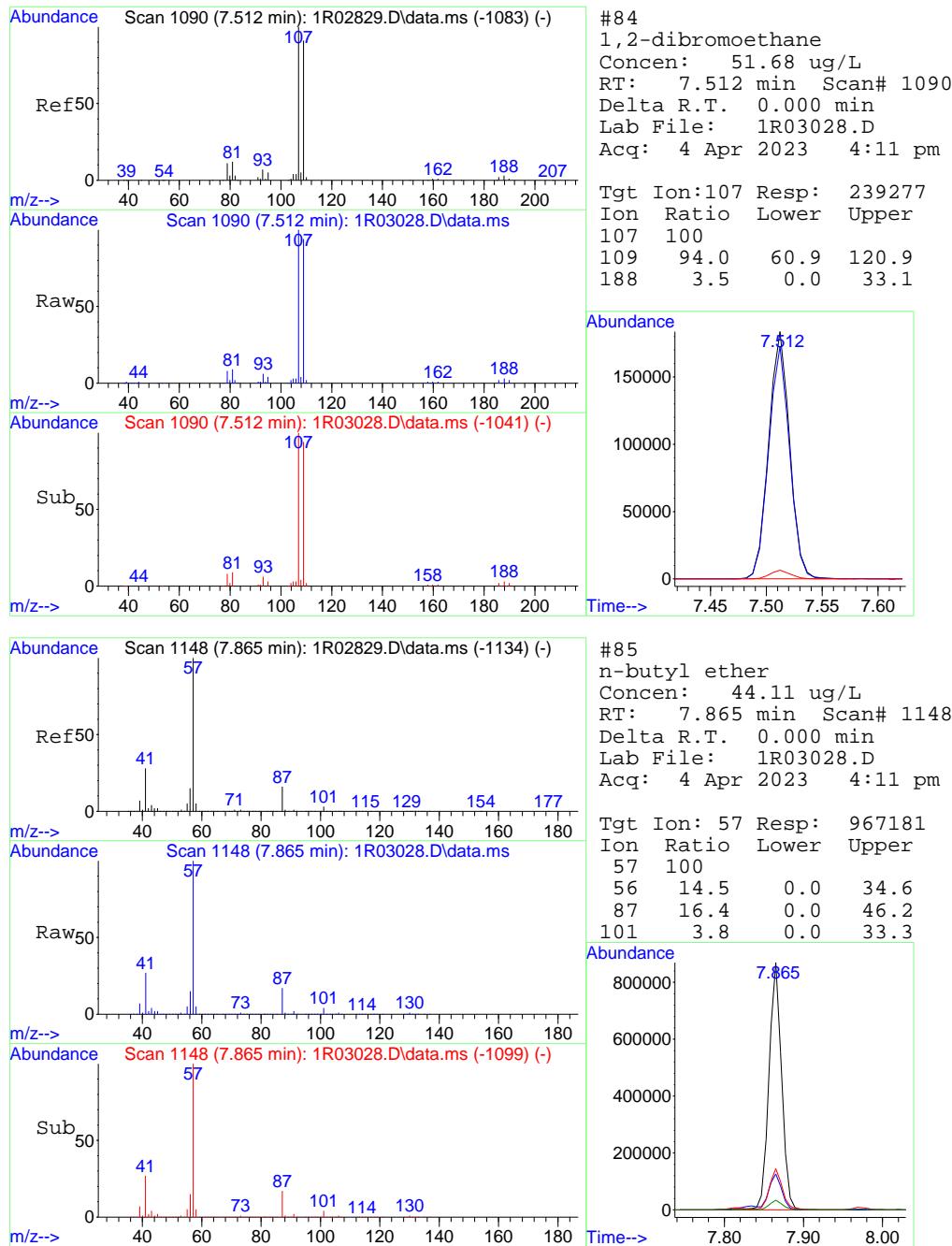


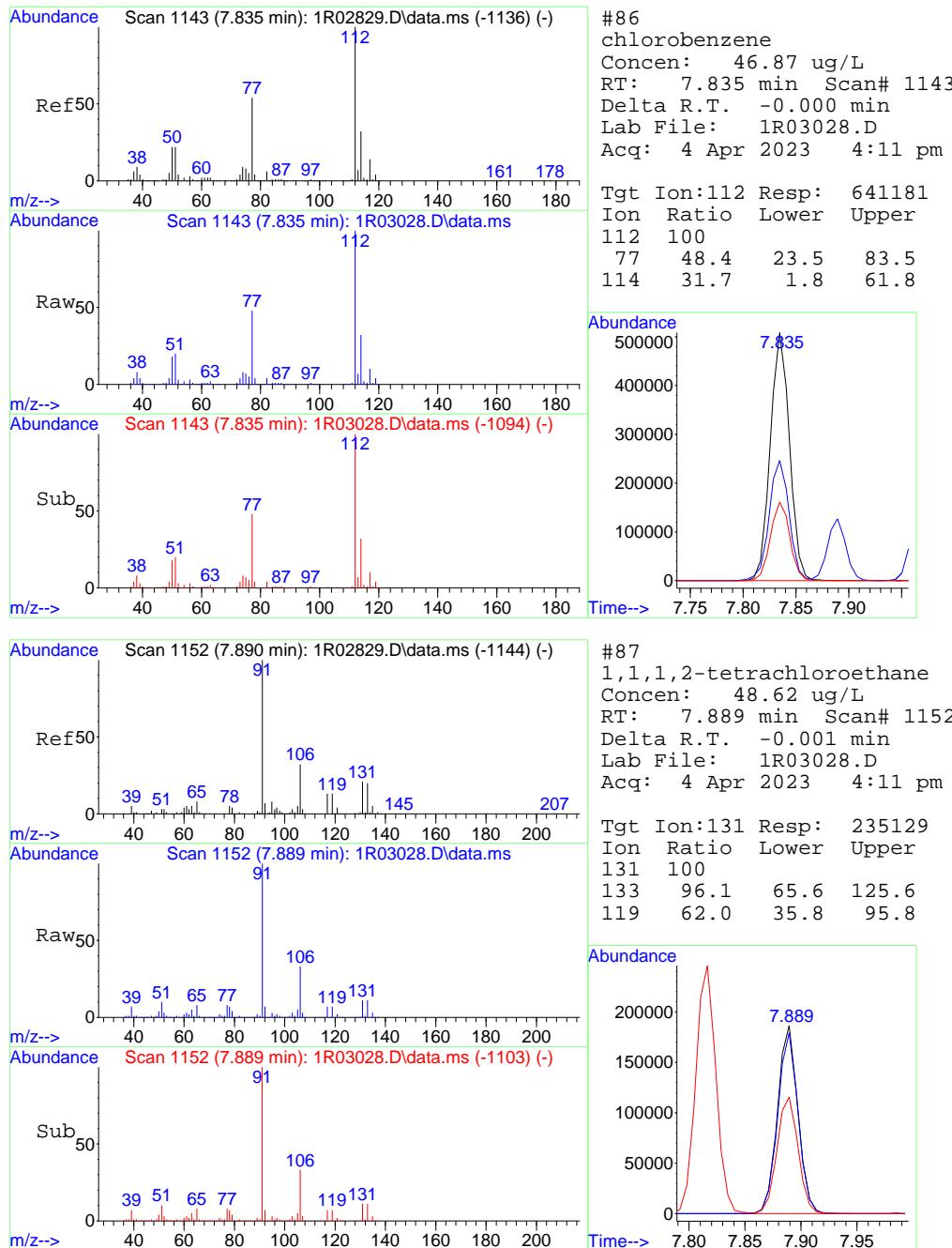


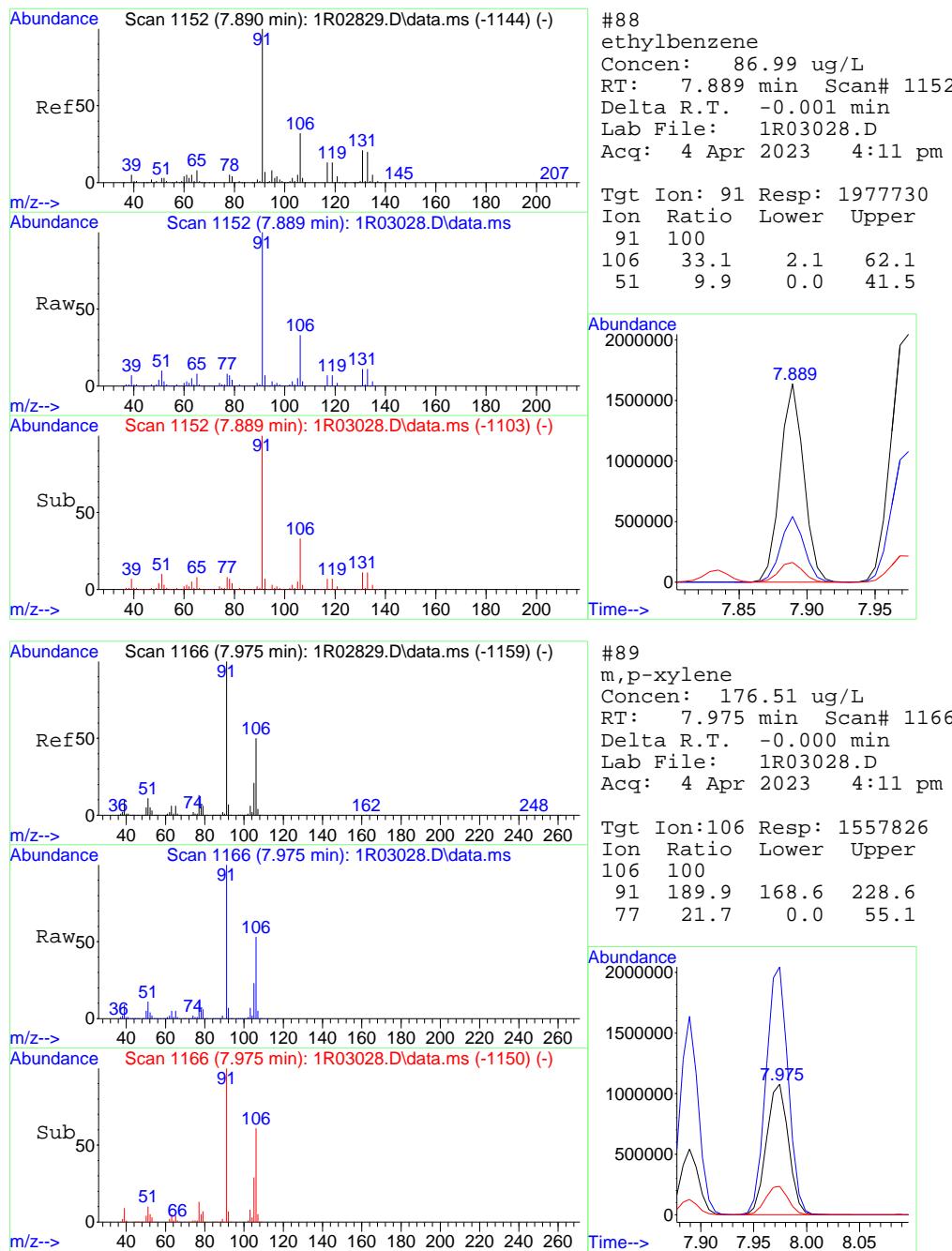


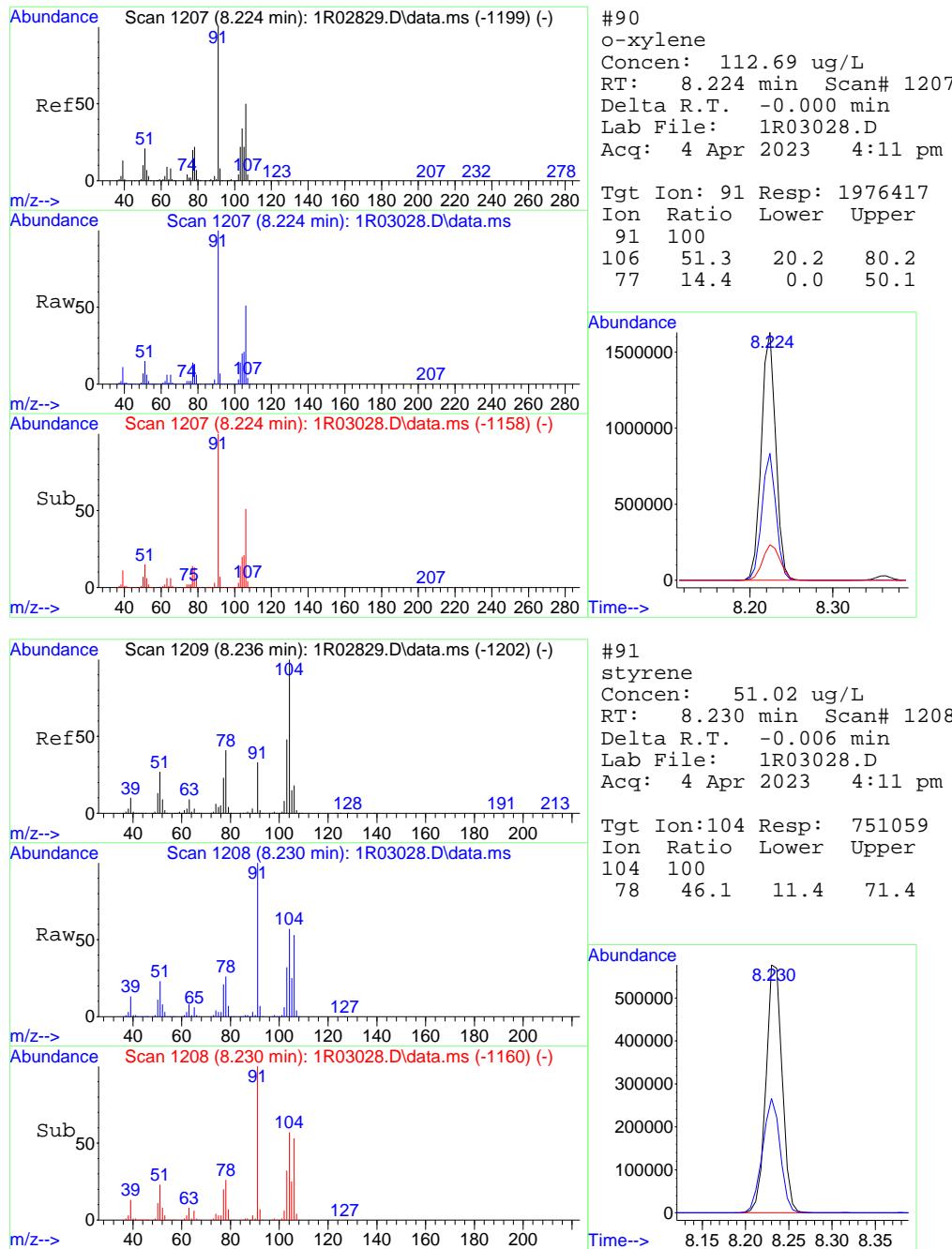


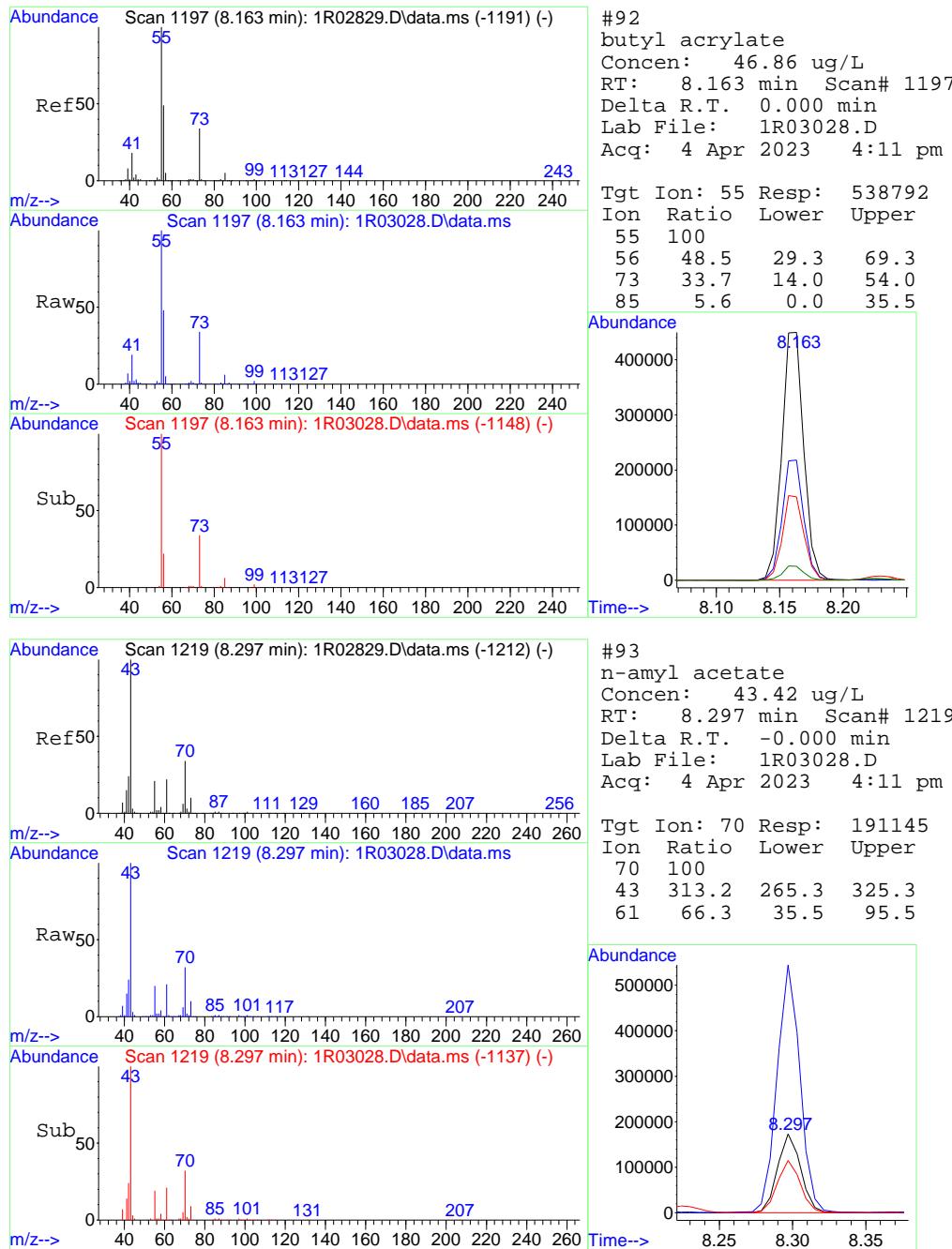


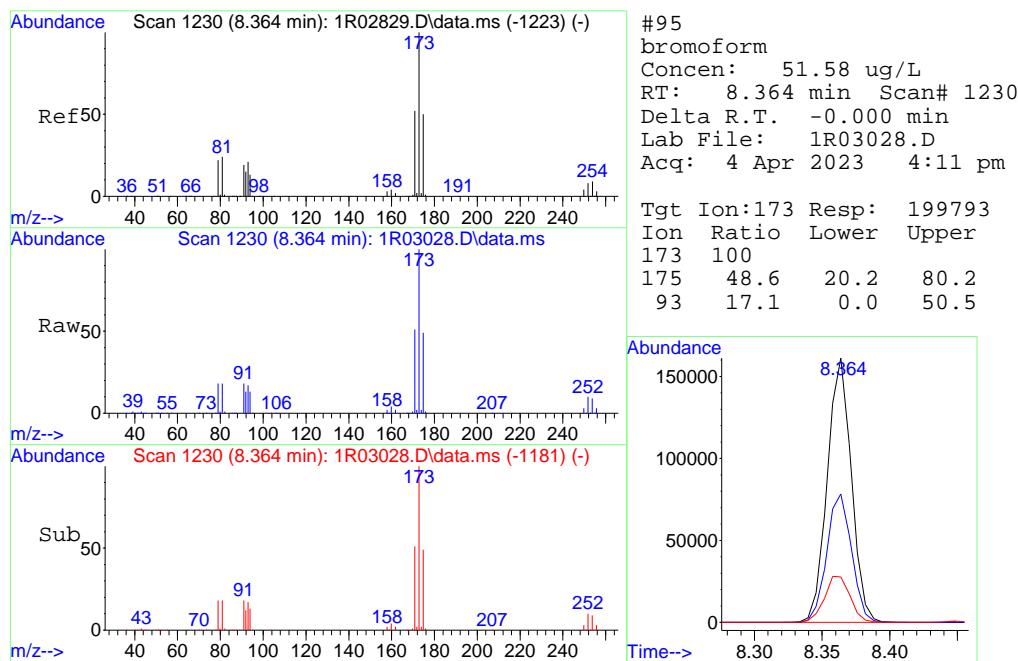
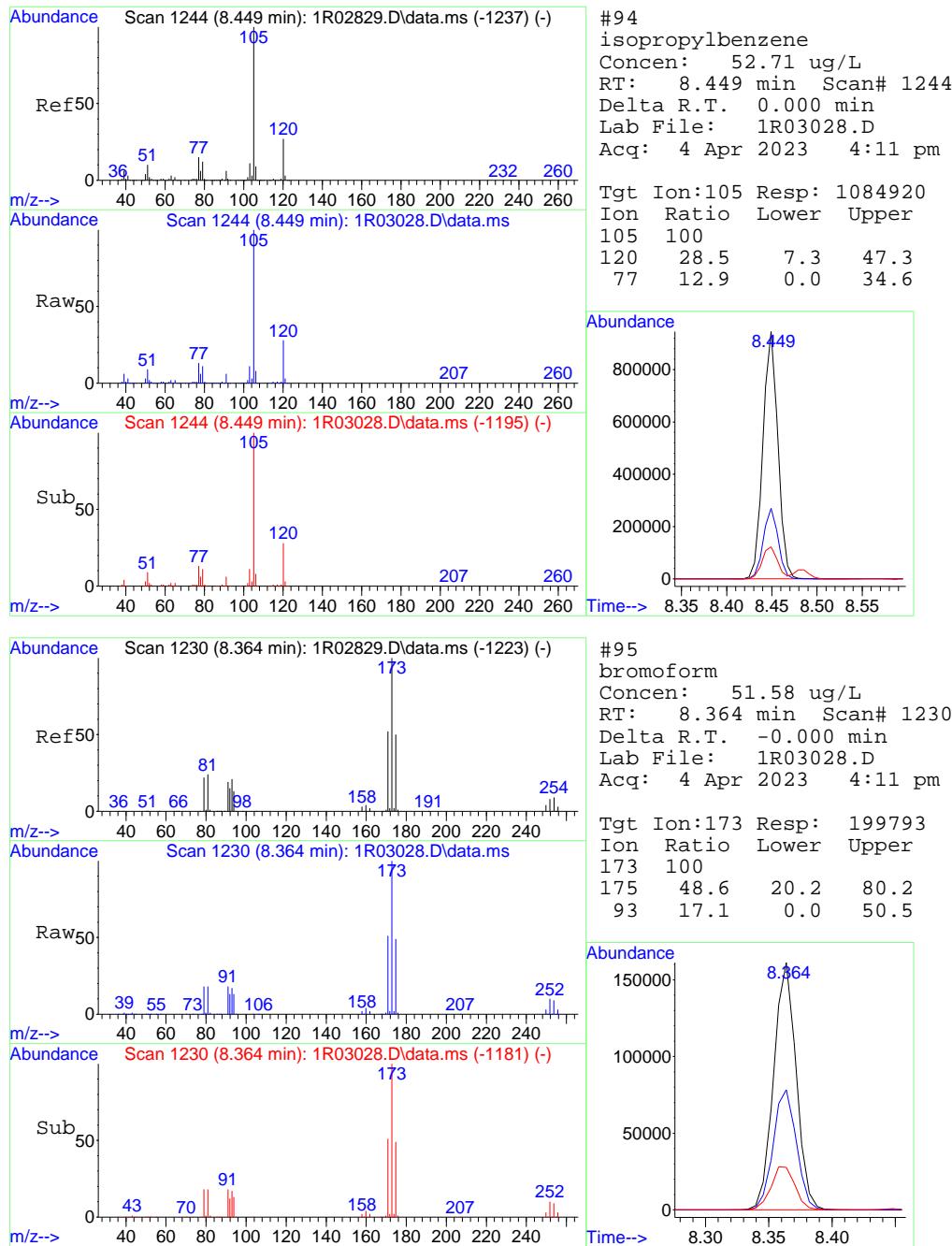


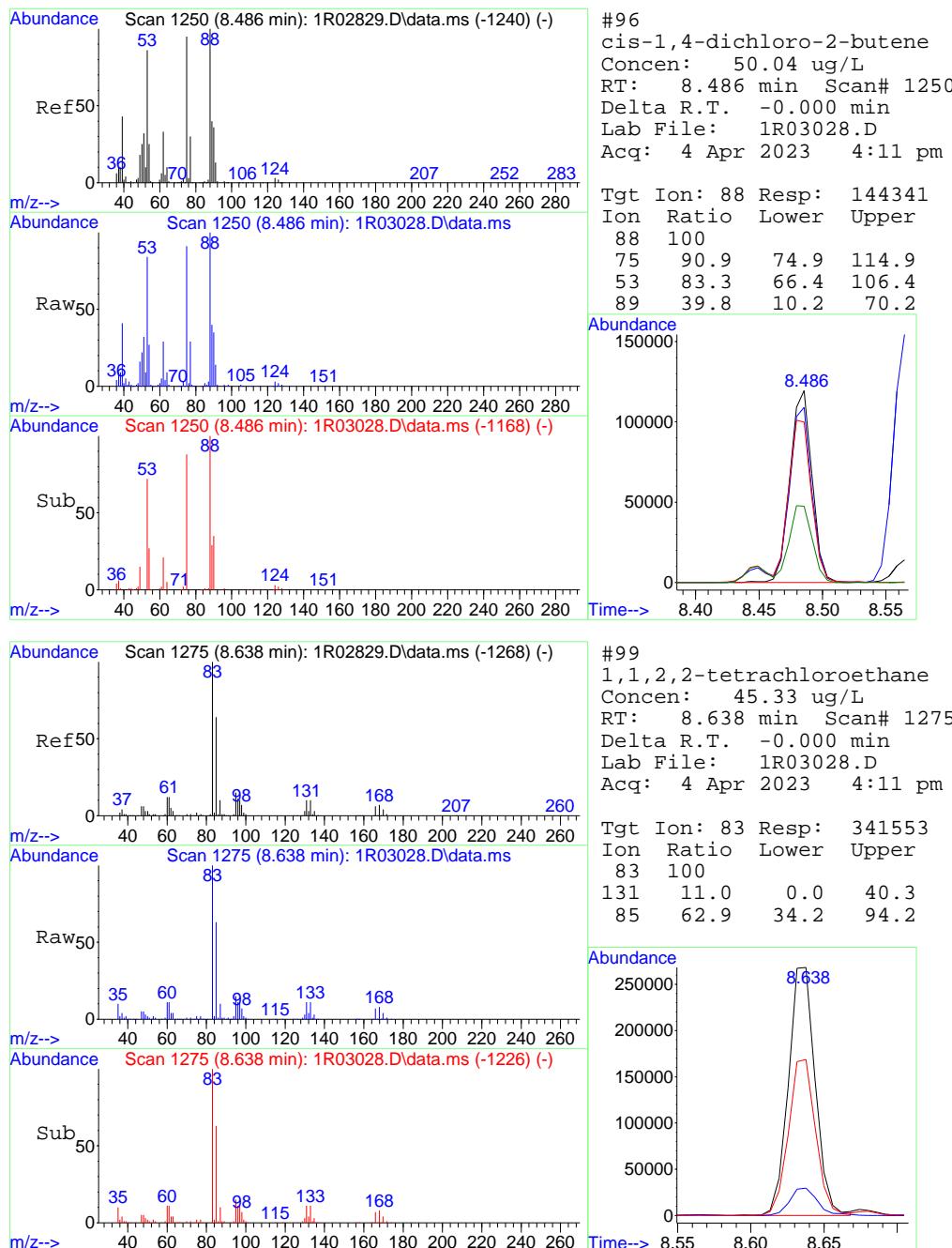


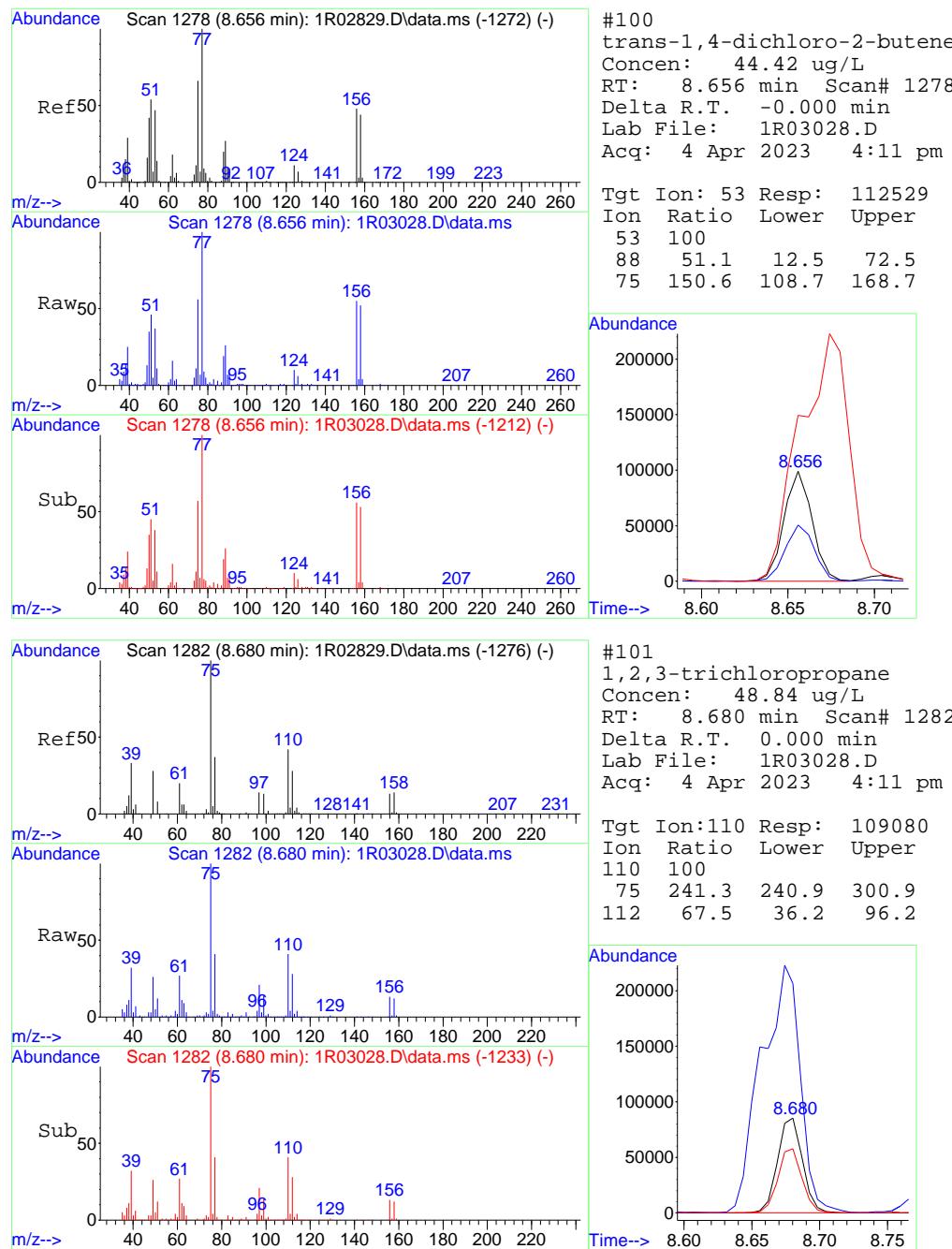


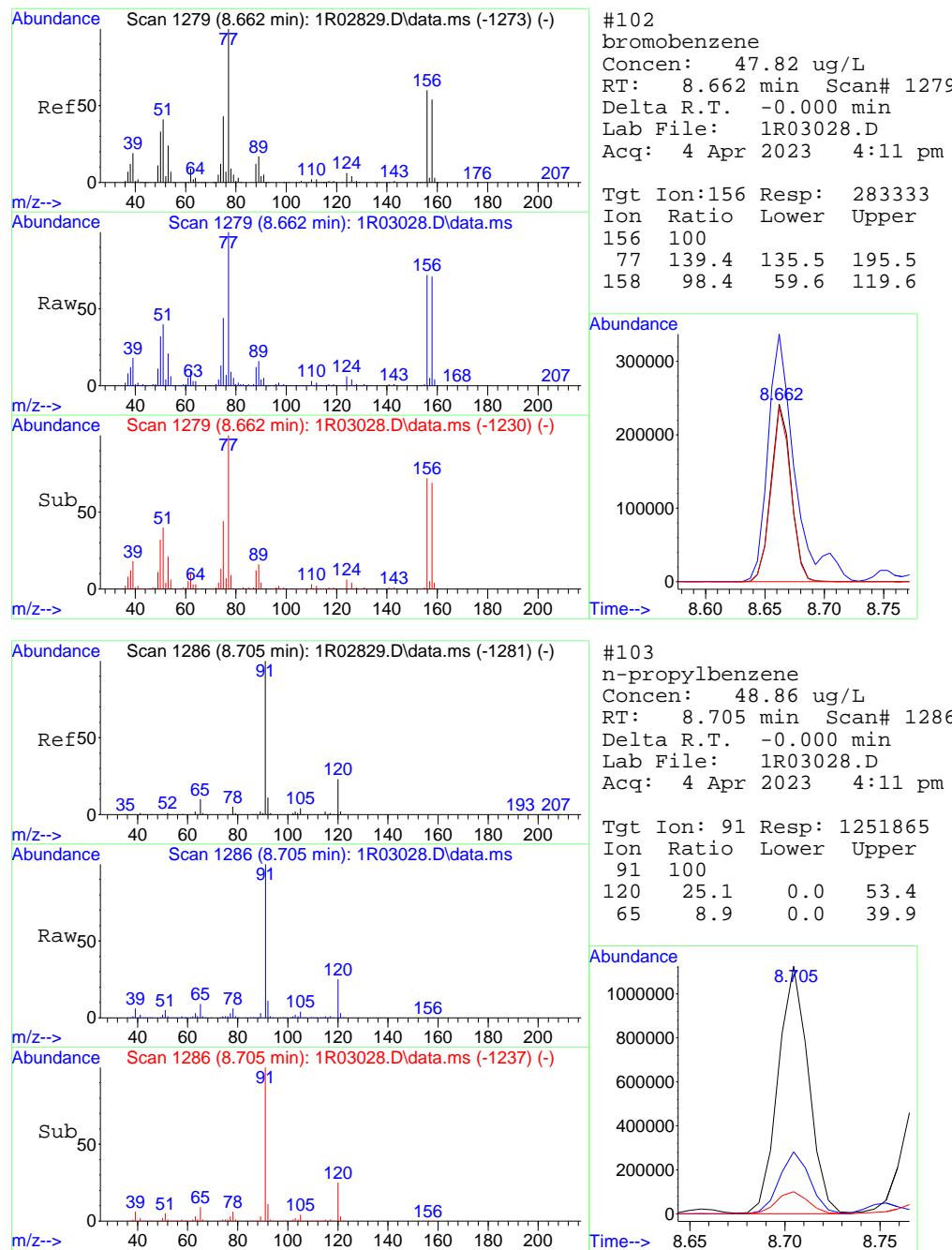


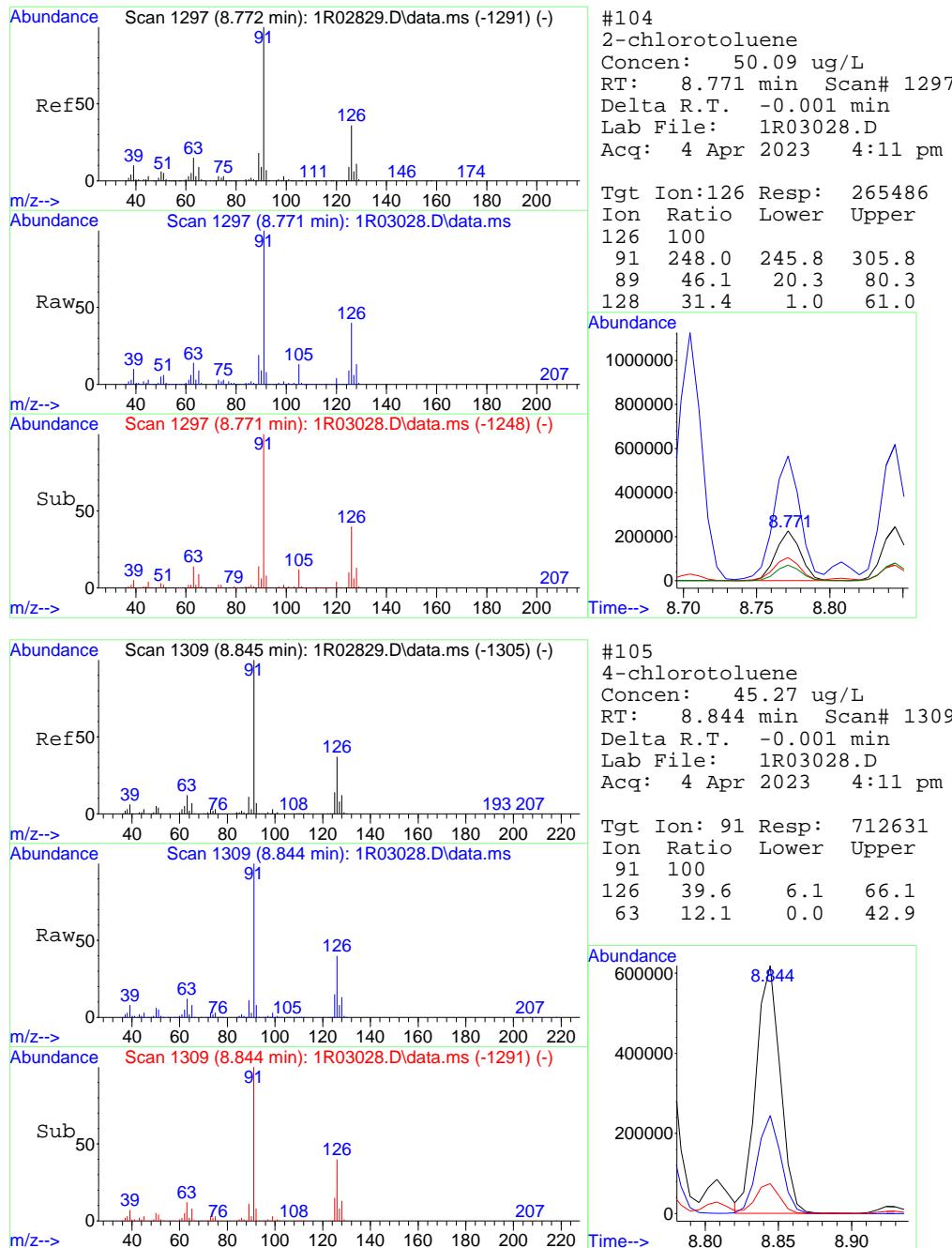


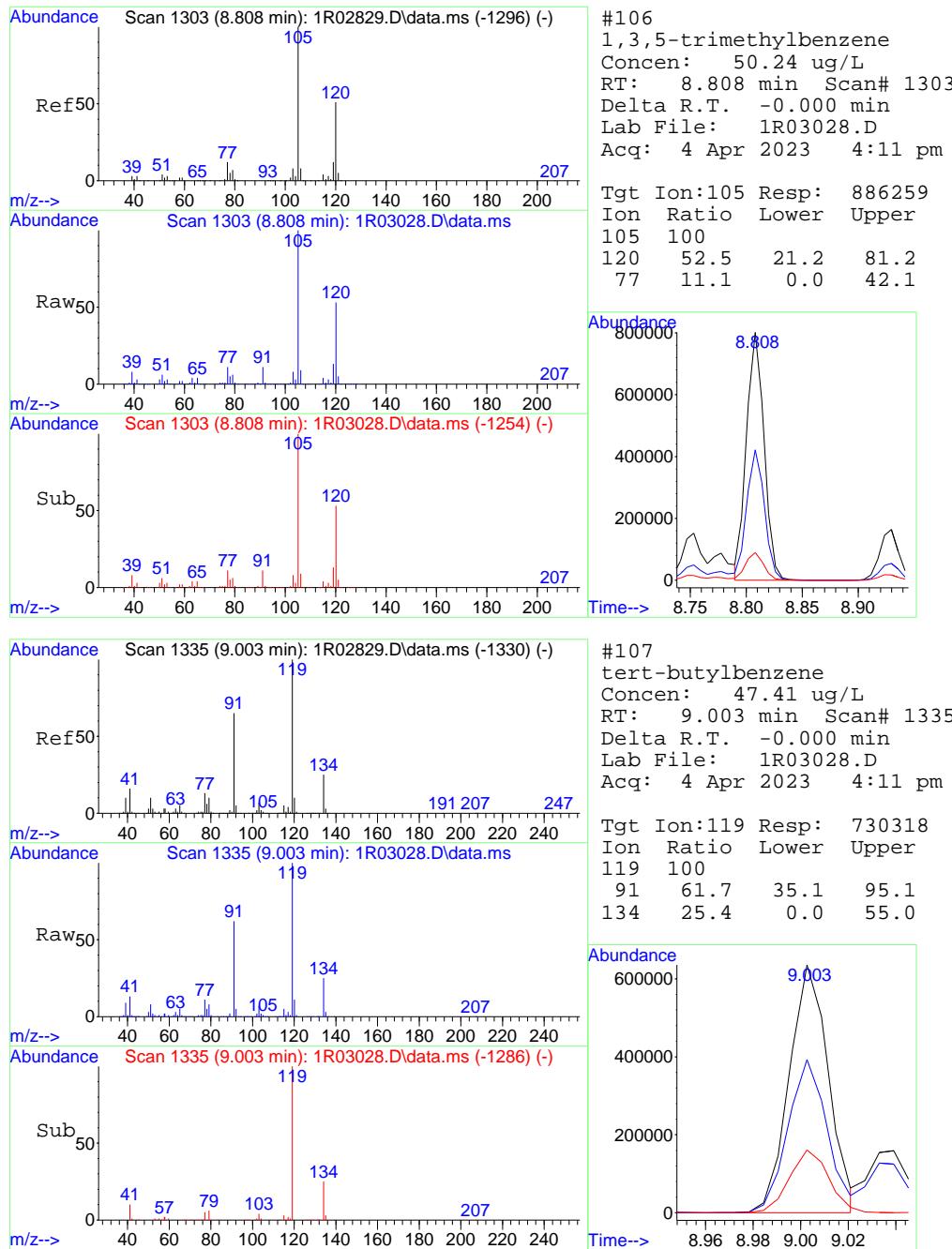


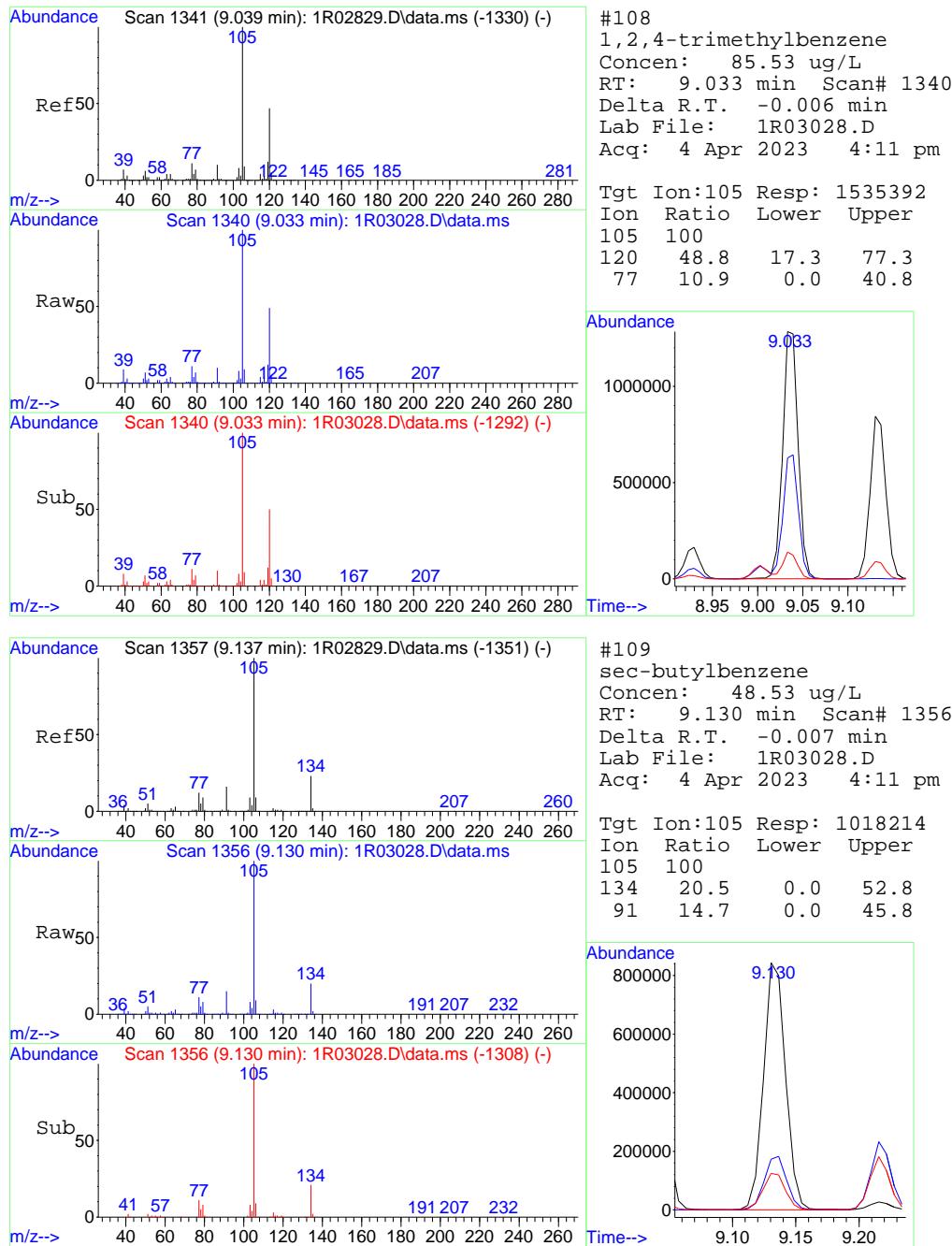


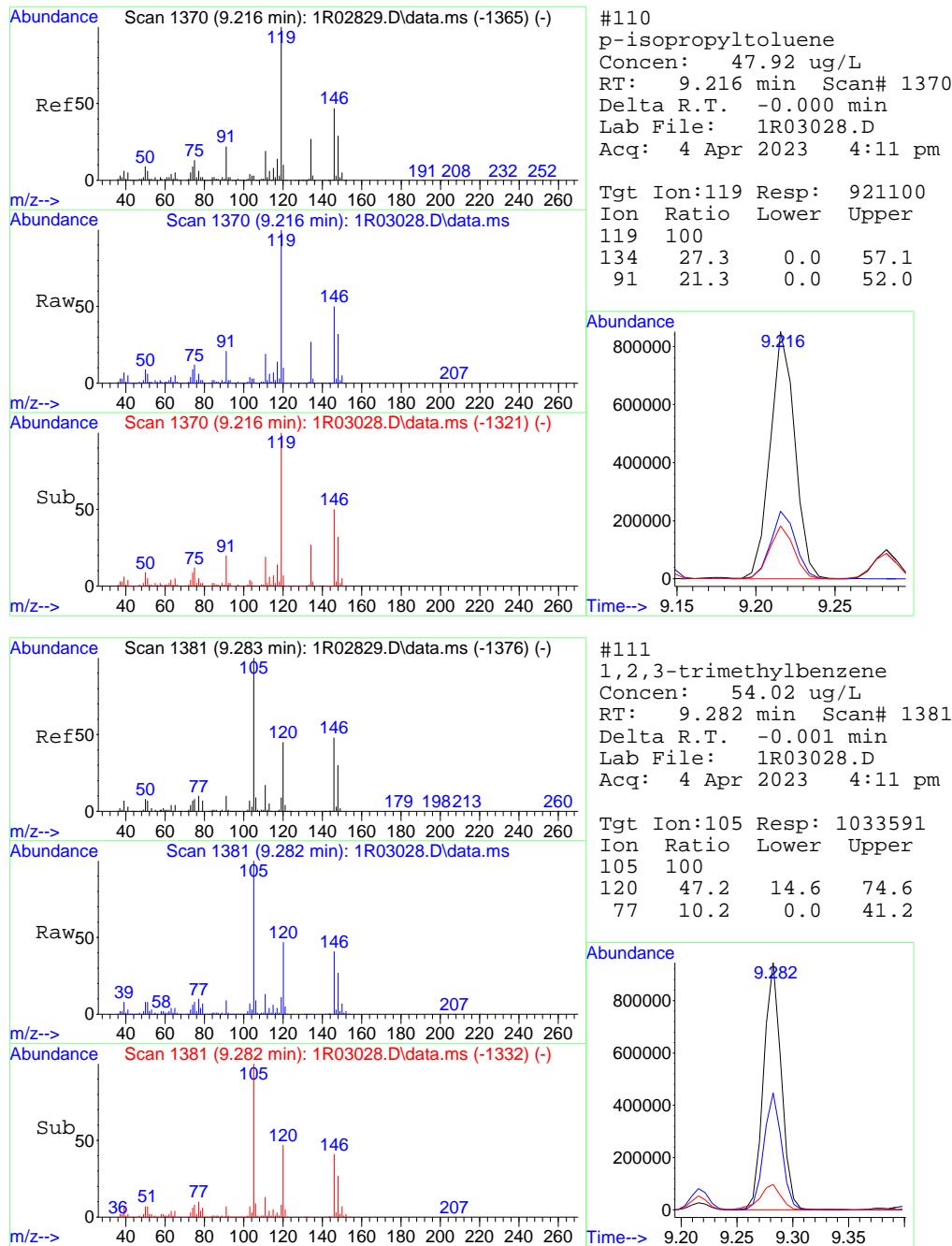


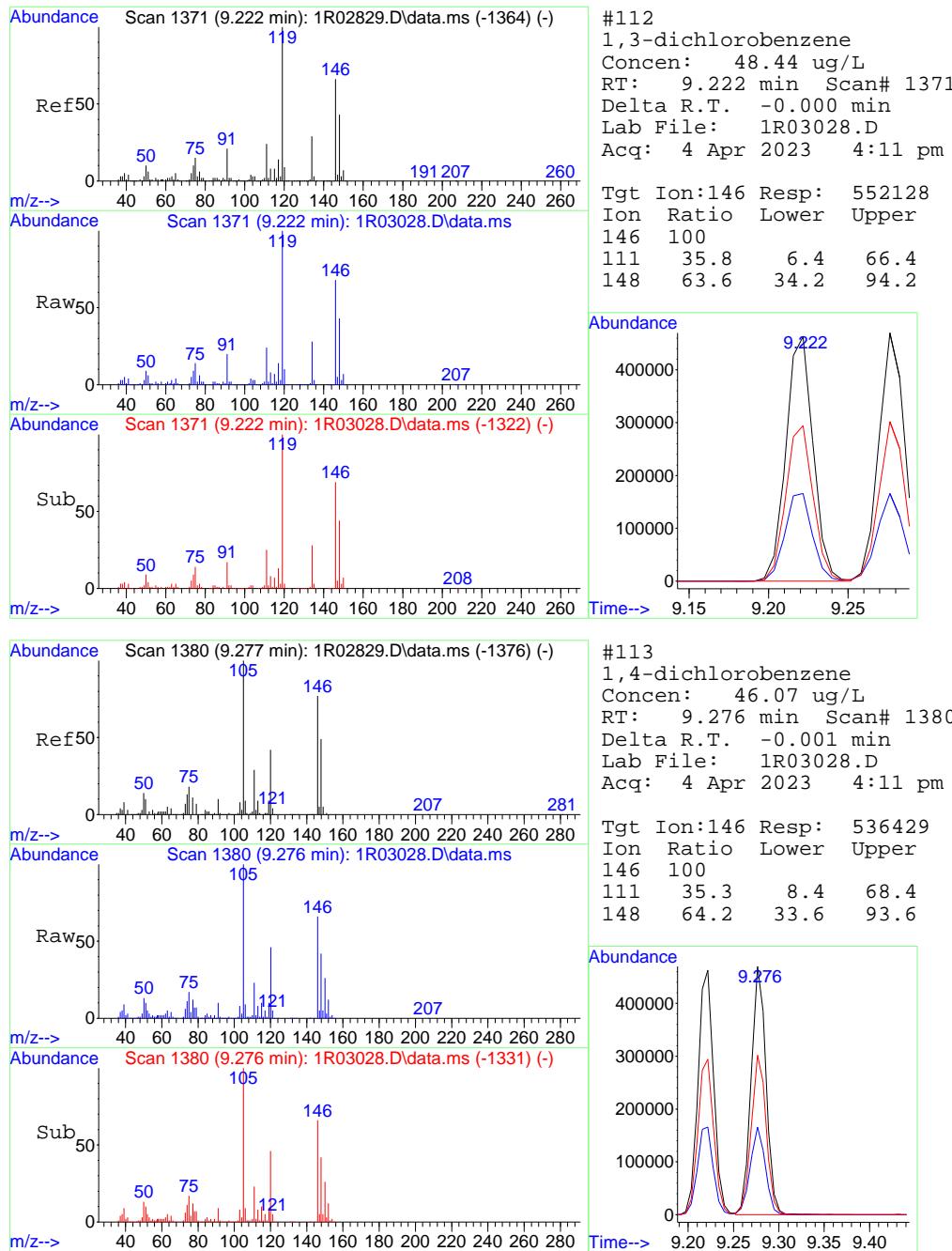


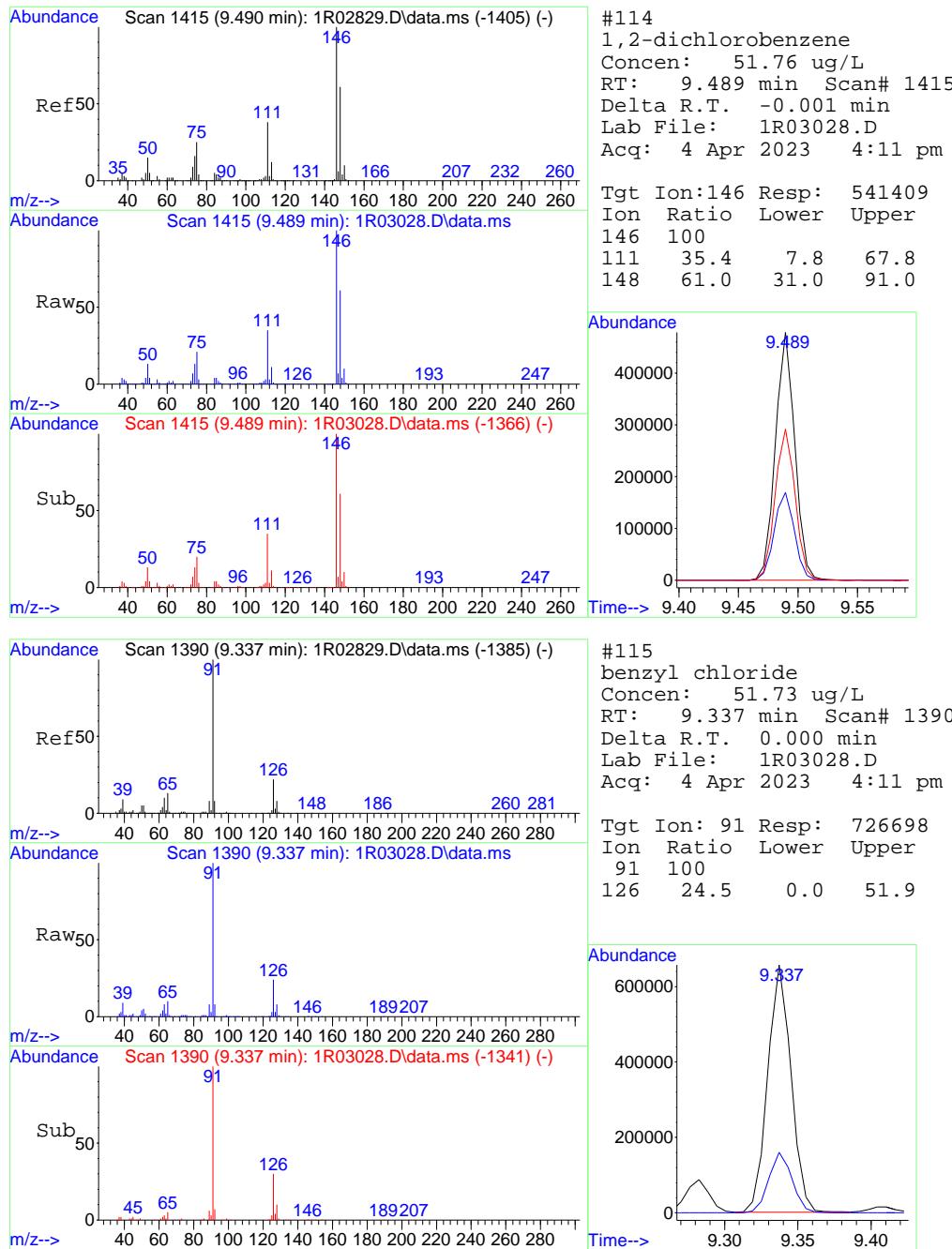


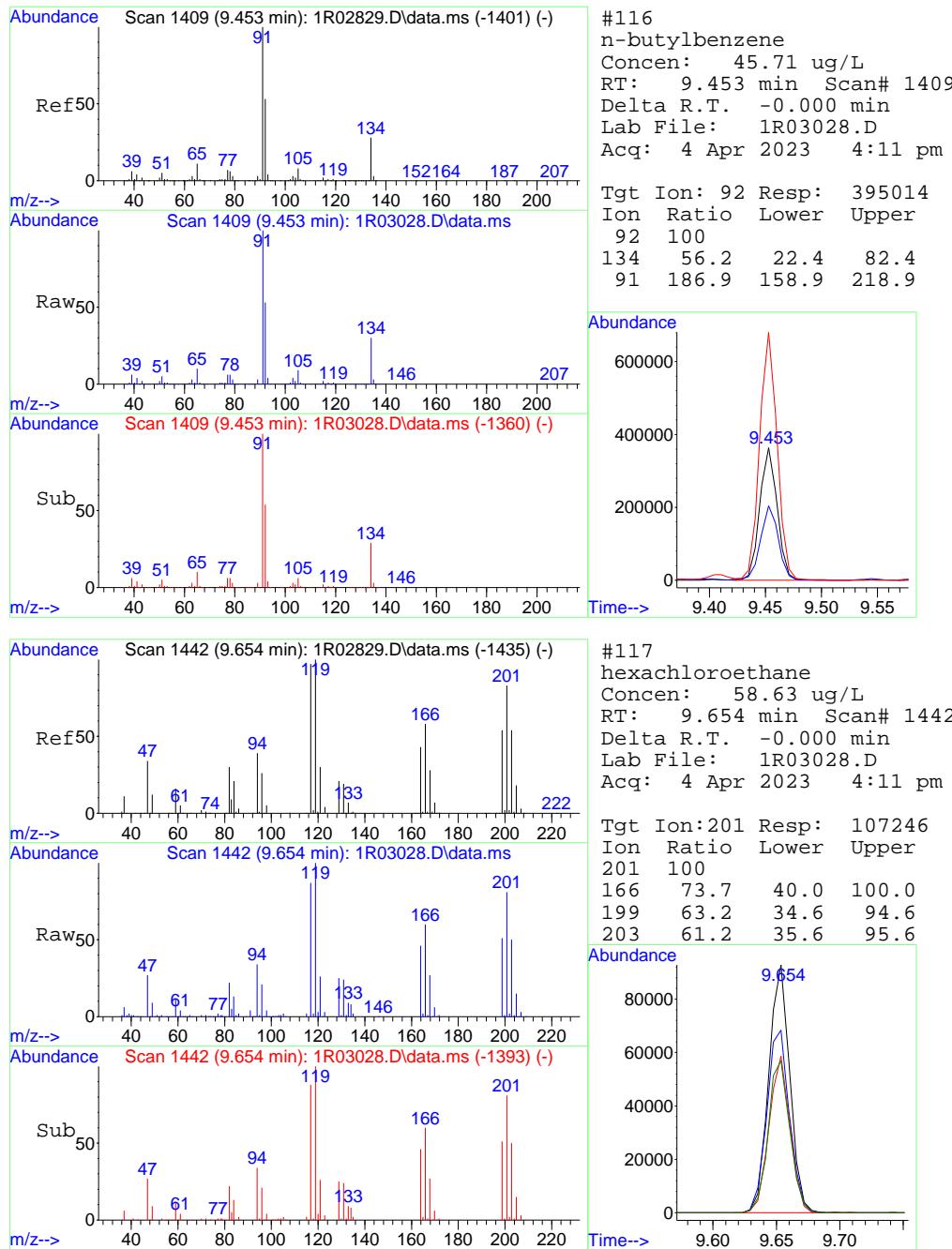


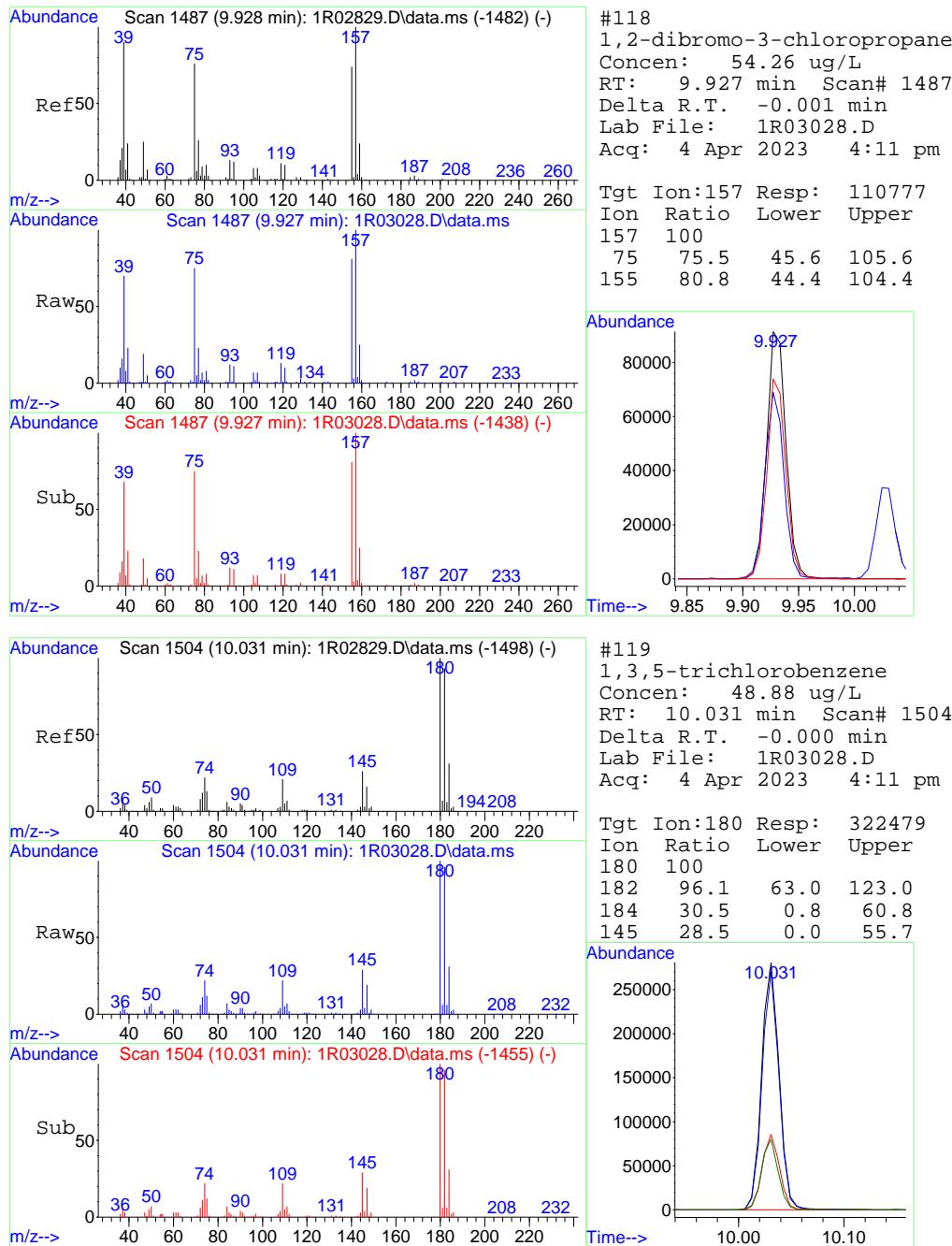


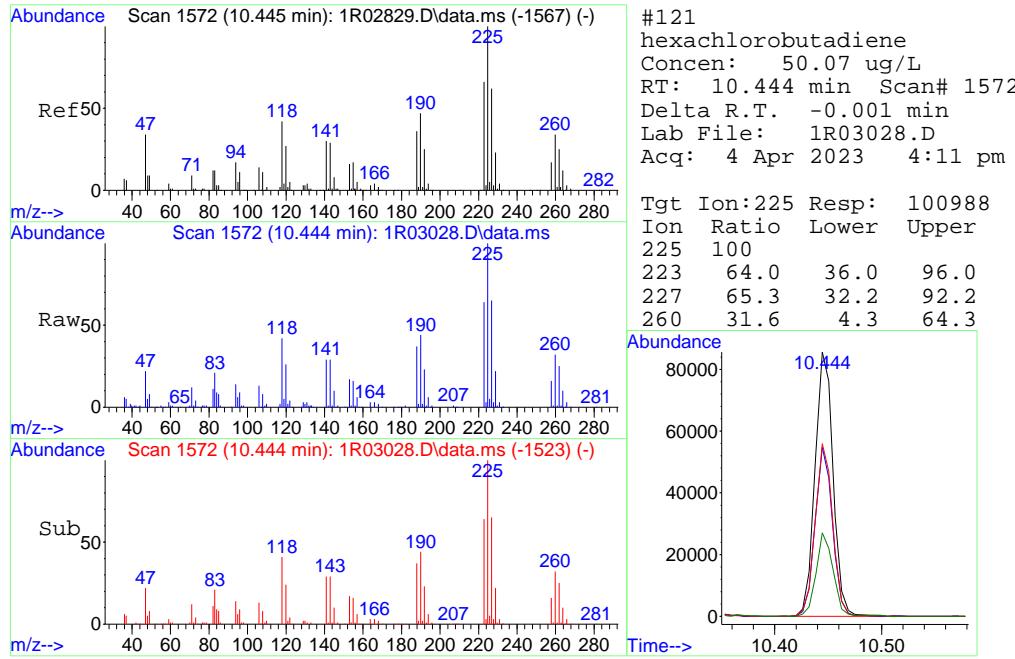
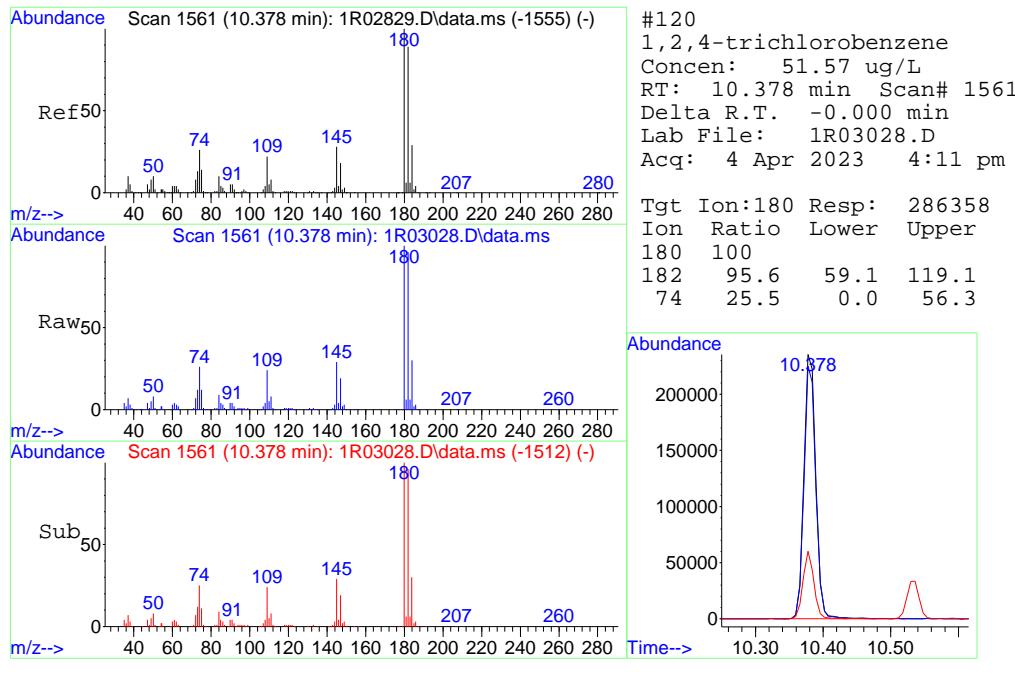


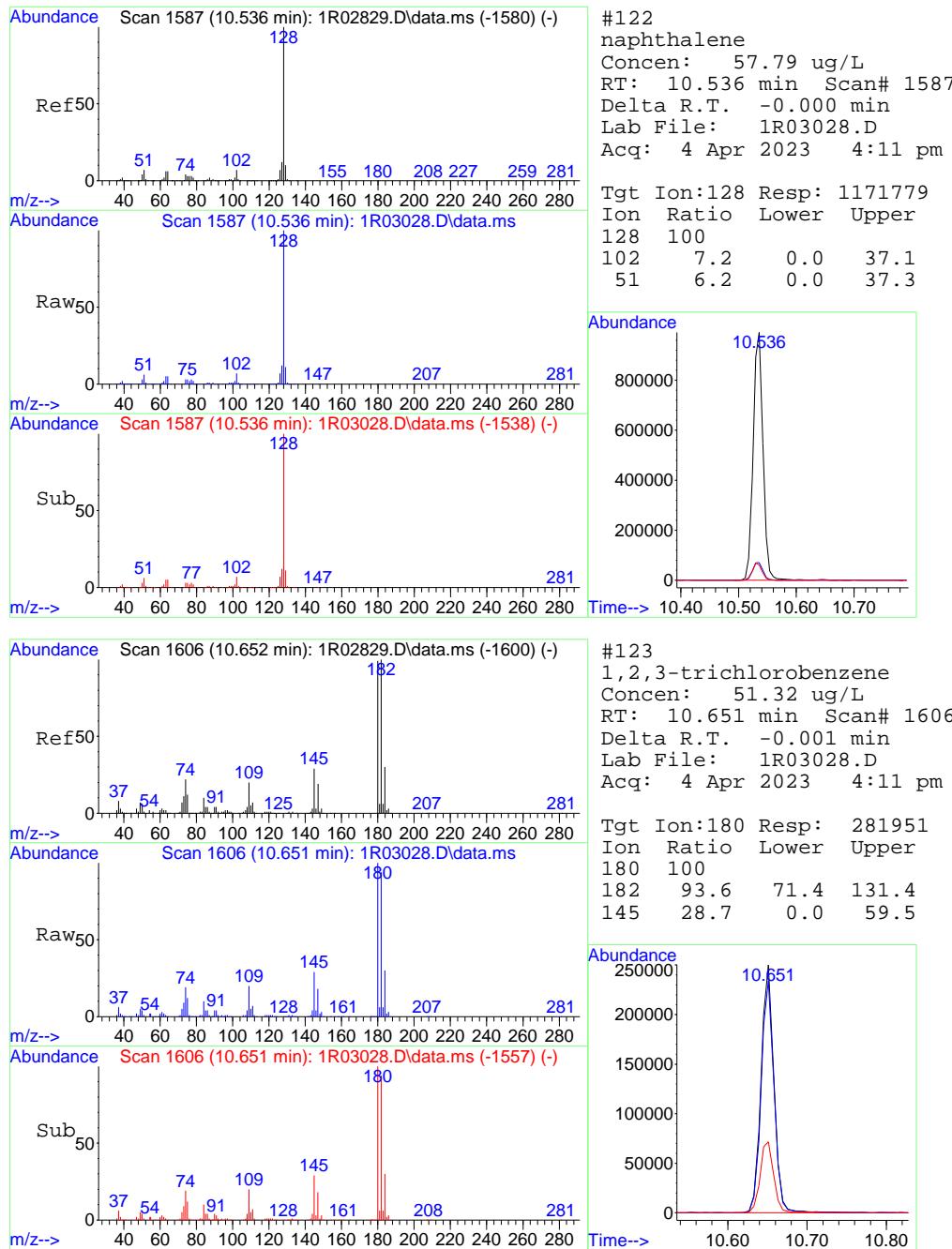


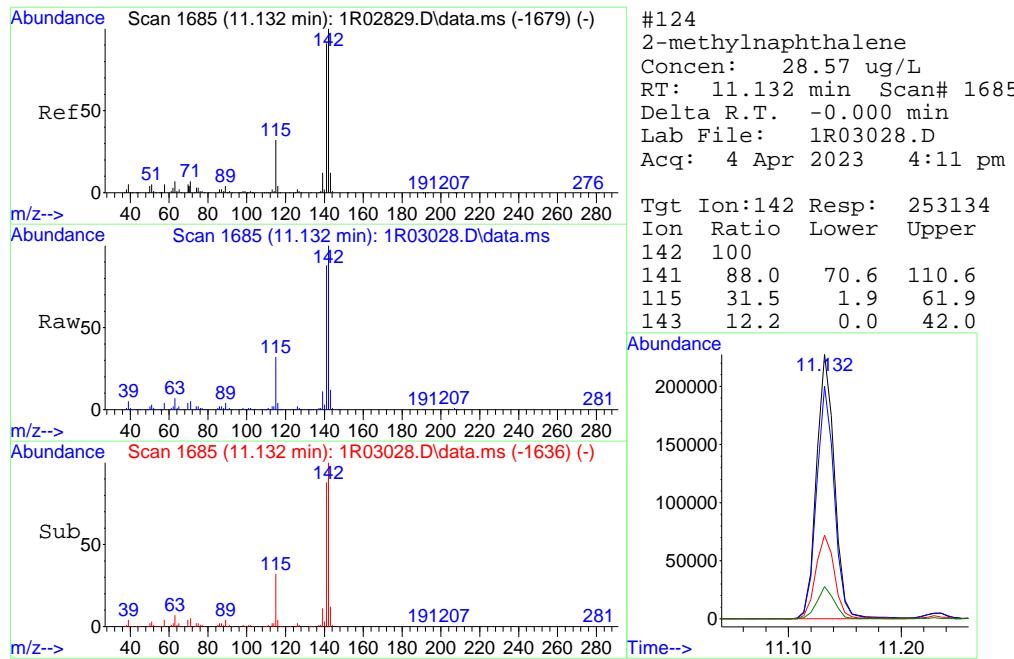












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03029.D  
 Acq On : 4 Apr 2023 4:36 pm  
 Operator : nickw  
 Sample : jd62953-1msd  
 Misc : MS67936,V1R0100,5,,,4  
 ALS Vial : 18 Sample Multiplier: 1

Quant Time: Apr 05 09:50:01 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.491	65	584603	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	697532	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	1038223	50.00	ug/L	0.00
73) chlorobenzene-d5	7.817	117	1013024	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	549087	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) dibromofluoromethane (s)	5.018	113	317667	52.40	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 104.80%			
53) 1,2-dichloroethane-d4 (s)	5.304	65	298996	49.53	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery = 99.06%			
74) toluene-d8 (s)	6.801	98	1292044	46.19	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 92.38%			
98) 4-bromofluorobenzene (s)	8.565	174	442538	49.25	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 98.50%			

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) ethanol	2.688	45	521423	4944.43	ug/L	96
3) tertiary butyl alcohol	3.570	59	351576	239.73	ug/L	99
4) 1,4-dioxane	6.119	88	122594	1121.53	ug/L	98
6) chlorodifluoromethane	1.654	51	246055	46.23	ug/L	97
7) dichlorodifluoromethane	1.642	85	269988	54.88	ug/L	97
8) chloromethane	1.806	50	269087	35.17	ug/L	99
9) vinyl chloride	1.897	62	250761	31.22	ug/L	99
10) 1,3-butadiene	1.934	54	207430	34.56	ug/L	94
11) bromomethane	2.189	94	96286	21.42	ug/L	99
12) chloroethane	2.293	64	150452	37.19	ug/L	97
13) trichlorofluoromethane	2.536	101	338779	55.03	ug/L	94
14) ethyl ether	2.804	74	122695	48.37	ug/L	86
15) acrolein	2.938	56	65693	45.00	ug/L	80
16) freon 113	3.041	151	166822	48.07	ug/L	98
17) 1,1-dichloroethene	3.035	96	186872	47.38	ug/L	95
18) acetone	3.071	58	161276	169.23	ug/L	100
19) acetonitrile	3.333	41	624500	486.44	ug/L	95
20) iodomethane	3.181	142	184312	61.58	ug/L	97
21) carbon disulfide	3.254	76	510918	44.11	ug/L	99
22) methylene chloride	3.497	84	213388	49.62	ug/L	99
23) methyl acetate	3.357	74	64151	49.92	ug/L #	77
24) methyl tert butyl ether	3.728	73	610560	49.71	ug/L	99
25) trans-1,2-dichloroethene	3.741	96	210969	48.18	ug/L	95
26) hexane	3.984	56	171032	46.08	ug/L	97
27) di-isopropyl ether	4.142	45	802316	46.04	ug/L	99
28) 2-butanone	4.592	72	205208	183.20	ug/L	96
29) 1,1-dichloroethane	4.130	63	375627	45.70	ug/L	97
30) chloroprene	4.191	53	347584	49.08	ug/L	97
31) acrylonitrile	3.704	53	142571	44.81	ug/L	99
32) vinyl acetate	4.118	86	57370	51.16	ug/L #	81
33) ethyl tert-butyl ether	4.452	59	677746	49.24	ug/L	98
34) ethyl acetate	4.623	45	74087	47.16	ug/L #	70
35) 2,2-dichloropropane	4.623	77	303016	50.37	ug/L	88
36) cis-1,2-dichloroethene	4.611	96	231912	47.38	ug/L	87
37) propionitrile	4.653	54	676758	468.12	ug/L	95
38) methyl acrylate	4.665	85	56527	47.82	ug/L #	84
39) methacrylonitrile	4.775	67	142687	45.66	ug/L	90
40) bromochloromethane	4.811	128	117510	50.66	ug/L #	84
41) tetrahydrofuran	4.823	42	171577	44.03	ug/L	95
42) chloroform	4.890	83	381233	47.40	ug/L	96
43) tert-Butyl Formate	4.909	59	189820	43.32	ug/L	96

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03029.D  
 Acq On : 4 Apr 2023 4:36 pm  
 Operator : nickw  
 Sample : jd62953-1msd  
 Misc : MS67936,V1R0100,5,,,4  
 ALS Vial : 18 Sample Multiplier: 1

Quant Time: Apr 05 09:50:01 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1,1-trichloroethane	5.036	97	329849	47.58	ug/L	92
46) cyclohexane	5.097	84	301574	46.86	ug/L	86
47) isobutyl alcohol	5.213	43	305889	520.11	ug/L	98
48) 1,1-dichloropropene	5.164	75	275840	47.49	ug/L	95
49) carbon tetrachloride	5.164	117	304665	52.72	ug/L	94
50) tert-amyl alcohol	5.310	73	189967	272.30	ug/L	96
51) isopropyl acetate	5.334	87	74200	47.79	ug/L	95
54) n-butyl alcohol	5.742	56	937499	2405.38	ug/L	97
55) 2,2,4-trimethylpentane	5.407	57	610013	42.35	ug/L	97
56) benzene	5.328	78	921783	45.03	ug/L	99
57) tert-amyl methyl ether	5.414	73	655286	45.03	ug/L	98
58) heptane	5.541	57	130646	40.13	ug/L	95
59) 1,2-dichloroethane	5.365	62	294887	45.23	ug/L	94
60) ethyl acrylate	5.894	55	440311	41.75	ug/L	99
61) trichloroethene	5.852	95	234945	45.50	ug/L	99
62) 2-chloroethyl vinyl ether	6.466	63	6618	1.49	ug/L	84
63) methyl methacrylate	6.083	100	82932	47.50	ug/L	#
64) methylcyclohexane	6.046	83	350791	42.88	ug/L	97
65) 1,2-dichloropropane	6.064	63	226354	40.15	ug/L	98
66) dibromomethane	6.131	93	151679	43.54	ug/L	99
67) bromodichloromethane	6.265	83	283311	45.61	ug/L	99
68) 2-nitropropane	6.430	41	131984	45.09	ug/L	95
69) epichlorohydrin	6.509	57	247209	233.65	ug/L	99
70) cis-1,3-dichloropropene	6.594	75	357181	44.69	ug/L	98
71) 4-methyl-2-pentanone	6.691	58	661762	170.10	ug/L	96
72) isoamyl alcohol	6.721	70	334574	923.25	ug/L	96
75) toluene	6.849	92	3335999	254.22	ug/L	97
76) ethyl methacrylate	7.026	69	337334	47.12	ug/L	99
77) trans-1,3-dichloropropene	7.007	75	333844	47.55	ug/L	91
78) 1,1,2-trichloroethane	7.153	83	179872	43.83	ug/L	94
79) tetrachloroethene	7.220	164	233898	45.25	ug/L	96
80) 2-hexanone	7.287	58	741018	178.77	ug/L	99
81) 1,3-dichloropropane	7.269	76	333731	45.53	ug/L	97
82) butyl acetate	7.354	56	240100	44.22	ug/L	93
83) dibromochloromethane	7.421	129	268967	52.03	ug/L	100
84) 1,2-dibromoethane	7.512	107	250696	51.05	ug/L	98
85) n-butyl ether	7.865	57	989980	42.57	ug/L	99
86) chlorobenzene	7.835	112	669874	46.17	ug/L	96
87) 1,1,1,2-tetrachloroethane	7.890	131	250801	48.90	ug/L	96
88) ethylbenzene	7.890	91	2060723	85.45	ug/L	98
89) m,p-xylene	7.975	106	1626360	173.74	ug/L	92
90) o-xylene	8.224	91	2059600	110.72	ug/L	95
91) styrene	8.236	104	795496	50.95	ug/L	97
92) butyl acrylate	8.163	55	561607	46.06	ug/L	97
93) n-amyl acetate	8.297	70	198717	42.56	ug/L	89
94) isopropylbenzene	8.449	105	1131272	51.82	ug/L	96
95) bromoform	8.364	173	220027	53.56	ug/L	95
96) cis-1,4-dichloro-2-butene	8.480	88	151453	49.51	ug/L	95
99) 1,1,2,2-tetrachloroethane	8.638	83	354239	42.92	ug/L	99
100) trans-1,4-dichloro-2-b...	8.656	53	115518	41.63	ug/L	88
101) 1,2,3-trichloropropane	8.680	110	114692	46.88	ug/L	#
102) bromobenzene	8.662	156	298479	45.99	ug/L	85
103) n-propylbenzene	8.705	91	1284960	45.78	ug/L	96
104) 2-chlorotoluene	8.772	126	281772	48.54	ug/L	88
105) 4-chlorotoluene	8.845	91	737037	42.74	ug/L	94
106) 1,3,5-trimethylbenzene	8.808	105	952647	49.30	ug/L	99
107) tert-butylbenzene	9.003	119	780063	46.23	ug/L	97
108) 1,2,4-trimethylbenzene	9.033	105	1606513	81.70	ug/L	99

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03029.D  
 Acq On : 4 Apr 2023 4:36 pm  
 Operator : nickw  
 Sample : jd62953-1msd  
 Misc : MS67936,V1R0100,5,,,4  
 ALS Vial : 18 Sample Multiplier: 1

Quant Time: Apr 05 09:50:01 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration

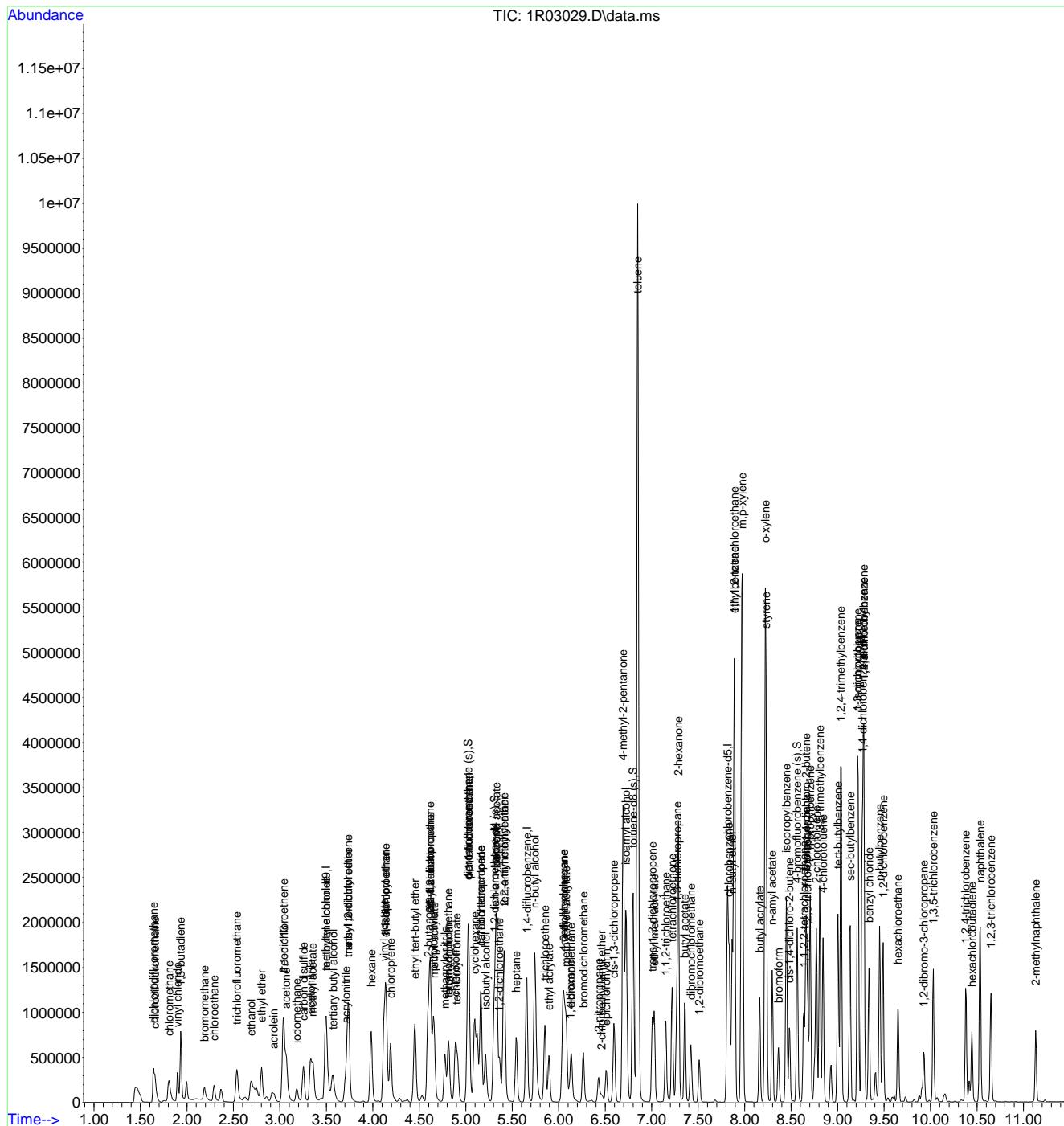
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
109) sec-butylbenzene	9.137	105	1070482	46.57	ug/L	98
110) p-isopropyltoluene	9.216	119	985734	46.82	ug/L	98
111) 1,2,3-trimethylbenzene	9.283	105	1086720	51.85	ug/L	98
112) 1,3-dichlorobenzene	9.222	146	590340	47.28	ug/L	99
113) 1,4-dichlorobenzene	9.277	146	577501	45.28	ug/L	98
114) 1,2-dichlorobenzene	9.490	146	575793	50.25	ug/L	95
115) benzyl chloride	9.337	91	734421	47.73	ug/L	93
116) n-butylbenzene	9.453	92	414684	43.81	ug/L	98
117) hexachloroethane	9.654	201	118033	58.91	ug/L	98
118) 1,2-dibromo-3-chloropr...	9.928	157	122616	54.83	ug/L	96
119) 1,3,5-trichlorobenzene	10.031	180	342386	47.38	ug/L	99
120) 1,2,4-trichlorobenzene	10.378	180	312107	51.31	ug/L	93
121) hexachlorobutadiene	10.445	225	108664	49.19	ug/L	93
122) naphthalene	10.536	128	1240380	55.85	ug/L	98
123) 1,2,3-trichlorobenzene	10.652	180	298596	49.61	ug/L	91
124) 2-methylnaphthalene	11.132	142	272894	28.11	ug/L	98

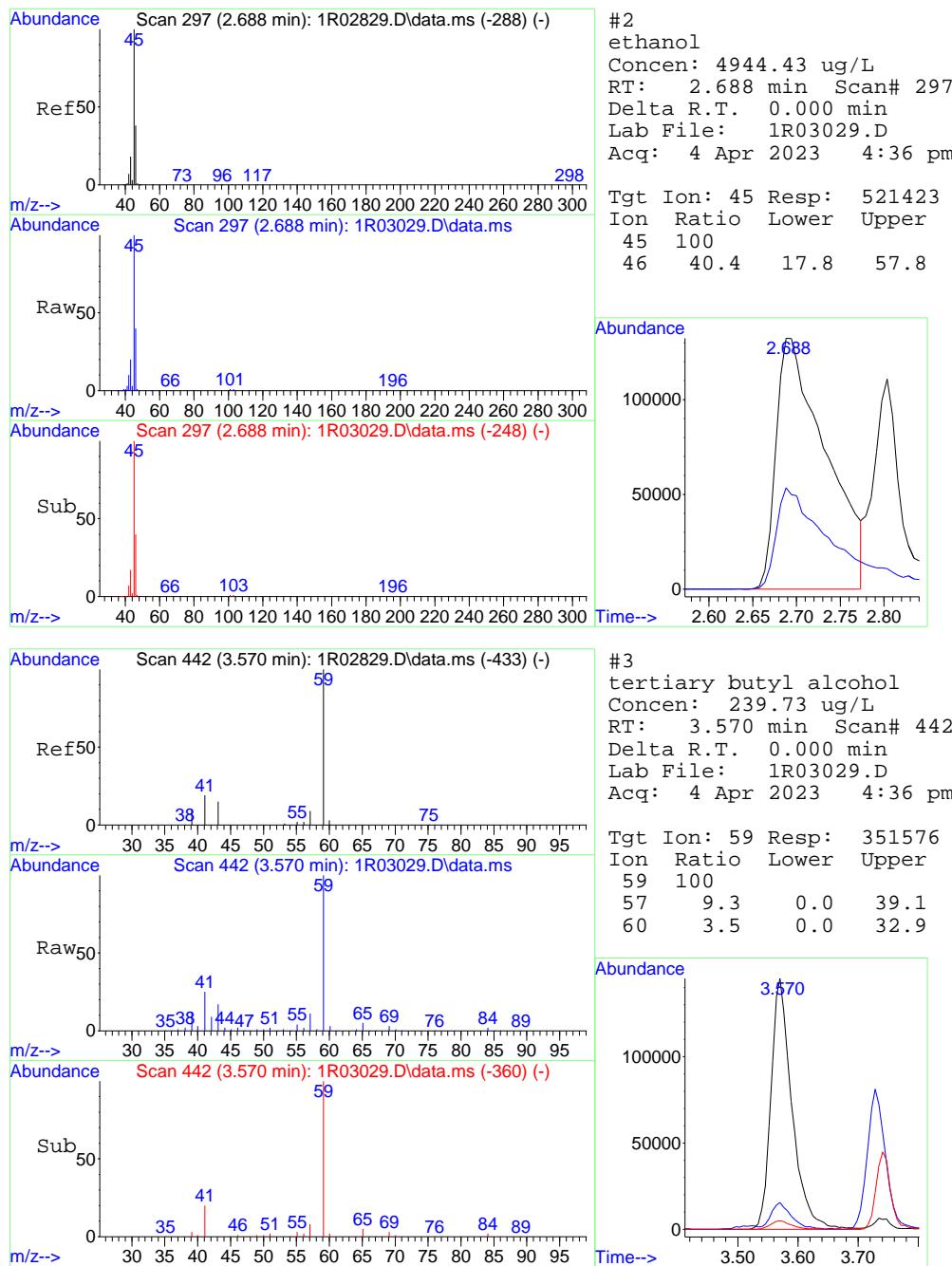
(#) = qualifier out of range (m) = manual integration (+) = signals summed

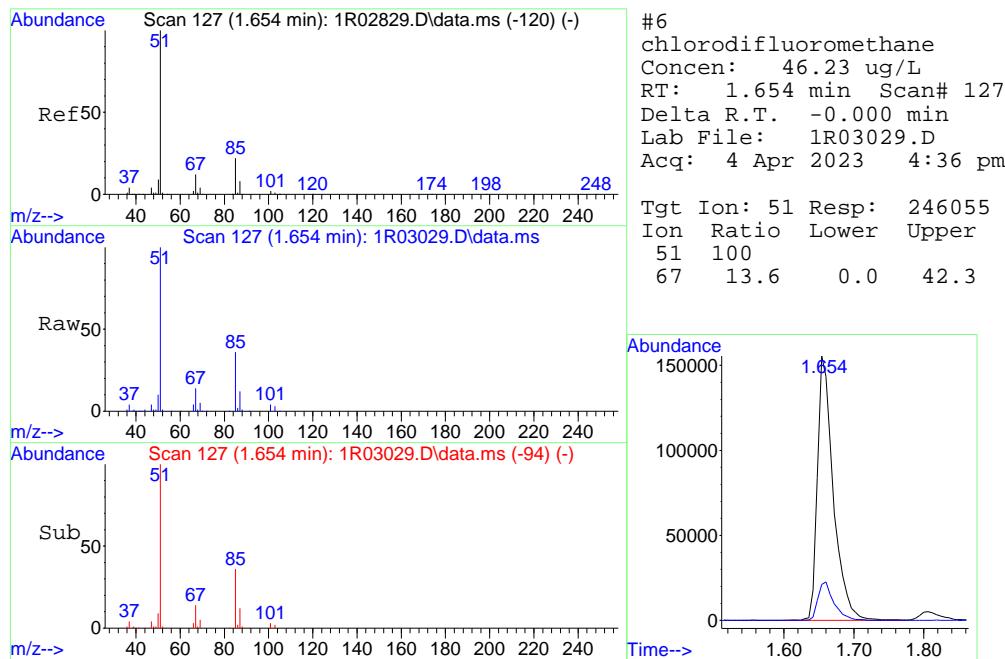
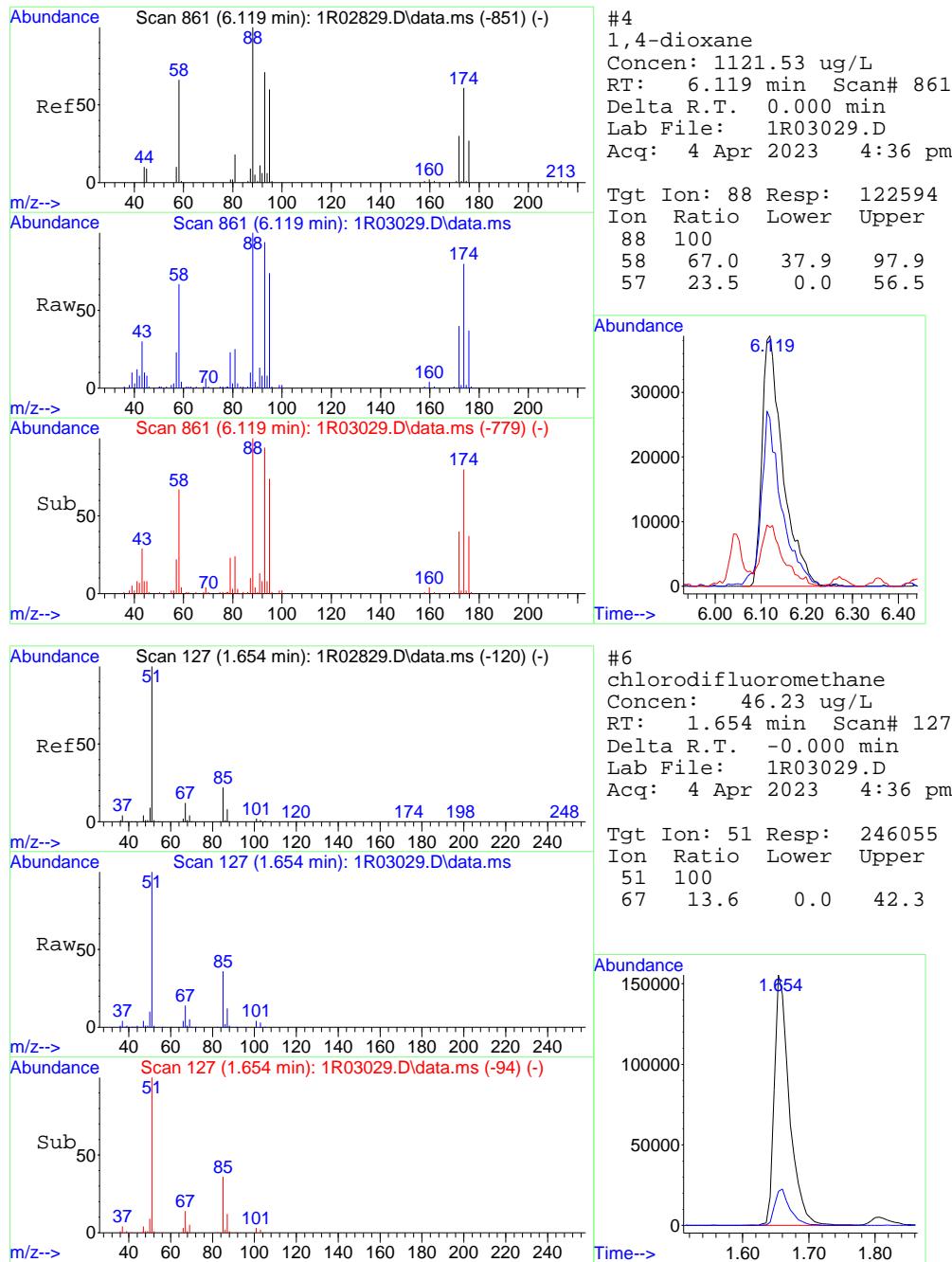
## Quantitation Report (QT Reviewed)

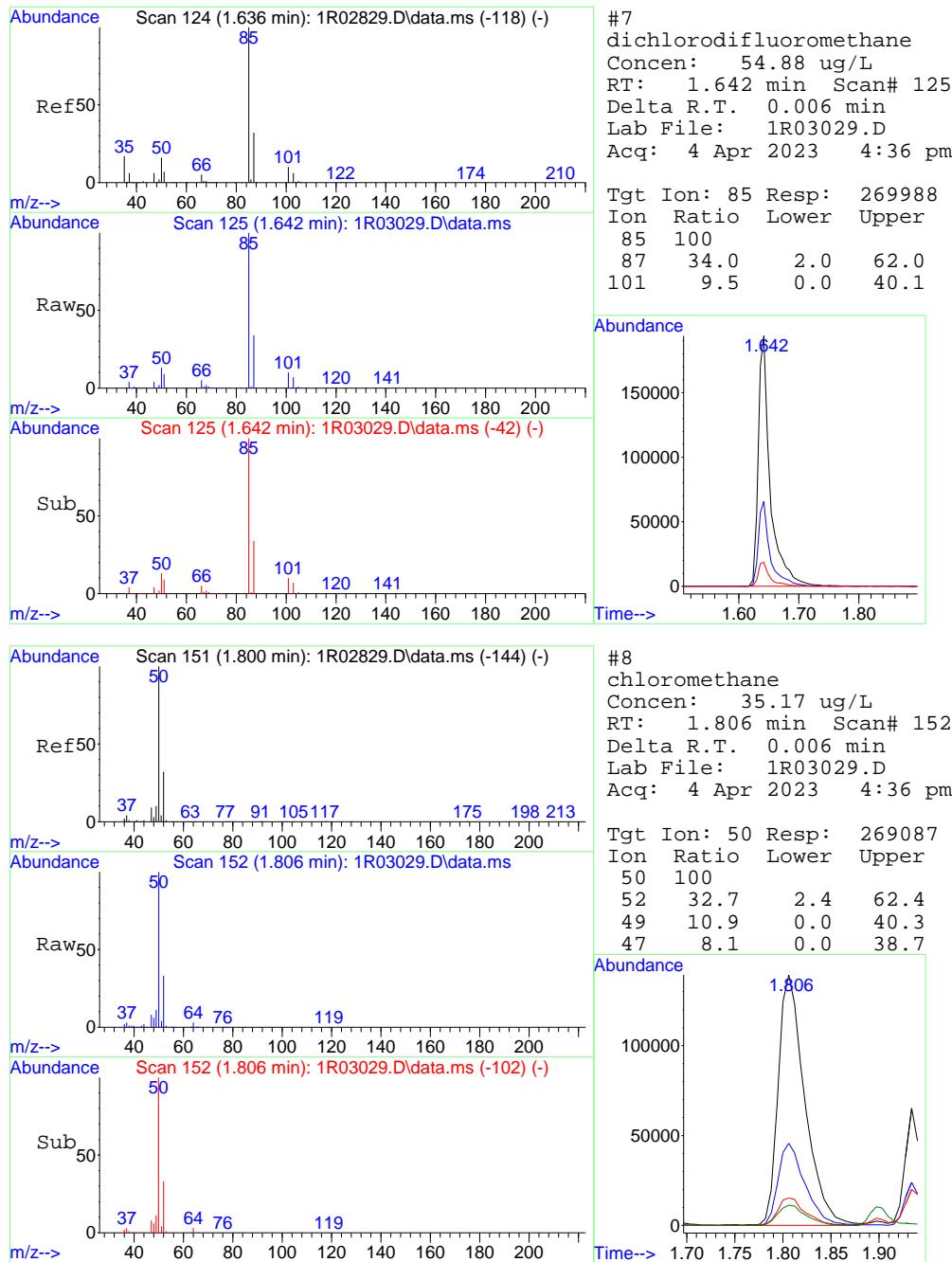
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Acq On : 4 Apr 2023 4:36 pm  
Operator : nickw  
Sample : jd62953-1msd  
Misc : MS67936,V1R0100,5,,,4  
ALS Vial : 18 Sample Multiplier: 1

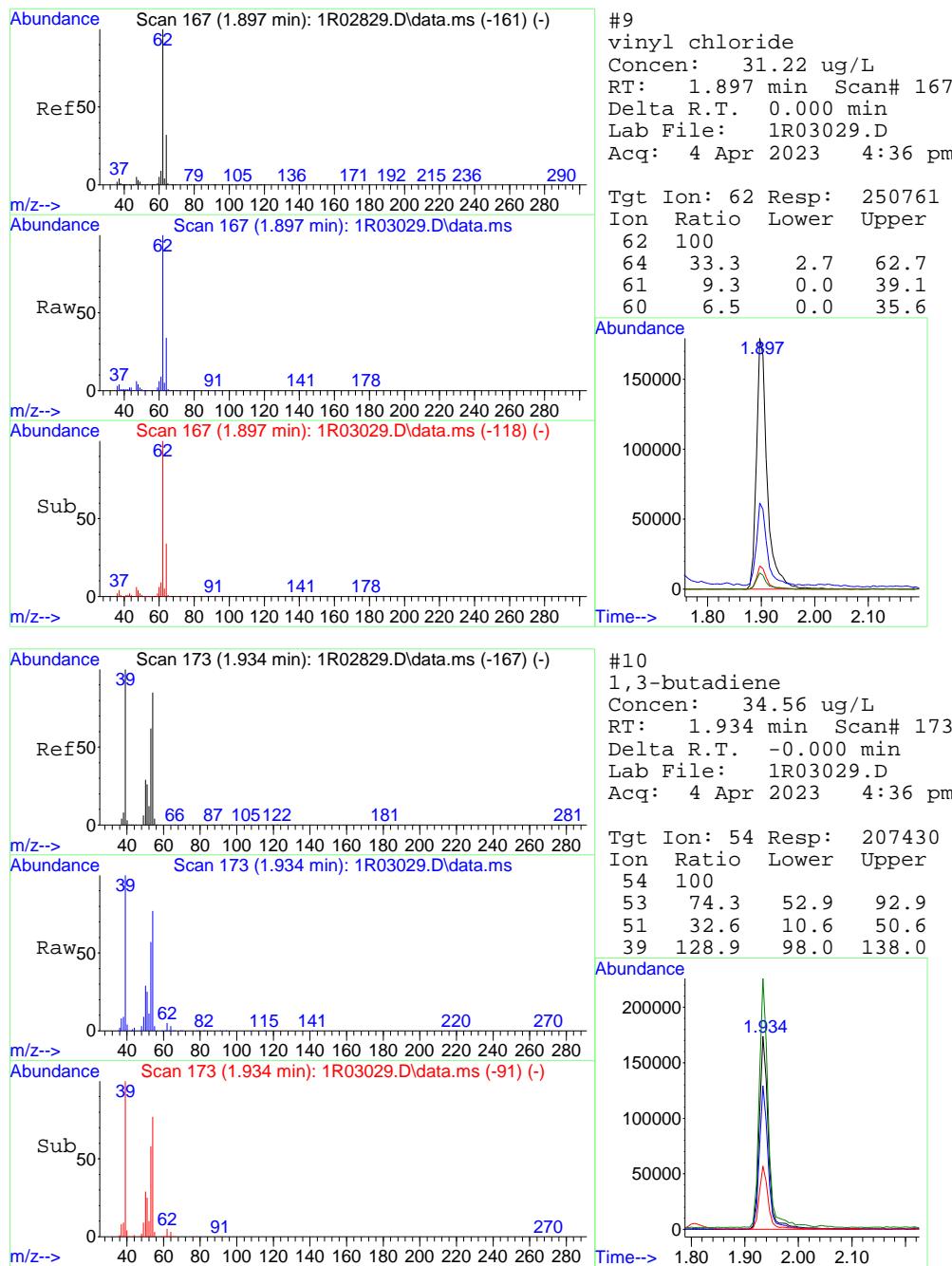
Quant Time: Apr 05 09:50:01 2023  
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
QLast Update : Thu Mar 30 09:13:25 2023  
Response via : Initial Calibration

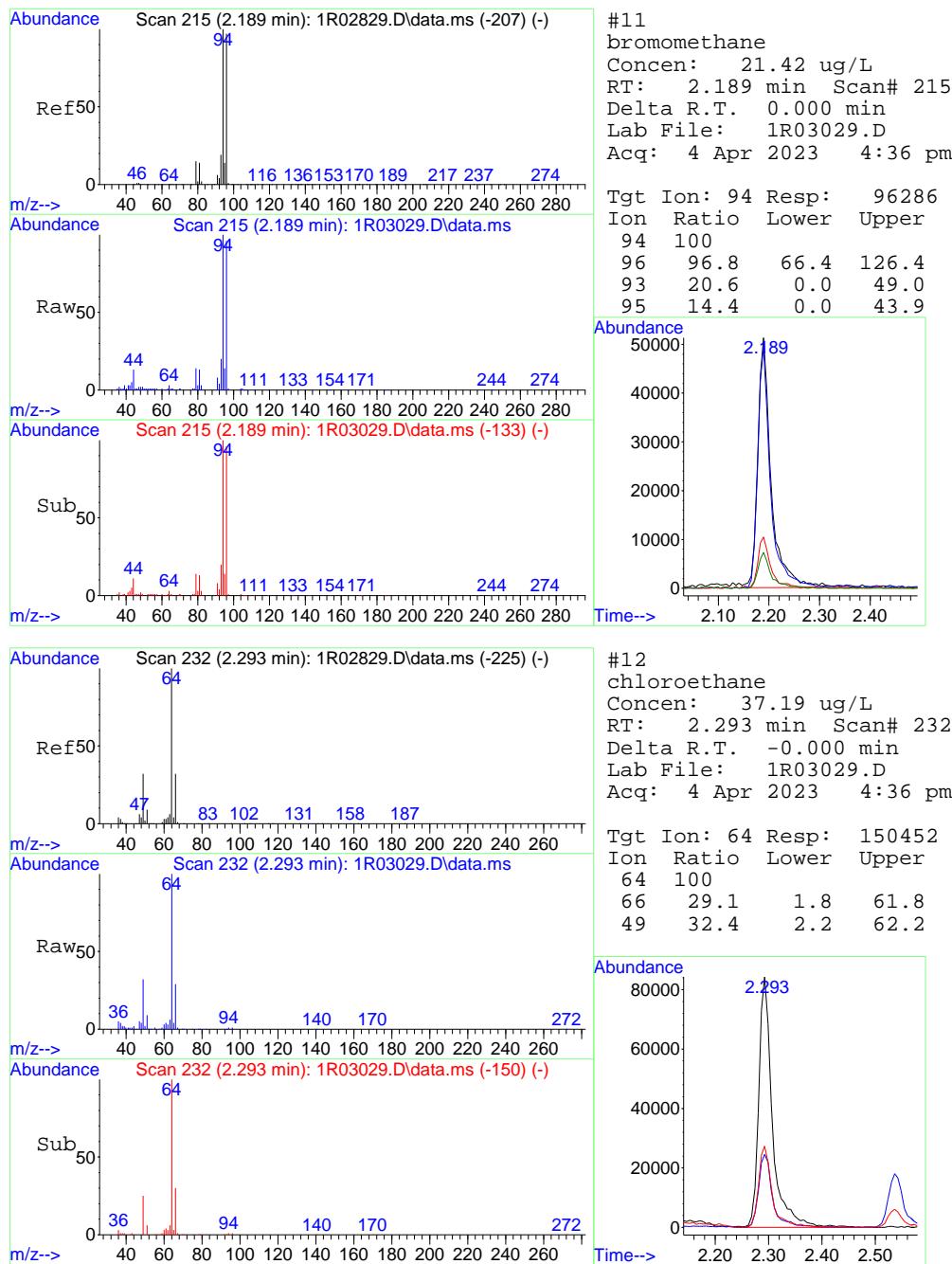


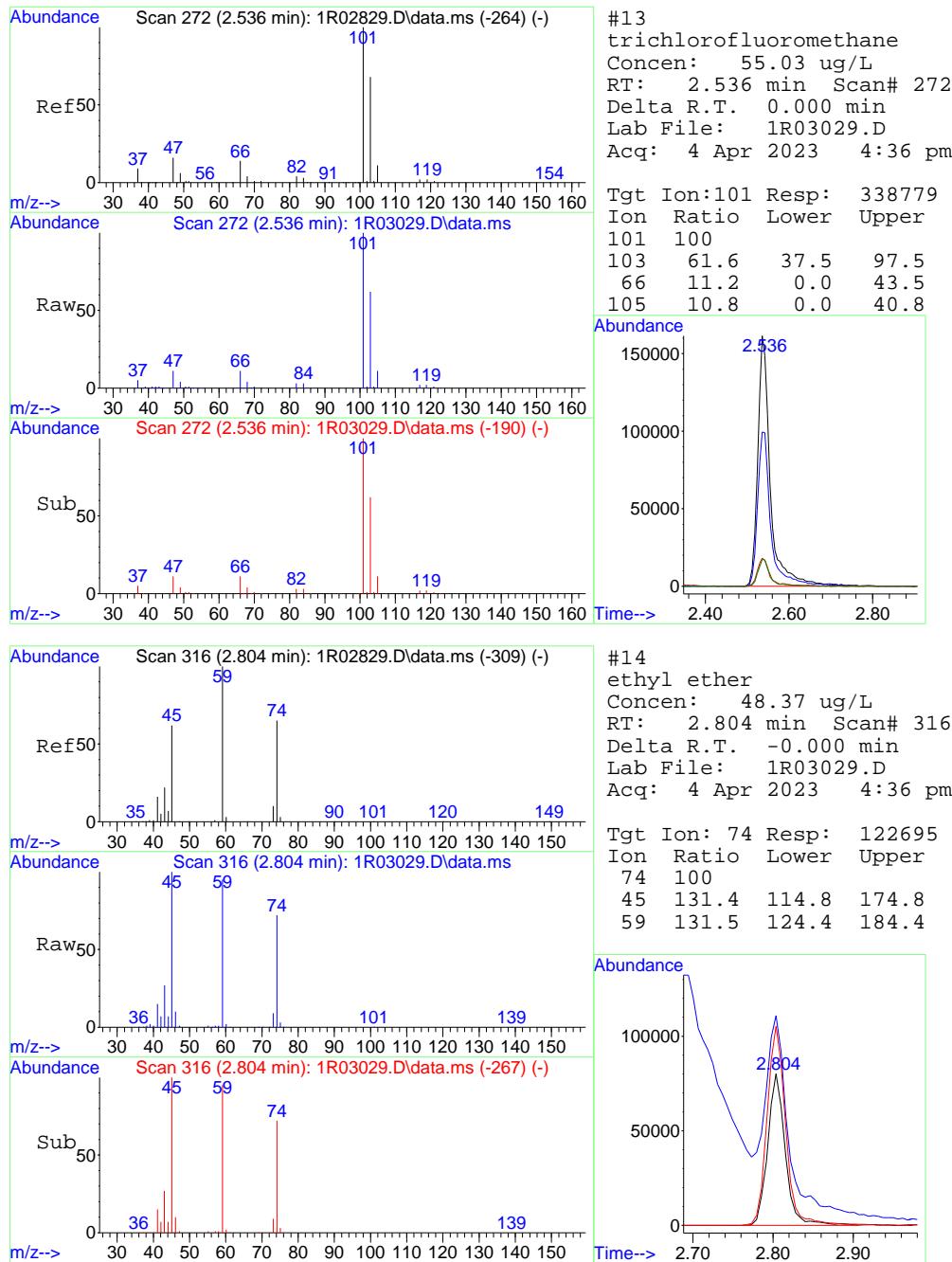


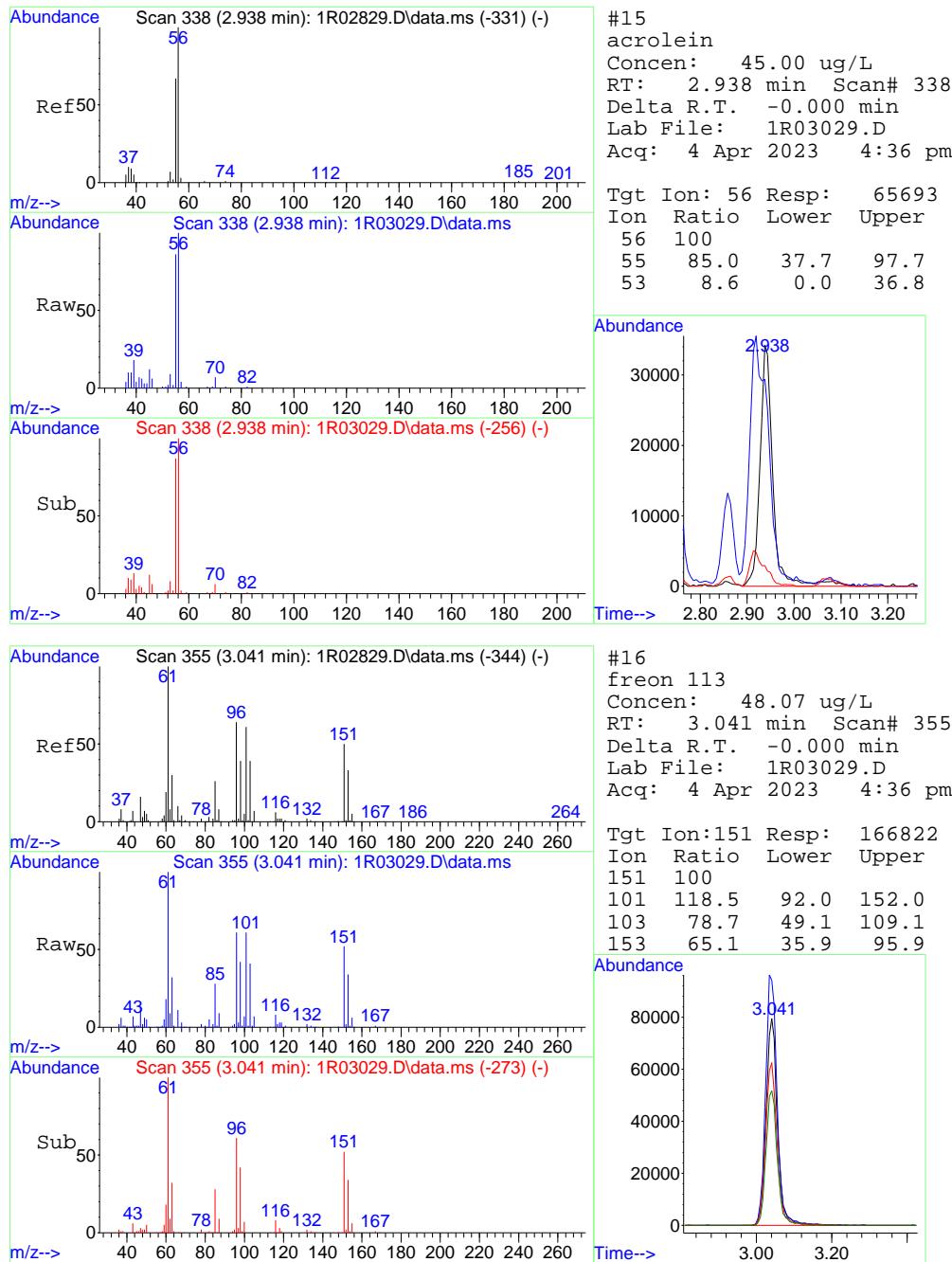


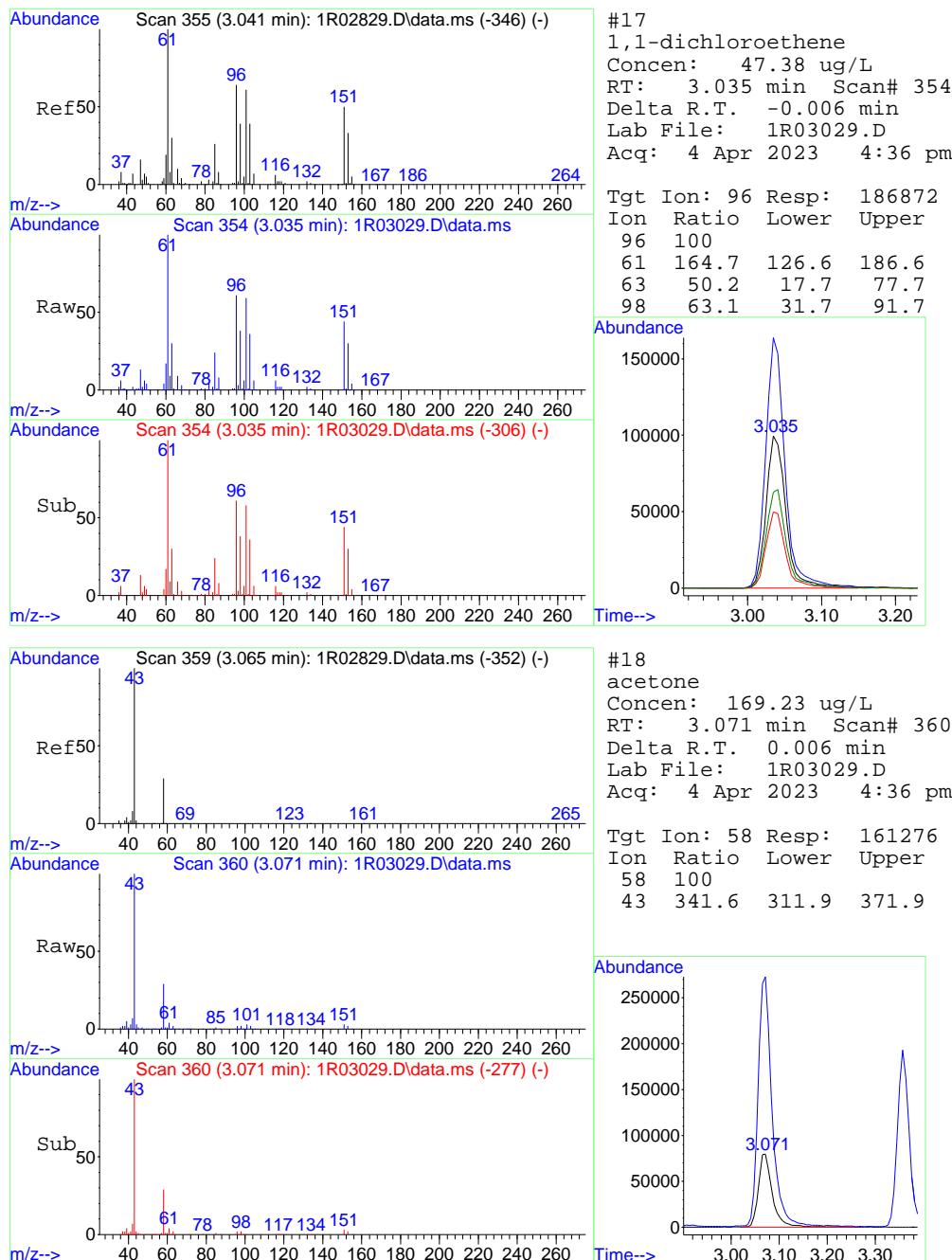


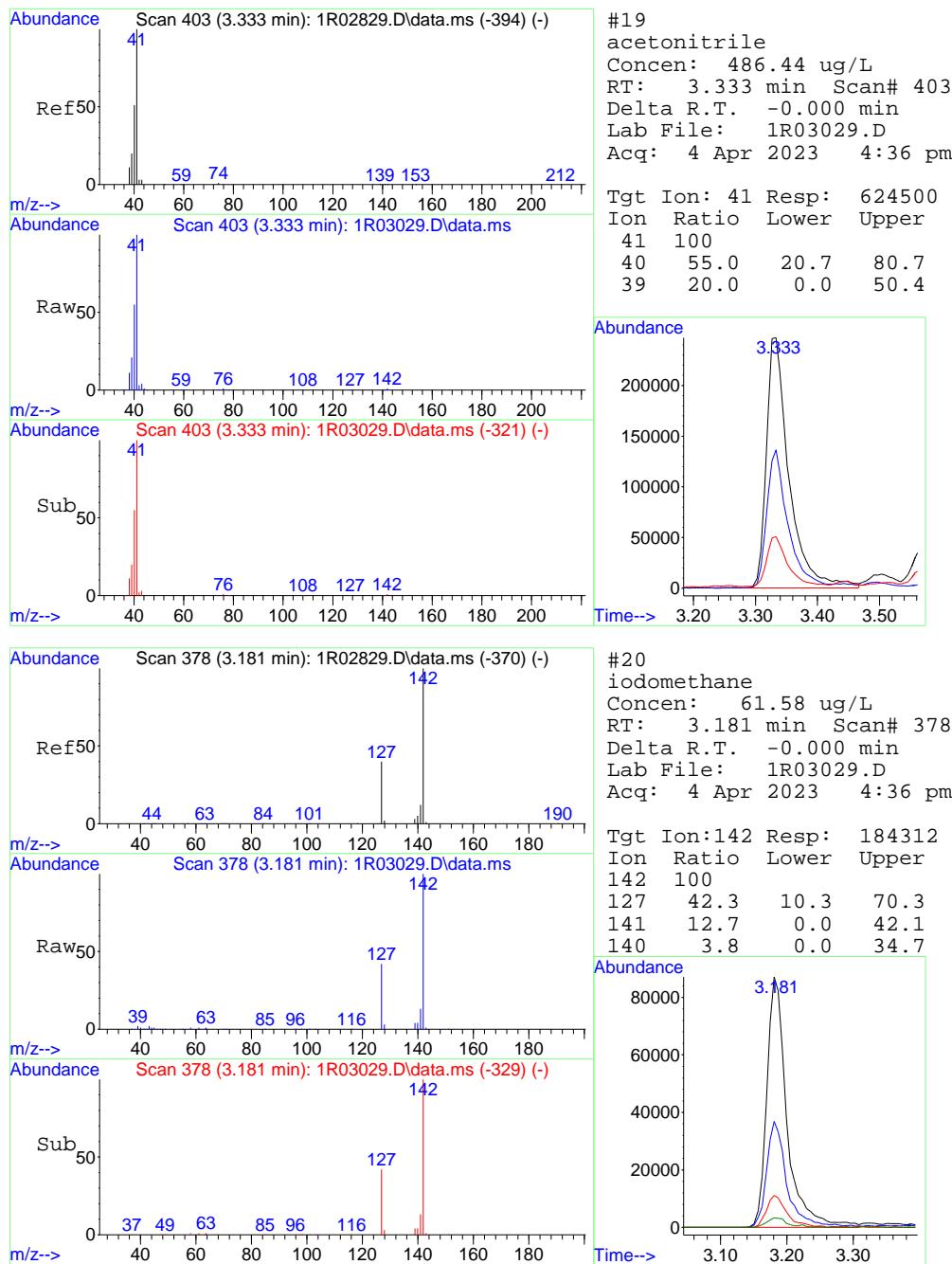


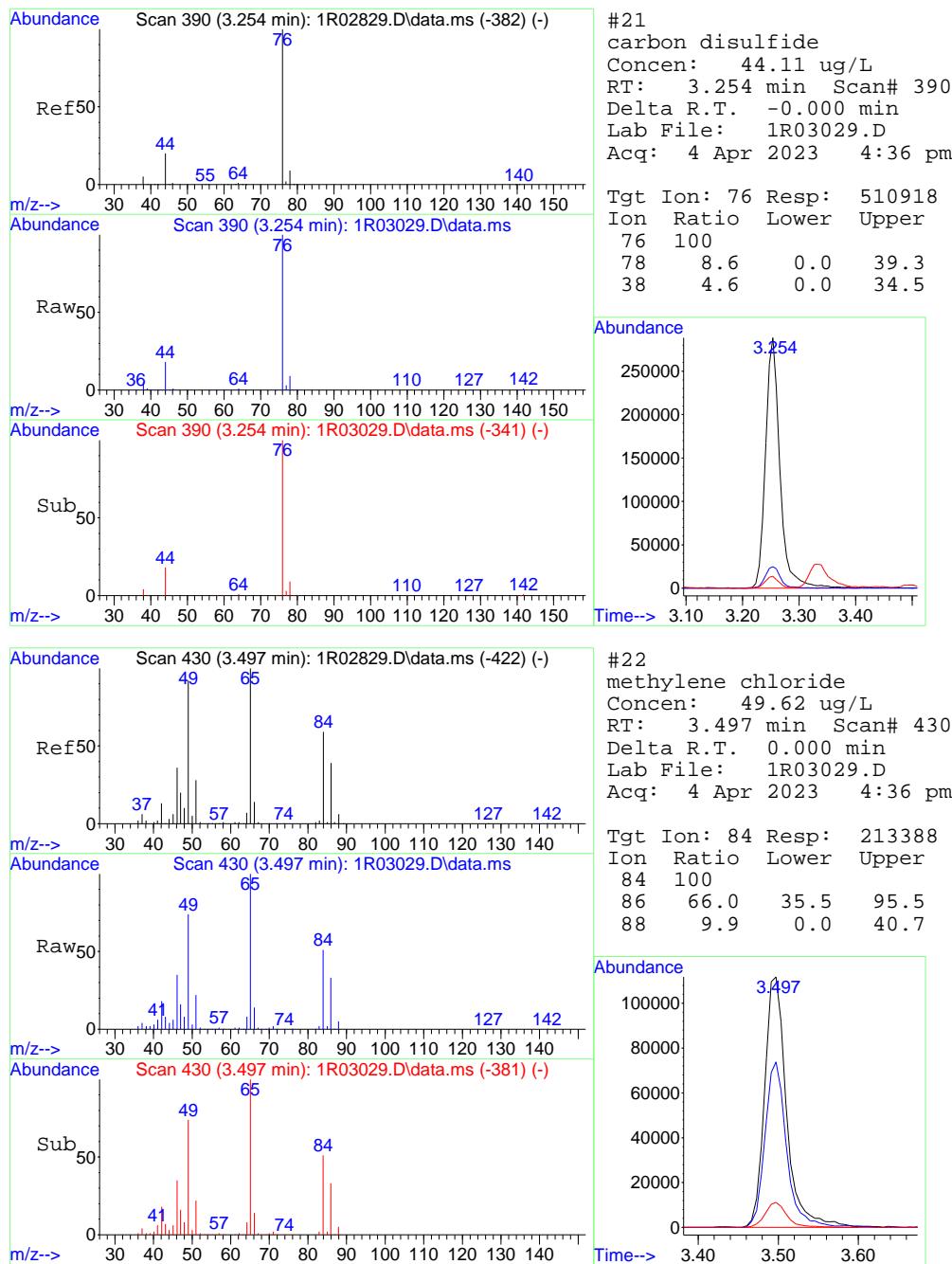


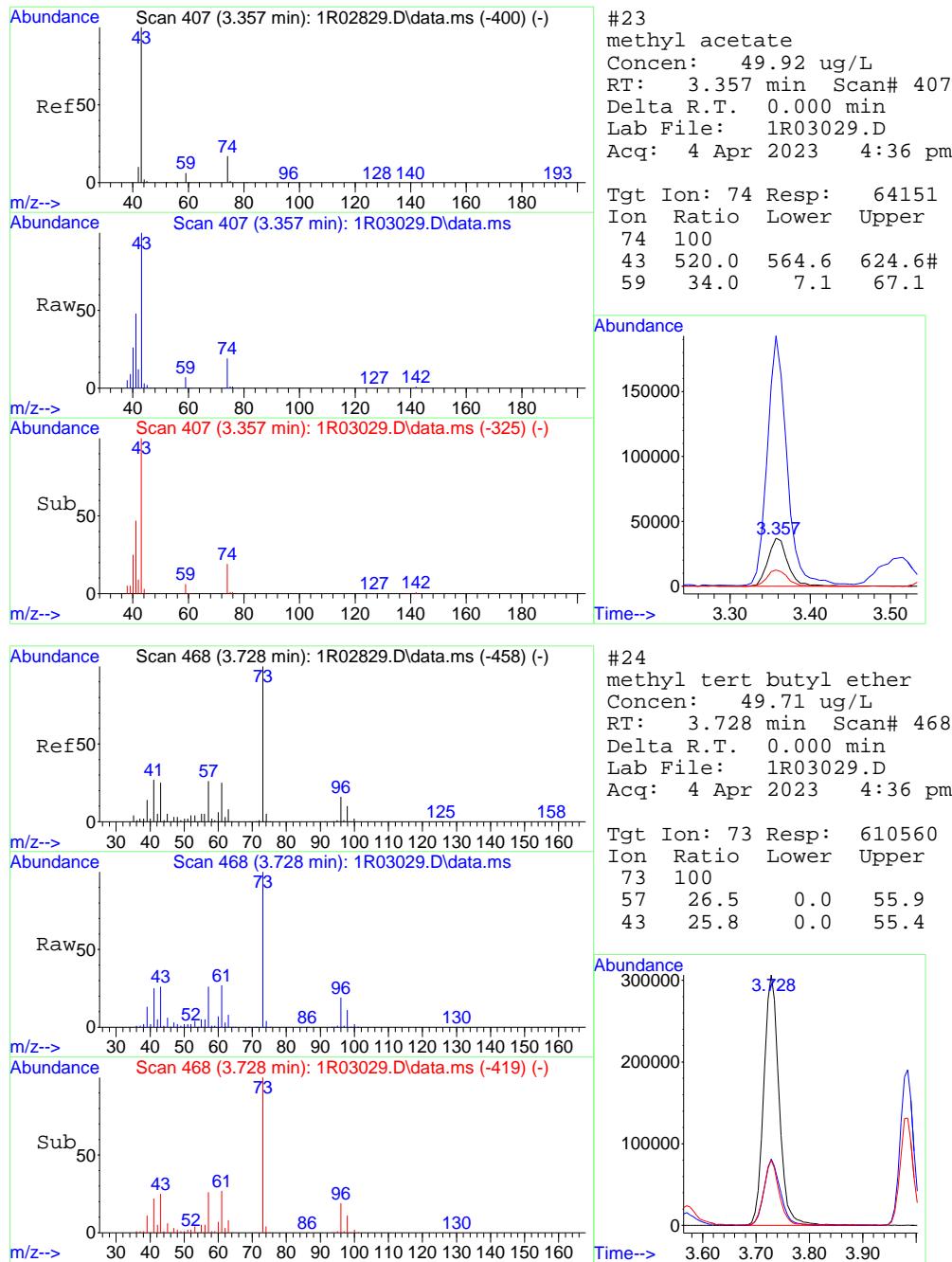


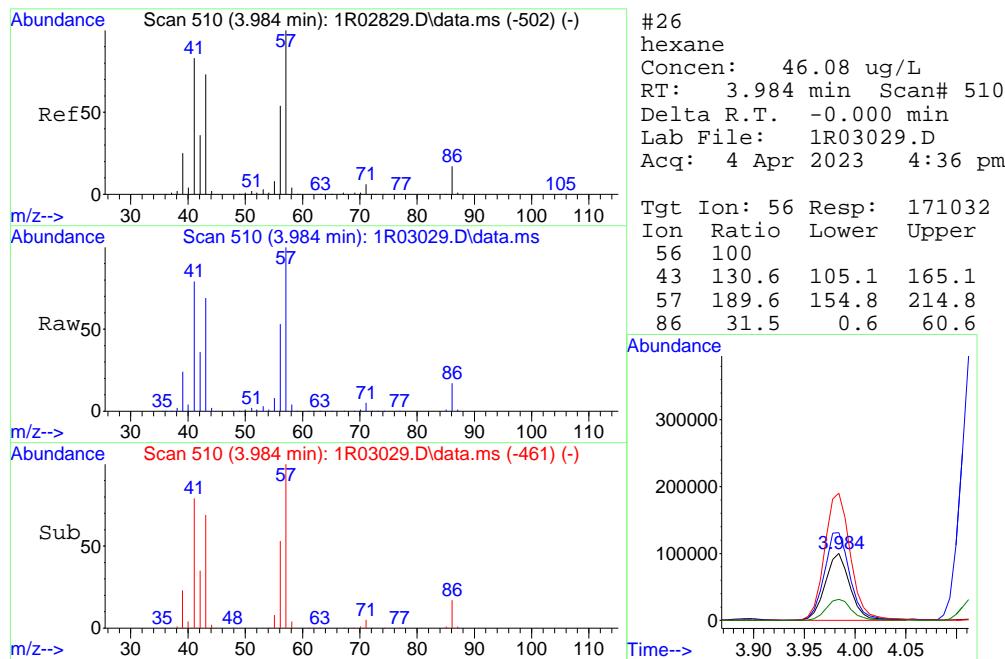
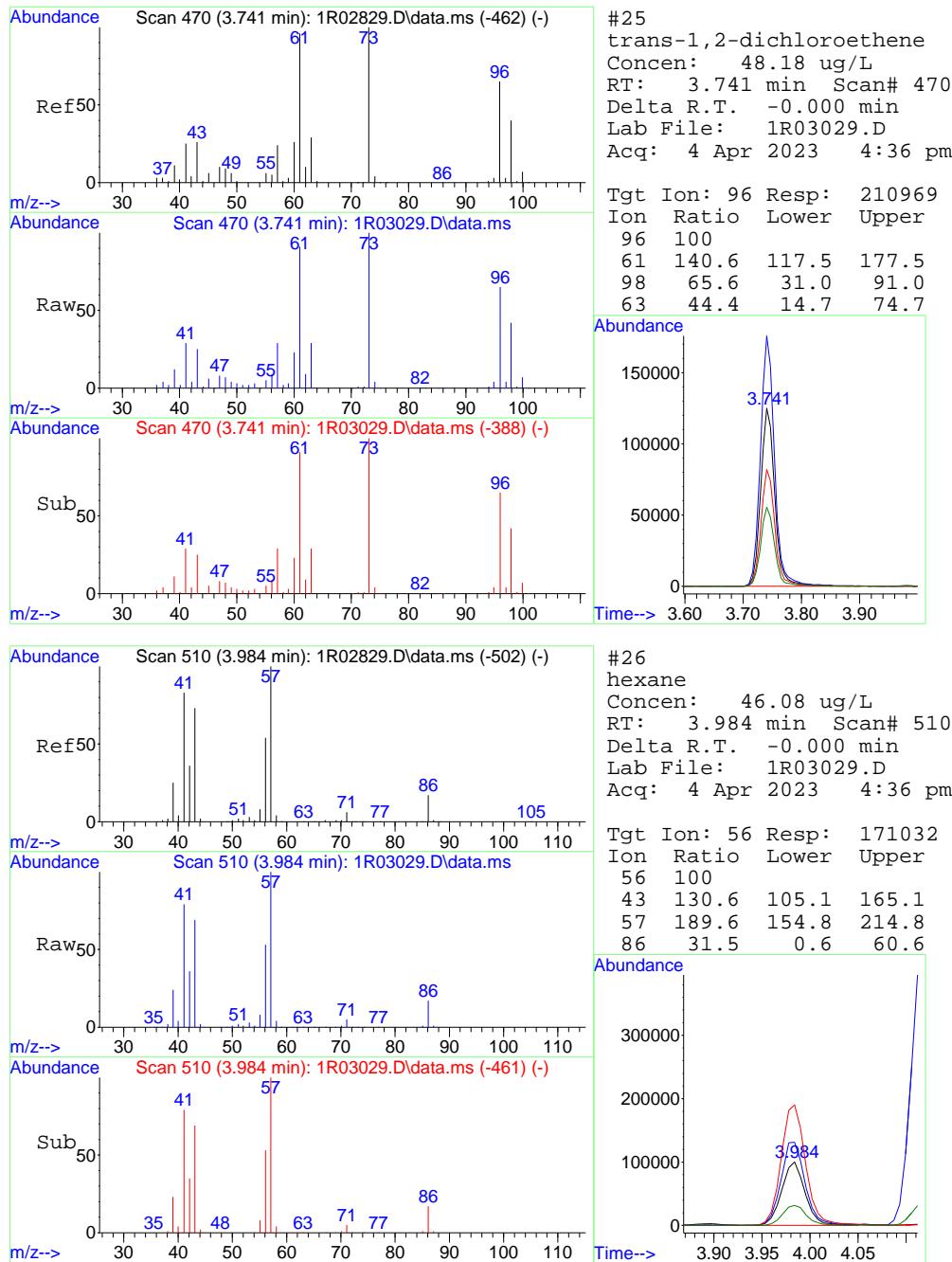


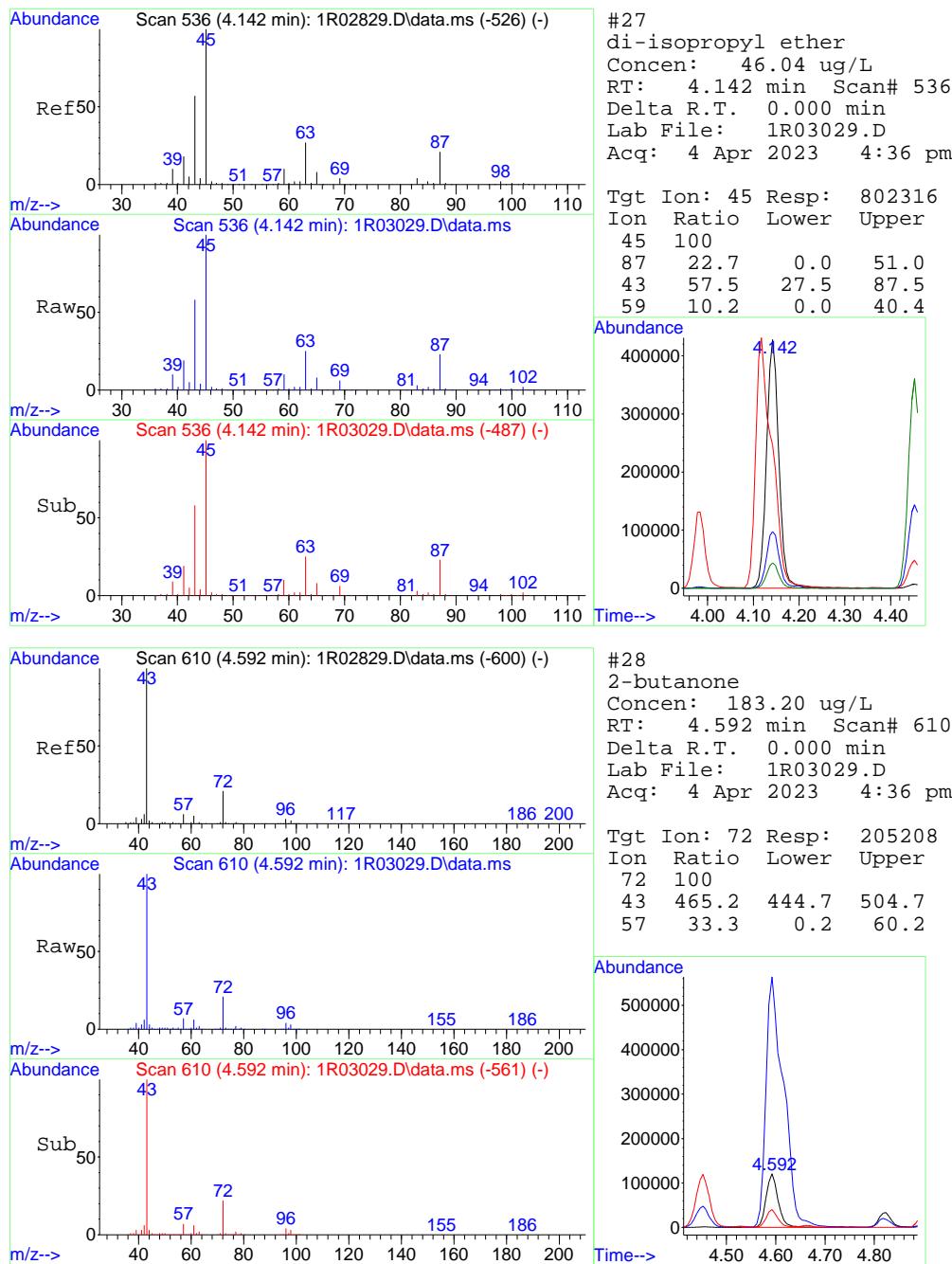


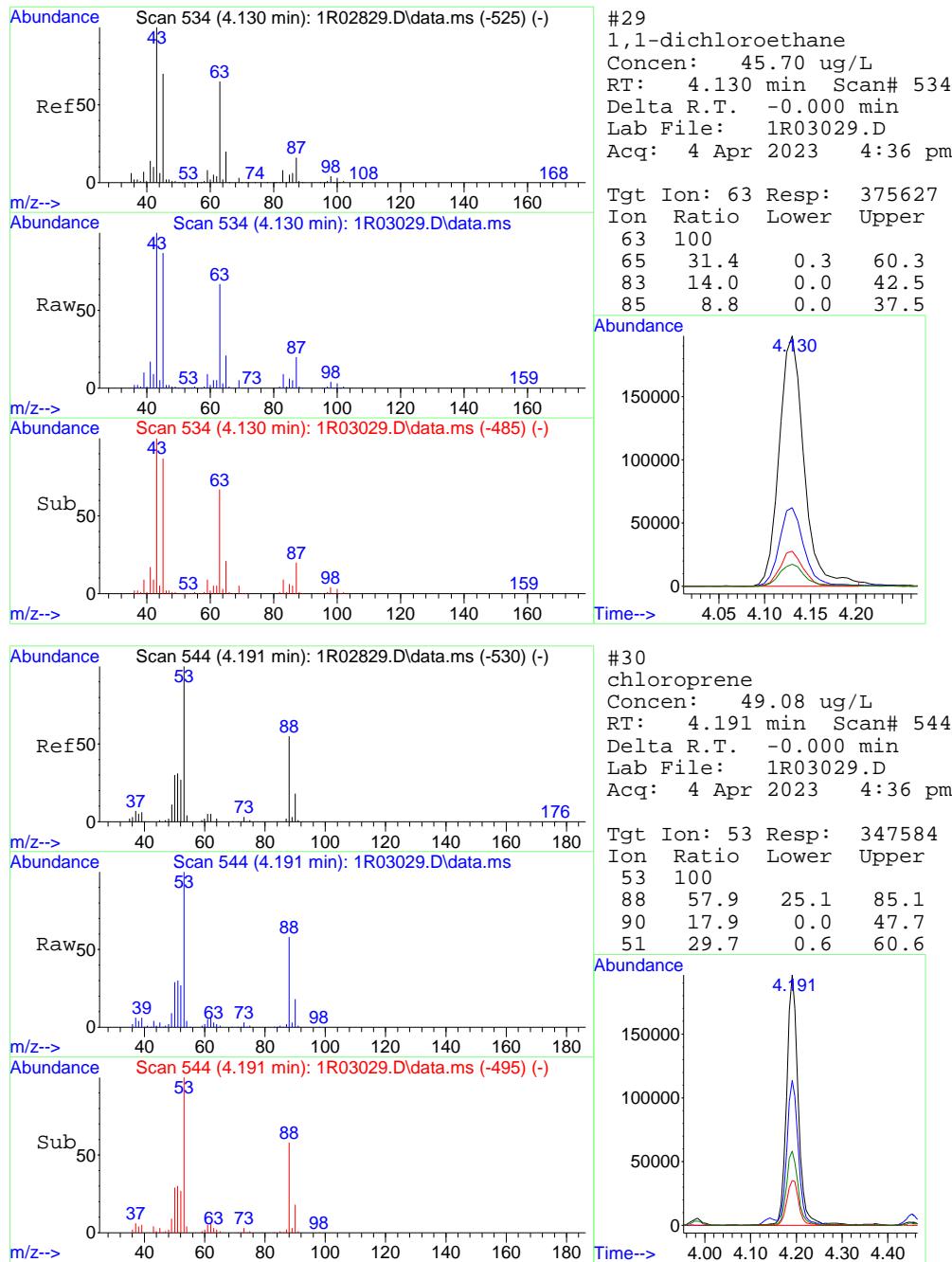


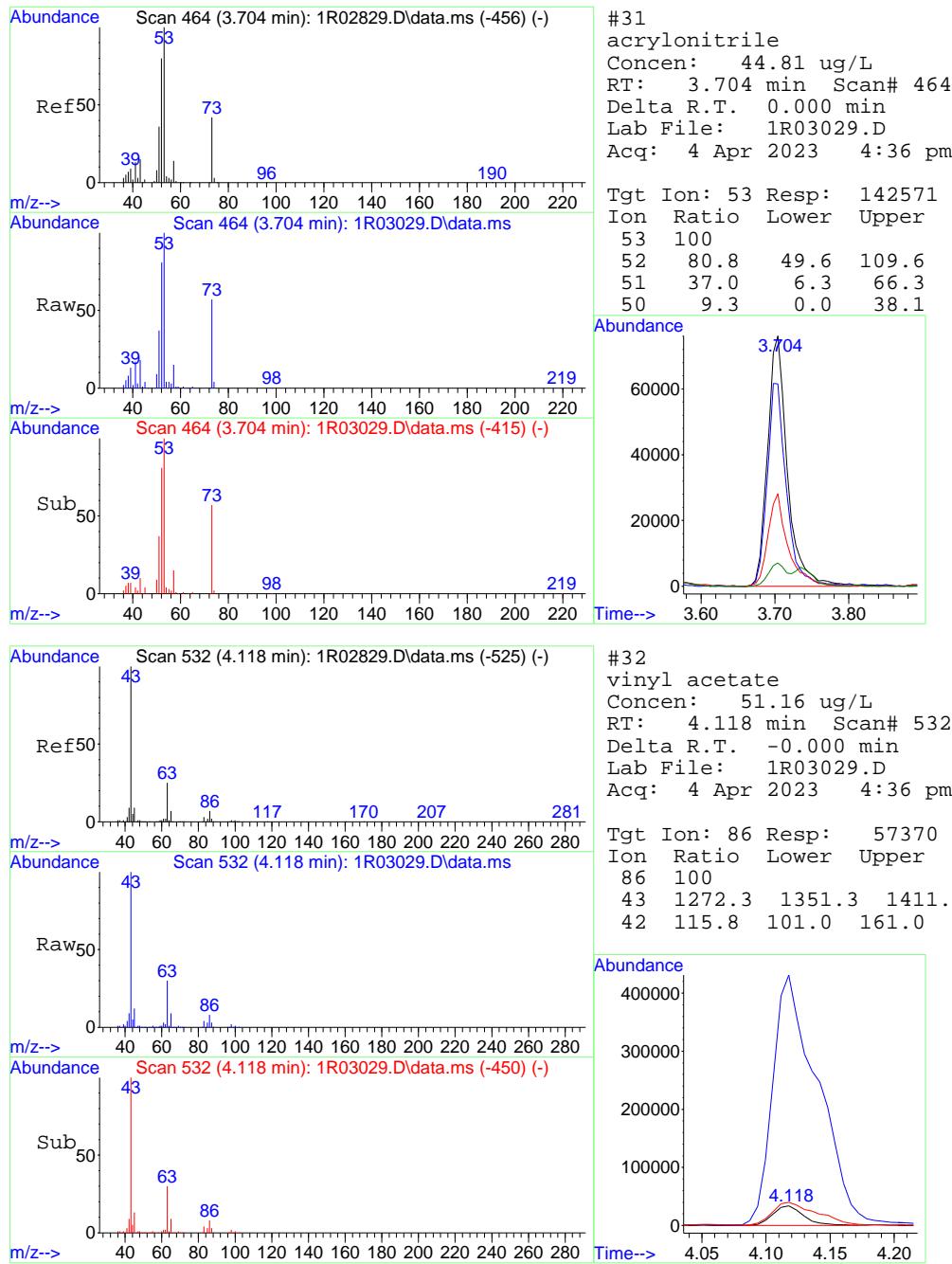


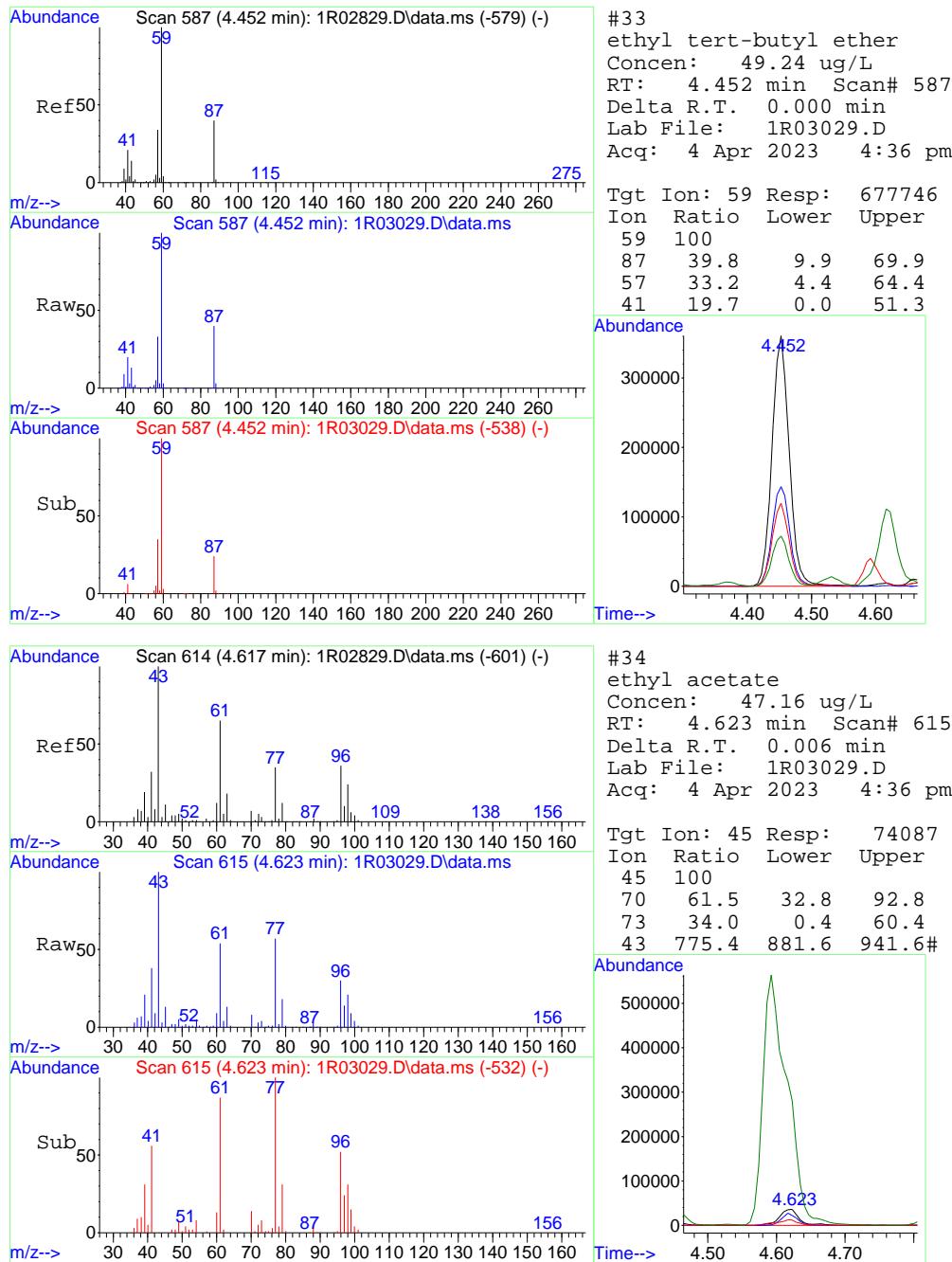


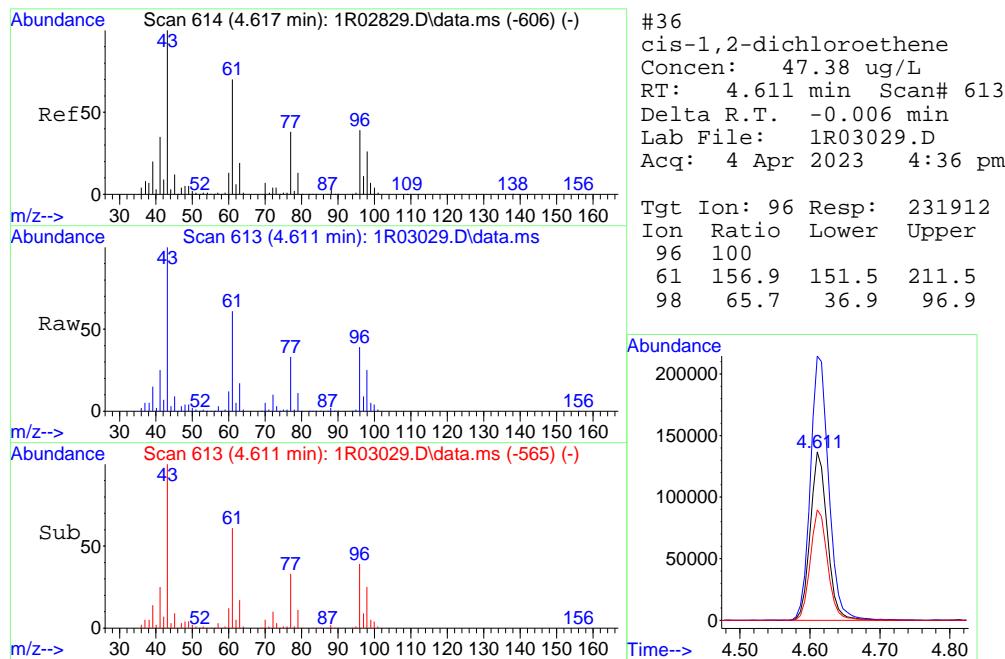
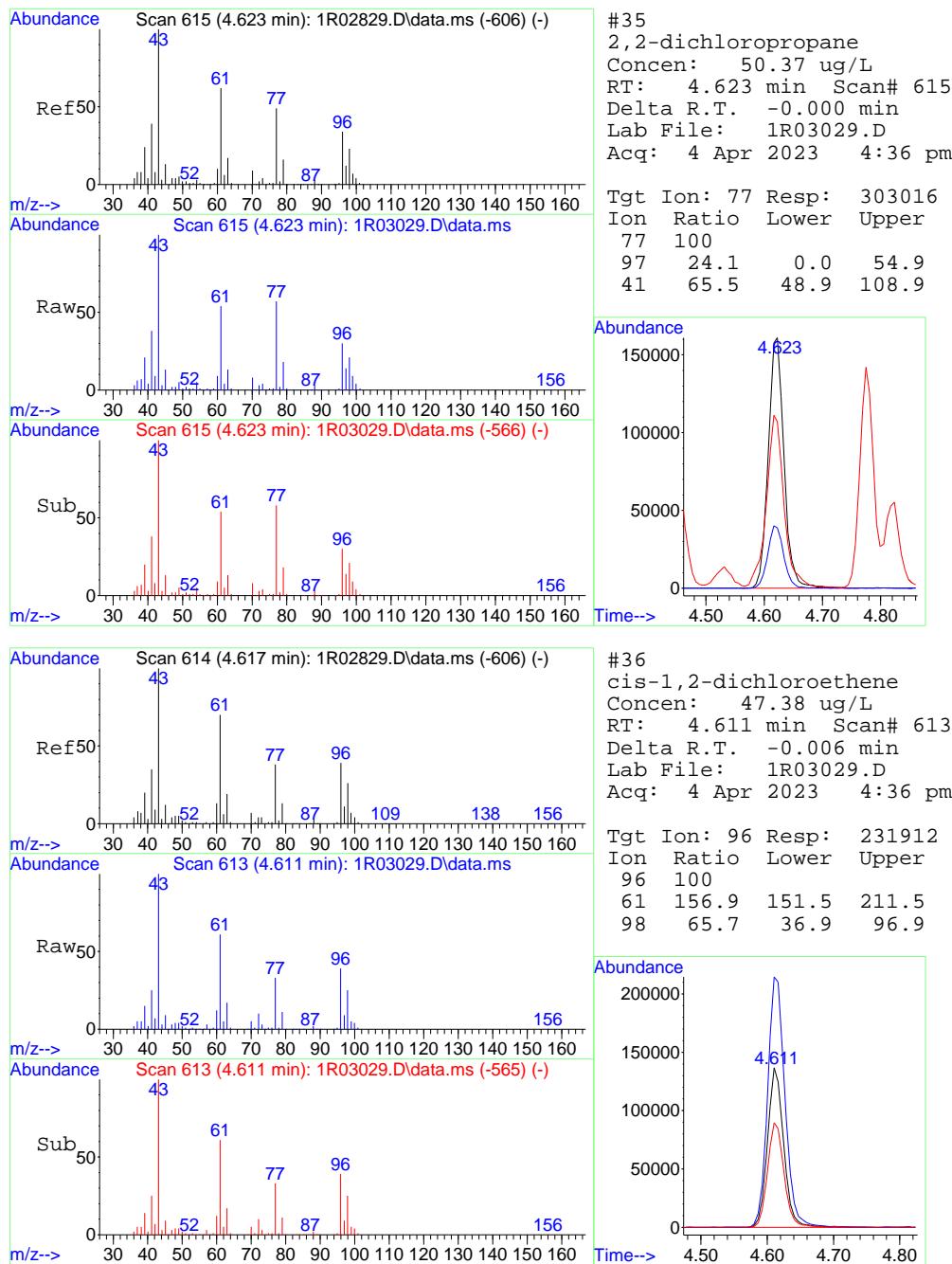


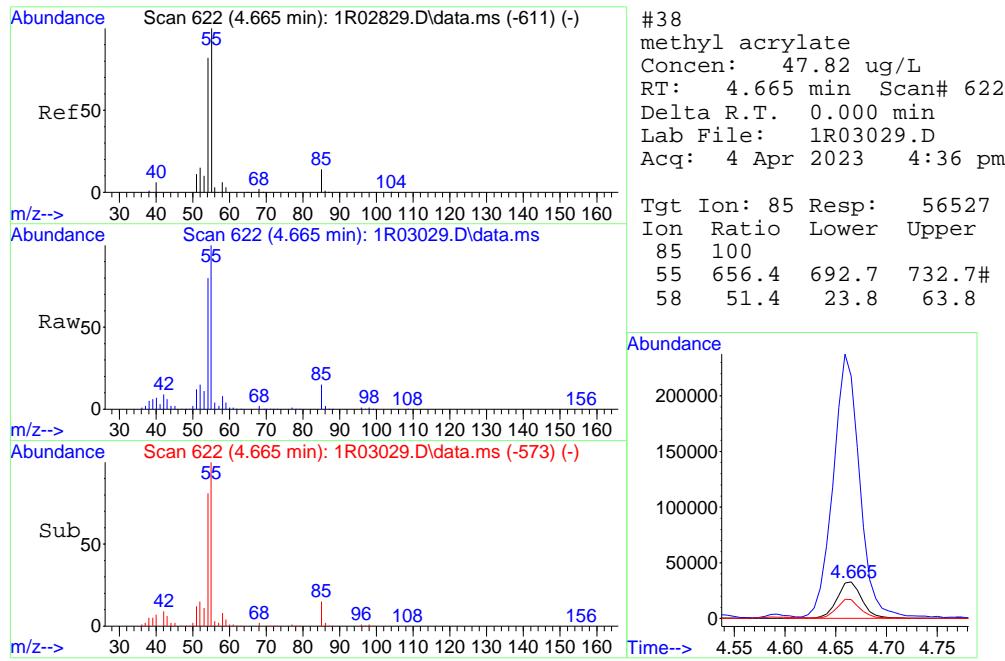
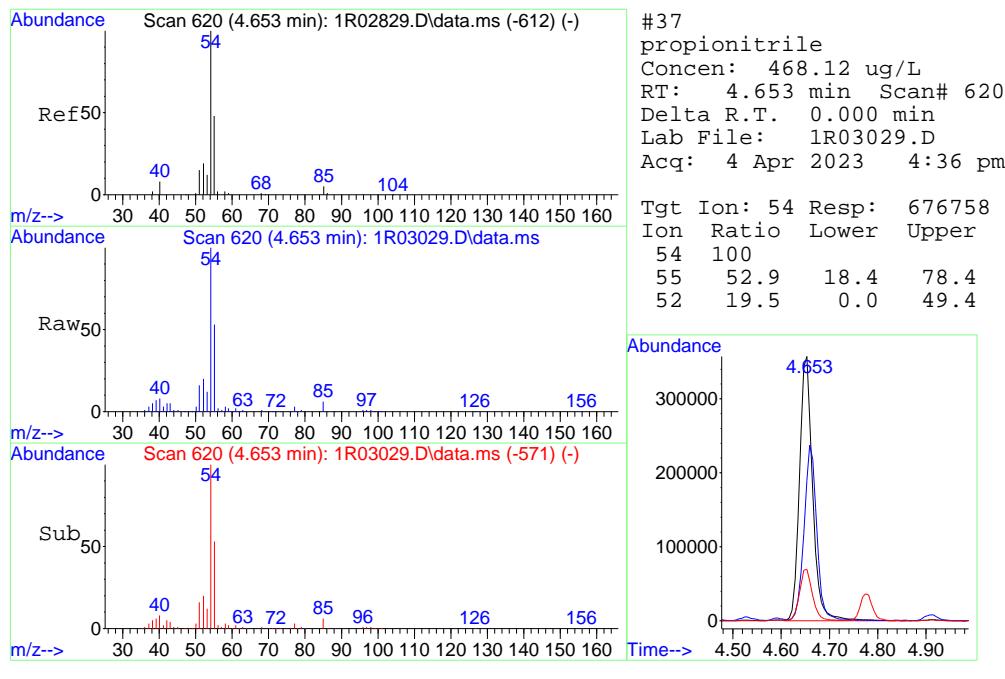


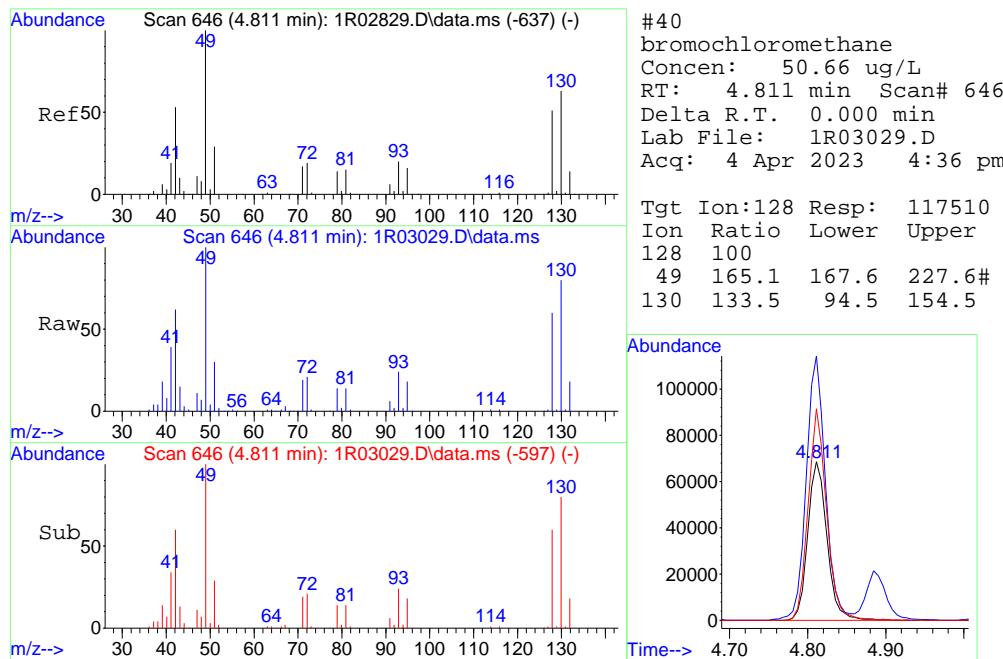
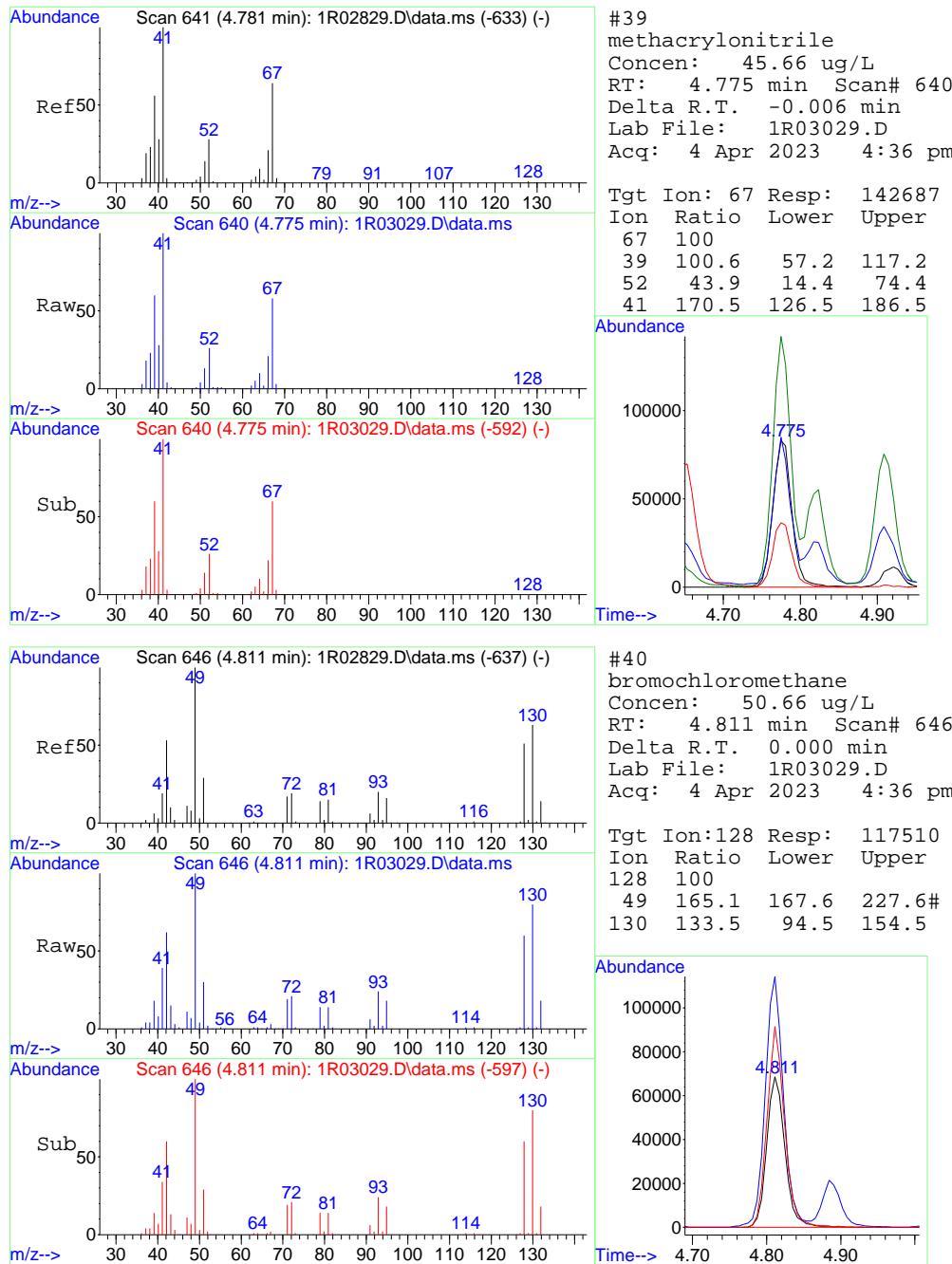


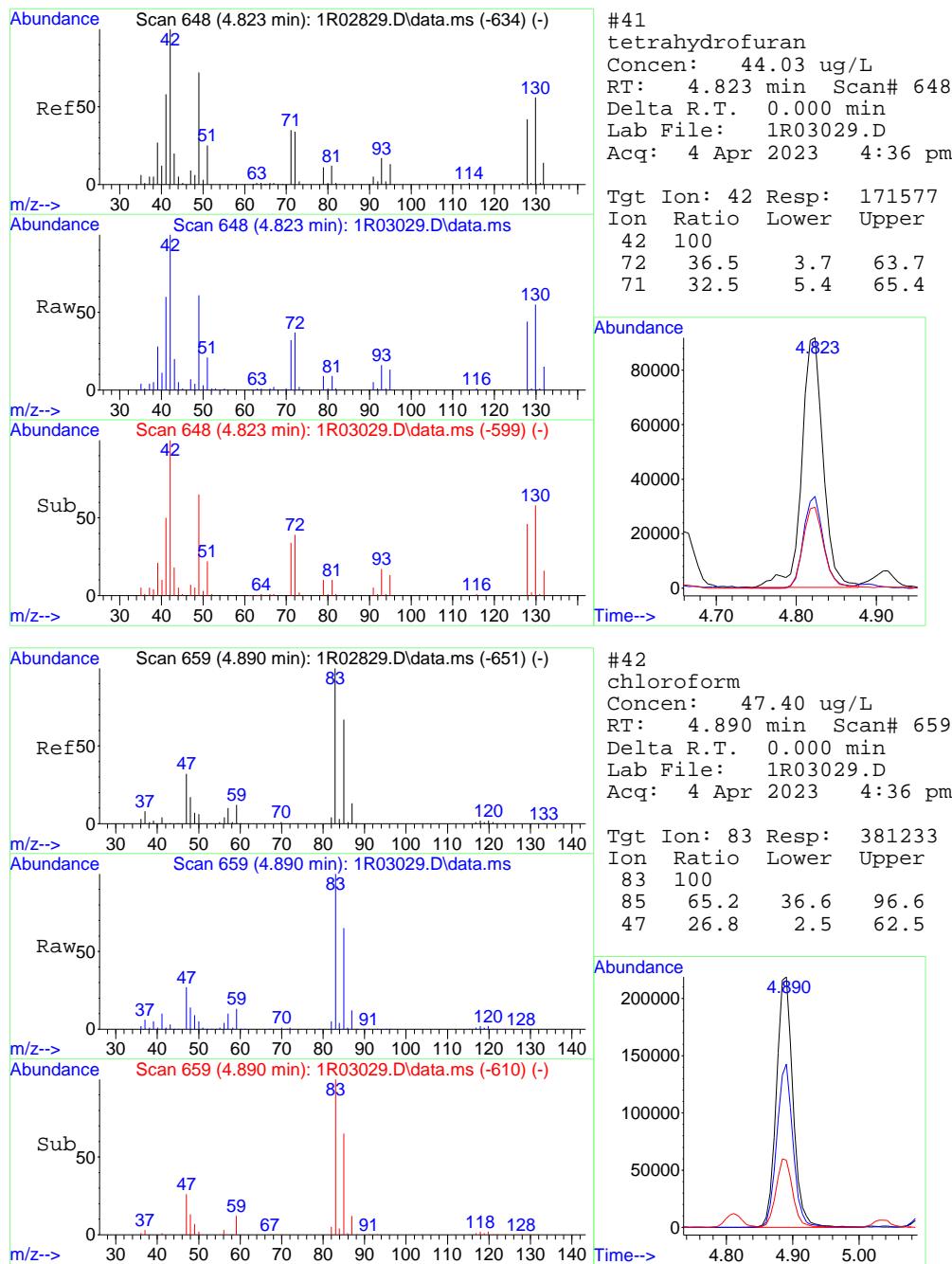


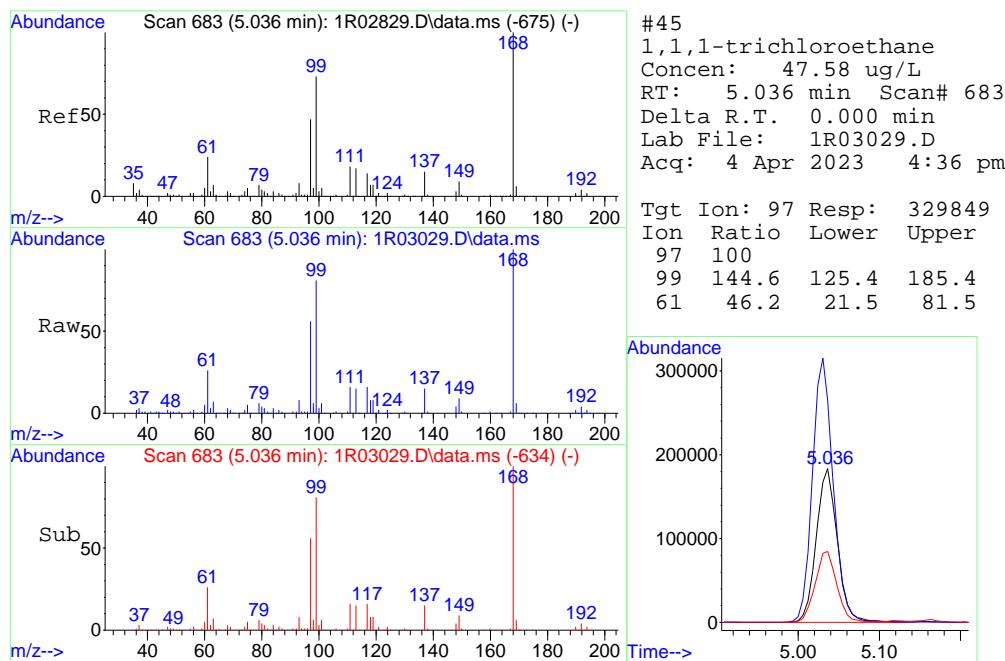
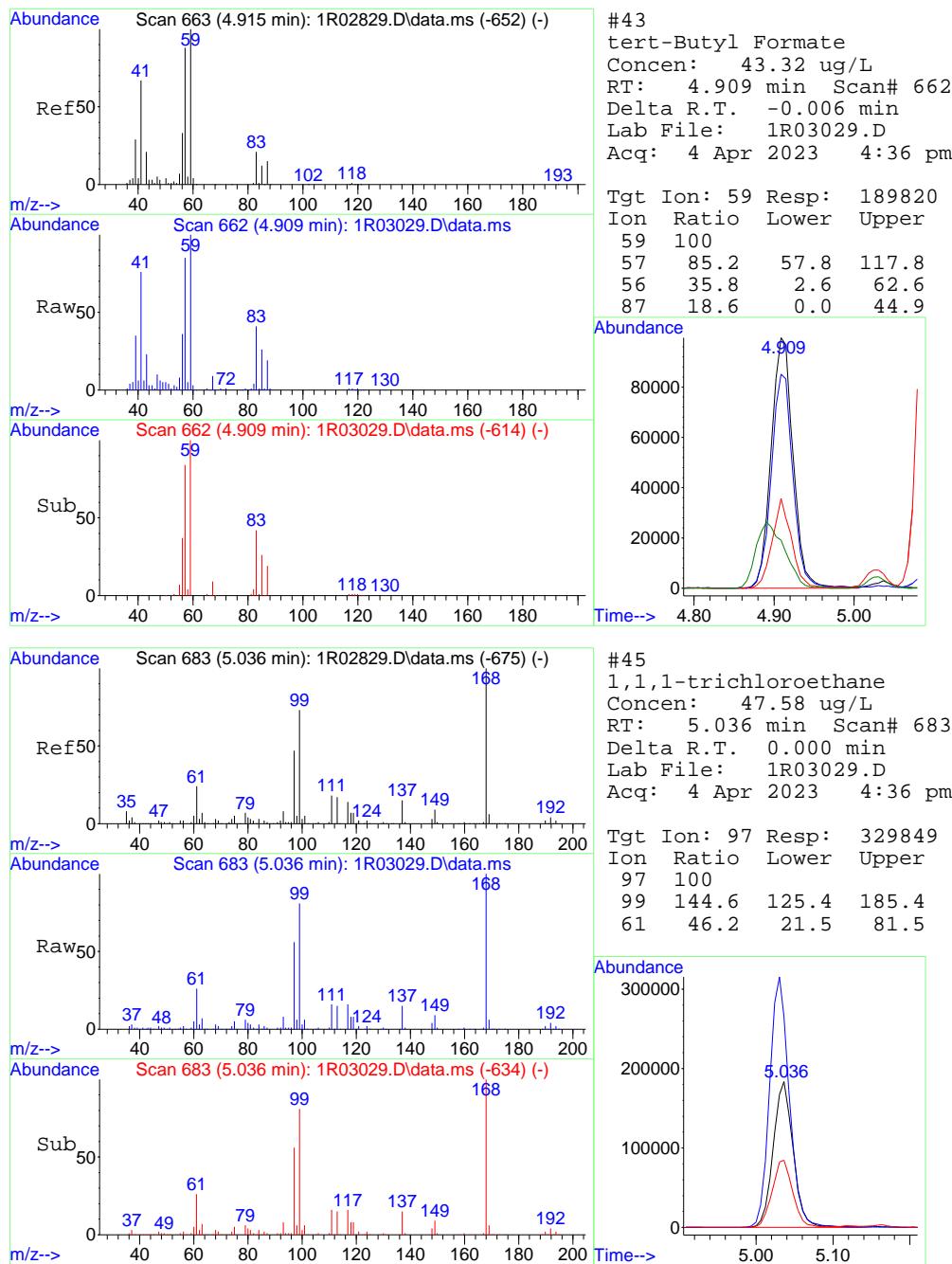


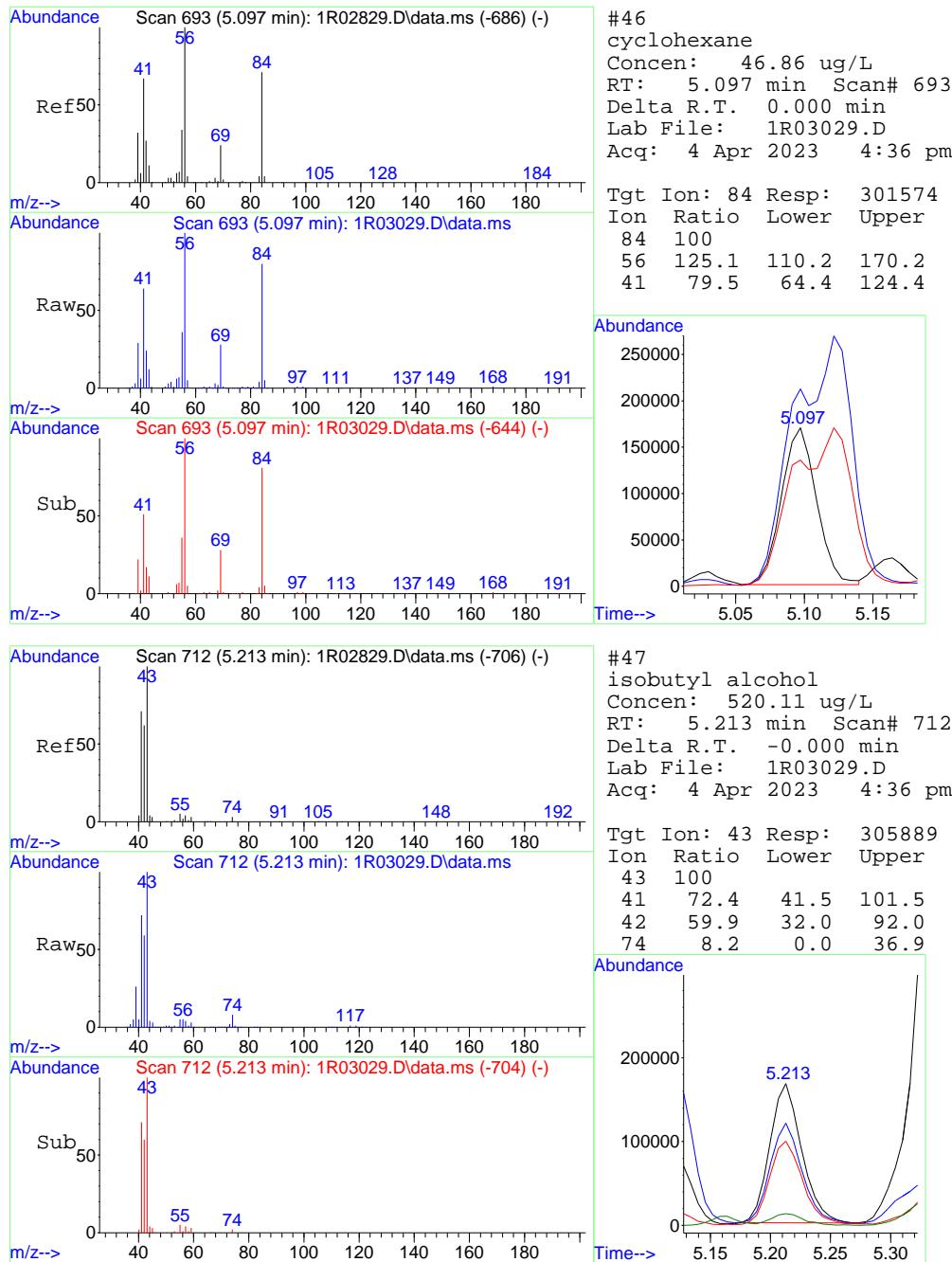


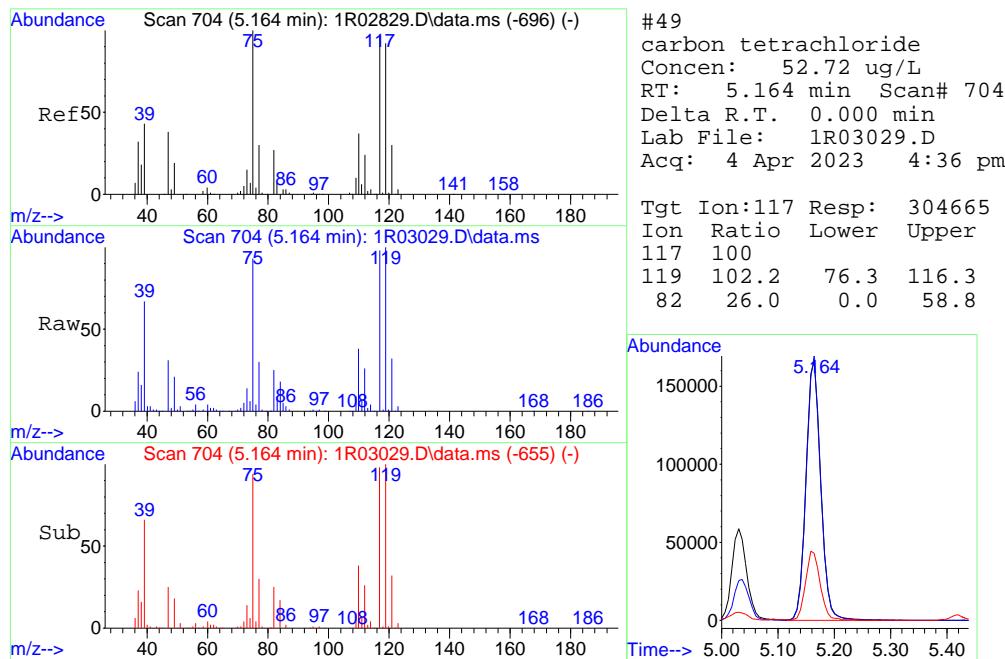
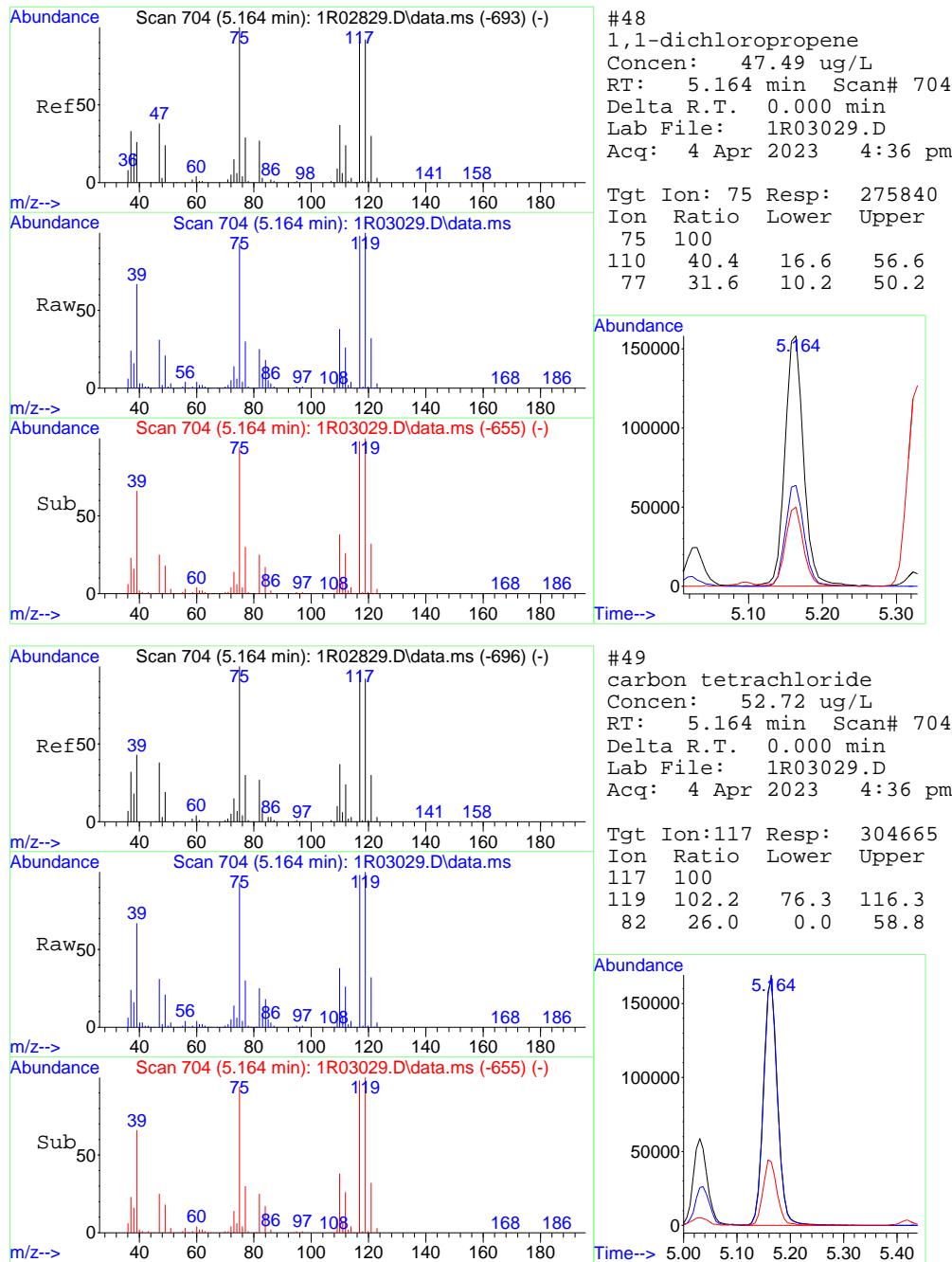


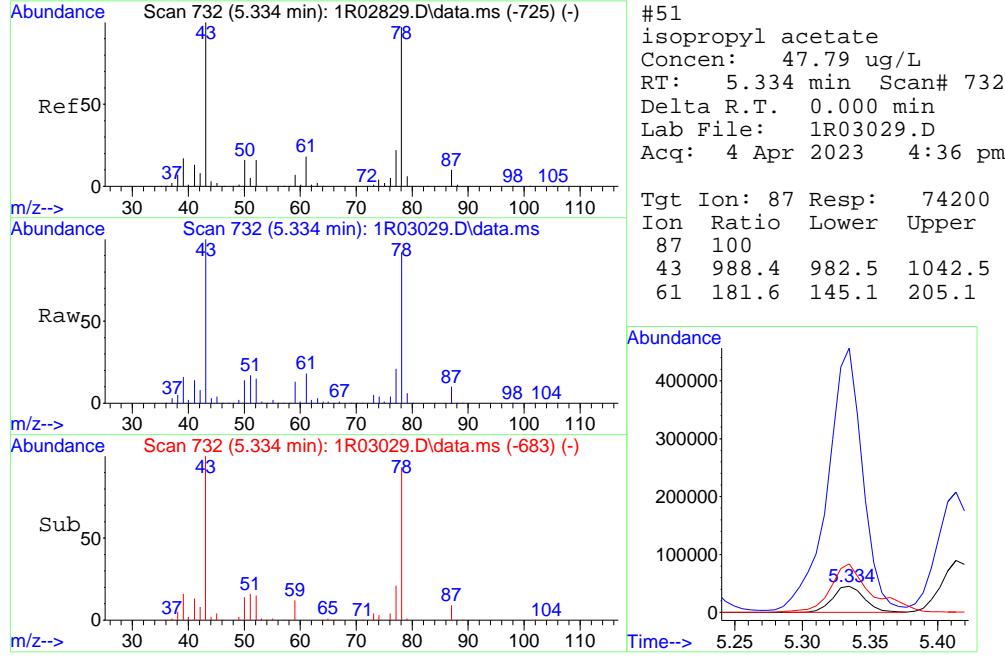
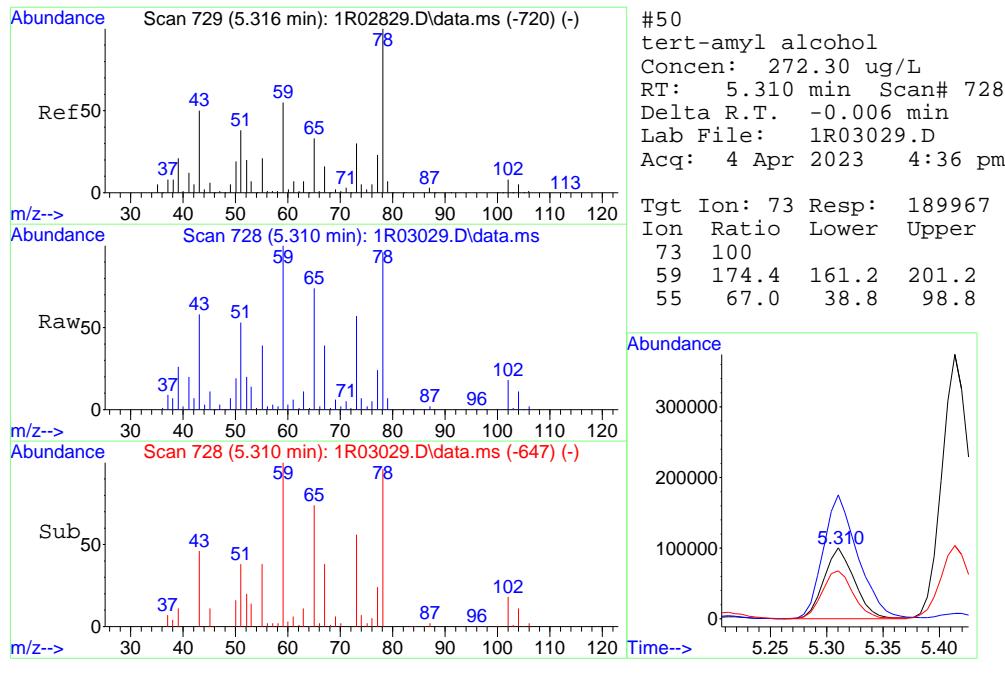


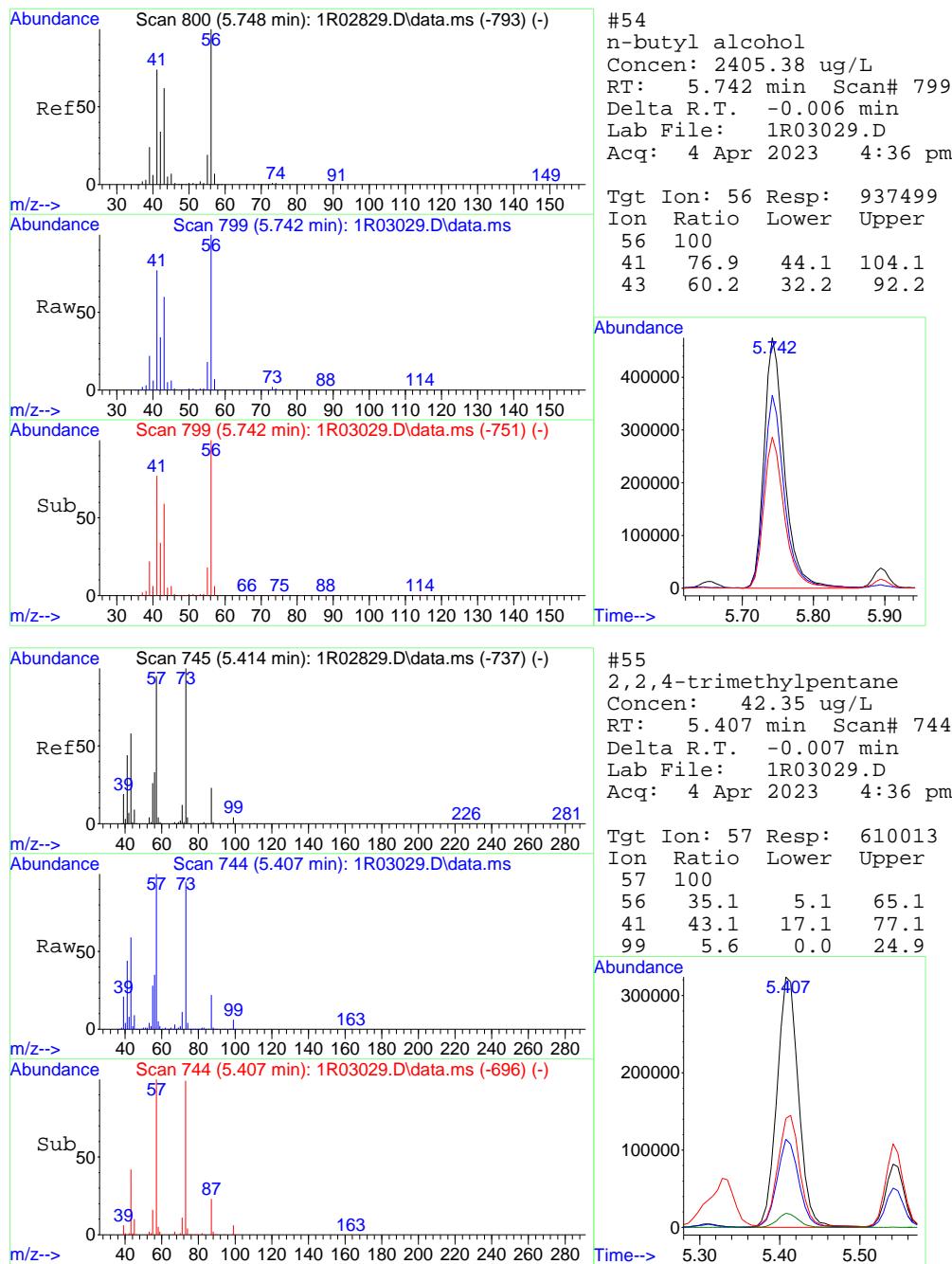


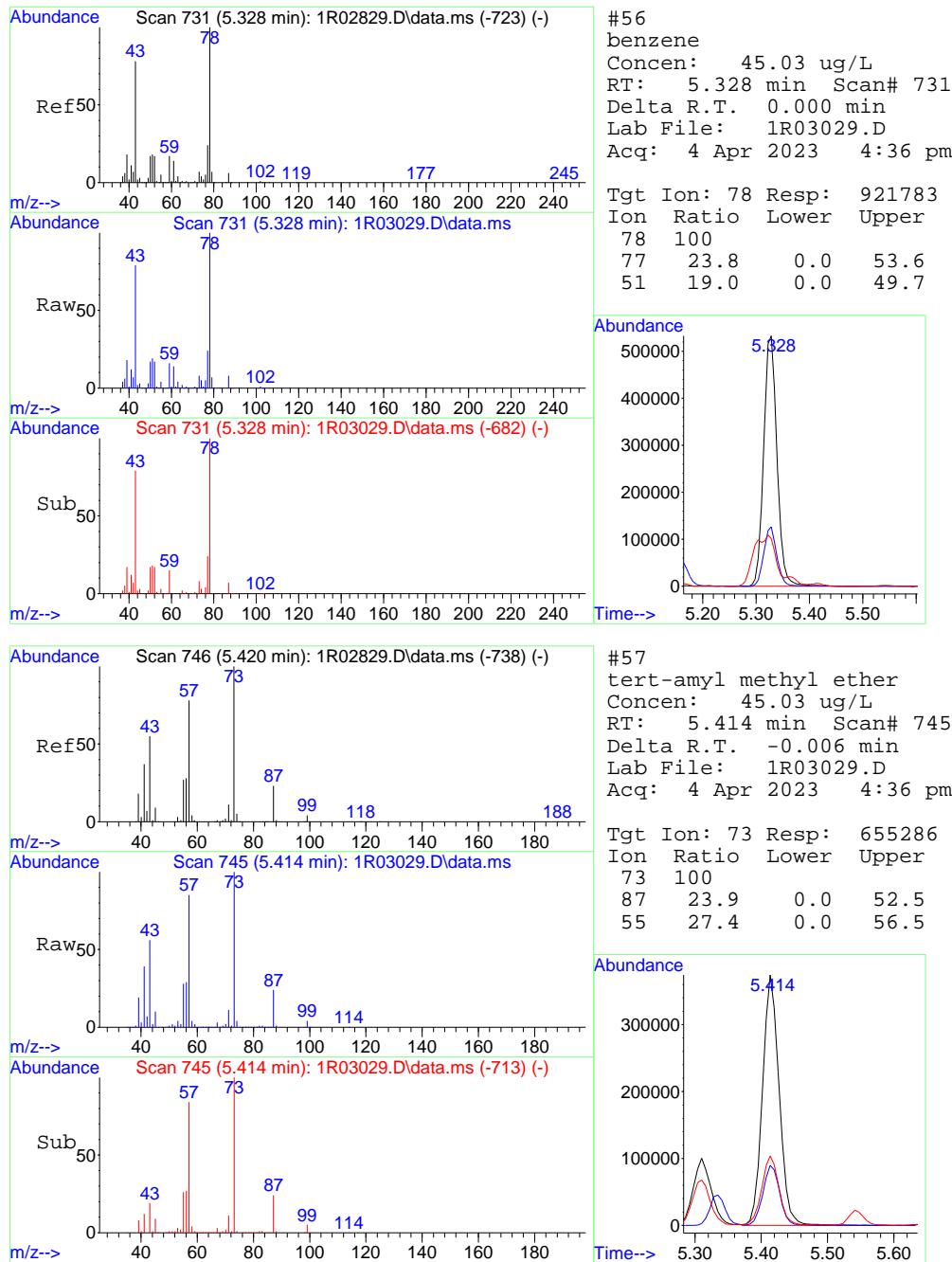


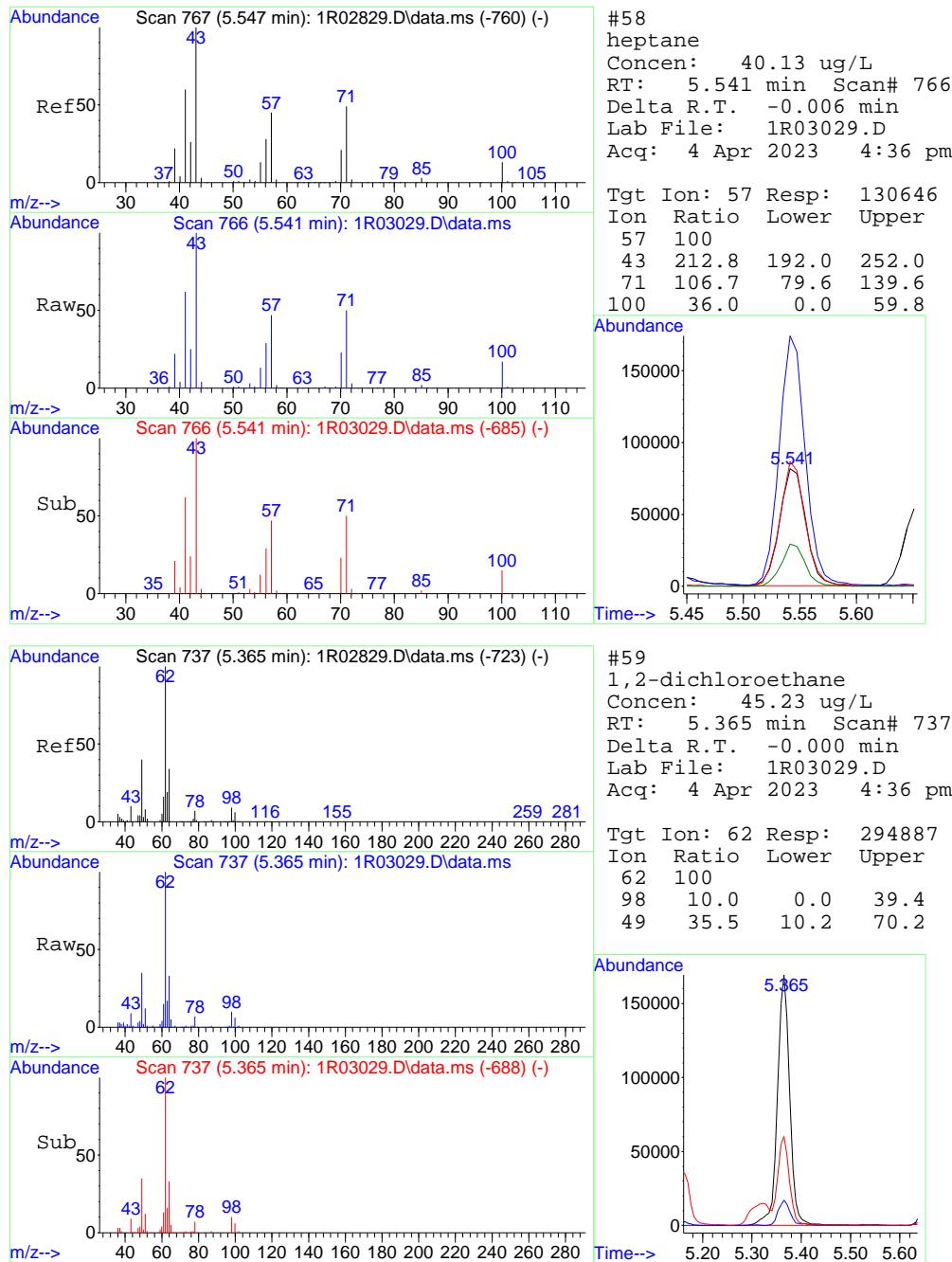


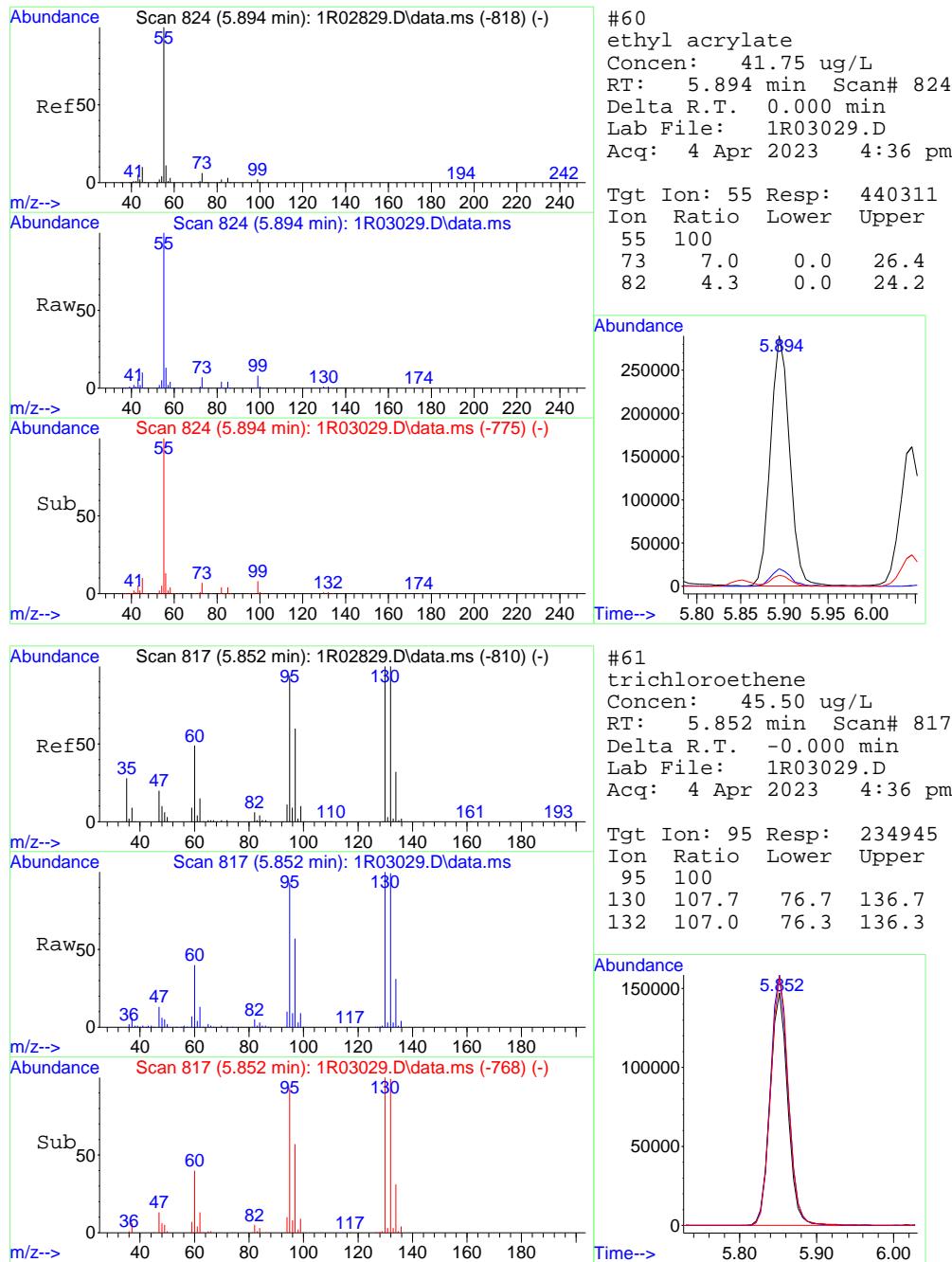


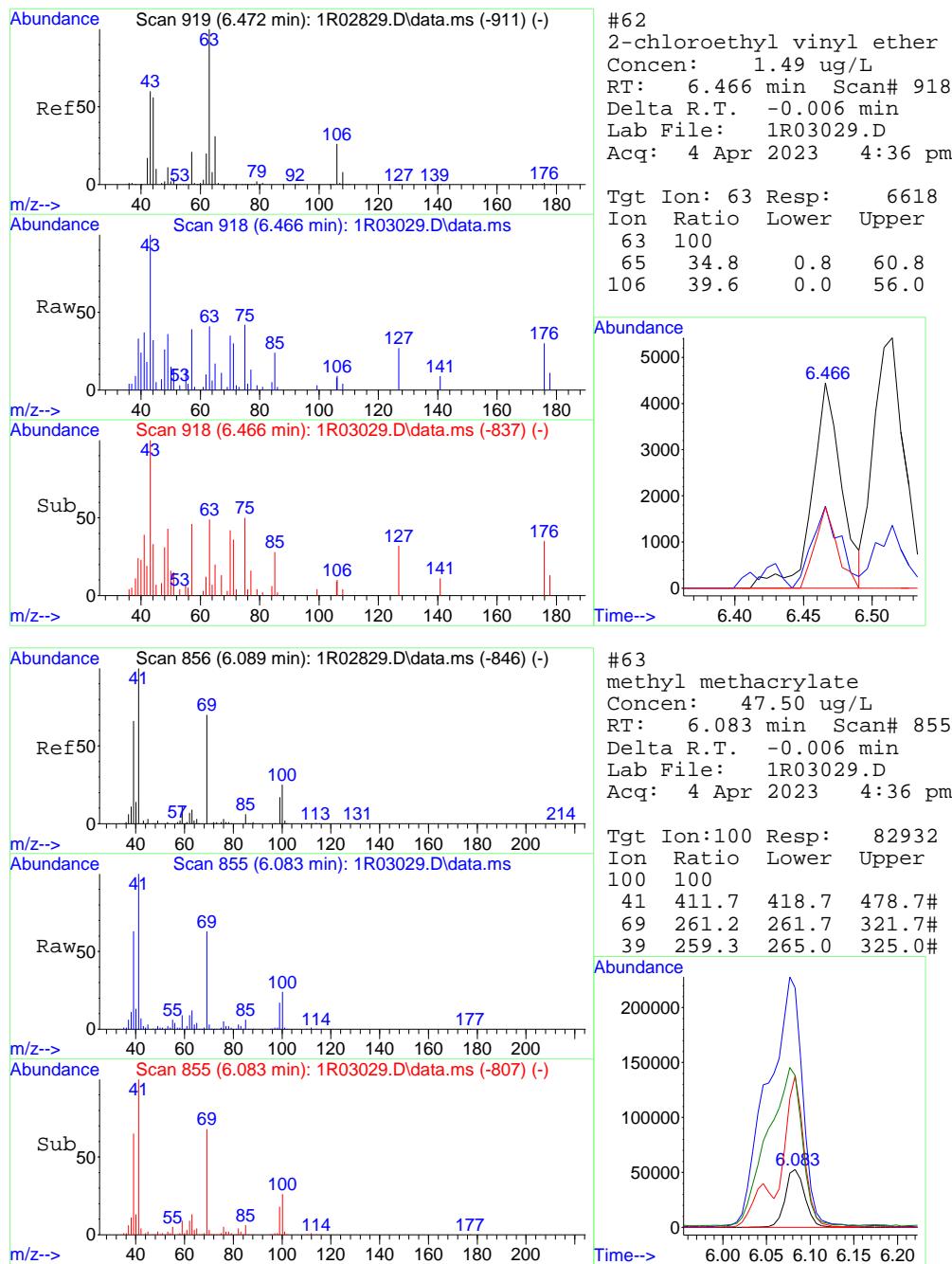


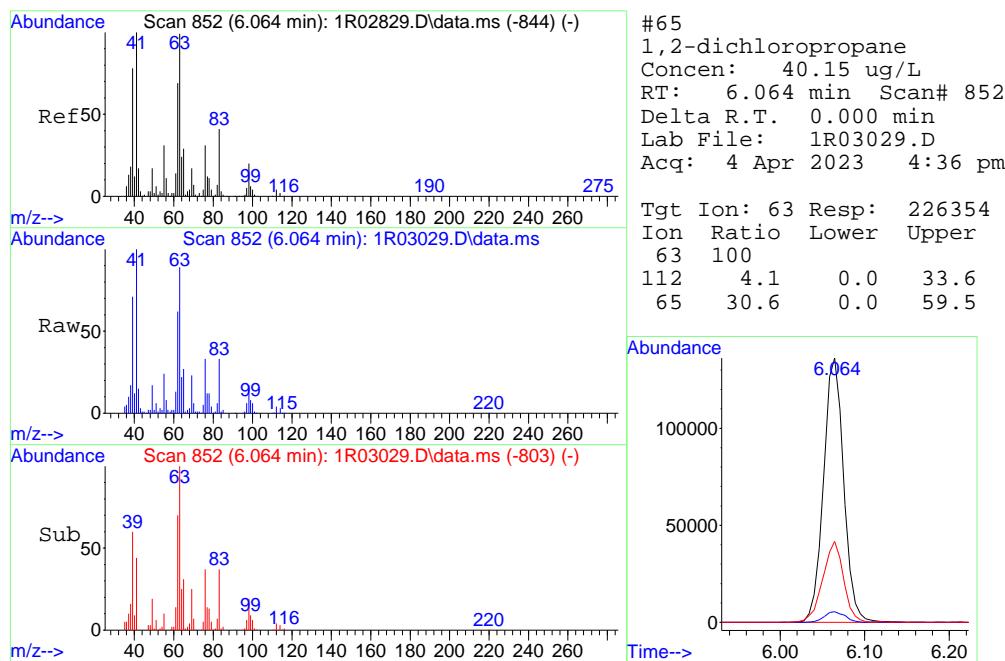
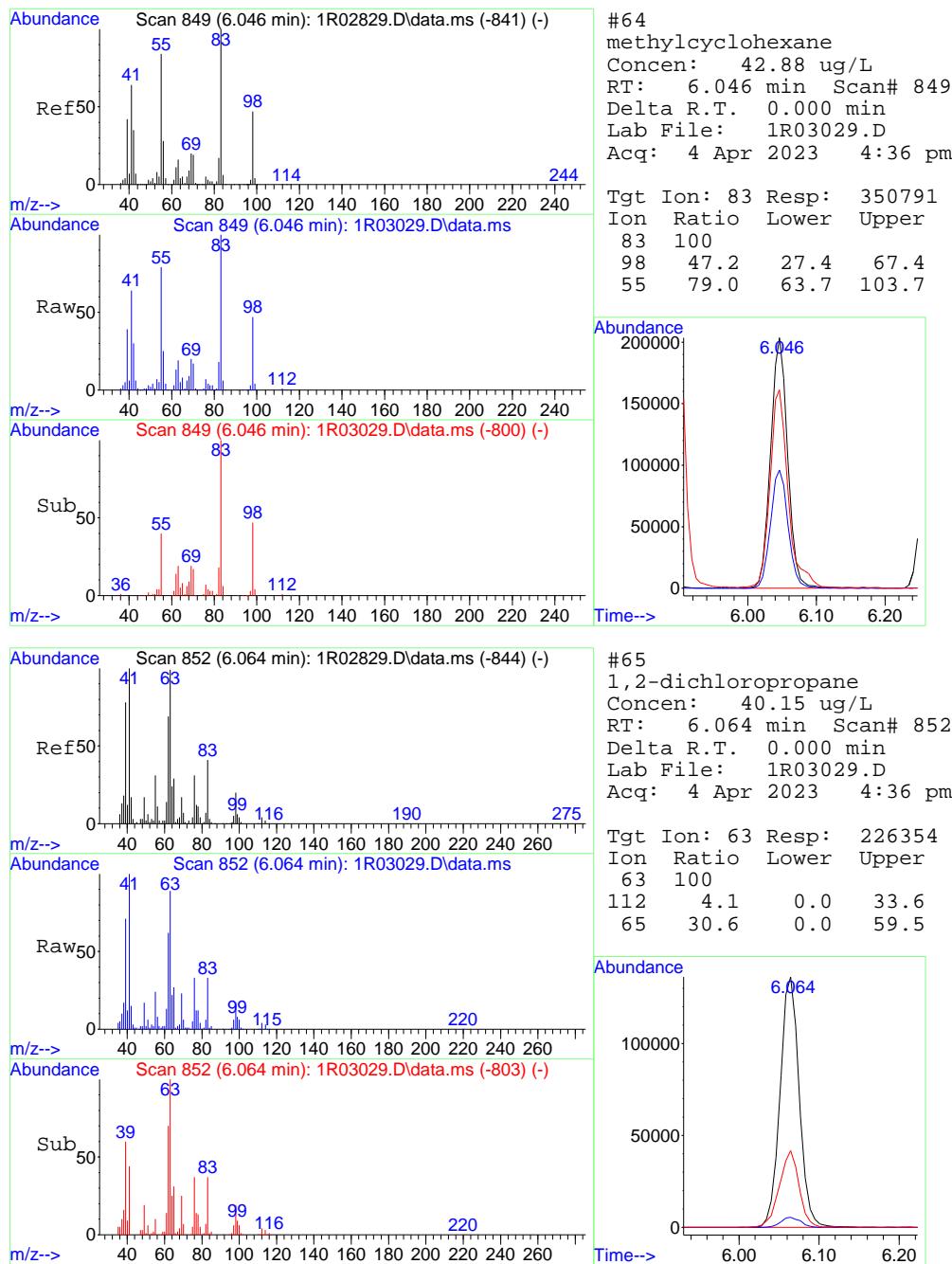


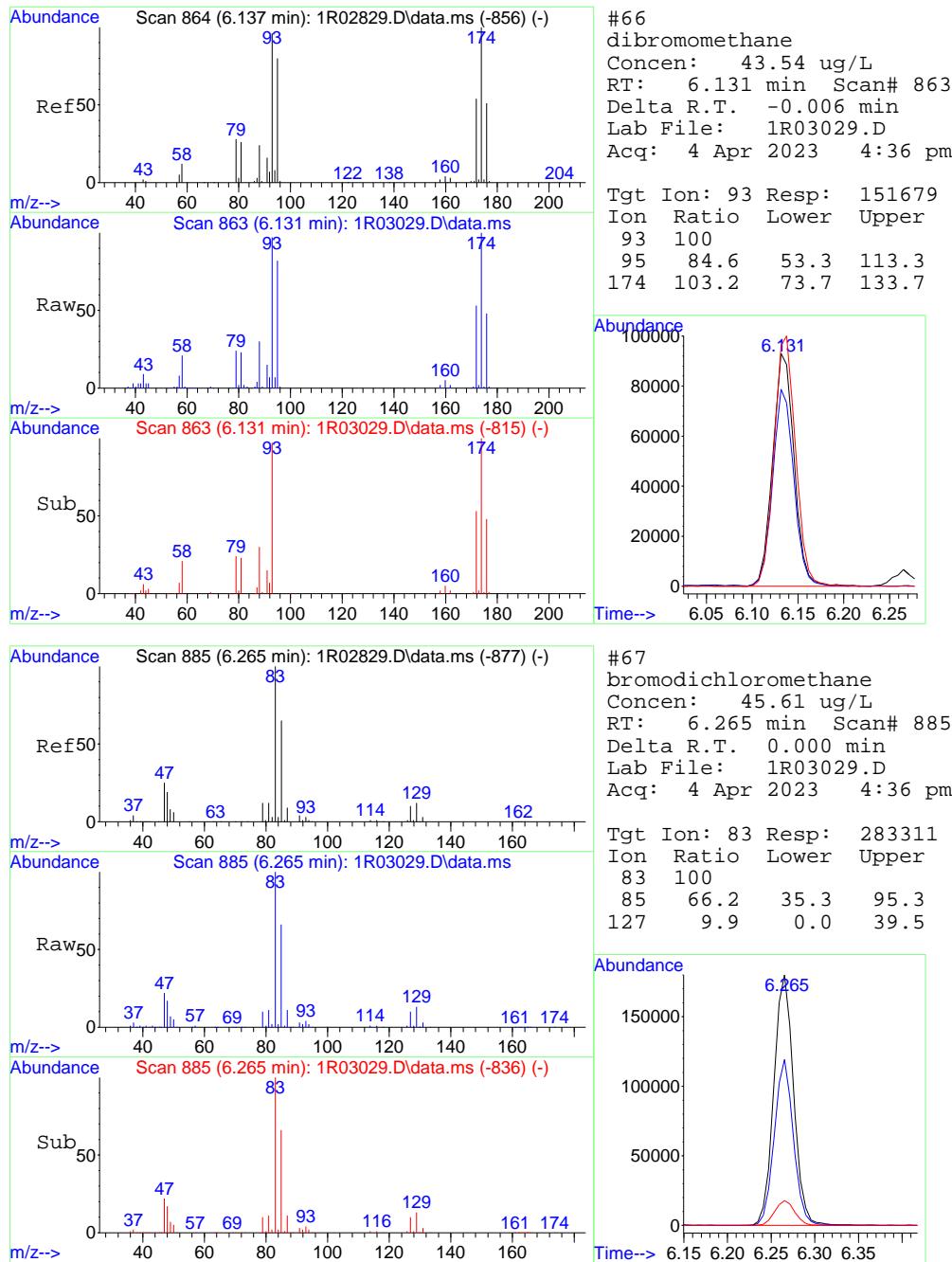


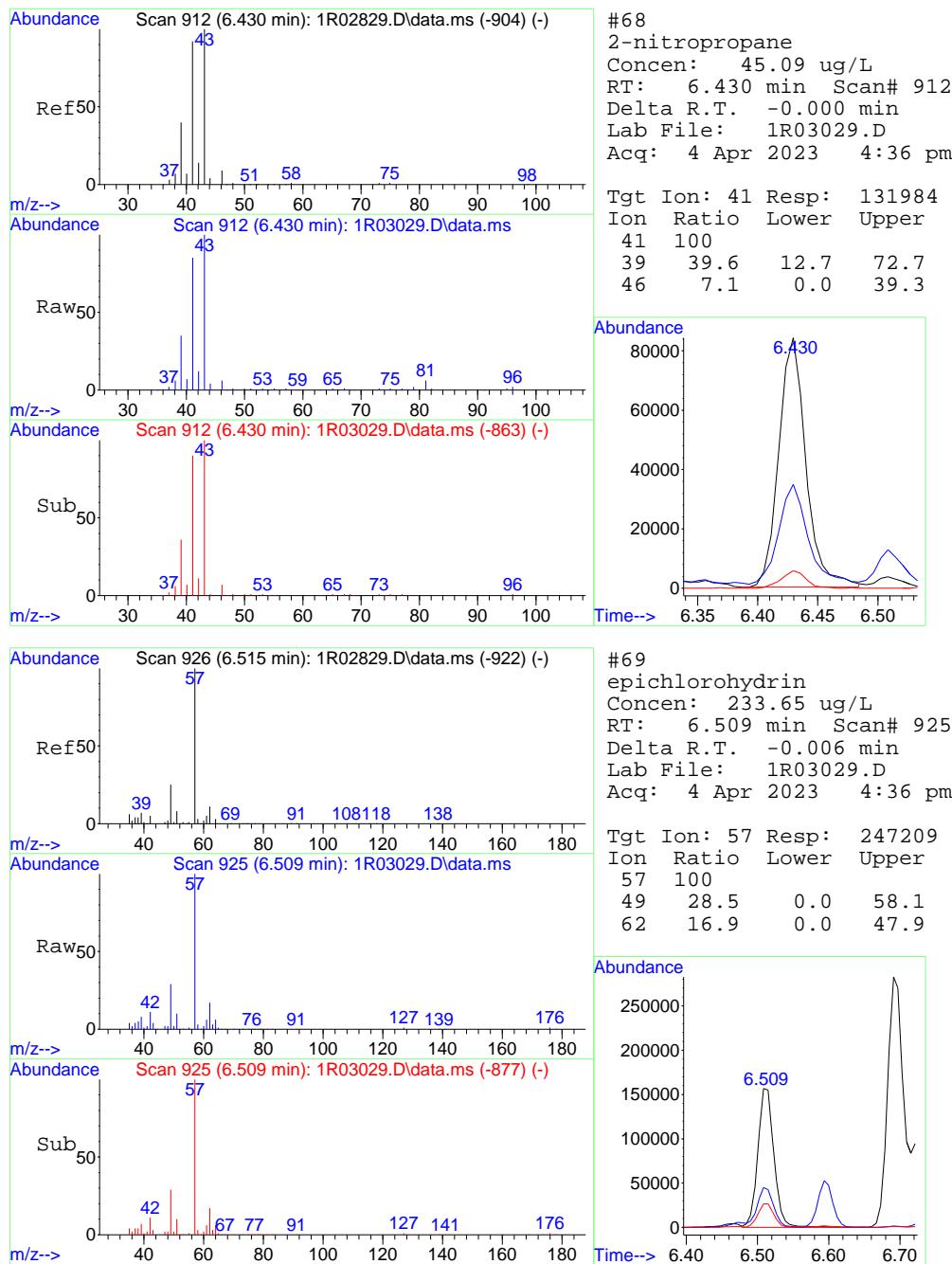


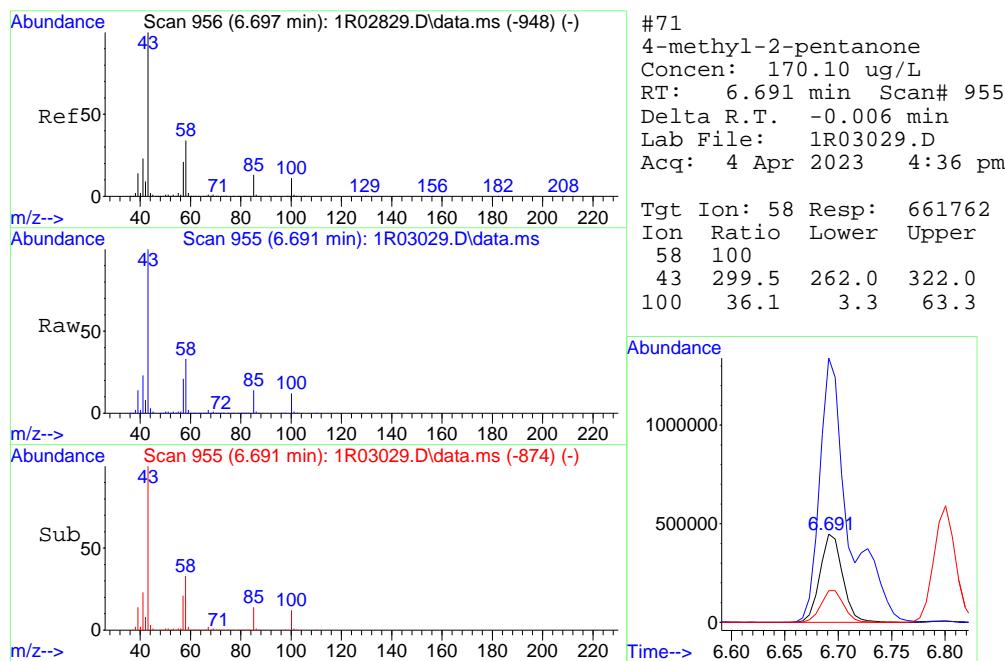
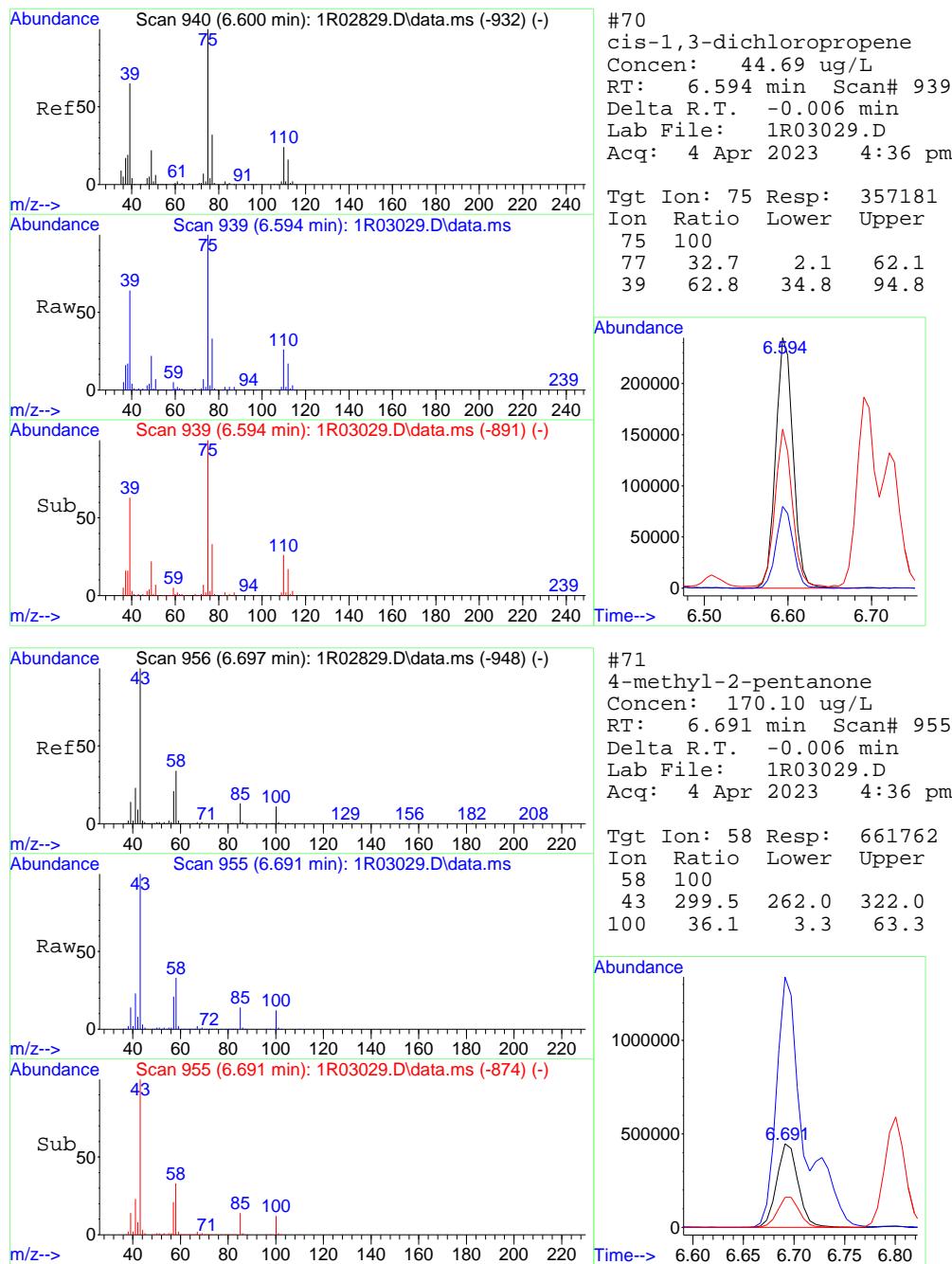


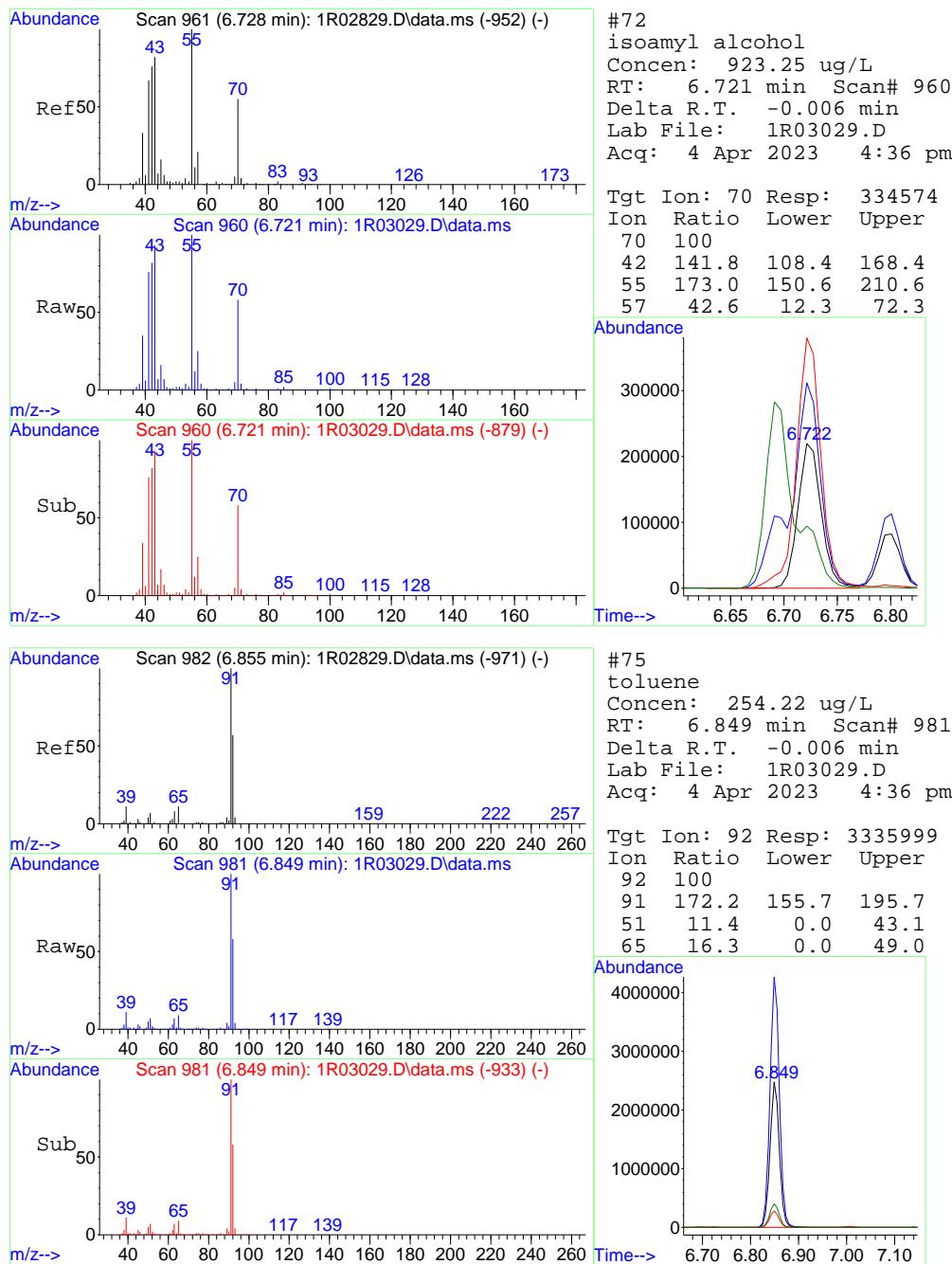


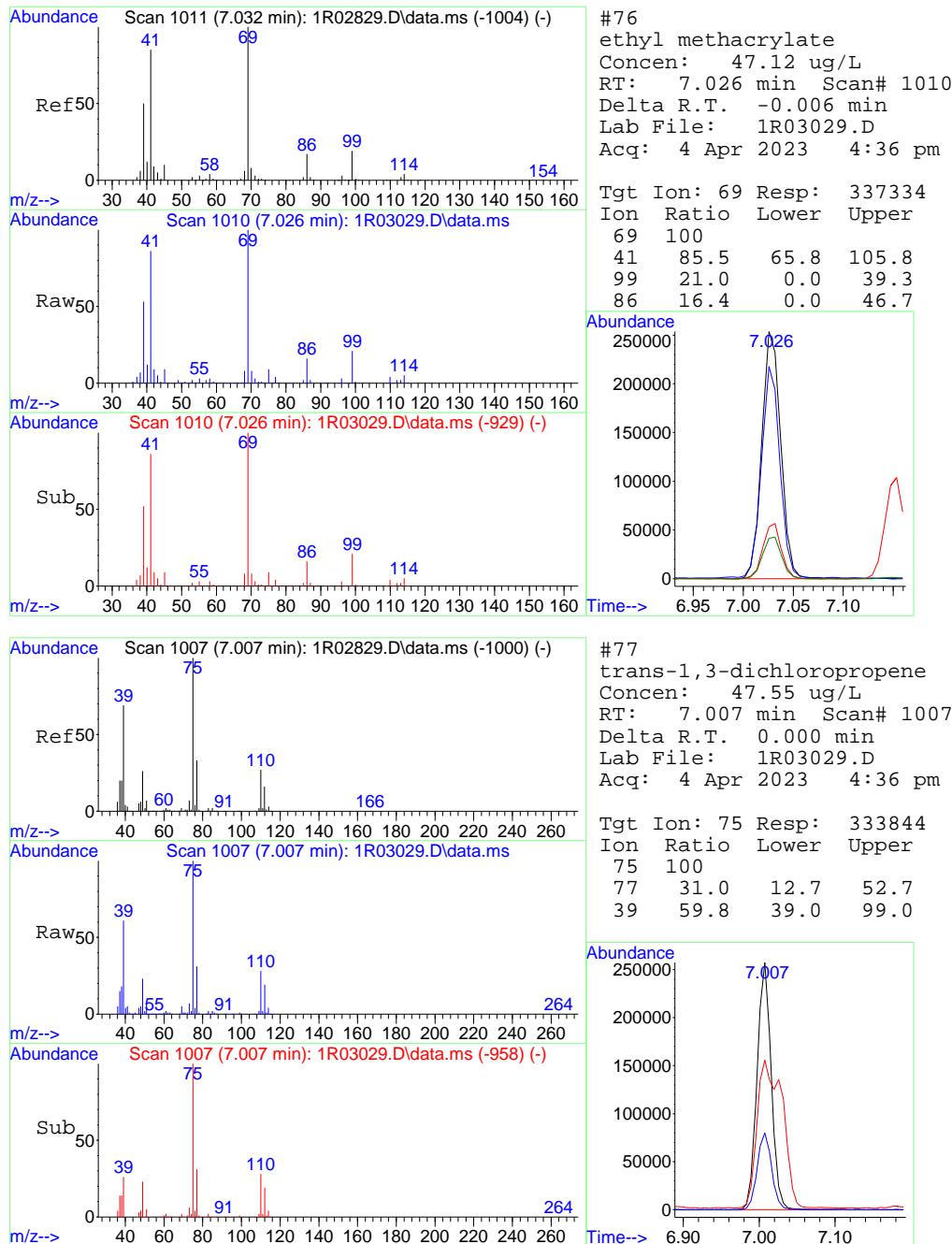


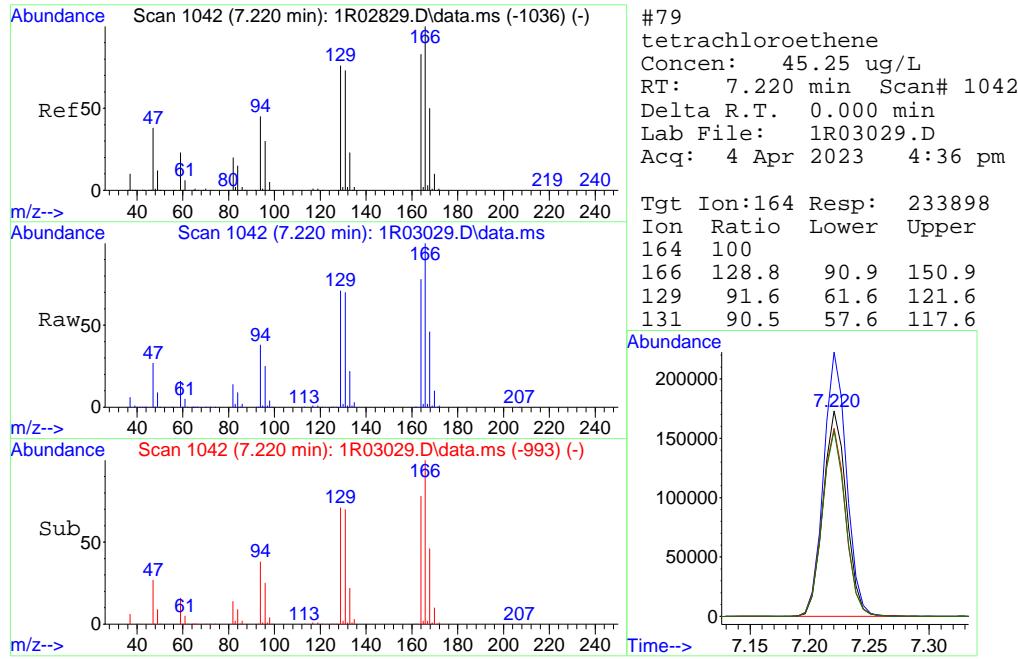
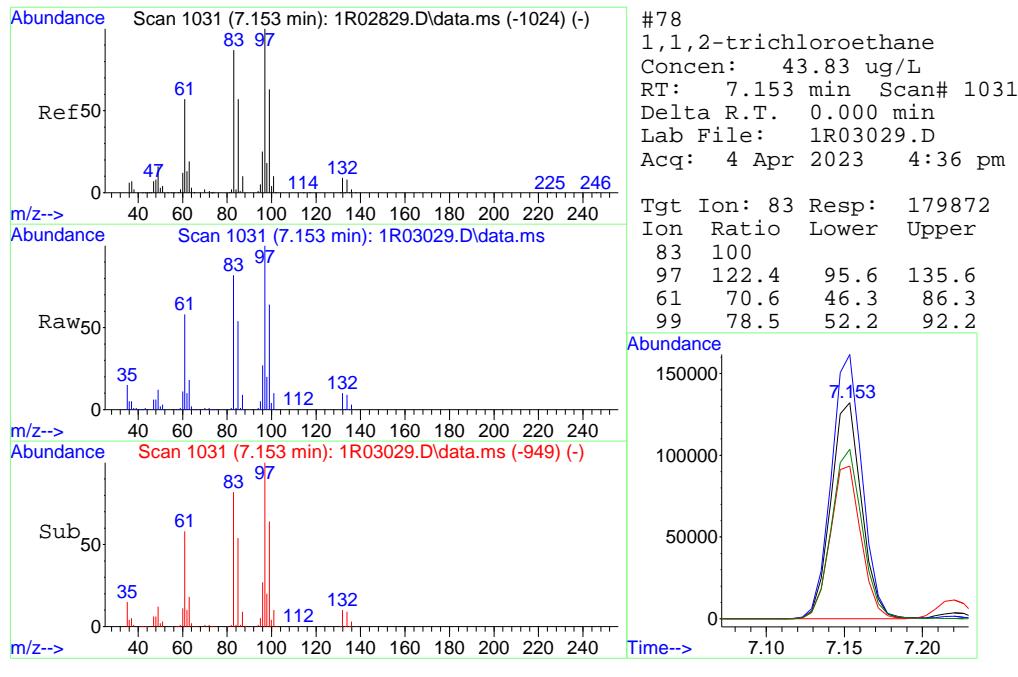


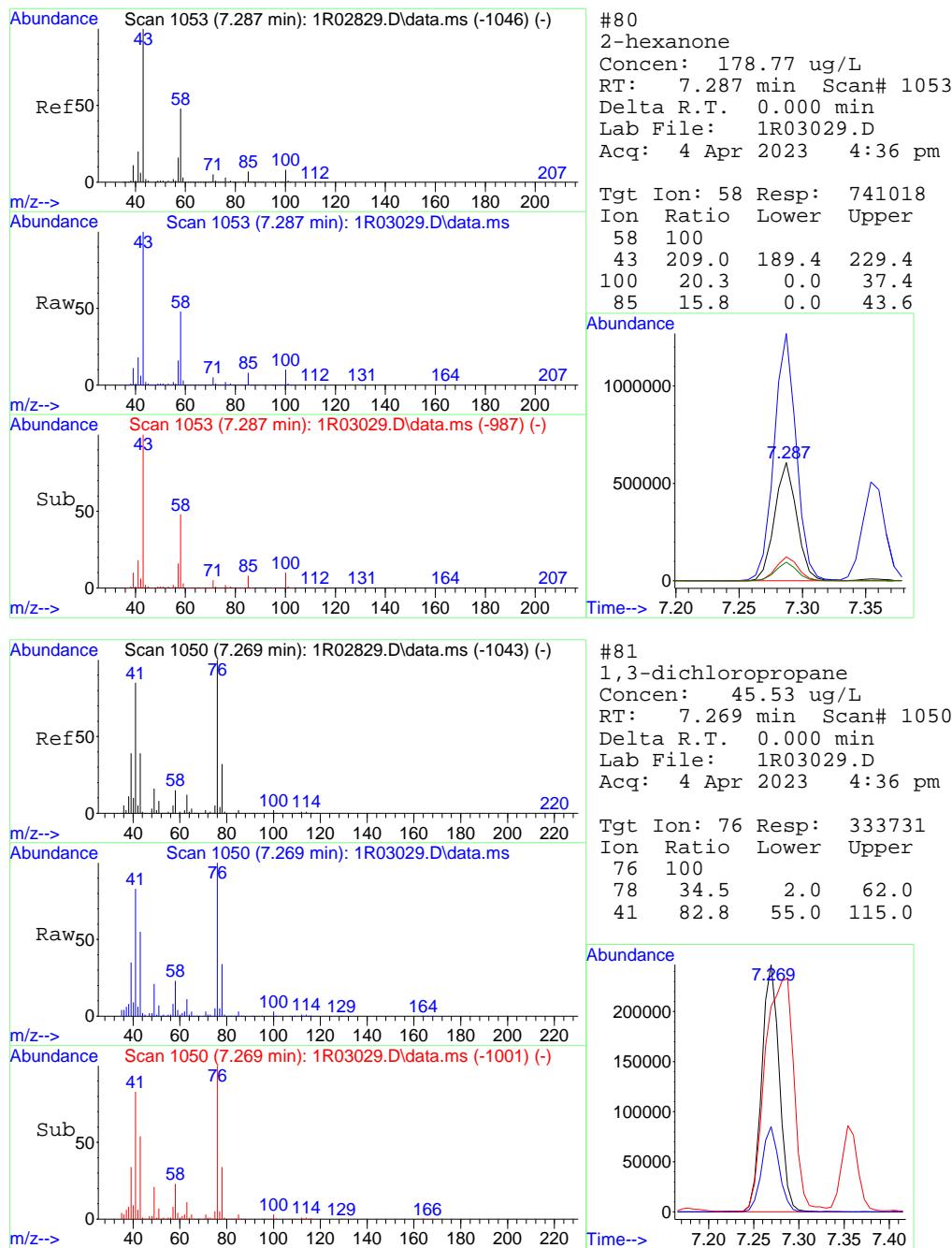


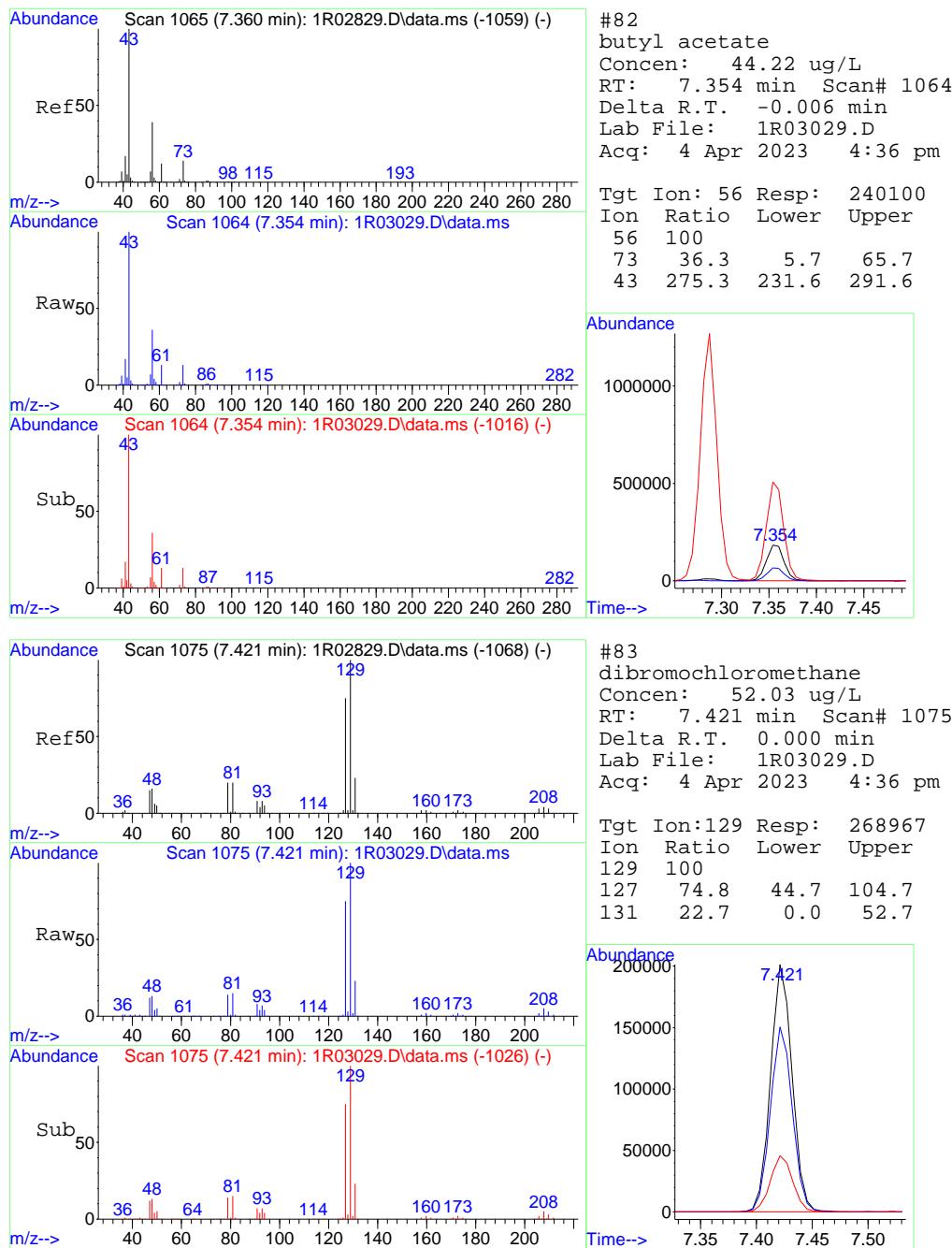


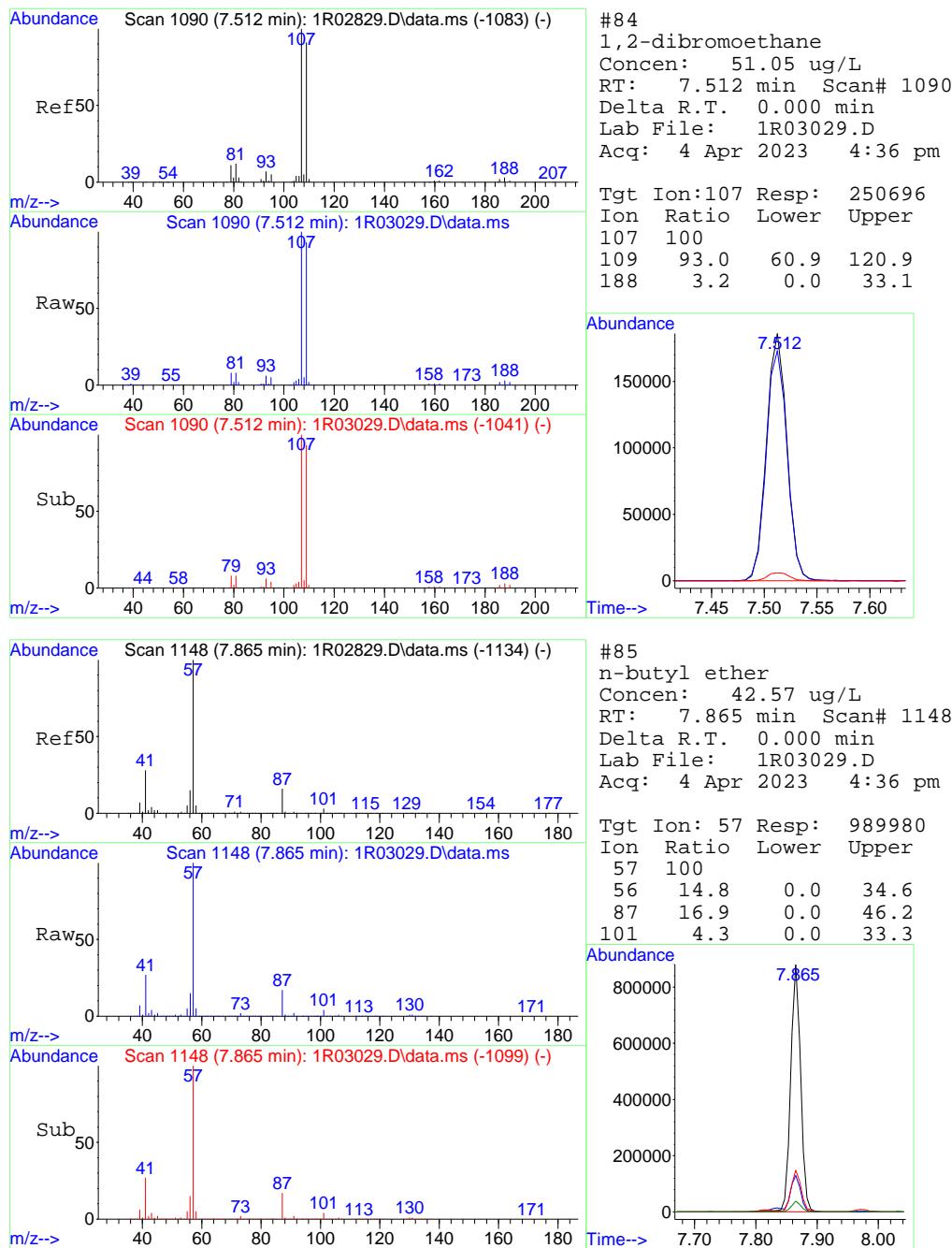


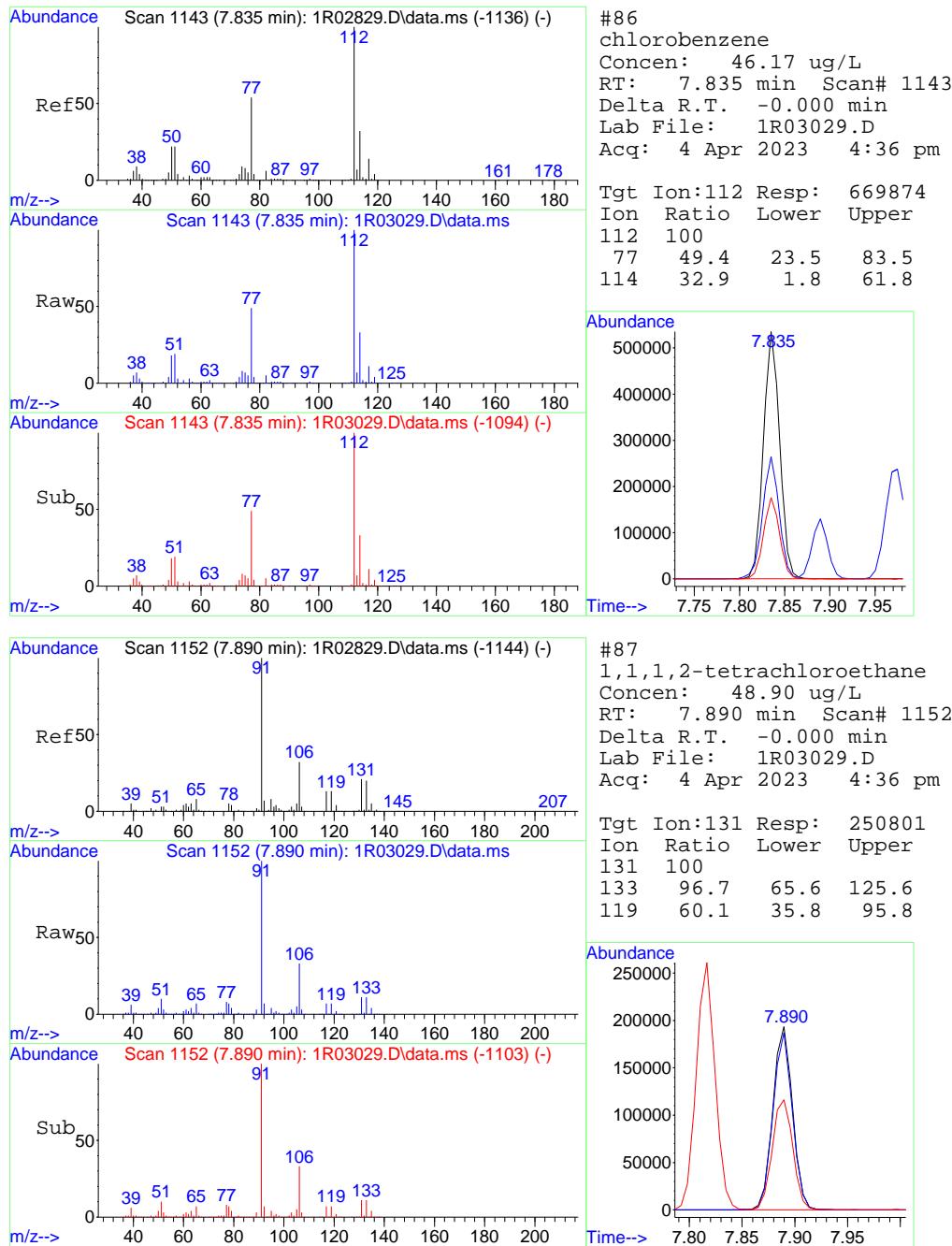


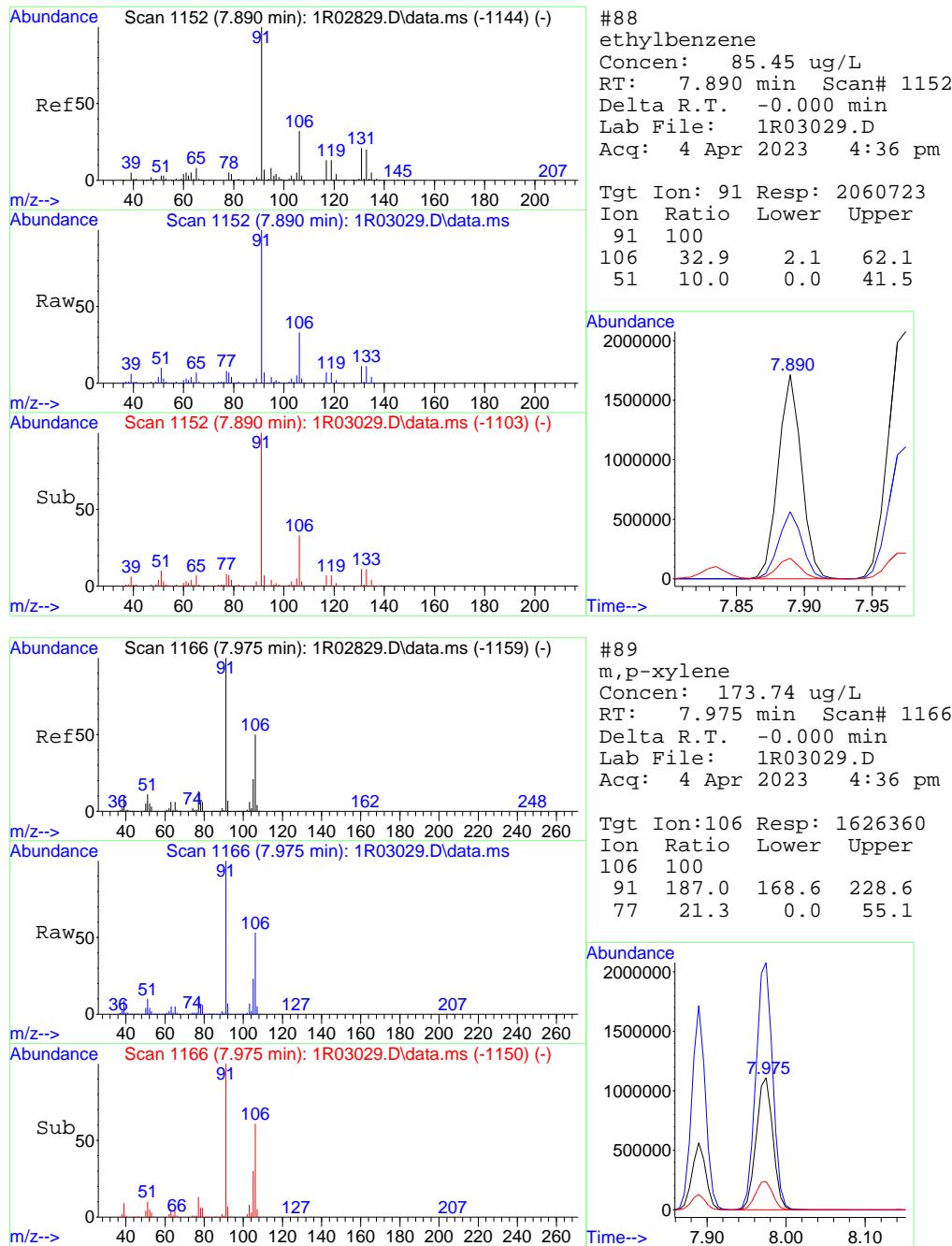


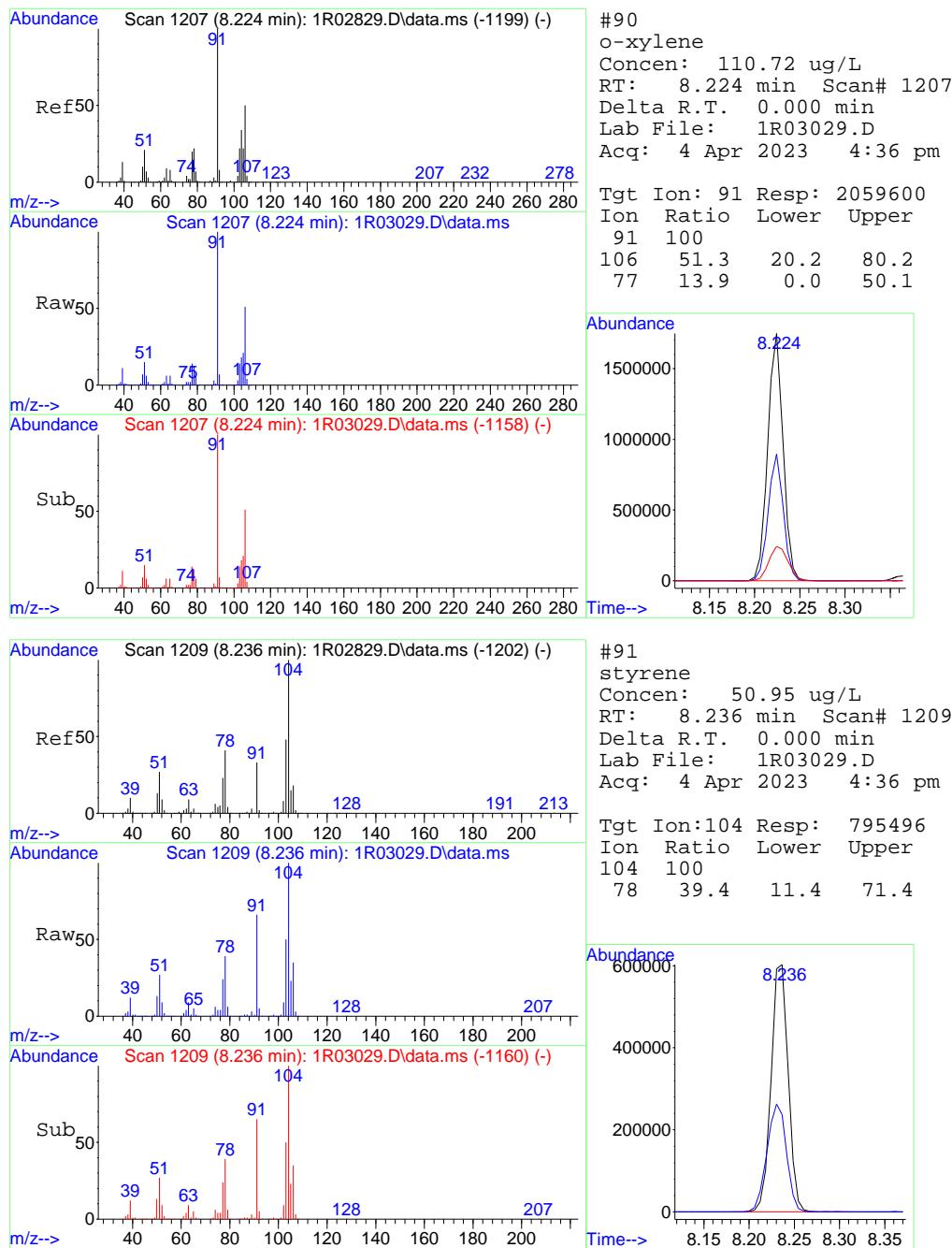


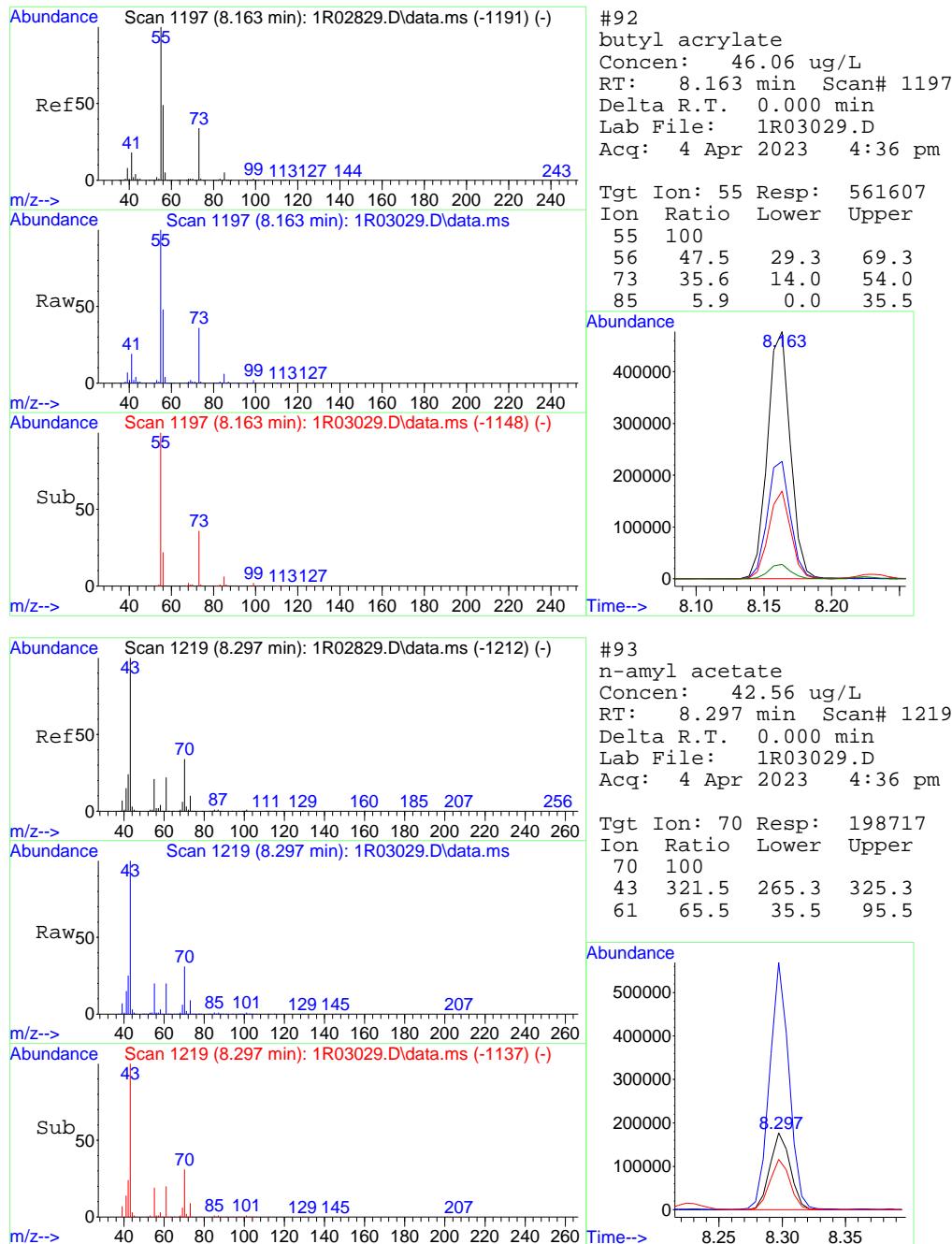


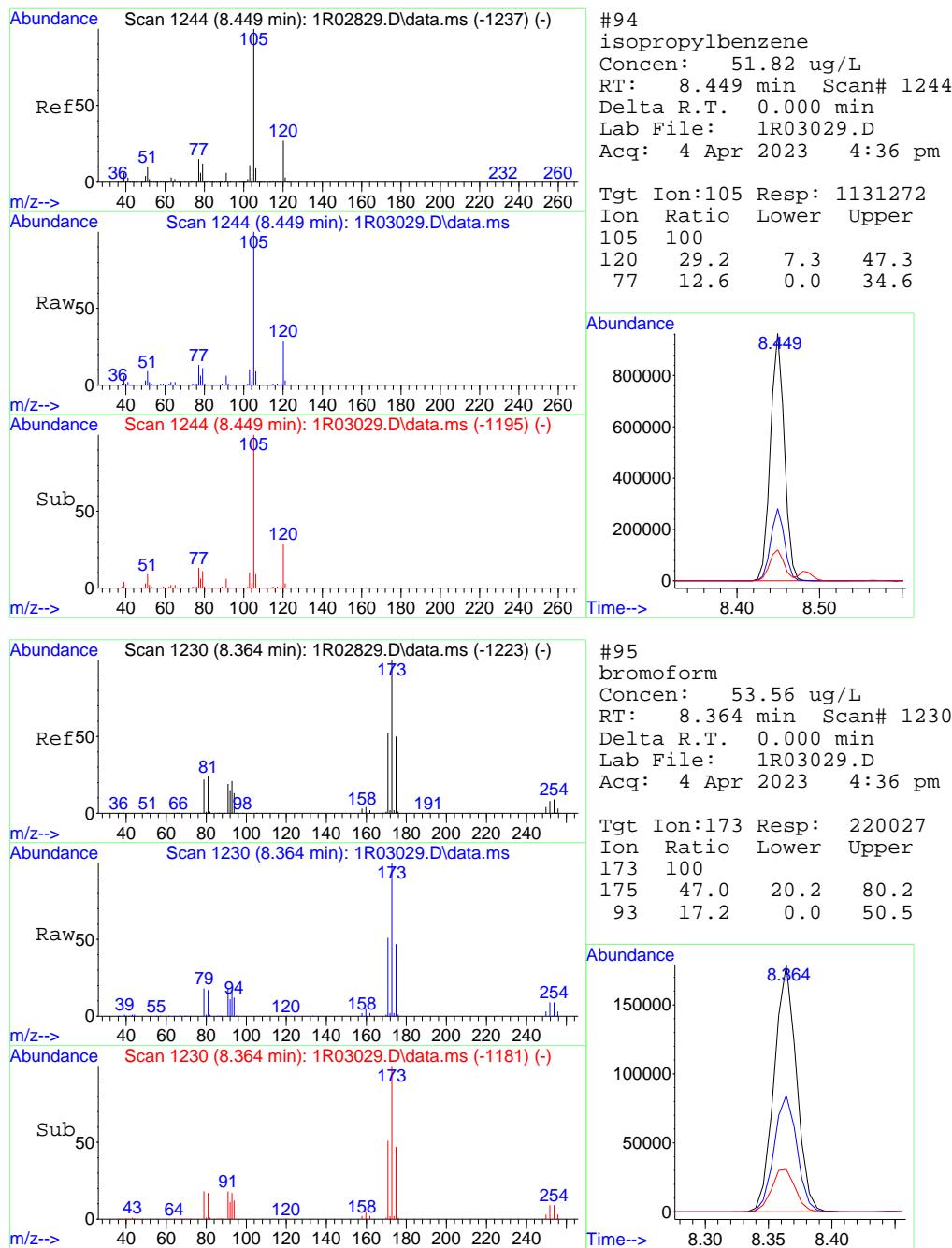


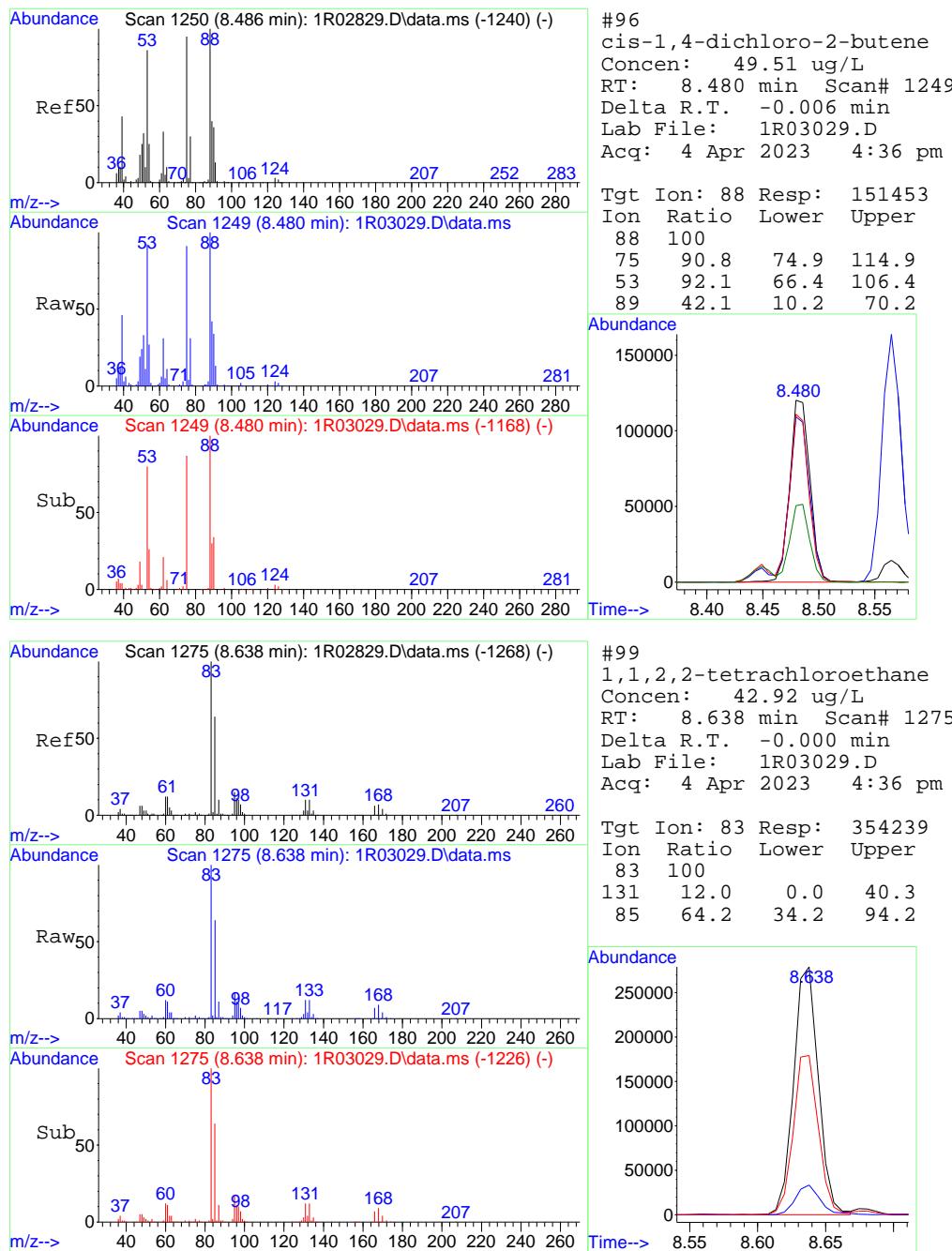


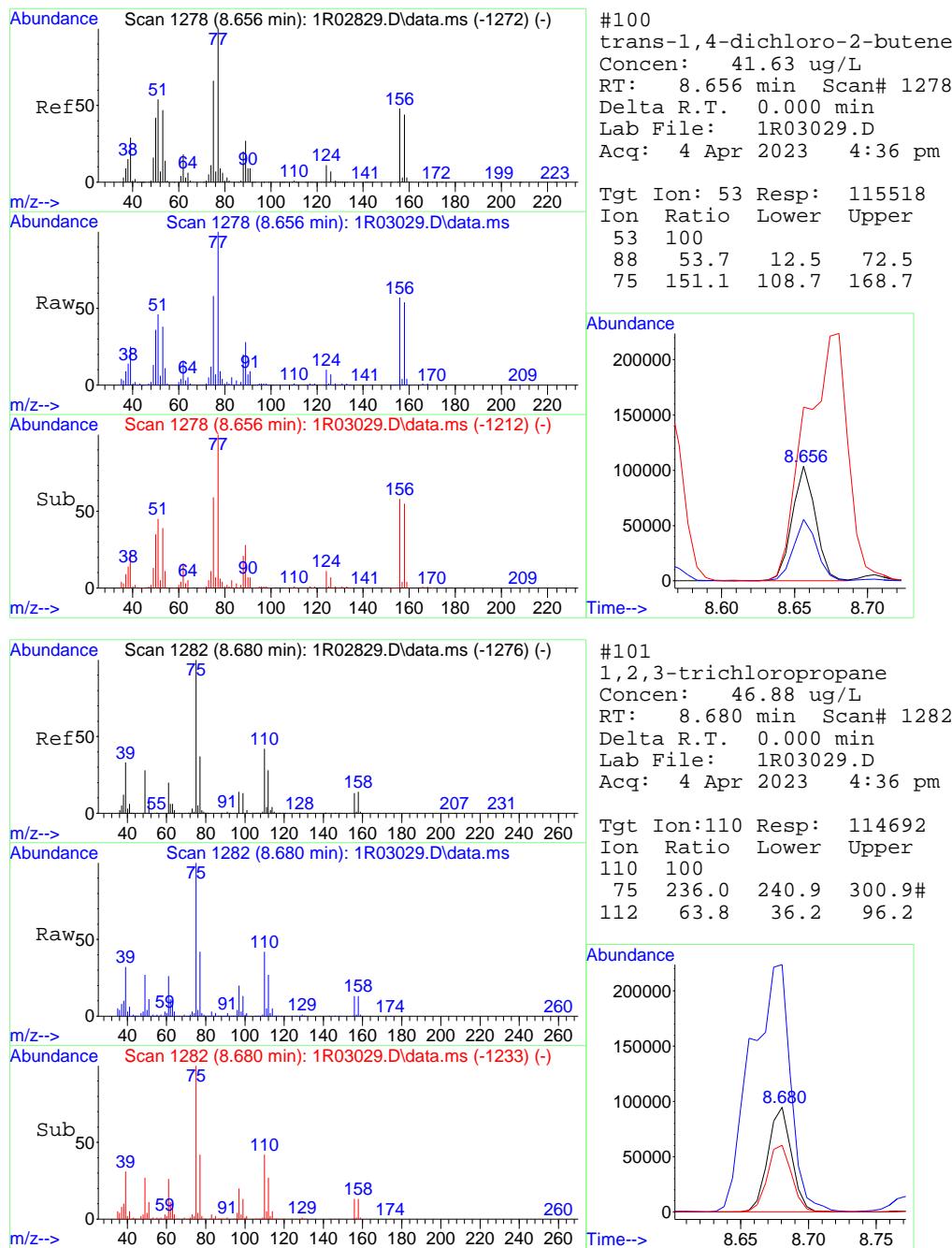


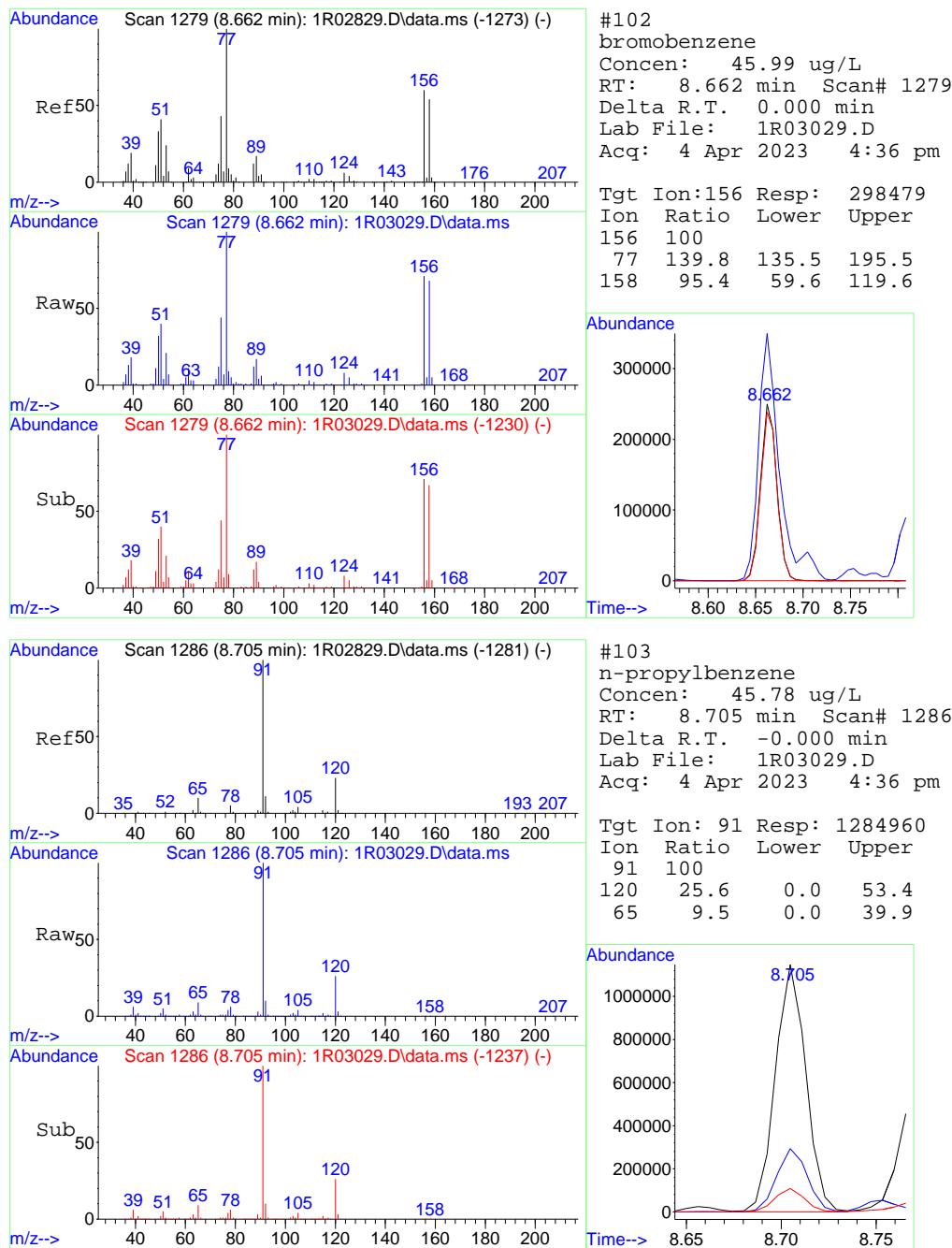


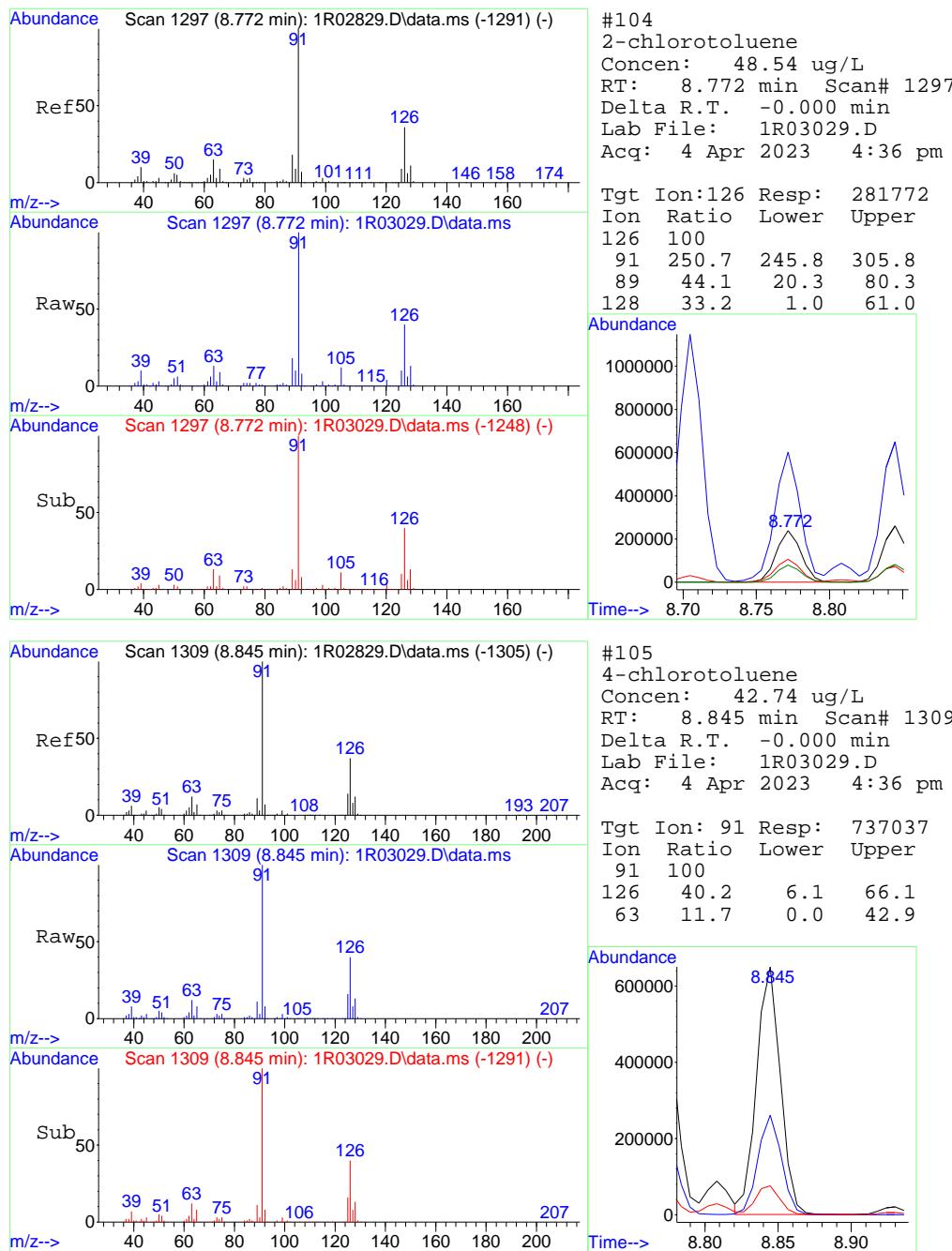


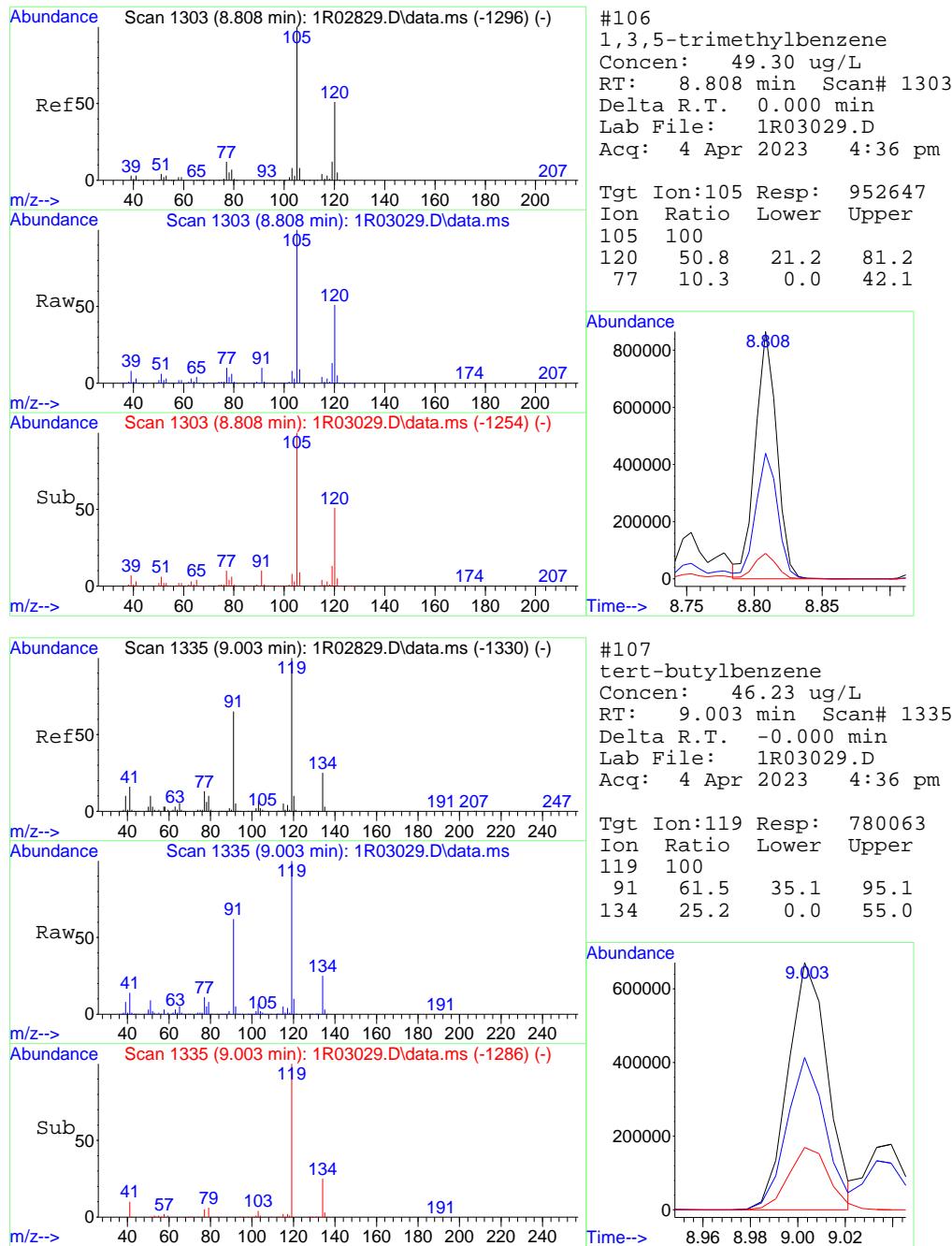


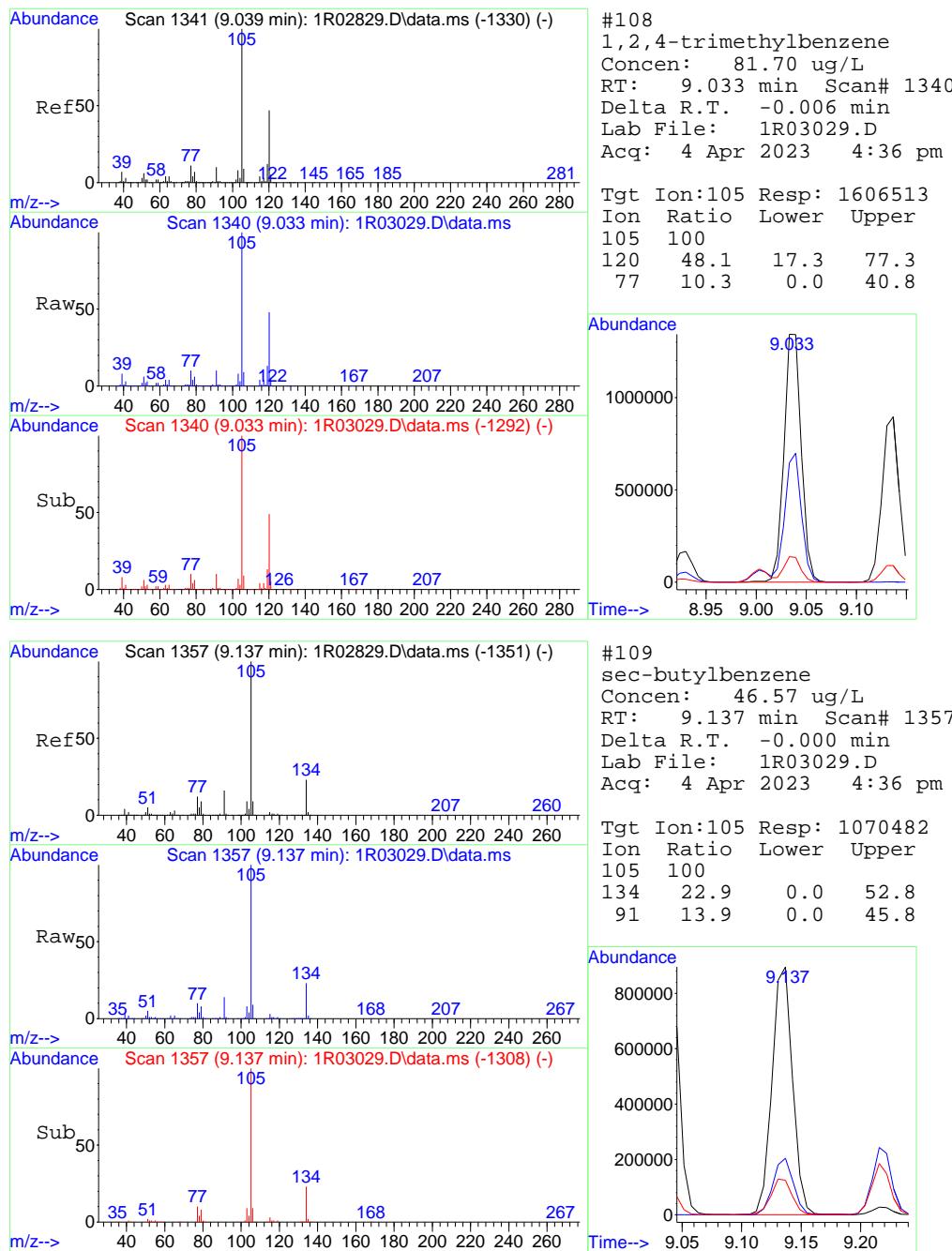


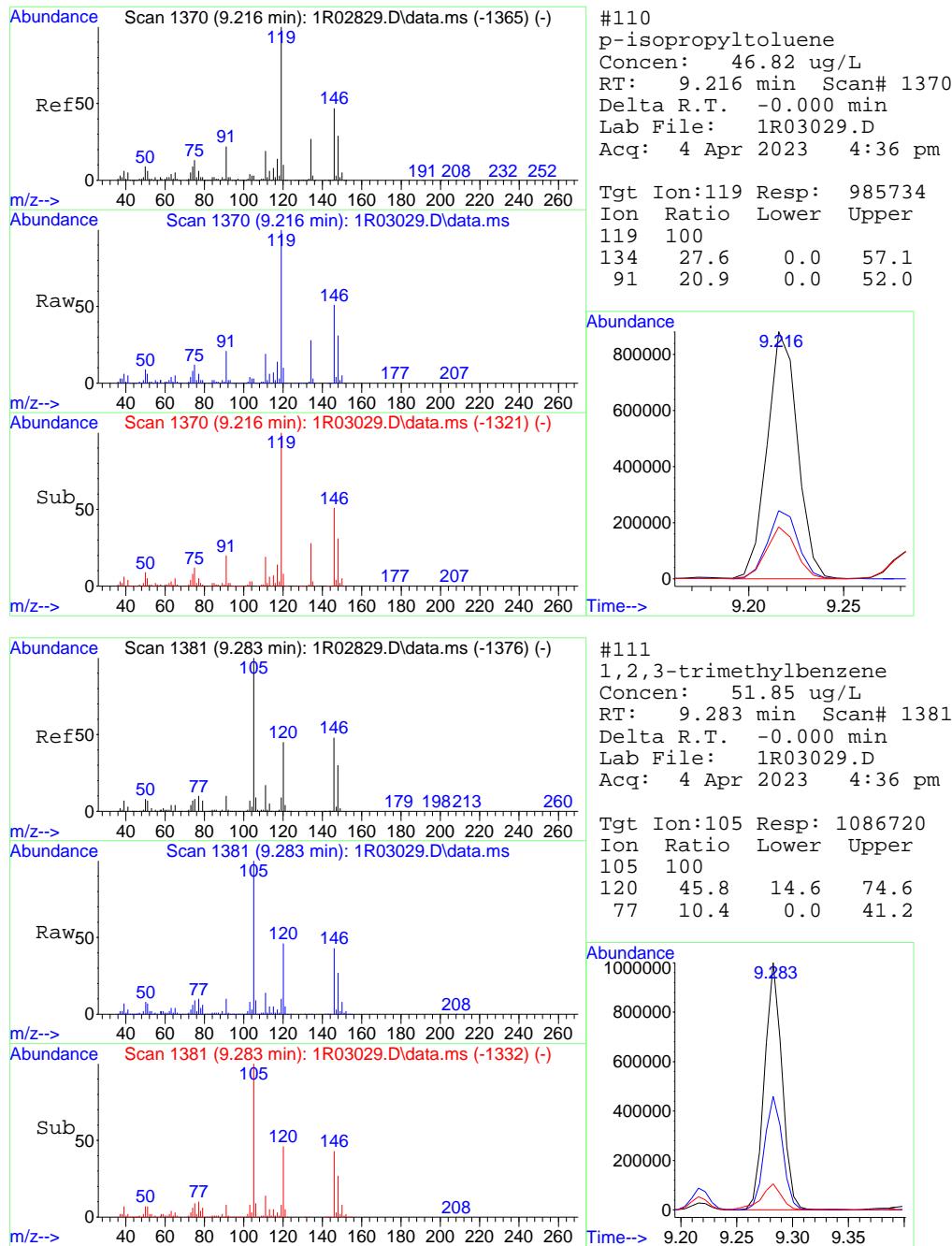


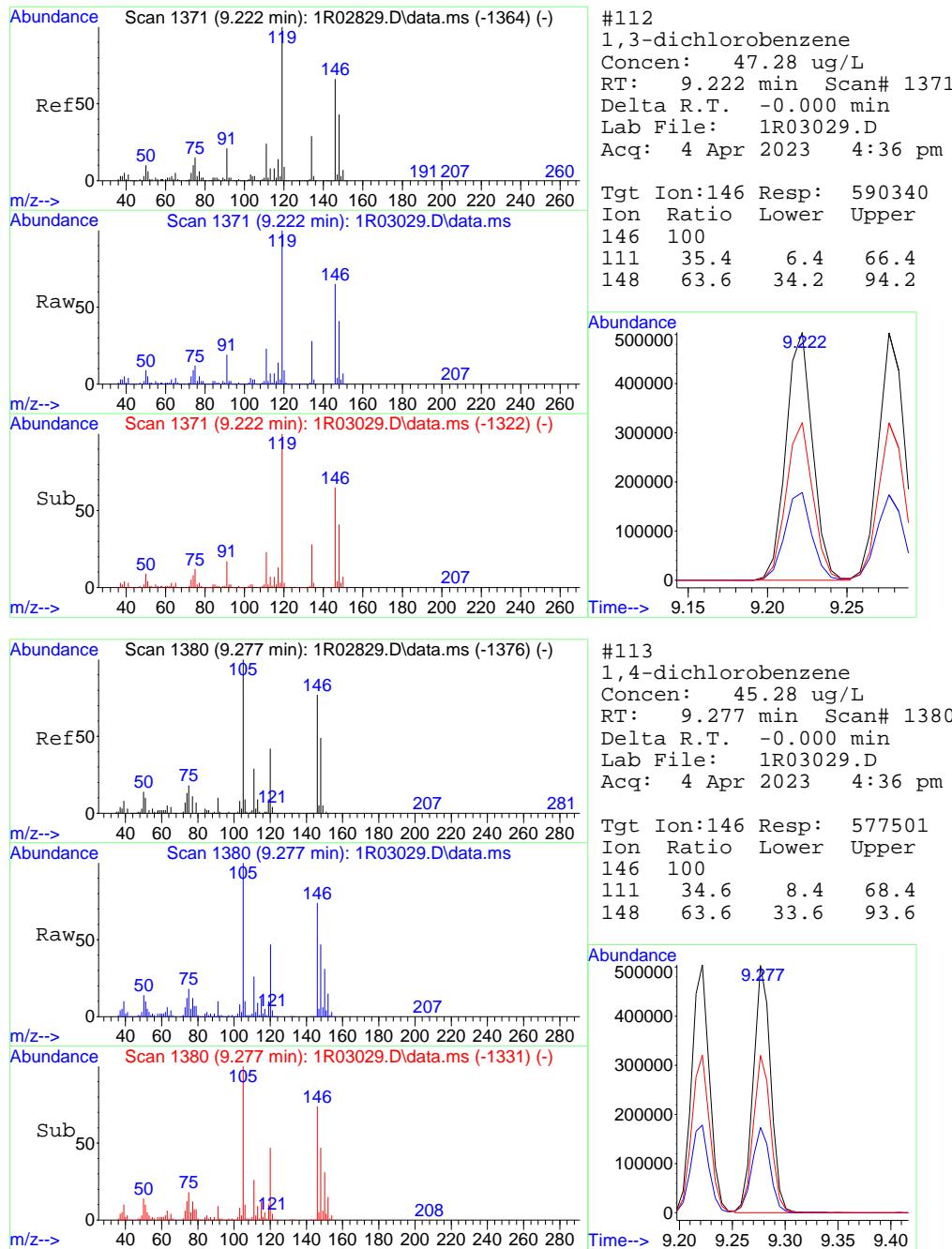


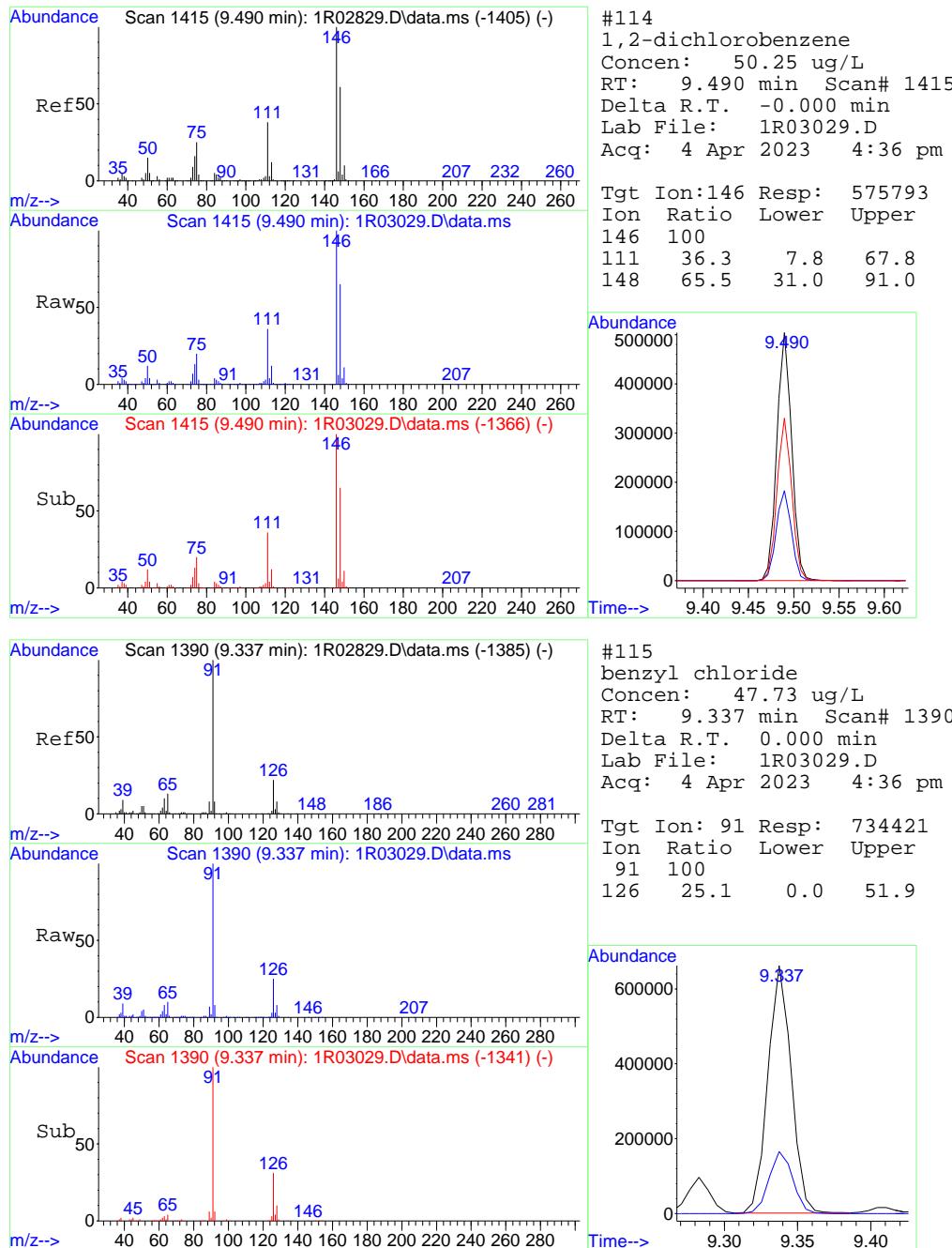


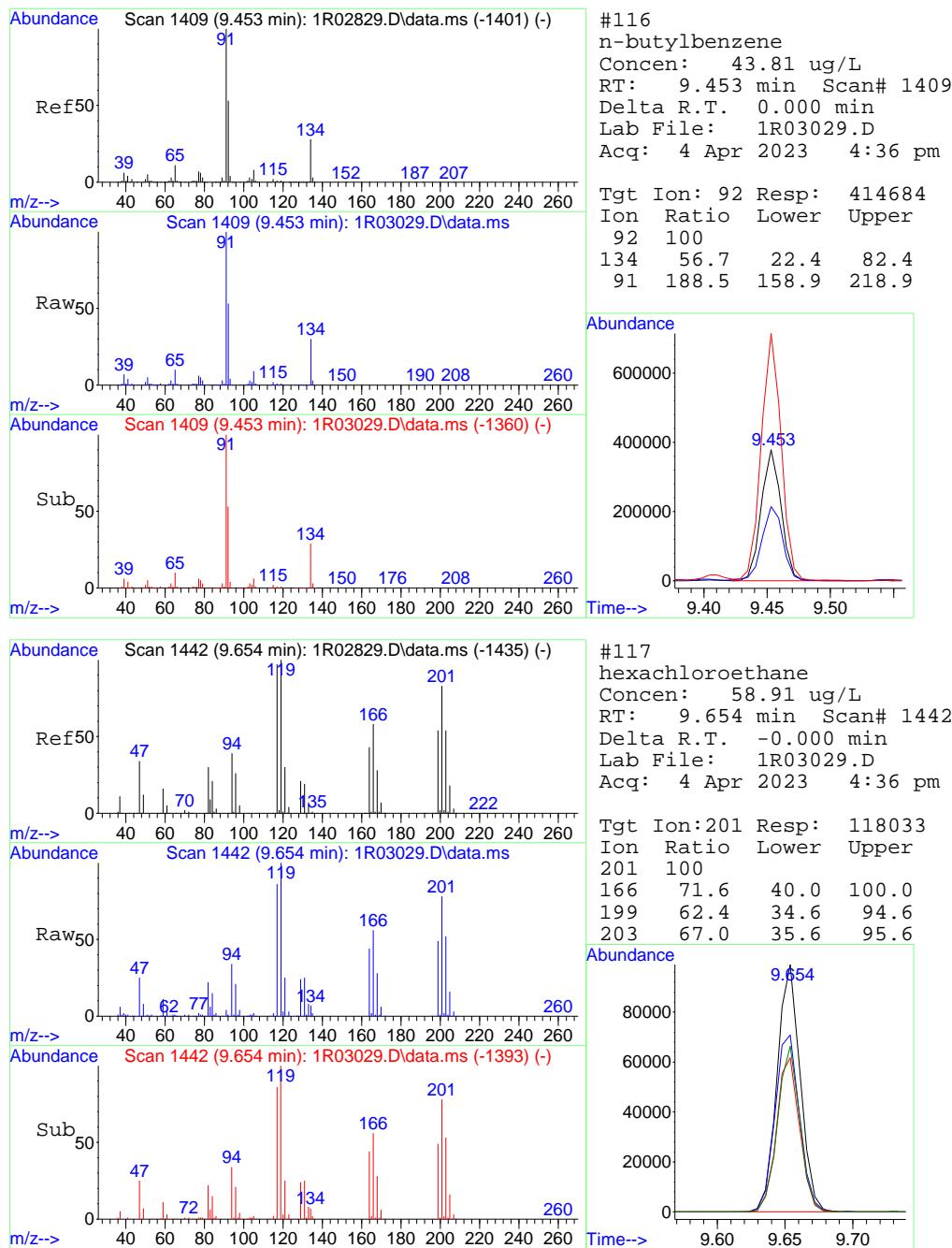


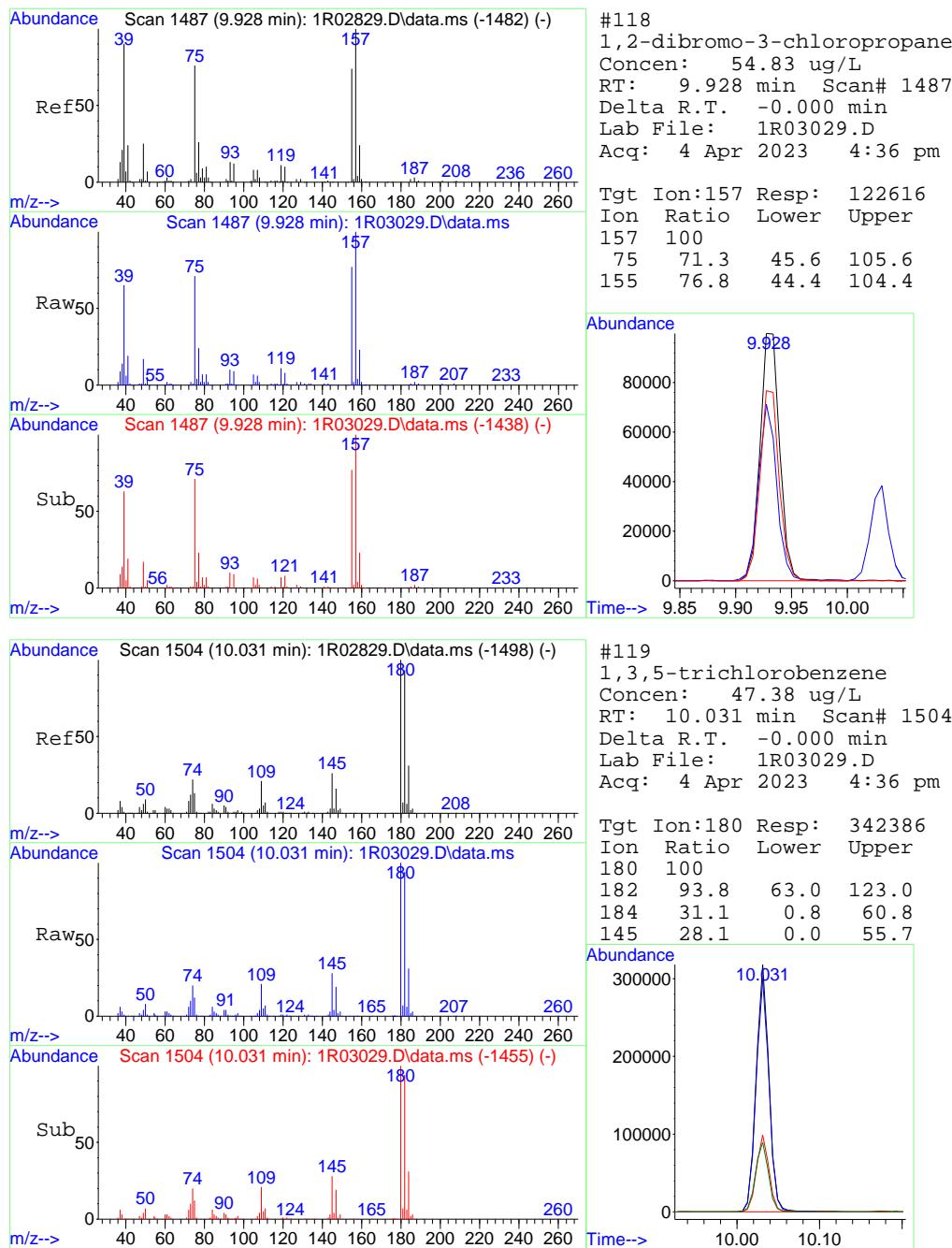


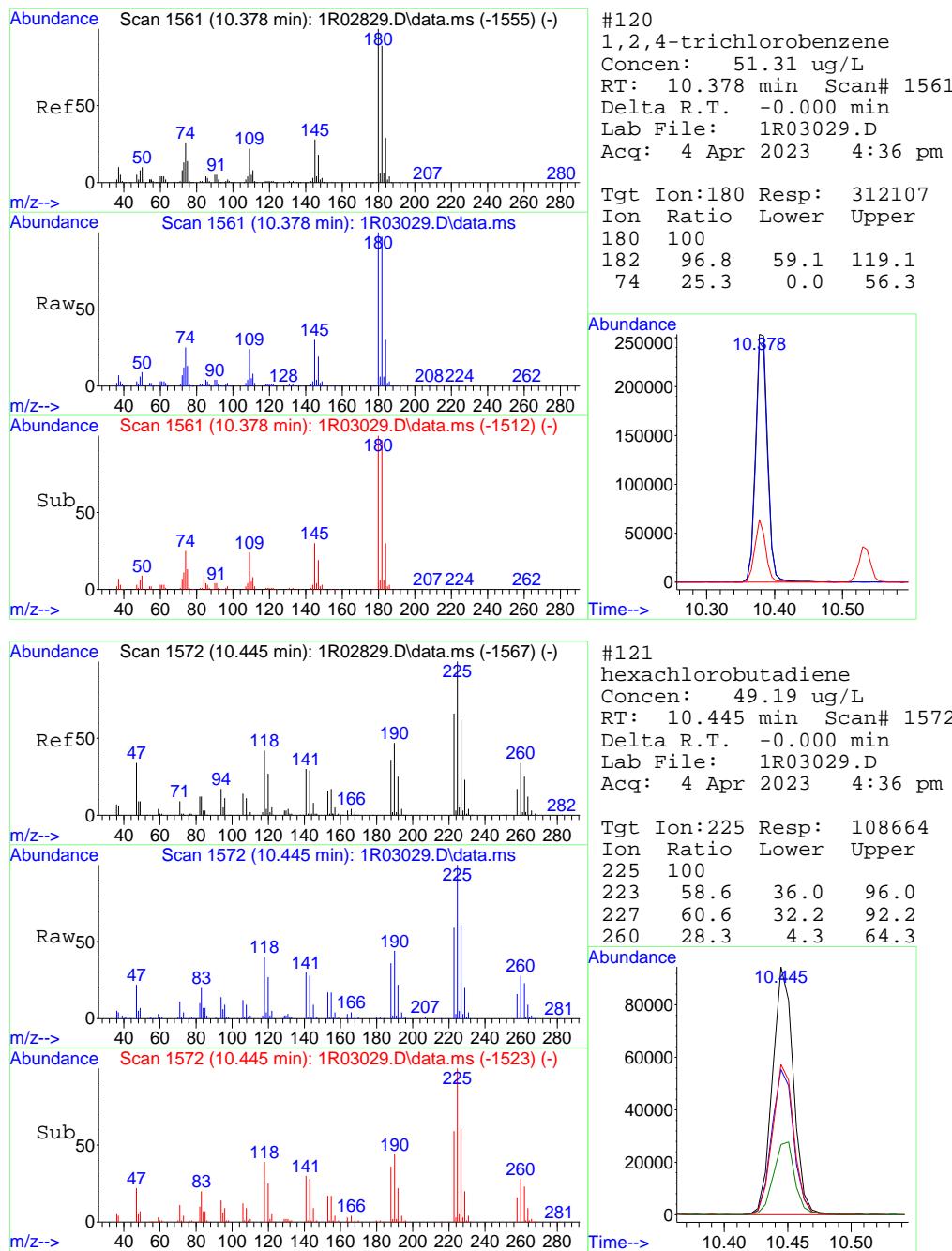


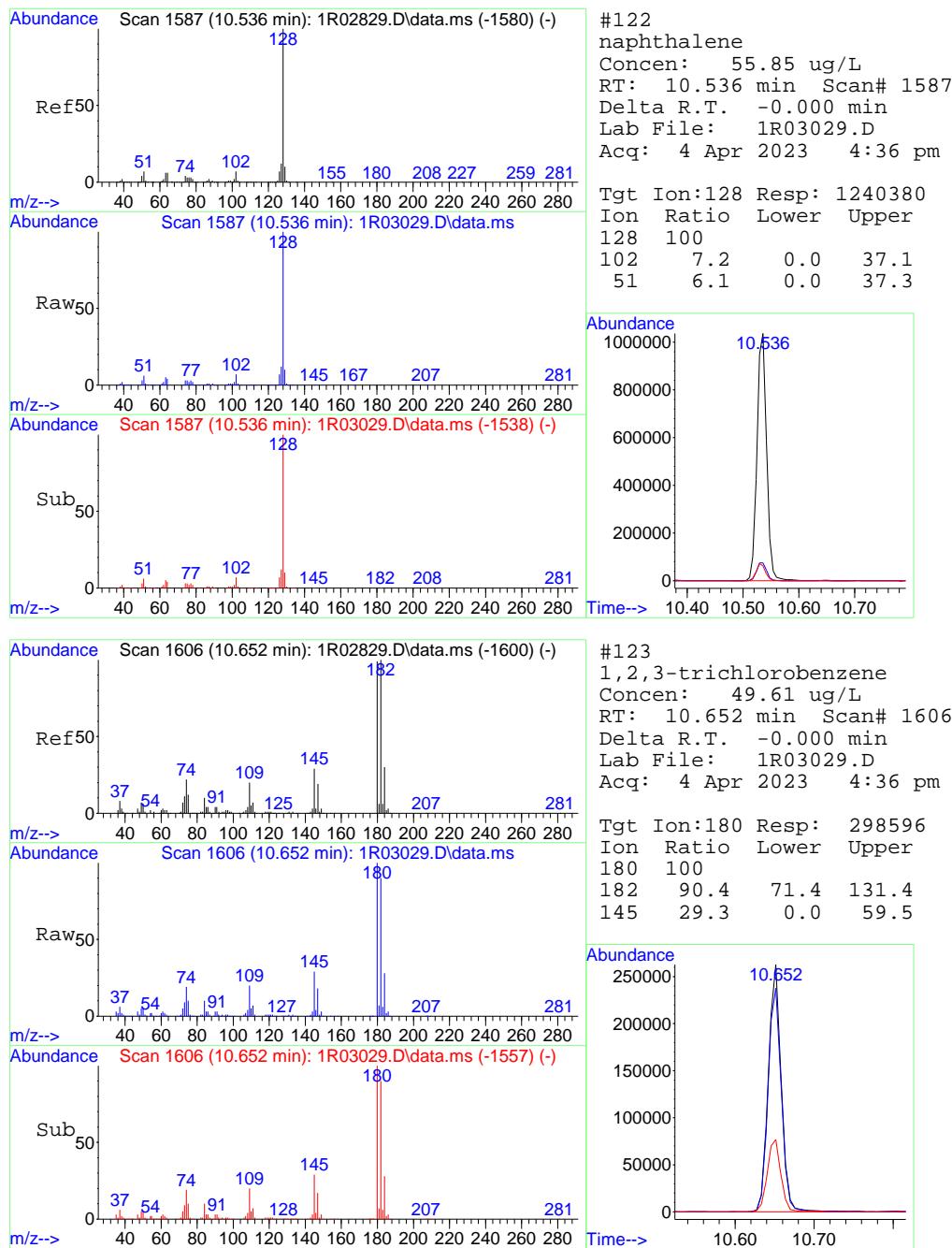


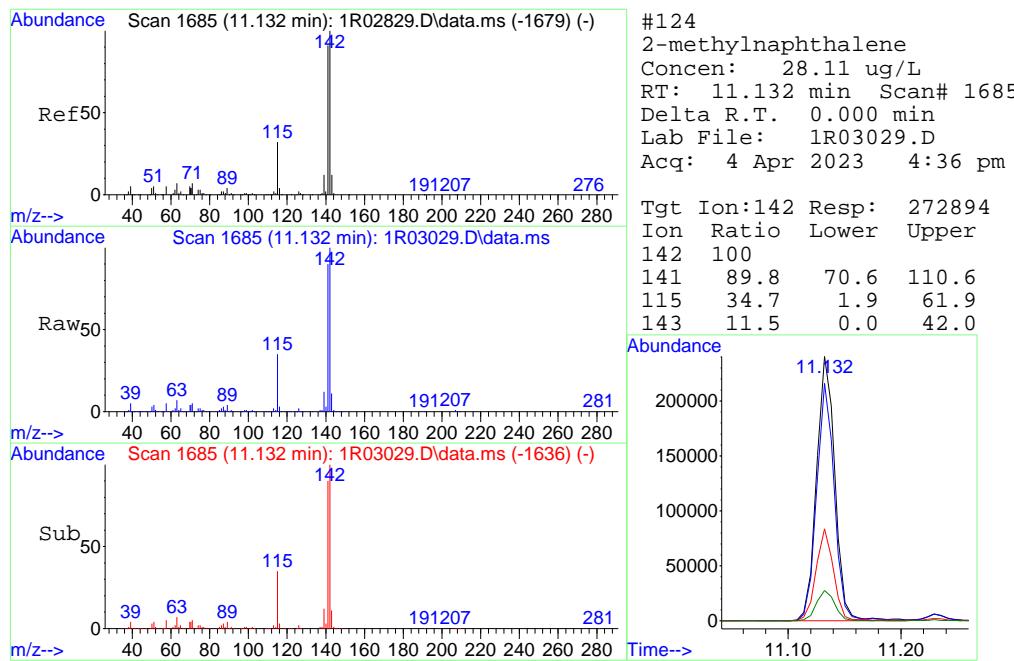








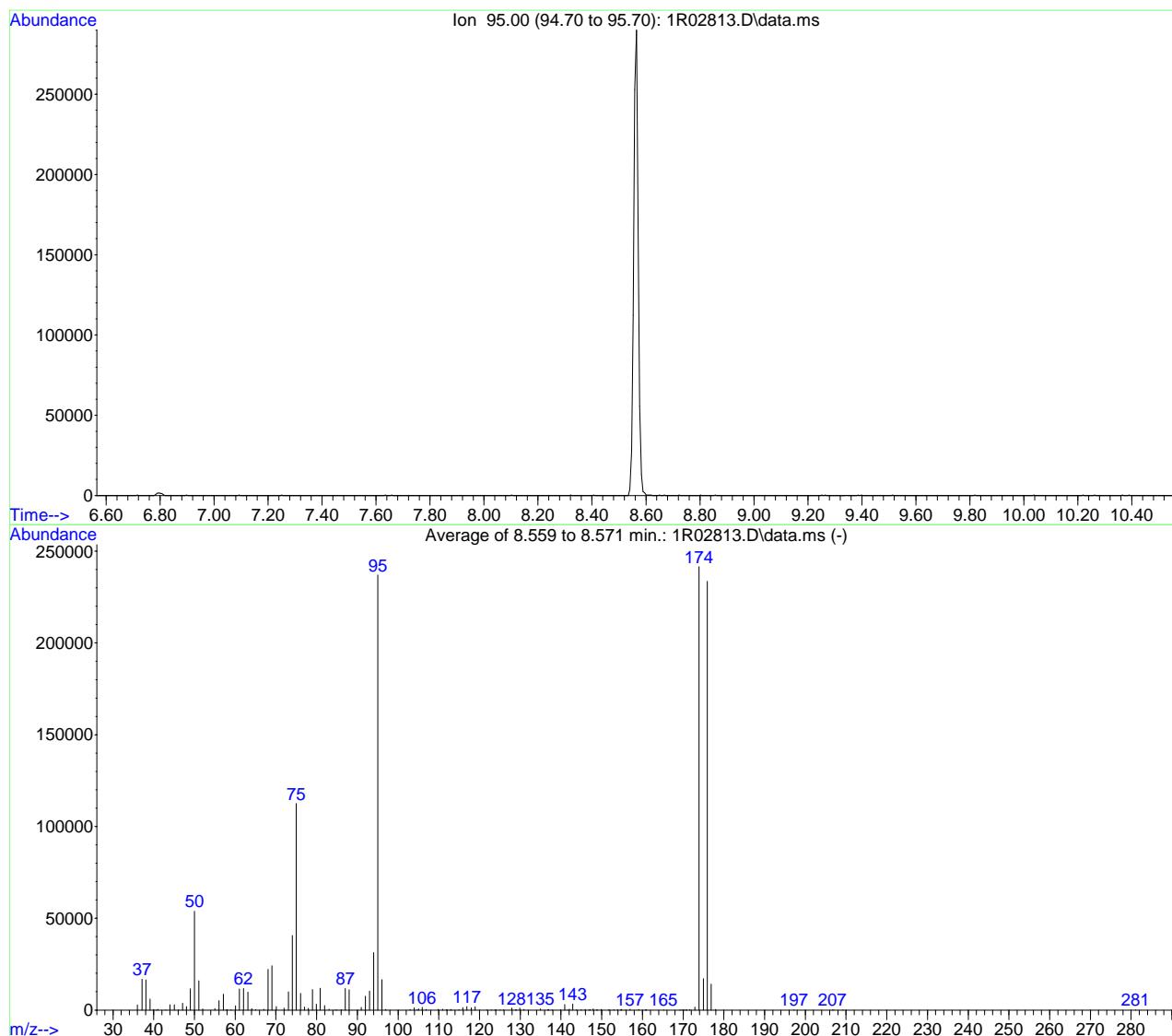




SW-846 Method 8260

Data File : C:\msdchem\1\data\V1r91\1R02813.D Vial: 1  
 Acq On : 28 Mar 2023 9:55 pm Operator: PrashanS  
 Sample : BFB Inst : GCMSR  
 Misc : MS67262,V1R0091,5,,,1 Multiplr: 1.00  
 MS Integration Params: rteint.p

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)  
 Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um



AutoFind: Scans 1262, 1263, 1264; Background Corrected with Scan 1256

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	22.7	53808	PASS
75	95	30	60	47.5	112563	PASS
95	95	100	100	100.0	237077	PASS
96	95	5	9	7.0	16559	PASS
173	174	0.00	2	0.7	1620	PASS
174	95	50	150	101.9	241536	PASS
175	174	5	9	7.1	17164	PASS
176	174	95	101	96.7	233536	PASS
177	176	5	9	6.0	14108	PASS

Average of 8.559 to 8.571 min.: 1R02813.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.00	2791	47.05	3690	59.00	82	72.00	1057
37.10	16822	48.00	1909	60.10	2352	73.05	9913
38.10	16400	49.00	11565	61.00	11439	74.00	40669
39.10	5986	50.00	53808	62.05	11712	75.00	112563
40.00	369	51.10	16037	63.10	9752	76.05	9157
40.70	82	52.05	709	64.05	811	77.00	1604
41.05	291	55.00	826	65.00	174	77.95	905
42.95	25	56.05	5206	67.00	474	79.00	11201
44.00	2866	57.10	8707	68.05	22211	79.95	3172
45.05	2924	57.90	209	69.00	24202	80.90	11862
46.05	148	58.05	210	70.10	1960	81.95	2419

Average of 8.559 to 8.571 min.: 1R02813.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
83.05	514	96.00	16559	112.10	58	124.00	193
84.60	58	96.95	368	112.90	332	125.10	124
85.00	65	102.95	383	113.10	128	125.90	73
86.00	267	103.95	1254	114.95	196	127.90	1121
87.00	11726	104.70	178	115.90	1395	128.80	150
88.00	11115	105.05	582	116.90	1874	129.00	203
90.95	1422	105.95	1542	117.95	1220	129.90	992
92.00	7508	106.90	314	118.90	1739	130.90	409
93.00	10304	109.85	321	121.80	89	134.95	600
94.00	31381	110.95	504	122.90	91	136.80	236
95.00	237077	111.90	233	123.80	72	136.95	335

Average of 8.559 to 8.571 min.: 1R02813.D\data.ms

BFB

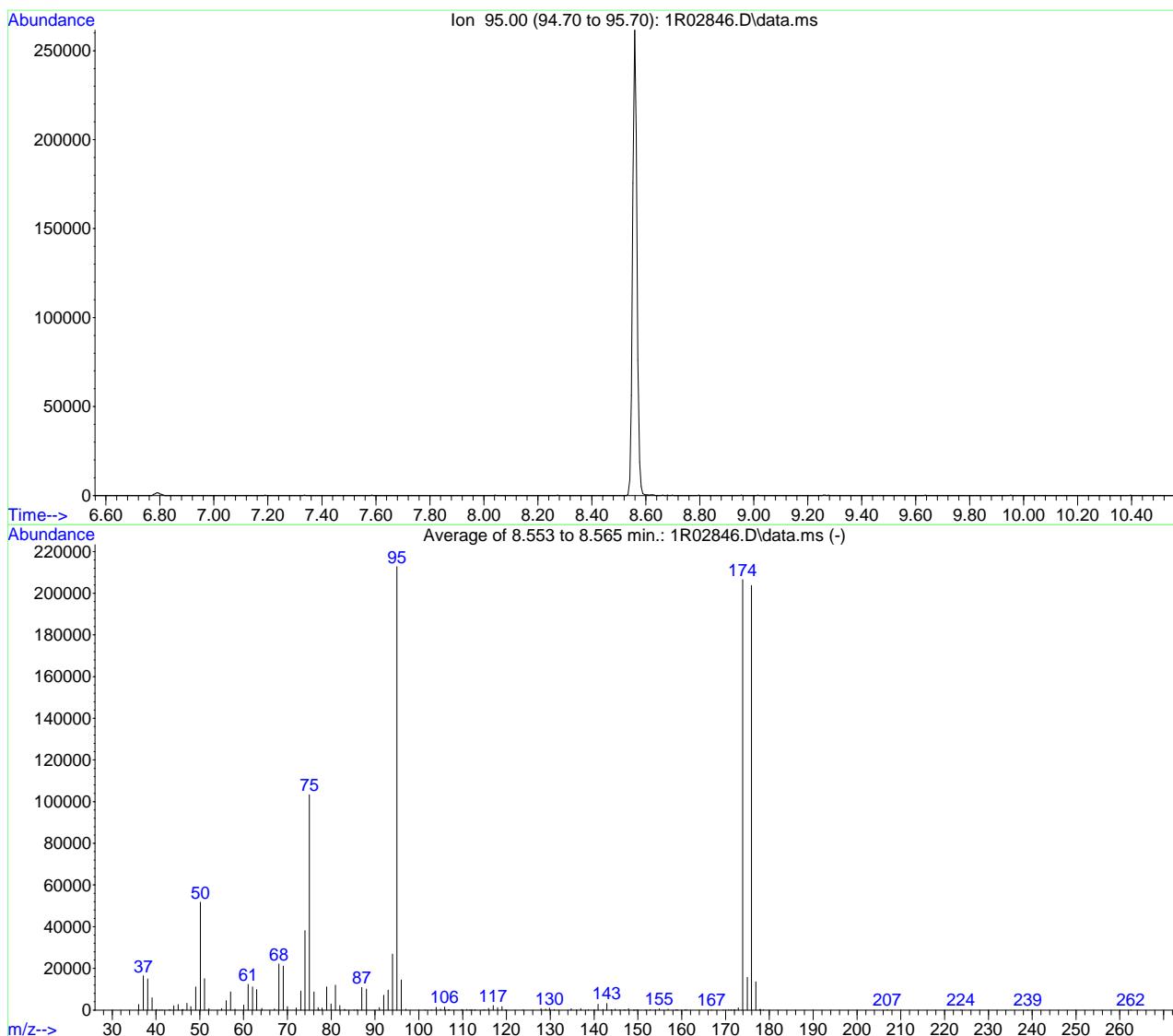
Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
137.80	72	147.90	593	156.95	534	172.90	1620
139.90	119	148.70	63	159.05	304	173.90	241536
140.95	3045	149.00	135	160.70	103	175.00	17164
141.60	67	149.85	310	160.95	238	175.90	233536
142.05	552	151.70	96	165.20	78	176.90	14108
142.90	3360	151.90	87	169.20	78	177.95	350
143.80	109	153.00	153	169.70	121	197.20	73
144.95	281	153.95	162	170.30	142	206.70	73
145.90	383	154.70	206	170.70	67	207.00	69
146.90	78	154.85	354	170.90	64	209.20	59
147.15	181	155.90	76	171.60	214	281.10	98

## SW-846 Method 8260

Data File : C:\msdchem\1\data\V1r91\1R02846.D Vial: 18  
 Acq On : 29 Mar 2023 4:59 pm Operator: PrashanS  
 Sample : bfb2 Inst : GCMSR  
 Misc : MS67262,V2R0091,5,,,,1 Multiplr: 1.00  
 MS Integration Params: rteint.p

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)  
 Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um



AutoFind: Scans 1261, 1262, 1263; Background Corrected with Scan 1255

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	24.3	51715	PASS
75	95	30	60	48.6	103344	PASS
95	95	100	100	100.0	212736	PASS
96	95	5	9	6.8	14480	PASS
173	174	0.00	2	0.5	1103	PASS
174	95	50	150	97.1	206571	PASS
175	174	5	9	7.6	15660	PASS
176	174	95	101	98.6	203776	PASS
177	176	5	9	6.6	13543	PASS

Average of 8.553 to 8.565 min.: 1R02846.D\data.ms

bfb2

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.05	2642	45.90	59	58.05	402	73.05	9074
37.10	16368	46.15	270	60.05	2434	74.00	38075
38.10	14893	47.05	3271	61.05	12360	75.00	103344
39.10	5925	47.95	1677	62.05	11067	76.05	8786
40.10	268	49.05	11081	63.00	9922	77.00	1184
40.80	59	50.10	51715	64.15	834	77.20	236
41.40	64	51.05	15048	67.05	553	77.90	963
43.05	195	52.05	694	68.05	22136	78.95	11107
43.30	63	54.95	726	69.05	21203	79.95	2913
44.05	2095	56.00	4589	70.00	1711	80.95	11908
45.05	2728	57.05	8765	72.00	1079	81.95	2230

Average of 8.553 to 8.565 min.: 1R02846.D\data.ms

bfb2

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
83.00	273	96.00	14480	111.65	163	123.80	63
83.20	118	97.05	439	112.10	147	127.90	741
85.20	59	101.40		62	112.70	108	128.95
85.85	230	102.90		85	112.90	90	129.85
86.95	10793	103.95		1305	114.60	63	130.90
88.00	10142	104.95		415	115.05	158	134.80
90.95	1230	105.90		1413	115.90	843	136.00
92.00	7094	106.70		63	116.95	2123	136.95
93.00	9668	106.95		235	117.95	1168	138.75
94.00	26891	109.80		340	118.95	1700	140.10
95.00	212736	110.95		264	119.80	72	140.90
							2826

Average of 8.553 to 8.565 min.: 1R02846.D\data.ms

bfb2

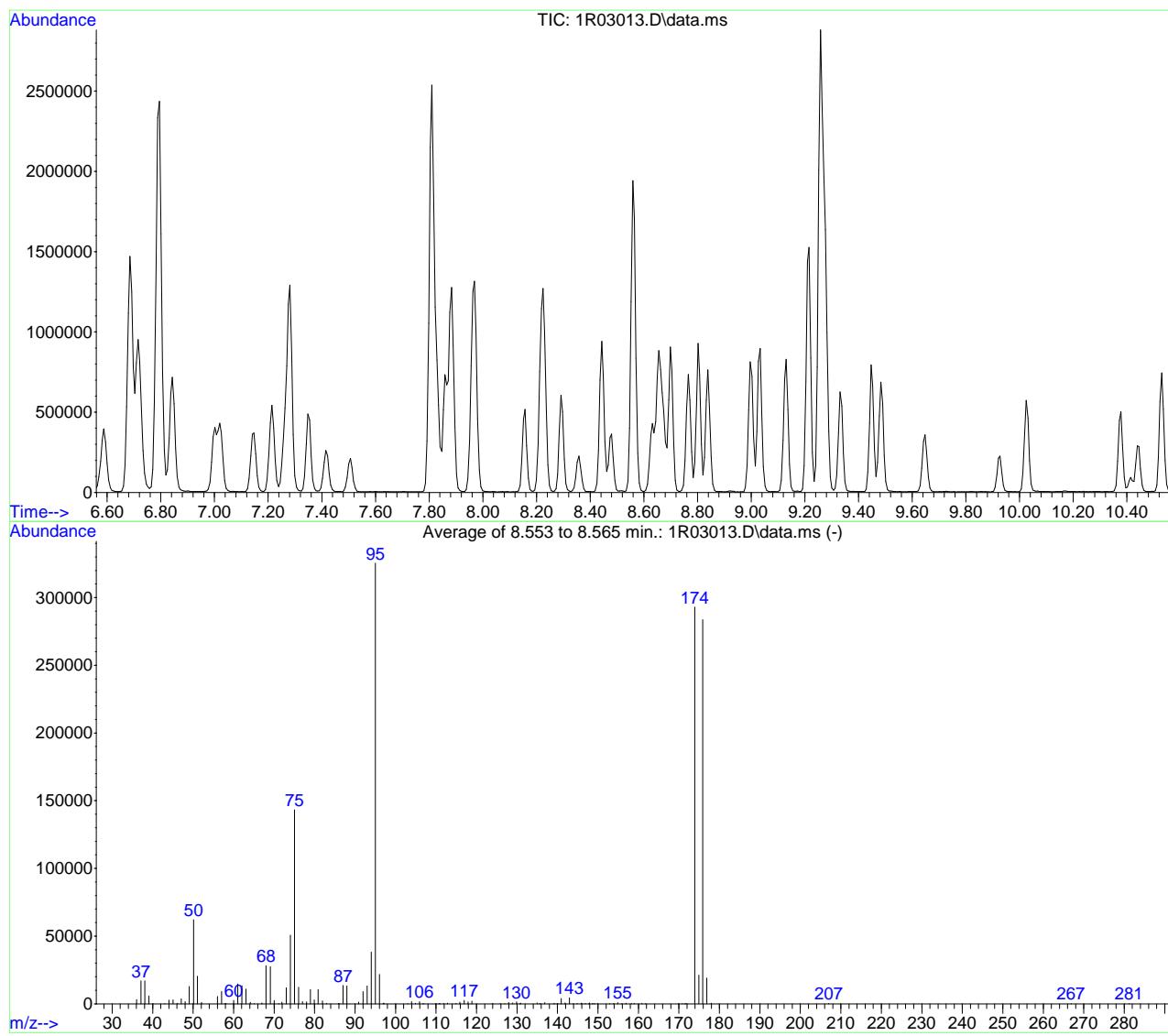
Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
142.05	259	151.80	70	170.70	95	178.00	291
142.90	3166	152.80	141	171.20	84	178.70	60
143.95	154	154.00	145	171.40	100	206.95	225
144.80	436	154.85	506	171.95	198	223.70	68
145.80	35	156.90	379	172.20	68	239.00	70
146.80	73	158.70	150	172.90	1103	262.50	69
147.10	58	159.00	132	173.90	206571		
147.85	723	160.75	206	174.95	15660		
148.95	6	161.15	149	175.90	203776		
149.50	69	166.80	58	176.90	13543		
149.90	233	169.50	82	177.80	261		

## SW-846 Method 8260

Data File : M:\MSVOA\GCMSR\VR0100\1R03013.D Vial: 2  
 Acq On : 4 Apr 2023 9:38 am Operator: nickw  
 Sample : bfb Inst : GCMSR  
 Misc : MS67909,V1R0100,5,,,1 Multiplr: 1.00  
 MS Integration Params: rteint.p

Method : C:\MSDCHEM\1\METHODS\M1R0091.M (RTE Integrator)  
 Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um



AutoFind: Scans 1261, 1262, 1263; Background Corrected with Scan 1254

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	19.1	62083	PASS
75	95	30	60	44.1	143333	PASS
95	95	100	100	100.0	325376	PASS
96	95	5	9	6.7	21747	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	150	90.1	293013	PASS
175	174	5	9	7.2	21152	PASS
176	174	95	101	96.8	283755	PASS
177	176	5	9	6.7	19014	PASS

Average of 8.553 to 8.565 min.: 1R03013.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.05	3060	49.05	12828	63.05	10893	75.05	143333
37.10	16964	50.10	62083	64.10	1105	76.05	12188
38.05	16939	51.05	20438	65.10	68	77.00	1604
39.10	5872	52.05	1035	67.00	685	78.00	1422
40.10	202	56.00	5337	68.00	28133	78.95	10500
43.00	151	57.00	9147	69.05	27416	79.95	3000
44.00	2753	57.80	83	70.05	2312	80.90	10565
45.05	2984	58.10	308	71.20	68	81.95	2249
46.15	308	60.00	2493	71.95	1300	82.95	232
47.05	3791	61.00	14655	73.00	11944	85.95	397
48.00	1601	62.05	13296	74.00	50715	86.20	149

Average of 8.553 to 8.565 min.: 1R03013.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
87.00	13500	103.95	1591	115.85	1303	130.90	452
88.00	13152	104.90	386	116.95	2556	134.95	702
90.90	1252	105.90	1709	117.90	1398	135.60	93
92.00	9066	106.80	126	118.85	2017	135.80	71
93.00	13115	107.00	199	124.00	279	136.85	867
94.00	38128	109.95	208	125.90	91	139.00	104
95.00	325376	110.75	222	127.00	67	139.30	62
96.00	21747	111.00	176	127.90	1120	139.90	193
96.95	457	111.90	356	128.95	636	140.10	89
97.20	230	112.85	451	129.90	1492	140.90	3905
103.00	191	114.95	397	130.70	66	141.90	603

Average of 8.553 to 8.565 min.: 1R03013.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
142.95	4407	149.95	191	170.80	140	281.20	68
143.85	187	152.90	114	171.40	162		
144.85	371	153.85	333	171.70	124		
145.85	452	154.90	876	172.00	119		
146.10	100	155.90	60	173.90	293013		
146.90	219	156.90	486	174.90	21152		
147.70	169	158.30	59	175.90	283755		
147.90	643	158.95	306	176.90	19014		
148.60	82	160.85	352	177.95	551		
148.90	142	161.70	65	207.00	208		
149.20	88	169.90	77	266.90	70		

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02815.D  
 Acq On : 28 Mar 2023 10:46 pm  
 Operator : PrashanS  
 Sample : IC0091-0.2  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Mar 30 09:05:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:12:11 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.485	65	524695	500.00	ug/L	0.00
5) pentafluorobenzene	5.024	168	631778	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.651	114	885203	50.00	ug/L	0.00
73) chlorobenzene-d5	7.810	117	816563	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.258	152	376952	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.012	113	273684	49.34	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 98.68%			
53) 1,2-dichloroethane-d4 (s)	5.298	65	236768	45.58	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery = 91.16%			
74) toluene-d8 (s)	6.794	98	1114103	44.51	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 89.02%			
98) 4-bromofluorobenzene (s)	8.565	174	328459	63.47	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 126.94%#			
<b>Target Compounds</b>						
				Qvalue		
7) dichlorodifluoromethane	1.642	85	861	0.21	ug/L	84
13) trichlorofluoromethane	2.536	101	1096	0.21	ug/L	66
17) 1,1-dichloroethene	3.035	96	729	0.23	ug/L	90
24) methyl tert butyl ether	3.728	73	1874	0.18	ug/L	57
27) di-isopropyl ether	4.136	45	2736	0.18	ug/L #	75
29) 1,1-dichloroethane	4.118	63	1444	0.21	ug/L	72
30) chloroprene	4.191	53	1205	0.19	ug/L	69
33) ethyl tert-butyl ether	4.440	59	2452	0.20	ug/L	92
36) cis-1,2-dichloroethene	4.610	96	944	0.23	ug/L #	59
37) propionitrile	4.647	54	2682	2.17	ug/L	79
40) bromochloromethane	4.805	128	460	0.23	ug/L #	28
55) 2,2,4-trimethylpentane	5.401	57	2571	0.22	ug/L	84
56) benzene	5.316	78	3576	0.21	ug/L #	1
57) tert-amyl methyl ether	5.407	73	2192	0.18	ug/L	77
59) 1,2-dichloroethane	5.359	62	1167	0.22	ug/L	84
61) trichloroethene	5.851	95	919	0.22	ug/L	82
62) 2-chloroethyl vinyl ether	6.466	63	2988	0.72	ug/L	83
64) methylcyclohexane	6.034	83	1559	0.24	ug/L #	67
70) cis-1,3-dichloropropene	6.594	75	1552	0.23	ug/L	83
71) 4-methyl-2-pentanone	6.685	58	2610	0.71	ug/L #	55
75) toluene	6.843	92	2487	0.23	ug/L #	77
77) trans-1,3-dichloropropene	7.001	75	957	0.17	ug/L #	25
78) 1,1,2-trichloroethane	7.147	83	703	0.23	ug/L #	80
79) tetrachloroethene	7.220	164	881	0.25	ug/L #	55
80) 2-hexanone	7.287	58	3063	0.91	ug/L #	71
81) 1,3-dichloropropane	7.263	76	1355	0.24	ug/L #	72
85) n-butyl ether	7.859	57	4019	0.21	ug/L	92
86) chlorobenzene	7.829	112	2360	0.21	ug/L #	31
88) ethylbenzene	7.890	91	4064	0.20	ug/L	89
89) m,p-xylene	7.963	106	2673	0.32	ug/L	89
90) o-xylene	8.224	91	3489	0.20	ug/L	87
92) butyl acrylate	8.157	55	1695	0.17	ug/L	95
95) bromoform	8.358	173	720	0.19	ug/L #	36
102) bromobenzene	8.662	156	697	0.19	ug/L #	77
104) 2-chlorotoluene	8.772	126	762	0.24	ug/L #	86
105) 4-chlorotoluene	8.839	91	2733	0.25	ug/L	75
106) 1,3,5-trimethylbenzene	8.808	105	2660	0.22	ug/L	93
108) 1,2,4-trimethylbenzene	9.033	105	2175	0.16	ug/L	92
109) sec-butylbenzene	9.131	105	2600	0.17	ug/L	86
110) p-isopropyltoluene	9.210	119	2970	0.20	ug/L	79

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02815.D  
 Acq On : 28 Mar 2023 10:46 pm  
 Operator : PrashanS  
 Sample : IC0091-0.2  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Mar 30 09:05:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:12:11 2023  
 Response via : Initial Calibration

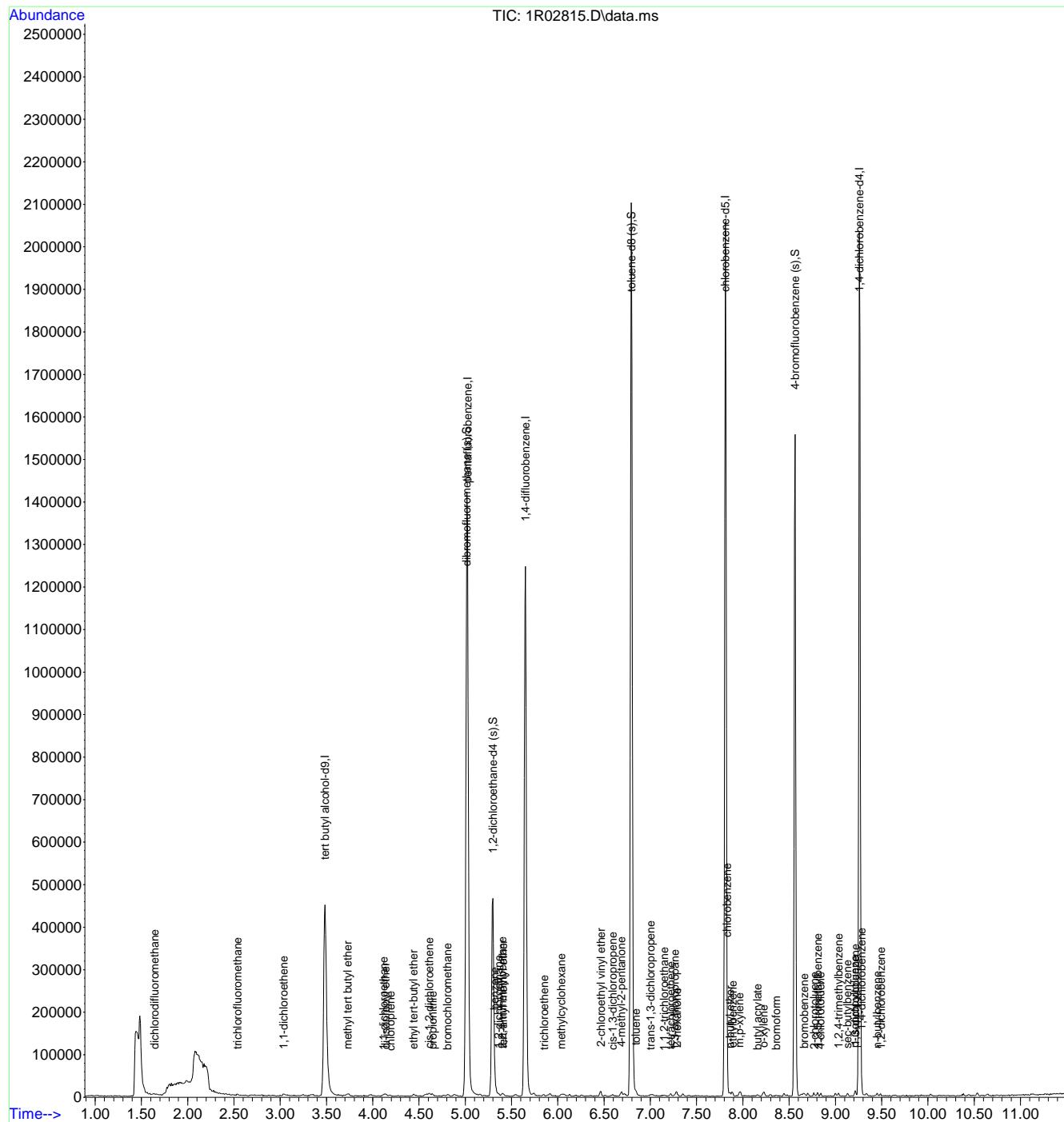
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
112) 1,3-dichlorobenzene	9.216	146	1659	0.19	ug/L	86
113) 1,4-dichlorobenzene	9.283	146	1867	0.22	ug/L	86
114) 1,2-dichlorobenzene	9.489	146	1404	0.18	ug/L	82
116) n-butylbenzene	9.453	92	1206	0.19	ug/L	88

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02815.D  
 Acq On : 28 Mar 2023 10:46 pm  
 Operator : PrashanS  
 Sample : IC0091-0.2  
 Misc : MS67262,V1R0091,5,,,.1  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Mar 30 09:05:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:12:11 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02817.D  
 Acq On : 28 Mar 2023 11:19 pm  
 Operator : PrashanS  
 Sample : IC0091-0.5  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 30 09:06:19 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:20:15 2023  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED  
 (compounds with "m" flag)

Kanya Veerawat  
 03/30/23 12:39

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) tert butyl alcohol-d9	3.491	65	518421	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	627743	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	870680	50.00	ug/L	0.00
73) chlorobenzene-d5	7.817	117	821774	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	371336	50.00	ug/L	0.00
System Monitoring Compounds						
44) dibromofluoromethane (s)	5.024	113	264721	48.35	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	96.70%		
53) 1,2-dichloroethane-d4 (s)	5.304	65	236206	48.37	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	96.74%		
74) toluene-d8 (s)	6.801	98	1096721	46.06	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	92.12%		
98) 4-bromofluorobenzene (s)	8.565	174	321812	55.63	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	111.26%		
Target Compounds						
				Qvalue		
2) ethanol	2.700	45	4952	56.75	ug/L	81
3) tertiary butyl alcohol	3.564	59	3469	2.92	ug/L	75
7) dichlorodifluoromethane	1.636	85	2561	0.61	ug/L	59
9) vinyl chloride	1.897	62	4107	0.64	ug/L	89
10) 1,3-butadiene	1.928	54	2574	0.52	ug/L	# 69
13) trichlorofluoromethane	2.542	101	2306	0.43	ug/L	84
14) ethyl ether	2.810	74	1119	0.52	ug/L	# 52
16) freon 113	3.047	151	1718	0.60	ug/L	# 73
17) 1,1-dichloroethene	3.035	96	1812	0.53	ug/L	# 56
18) acetone	3.065	58	1992	2.93	ug/L	# 54
21) carbon disulfide	3.260	76	5372	0.59	ug/L	92
22) methylene chloride	3.491	84	1764	0.48	ug/L	84
24) methyl tert butyl ether	3.728	73	5546	0.56	ug/L	87
25) trans-1,2-dichloroethene	3.747	96	1951	0.53	ug/L	86
26) hexane	3.984	56	1921	0.64	ug/L	# 67
27) di-isopropyl ether	4.142	45	7804	0.54	ug/L	95
28) 2-butanone	4.598	72	2156	2.38	ug/L	# 38
29) 1,1-dichloroethane	4.130	63	3740	0.53	ug/L	82
30) chloroprene	4.191	53	2867	0.47	ug/L	94
31) acrylonitrile	3.698	53	1253	0.43	ug/L	# 68
33) ethyl tert-butyl ether	4.452	59	5059	0.42	ug/L	94
35) 2,2-dichloropropane	4.623	77	2479	0.52	ug/L	74
36) cis-1,2-dichloroethene	4.617	96	1955	0.45	ug/L	# 75
37) propionitrile	4.653	54	5866	4.58	ug/L	91
40) bromochloromethane	4.817	128	826	0.39	ug/L	# 53
42) chloroform	4.896	83	4295	0.64	ug/L	72
43) tert-Butyl Formate	4.902	59	1680	0.40	ug/L	80
45) 1,1,1-trichloroethane	5.030	97	3192	0.56	ug/L	# 1
46) cyclohexane	5.103	84	3243	0.57	ug/L	# 25
48) 1,1-dichloropropene	5.164	75	2482	0.49	ug/L	79
49) carbon tetrachloride	5.170	117	2275	0.45	ug/L	82
55) 2,2,4-trimethylpentane	5.407	57	5470	0.45	ug/L	84
56) benzene	5.322	78	7904m	0.46	ug/L	
57) tert-amyl methyl ether	5.426	73	5569	0.49	ug/L	90
58) heptane	5.547	57	1214	0.47	ug/L	# 70
59) 1,2-dichloroethane	5.371	62	2957	0.54	ug/L	98
60) ethyl acrylate	5.894	55	4310	0.49	ug/L	88
61) trichloroethene	5.852	95	1650	0.38	ug/L	# 69
62) 2-chloroethyl vinyl ether	6.466	63	7689	2.19	ug/L	78
64) methylcyclohexane	6.052	83	2781	0.39	ug/L	# 84

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02817.D  
 Acq On : 28 Mar 2023 11:19 pm  
 Operator : PrashanS  
 Sample : IC0091-0.5  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 30 09:06:19 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:20:15 2023  
 Response via : Initial Calibration

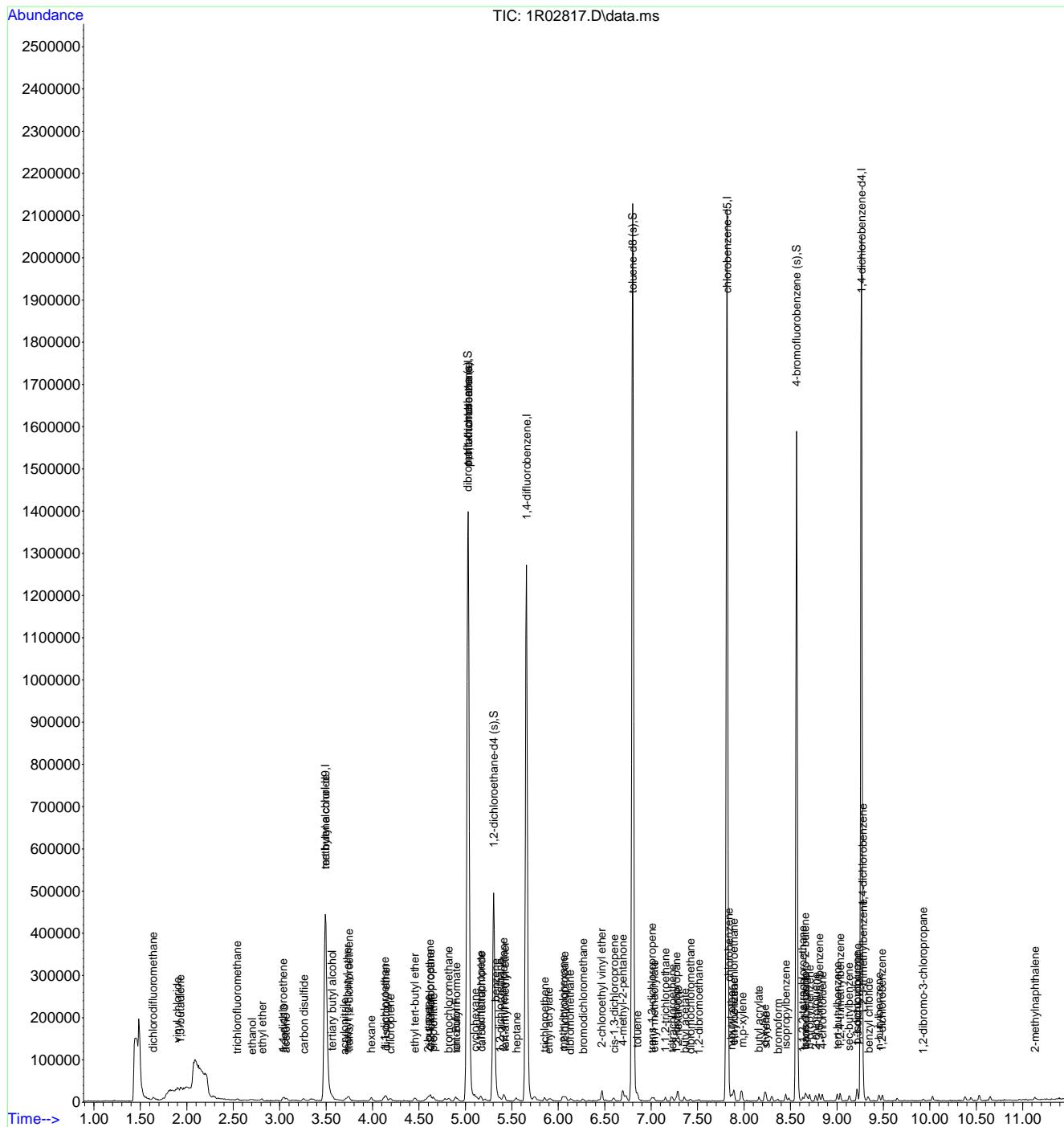
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
65) 1,2-dichloropropane	6.064	63	2029	0.44	ug/L	80
66) dibromomethane	6.125	93	1622	0.58	ug/L #	45
67) bromodichloromethane	6.265	83	2619	0.50	ug/L	71
70) cis-1,3-dichloropropene	6.600	75	2710	0.38	ug/L	82
71) 4-methyl-2-pentanone	6.691	58	5571	1.63	ug/L #	64
75) toluene	6.849	92	4042	0.34	ug/L	96
76) ethyl methacrylate	7.032	69	2480	0.42	ug/L	90
77) trans-1,3-dichloropropene	7.007	75	2679	0.51	ug/L	86
78) 1,1,2-trichloroethane	7.153	83	1530	0.46	ug/L	84
79) tetrachloroethene	7.226	164	2255	0.57	ug/L #	76
80) 2-hexanone	7.287	58	5671	1.56	ug/L #	85
81) 1,3-dichloropropane	7.269	76	2442	0.39	ug/L	88
82) butyl acetate	7.360	56	2354	0.55	ug/L #	41
83) dibromochloromethane	7.421	129	1687	0.41	ug/L	76
84) 1,2-dibromoethane	7.512	107	1809	0.47	ug/L	96
85) n-butyl ether	7.865	57	8181	0.42	ug/L	86
86) chlorobenzene	7.835	112	5342	0.46	ug/L	76
87) 1,1,1,2-tetrachloroethane	7.890	131	1655	0.37	ug/L #	80
88) ethylbenzene	7.890	91	9203	0.45	ug/L	87
89) m,p-xylene	7.981	106	6760	0.89	ug/L #	62
90) o-xylene	8.224	91	5750	0.33	ug/L	85
91) styrene	8.236	104	5106	0.32	ug/L	80
92) butyl acrylate	8.157	55	3922	0.42	ug/L	95
94) isopropylbenzene	8.449	105	6962	0.33	ug/L	96
95) bromoform	8.364	173	1370	0.37	ug/L	86
99) 1,1,2,2-tetrachloroethane	8.632	83	2771	0.68	ug/L	96
100) trans-1,4-dichloro-2-b...	8.656	53	968	0.60	ug/L #	52
102) bromobenzene	8.662	156	2294	0.66	ug/L #	77
103) n-propylbenzene	8.705	91	9575	0.61	ug/L	94
104) 2-chlorotoluene	8.772	126	1499	0.44	ug/L	92
105) 4-chlorotoluene	8.845	91	6080	0.50	ug/L	83
106) 1,3,5-trimethylbenzene	8.808	105	5720	0.46	ug/L	89
107) tert-butylbenzene	9.003	119	5306	0.57	ug/L	78
108) 1,2,4-trimethylbenzene	9.033	105	6270	0.52	ug/L	100
109) sec-butylbenzene	9.131	105	7615	0.55	ug/L	96
110) p-isopropyltoluene	9.216	119	6012	0.41	ug/L	96
111) 1,2,3-trimethylbenzene	9.283	105	6700	0.47	ug/L	90
112) 1,3-dichlorobenzene	9.222	146	4053	0.49	ug/L	90
113) 1,4-dichlorobenzene	9.277	146	4833	0.55	ug/L	73
114) 1,2-dichlorobenzene	9.483	146	3250	0.44	ug/L	84
115) benzyl chloride	9.344	91	4893	0.49	ug/L	98
116) n-butylbenzene	9.453	92	3442	0.57	ug/L	80
118) 1,2-dibromo-3-chloropr...	9.928	157	756	0.68	ug/L #	45
124) 2-methylnaphthalene	11.132	142	1843	0.40	ug/L #	67

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02817.D  
 Acq On : 28 Mar 2023 11:19 pm  
 Operator : PrashanS  
 Sample : IC0091-0.5  
 Misc : MS67262,V1R0091,5,,,.1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 30 09:06:19 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:20:15 2023  
 Response via : Initial Calibration



# Manual Integration Approval Summary

Page 1 of 1

Sample Number: V1R91-IC0091  
Lab FileID: 1R02817.D  
Injection Time: 03/28/23 23:19

Method: SW846 8260D  
Analyst approved: 03/30/23 10:37 MoHui Huang  
Supervisor approved: 03/30/23 12:39 Kanya Veerawat

Parameter	CAS	Sig#	R.T. (min.)	Reason
Benzene	71-43-2		5.32	Missed peak

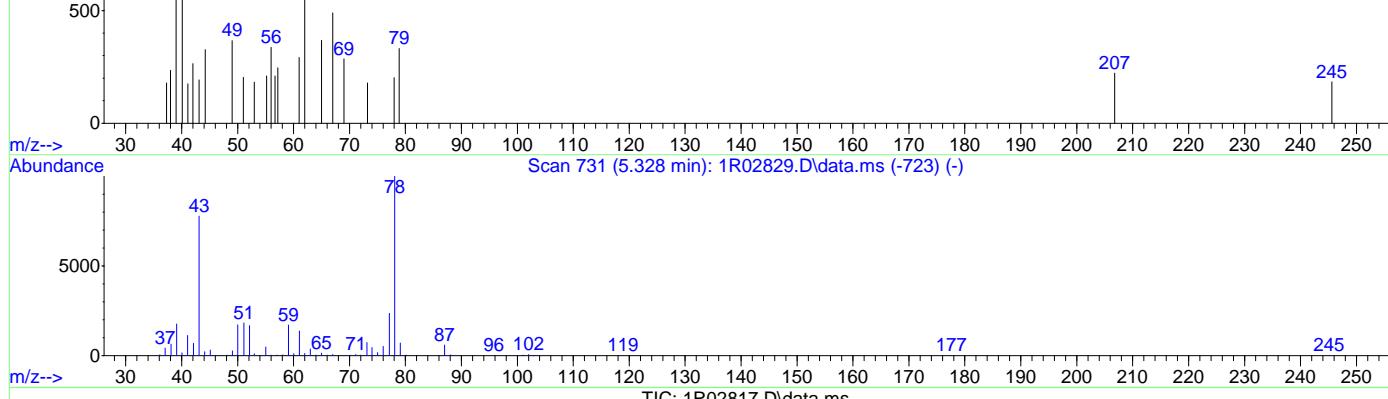
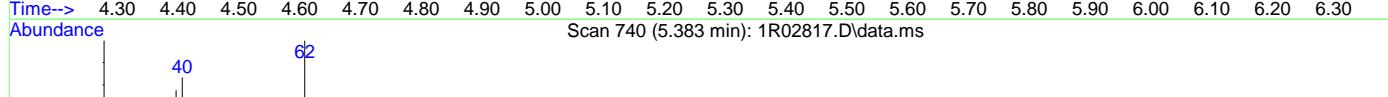
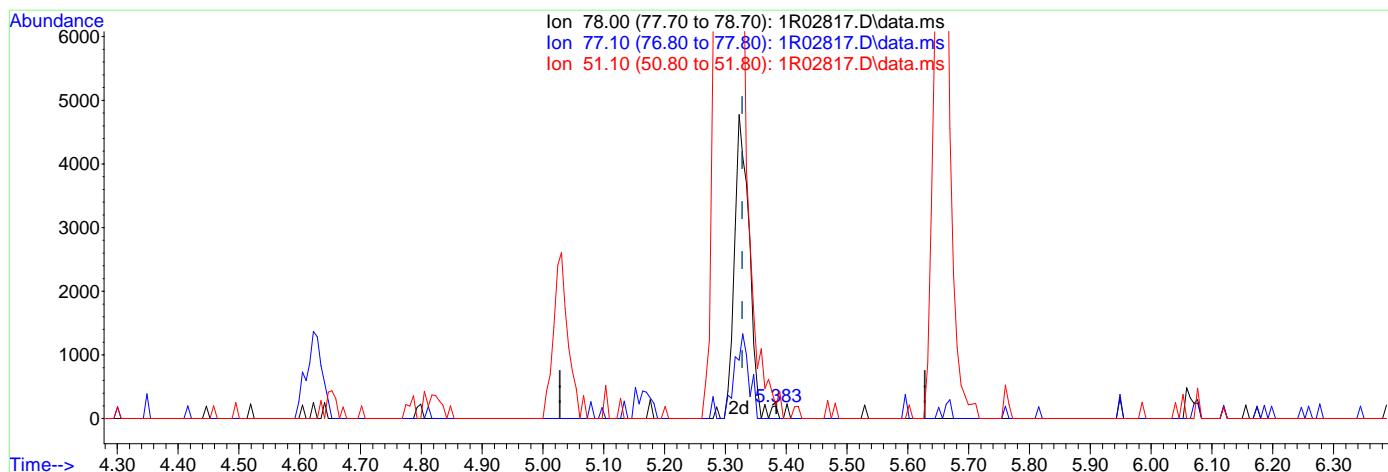
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7

## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\VR0091-raw\  
 Data File : 1R02817.D  
 Acq On : 28 Mar 2023 11:19 pm  
 Operator : PrashanS  
 Sample : IC0091-0.5  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 28 23:31:13 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\MR0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Jan 11 14:18:02 2023  
 Response via : Initial Calibration



(56) benzene

5.383min (+0.055) 0.01ug/L

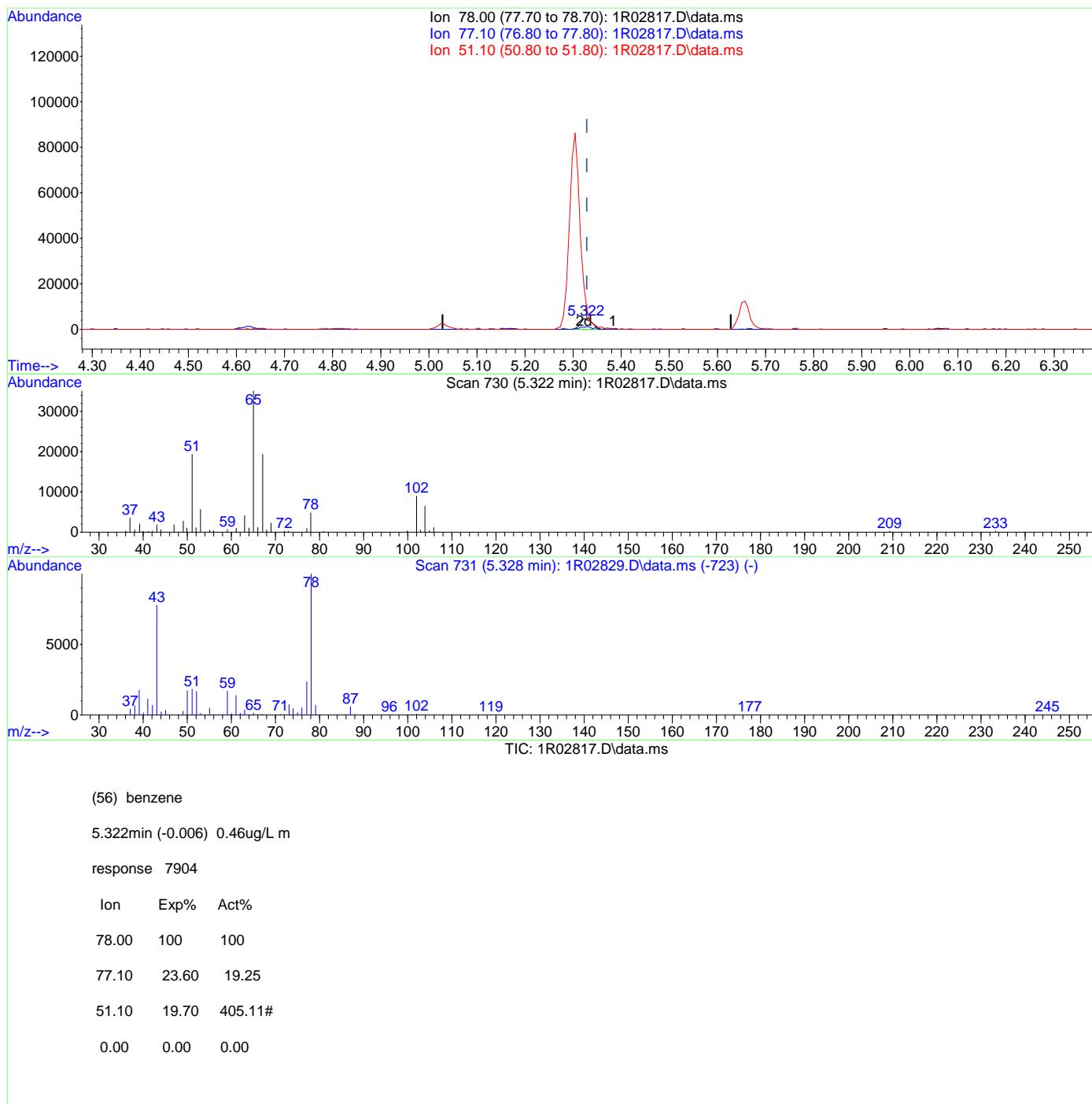
response 141

Ion	Exp%	Act%
78.00	100	100
77.10	23.30	0.00
51.10	21.00	0.00
0.00	0.00	0.00

## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02817.D  
 Acq On : 28 Mar 2023 11:19 pm  
 Operator : PrashanS  
 Sample : IC0091-0.5  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 30 09:06:19 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:20:15 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 29 12:43:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:26:40 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) tert butyl alcohol-d9	3.491	65	478994	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	610388	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	848858	50.00	ug/L	0.00
73) chlorobenzene-d5	7.817	117	819932	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	373683	50.00	ug/L	0.00
System Monitoring Compounds						
44) dibromofluoromethane (s)	5.024	113	261498	49.67	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 99.34%			
53) 1,2-dichloroethane-d4 (s)	5.304	65	234709	49.84	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery = 99.68%			
74) toluene-d8 (s)	6.801	98	1088777	47.07	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 94.14%			
98) 4-bromofluorobenzene (s)	8.565	174	311507	51.58	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 103.16%			
Target Compounds						
				Qvalue		
2) ethanol	2.682	45	8307	96.52	ug/L	89
3) tertiary butyl alcohol	3.564	59	6699	5.63	ug/L	96
4) 1,4-dioxane	6.113	88	2572	32.01	ug/L #	56
7) dichlorodifluoromethane	1.636	85	4339	0.99	ug/L	92
8) chloromethane	1.794	50	10408	1.66	ug/L	97
9) vinyl chloride	1.891	62	7819	1.10	ug/L	78
10) 1,3-butadiene	1.934	54	6392	1.30	ug/L #	83
12) chloroethane	2.299	64	4125	1.25	ug/L	82
13) trichlorofluoromethane	2.536	101	5613	1.14	ug/L	94
14) ethyl ether	2.798	74	2248	1.05	ug/L #	64
16) freon 113	3.041	151	3335	1.09	ug/L #	76
17) 1,1-dichloroethene	3.041	96	3645	1.08	ug/L #	77
18) acetone	3.071	58	4054	4.97	ug/L #	53
21) carbon disulfide	3.248	76	11667	1.21	ug/L	95
22) methylene chloride	3.497	84	3496	1.00	ug/L	98
23) methyl acetate	3.357	74	1248	1.27	ug/L #	85
24) methyl tert butyl ether	3.728	73	12104	1.21	ug/L	96
25) trans-1,2-dichloroethene	3.747	96	4221	1.15	ug/L	80
26) hexane	3.984	56	3595	1.08	ug/L #	79
27) di-isopropyl ether	4.142	45	16265	1.13	ug/L	98
28) 2-butanone	4.586	72	3996	4.14	ug/L #	47
29) 1,1-dichloroethane	4.136	63	7278	1.04	ug/L	98
30) chloroprene	4.185	53	7119	1.23	ug/L	77
31) acrylonitrile	3.704	53	2588	0.98	ug/L #	61
32) vinyl acetate	4.118	86	1125	1.18	ug/L #	1
33) ethyl tert-butyl ether	4.452	59	12548	1.12	ug/L	92
34) ethyl acetate	4.617	45	1541	1.18	ug/L #	16
35) 2,2-dichloropropane	4.629	77	5569	1.17	ug/L	81
36) cis-1,2-dichloroethene	4.617	96	4601	1.12	ug/L #	71
37) propionitrile	4.647	54	13815	11.41	ug/L	69
38) methyl acrylate	4.659	85	1094	1.07	ug/L #	19
39) methacrylonitrile	4.775	67	3000	1.12	ug/L #	57
40) bromochloromethane	4.811	128	2033	1.06	ug/L #	77
41) tetrahydrofuran	4.829	42	3989	1.26	ug/L	77
42) chloroform	4.884	83	7401	1.00	ug/L	87
43) tert-Butyl Formate	4.902	59	3635	0.99	ug/L	88
45) 1,1,1-trichloroethane	5.036	97	6610	1.12	ug/L #	1
46) cyclohexane	5.103	84	5699	0.97	ug/L #	67
48) 1,1-dichloropropene	5.158	75	5360	1.10	ug/L	80
49) carbon tetrachloride	5.170	117	5178	1.11	ug/L	97

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 29 12:43:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:26:40 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) tert-amyl alcohol	5.304	73	3043	5.36	ug/L	#
51) isopropyl acetate	5.340	87	1434	1.13	ug/L	#
54) n-butyl alcohol	5.736	56	14508	45.95	ug/L	95
55) 2,2,4-trimethylpentane	5.414	57	13111	1.15	ug/L	81
56) benzene	5.328	78	16586m	1.02	ug/L	
57) tert-amyl methyl ether	5.414	73	11878	1.08	ug/L	89
58) heptane	5.541	57	2923	1.20	ug/L	#
59) 1,2-dichloroethane	5.371	62	5167	0.94	ug/L	82
60) ethyl acrylate	5.894	55	9052	1.07	ug/L	98
61) trichloroethene	5.852	95	4497	1.16	ug/L	89
62) 2-chloroethyl vinyl ether	6.472	63	17439	5.32	ug/L	96
63) methyl methacrylate	6.083	100	1494	1.14	ug/L	#
64) methylcyclohexane	6.046	83	7094	1.11	ug/L	89
65) 1,2-dichloropropane	6.064	63	5134	1.22	ug/L	97
66) dibromomethane	6.131	93	2926	1.00	ug/L	83
67) bromodichloromethane	6.271	83	5503	1.08	ug/L	91
69) epichlorohydrin	6.515	57	3816	4.36	ug/L	72
70) cis-1,3-dichloropropene	6.594	75	6374	0.99	ug/L	93
71) 4-methyl-2-pentanone	6.691	58	11579	3.71	ug/L	98
75) toluene	6.849	92	10938	1.04	ug/L	#
76) ethyl methacrylate	7.026	69	5558	1.02	ug/L	88
77) trans-1,3-dichloropropene	7.007	75	5601	1.07	ug/L	88
78) 1,1,2-trichloroethane	7.160	83	3336	1.03	ug/L	#
79) tetrachloroethene	7.226	164	4845	1.17	ug/L	86
80) 2-hexanone	7.287	58	13068	3.89	ug/L	92
81) 1,3-dichloropropane	7.269	76	6058	1.05	ug/L	89
82) butyl acetate	7.360	56	4367	0.98	ug/L	94
83) dibromochloromethane	7.421	129	4490	1.21	ug/L	93
84) 1,2-dibromoethane	7.506	107	4375	1.18	ug/L	91
85) n-butyl ether	7.865	57	19898	1.08	ug/L	94
86) chlorobenzene	7.835	112	11897	1.06	ug/L	96
87) 1,1,1,2-tetrachloroethane	7.883	131	3992	1.03	ug/L	95
88) ethylbenzene	7.890	91	19173	0.97	ug/L	87
89) m,p-xylene	7.969	106	15648	2.14	ug/L	#
90) o-xylene	8.224	91	14547	0.94	ug/L	91
91) styrene	8.236	104	11326	0.86	ug/L	93
92) butyl acrylate	8.163	55	11328	1.29	ug/L	95
93) n-amyl acetate	8.297	70	3496	0.76	ug/L	86
94) isopropylbenzene	8.449	105	17048	0.97	ug/L	99
95) bromoform	8.364	173	2881	0.85	ug/L	91
96) cis-1,4-dichloro-2-butene	8.486	88	2245	0.72	ug/L	97
99) 1,1,2,2-tetrachloroethane	8.638	83	6702	1.39	ug/L	84
100) trans-1,4-dichloro-2-b...	8.650	53	1910	1.07	ug/L	#
101) 1,2,3-trichloropropane	8.674	110	1769	1.40	ug/L	#
102) bromobenzene	8.668	156	5686	1.47	ug/L	#
103) n-propylbenzene	8.705	91	21105	1.20	ug/L	92
104) 2-chlorotoluene	8.772	126	4492m	1.35	ug/L	
105) 4-chlorotoluene	8.845	91	11589	0.95	ug/L	99
106) 1,3,5-trimethylbenzene	8.808	105	13768	1.12	ug/L	93
107) tert-butylbenzene	9.009	119	12641m	1.26	ug/L	
108) 1,2,4-trimethylbenzene	9.039	105	14740	1.20	ug/L	89
109) sec-butylbenzene	9.137	105	17021	1.19	ug/L	94
110) p-isopropyltoluene	9.216	119	14494	1.05	ug/L	91
111) 1,2,3-trimethylbenzene	9.283	105	14902	1.07	ug/L	95
112) 1,3-dichlorobenzene	9.216	146	8257	1.00	ug/L	94
113) 1,4-dichlorobenzene	9.277	146	8297	0.91	ug/L	83
114) 1,2-dichlorobenzene	9.489	146	8195	1.16	ug/L	85
115) benzyl chloride	9.337	91	10988	1.10	ug/L	77

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 29 12:43:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:26:40 2023  
 Response via : Initial Calibration

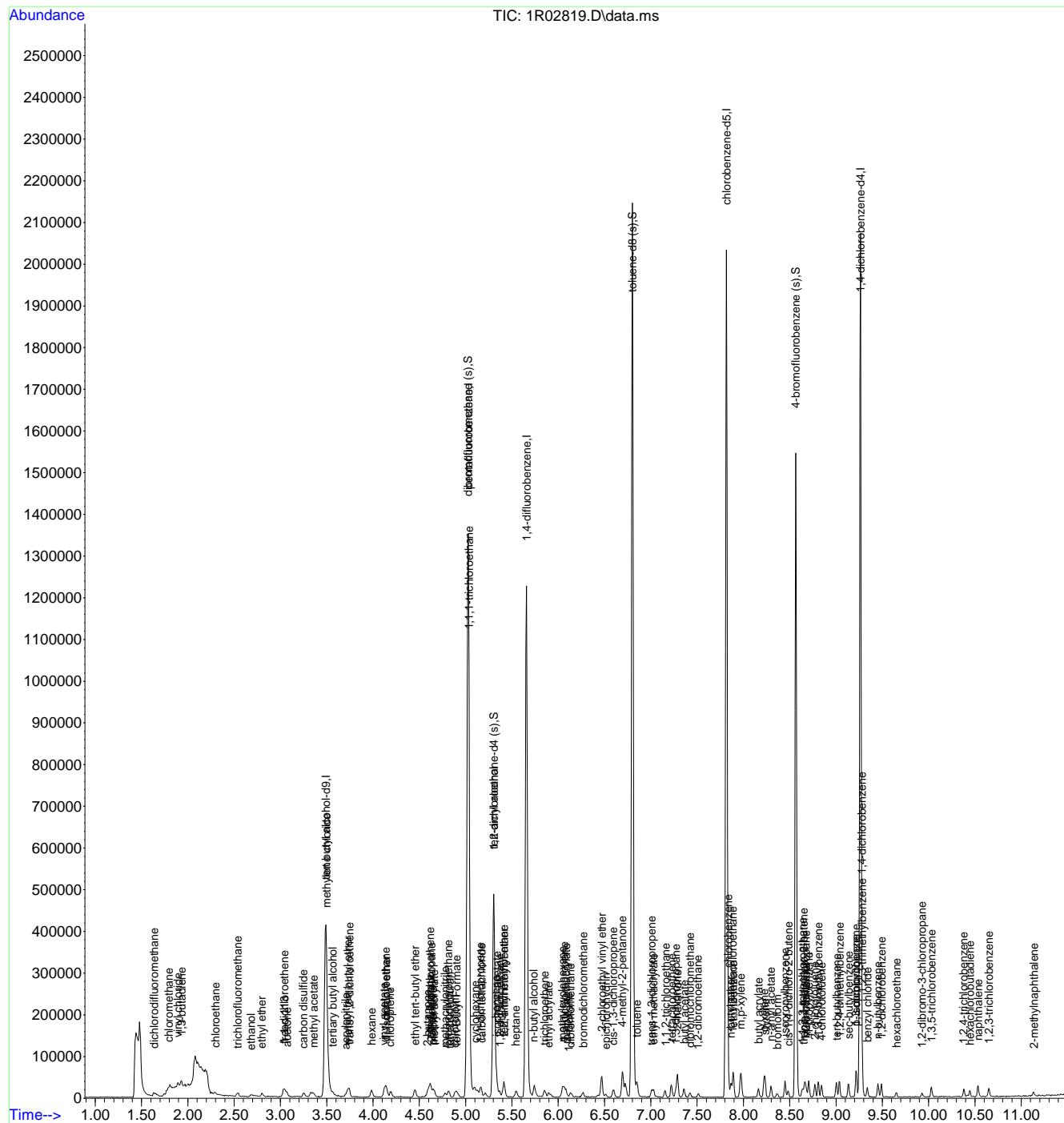
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
116) n-butylbenzene	9.453	92	6554	1.03	ug/L	82
117) hexachloroethane	9.654	201	1071	0.92	ug/L	71
118) 1,2-dibromo-3-chloropr...	9.928	157	1528	1.15	ug/L	91
119) 1,3,5-trichlorobenzene	10.031	180	5175	1.22	ug/L	96
120) 1,2,4-trichlorobenzene	10.378	180	4315	1.09	ug/L	94
121) hexachlorobutadiene	10.445	225	1747	1.41	ug/L #	55
122) naphthalene	10.536	128	15257	1.38	ug/L	96
123) 1,2,3-trichlorobenzene	10.651	180	4886	1.35	ug/L #	67
124) 2-methylnaphthalene	11.138	142	3508	0.58	ug/L #	70

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 29 12:43:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:26:40 2023  
 Response via : Initial Calibration



# Manual Integration Approval Summary

Page 1 of 1

Sample Number: V1R91-IC0091  
Lab FileID: 1R02819.D  
Injection Time: 03/28/23 23:52

Method: SW846 8260D  
Analyst approved: 03/30/23 10:37 MoHui Huang  
Supervisor approved: 03/30/23 12:39 Kanya Veerawat

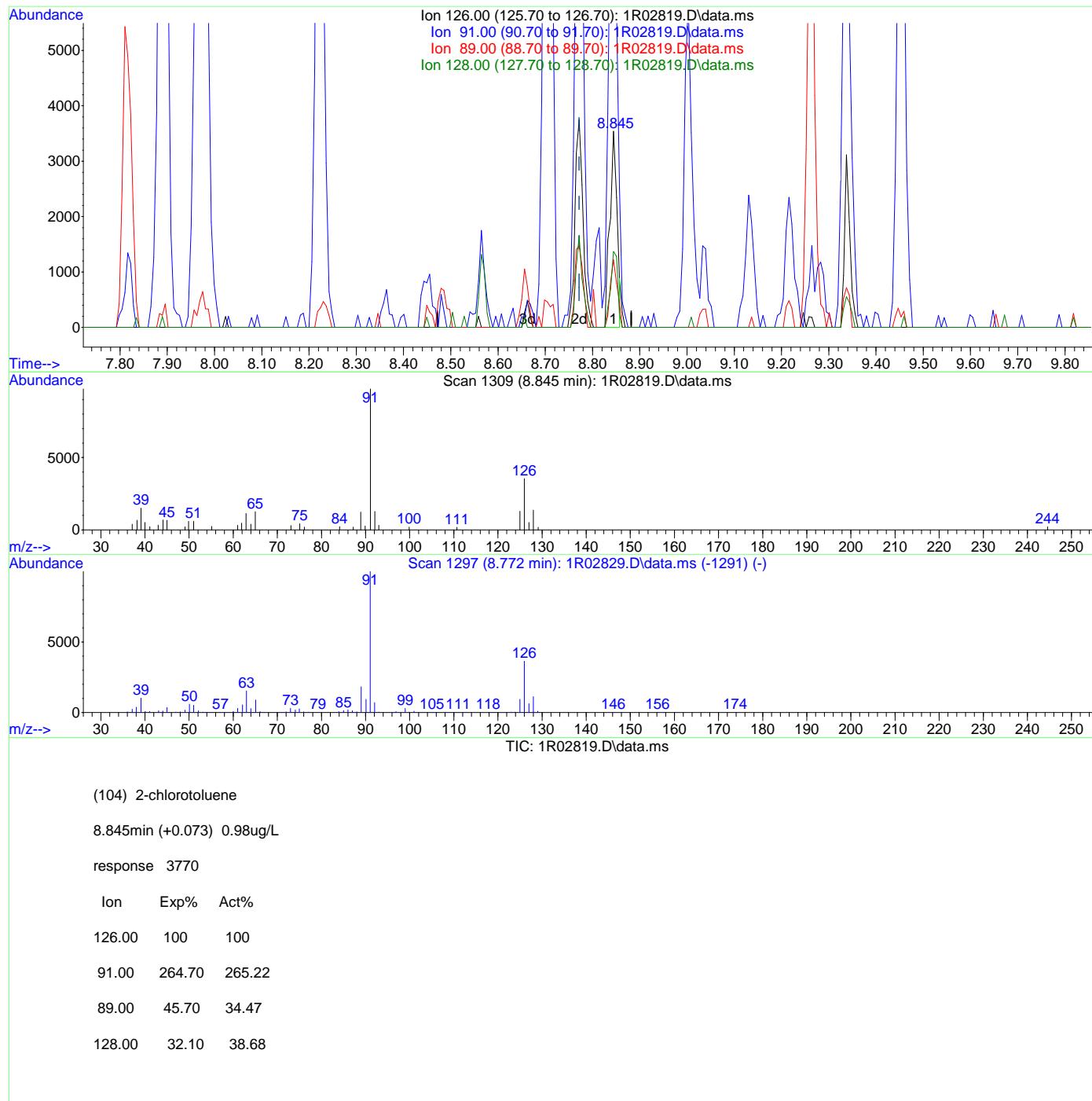
Parameter	CAS	Sig#	R.T. (min.)	Reason
Benzene	71-43-2		5.33	Missed peak
o-Chlorotoluene	95-49-8		8.77	Missed peak
tert-Butylbenzene	98-06-6		9.01	Missed peak

7.6.3.1  
7

## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\VR0091-raw\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

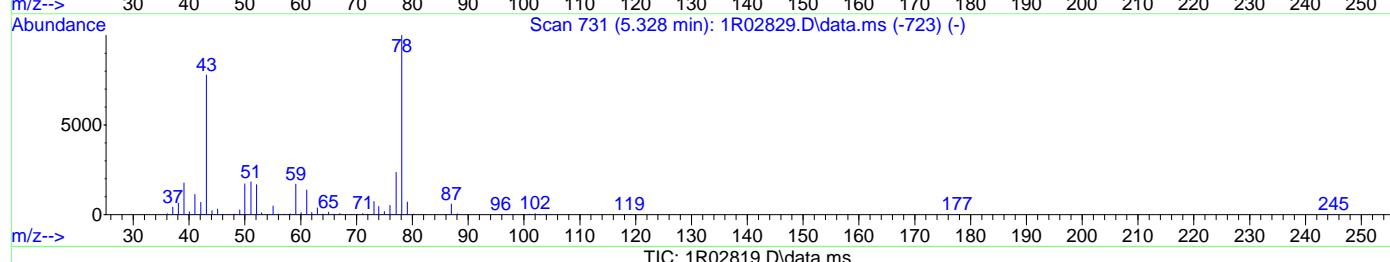
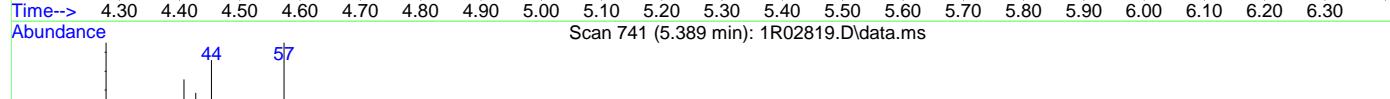
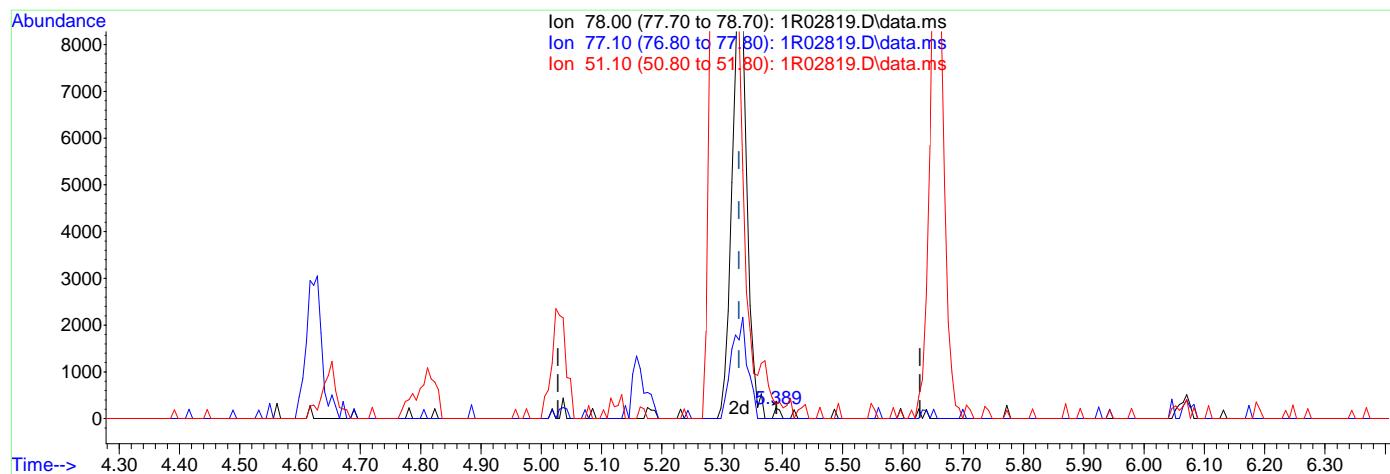
Quant Time: Mar 29 00:04:10 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\MR0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Jan 11 14:18:02 2023  
 Response via : Initial Calibration



## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\VR0091-raw\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 29 00:04:10 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\MR0091.M  
 Quant Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Jan 11 14:18:02 2023  
 Response via : Initial Calibration



(56) benzene

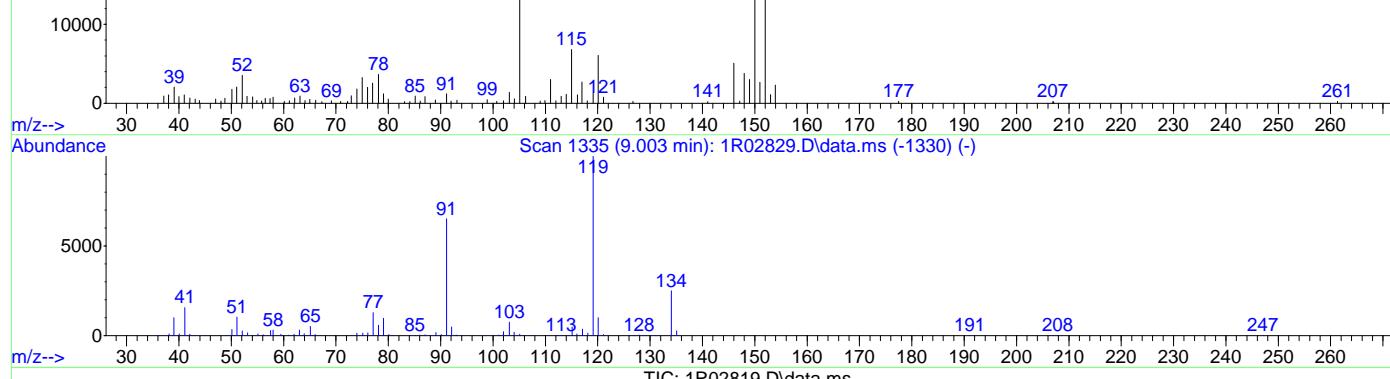
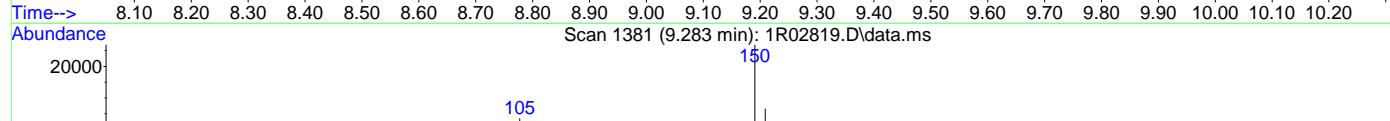
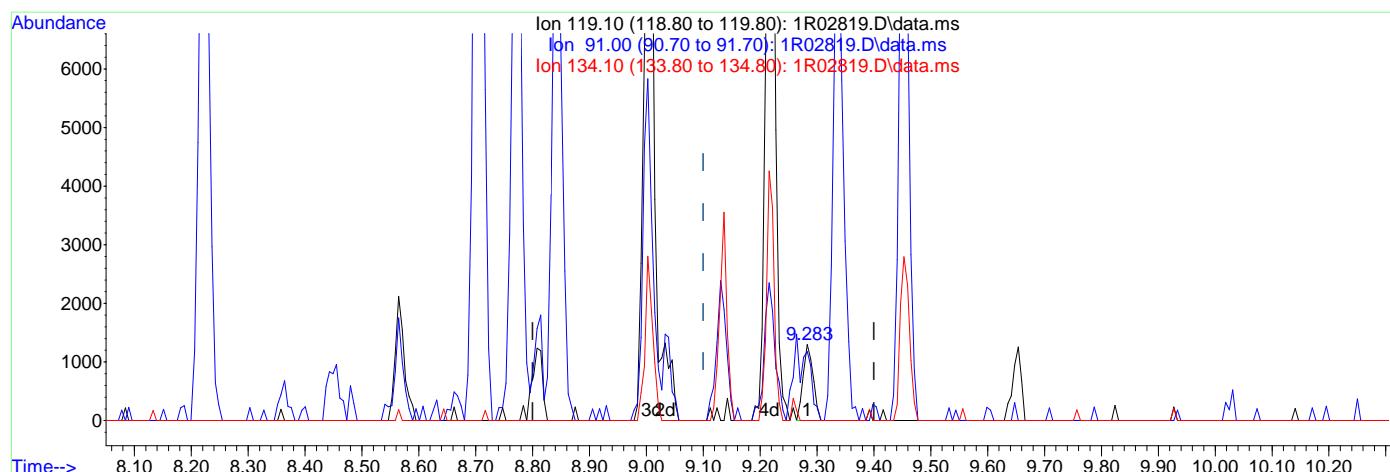
5.389min (+0.061) 0.01ug/L

response 148

Ion	Exp%	Act%
78.00	100	100
77.10	23.30	0.00
51.10	21.00	12.95
0.00	0.00	0.00

## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\VR0091-raw\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,,1  
 ALS Vial : 4 Sample Multiplier: 1  
 Quant Time: Mar 30 10:35:59 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Thu Mar 30 09:13:25 2023  
 Response via : Initial Calibration



(107) tert-butylbenzene

9.283min (+0.183) 0.14ug/L

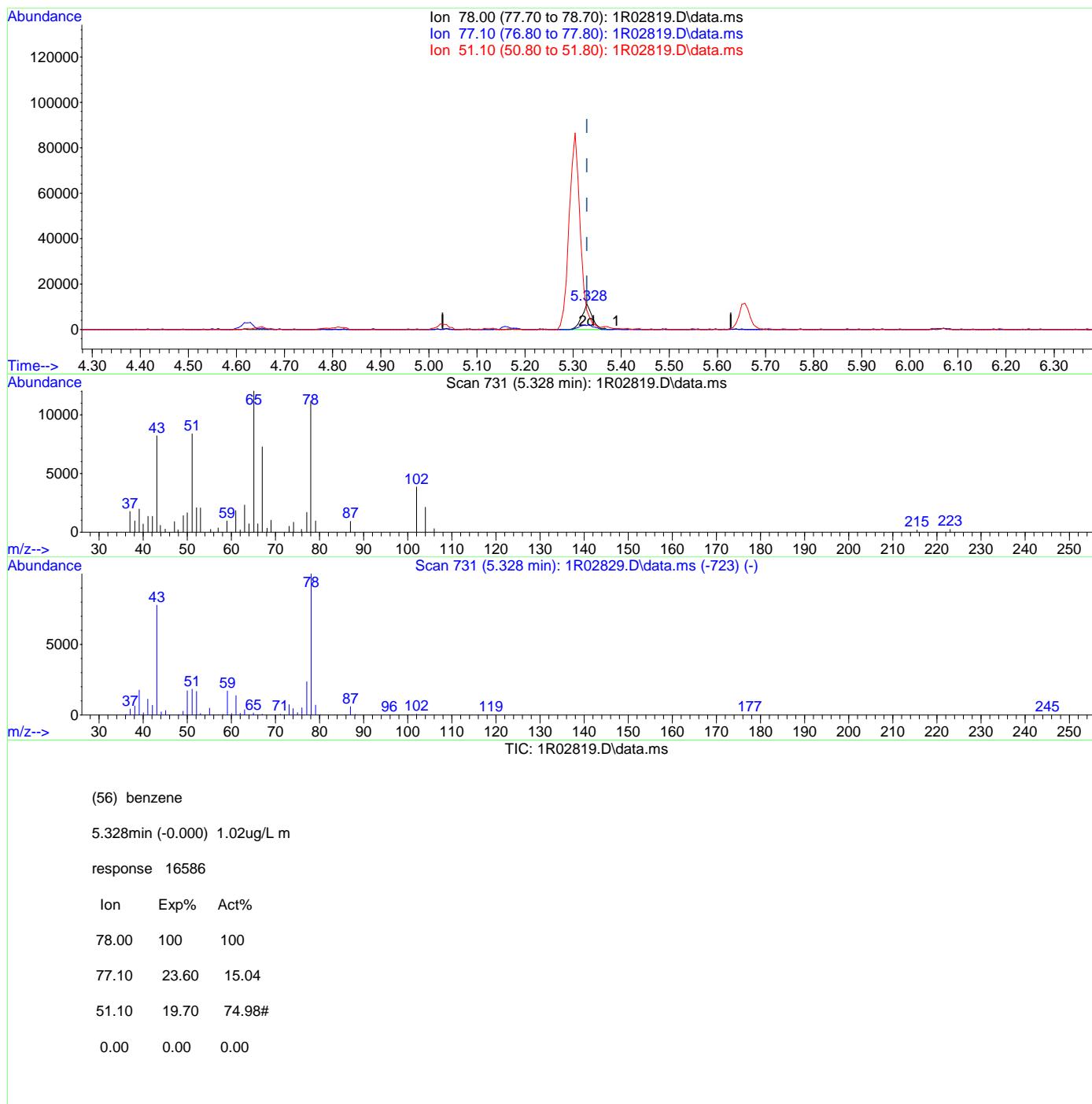
response 1601

Ion	Exp%	Act%
119.10	100	100
91.00	65.10	91.05
134.10	25.00	0.00
0.00	0.00	0.00

## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

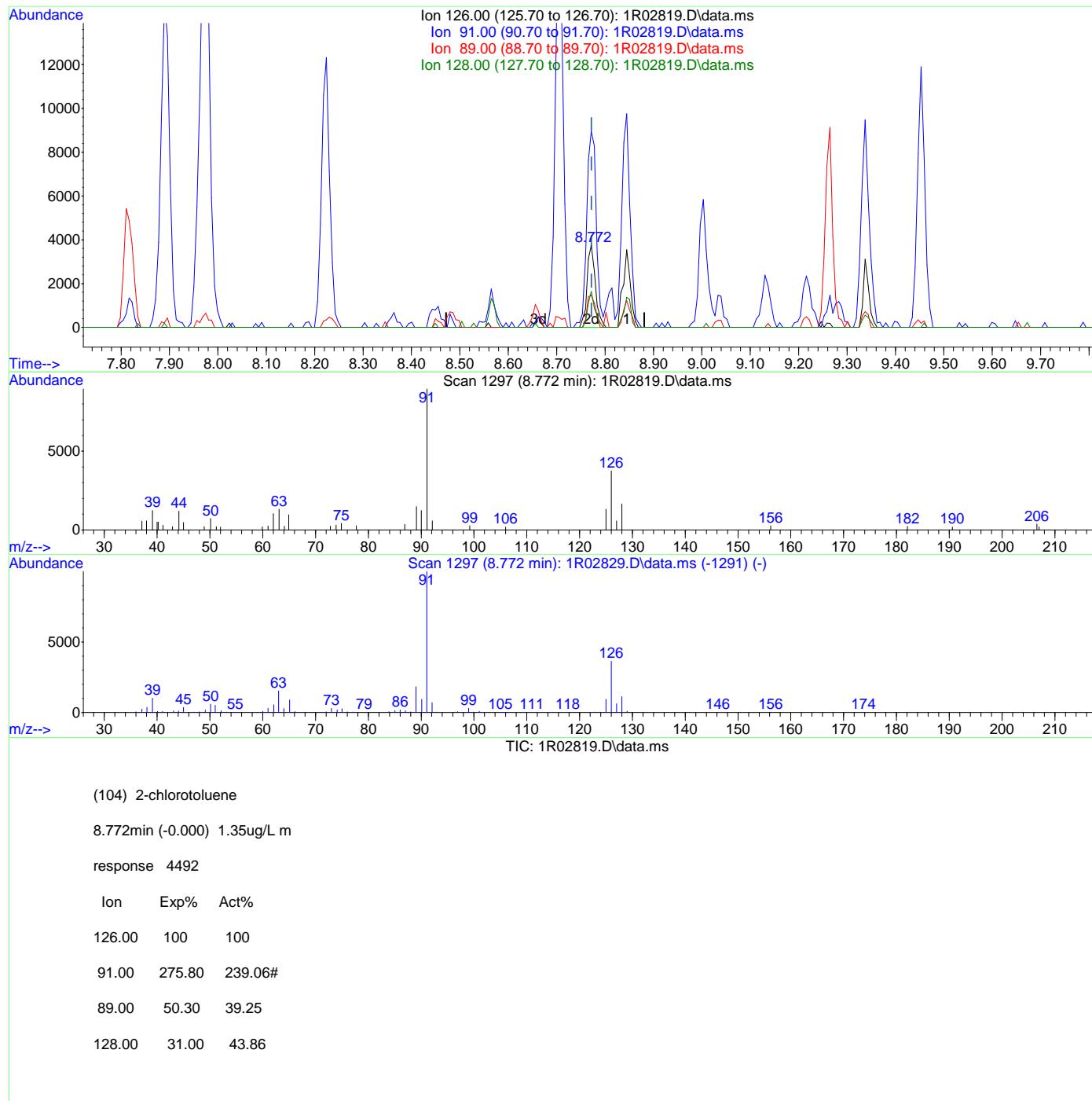
Quant Time: Mar 29 12:43:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:26:40 2023  
 Response via : Initial Calibration



## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

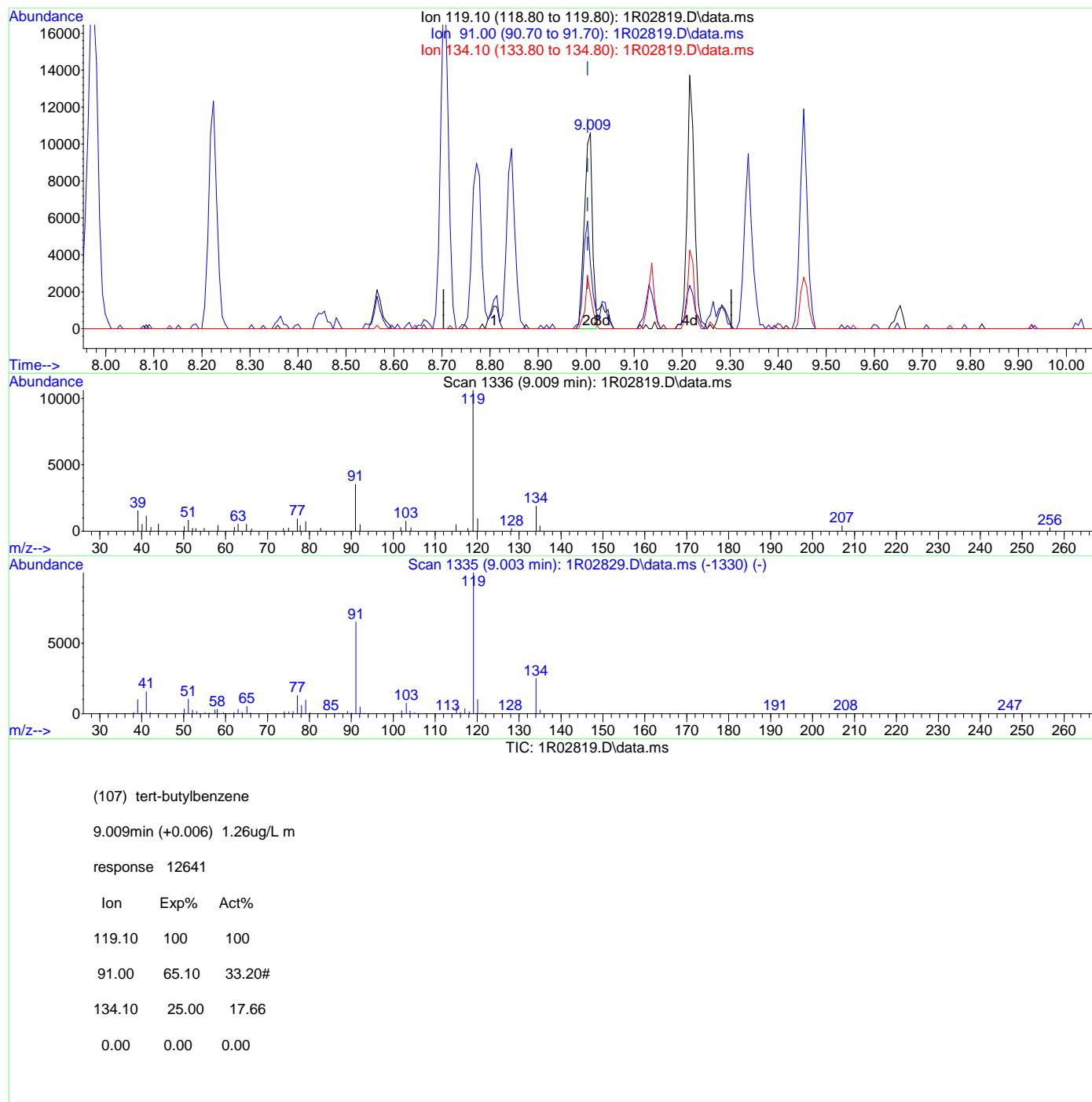
Quant Time: Mar 29 12:43:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:26:40 2023  
 Response via : Initial Calibration



## Quantitation Report (Qedit)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02819.D  
 Acq On : 28 Mar 2023 11:52 pm  
 Operator : PrashanS  
 Sample : IC0091-1  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 29 12:43:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rxi624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:26:40 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02821.D  
 Acq On : 29 Mar 2023 12:25 am  
 Operator : PrashanS  
 Sample : IC0091-2  
 Misc : MS67262,V1R0091,5,,,,1  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Mar 29 12:45:53 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:32:45 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.491	65	465351	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	591860	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	826134	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	787945	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	360724	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.018	113	255997	50.23	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 100.46%			
53) 1,2-dichloroethane-d4 (s)	5.304	65	229877	50.20	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery = 100.40%			
74) toluene-d8 (s)	6.800	98	1067225	48.72	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 97.44%			
98) 4-bromofluorobenzene (s)	8.565	174	313025	53.27	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 106.54%			
<b>Target Compounds</b>						
				Qvalue		
2) ethanol	2.682	45	22124	267.70	ug/L	95
3) tertiary butyl alcohol	3.564	59	13025	10.81	ug/L	96
4) 1,4-dioxane	6.113	88	5034	56.56	ug/L	84
6) chlorodifluoromethane	1.654	51	11278	2.17	ug/L	86
7) dichlorodifluoromethane	1.635	85	9782	2.31	ug/L	75
8) chloromethane	1.806	50	17520	2.17	ug/L	95
9) vinyl chloride	1.897	62	16699	2.35	ug/L	90
10) 1,3-butadiene	1.927	54	11487	2.19	ug/L #	80
11) bromomethane	2.189	94	9718	2.22	ug/L	75
12) chloroethane	2.292	64	8496	2.36	ug/L	95
13) trichlorofluoromethane	2.542	101	11873	2.40	ug/L	89
14) ethyl ether	2.797	74	4707	2.24	ug/L	89
15) acrolein	2.943	56	3121	2.69	ug/L	87
16) freon 113	3.047	151	6377	2.08	ug/L	82
17) 1,1-dichloroethene	3.035	96	7512	2.24	ug/L	82
18) acetone	3.065	58	8183	9.58	ug/L	90
19) acetonitrile	3.333	41	27581	27.66	ug/L	95
20) iodomethane	3.181	142	2623	1.04	ug/L	91
21) carbon disulfide	3.254	76	23615	2.36	ug/L	92
22) methylene chloride	3.491	84	8822	2.60	ug/L	74
23) methyl acetate	3.357	74	2578	2.39	ug/L #	78
24) methyl tert butyl ether	3.728	73	24539	2.41	ug/L	92
25) trans-1,2-dichloroethene	3.740	96	8585	2.29	ug/L	95
26) hexane	3.984	56	7271	2.19	ug/L	92
27) di-isopropyl ether	4.142	45	32838	2.28	ug/L	96
28) 2-butanone	4.598	72	8721	9.21	ug/L	97
29) 1,1-dichloroethane	4.136	63	15614	2.28	ug/L	95
30) chloroprene	4.191	53	13166	2.22	ug/L	87
31) acrylonitrile	3.704	53	6008	2.37	ug/L	84
32) vinyl acetate	4.124	86	2107	2.09	ug/L #	6
33) ethyl tert-butyl ether	4.452	59	26943	2.41	ug/L	95
34) ethyl acetate	4.622	45	2373	1.72	ug/L	96
35) 2,2-dichloropropane	4.616	77	10578	2.17	ug/L	99
36) cis-1,2-dichloroethene	4.610	96	8569	2.09	ug/L	84
37) propionitrile	4.647	54	28621	23.55	ug/L	72
38) methyl acrylate	4.665	85	2013	1.96	ug/L #	89
39) methacrylonitrile	4.775	67	5765	2.10	ug/L #	82
40) bromochloromethane	4.817	128	4285	2.27	ug/L	79
41) tetrahydrofuran	4.817	42	7695	2.22	ug/L	85
42) chloroform	4.890	83	15200	2.11	ug/L	89

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02821.D  
 Acq On : 29 Mar 2023 12:25 am  
 Operator : PrashanS  
 Sample : IC0091-2  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Mar 29 12:45:53 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:32:45 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) tert-Butyl Formate	4.908	59	8053	2.28	ug/L	91
45) 1,1,1-trichloroethane	5.036	97	13835	2.33	ug/L #	1
46) cyclohexane	5.097	84	12011	2.12	ug/L	87
47) isobutyl alcohol	5.213	43	12160	25.38	ug/L	81
48) 1,1-dichloropropene	5.164	75	10979	2.25	ug/L	93
49) carbon tetrachloride	5.164	117	10173	2.16	ug/L	86
50) tert-amyl alcohol	5.310	73	6040	10.59	ug/L #	76
51) isopropyl acetate	5.334	87	3152	2.40	ug/L #	93
54) n-butyl alcohol	5.742	56	32616	110.63	ug/L	98
55) 2,2,4-trimethylpentane	5.413	57	24900	2.16	ug/L	96
56) benzene	5.334	78	36955	2.33	ug/L	85
57) tert-amyl methyl ether	5.419	73	25076	2.29	ug/L	96
58) heptane	5.547	57	6414	2.53	ug/L #	76
59) 1,2-dichloroethane	5.365	62	11137	2.11	ug/L	97
60) ethyl acrylate	5.900	55	17914	2.13	ug/L	91
61) trichloroethene	5.851	95	8911	2.27	ug/L #	79
62) 2-chloroethyl vinyl ether	6.472	63	38247	11.80	ug/L	99
63) methyl methacrylate	6.089	100	2720	1.99	ug/L #	69
64) methylcyclohexane	6.052	83	13489	2.11	ug/L	88
65) 1,2-dichloropropane	6.064	63	9793	2.23	ug/L	94
66) dibromomethane	6.137	93	5264	1.84	ug/L	97
67) bromodichloromethane	6.265	83	10279	2.02	ug/L	94
69) epichlorohydrin	6.514	57	8433	10.58	ug/L	92
70) cis-1,3-dichloropropene	6.600	75	13278	2.13	ug/L	93
71) 4-methyl-2-pentanone	6.691	58	27474	9.21	ug/L #	84
72) isoamyl alcohol	6.721	70	10546	31.88	ug/L #	81
75) toluene	6.849	92	21986	2.15	ug/L #	82
76) ethyl methacrylate	7.032	69	12129	2.31	ug/L	92
77) trans-1,3-dichloropropene	7.007	75	12033	2.35	ug/L	95
78) 1,1,2-trichloroethane	7.153	83	8109	2.59	ug/L	91
79) tetrachloroethene	7.226	164	9430	2.27	ug/L	90
80) 2-hexanone	7.287	58	29851	9.32	ug/L	99
81) 1,3-dichloropropane	7.269	76	13502	2.40	ug/L	84
82) butyl acetate	7.360	56	9629	2.26	ug/L #	83
83) dibromochloromethane	7.421	129	9022	2.36	ug/L	94
84) 1,2-dibromoethane	7.512	107	8899	2.36	ug/L	89
85) n-butyl ether	7.865	57	40236	2.23	ug/L	95
86) chlorobenzene	7.835	112	26461	2.42	ug/L	97
87) 1,1,1,2-tetrachloroethane	7.889	131	8092	2.15	ug/L	95
88) ethylbenzene	7.889	91	38788	2.06	ug/L	90
89) m,p-xylene	7.974	106	30462	4.27	ug/L	88
90) o-xylene	8.224	91	29290	2.00	ug/L	97
91) styrene	8.236	104	24550	2.04	ug/L	89
92) butyl acrylate	8.163	55	21947	2.43	ug/L	97
93) n-amyl acetate	8.297	70	8371	2.15	ug/L	95
94) isopropylbenzene	8.449	105	33737	2.02	ug/L	97
95) bromoform	8.364	173	6160	1.96	ug/L	95
96) cis-1,4-dichloro-2-butene	8.486	88	4652	1.80	ug/L	83
99) 1,1,2,2-tetrachloroethane	8.638	83	13936	2.64	ug/L	99
100) trans-1,4-dichloro-2-b...	8.656	53	4363	2.48	ug/L	91
101) 1,2,3-trichloropropane	8.680	110	3912	2.67	ug/L	91
102) bromobenzene	8.662	156	9897	2.37	ug/L	90
103) n-propylbenzene	8.705	91	42752	2.36	ug/L	92
104) 2-chlorotoluene	8.771	126	9869	2.83	ug/L #	80
105) 4-chlorotoluene	8.844	91	25284	2.18	ug/L	96
106) 1,3,5-trimethylbenzene	8.808	105	28568	2.34	ug/L	94
107) tert-butylbenzene	9.003	119	26131	2.48	ug/L	95
108) 1,2,4-trimethylbenzene	9.039	105	27803	2.23	ug/L	94

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02821.D  
 Acq On : 29 Mar 2023 12:25 am  
 Operator : PrashanS  
 Sample : IC0091-2  
 Misc : MS67262,V1R0091,5,,,,1  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Mar 29 12:45:53 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:32:45 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
109) sec-butylbenzene	9.136	105	34583	2.39	ug/L	97
110) p-isopropyltoluene	9.222	119	30420	2.26	ug/L	93
111) 1,2,3-trimethylbenzene	9.282	105	31402	2.28	ug/L	98
112) 1,3-dichlorobenzene	9.222	146	18778	2.35	ug/L	93
113) 1,4-dichlorobenzene	9.276	146	17212	2.00	ug/L	88
114) 1,2-dichlorobenzene	9.489	146	17149	2.41	ug/L	95
115) benzyl chloride	9.337	91	23787	2.39	ug/L	93
116) n-butylbenzene	9.453	92	14153	2.29	ug/L	90
117) hexachloroethane	9.647	201	2401	2.19	ug/L #	82
118) 1,2-dibromo-3-chloropr...	9.927	157	3450	2.57	ug/L	79
119) 1,3,5-trichlorobenzene	10.031	180	11241	2.56	ug/L	97
120) 1,2,4-trichlorobenzene	10.384	180	9990	2.56	ug/L	94
121) hexachlorobutadiene	10.451	225	3490	2.57	ug/L	89
122) naphthalene	10.536	128	35435	2.79	ug/L	94
123) 1,2,3-trichlorobenzene	10.651	180	9178	2.35	ug/L	95
124) 2-methylnaphthalene	11.132	142	7398	1.21	ug/L	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

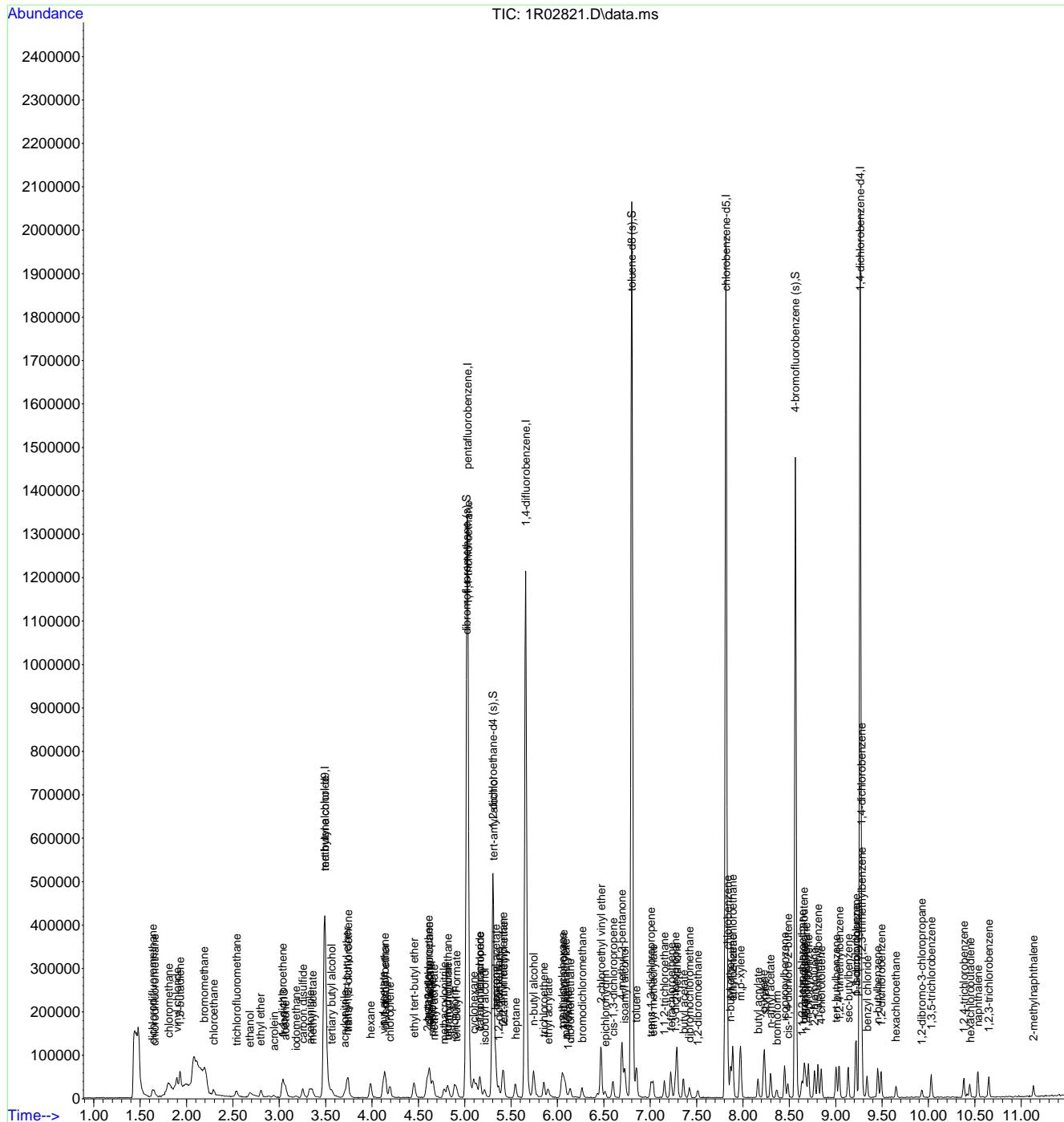
7.6.4

7

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02821.D  
 Acq On : 29 Mar 2023 12:25 am  
 Operator : PrashanS  
 Sample : IC0091-2  
 Misc : MS67262,V1R0091,5,,,.1  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Mar 29 12:45:53 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:32:45 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02823.D  
 Acq On : 29 Mar 2023 12:58 am  
 Operator : PrashanS  
 Sample : IC0091-4  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Mar 29 12:34:27 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:34:20 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.491	65	458621	500.00	ug/L	0.00
5) pentafluorobenzene	5.036	168	590305	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	812769	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	794623	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	375187	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.024	113	260799	51.26	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	102.52%		
53) 1,2-dichloroethane-d4 (s)	5.304	65	228555	50.69	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	101.38%		
74) toluene-d8 (s)	6.800	98	1063624	48.40	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	96.80%		
98) 4-bromofluorobenzene (s)	8.571	174	320601	51.78	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	103.56%		
<b>Target Compounds</b>						
				Qvalue		
2) ethanol	2.682	45	35894	406.30	ug/L	91
3) tertiary butyl alcohol	3.570	59	23684	19.55	ug/L	95
4) 1,4-dioxane	6.113	88	7948	86.81	ug/L	96
6) chlorodifluoromethane	1.660	51	20659	3.88	ug/L	92
7) dichlorodifluoromethane	1.642	85	17929	4.12	ug/L	98
8) chloromethane	1.806	50	30641	3.70	ug/L	92
9) vinyl chloride	1.897	62	31070	4.20	ug/L	89
10) 1,3-butadiene	1.934	54	23218	4.33	ug/L	86
11) bromomethane	2.195	94	13924	3.08	ug/L	88
12) chloroethane	2.299	64	13258	3.48	ug/L	95
13) trichlorofluoromethane	2.548	101	21416	4.17	ug/L	96
14) ethyl ether	2.803	74	9068	4.19	ug/L #	69
15) acrolein	2.943	56	4687	3.45	ug/L	90
16) freon 113	3.047	151	11367	3.69	ug/L	93
17) 1,1-dichloroethene	3.047	96	13305	3.89	ug/L	92
18) acetone	3.077	58	10849	12.13	ug/L	98
19) acetonitrile	3.333	41	47394	40.00	ug/L	99
20) iodomethane	3.187	142	4404	2.31	ug/L	85
21) carbon disulfide	3.260	76	38796	3.72	ug/L	97
22) methylene chloride	3.503	84	15733	4.32	ug/L	97
23) methyl acetate	3.369	74	4375	3.82	ug/L #	49
24) methyl tert butyl ether	3.734	73	40683	3.84	ug/L	97
25) trans-1,2-dichloroethene	3.753	96	13803	3.57	ug/L	94
26) hexane	3.990	56	13014	3.84	ug/L	87
27) di-isopropyl ether	4.148	45	60783	4.12	ug/L	94
28) 2-butanone	4.598	72	13326	13.60	ug/L #	76
29) 1,1-dichloroethane	4.136	63	27232	3.88	ug/L	91
30) chloroprene	4.197	53	22227	3.67	ug/L	94
31) acrylonitrile	3.704	53	11309	4.27	ug/L	85
32) vinyl acetate	4.118	86	3484	3.42	ug/L #	72
33) ethyl tert-butyl ether	4.458	59	45569	3.93	ug/L	92
34) ethyl acetate	4.622	45	5705	4.35	ug/L #	85
35) 2,2-dichloropropane	4.629	77	20128	4.06	ug/L	97
36) cis-1,2-dichloroethene	4.616	96	16316	3.96	ug/L #	83
37) propionitrile	4.653	54	46749	37.25	ug/L	90
38) methyl acrylate	4.677	85	3454	3.40	ug/L #	40
39) methacrylonitrile	4.775	67	10179	3.65	ug/L	94
40) bromochloromethane	4.811	128	7969	4.12	ug/L	85
41) tetrahydrofuran	4.823	42	13617	3.79	ug/L	96
42) chloroform	4.890	83	24699	3.40	ug/L	97

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02823.D  
 Acq On : 29 Mar 2023 12:58 am  
 Operator : PrashanS  
 Sample : IC0091-4  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Mar 29 12:34:27 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:34:20 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) tert-Butyl Formate	4.914	59	14751	4.04	ug/L	96
45) 1,1,1-trichloroethane	5.036	97	22024	3.57	ug/L #	1
46) cyclohexane	5.103	84	20501	3.58	ug/L	93
47) isobutyl alcohol	5.213	43	20639	38.07	ug/L	87
48) 1,1-dichloropropene	5.164	75	19254	3.84	ug/L	91
49) carbon tetrachloride	5.170	117	18733	3.92	ug/L	84
50) tert-amyl alcohol	5.316	73	12355	21.30	ug/L #	83
51) isopropyl acetate	5.340	87	4664	3.34	ug/L #	82
54) n-butyl alcohol	5.742	56	55934	186.24	ug/L	94
55) 2,2,4-trimethylpentane	5.413	57	43093	3.74	ug/L	94
56) benzene	5.334	78	63164	3.92	ug/L	89
57) tert-amyl methyl ether	5.419	73	45727	4.13	ug/L	96
58) heptane	5.541	57	10616	3.99	ug/L #	63
59) 1,2-dichloroethane	5.371	62	19030	3.63	ug/L	86
60) ethyl acrylate	5.900	55	31543	3.74	ug/L	93
61) trichloroethene	5.857	95	14793	3.73	ug/L	94
62) 2-chloroethyl vinyl ether	6.472	63	68716	20.80	ug/L	99
63) methyl methacrylate	6.089	100	4968	3.70	ug/L #	86
64) methylcyclohexane	6.046	83	26700	4.19	ug/L	88
65) 1,2-dichloropropane	6.064	63	17535	3.94	ug/L	99
66) dibromomethane	6.131	93	10892	3.96	ug/L	84
67) bromodichloromethane	6.265	83	18263	3.64	ug/L	98
68) 2-nitropropane	6.429	41	9947	3.30	ug/L	83
69) epichlorohydrin	6.508	57	16637	20.81	ug/L	91
70) cis-1,3-dichloropropene	6.600	75	23139	3.72	ug/L	93
71) 4-methyl-2-pentanone	6.697	58	45351	15.00	ug/L	97
72) isoamyl alcohol	6.721	70	19635	67.15	ug/L	86
75) toluene	6.855	92	39392	3.76	ug/L	96
76) ethyl methacrylate	7.032	69	22242	4.04	ug/L	97
77) trans-1,3-dichloropropene	7.013	75	21594	4.04	ug/L	92
78) 1,1,2-trichloroethane	7.153	83	12841	3.84	ug/L	96
79) tetrachloroethylene	7.226	164	15793	3.67	ug/L	93
80) 2-hexanone	7.287	58	51668	15.48	ug/L	94
81) 1,3-dichloropropane	7.275	76	22353	3.79	ug/L	86
82) butyl acetate	7.354	56	17291	3.90	ug/L	91
83) dibromochloromethane	7.427	129	15061	3.74	ug/L	93
84) 1,2-dibromoethane	7.512	107	15052	3.78	ug/L	97
85) n-butyl ether	7.865	57	72101	3.87	ug/L	96
86) chlorobenzene	7.835	112	42787	3.72	ug/L	96
87) 1,1,1,2-tetrachloroethane	7.889	131	14648	3.79	ug/L	97
88) ethylbenzene	7.889	91	71429	3.73	ug/L	96
89) m,p-xylene	7.974	106	54076	7.41	ug/L	87
90) o-xylene	8.224	91	51731	3.51	ug/L	96
91) styrene	8.236	104	43786	3.59	ug/L	98
92) butyl acrylate	8.163	55	38213	4.02	ug/L	95
93) n-amyl acetate	8.297	70	12985	3.23	ug/L	94
94) isopropylbenzene	8.449	105	62417	3.69	ug/L	96
95) bromoform	8.364	173	10428	3.30	ug/L	92
96) cis-1,4-dichloro-2-butene	8.486	88	7522	2.99	ug/L #	77
99) 1,1,2,2-tetrachloroethane	8.638	83	21607	3.65	ug/L	89
100) trans-1,4-dichloro-2-b...	8.656	53	7749	3.99	ug/L	89
101) 1,2,3-trichloropropane	8.680	110	7324	4.32	ug/L #	75
102) bromobenzene	8.662	156	17218	3.82	ug/L	94
103) n-propylbenzene	8.705	91	78314	3.98	ug/L	97
104) 2-chlorotoluene	8.771	126	16561	4.22	ug/L	88
105) 4-chlorotoluene	8.844	91	45770	3.73	ug/L	93
106) 1,3,5-trimethylbenzene	8.808	105	52960	4.03	ug/L	97
107) tert-butylbenzene	9.003	119	45763	3.94	ug/L	98

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02823.D  
 Acq On : 29 Mar 2023 12:58 am  
 Operator : PrashanS  
 Sample : IC0091-4  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Mar 29 12:34:27 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:34:20 2023  
 Response via : Initial Calibration

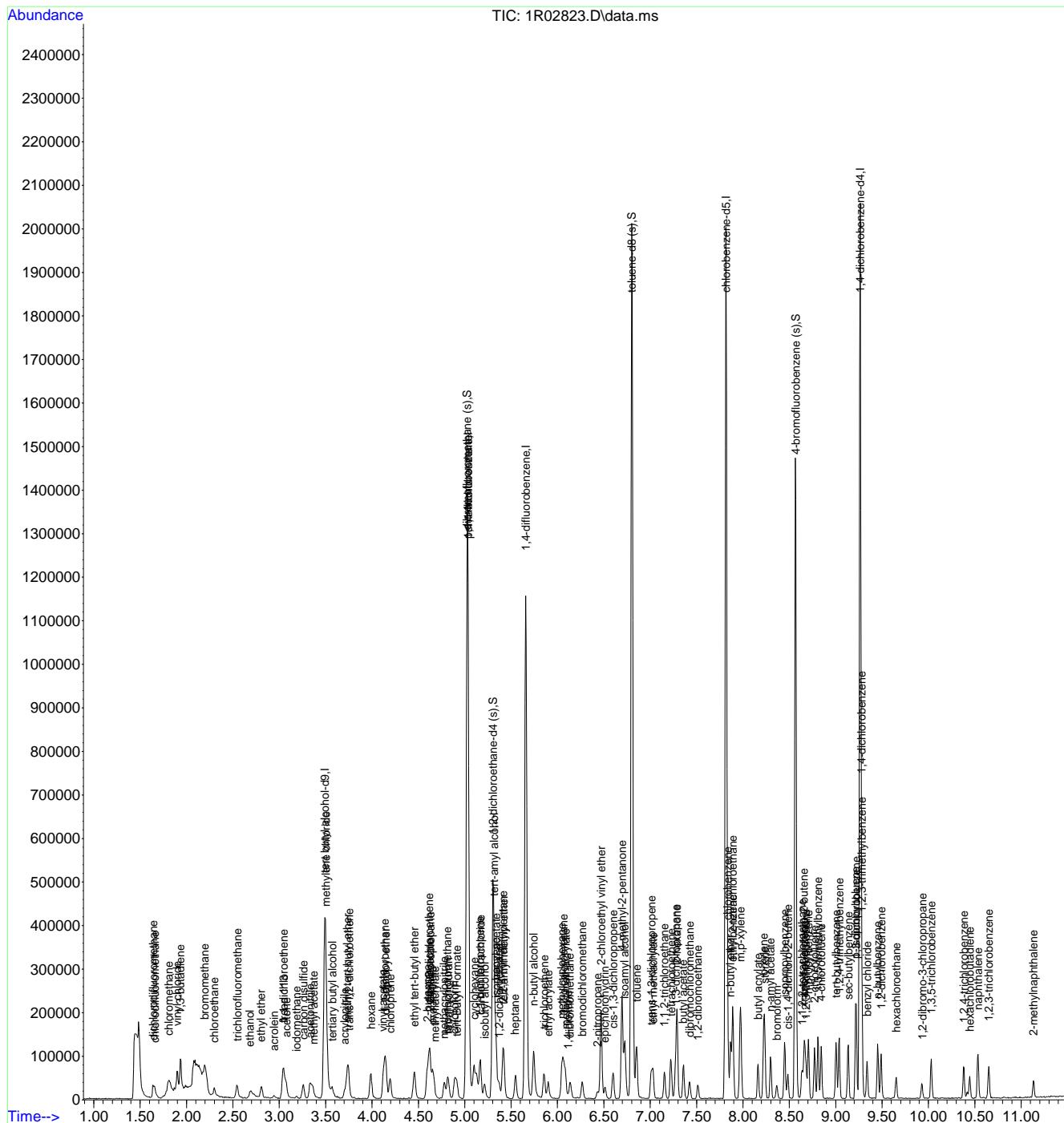
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
108) 1,2,4-trimethylbenzene	9.039	105	52099	3.93	ug/L	92
109) sec-butylbenzene	9.136	105	62018	3.97	ug/L	100
110) p-isopropyltoluene	9.216	119	54675	3.80	ug/L	95
111) 1,2,3-trimethylbenzene	9.282	105	52885	3.57	ug/L	96
112) 1,3-dichlorobenzene	9.222	146	31654	3.68	ug/L	91
113) 1,4-dichlorobenzene	9.276	146	33742	3.77	ug/L	99
114) 1,2-dichlorobenzene	9.489	146	30151	3.92	ug/L	95
115) benzyl chloride	9.337	91	40626	3.75	ug/L	96
116) n-butylbenzene	9.453	92	26042	3.93	ug/L	94
117) hexachloroethane	9.654	201	5038	4.32	ug/L	90
118) 1,2-dibromo-3-chloropr...	9.933	157	5961	3.98	ug/L	87
119) 1,3,5-trichlorobenzene	10.031	180	20522	4.20	ug/L	97
120) 1,2,4-trichlorobenzene	10.384	180	16731	3.90	ug/L	94
121) hexachlorobutadiene	10.444	225	6072	4.01	ug/L	95
122) naphthalene	10.536	128	60621	4.06	ug/L	99
123) 1,2,3-trichlorobenzene	10.651	180	16489	3.89	ug/L	94
124) 2-methylnaphthalene	11.132	142	13544	2.02	ug/L #	79

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02823.D  
 Acq On : 29 Mar 2023 12:58 am  
 Operator : PrashanS  
 Sample : IC0091-4  
 Misc : MS67262,V1R0091,5,,,.1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Mar 29 12:34:27 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:34:20 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02825.D  
 Acq On : 29 Mar 2023 1:31 am  
 Operator : PrashanS  
 Sample : IC0091-8  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Mar 29 12:35:33 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:35:26 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.485	65	439062	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	587607	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	819707	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	801975	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	367914	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.018	113	253711	49.88	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 99.76%			
53) 1,2-dichloroethane-d4 (s)	5.304	65	227226	49.85	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery = 99.70%			
74) toluene-d8 (s)	6.801	98	1065702	48.31	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 96.62%			
98) 4-bromofluorobenzene (s)	8.565	174	313910	51.40	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 102.80%			
<b>Target Compounds</b>						
				Qvalue		
2) ethanol	2.682	45	70404	829.82	ug/L	94
3) tertiary butyl alcohol	3.564	59	43680	37.84	ug/L	94
4) 1,4-dioxane	6.113	88	15620	184.29	ug/L	87
6) chlorodifluoromethane	1.648	51	38281	7.28	ug/L	95
7) dichlorodifluoromethane	1.629	85	35574	8.16	ug/L	99
8) chloromethane	1.800	50	58992	7.29	ug/L	99
9) vinyl chloride	1.891	62	58604	7.88	ug/L	98
10) 1,3-butadiene	1.928	54	44288	8.17	ug/L	97
11) bromomethane	2.189	94	26311	6.20	ug/L	96
12) chloroethane	2.293	64	25092	6.84	ug/L	94
13) trichlorofluoromethane	2.542	101	41476	8.06	ug/L	92
14) ethyl ether	2.798	74	17352	7.99	ug/L	95
15) acrolein	2.944	56	9579	7.43	ug/L	86
16) freon 113	3.035	151	22783	7.54	ug/L	89
17) 1,1-dichloroethene	3.035	96	27830	8.21	ug/L	95
18) acetone	3.065	58	24843	29.32	ug/L	87
19) acetonitrile	3.327	41	87044	73.80	ug/L	96
20) iodomethane	3.181	142	13702	8.40	ug/L	95
21) carbon disulfide	3.254	76	78247	7.64	ug/L	99
22) methylene chloride	3.497	84	31102	8.45	ug/L	94
23) methyl acetate	3.357	74	8425	7.47	ug/L	94
24) methyl tert butyl ether	3.728	73	78098	7.46	ug/L	97
25) trans-1,2-dichloroethene	3.740	96	28595	7.59	ug/L	95
26) hexane	3.978	56	25003	7.48	ug/L	94
27) di-isopropyl ether	4.142	45	116235	7.88	ug/L	96
28) 2-butanone	4.592	72	30144	31.85	ug/L	94
29) 1,1-dichloroethane	4.130	63	54473	7.83	ug/L	95
30) chloroprene	4.197	53	44728	7.53	ug/L	92
31) acrylonitrile	3.704	53	22719	8.51	ug/L	90
32) vinyl acetate	4.124	86	7386	7.55	ug/L #	60
33) ethyl tert-butyl ether	4.452	59	91492	7.95	ug/L	94
34) ethyl acetate	4.617	45	10920	8.18	ug/L #	81
35) 2,2-dichloropropane	4.617	77	37059	7.48	ug/L	96
36) cis-1,2-dichloroethene	4.617	96	30741	7.51	ug/L	94
37) propionitrile	4.653	54	92561	74.94	ug/L	97
38) methyl acrylate	4.665	85	7871	8.09	ug/L #	75
39) methacrylonitrile	4.781	67	19614	7.23	ug/L	92
40) bromochloromethane	4.817	128	14052	7.26	ug/L	88
41) tetrahydrofuran	4.823	42	26099	7.40	ug/L	92
42) chloroform	4.890	83	49440	7.04	ug/L	93

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02825.D  
 Acq On : 29 Mar 2023 1:31 am  
 Operator : PrashanS  
 Sample : IC0091-8  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Mar 29 12:35:33 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:35:26 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) tert-Butyl Formate	4.915	59	28314	7.78	ug/L	95
45) 1,1,1-trichloroethane	5.036	97	43655	7.27	ug/L #	1
46) cyclohexane	5.097	84	40112	7.19	ug/L	95
47) isobutyl alcohol	5.207	43	40068	75.46	ug/L	98
48) 1,1-dichloropropene	5.164	75	39731	8.03	ug/L	99
49) carbon tetrachloride	5.164	117	37420	7.89	ug/L	96
50) tert-amyl alcohol	5.310	73	23729	40.44	ug/L	94
51) isopropyl acetate	5.334	87	10134	7.61	ug/L	95
54) n-butyl alcohol	5.742	56	118314	397.44	ug/L	97
55) 2,2,4-trimethylpentane	5.413	57	88457	7.69	ug/L	96
56) benzene	5.328	78	117813	7.27	ug/L	93
57) tert-amyl methyl ether	5.420	73	88046	7.84	ug/L	95
58) heptane	5.547	57	20241	7.55	ug/L	97
59) 1,2-dichloroethane	5.365	62	37740	7.24	ug/L	97
60) ethyl acrylate	5.894	55	63303	7.55	ug/L	98
61) trichloroethene	5.851	95	30794	7.78	ug/L	95
62) 2-chloroethyl vinyl ether	6.466	63	133066	39.68	ug/L	96
63) methyl methacrylate	6.089	100	10042	7.56	ug/L #	73
64) methylcyclohexane	6.052	83	49131	7.59	ug/L	95
65) 1,2-dichloropropane	6.064	63	33526	7.49	ug/L	96
66) dibromomethane	6.137	93	19463	7.02	ug/L	85
67) bromodichloromethane	6.265	83	35506	7.15	ug/L	99
68) 2-nitropropane	6.429	41	18333	6.41	ug/L	96
69) epichlorohydrin	6.509	57	32912	40.41	ug/L	95
70) cis-1,3-dichloropropene	6.600	75	47566	7.68	ug/L	97
71) 4-methyl-2-pentanone	6.697	58	92596	30.69	ug/L	97
72) isoamyl alcohol	6.721	70	41048	147.07	ug/L	93
75) toluene	6.849	92	77915	7.45	ug/L	92
76) ethyl methacrylate	7.032	69	43912	7.88	ug/L	97
77) trans-1,3-dichloropropene	7.007	75	43284	8.00	ug/L	99
78) 1,1,2-trichloroethane	7.153	83	26329	7.86	ug/L	94
79) tetrachloroethylene	7.220	164	32150	7.51	ug/L	94
80) 2-hexanone	7.287	58	102741	30.67	ug/L	98
81) 1,3-dichloropropane	7.269	76	47376	8.03	ug/L	96
82) butyl acetate	7.360	56	33482	7.52	ug/L	91
83) dibromochloromethane	7.427	129	29651	7.40	ug/L	96
84) 1,2-dibromoethane	7.512	107	29349	7.39	ug/L	98
85) n-butyl ether	7.865	57	143165	7.65	ug/L	99
86) chlorobenzene	7.835	112	85599	7.46	ug/L	96
87) 1,1,1,2-tetrachloroethane	7.889	131	29324	7.59	ug/L	96
88) ethylbenzene	7.889	91	142287	7.45	ug/L	98
89) m,p-xylene	7.975	106	111044	15.27	ug/L	93
90) o-xylene	8.224	91	104788	7.18	ug/L	96
91) styrene	8.236	104	86090	7.14	ug/L	93
92) butyl acrylate	8.163	55	77569	8.08	ug/L	99
93) n-amyl acetate	8.297	70	26842	6.95	ug/L #	79
94) isopropylbenzene	8.449	105	126746	7.55	ug/L	96
95) bromoform	8.364	173	23747	7.67	ug/L	94
96) cis-1,4-dichloro-2-butene	8.486	88	16471	6.92	ug/L #	84
99) 1,1,2,2-tetrachloroethane	8.638	83	50824	8.91	ug/L	97
100) trans-1,4-dichloro-2-b...	8.656	53	15739	8.27	ug/L	97
101) 1,2,3-trichloropropane	8.680	110	13325	7.86	ug/L	97
102) bromobenzene	8.662	156	35952	8.20	ug/L	90
103) n-propylbenzene	8.705	91	155426	8.07	ug/L	98
104) 2-chlorotoluene	8.772	126	32227	8.30	ug/L	98
105) 4-chlorotoluene	8.845	91	88833	7.46	ug/L	97
106) 1,3,5-trimethylbenzene	8.808	105	107954	8.37	ug/L	99
107) tert-butylbenzene	9.003	119	91177	8.03	ug/L	95

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02825.D  
 Acq On : 29 Mar 2023 1:31 am  
 Operator : PrashanS  
 Sample : IC0091-8  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Mar 29 12:35:33 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:35:26 2023  
 Response via : Initial Calibration

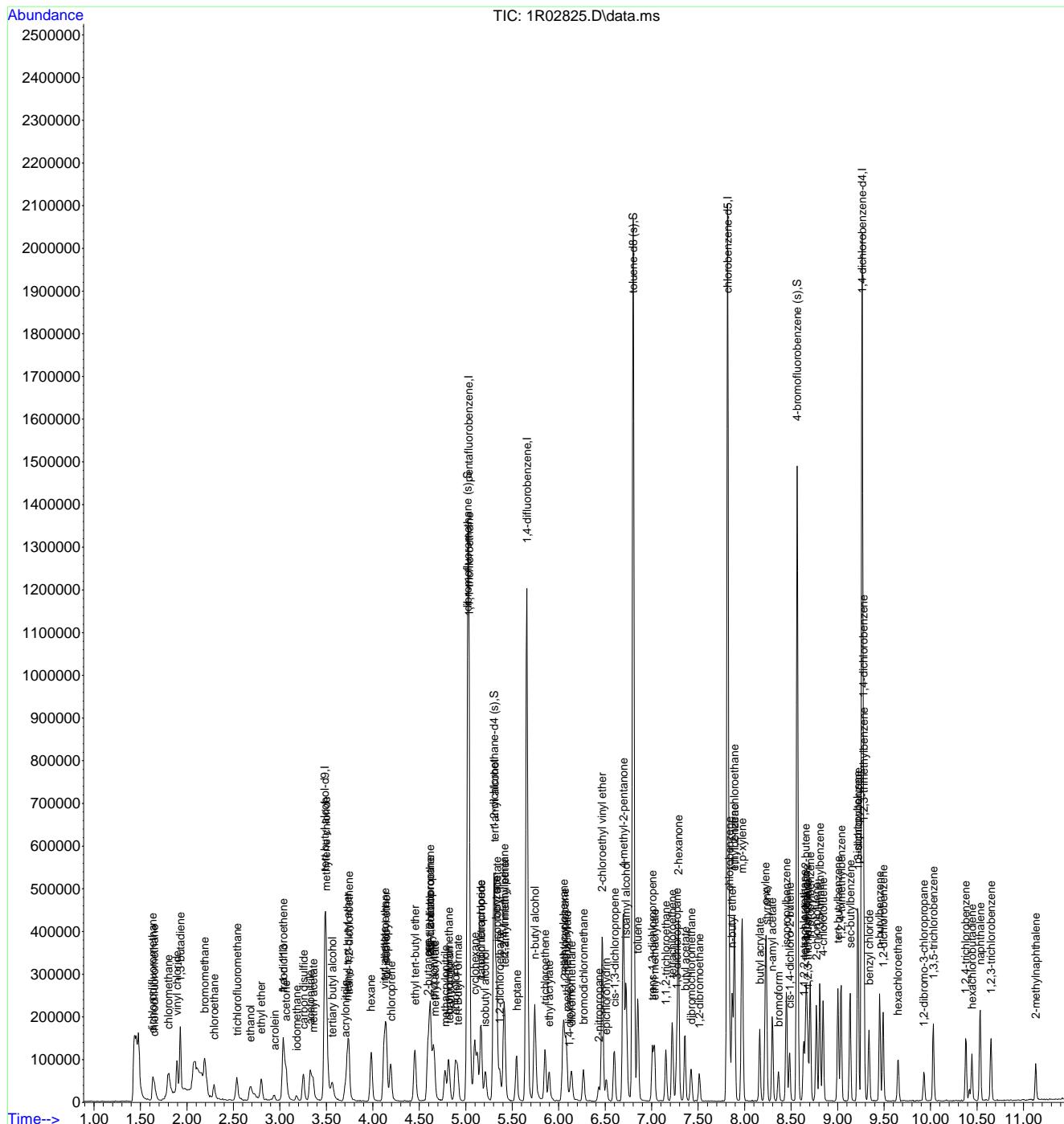
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
108) 1,2,4-trimethylbenzene	9.039	105	108936	8.41	ug/L	98
109) sec-butylbenzene	9.137	105	127218	8.31	ug/L	95
110) p-isopropyltoluene	9.216	119	109816	7.85	ug/L	96
111) 1,2,3-trimethylbenzene	9.283	105	108681	7.65	ug/L	98
112) 1,3-dichlorobenzene	9.222	146	63177	7.60	ug/L	94
113) 1,4-dichlorobenzene	9.277	146	61804	7.11	ug/L	87
114) 1,2-dichlorobenzene	9.489	146	63434	8.43	ug/L	99
115) benzyl chloride	9.337	91	79081	7.53	ug/L	93
116) n-butylbenzene	9.453	92	50011	7.72	ug/L	99
117) hexachloroethane	9.648	201	12028	10.35	ug/L	82
118) 1,2-dibromo-3-chloropr...	9.927	157	13361	9.11	ug/L	93
119) 1,3,5-trichlorobenzene	10.031	180	40488	8.37	ug/L	97
120) 1,2,4-trichlorobenzene	10.378	180	32689	7.80	ug/L	86
121) hexachlorobutadiene	10.445	225	12653	8.52	ug/L	94
122) naphthalene	10.536	128	122432	8.32	ug/L	99
123) 1,2,3-trichlorobenzene	10.651	180	32241	7.81	ug/L	99
124) 2-methylnaphthalene	11.132	142	27221	4.14	ug/L	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
Data File : 1R02825.D  
Acq On : 29 Mar 2023 1:31 am  
Operator : PrashanS  
Sample : IC0091-8  
Misc : MS67262,V1R0091,5,,,1  
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Mar 29 12:35:33 2023  
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
QLast Update : Wed Mar 29 12:35:26 2023  
Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02827.D  
 Acq On : 29 Mar 2023 2:04 am  
 Operator : PrashanS  
 Sample : IC0091-20  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Mar 29 12:36:29 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:36:23 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.491	65	497795	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	614482	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	876260	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	843977	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	451452	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.018	113	266431	50.11	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	100.22%		
53) 1,2-dichloroethane-d4 (s)	5.304	65	251179	51.57	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	103.14%		
74) toluene-d8 (s)	6.800	98	1176684	50.93	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	101.86%		
98) 4-bromofluorobenzene (s)	8.565	174	347191	46.14	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	92.28%		
<b>Target Compounds</b>						
				Qvalue		
2) ethanol	2.682	45	179853	1858.20	ug/L	99
3) tertiary butyl alcohol	3.564	59	111393	85.89	ug/L	96
4) 1,4-dioxane	6.125	88	40838	431.75	ug/L	94
6) chlorodifluoromethane	1.648	51	92801	17.18	ug/L	97
7) dichlorodifluoromethane	1.629	85	82414	18.03	ug/L	98
8) chloromethane	1.800	50	137663	16.56	ug/L	96
9) vinyl chloride	1.891	62	134620	17.35	ug/L	95
10) 1,3-butadiene	1.927	54	102338	17.99	ug/L	95
11) bromomethane	2.183	94	71986	16.99	ug/L	90
12) chloroethane	2.293	64	66836	17.95	ug/L	95
13) trichlorofluoromethane	2.536	101	101798	18.90	ug/L	92
14) ethyl ether	2.804	74	41935	18.46	ug/L	90
15) acrolein	2.937	56	24389	18.41	ug/L	98
16) freon 113	3.041	151	53298	17.03	ug/L	95
17) 1,1-dichloroethene	3.035	96	61863	17.39	ug/L	91
18) acetone	3.065	58	65777	75.30	ug/L	93
19) acetonitrile	3.327	41	219744	181.68	ug/L	96
20) iodomethane	3.181	142	41106	23.80	ug/L	98
21) carbon disulfide	3.248	76	184641	17.36	ug/L	99
22) methylene chloride	3.497	84	70793	18.21	ug/L	98
23) methyl acetate	3.351	74	20981	18.02	ug/L #	80
24) methyl tert butyl ether	3.728	73	193236	17.83	ug/L	96
25) trans-1,2-dichloroethene	3.740	96	69163	17.70	ug/L	91
26) hexane	3.984	56	57166	16.52	ug/L	91
27) di-isopropyl ether	4.142	45	275141	17.87	ug/L	95
28) 2-butanone	4.592	72	76219	77.08	ug/L	98
29) 1,1-dichloroethane	4.130	63	130156	17.95	ug/L	98
30) chloroprene	4.191	53	105985	17.20	ug/L	94
31) acrylonitrile	3.698	53	55667	19.73	ug/L	96
32) vinyl acetate	4.118	86	16353	16.17	ug/L #	90
33) ethyl tert-butyl ether	4.452	59	220761	18.36	ug/L	98
34) ethyl acetate	4.616	45	25969	18.53	ug/L #	79
35) 2,2-dichloropropane	4.622	77	90094	17.59	ug/L	96
36) cis-1,2-dichloroethene	4.610	96	76245	17.97	ug/L	89
37) propionitrile	4.647	54	238563	186.39	ug/L	84
38) methyl acrylate	4.665	85	18794	18.42	ug/L #	72
39) methacrylonitrile	4.781	67	47487	17.07	ug/L	96
40) bromochloromethane	4.811	128	34399	17.22	ug/L	90
41) tetrahydrofuran	4.817	42	62017	17.07	ug/L	92
42) chloroform	4.890	83	120077	16.69	ug/L	96

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02827.D  
 Acq On : 29 Mar 2023 2:04 am  
 Operator : PrashanS  
 Sample : IC0091-20  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Mar 29 12:36:29 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:36:23 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) tert-Butyl Formate	4.908	59	68289	18.02	ug/L	95
45) 1,1,1-trichloroethane	5.036	97	108177	17.49	ug/L #	24
46) cyclohexane	5.097	84	105191	18.33	ug/L	88
47) isobutyl alcohol	5.213	43	103596	189.25	ug/L	99
48) 1,1-dichloropropene	5.164	75	95270	18.40	ug/L	95
49) carbon tetrachloride	5.164	117	93168	18.83	ug/L	100
50) tert-amyl alcohol	5.310	73	61719	100.36	ug/L	97
51) isopropyl acetate	5.328	87	24639	17.86	ug/L #	88
54) n-butyl alcohol	5.742	56	311874	981.29	ug/L	95
55) 2,2,4-trimethylpentane	5.407	57	215360	17.61	ug/L	97
56) benzene	5.328	78	301819	17.66	ug/L	97
57) tert-amyl methyl ether	5.419	73	219524	18.33	ug/L	98
58) heptane	5.541	57	47789	16.84	ug/L	86
59) 1,2-dichloroethane	5.365	62	91444	16.64	ug/L	94
60) ethyl acrylate	5.894	55	153576	17.29	ug/L	98
61) trichloroethene	5.851	95	78072	18.52	ug/L	95
62) 2-chloroethyl vinyl ether	6.466	63	338721	94.60	ug/L	99
63) methyl methacrylate	6.089	100	24640	17.55	ug/L #	81
64) methylcyclohexane	6.046	83	116778	17.00	ug/L	98
65) 1,2-dichloropropane	6.064	63	82515	17.44	ug/L	99
66) dibromomethane	6.137	93	48396	16.68	ug/L	93
67) bromodichloromethane	6.265	83	89336	17.13	ug/L	96
68) 2-nitropropane	6.429	41	41820	14.39	ug/L	99
69) epichlorohydrin	6.515	57	79382	90.99	ug/L	97
70) cis-1,3-dichloropropene	6.594	75	119090	18.08	ug/L	97
71) 4-methyl-2-pentanone	6.691	58	241576	75.34	ug/L	97
72) isoamyl alcohol	6.721	70	113568	388.48	ug/L	95
75) toluene	6.849	92	203852	18.70	ug/L	97
76) ethyl methacrylate	7.032	69	116401	19.91	ug/L	99
77) trans-1,3-dichloropropene	7.007	75	111164	19.53	ug/L	97
78) 1,1,2-trichloroethane	7.153	83	61126	17.39	ug/L	97
79) tetrachloroethylene	7.220	164	75281	16.85	ug/L	95
80) 2-hexanone	7.287	58	267472	76.32	ug/L	100
81) 1,3-dichloropropane	7.269	76	114026	18.35	ug/L	98
82) butyl acetate	7.360	56	84511	18.21	ug/L	96
83) dibromochloromethane	7.421	129	77858	18.69	ug/L	95
84) 1,2-dibromoethane	7.512	107	76692	18.58	ug/L	93
85) n-butyl ether	7.865	57	372395	19.04	ug/L	100
86) chlorobenzene	7.835	112	220500	18.45	ug/L	98
87) 1,1,1,2-tetrachloroethane	7.889	131	77784	19.30	ug/L	97
88) ethylbenzene	7.889	91	372338	18.71	ug/L	98
89) m,p-xylene	7.975	106	287122	37.76	ug/L	95
90) o-xylene	8.224	91	282162	18.65	ug/L	99
91) styrene	8.236	104	241245	19.35	ug/L	98
92) butyl acrylate	8.163	55	196439	19.41	ug/L	98
93) n-amyl acetate	8.297	70	75470	19.06	ug/L	96
94) isopropylbenzene	8.449	105	334148	19.09	ug/L	99
95) bromoform	8.364	173	58441	18.04	ug/L	96
96) cis-1,4-dichloro-2-butene	8.486	88	43594	17.88	ug/L	89
99) 1,1,2,2-tetrachloroethane	8.638	83	126364	17.71	ug/L	97
100) trans-1,4-dichloro-2-b...	8.656	53	44652	19.02	ug/L	93
101) 1,2,3-trichloropropane	8.680	110	36715	17.72	ug/L	88
102) bromobenzene	8.662	156	91212	16.89	ug/L	97
103) n-propylbenzene	8.705	91	422613	17.85	ug/L	98
104) 2-chlorotoluene	8.771	126	87376	18.23	ug/L	91
105) 4-chlorotoluene	8.845	91	250851	17.33	ug/L	99
106) 1,3,5-trimethylbenzene	8.808	105	287897	18.08	ug/L	95
107) tert-butylbenzene	9.003	119	245547	17.61	ug/L	99

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02827.D  
 Acq On : 29 Mar 2023 2:04 am  
 Operator : PrashanS  
 Sample : IC0091-20  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Mar 29 12:36:29 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:36:23 2023  
 Response via : Initial Calibration

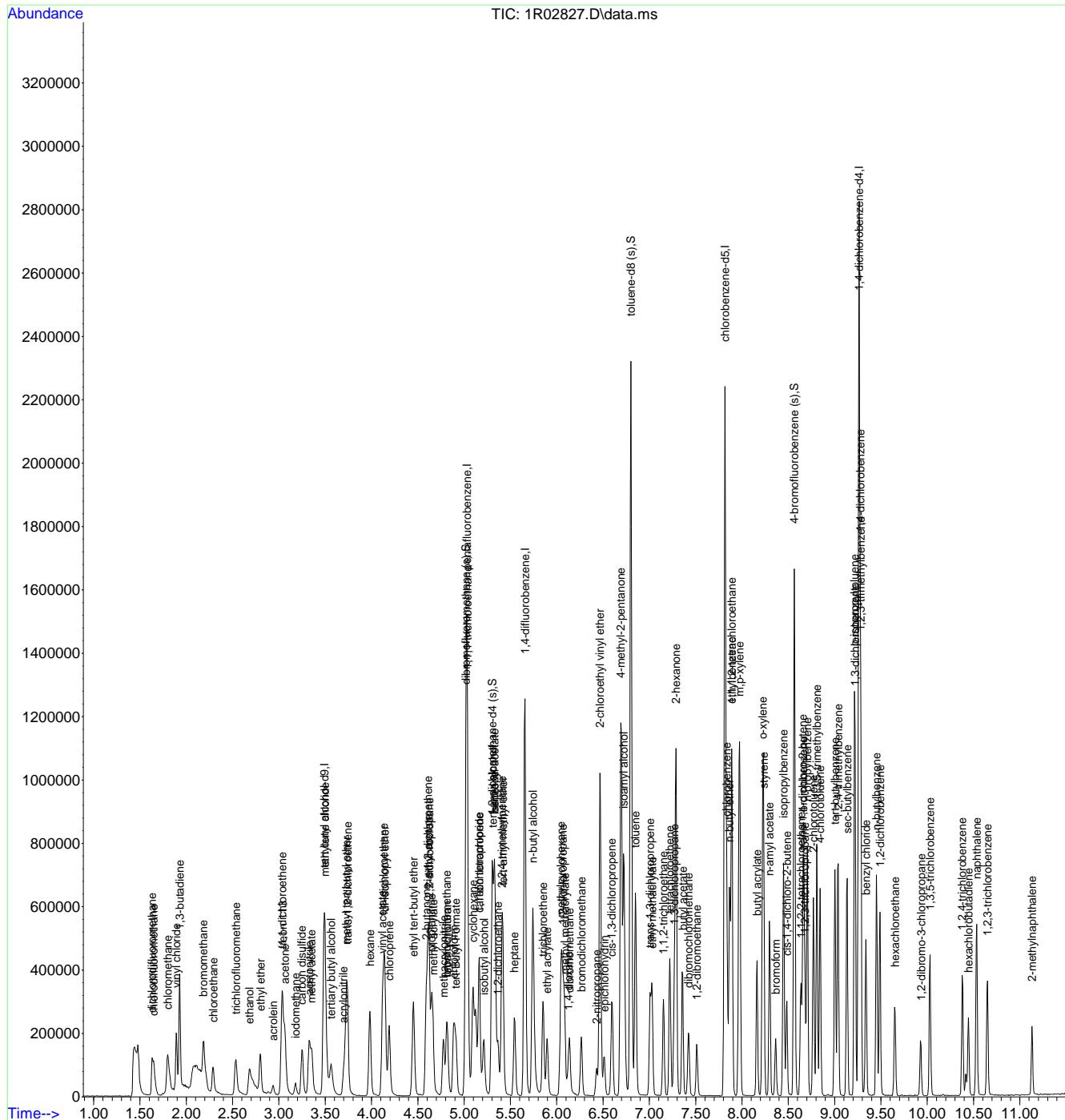
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
108) 1,2,4-trimethylbenzene	9.039	105	294964	18.41	ug/L	98
109) sec-butylbenzene	9.137	105	346094	18.32	ug/L	97
110) p-isopropyltoluene	9.216	119	298758	17.45	ug/L	98
111) 1,2,3-trimethylbenzene	9.283	105	306839	17.73	ug/L	95
112) 1,3-dichlorobenzene	9.222	146	171852	16.96	ug/L	99
113) 1,4-dichlorobenzene	9.276	146	180055	17.16	ug/L	99
114) 1,2-dichlorobenzene	9.489	146	167983	18.06	ug/L	100
115) benzyl chloride	9.337	91	233843	18.33	ug/L	96
116) n-butylbenzene	9.453	92	147689	18.68	ug/L	95
117) hexachloroethane	9.654	201	32309	21.59	ug/L	98
118) 1,2-dibromo-3-chloropr...	9.933	157	33227	18.05	ug/L	88
119) 1,3,5-trichlorobenzene	10.031	180	101824	17.03	ug/L	98
120) 1,2,4-trichlorobenzene	10.378	180	86348	16.86	ug/L	96
121) hexachlorobutadiene	10.444	225	33423	18.14	ug/L	96
122) naphthalene	10.536	128	314673	17.29	ug/L	99
123) 1,2,3-trichlorobenzene	10.651	180	85523	16.94	ug/L	93
124) 2-methylnaphthalene	11.132	142	71181	8.76	ug/L	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
Data File : 1R02827.D  
Acq On : 29 Mar 2023 2:04 am  
Operator : PrashanS  
Sample : IC0091-20  
Misc : MS67262,V1R0091,5,,,1  
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Mar 29 12:36:29 2023  
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
QLast Update: Wed Mar 29 12:36:23 2023  
Response via: Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02829.D  
 Acq On : 29 Mar 2023 2:37 am  
 Operator : PrashanS  
 Sample : ICC0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Mar 29 12:11:30 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:11:26 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.491	65	501834	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	694530	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	940928	50.00	ug/L	0.00
73) chlorobenzene-d5	7.817	117	1005733	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	682662	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.018	113	304894	50.00	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 100.00%			
53) 1,2-dichloroethane-d4 (s)	5.304	65	276087	50.00	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery = 100.00%			
74) toluene-d8 (s)	6.801	98	1541613	50.00	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 100.00%			
98) 4-bromofluorobenzene (s)	8.565	174	468613	50.00	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery = 100.00%			
<b>Target Compounds</b>						
				Qvalue		
2) ethanol	2.688	45	422373	5000.00	ug/L	100
3) tertiary butyl alcohol	3.570	59	287426	250.00	ug/L	100
4) 1,4-dioxane	6.119	88	105218	1250.00	ug/L	100
6) chlorodifluoromethane	1.654	51	246216	50.00	ug/L	100
7) dichlorodifluoromethane	1.636	85	227725	50.00	ug/L	100
8) chloromethane	1.800	50	357107	50.00	ug/L	100
9) vinyl chloride	1.897	62	352598	50.14	ug/L	100
10) 1,3-butadiene	1.934	54	275337	50.00	ug/L	100
11) bromomethane	2.189	94	216280	50.00	ug/L	100
12) chloroethane	2.293	64	188031	50.00	ug/L	100
13) trichlorofluoromethane	2.536	101	285374	50.00	ug/L	100
14) ethyl ether	2.804	74	118894	50.00	ug/L	100
15) acrolein	2.938	56	68155	50.00	ug/L	100
16) freon 113	3.041	151	158855	50.05	ug/L	100
17) 1,1-dichloroethene	3.041	96	177652	50.00	ug/L	100
18) acetone	3.065	58	150660	200.00	ug/L	100
19) acetonitrile	3.333	41	584973	500.00	ug/L	100
20) iodomethane	3.181	142	147568	50.00	ug/L	100
21) carbon disulfide	3.254	76	506315	50.00	ug/L	100
22) methylene chloride	3.497	84	203993	50.00	ug/L	100
23) methyl acetate	3.357	74	55706	50.00	ug/L	100
24) methyl tert butyl ether	3.728	73	575169	50.00	ug/L	100
25) trans-1,2-dichloroethene	3.741	96	202690	50.00	ug/L	100
26) hexane	3.984	56	166590	50.00	ug/L	100
27) di-isopropyl ether	4.142	45	833277	50.00	ug/L	100
28) 2-butanone	4.592	72	200858	200.00	ug/L	100
29) 1,1-dichloroethane	4.130	63	382464	50.16	ug/L	100
30) chloroprene	4.191	53	340489	50.00	ug/L	100
31) acrylonitrile	3.704	53	160405	50.00	ug/L	100
32) vinyl acetate	4.118	86	54080	50.00	ug/L	100
33) ethyl tert-butyl ether	4.452	59	673197	50.00	ug/L	100
34) ethyl acetate	4.617	45	74209	50.00	ug/L	100
35) 2,2-dichloropropane	4.623	77	265878	50.00	ug/L	100
36) cis-1,2-dichloroethene	4.617	96	222967	50.00	ug/L	100
37) propionitrile	4.653	54	680248	500.00	ug/L	100
38) methyl acrylate	4.665	85	58021	50.00	ug/L	100
39) methacrylonitrile	4.781	67	151882	50.00	ug/L	100
40) bromochloromethane	4.811	128	110133	50.00	ug/L	100
41) tetrahydrofuran	4.823	42	180696	50.00	ug/L	100
42) chloroform	4.890	83	369311	50.00	ug/L	100

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02829.D  
 Acq On : 29 Mar 2023 2:37 am  
 Operator : PrashanS  
 Sample : ICC0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Mar 29 12:11:30 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:11:26 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) tert-Butyl Formate	4.915	59	230067	50.00	ug/L	100
45) 1,1,1-trichloroethane	5.036	97	315681	50.00	ug/L	100
46) cyclohexane	5.097	84	312714	50.00	ug/L	100
47) isobutyl alcohol	5.213	43	281153	500.00	ug/L	100
48) 1,1-dichloropropene	5.164	75	277543	50.00	ug/L	100
49) carbon tetrachloride	5.164	117	281010	50.00	ug/L	100
50) tert-amyl alcohol	5.316	73	161535	250.00	ug/L	100
51) isopropyl acetate	5.334	87	72291	50.00	ug/L	100
54) n-butyl alcohol	5.748	56	874923	2500.00	ug/L	100
55) 2,2,4-trimethylpentane	5.414	57	626965	50.00	ug/L	100
56) benzene	5.328	78	887008	50.00	ug/L	100
57) tert-amyl methyl ether	5.420	73	651152	50.08	ug/L	100
58) heptane	5.547	57	139765	50.00	ug/L	100
59) 1,2-dichloroethane	5.365	62	285924	50.00	ug/L	100
60) ethyl acrylate	5.894	55	472754	50.00	ug/L	100
61) trichloroethene	5.852	95	223654	50.00	ug/L	100
62) 2-chloroethyl vinyl ether	6.472	63	1099080	250.00	ug/L	100
63) methyl methacrylate	6.089	100	72620	50.00	ug/L	100
64) methylcyclohexane	6.046	83	350060	50.00	ug/L	100
65) 1,2-dichloropropane	6.064	63	247629	50.00	ug/L	100
66) dibromomethane	6.137	93	150195	50.00	ug/L	100
67) bromodichloromethane	6.265	83	280490	50.00	ug/L	100
68) 2-nitropropane	6.430	41	130218	49.81	ug/L	99
69) epichlorohydrin	6.515	57	242549	250.00	ug/L	100
70) cis-1,3-dichloropropene	6.600	75	362576	50.00	ug/L	100
71) 4-methyl-2-pentanone	6.697	58	779840	200.00	ug/L	100
72) isoamyl alcohol	6.728	70	376734	1000.00	ug/L	100
75) toluene	6.855	92	679358	50.00	ug/L	100
76) ethyl methacrylate	7.032	69	362878	50.00	ug/L	100
77) trans-1,3-dichloropropene	7.007	75	342756	50.00	ug/L	100
78) 1,1,2-trichloroethane	7.153	83	189705	50.00	ug/L	100
79) tetrachloroethylene	7.220	164	216040	50.00	ug/L	100
80) 2-hexanone	7.287	58	832743	200.00	ug/L	100
81) 1,3-dichloropropane	7.269	76	347870	50.00	ug/L	100
82) butyl acetate	7.360	56	259973	50.00	ug/L	100
83) dibromochloromethane	7.421	129	248921	50.00	ug/L	100
84) 1,2-dibromoethane	7.512	107	233677	50.00	ug/L	100
85) n-butyl ether	7.865	57	1153362	50.00	ug/L	100
86) chlorobenzene	7.835	112	683279	50.00	ug/L	100
87) 1,1,1,2-tetrachloroethane	7.890	131	273322	50.00	ug/L	100
88) ethylbenzene	7.890	91	1261421	50.00	ug/L	100
89) m,p-xylene	7.975	106	1035613	100.00	ug/L	100
90) o-xylene	8.224	91	1064079	50.00	ug/L	100
91) styrene	8.236	104	986379	50.00	ug/L	100
92) butyl acrylate	8.163	55	611847	50.00	ug/L	100
93) n-amyl acetate	8.297	70	282318	50.00	ug/L	100
94) isopropylbenzene	8.449	105	1302729	50.00	ug/L	100
95) bromoform	8.364	173	237514	50.00	ug/L	100
96) cis-1,4-dichloro-2-butene	8.486	88	191989	50.00	ug/L	100
99) 1,1,2,2-tetrachloroethane	8.638	83	374412	50.00	ug/L	100
100) trans-1,4-dichloro-2-b...	8.656	53	147572	50.00	ug/L	100
101) 1,2,3-trichloropropane	8.680	110	115549	50.00	ug/L	100
102) bromobenzene	8.662	156	324520	50.00	ug/L	100
103) n-propylbenzene	8.705	91	1446533	50.00	ug/L	100
104) 2-chlorotoluene	8.772	126	288159	50.00	ug/L	100
105) 4-chlorotoluene	8.845	91	977811	50.00	ug/L	100
106) 1,3,5-trimethylbenzene	8.808	105	1106738	50.00	ug/L	100
107) tert-butylbenzene	9.003	119	859101	50.00	ug/L	100

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02829.D  
 Acq On : 29 Mar 2023 2:37 am  
 Operator : PrashanS  
 Sample : ICC0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Mar 29 12:11:30 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:11:26 2023  
 Response via : Initial Calibration

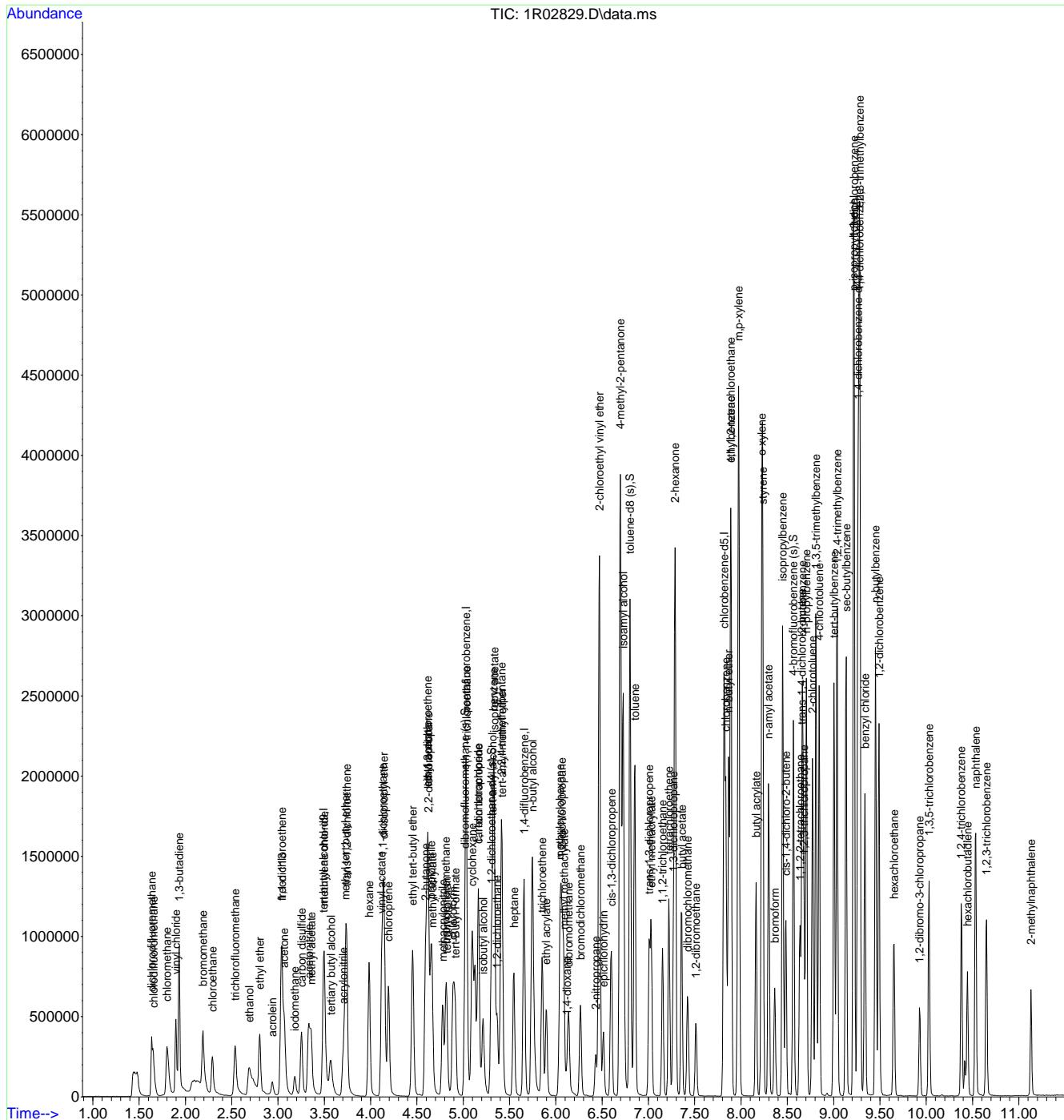
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
108) 1,2,4-trimethylbenzene	9.039	105	1229963	50.00	ug/L	100
109) sec-butylbenzene	9.137	105	1346439	50.00	ug/L	100
110) p-isopropyltoluene	9.216	119	1329146	50.00	ug/L	100
111) 1,2,3-trimethylbenzene	9.283	105	1309487	50.00	ug/L	100
112) 1,3-dichlorobenzene	9.222	146	771446	50.00	ug/L	100
113) 1,4-dichlorobenzene	9.277	146	762393	50.00	ug/L	100
114) 1,2-dichlorobenzene	9.490	146	709701	50.00	ug/L	100
115) benzyl chloride	9.337	91	918856	50.00	ug/L	100
116) n-butylbenzene	9.453	92	565553	50.00	ug/L	100
117) hexachloroethane	9.654	201	116356	50.00	ug/L	100
118) 1,2-dibromo-3-chloropr...	9.928	157	102775	50.00	ug/L	100
119) 1,3,5-trichlorobenzene	10.031	180	312817	50.00	ug/L	100
120) 1,2,4-trichlorobenzene	10.378	180	277477	50.00	ug/L	100
121) hexachlorobutadiene	10.445	225	94479	50.00	ug/L	100
122) naphthalene	10.536	128	1009565	50.00	ug/L	100
123) 1,2,3-trichlorobenzene	10.652	180	256799	50.00	ug/L	100
124) 2-methylnaphthalene	11.132	142	209203	25.00	ug/L	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
Data File : 1R02829.D  
Acq On : 29 Mar 2023 2:37 am  
Operator : PrashanS  
Sample : ICC0091-50  
Misc : MS67262,VLR0091,5,,,1  
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Mar 29 12:11:30 2023  
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
QLast Update: Wed Mar 29 12:11:26 2023  
Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02831.D  
 Acq On : 29 Mar 2023 3:10 am  
 Operator : PrashanS  
 Sample : IC0091-100  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Mar 29 12:37:33 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:37:22 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.497	65	580217	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	826610	50.00	ug/L	# 0.00
52) 1,4-difluorobenzene	5.657	114	1081223	50.00	ug/L	0.00
73) chlorobenzene-d5	7.817	117	1320491	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	764898	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.024	113	374444	52.34	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	104.68%		
53) 1,2-dichloroethane-d4 (s)	5.310	65	359641	59.61	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	119.22%		
74) toluene-d8 (s)	6.801	98	1960065	54.10	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	108.20%		
98) 4-bromofluorobenzene (s)	8.571	174	612793	48.54	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	97.08%		
<b>Target Compounds</b>						
				Qvalue		
2) ethanol	2.700	45	925310	8285.95	ug/L	97
3) tertiary butyl alcohol	3.576	59	718419	485.00	ug/L	98
4) 1,4-dioxane	6.125	88	292363	2713.62	ug/L	97
6) chlorodifluoromethane	1.654	51	581613	81.98	ug/L	100
7) dichlorodifluoromethane	1.636	85	564712	92.99	ug/L	97
8) chloromethane	1.806	50	793784	73.06	ug/L	98
9) vinyl chloride	1.897	62	814745	79.55	ug/L	97
10) 1,3-butadiene	1.934	54	650103	86.17	ug/L	99
11) bromomethane	2.183	94	568909	102.40	ug/L	99
12) chloroethane	2.293	64	504842	102.53	ug/L	100
13) trichlorofluoromethane	2.536	101	734929	102.13	ug/L	97
14) ethyl ether	2.804	74	303344	100.38	ug/L	91
15) acrolein	2.944	56	172633	98.45	ug/L	98
16) freon 113	3.041	151	389034	94.41	ug/L	95
17) 1,1-dichloroethene	3.041	96	454166	96.47	ug/L	94
18) acetone	3.071	58	442272	379.55	ug/L	98
19) acetonitrile	3.333	41	1411840	883.92	ug/L	97
20) iodomethane	3.187	142	405218	168.00	ug/L	99
21) carbon disulfide	3.254	76	1298888	92.54	ug/L	100
22) methylene chloride	3.497	84	501151	97.08	ug/L	95
23) methyl acetate	3.363	74	148603	96.49	ug/L	# 75
24) methyl tert butyl ether	3.728	73	1581084	109.93	ug/L	98
25) trans-1,2-dichloroethene	3.747	96	548252	106.06	ug/L	98
26) hexane	3.984	56	405380	89.32	ug/L	89
27) di-isopropyl ether	4.142	45	2410798	117.98	ug/L	98
28) 2-butanone	4.598	72	616497	465.87	ug/L	97
29) 1,1-dichloroethane	4.130	63	1146351	119.06	ug/L	98
30) chloroprene	4.191	53	1026073	126.00	ug/L	97
31) acrylonitrile	3.704	53	395669	104.45	ug/L	96
32) vinyl acetate	4.124	86	142995	108.60	ug/L	# 82
33) ethyl tert-butyl ether	4.452	59	1835242	114.63	ug/L	96
34) ethyl acetate	4.623	45	212133	113.90	ug/L	# 89
35) 2,2-dichloropropane	4.623	77	946746	139.79	ug/L	91
36) cis-1,2-dichloroethene	4.617	96	713274	126.55	ug/L	93
37) propionitrile	4.659	54	1987901	1164.48	ug/L	81
38) methyl acrylate	4.665	85	176664	130.45	ug/L	94
39) methacrylonitrile	4.781	67	436639	119.57	ug/L	97
40) bromochloromethane	4.817	128	348106	131.82	ug/L	# 83
41) tetrahydrofuran	4.823	42	482225	101.15	ug/L	95
42) chloroform	4.890	83	1055514	111.68	ug/L	92

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02831.D  
 Acq On : 29 Mar 2023 3:10 am  
 Operator : PrashanS  
 Sample : IC0091-100  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Mar 29 12:37:33 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:37:22 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) tert-Butyl Formate	4.915	59	603718	120.12	ug/L	96
45) 1,1,1-trichloroethane	5.042	97	842780	103.14	ug/L #	62
46) cyclohexane	5.097	84	804929	105.52	ug/L	90
47) isobutyl alcohol	5.219	43	650128	892.47	ug/L	98
48) 1,1-dichloropropene	5.164	75	725990	105.43	ug/L	97
49) carbon tetrachloride	5.164	117	745011	112.88	ug/L	99
50) tert-amyl alcohol	5.316	73	439659	531.14	ug/L	97
51) isopropyl acetate	5.334	87	205837	112.93	ug/L #	92
54) n-butyl alcohol	5.754	56	2396184	6129.32	ug/L	98
55) 2,2,4-trimethylpentane	5.414	57	1639193	110.30	ug/L	98
56) benzene	5.328	78	2524955	121.53	ug/L	99
57) tert-amyl methyl ether	5.420	73	1824924	124.79	ug/L	98
58) heptane	5.547	57	336740	98.36	ug/L	87
59) 1,2-dichloroethane	5.365	62	744969	112.23	ug/L	98
60) ethyl acrylate	5.900	55	1295793	120.58	ug/L	99
61) trichloroethene	5.852	95	652692	126.66	ug/L	98
62) 2-chloroethyl vinyl ether	6.472	63	3054845	696.12	ug/L	98
63) methyl methacrylate	6.089	100	239961	141.43	ug/L #	82
64) methylcyclohexane	6.046	83	993317	119.46	ug/L	98
65) 1,2-dichloropropane	6.064	63	713812	124.52	ug/L	97
66) dibromomethane	6.137	93	458720	131.23	ug/L	94
67) bromodichloromethane	6.265	83	701235	111.24	ug/L	98
68) 2-nitropropane	6.429	41	342779	101.28	ug/L	93
69) epichlorohydrin	6.515	57	660846	623.24	ug/L	98
70) cis-1,3-dichloropropene	6.600	75	1004197	125.07	ug/L	94
71) 4-methyl-2-pentanone	6.697	58	2041774	519.87	ug/L	99
72) isoamyl alcohol	6.728	70	928768	2589.70	ug/L	98
75) toluene	6.855	92	1857330	109.80	ug/L	99
76) ethyl methacrylate	7.032	69	1089514	119.17	ug/L	93
77) trans-1,3-dichloropropene	7.007	75	1083074	121.99	ug/L	91
78) 1,1,2-trichloroethane	7.153	83	523826	96.81	ug/L	98
79) tetrachloroethylene	7.220	164	580654	84.74	ug/L	95
80) 2-hexanone	7.287	58	2219778	407.17	ug/L	97
81) 1,3-dichloropropane	7.269	76	939405	97.64	ug/L	96
82) butyl acetate	7.360	56	734720	102.49	ug/L	97
83) dibromochloromethane	7.421	129	744276	115.26	ug/L	98
84) 1,2-dibromoethane	7.512	107	614495	96.13	ug/L	98
85) n-butyl ether	7.865	57	3193299	104.96	ug/L	100
86) chlorobenzene	7.835	112	2100361	113.40	ug/L	96
87) 1,1,1,2-tetrachloroethane	7.890	131	827877	131.97	ug/L	97
88) ethylbenzene	7.890	91	3548587	114.92	ug/L	98
89) m,p-xylene	7.975	106	2792845	236.38	ug/L	95
90) o-xylene	8.224	91	2861186	121.92	ug/L	98
91) styrene	8.236	104	2448886	126.13	ug/L	95
92) butyl acrylate	8.163	55	1756277	111.32	ug/L	98
93) n-amyl acetate	8.297	70	645058	104.94	ug/L	98
94) isopropylbenzene	8.449	105	3466020	127.38	ug/L	98
95) bromoform	8.364	173	702390	140.29	ug/L	98
96) cis-1,4-dichloro-2-butene	8.486	88	513020	136.92	ug/L	97
99) 1,1,2,2-tetrachloroethane	8.638	83	1027470	86.42	ug/L	99
100) trans-1,4-dichloro-2-b...	8.656	53	364644	92.31	ug/L	87
101) 1,2,3-trichloropropane	8.680	110	338060	98.16	ug/L	87
102) bromobenzene	8.662	156	916762	102.19	ug/L	88
103) n-propylbenzene	8.705	91	3871406	98.01	ug/L	99
104) 2-chlorotoluene	8.772	126	831391	103.54	ug/L	97
105) 4-chlorotoluene	8.845	91	2391026	99.17	ug/L	97
106) 1,3,5-trimethylbenzene	8.808	105	2830885	106.19	ug/L	98
107) tert-butylbenzene	9.003	119	2436622	104.95	ug/L	97

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02831.D  
 Acq On : 29 Mar 2023 3:10 am  
 Operator : PrashanS  
 Sample : IC0091-100  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Mar 29 12:37:33 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:37:22 2023  
 Response via : Initial Calibration

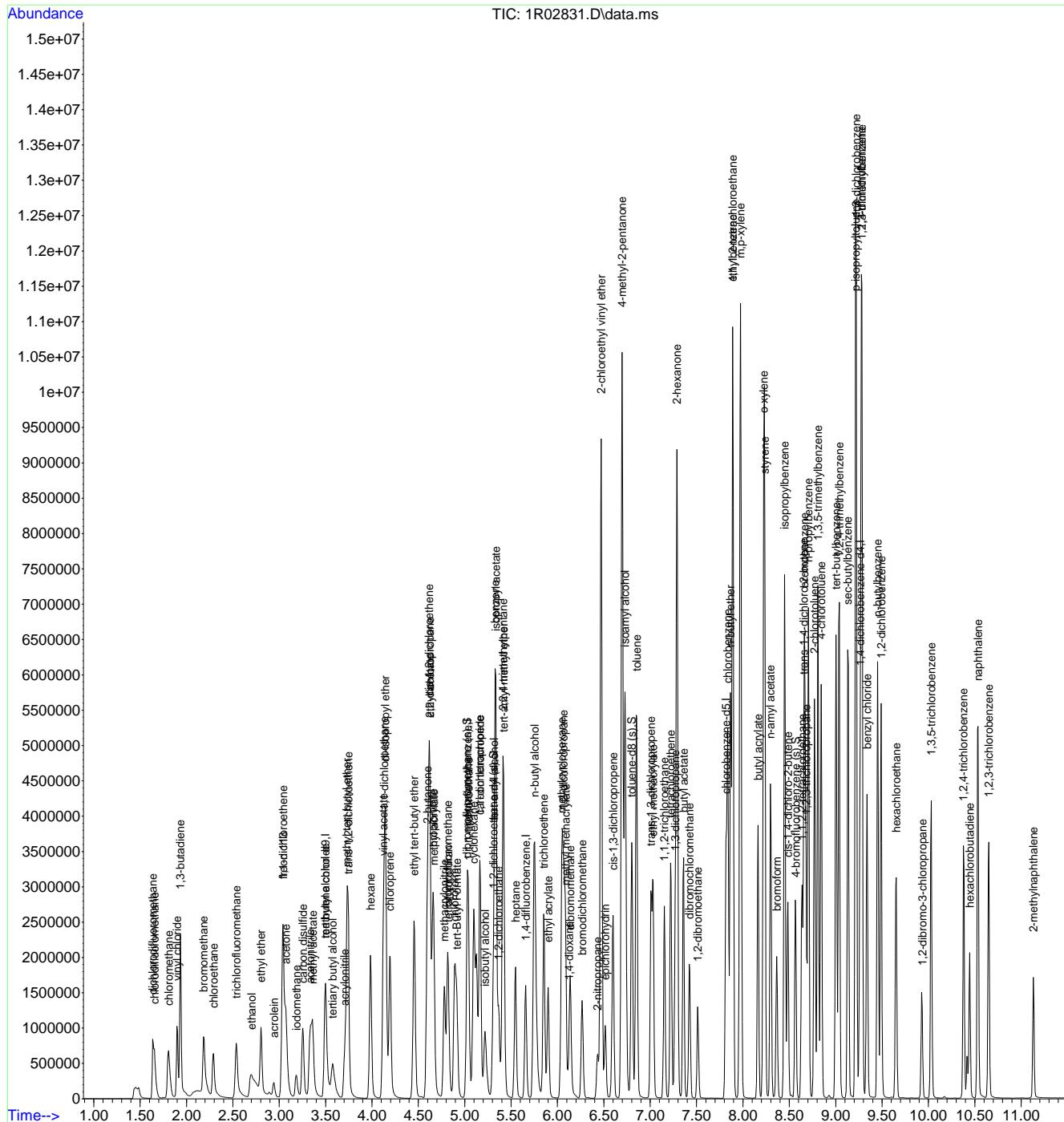
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
108) 1,2,4-trimethylbenzene	9.039	105	2993354	111.40	ug/L	98
109) sec-butylbenzene	9.131	105	3399213	107.31	ug/L	96
110) p-isopropyltoluene	9.216	119	3219822	112.82	ug/L	99
111) 1,2,3-trimethylbenzene	9.283	105	3029099	105.00	ug/L	99
112) 1,3-dichlorobenzene	9.222	146	1885416	111.95	ug/L	98
113) 1,4-dichlorobenzene	9.277	146	1826198	104.59	ug/L	98
114) 1,2-dichlorobenzene	9.490	146	1707172	109.63	ug/L	98
115) benzyl chloride	9.337	91	2137870	100.08	ug/L	99
116) n-butylbenzene	9.453	92	1265490	95.25	ug/L	96
117) hexachloroethane	9.654	201	376184	146.73	ug/L	96
118) 1,2-dibromo-3-chloropr...	9.934	157	298421	97.03	ug/L	90
119) 1,3,5-trichlorobenzene	10.031	180	967568	97.57	ug/L	94
120) 1,2,4-trichlorobenzene	10.378	180	796743	93.65	ug/L	98
121) hexachlorobutadiene	10.445	225	247951	80.48	ug/L	96
122) naphthalene	10.536	128	3170506	105.22	ug/L	97
123) 1,2,3-trichlorobenzene	10.651	180	809286	96.75	ug/L	90
124) 2-methylnaphthalene	11.132	142	569499	42.13	ug/L	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02831.D  
 Acq On : 29 Mar 2023 3:10 am  
 Operator : Prashans  
 Sample : IC0091-100  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Mar 29 12:37:33 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:37:22 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02833.D  
 Acq On : 29 Mar 2023 3:42 am  
 Operator : PrashanS  
 Sample : IC0091-200  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 30 09:19:40 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) tert butyl alcohol-d9	3.509	65	882700	500.00	ug/L	0.02
5) pentafluorobenzene	5.030	168	1195366	50.00	ug/L	# 0.00
52) 1,4-difluorobenzene	5.657	114	1507228	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	1606818	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	850440	50.00	ug/L	0.00
<b>System Monitoring Compounds</b>						
44) dibromofluoromethane (s)	5.018	113	515145	49.58	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	99.16%		
53) 1,2-dichloroethane-d4 (s)	5.304	65	516038	58.89	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	117.78%		
74) toluene-d8 (s)	6.800	98	2165015	48.79	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	97.58%		
98) 4-bromofluorobenzene (s)	8.565	174	661717	47.55	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	95.10%		
<b>Target Compounds</b>						
				Qvalue		
2) ethanol	2.700	45	2061487	12946.57	ug/L	99
3) tertiary butyl alcohol	3.582	59	1968848	889.13	ug/L	99
4) 1,4-dioxane	6.125	88	795174	4817.83	ug/L	96
6) chlorodifluoromethane	1.648	51	1269872	139.23	ug/L	98
7) dichlorodifluoromethane	1.629	85	1176988	139.61	ug/L	98
8) chloromethane	1.794	50	1667723	196.67	ug/L	99
9) vinyl chloride	1.891	62	1680007	122.07	ug/L	97
10) 1,3-butadiene	1.927	54	1359352	132.17	ug/L	97
12) chloroethane	2.268	64	1086712	156.76	ug/L	95
13) trichlorofluoromethane	2.518	101	2326690	220.52	ug/L	96
14) ethyl ether	2.797	74	829133	190.73	ug/L	90
15) acrolein	2.931	56	467658	186.92	ug/L	99
16) freon 113	3.029	151	1237823	208.12	ug/L	96
17) 1,1-dichloroethene	3.029	96	1309964	193.83	ug/L	98
18) acetone	3.065	58	1042887	638.58	ug/L	98
19) acetonitrile	3.333	41	3619820	1645.32	ug/L	98
21) carbon disulfide	3.241	76	3555285	179.12	ug/L	98
22) methylene chloride	3.491	84	1365790	185.32	ug/L	99
23) methyl acetate	3.357	74	421333	191.31	ug/L	# 72
24) methyl tert butyl ether	3.728	73	4246623	201.76	ug/L	96
25) trans-1,2-dichloroethene	3.740	96	1454709	193.87	ug/L	93
26) hexane	3.978	56	1089137	171.22	ug/L	92
27) di-isopropyl ether	4.142	45	5479635	183.48	ug/L	95
28) 2-butanone	4.598	72	1321722	688.53	ug/L	94
29) 1,1-dichloroethane	4.124	63	2576864	182.95	ug/L	98
30) chloroprene	4.184	53	2465246	203.11	ug/L	99
31) acrylonitrile	3.704	53	1001148	183.61	ug/L	99
32) vinyl acetate	4.118	86	375969	195.66	ug/L	# 75
33) ethyl tert-butyl ether	4.452	59	4824731	204.53	ug/L	96
34) ethyl acetate	4.616	45	462052	171.63	ug/L	# 37
35) 2,2-dichloropropane	4.616	77	2107208	204.40	ug/L	90
36) cis-1,2-dichloroethene	4.610	96	1671784	199.30	ug/L	87
37) propionitrile	4.659	54	4301422	1736.21	ug/L	70
38) methyl acrylate	4.665	85	379839	187.52	ug/L	99
39) methacrylonitrile	4.781	67	967997	180.77	ug/L	98
40) bromochloromethane	4.811	128	833154	209.60	ug/L	# 80
41) tetrahydrofuran	4.823	42	1016764	152.26	ug/L	96
42) chloroform	4.890	83	2625003	190.44	ug/L	93
43) tert-Butyl Formate	4.908	59	1594133	212.30	ug/L	92
45) 1,1,1-trichloroethane	5.030	97	2404621	202.41	ug/L	# 59

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02833.D  
 Acq On : 29 Mar 2023 3:42 am  
 Operator : PrashanS  
 Sample : IC0091-200  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 30 09:19:40 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cyclohexane	5.097	84	2080024	188.61	ug/L	86
47) isobutyl alcohol	5.225	43	1695638	1682.40	ug/L	99
48) 1,1-dichloropropene	5.158	75	1876987	188.56	ug/L	96
49) carbon tetrachloride	5.158	117	2310185	233.26	ug/L	98
50) tert-amyl alcohol	5.322	73	1110721	929.04	ug/L	97
51) isopropyl acetate	5.334	87	499172	187.60	ug/L #	93
54) n-butyl alcohol	5.760	56	5886643	10403.80	ug/L	96
55) 2,2,4-trimethylpentane	5.407	57	4103737	196.25	ug/L	93
56) benzene	5.328	78	6049069	203.57	ug/L	95
57) tert-amyl methyl ether	5.413	73	4551800	215.44	ug/L	98
58) heptane	5.541	57	886462	187.58	ug/L	92
59) 1,2-dichloroethane	5.365	62	2047965	216.37	ug/L	91
60) ethyl acrylate	5.900	55	2971411	194.09	ug/L	99
61) trichloroethene	5.851	95	1681215	224.25	ug/L	92
62) 2-chloroethyl vinyl ether	6.472	63	6952646	1079.45	ug/L	96
63) methyl methacrylate	6.083	100	548056	216.24	ug/L #	80
64) methylcyclohexane	6.046	83	2438900	205.34	ug/L	96
65) 1,2-dichloropropane	6.064	63	1553088	189.77	ug/L	98
66) dibromomethane	6.137	93	1003123	198.33	ug/L	89
67) bromodichloromethane	6.265	83	1972872	218.80	ug/L	96
68) 2-nitropropane	6.435	41	825362	194.21	ug/L	91
69) epichlorohydrin	6.514	57	1576698	1026.51	ug/L	99
70) cis-1,3-dichloropropene	6.600	75	2496310	215.16	ug/L	91
71) 4-methyl-2-pentanone	6.697	58	4450273	787.97	ug/L #	72
72) isoamyl alcohol	6.733	70	2142268	4072.04	ug/L	98
75) toluene	6.849	92	4174301	200.55	ug/L	98
76) ethyl methacrylate	7.032	69	2212076	194.82	ug/L	93
77) trans-1,3-dichloropropene	7.007	75	2353177	211.29	ug/L	88
78) 1,1,2-trichloroethane	7.153	83	1208655	185.69	ug/L	95
79) tetrachloroethene	7.220	164	1643119	200.42	ug/L	97
80) 2-hexanone	7.287	58	4740215	720.96	ug/L	89
81) 1,3-dichloropropane	7.269	76	2233948	192.16	ug/L	91
82) butyl acetate	7.360	56	1503328	174.54	ug/L	97
83) dibromochloromethane	7.427	129	1946621	237.42	ug/L	96
84) 1,2-dibromoethane	7.512	107	1635385	209.96	ug/L	99
85) n-butyl ether	7.865	57	6767772	183.48	ug/L	100
86) chlorobenzene	7.835	112	4860787	211.19	ug/L	93
87) 1,1,2-tetrachloroethane	7.889	131	1942330	238.75	ug/L	97
88) ethylbenzene	7.889	91	7745064	202.49	ug/L	94
89) m,p-xylene	7.975	106	6588021	443.70	ug/L #	74
90) o-xylene	8.224	91	6360791	215.57	ug/L	97
91) styrene	8.236	104	5546974	223.96	ug/L	94
92) butyl acrylate	8.163	55	3692312	190.90	ug/L	98
93) n-amyl acetate	8.297	70	1305004	176.20	ug/L	95
94) isopropylbenzene	8.449	105	7544342	217.86	ug/L	96
95) bromoform	8.364	173	1574479	241.64	ug/L	97
96) cis-1,4-dichloro-2-butene	8.486	88	1044283	215.21	ug/L	95
99) 1,1,2,2-tetrachloroethane	8.638	83	2218826	173.57	ug/L	98
100) trans-1,4-dichloro-2-b...	8.656	53	776697	180.71	ug/L	90
101) 1,2,3-trichloropropane	8.680	110	716248	189.02	ug/L	87
102) bromobenzene	8.662	156	2106140	209.52	ug/L	85
103) n-propylbenzene	8.705	91	8153075	187.56	ug/L	95
104) 2-chlorotoluene	8.771	126	1849722	205.72	ug/L	92
105) 4-chlorotoluene	8.844	91	5209191	195.04	ug/L	96
106) 1,3,5-trimethylbenzene	8.808	105	6142565	205.23	ug/L	98
107) tert-butylbenzene	9.003	119	5485238	209.89	ug/L	95
108) 1,2,4-trimethylbenzene	9.039	105	6457896	212.03	ug/L	95
109) sec-butylbenzene	9.136	105	7290379	204.78	ug/L	98

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02833.D  
 Acq On : 29 Mar 2023 3:42 am  
 Operator : PrashanS  
 Sample : IC0091-200  
 Misc : MS67262,V1R0091,5,,,,1  
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 30 09:19:40 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

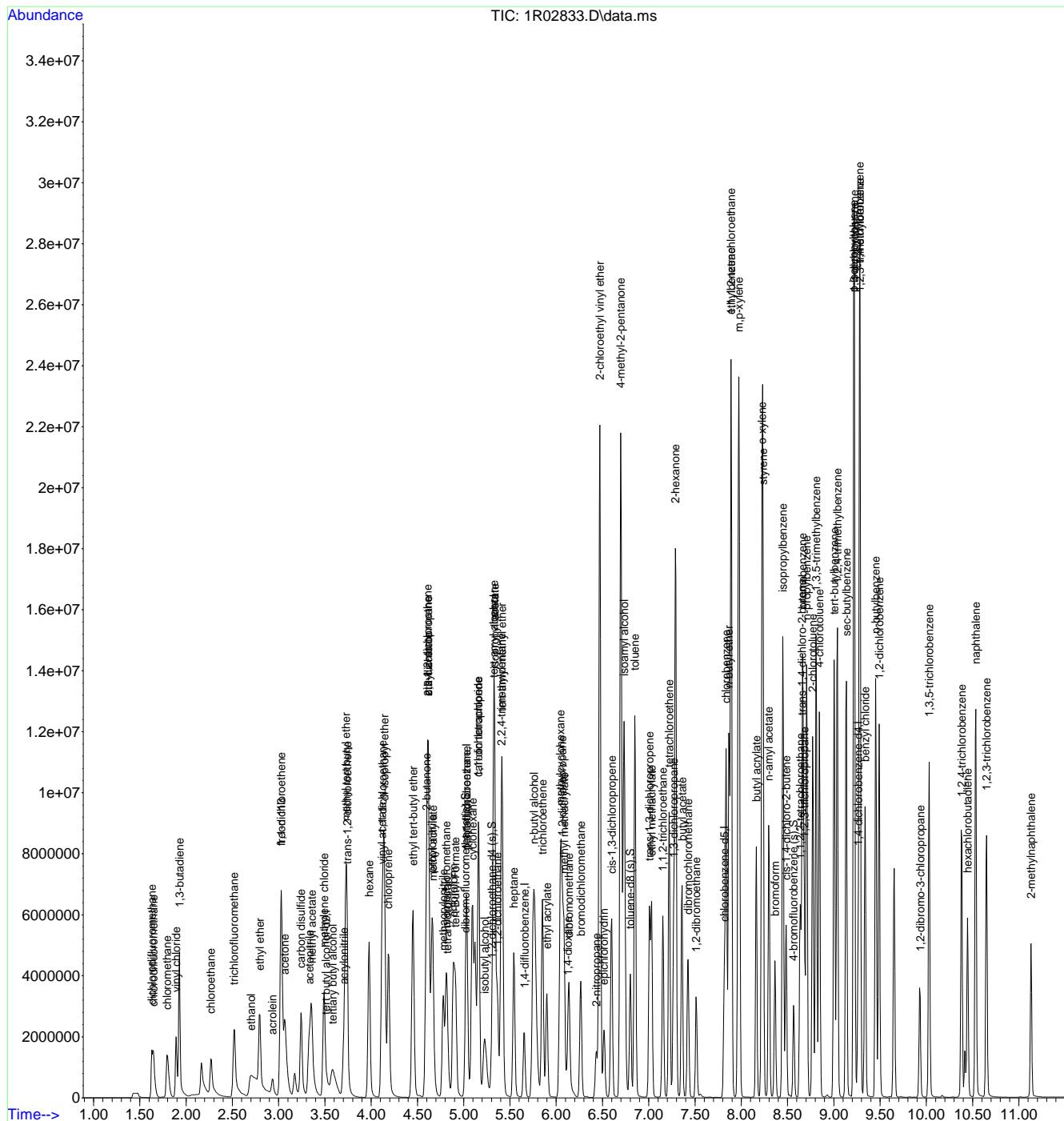
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
110) p-isopropyltoluene	9.216	119	7291852	223.60	ug/L	98
111) 1,2,3-trimethylbenzene	9.282	105	6791719	209.22	ug/L	94
112) 1,3-dichlorobenzene	9.222	146	4527234	234.08	ug/L	95
113) 1,4-dichlorobenzene	9.276	146	4389179	222.18	ug/L	95
114) 1,2-dichlorobenzene	9.489	146	3994504	225.09	ug/L	98
115) benzyl chloride	9.337	91	4896143	205.43	ug/L	98
116) n-butylbenzene	9.453	92	2870969	195.83	ug/L	99
118) 1,2-dibromo-3-chloropr...	9.933	157	783950	226.33	ug/L	92
119) 1,3,5-trichlorobenzene	10.031	180	2607797	232.42	ug/L	97
120) 1,2,4-trichlorobenzene	10.384	180	2187365	229.12	ug/L	91
121) hexachlorobutadiene	10.444	225	785621	230.59	ug/L	96
122) naphthalene	10.536	128	7786221	226.35	ug/L	96
123) 1,2,3-trichlorobenzene	10.651	180	2104372	223.81	ug/L	95
124) 2-methylnaphthalene	11.132	142	1743472	115.97	ug/L	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\v1r91\  
Data File : 1R02833.D  
Acq On : 29 Mar 2023 3:42 am  
Operator : PrashanS  
Sample : IC0091-200  
Misc : MS67262,V1R0091,5,,,1  
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Mar 30 09:19:40 2023  
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
QLast Update: Wed Mar 29 12:39:36 2023  
Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02839.D  
 Acq On : 29 Mar 2023 5:20 am  
 Operator : PrashanS  
 Sample : ICV0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 30 09:17:23 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) tert butyl alcohol-d9	3.491	65	541895	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	704790	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.651	114	989846	50.00	ug/L	0.00
73) chlorobenzene-d5	7.816	117	1009852	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	680734	50.00	ug/L	0.00
<hr/>						
System Monitoring Compounds						
44) dibromofluoromethane (s)	5.018	113	319382	52.14	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	104.28%		
53) 1,2-dichloroethane-d4 (s)	5.298	65	293864	51.06	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	102.12%		
74) toluene-d8 (s)	6.801	98	1544616	55.39	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	110.78%		
98) 4-bromofluorobenzene (s)	8.565	174	438448	39.36	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	78.72%#		
<hr/>						
Target Compounds				Qvalue		
2) ethanol	2.688	45	459363	4699.24	ug/L	97
3) tertiary butyl alcohol	3.564	59	303305	223.12	ug/L	99
4) 1,4-dioxane	6.113	88	119496	1179.34	ug/L	96
7) dichlorodifluoromethane	1.636	85	183816	36.98	ug/L	98
8) chloromethane	1.800	50	283805	36.87	ug/L	99
9) vinyl chloride	1.897	62	295757	36.45	ug/L	97
10) 1,3-butadiene	1.928	54	314988	51.94	ug/L	97
11) bromomethane	2.183	94	155555	34.24	ug/L	95
12) chloroethane	2.287	64	157332	38.49	ug/L	98
13) trichlorofluoromethane	2.536	101	252434	40.58	ug/L	99
14) ethyl ether	2.798	74	125545	48.98	ug/L	92
15) acrolein	2.937	56	68482	46.42	ug/L	97
16) freon 113	3.041	151	162126	46.23	ug/L	91
17) 1,1-dichloroethene	3.035	96	182280	45.74	ug/L	92
18) acetone	3.065	58	175080	181.83	ug/L	98
19) acetonitrile	3.327	41	512809	395.33	ug/L	99
20) iodomethane	3.181	142	124254	44.22	ug/L	98
21) carbon disulfide	3.254	76	554401	47.37	ug/L	99
22) methylene chloride	3.497	84	216157	49.74	ug/L	98
23) methyl acetate	3.357	74	60904	46.90	ug/L	91
24) methyl tert butyl ether	3.728	73	614541	49.52	ug/L	99
25) trans-1,2-dichloroethene	3.740	96	215877	48.80	ug/L	97
26) hexane	3.984	56	122035	32.54	ug/L	90
27) di-isopropyl ether	4.142	45	889753	50.53	ug/L	98
28) 2-butanone	4.592	72	227836	201.30	ug/L	99
29) 1,1-dichloroethane	4.130	63	406879	48.99	ug/L	97
30) chloroprene	4.191	53	353805	49.44	ug/L	98
31) acrylonitrile	3.704	53	160214	49.84	ug/L	98
32) vinyl acetate	4.118	86	47286	41.74	ug/L #	91
33) ethyl tert-butyl ether	4.452	59	681784	49.02	ug/L	98
34) ethyl acetate	4.617	45	81057	51.07	ug/L #	90
35) 2,2-dichloropropane	4.623	77	281864	46.37	ug/L	97
36) cis-1,2-dichloroethene	4.610	96	243401	49.21	ug/L	88
37) propionitrile	4.653	54	716594	490.57	ug/L	91
38) methyl acrylate	4.665	85	63348	53.04	ug/L #	86
39) methacrylonitrile	4.775	67	153561	48.64	ug/L	88
40) bromochloromethane	4.811	128	109147	46.57	ug/L	94
41) tetrahydrofuran	4.817	42	184216	46.79	ug/L	98
42) chloroform	4.890	83	379905	46.75	ug/L	97
43) tert-Butyl Formate	4.909	59	168449	38.05	ug/L	97

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02839.D  
 Acq On : 29 Mar 2023 5:20 am  
 Operator : PrashanS  
 Sample : ICV0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 30 09:17:23 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1,1-trichloroethane	5.036	97	333275	47.58	ug/L	90
46) cyclohexane	5.097	84	272601	41.92	ug/L	93
47) isobutyl alcohol	5.213	43	287877	484.44	ug/L	97
48) 1,1-dichloropropene	5.164	75	292114	49.77	ug/L	99
49) carbon tetrachloride	5.164	117	291631	49.94	ug/L	98
50) tert-amyl alcohol	5.310	73	170901	242.44	ug/L	97
51) isopropyl acetate	5.334	87	76293	48.63	ug/L	96
54) n-butyl alcohol	5.742	56	931242	2506.10	ug/L	98
55) 2,2,4-trimethylpentane	5.407	57	682261	49.68	ug/L	98
56) benzene	5.328	78	929854	47.65	ug/L	97
57) tert-amyl methyl ether	5.413	73	656387	47.30	ug/L	98
58) heptane	5.541	57	140533	45.28	ug/L	94
59) 1,2-dichloroethane	5.365	62	286001	46.01	ug/L	99
60) ethyl acrylate	5.894	55	494789	49.21	ug/L	98
61) trichloroethene	5.851	95	236609	48.06	ug/L	98
62) 2-chloroethyl vinyl ether	6.466	63	1116977	264.06	ug/L	98
63) methyl methacrylate	6.083	100	102114	61.35	ug/L	91
64) methylcyclohexane	6.046	83	401438	51.47	ug/L	91
65) 1,2-dichloropropane	6.064	63	266827	49.65	ug/L	97
66) dibromomethane	6.131	93	163151	49.12	ug/L	97
67) bromodichloromethane	6.265	83	275588	46.54	ug/L	98
68) 2-nitropropane	6.429	41	131744	47.20	ug/L	99
69) epichlorohydrin	6.515	57	243681	241.57	ug/L	96
70) cis-1,3-dichloropropene	6.594	75	365894	48.02	ug/L	97
71) 4-methyl-2-pentanone	6.691	58	807155	217.62	ug/L	94
72) isoamyl alcohol	6.721	70	402961	1166.31	ug/L	95
75) toluene	6.849	92	684351	52.32	ug/L	94
76) ethyl methacrylate	7.032	69	330579	46.32	ug/L	97
77) trans-1,3-dichloropropene	7.007	75	351480	50.21	ug/L	96
78) 1,1,2-trichloroethane	7.153	83	195104	47.69	ug/L	95
79) tetrachloroethene	7.220	164	307562	59.69	ug/L	96
80) 2-hexanone	7.287	58	843165	204.05	ug/L	97
81) 1,3-dichloropropane	7.269	76	357984	49.00	ug/L	96
82) butyl acetate	7.354	56	281792	52.06	ug/L	92
83) dibromochloromethane	7.421	129	253956	49.28	ug/L	92
84) 1,2-dibromoethane	7.512	107	240088	49.04	ug/L	99
85) n-butyl ether	7.865	57	1227617	52.96	ug/L	100
86) chlorobenzene	7.835	112	682909	47.21	ug/L	99
87) 1,1,1,2-tetrachloroethane	7.889	131	287665	56.26	ug/L	95
88) ethylbenzene	7.889	91	1321713	54.98	ug/L	99
89) m,p-xylene	7.975	106	1064697	114.10	ug/L	99
90) o-xylene	8.224	91	1084446	58.48	ug/L	100
91) styrene	8.236	104	864425	55.53	ug/L	98
92) butyl acrylate	8.163	55	639774	52.63	ug/L	98
93) n-amyl acetate	8.297	70	275375	59.16	ug/L	96
94) isopropylbenzene	8.449	105	1316290	60.48	ug/L	99
95) bromoform	8.364	173	234043	57.15	ug/L	98
96) cis-1,4-dichloro-2-butene	8.480	88	177504	58.21	ug/L	92
99) 1,1,2,2-tetrachloroethane	8.632	83	406159	39.69	ug/L	93
100) trans-1,4-dichloro-2-b...	8.656	53	144227	41.92	ug/L	89
101) 1,2,3-trichloropropane	8.674	110	120316	39.67	ug/L #	84
102) bromobenzene	8.662	156	301243	37.44	ug/L	95
103) n-propylbenzene	8.705	91	1426638	41.00	ug/L	98
104) 2-chlorotoluene	8.772	126	288795	40.13	ug/L	88
105) 4-chlorotoluene	8.845	91	972752	45.50	ug/L	97
106) 1,3,5-trimethylbenzene	8.808	105	1093153	45.63	ug/L	95
107) tert-butylbenzene	9.003	119	848816	40.58	ug/L	96
108) 1,2,4-trimethylbenzene	9.033	105	1228710	50.40	ug/L	98

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02839.D  
 Acq On : 29 Mar 2023 5:20 am  
 Operator : PrashanS  
 Sample : ICV0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 30 09:17:23 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

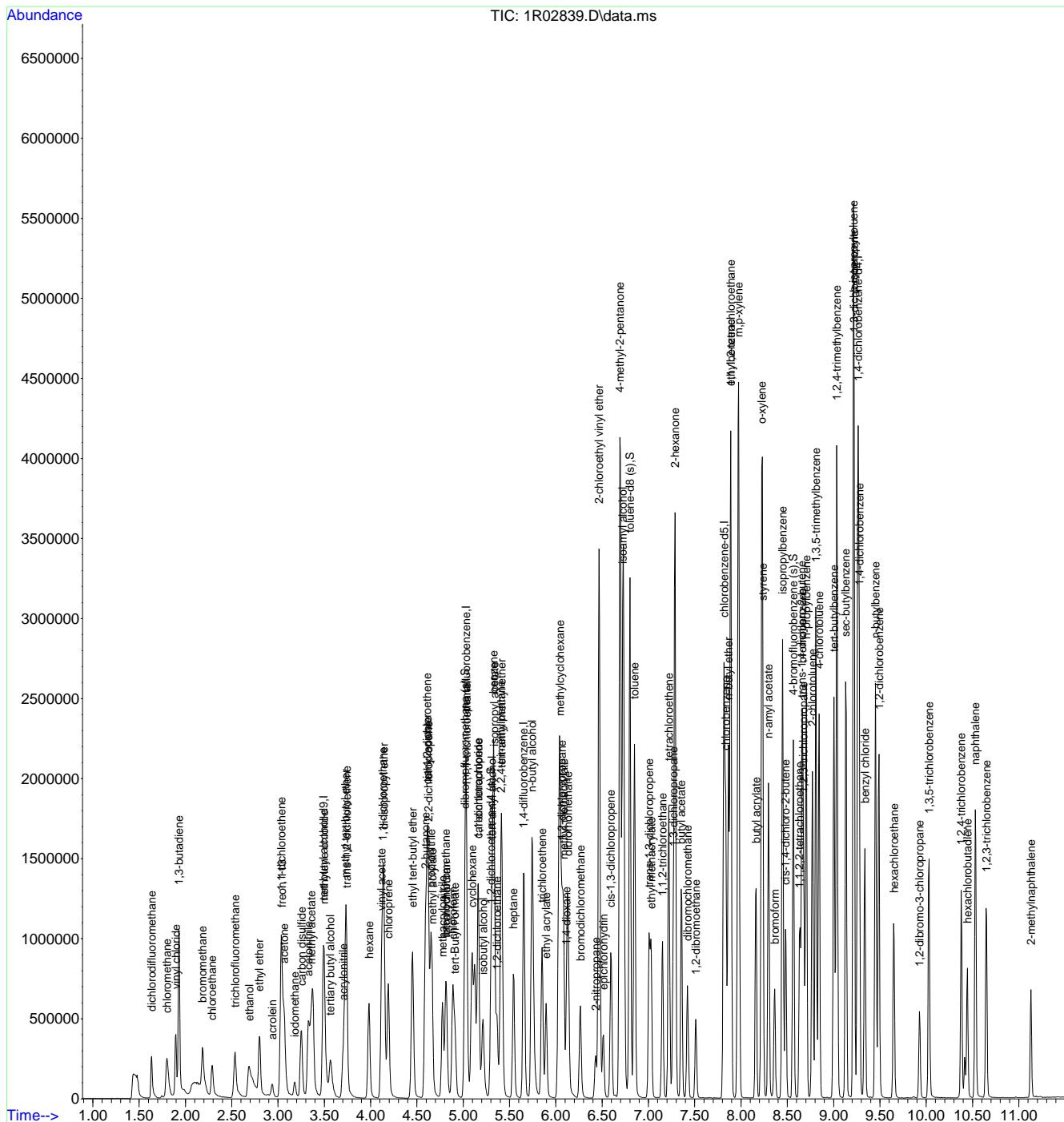
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
109) sec-butylbenzene	9.131	105	1291072	45.31	ug/L	97
110) p-isopropyltoluene	9.216	119	1299783	49.79	ug/L	99
112) 1,3-dichlorobenzene	9.222	146	750967	48.51	ug/L	99
113) 1,4-dichlorobenzene	9.277	146	750246	47.44	ug/L	97
114) 1,2-dichlorobenzene	9.489	146	663501	46.71	ug/L	99
115) benzyl chloride	9.337	91	774415	40.59	ug/L	99
116) n-butylbenzene	9.453	92	525347	44.77	ug/L	100
117) hexachloroethane	9.654	201	125448	50.50	ug/L	97
118) 1,2-dibromo-3-chloropr...	9.927	157	100467	36.24	ug/L	95
119) 1,3,5-trichlorobenzene	10.031	180	335863	37.49	ug/L	99
120) 1,2,4-trichlorobenzene	10.378	180	283395	37.58	ug/L	95
121) hexachlorobutadiene	10.445	225	100895	36.84	ug/L	95
122) naphthalene	10.530	128	1018149	36.98	ug/L	97
123) 1,2,3-trichlorobenzene	10.651	180	272270	36.49	ug/L	96
124) 2-methylnaphthalene	11.132	142	226899	18.86	ug/L	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\v1r91\  
Data File : 1R02839.D  
Acq On : 29 Mar 2023 5:20 am  
Operator : PrashanS  
Sample : ICV0091-50  
Misc : MS67262,V1R0091,5,,,1  
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 30 09:17:23 2023  
Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
QLast Update: Wed Mar 29 12:39:36 2023  
Response via: Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02841.D  
 Acq On : 29 Mar 2023 5:53 am  
 Operator : PrashanS  
 Sample : ICV0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 15 Sample Multiplier: 1

Quant Time: Mar 30 09:10:23 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

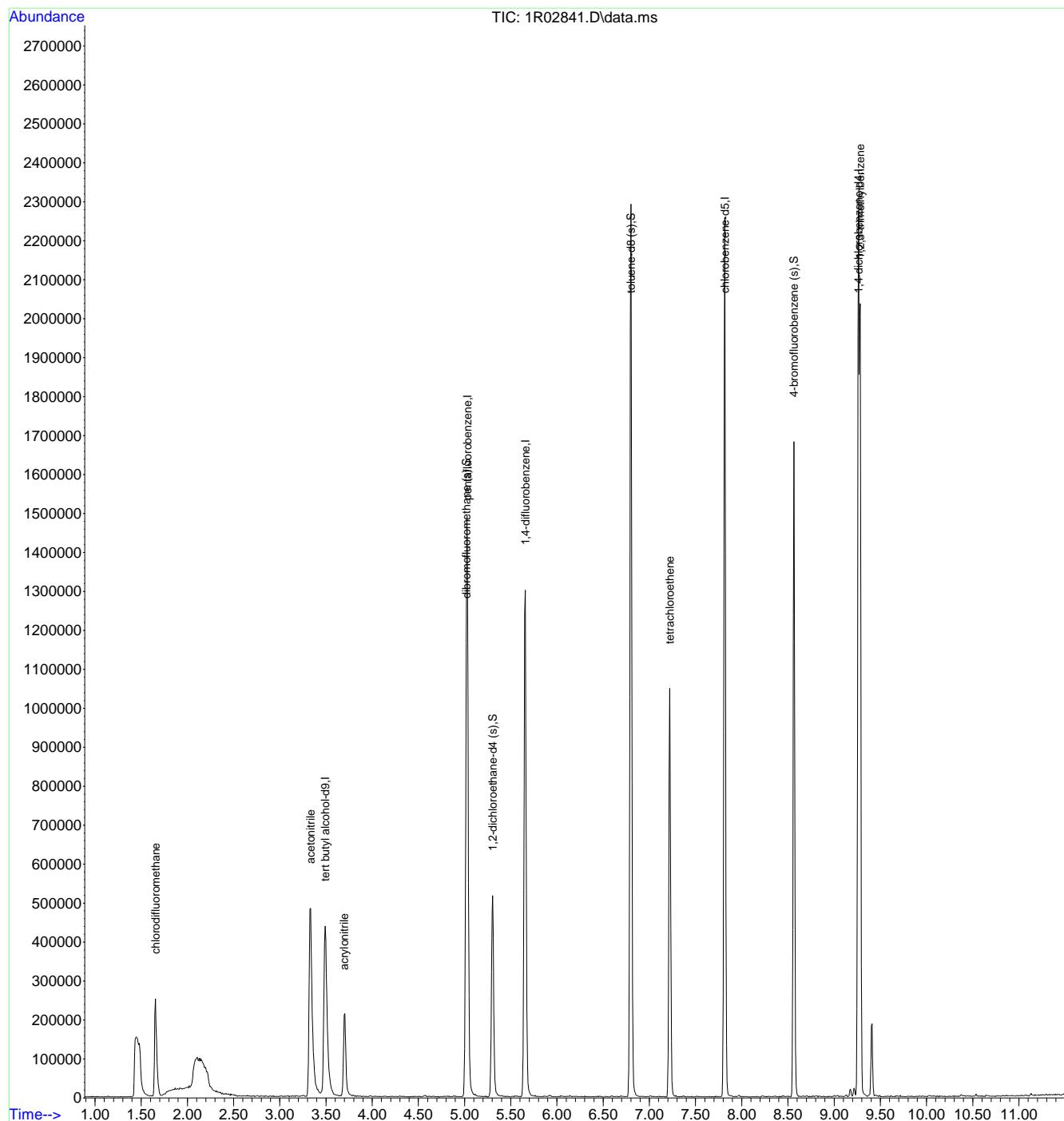
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) tert butyl alcohol-d9	3.491	65	565330	500.00	ug/L	0.00
5) pentafluorobenzene	5.030	168	654176	50.00	ug/L	0.00
52) 1,4-difluorobenzene	5.657	114	923767	50.00	ug/L	0.00
73) chlorobenzene-d5	7.817	117	886053	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.264	152	399541	50.00	ug/L	0.00
<hr/>						
System Monitoring Compounds						
44) dibromofluoromethane (s)	5.018	113	288893	50.81	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	101.62%		
53) 1,2-dichloroethane-d4 (s)	5.304	65	261660	48.72	ug/L	0.00
Spiked Amount 50.000	Range 81 - 124		Recovery =	97.44%		
74) toluene-d8 (s)	6.801	98	1182124	48.31	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	96.62%		
98) 4-bromofluorobenzene (s)	8.565	174	345539	52.85	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	105.70%		
<hr/>						
Target Compounds						
					Qvalue	
6) chlorodifluoromethane	1.654	51	282731	56.64	ug/L	99
19) acetonitrile	3.333	41	650255	540.07	ug/L	99
31) acrylonitrile	3.698	53	156426	52.42	ug/L	91
79) tetrachloroethene	7.220	164	175720	38.87	ug/L	96
111) 1,2,3-trimethylbenzene	9.283	105	786079	51.54	ug/L	98
<hr/>						

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02841.D  
 Acq On : 29 Mar 2023 5:53 am  
 Operator : Prashans  
 Sample : ICV0091-50  
 Misc : MS67262,V1R0091,5,,,1  
 ALS Vial : 15 Sample Multiplier: 1

Quant Time: Mar 30 09:10:23 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02847.D  
 Acq On : 29 Mar 2023 5:25 pm  
 Operator : PrashanS  
 Sample : icv0091-50  
 Misc : MS67262,V2R0091,5,,,1  
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Mar 30 09:22:51 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

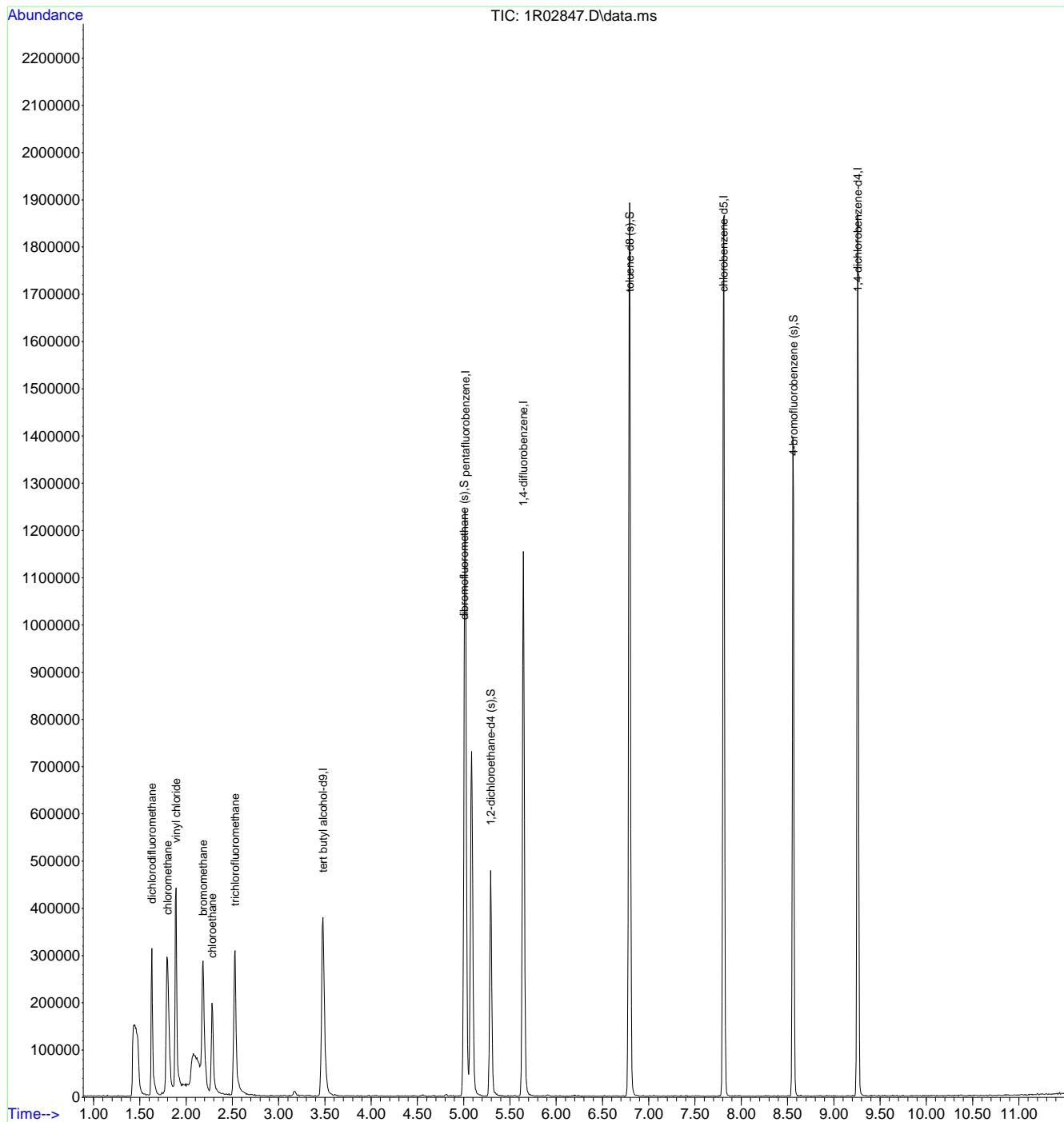
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) tert butyl alcohol-d9	3.479	65	445107	500.00	ug/L	-0.01
5) pentafluorobenzene	5.018	168	563508	50.00	ug/L	-0.01
52) 1,4-difluorobenzene	5.645	114	785396	50.00	ug/L	-0.01
73) chlorobenzene-d5	7.810	117	733917	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.258	152	342913	50.00	ug/L	0.00
<hr/>						
System Monitoring Compounds						
44) dibromofluoromethane (s)	5.006	113	244740	49.97	ug/L	-0.01
Spiked Amount 50.000	Range 80 - 120		Recovery =	99.94%		
53) 1,2-dichloroethane-d4 (s)	5.292	65	212176	46.47	ug/L	-0.01
Spiked Amount 50.000	Range 81 - 124		Recovery =	92.94%		
74) toluene-d8 (s)	6.794	98	980663	48.39	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	96.78%		
98) 4-bromofluorobenzene (s)	8.565	174	293509	52.31	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	104.62%		
<hr/>						
Target Compounds						
7) dichlorodifluoromethane	1.629	85	196088	49.34	ug/L	98
8) chloromethane	1.800	50	339197	58.02	ug/L	99
9) vinyl chloride	1.891	62	328739	50.67	ug/L	97
11) bromomethane	2.183	94	148058	40.77	ug/L	99
12) chloroethane	2.280	64	140189	42.90	ug/L	98
13) trichlorofluoromethane	2.530	101	234470	47.14	ug/L	97
<hr/>						

(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\V1r91\  
 Data File : 1R02847.D  
 Acq On : 29 Mar 2023 5:25 pm  
 Operator : Prashans  
 Sample : icv0091-50  
 Misc : MS67262,V2R0091,5,,,1  
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Mar 30 09:22:51 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration



## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03013.D  
 Acq On : 4 Apr 2023 9:38 am  
 Operator : nickw  
 Sample : cc91-20  
 Misc : MS67909,V1R0100,5,,,1  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 04 09:50:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	3.479	65	649271	500.00	ug/L	-0.01
5) pentafluorobenzene	5.018	168	688798	50.00	ug/L	-0.01
52) 1,4-difluorobenzene	5.645	114	1060871	50.00	ug/L	-0.01
73) chlorobenzene-d5	7.810	117	1047884	50.00	ug/L	0.00
97) 1,4-dichlorobenzene-d4	9.258	152	522573	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) dibromofluoromethane (s)	5.012	113	324490	54.20	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	108.40%		
53) 1,2-dichloroethane-d4 (s)	5.292	65	320005	51.88	ug/L	-0.01
Spiked Amount 50.000	Range 81 - 124		Recovery =	103.76%		
74) toluene-d8 (s)	6.794	98	1359764	46.99	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	93.98%		
98) 4-bromofluorobenzene (s)	8.559	174	435345	50.91	ug/L	0.00
Spiked Amount 50.000	Range 80 - 120		Recovery =	101.82%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) ethanol	2.676	45	249673	2131.73	ug/L	97
3) tertiary butyl alcohol	3.558	59	152459	93.60	ug/L	99
4) 1,4-dioxane	6.107	88	55787	459.52	ug/L	94
6) chlorodifluoromethane	1.648	51	86104	16.38	ug/L	97
7) dichlorodifluoromethane	1.629	85	101245	20.84	ug/L	99
8) chloromethane	1.794	50	130491	16.34	ug/L	98
9) vinyl chloride	1.891	62	112717	14.21	ug/L	99
10) 1,3-butadiene	1.928	54	72006	12.15	ug/L	95
11) bromomethane	2.183	94	66569	14.99	ug/L	98
12) chloroethane	2.287	64	64073	16.04	ug/L	98
13) trichlorofluoromethane	2.530	101	132897	21.86	ug/L	98
14) ethyl ether	2.791	74	49719	19.85	ug/L	90
15) acrolein	2.931	56	30412	21.10	ug/L	89
16) freon 113	3.029	151	64322	18.77	ug/L	98
17) 1,1-dichloroethene	3.029	96	77879	20.00	ug/L	97
18) acetone	3.059	58	96156	102.18	ug/L	95
19) acetonitrile	3.321	41	282983	223.22	ug/L	98
20) iodomethane	3.175	142	28127	12.53	ug/L	98
21) carbon disulfide	3.242	76	203823	17.82	ug/L	98
22) methylene chloride	3.485	84	88078	20.74	ug/L	96
23) methyl acetate	3.351	74	26223	20.66	ug/L #	81
24) methyl tert butyl ether	3.716	73	252377	20.81	ug/L	98
25) trans-1,2-dichloroethene	3.734	96	84681	19.59	ug/L	95
26) hexane	3.972	56	70014	19.10	ug/L	96
27) di-isopropyl ether	4.130	45	347350	20.18	ug/L	99
28) 2-butanone	4.580	72	102032	92.24	ug/L	99
29) 1,1-dichloroethane	4.118	63	157131	19.36	ug/L	97
30) chloroprene	4.178	53	140080	20.03	ug/L	98
31) acrylonitrile	3.692	53	66317	21.11	ug/L	98
32) vinyl acetate	4.105	86	23132	20.89	ug/L #	84
33) ethyl tert-butyl ether	4.440	59	293401	21.59	ug/L	95
34) ethyl acetate	4.604	45	29265	18.86	ug/L #	91
35) 2,2-dichloropropane	4.610	77	124268	20.92	ug/L	93
36) cis-1,2-dichloroethene	4.604	96	99584	20.60	ug/L	90
37) propionitrile	4.641	54	295196	206.78	ug/L	96
38) methyl acrylate	4.653	85	24121	20.67	ug/L #	82
39) methacrylonitrile	4.769	67	63022	20.42	ug/L	96
40) bromochloromethane	4.805	128	48635	21.23	ug/L #	84
41) tetrahydrofuran	4.811	42	68825	17.89	ug/L	97
42) chloroform	4.878	83	158879	20.00	ug/L	98
43) tert-Butyl Formate	4.902	59	94927	21.94	ug/L	97

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03013.D  
 Acq On : 4 Apr 2023 9:38 am  
 Operator : nickw  
 Sample : cc91-20  
 Misc : MS67909,V1R0100,5,,,1  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 04 09:50:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, Rx1624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1,1-trichloroethane	5.024	97	133252	19.47	ug/L #	5
46) cyclohexane	5.085	84	129351	20.35	ug/L	85
47) isobutyl alcohol	5.201	43	131780	226.91	ug/L	98
48) 1,1-dichloropropene	5.152	75	116349	20.28	ug/L	99
49) carbon tetrachloride	5.152	117	117315	20.56	ug/L	94
50) tert-amyl alcohol	5.298	73	79293	115.10	ug/L	96
51) isopropyl acetate	5.322	87	31551	20.58	ug/L #	91
54) n-butyl alcohol	5.730	56	404525	1015.75	ug/L	98
55) 2,2,4-trimethylpentane	5.401	57	249928	16.98	ug/L	97
56) benzene	5.316	78	357324	17.08	ug/L	97
57) tert-amyl methyl ether	5.407	73	280217	18.84	ug/L	98
58) heptane	5.535	57	55979	16.83	ug/L	89
59) 1,2-dichloroethane	5.353	62	124407	18.67	ug/L	93
60) ethyl acrylate	5.888	55	191145	17.74	ug/L	99
61) trichloroethene	5.839	95	97244	18.43	ug/L	100
62) 2-chloroethyl vinyl ether	6.460	63	402440	88.77	ug/L	100
63) methyl methacrylate	6.070	100	33845	18.97	ug/L #	90
64) methylcyclohexane	6.040	83	141682	16.95	ug/L	99
65) 1,2-dichloropropane	6.052	63	98162	17.04	ug/L	97
66) dibromomethane	6.125	93	63836	17.93	ug/L	98
67) bromodichloromethane	6.253	83	118250	18.63	ug/L	99
68) 2-nitropropane	6.423	41	53289	17.81	ug/L	92
69) epichlorohydrin	6.502	57	111968	103.57	ug/L	98
70) cis-1,3-dichloropropene	6.588	75	151093	18.50	ug/L	96
71) 4-methyl-2-pentanone	6.685	58	288058	72.46	ug/L	98
72) isoamyl alcohol	6.715	70	142804	385.65	ug/L	96
75) toluene	6.843	92	236969	17.46	ug/L	98
76) ethyl methacrylate	7.020	69	143509	19.38	ug/L	99
77) trans-1,3-dichloropropene	7.001	75	143450	19.75	ug/L	97
78) 1,1,2-trichloroethane	7.147	83	77870	18.34	ug/L	95
79) tetrachloroethene	7.214	164	94952	17.76	ug/L	95
80) 2-hexanone	7.281	58	323216	75.38	ug/L	95
81) 1,3-dichloropropane	7.263	76	142783	18.83	ug/L	98
82) butyl acetate	7.348	56	103452	18.42	ug/L	95
83) dibromochloromethane	7.415	129	105605	19.75	ug/L	99
84) 1,2-dibromoethane	7.506	107	106315	20.93	ug/L	96
85) n-butyl ether	7.859	57	410303	17.06	ug/L	99
86) chlorobenzene	7.829	112	276155	18.40	ug/L	95
87) 1,1,1,2-tetrachloroethane	7.883	131	96259	18.14	ug/L	99
88) ethylbenzene	7.883	91	447245	17.93	ug/L	99
89) m,p-xylene	7.969	106	349303	36.07	ug/L	97
90) o-xylene	8.218	91	350219	18.20	ug/L	98
91) styrene	8.230	104	297520	18.42	ug/L	95
92) butyl acrylate	8.157	55	237949	18.86	ug/L	98
93) n-amyl acetate	8.291	70	84897	17.58	ug/L	94
94) isopropylbenzene	8.443	105	435640	19.29	ug/L	98
95) bromoform	8.358	173	76525	18.01	ug/L	97
96) cis-1,4-dichloro-2-butene	8.480	88	60960	19.26	ug/L	94
99) 1,1,2,2-tetrachloroethane	8.632	83	149563	19.04	ug/L	96
100) trans-1,4-dichloro-2-b...	8.650	53	49796	18.85	ug/L	94
101) 1,2,3-trichloropropane	8.674	110	48595	20.87	ug/L	94
102) bromobenzene	8.656	156	121700	19.70	ug/L	93
103) n-propylbenzene	8.699	91	511242	19.14	ug/L	99
104) 2-chlorotoluene	8.765	126	113856	20.61	ug/L	93
105) 4-chlorotoluene	8.838	91	309613	18.87	ug/L	95
106) 1,3,5-trimethylbenzene	8.802	105	351890	19.13	ug/L	99
107) tert-butylbenzene	8.997	119	311545	19.40	ug/L	98
108) 1,2,4-trimethylbenzene	9.033	105	365862	19.55	ug/L	97

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03013.D  
 Acq On : 4 Apr 2023 9:38 am  
 Operator : nickw  
 Sample : cc91-20  
 Misc : MS67909,V1R0100,5,,,1  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 04 09:50:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
109) sec-butylbenzene	9.131	105	433894	19.83	ug/L	97
110) p-isopropyltoluene	9.210	119	378667	18.90	ug/L	99
111) 1,2,3-trimethylbenzene	9.277	105	369350	18.52	ug/L	99
112) 1,3-dichlorobenzene	9.216	146	221255	18.62	ug/L	98
113) 1,4-dichlorobenzene	9.270	146	217117	17.89	ug/L	99
114) 1,2-dichlorobenzene	9.483	146	217376	19.93	ug/L	99
115) benzyl chloride	9.331	91	315335	21.53	ug/L	99
116) n-butylbenzene	9.447	92	165718	18.40	ug/L	98
117) hexachloroethane	9.648	201	38944	20.42	ug/L	94
118) 1,2-dibromo-3-chloropr...	9.927	157	47591	22.36	ug/L	93
119) 1,3,5-trichlorobenzene	10.025	180	135311	19.67	ug/L	96
120) 1,2,4-trichlorobenzene	10.378	180	114984	19.86	ug/L	90
121) hexachlorobutadiene	10.445	225	39763	18.91	ug/L	98
122) naphthalene	10.530	128	432371	20.46	ug/L	99
123) 1,2,3-trichlorobenzene	10.645	180	113211	19.77	ug/L	92
124) 2-methylnaphthalene	11.132	142	96658	10.46	ug/L	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

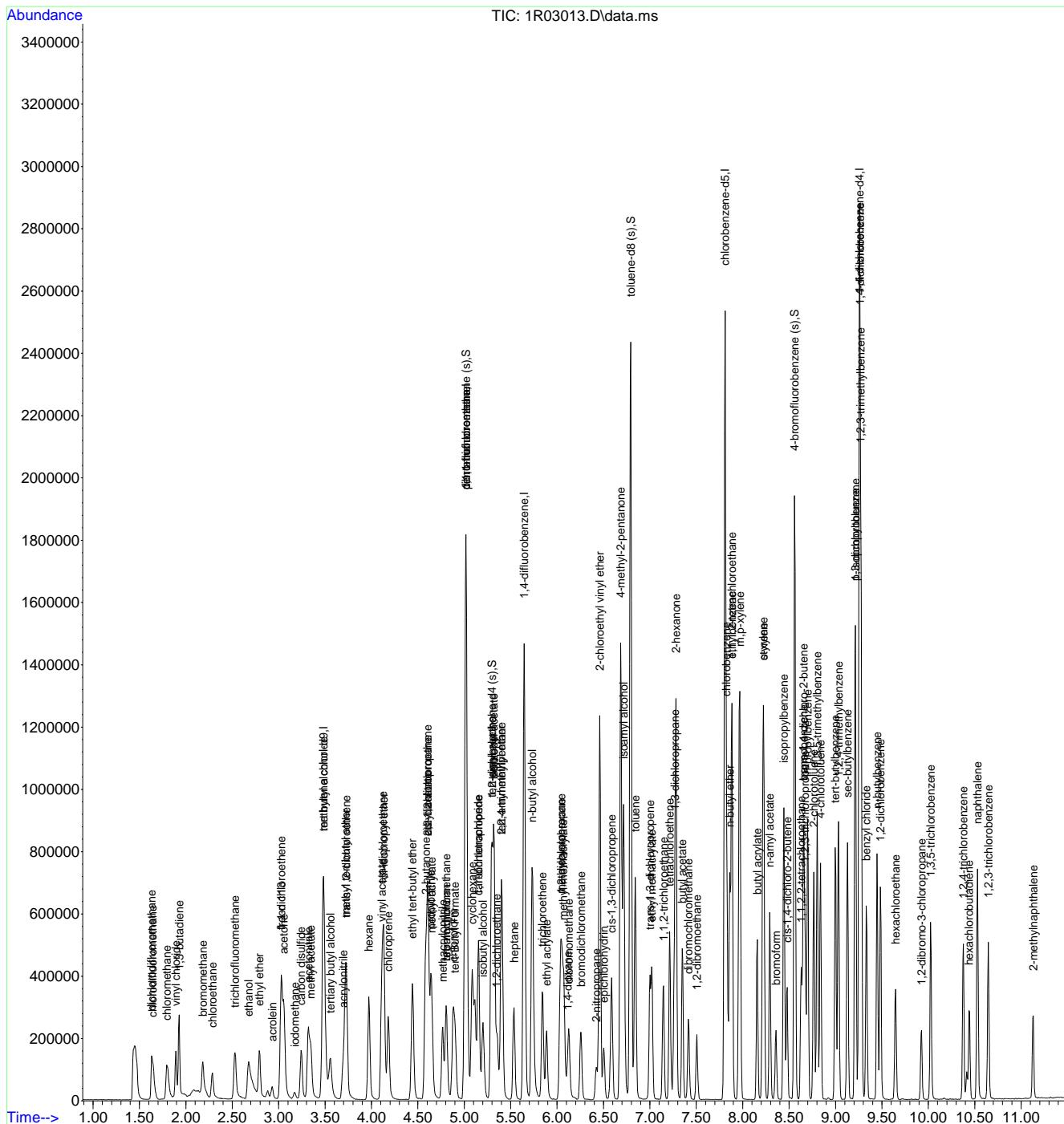
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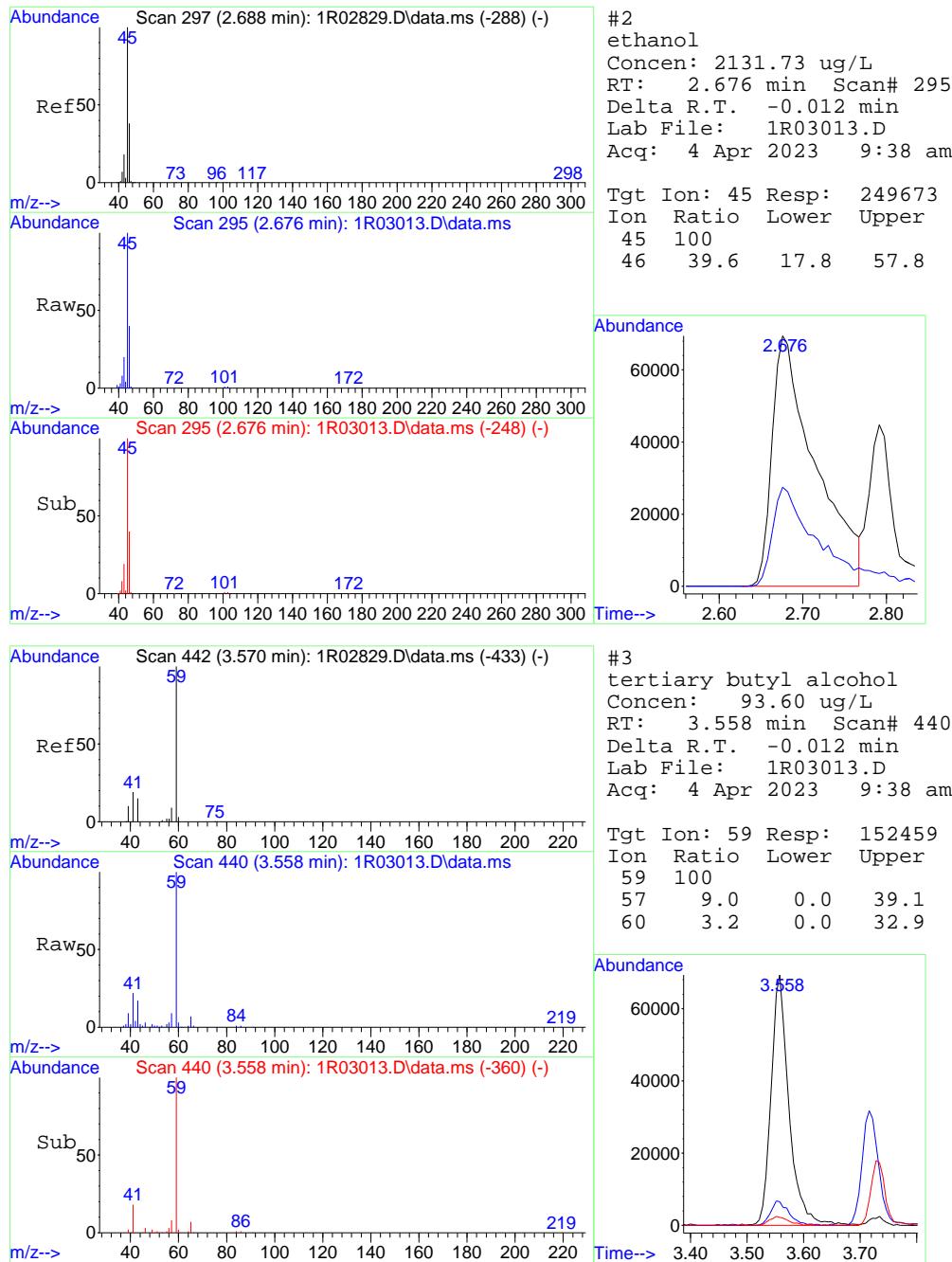
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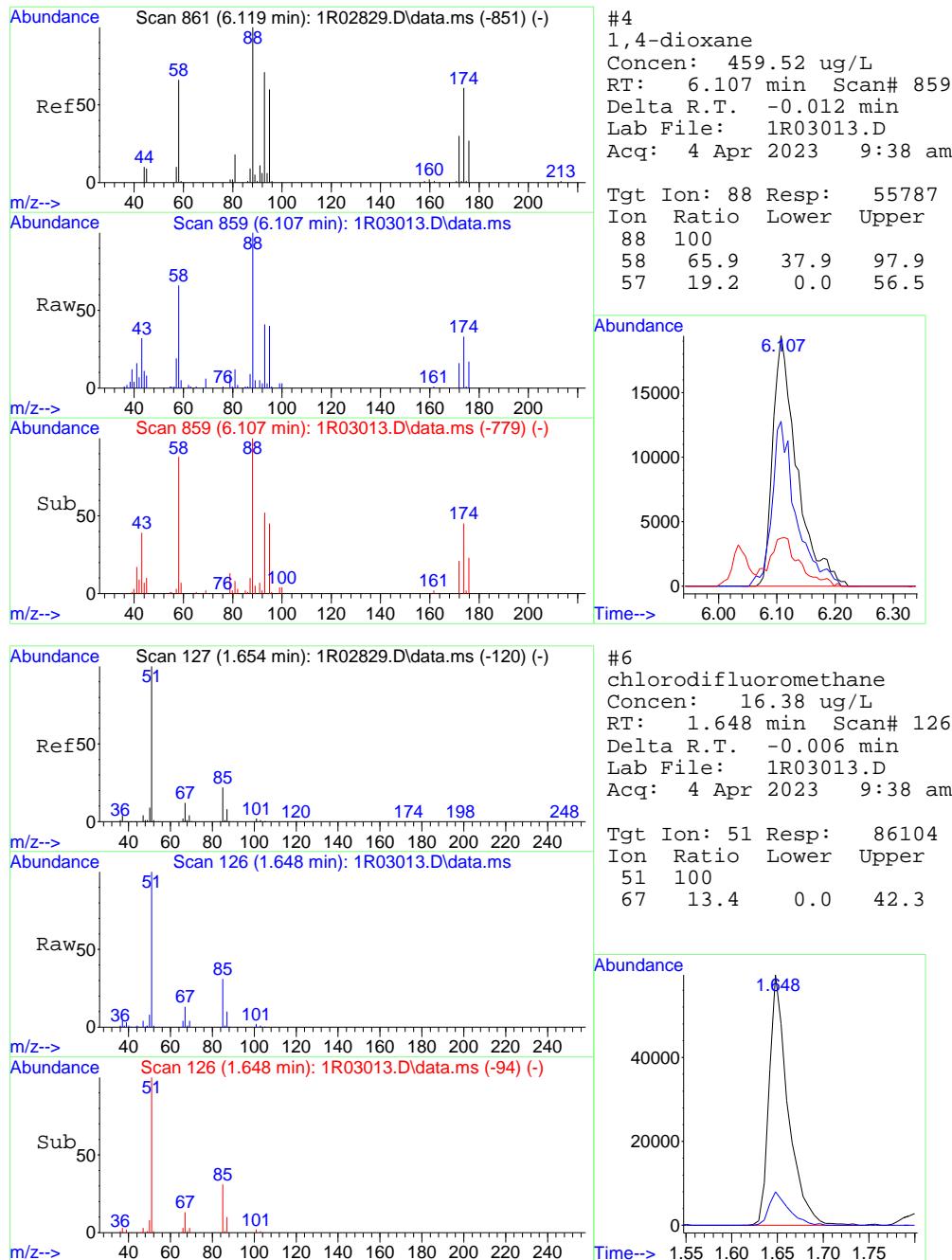
## Quantitation Report (QT Reviewed)

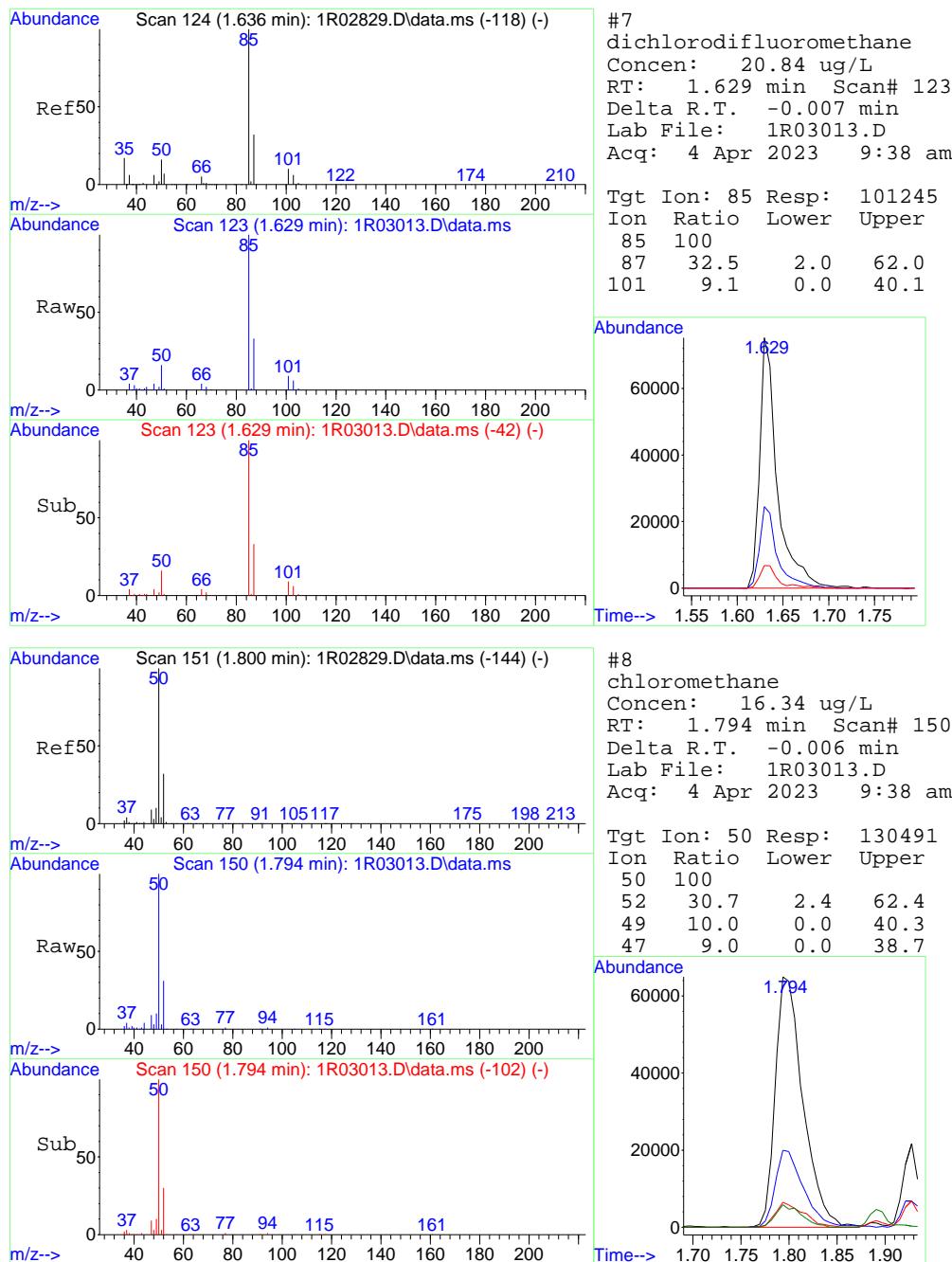
Data Path : C:\msdchem\1\data\Mei 04-06-2023\VR0100\  
 Data File : 1R03013.D  
 Acq On : 4 Apr 2023 9:38 am  
 Operator : nickw  
 Sample : cc91-20  
 Misc : MS67909,V1R0100,5,,,1  
 ALS Vial : 2 Sample Multiplier: 1

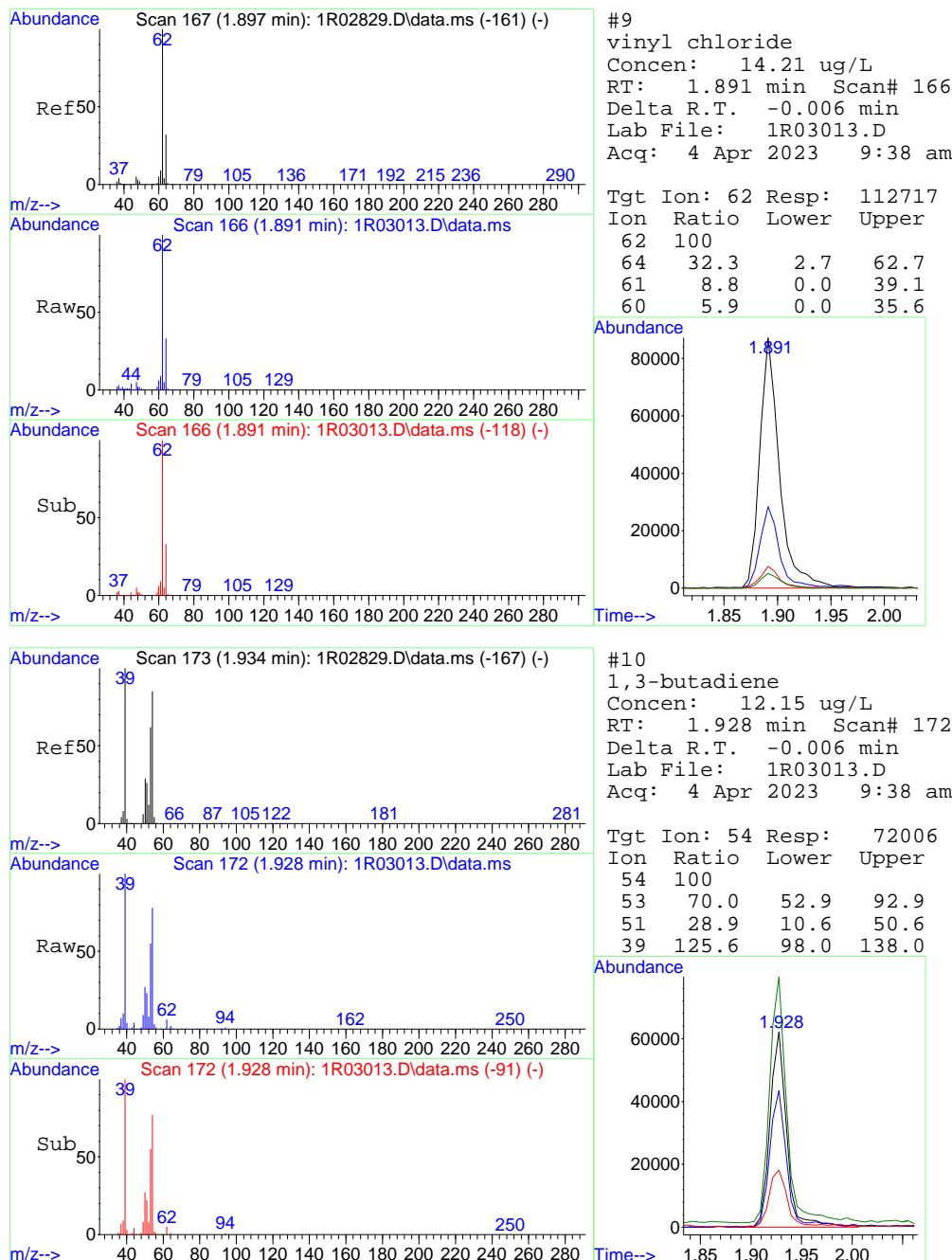
Quant Time: Apr 04 09:50:11 2023  
 Quant Method : C:\MSDCHEM\1\METHODS\M1R0091.M  
 Quant Title : SW846 8260D, RxI624Sil MS 60m x 0.25mm x 1.4um  
 QLast Update : Wed Mar 29 12:39:36 2023  
 Response via : Initial Calibration

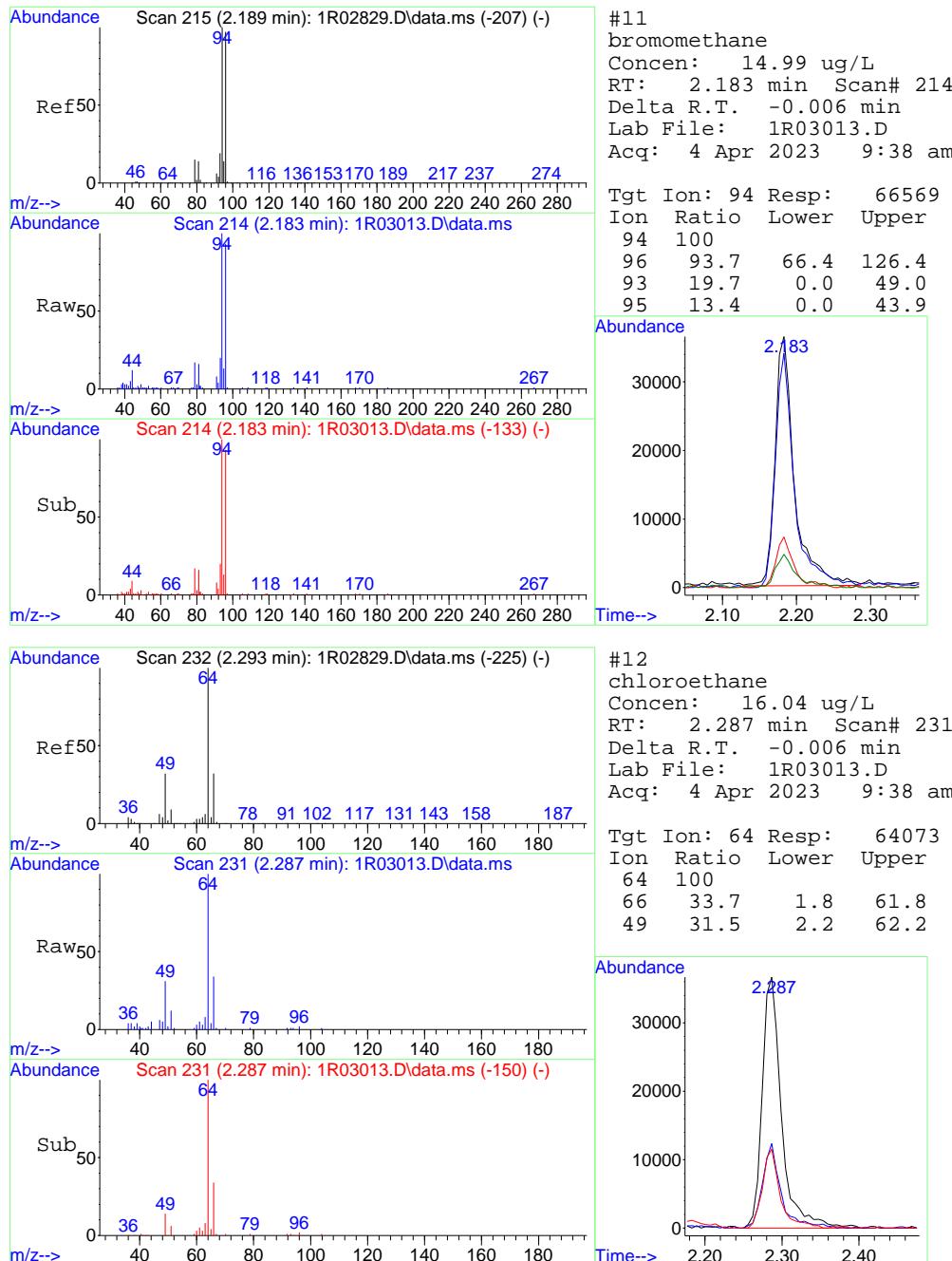


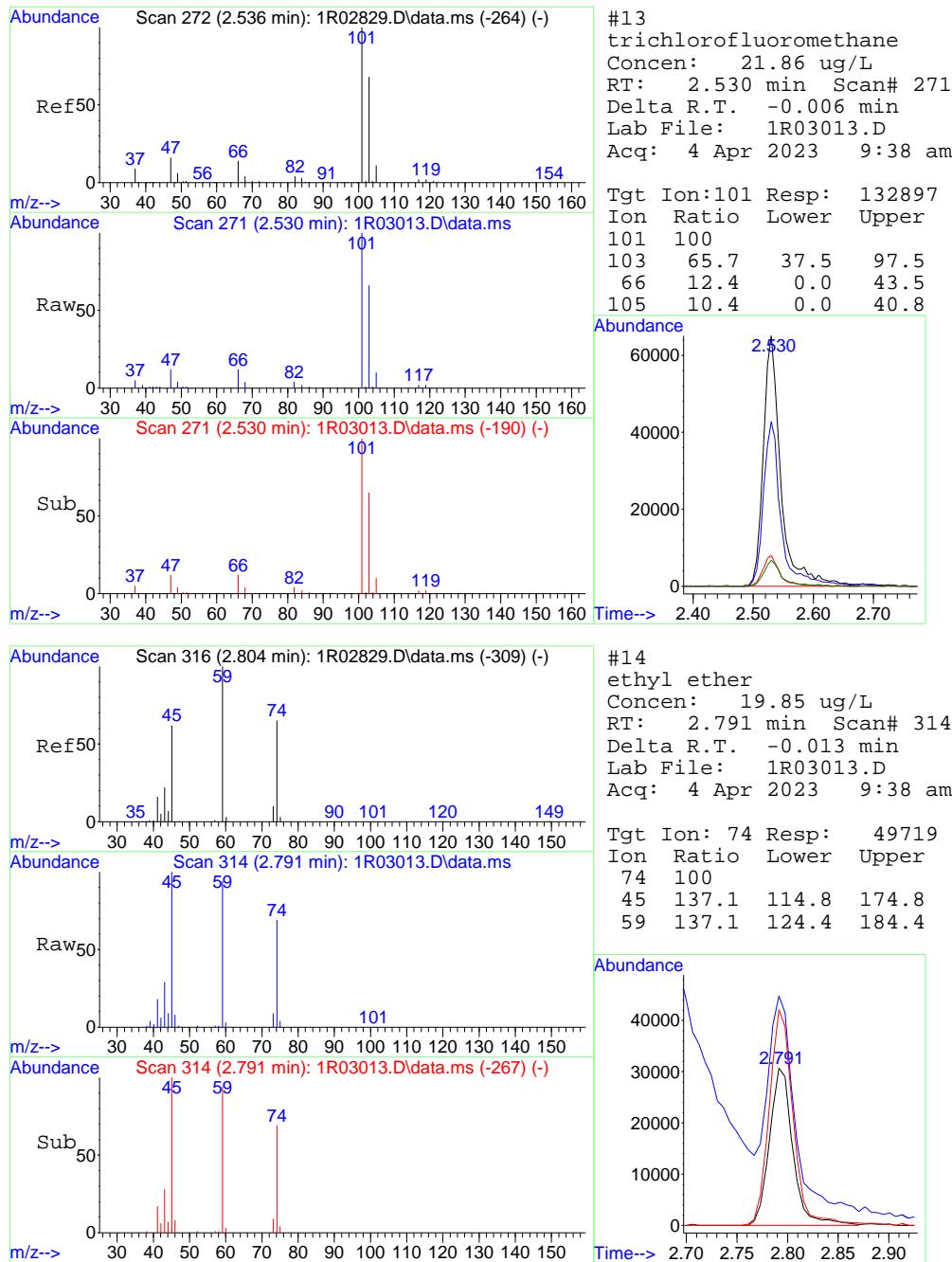


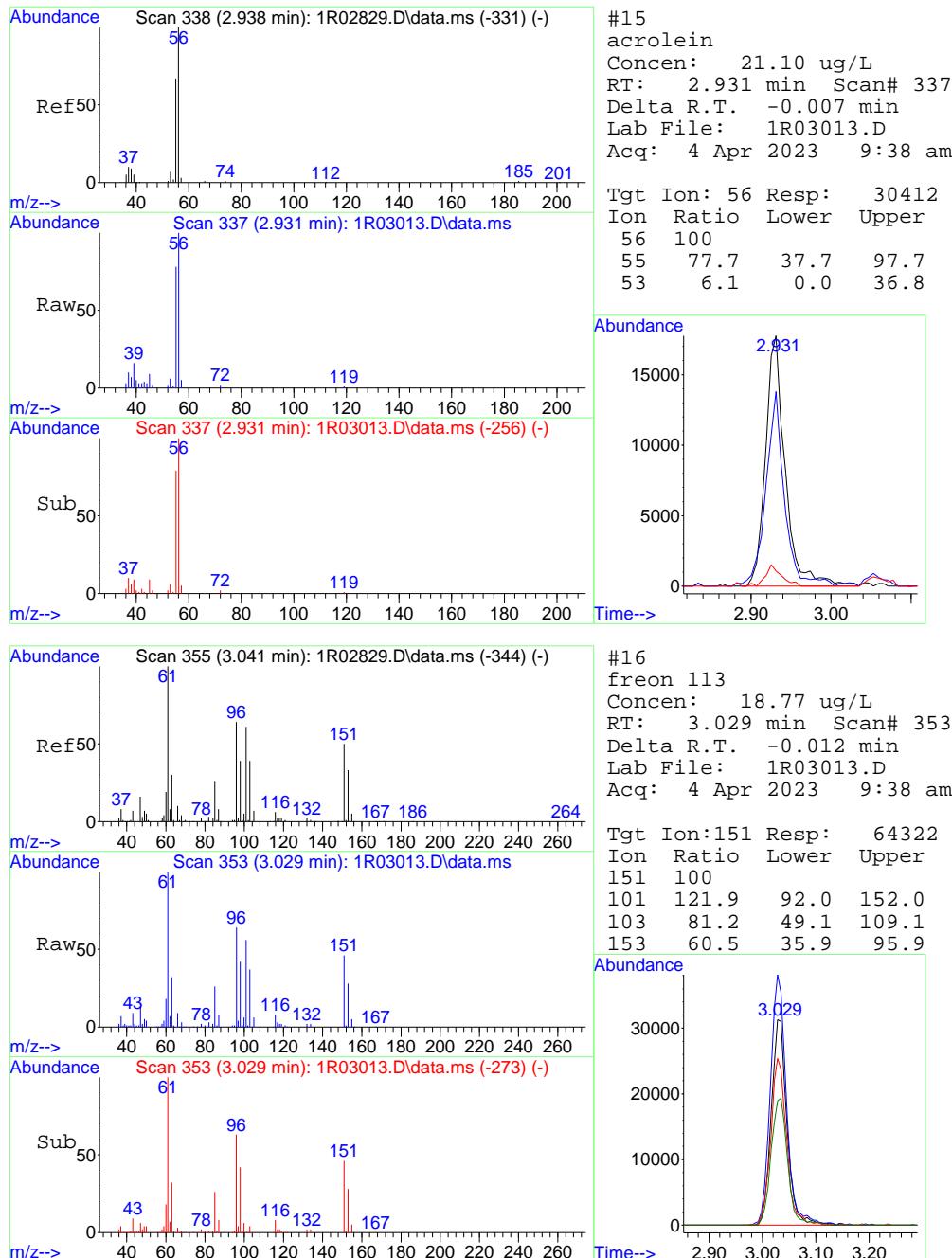


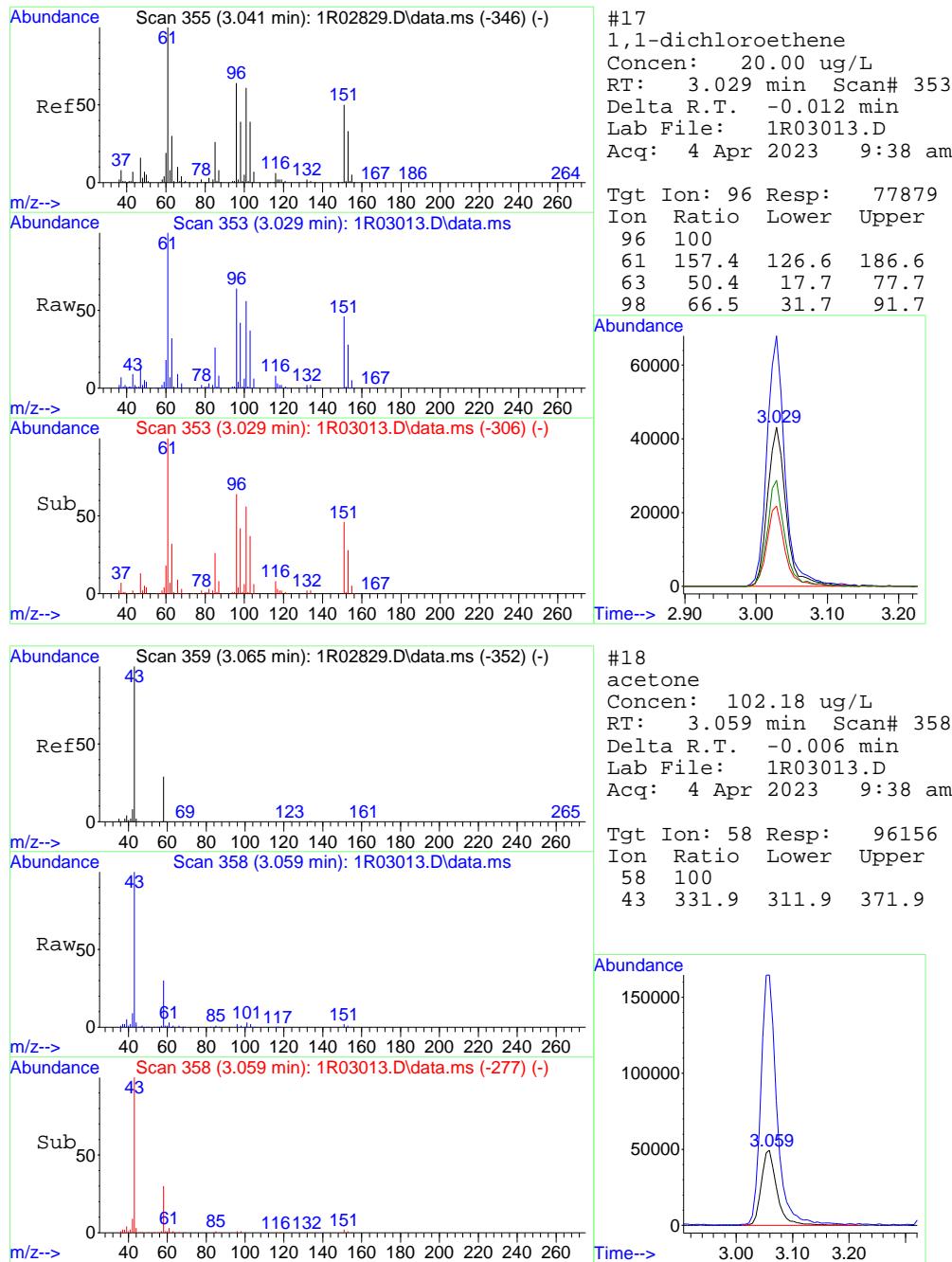


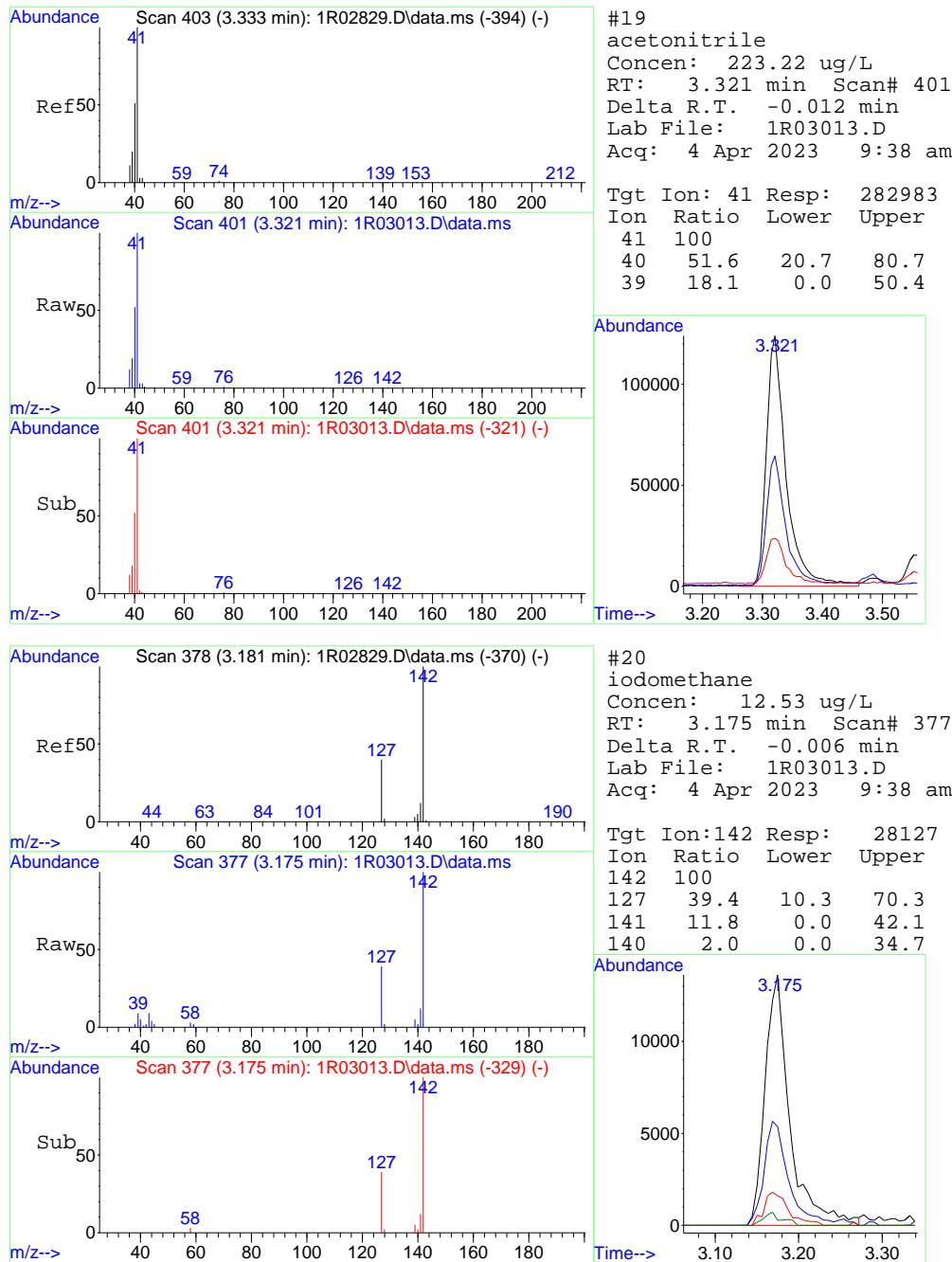


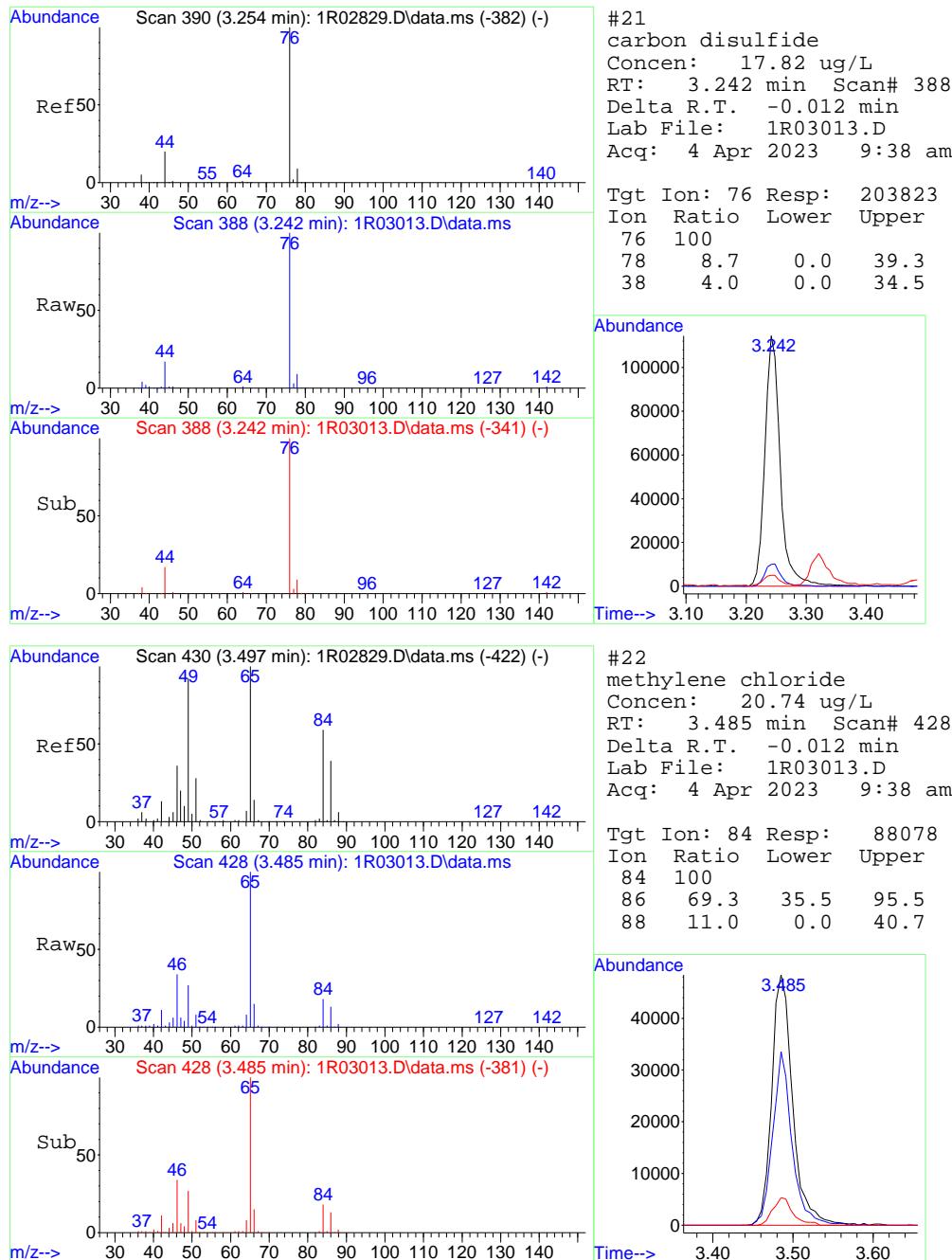


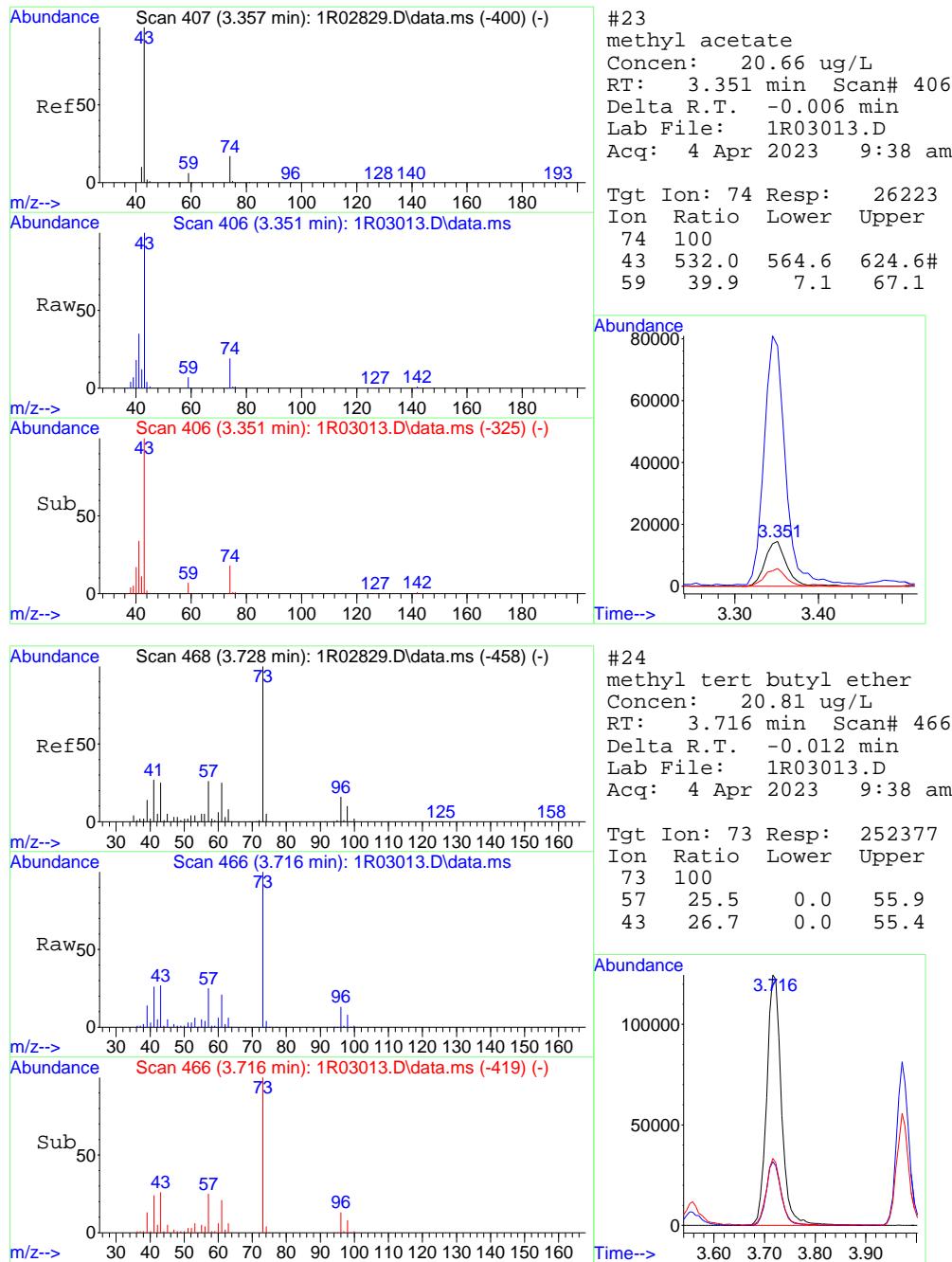


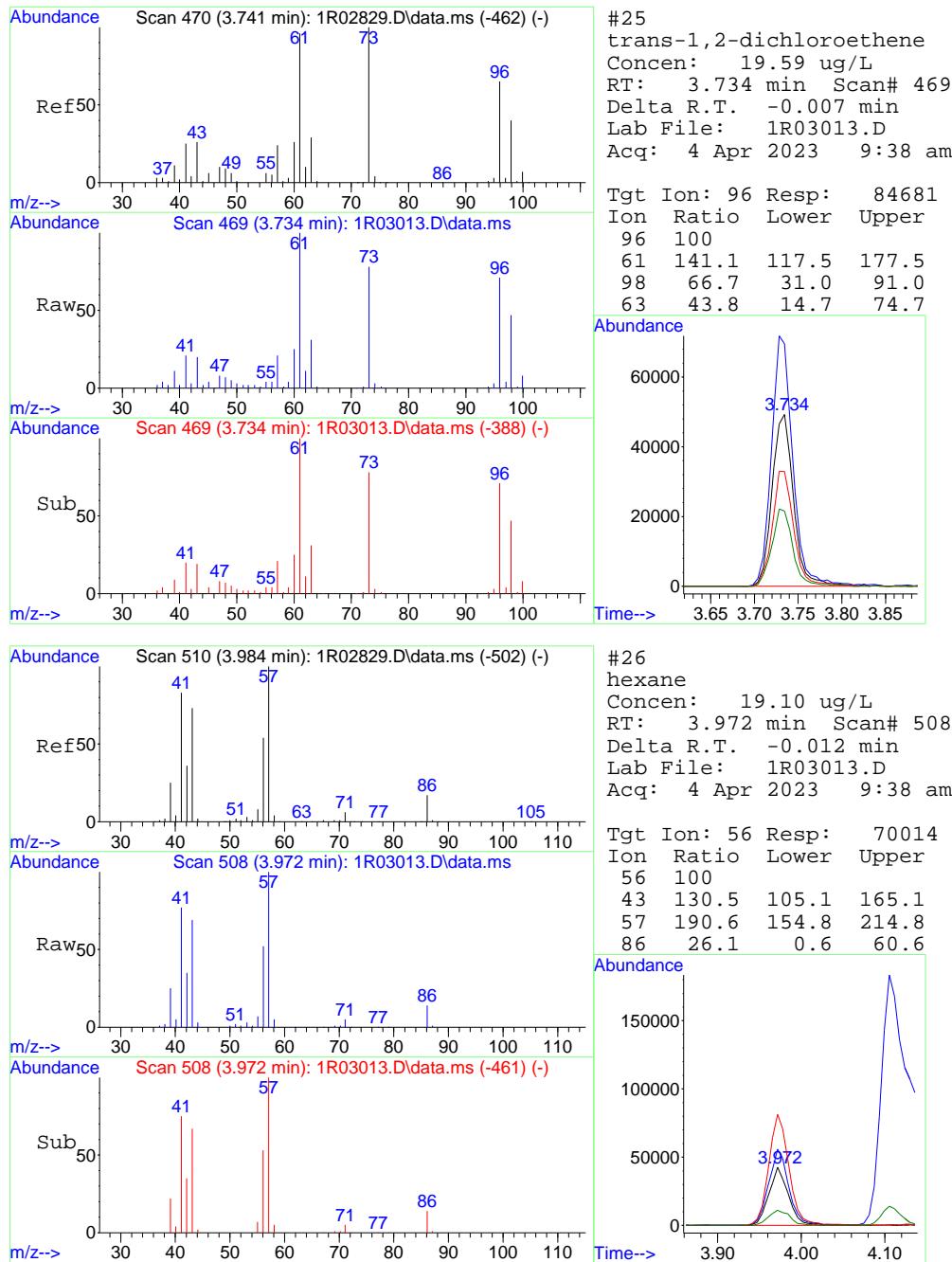


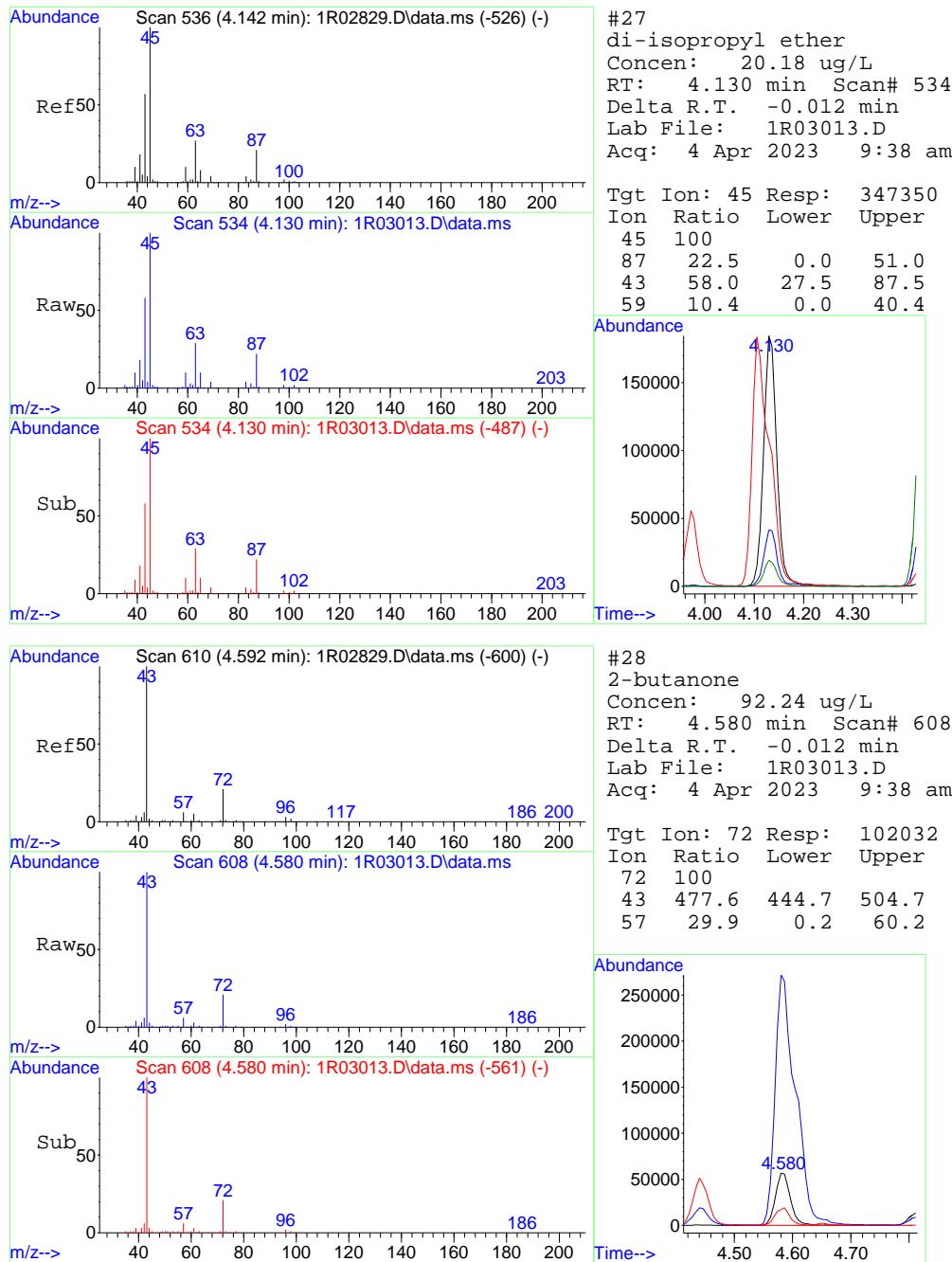


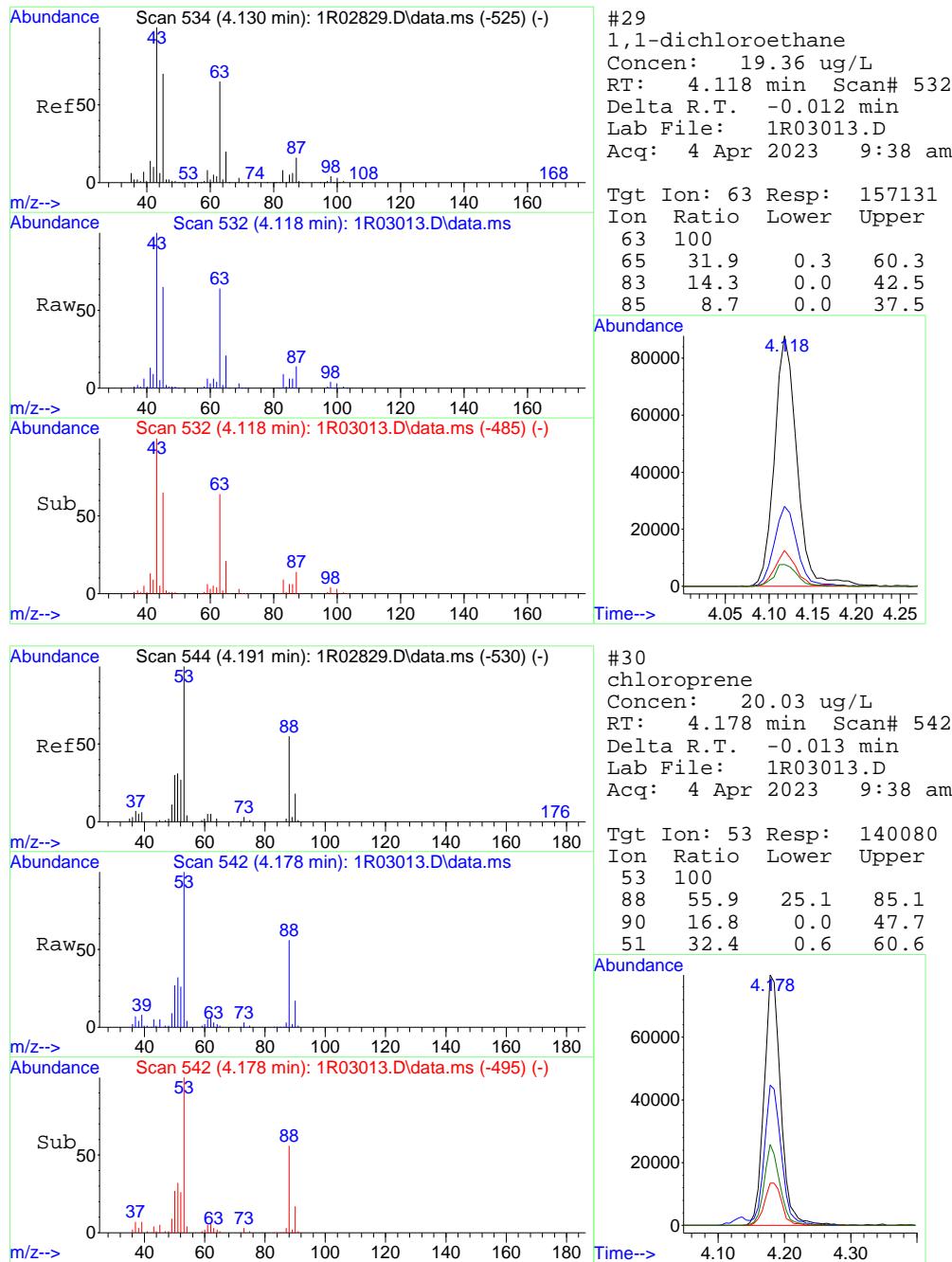


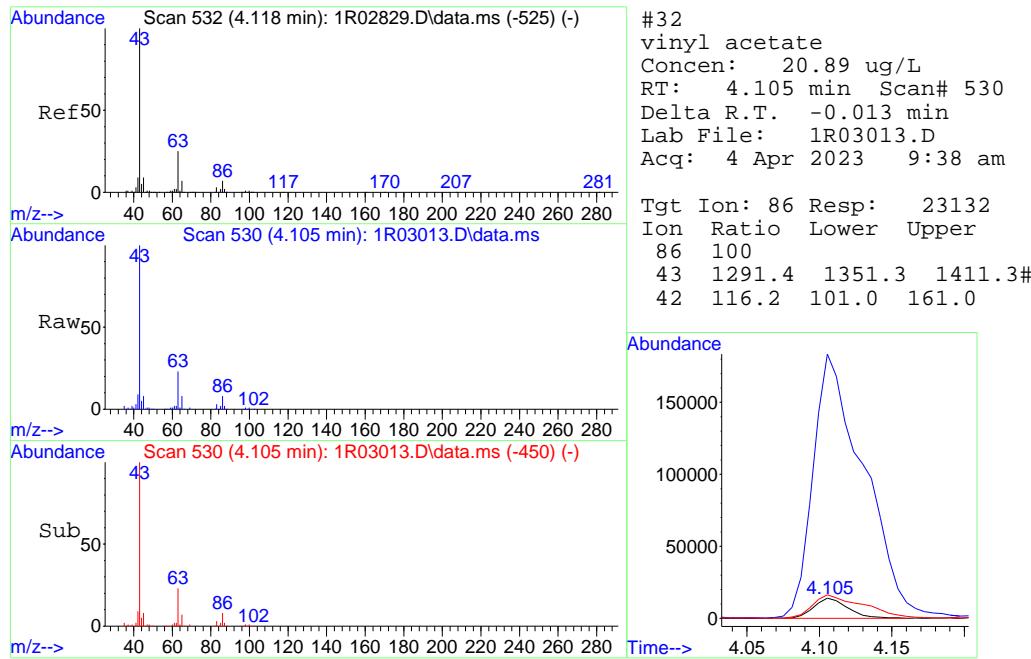
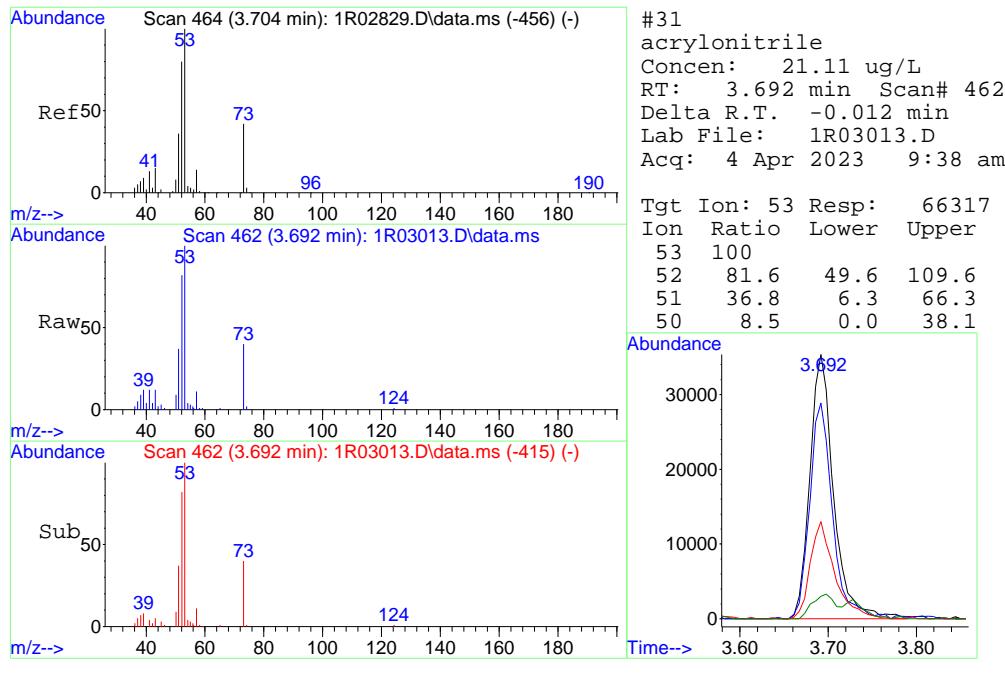


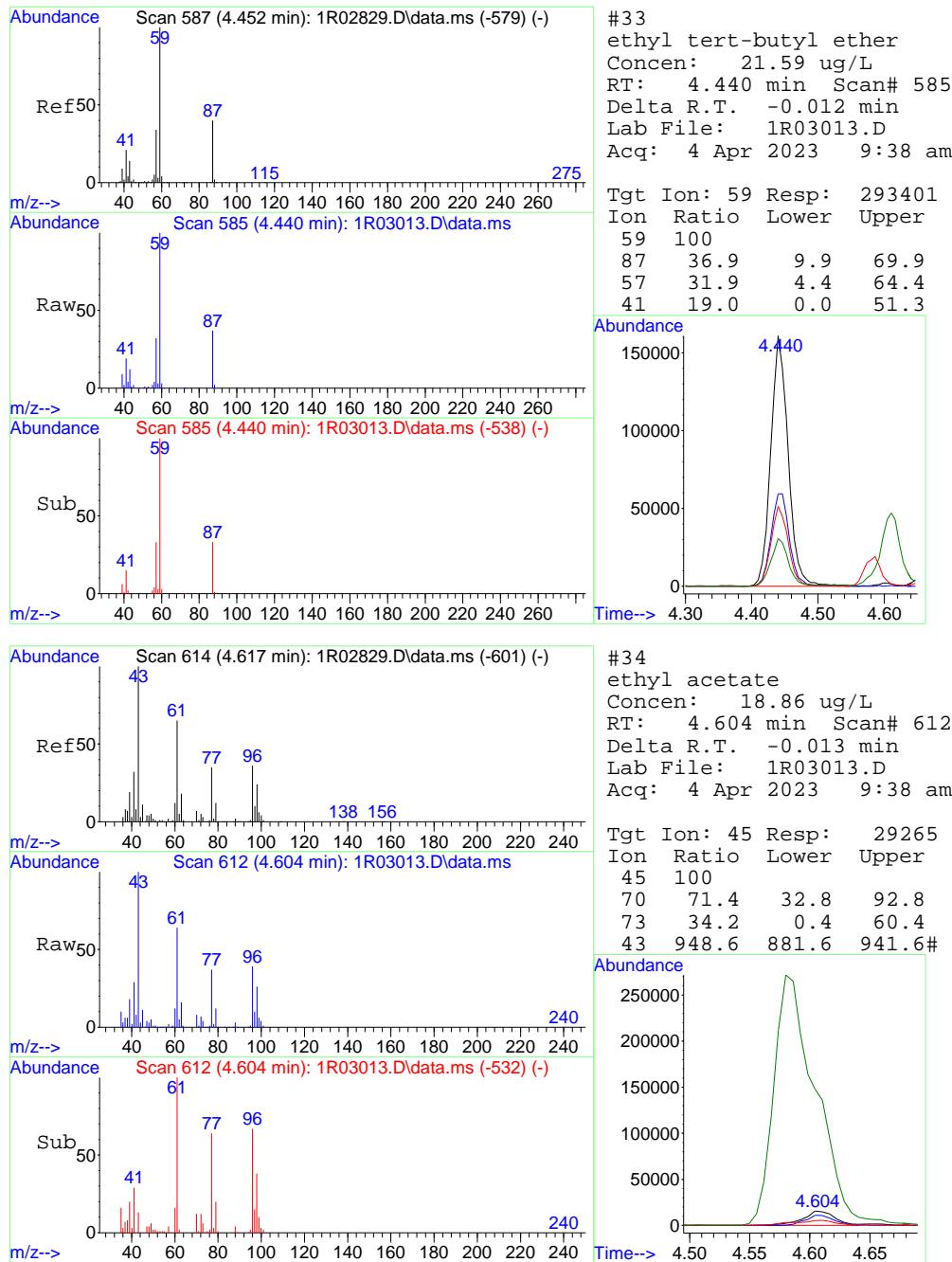


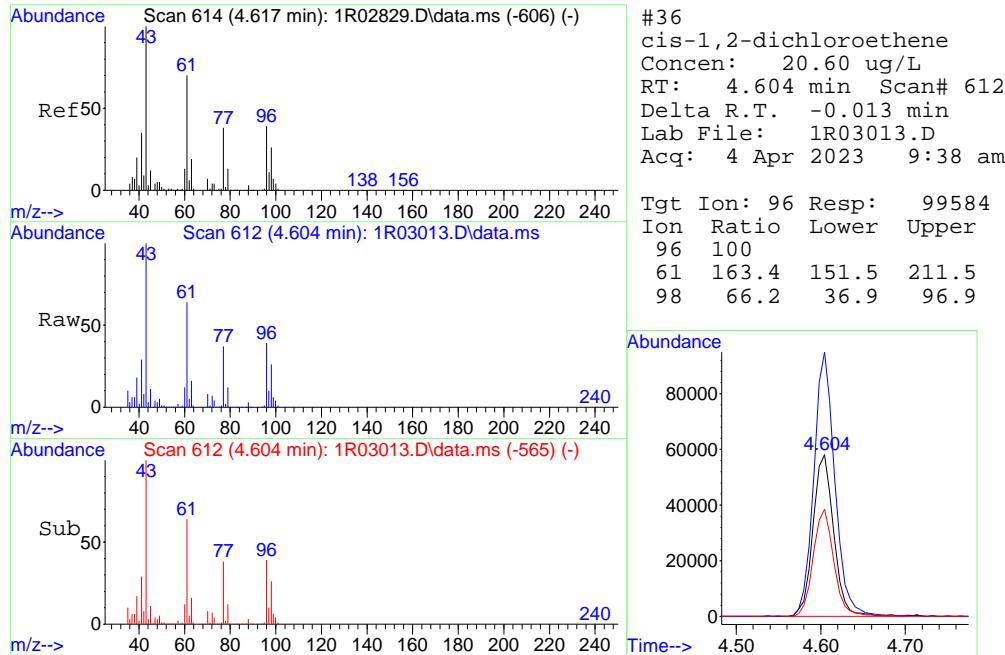
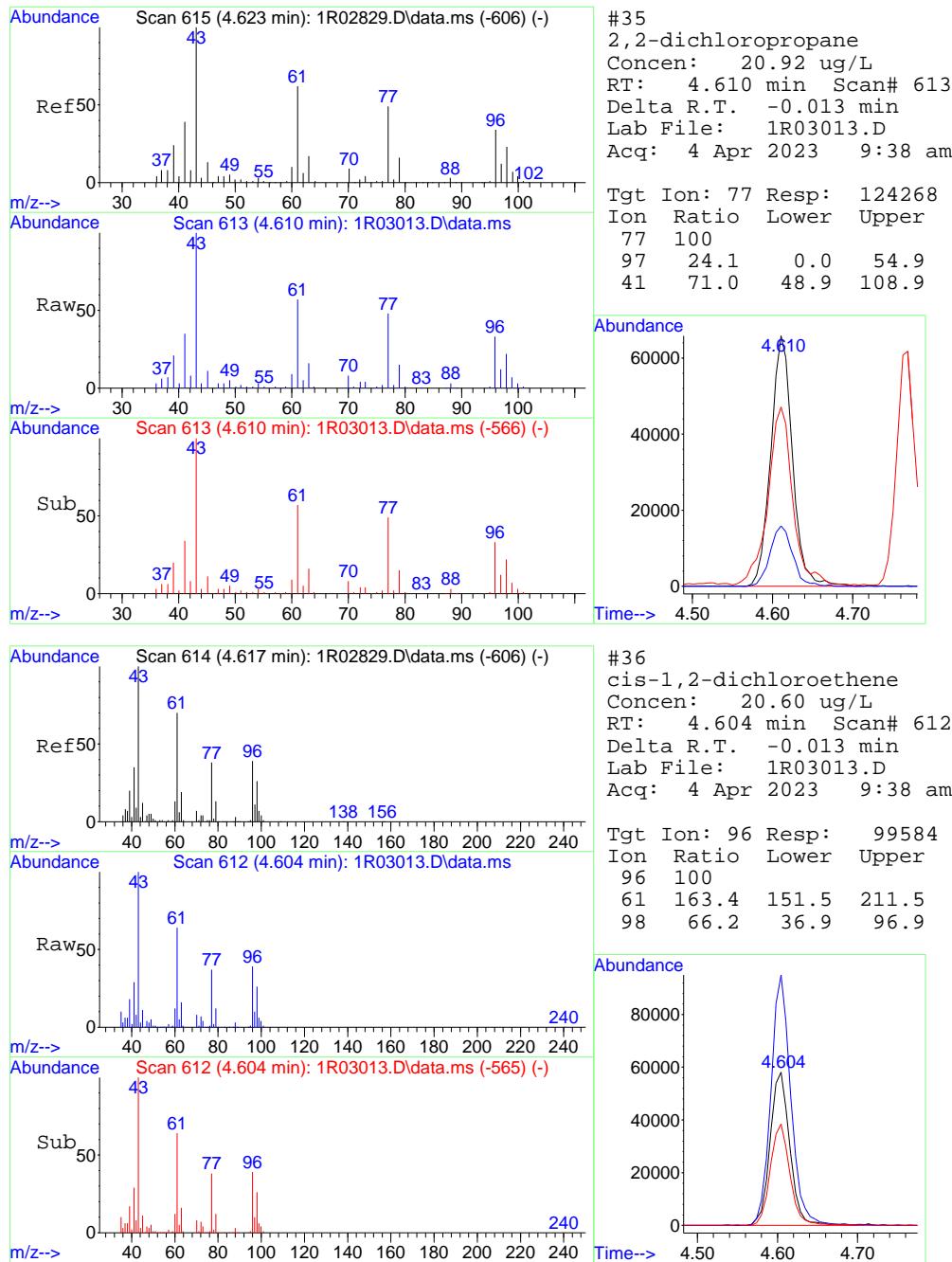


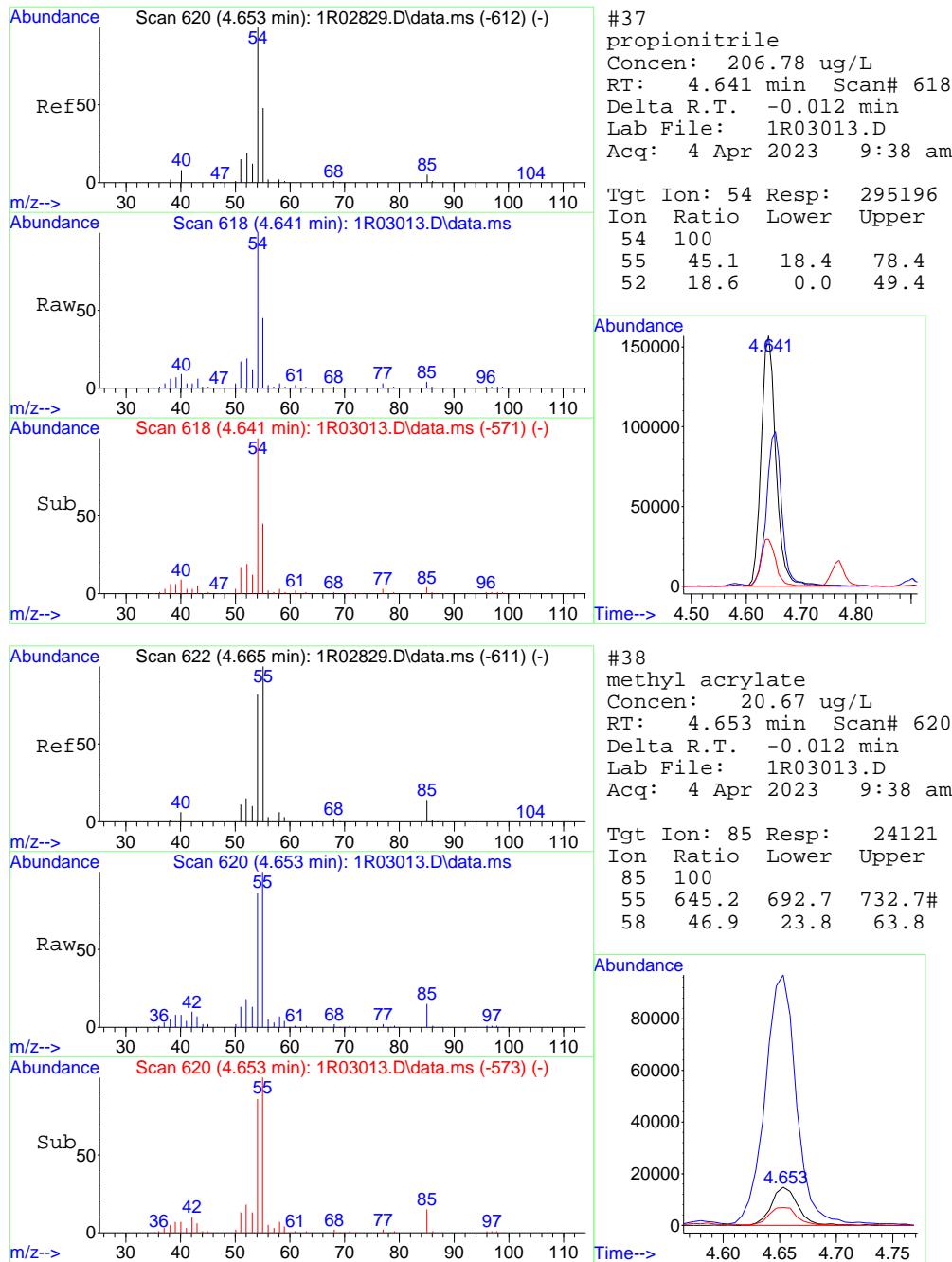


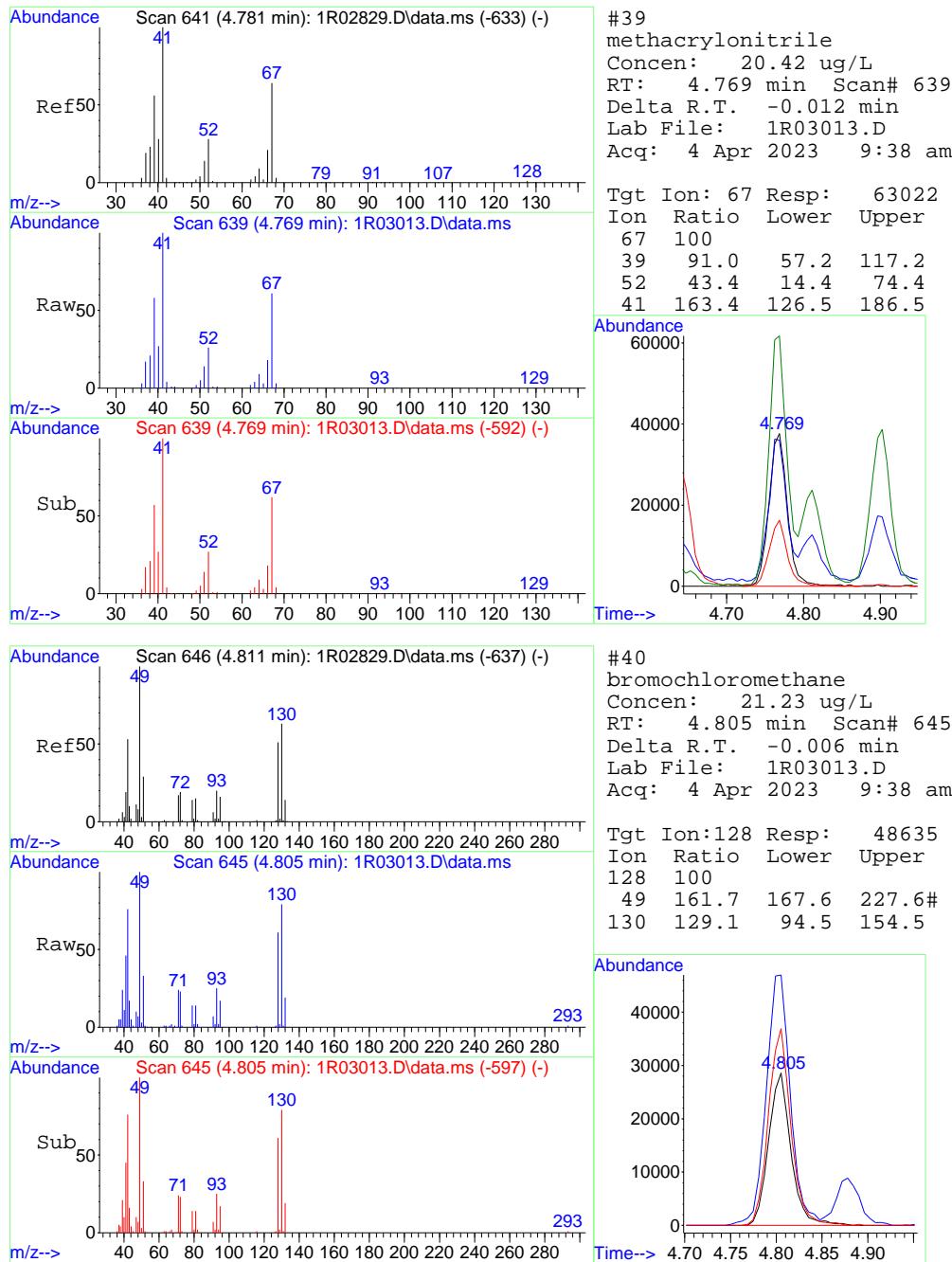


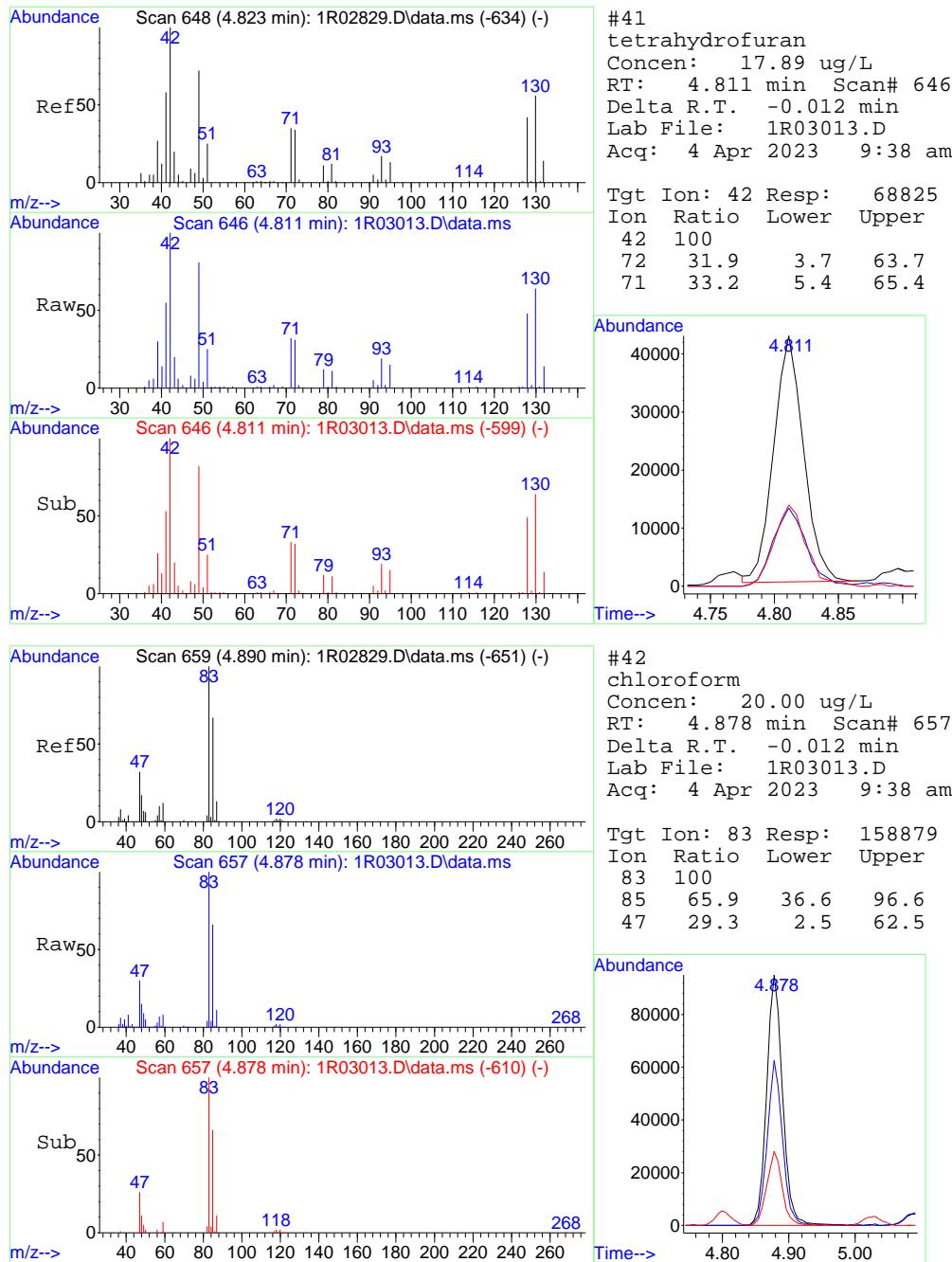


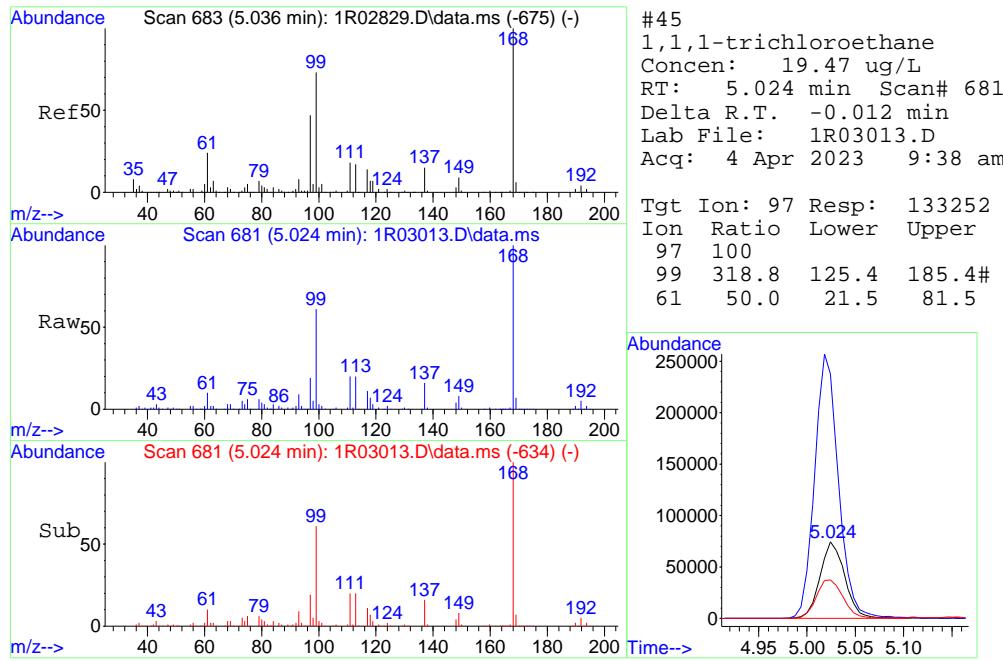
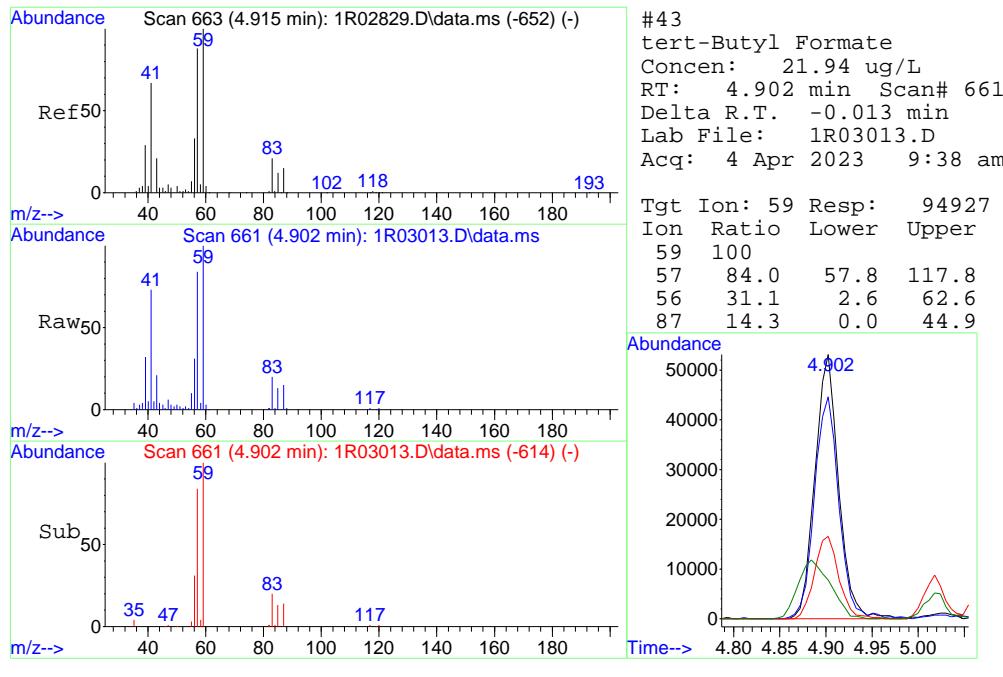


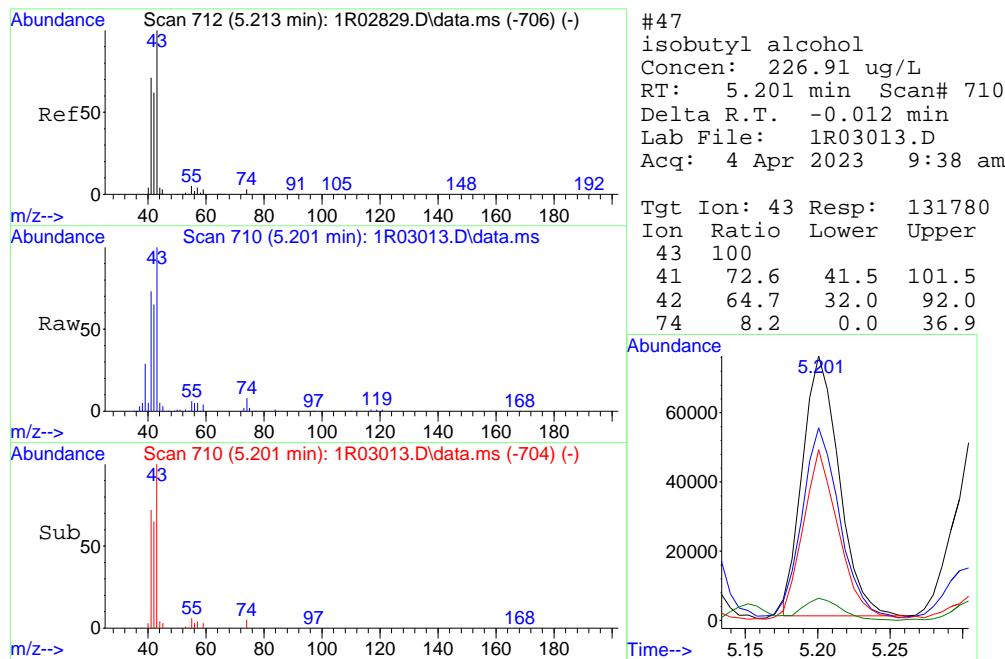
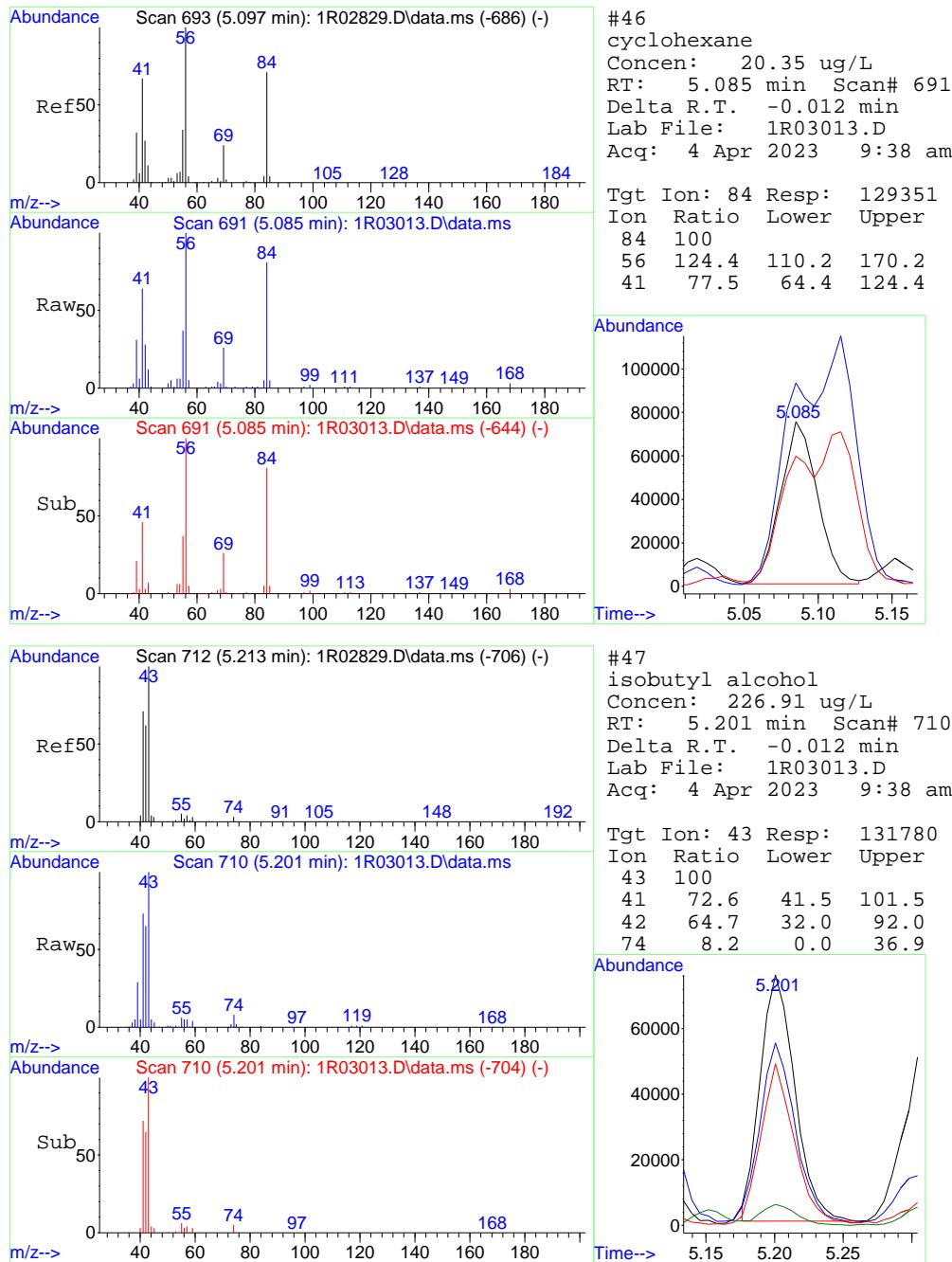


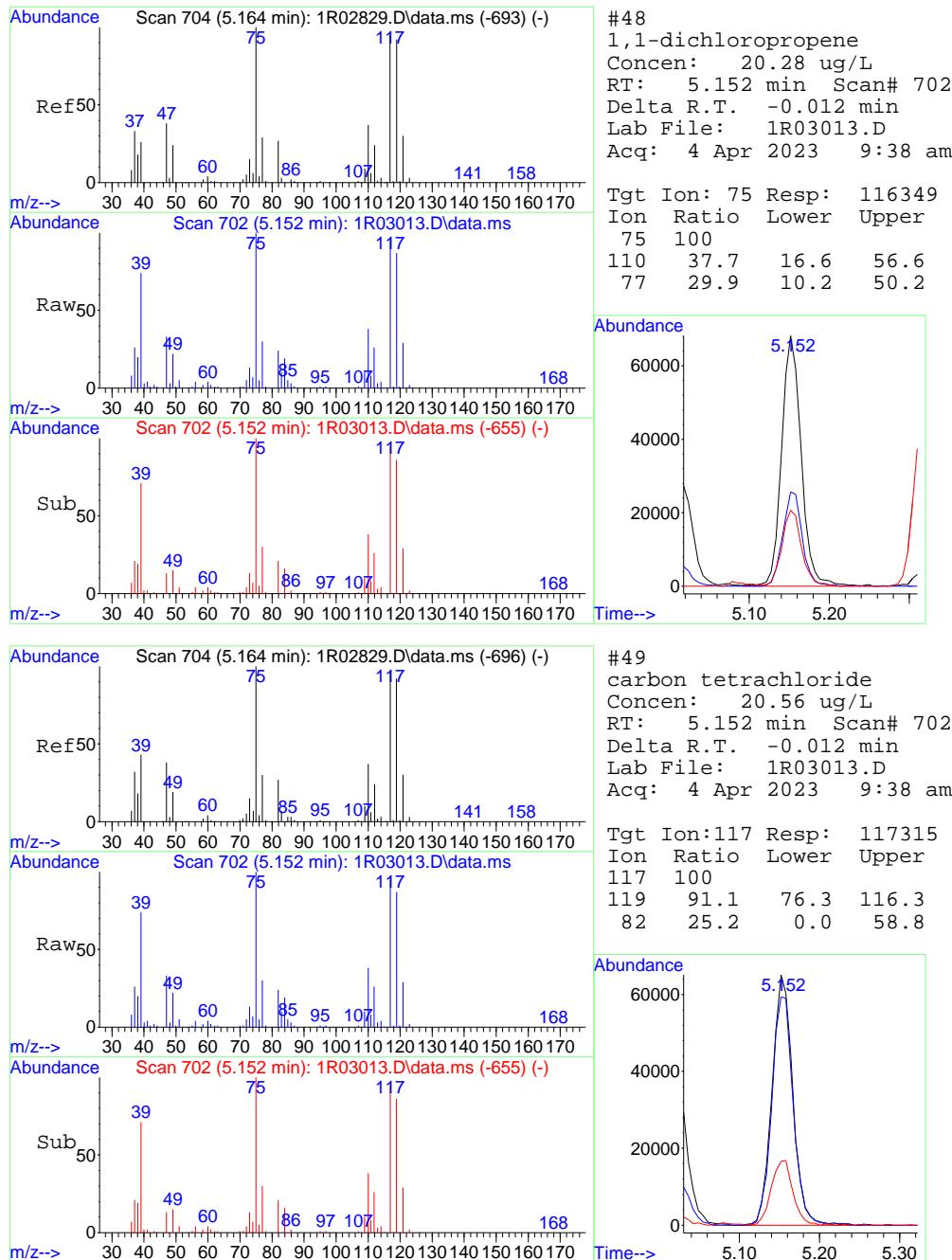


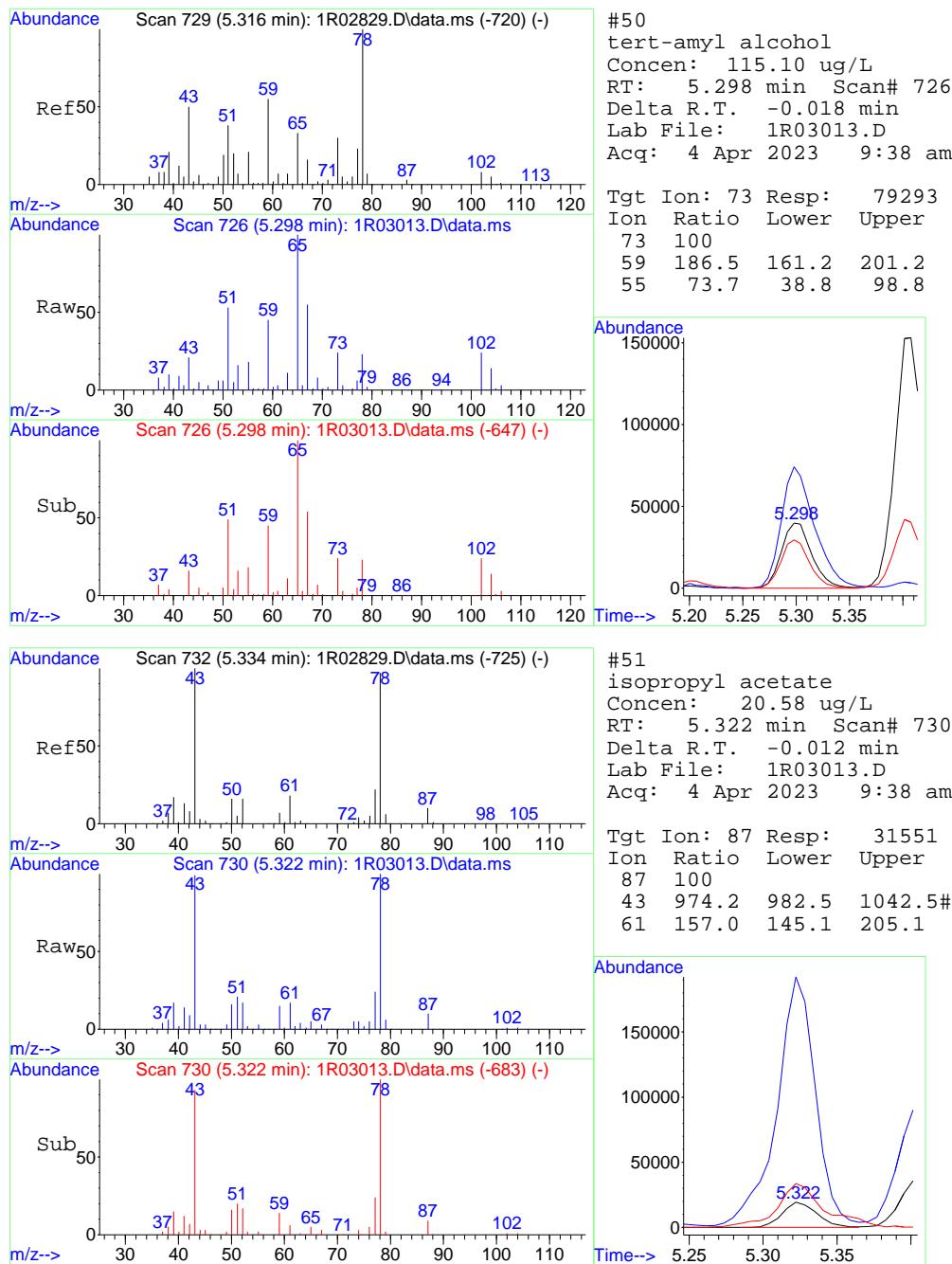


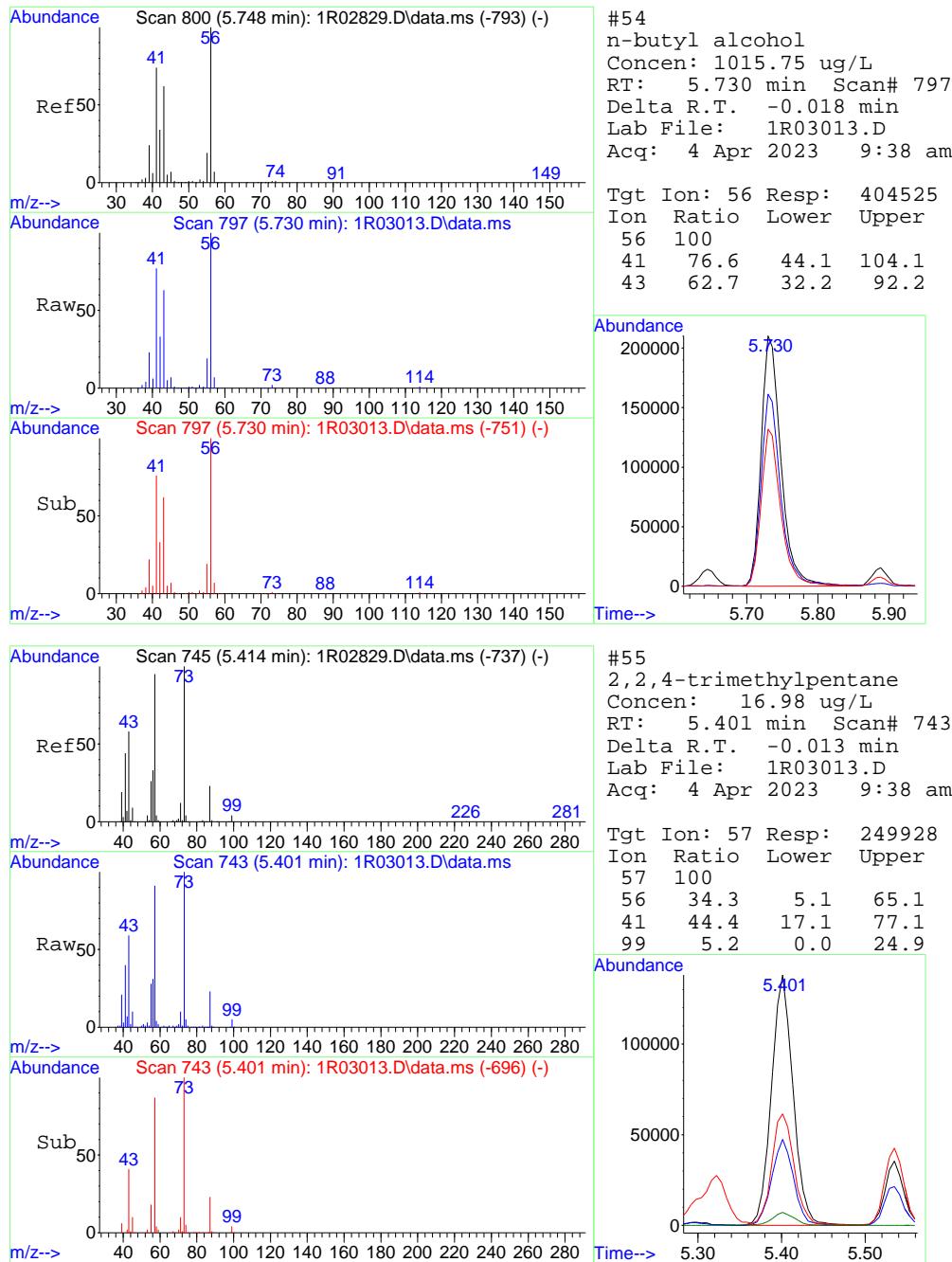


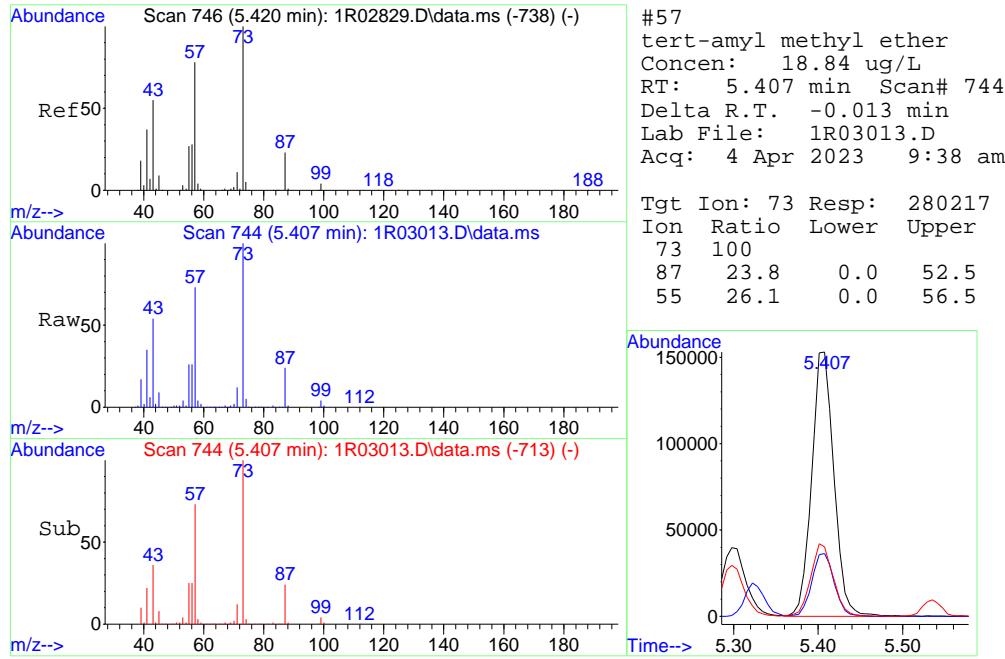
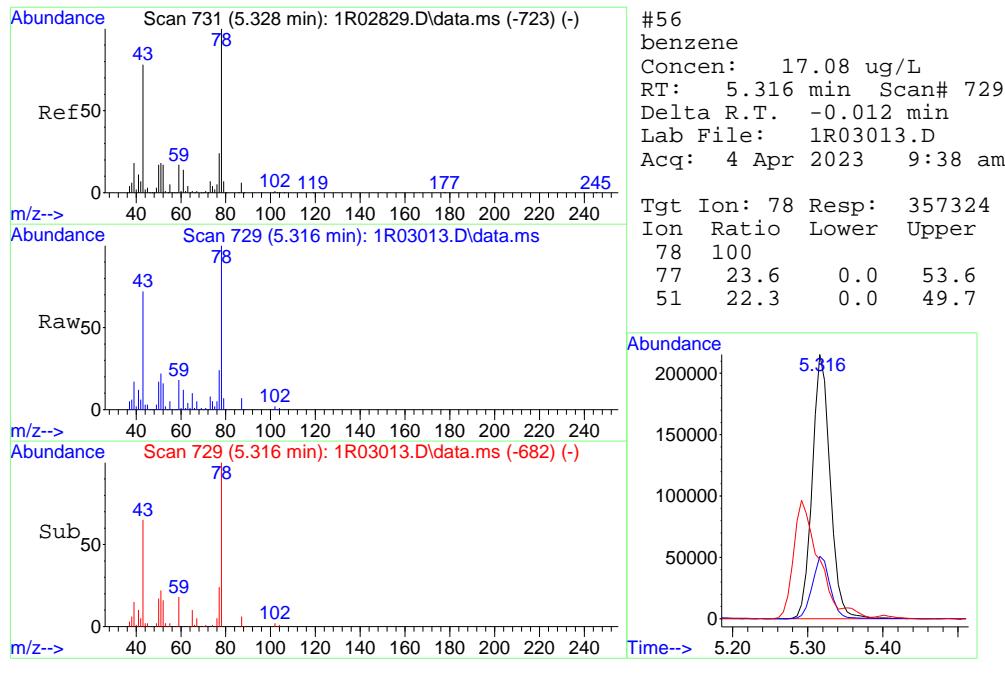




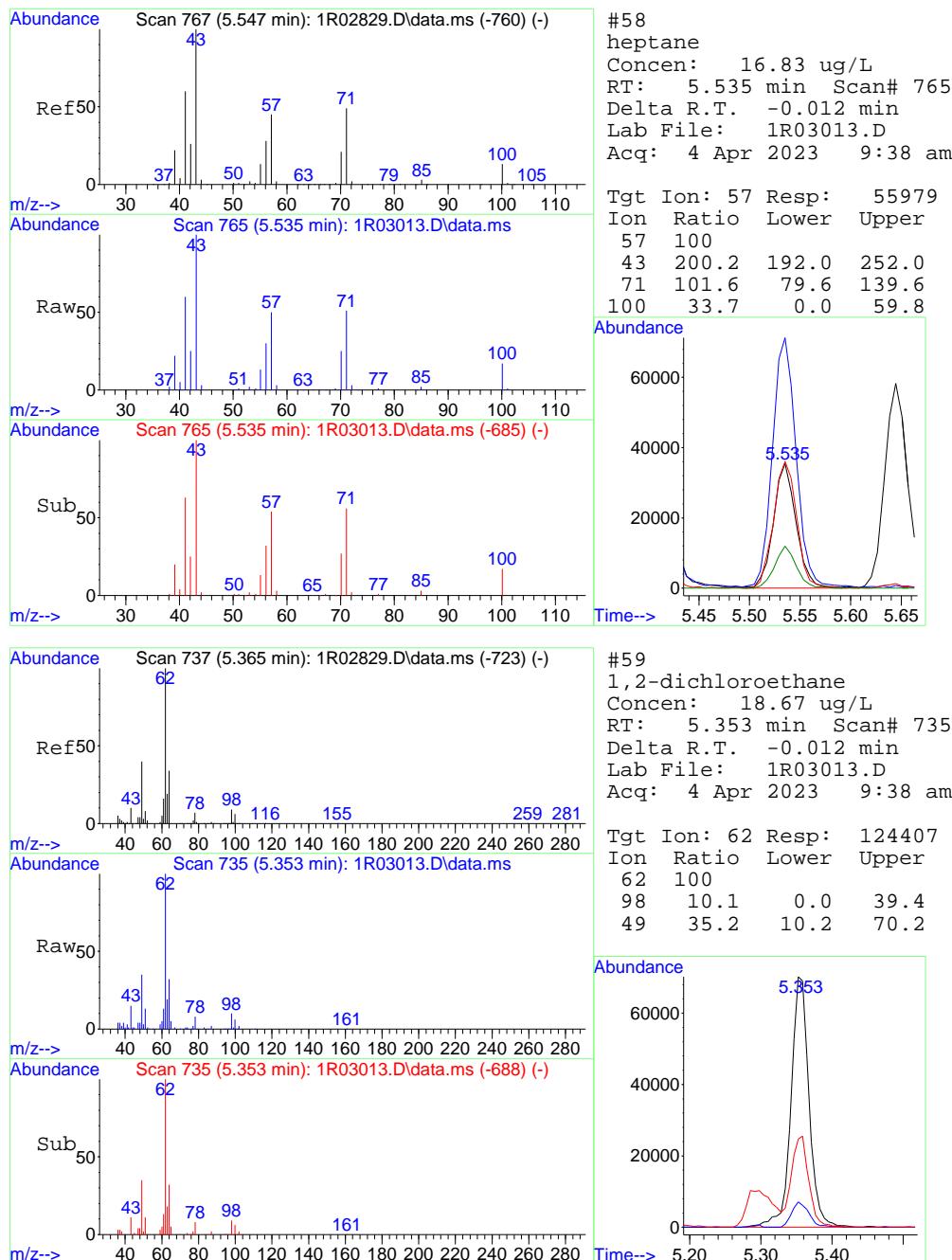


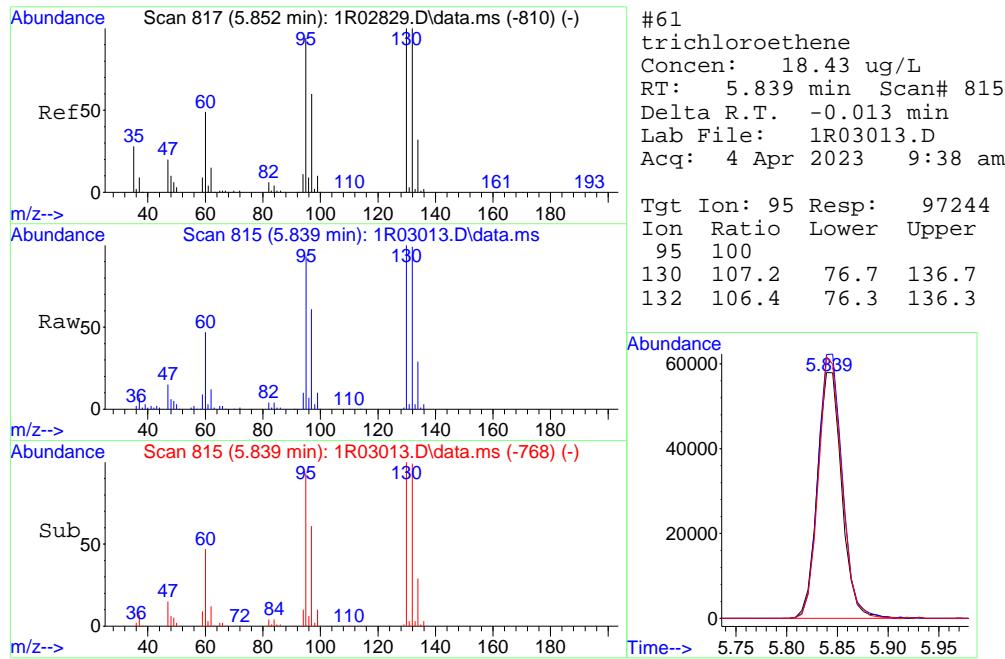
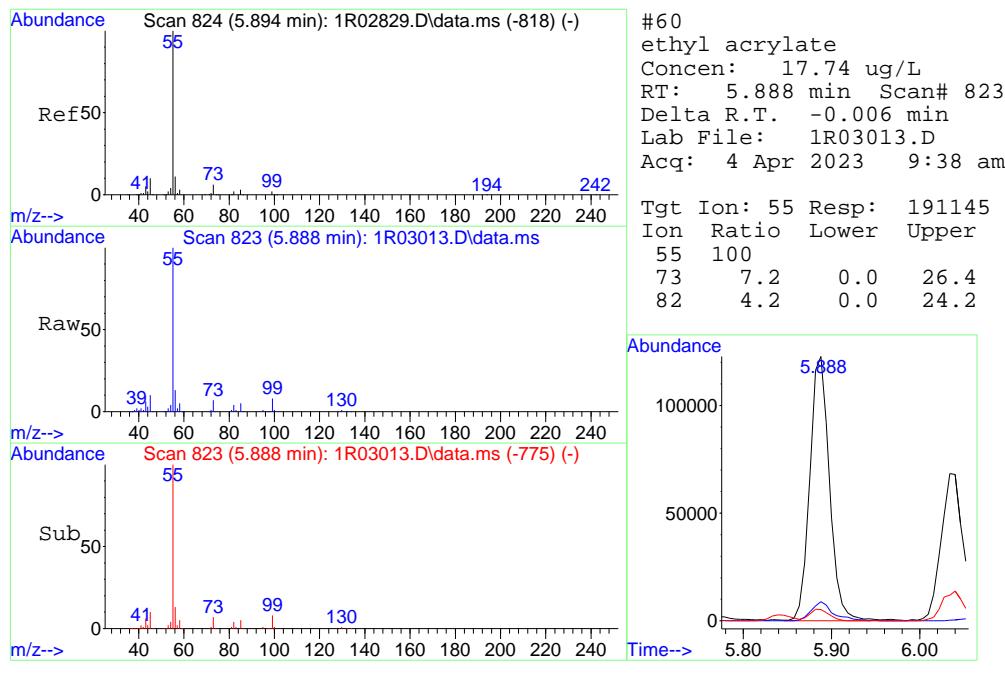


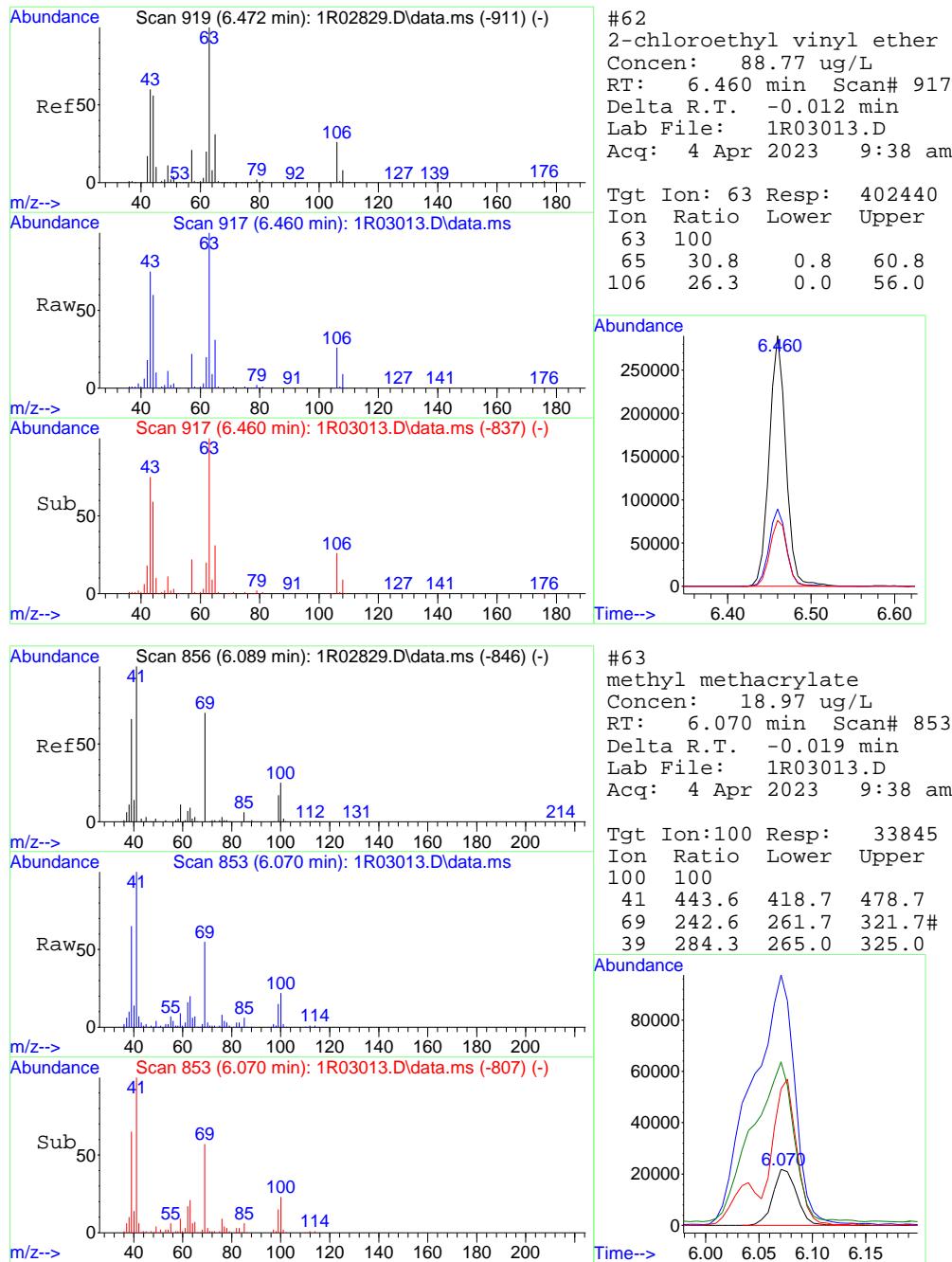


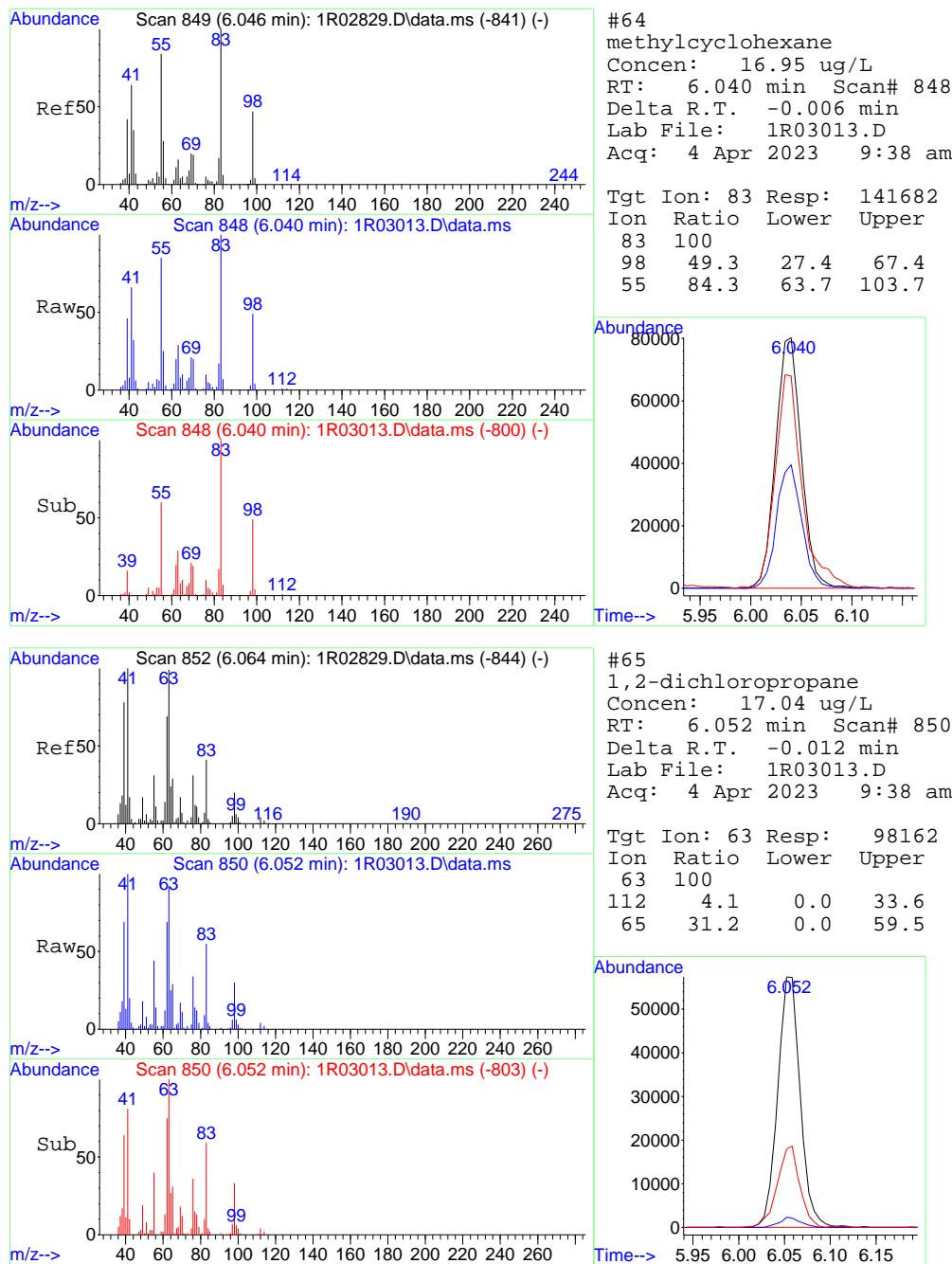


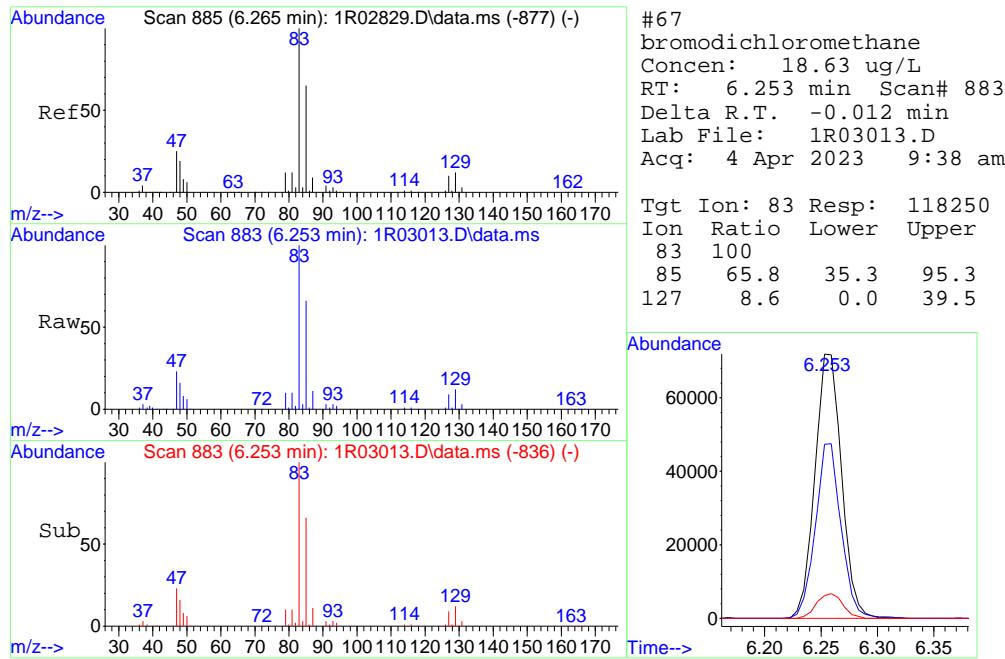
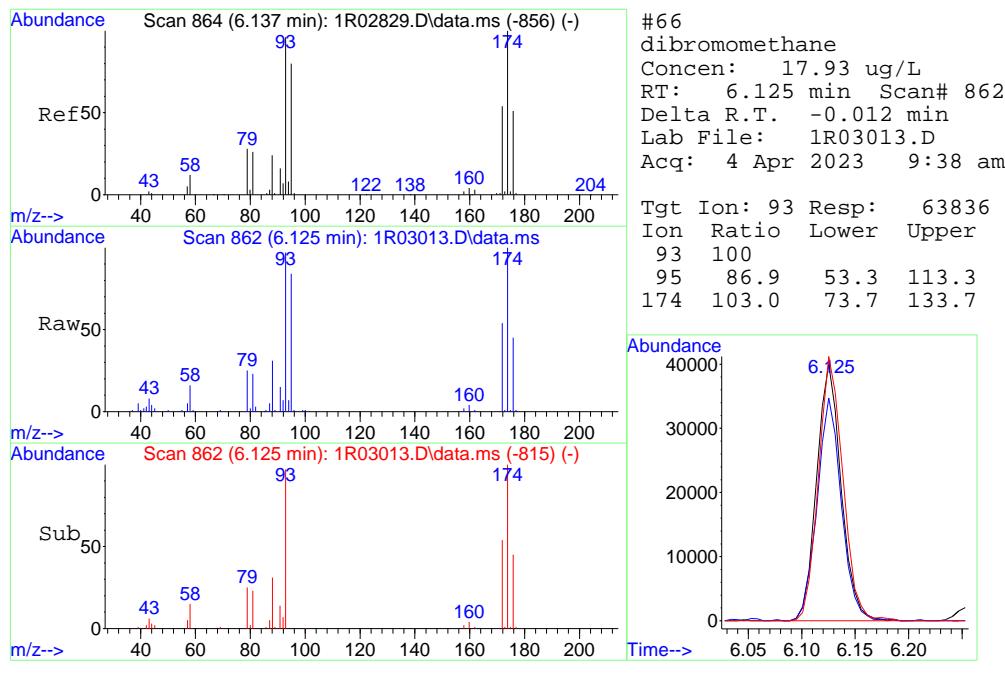
7.6.14

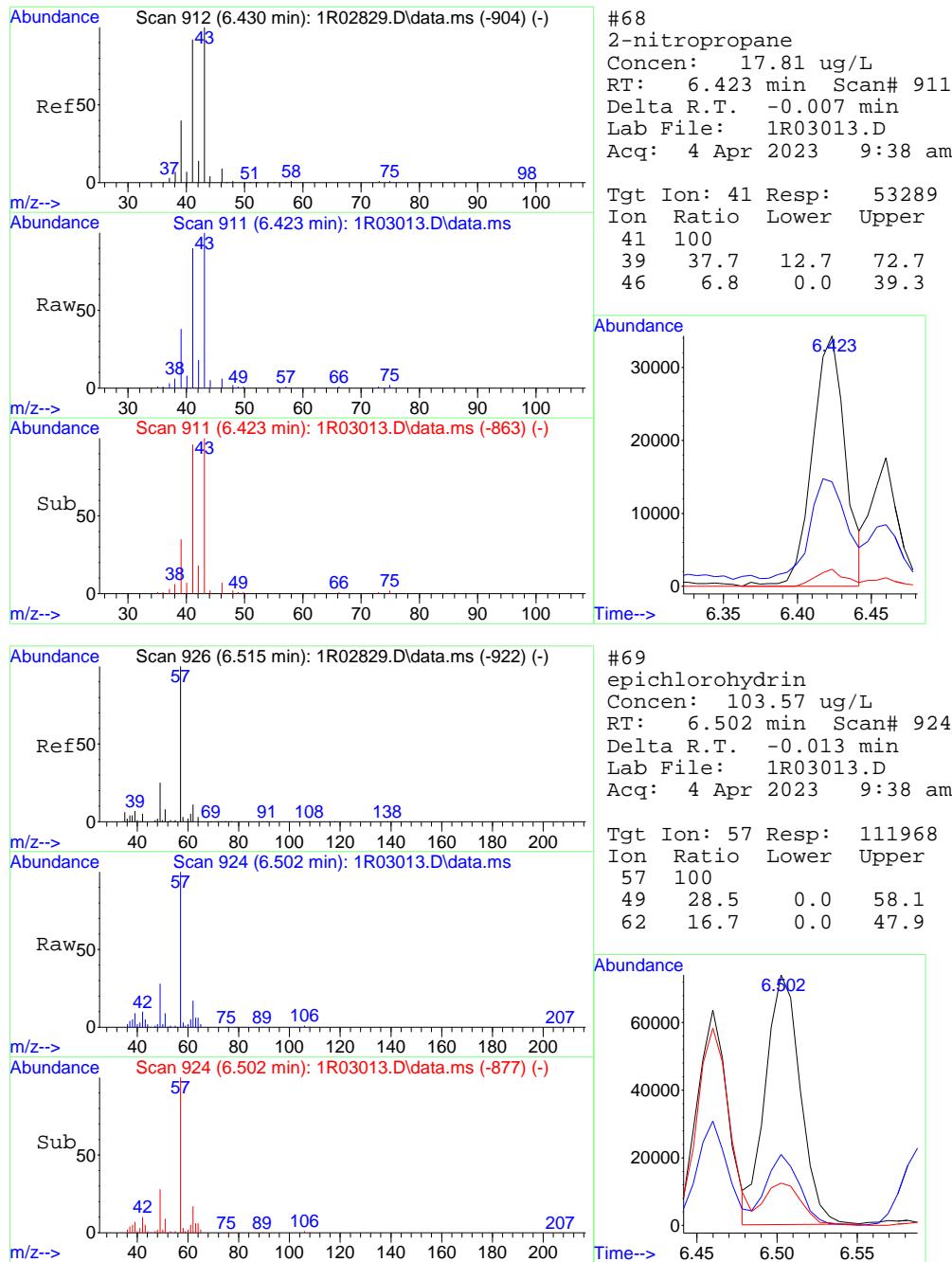


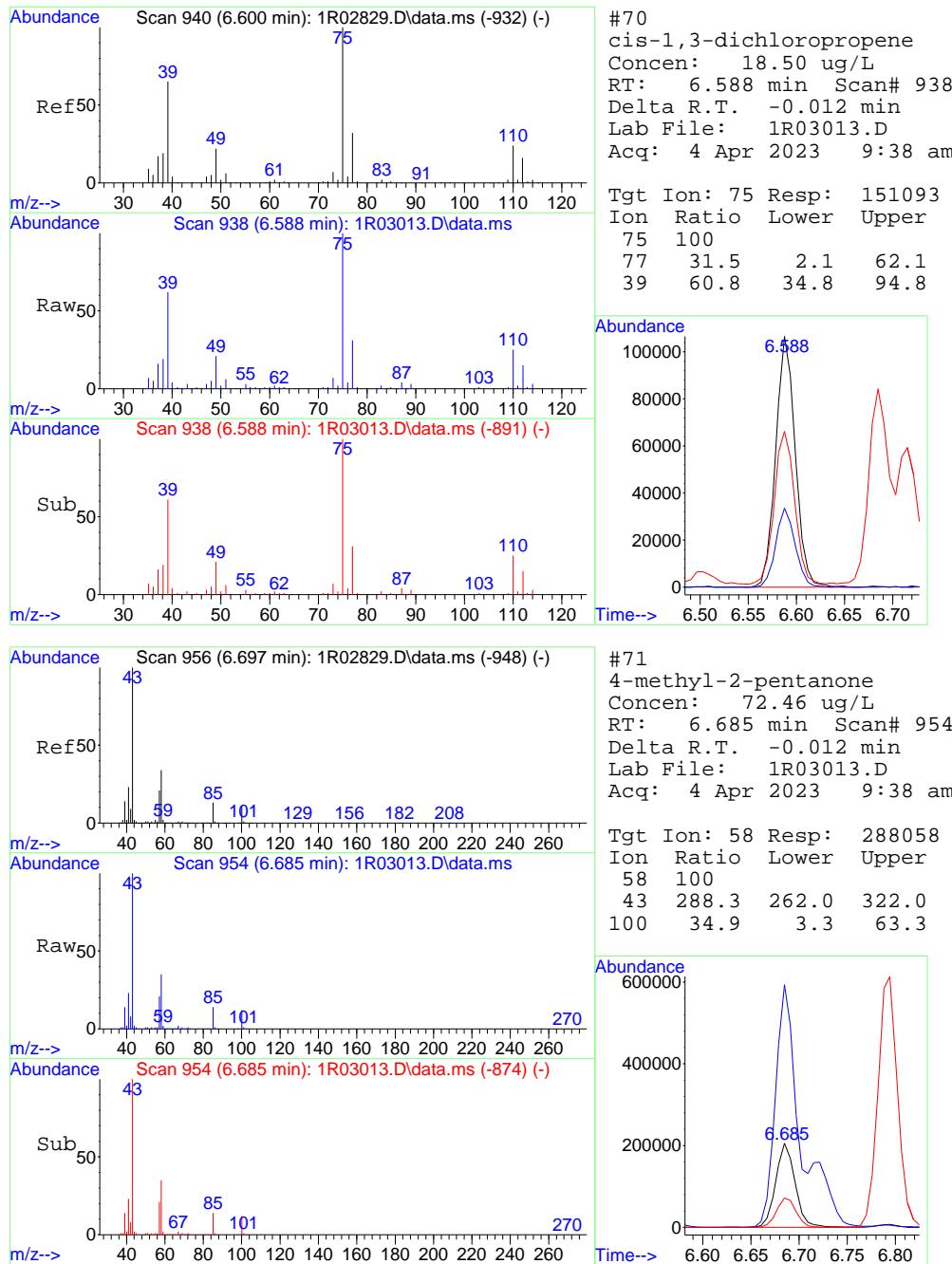


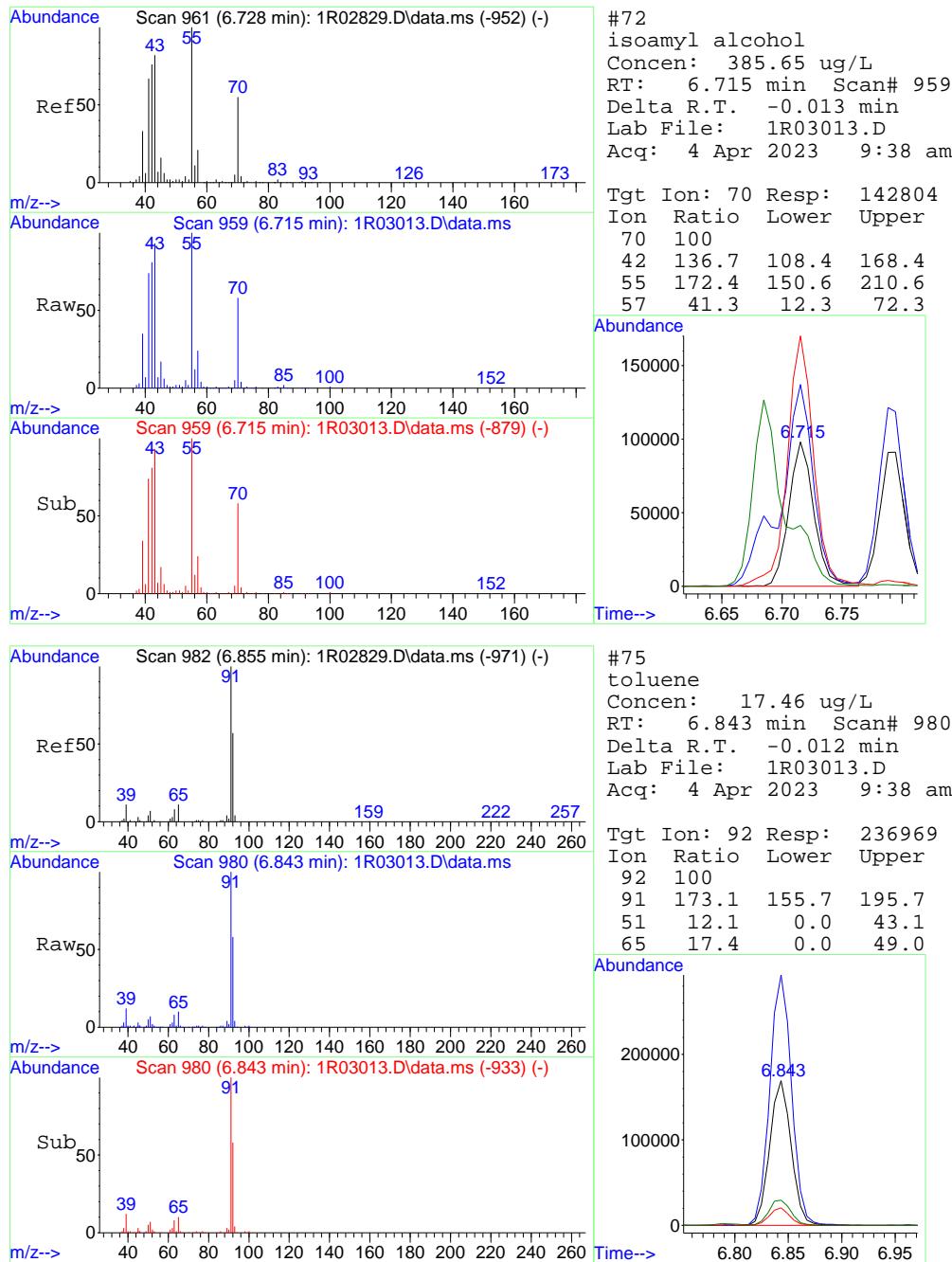


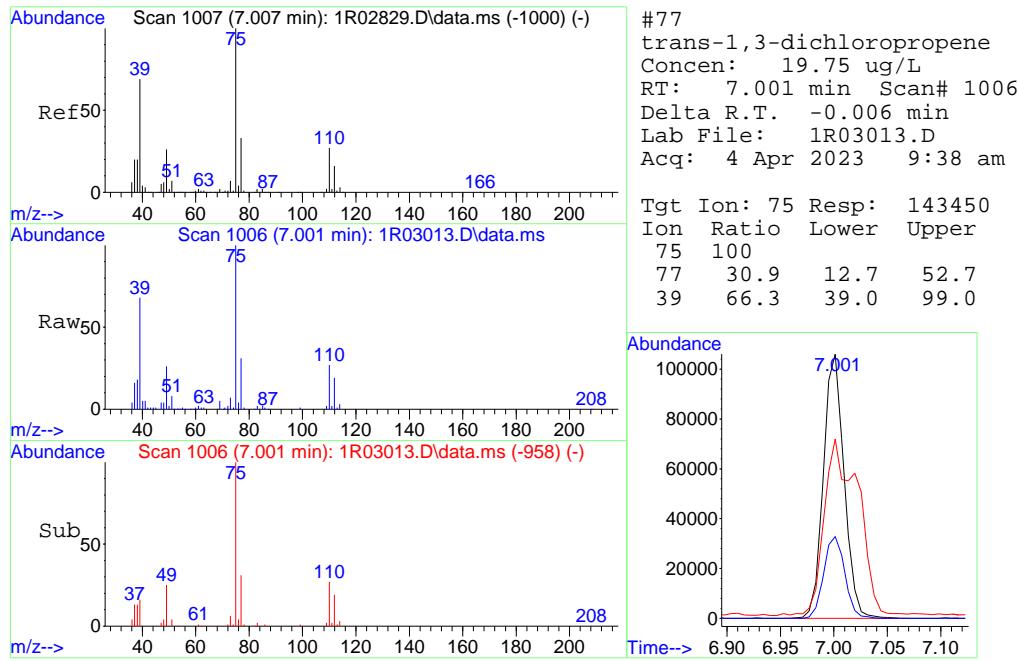
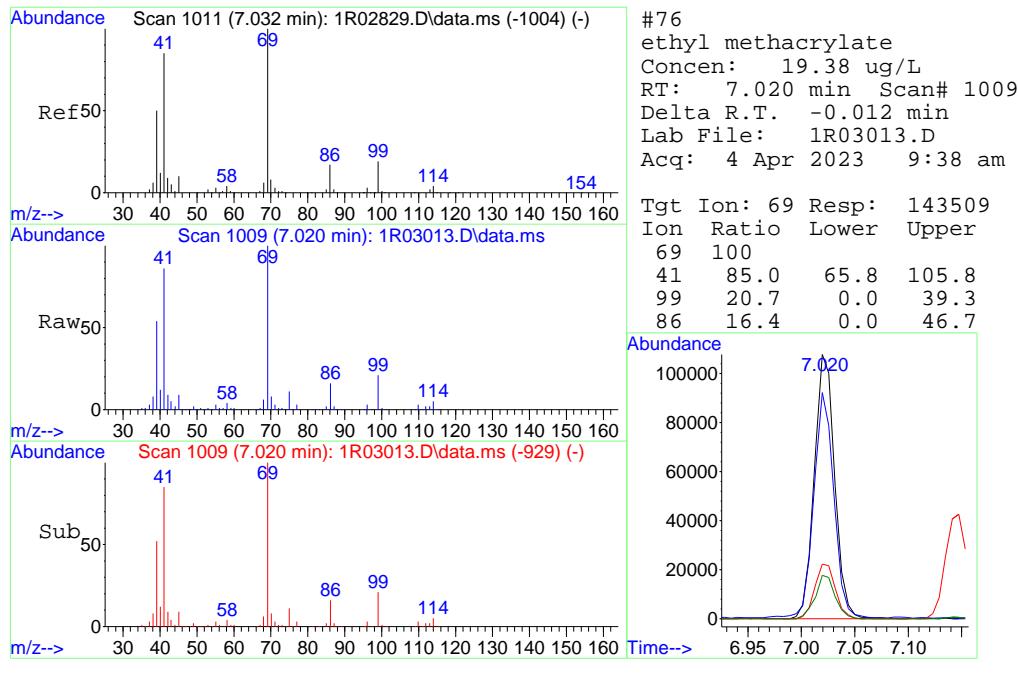


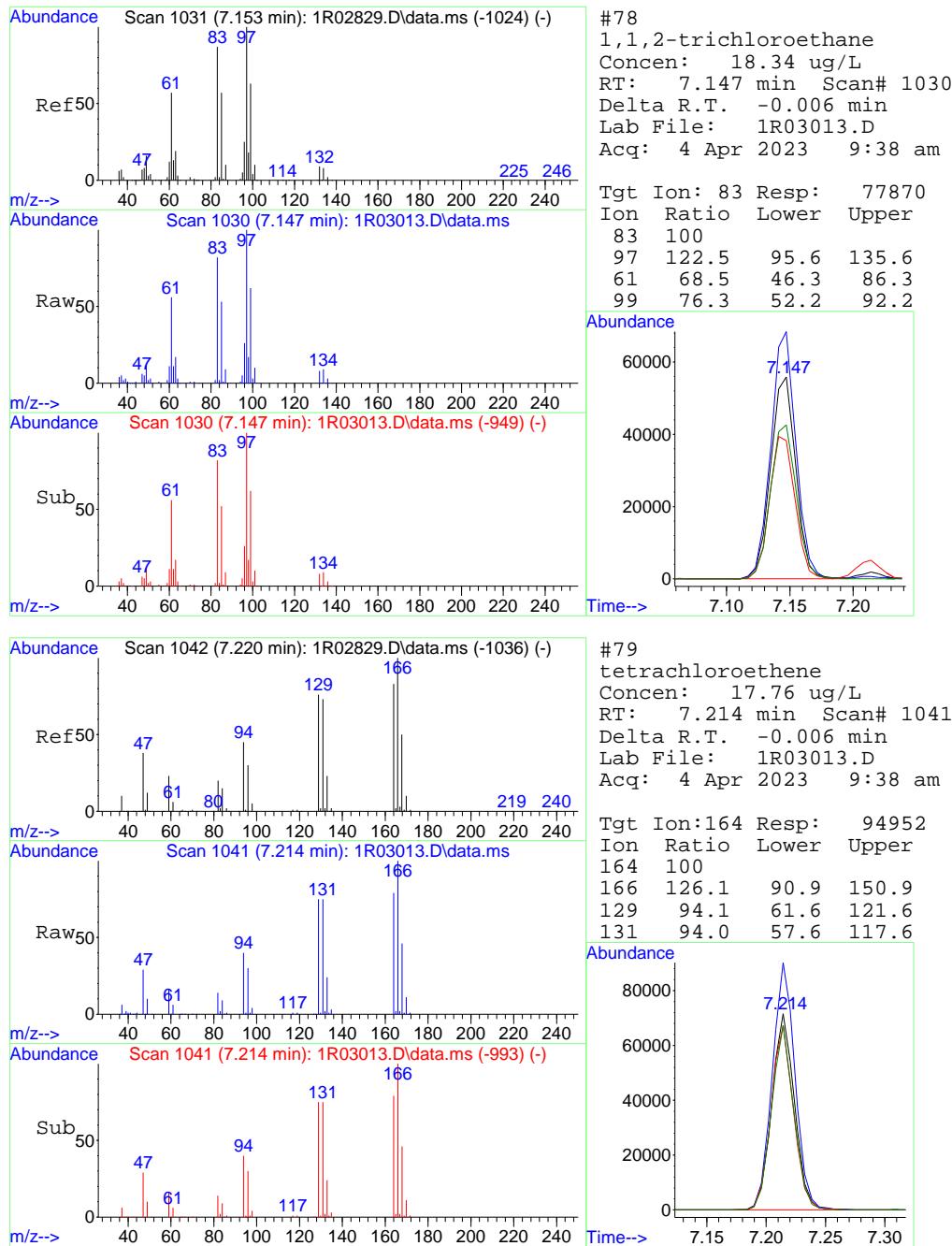


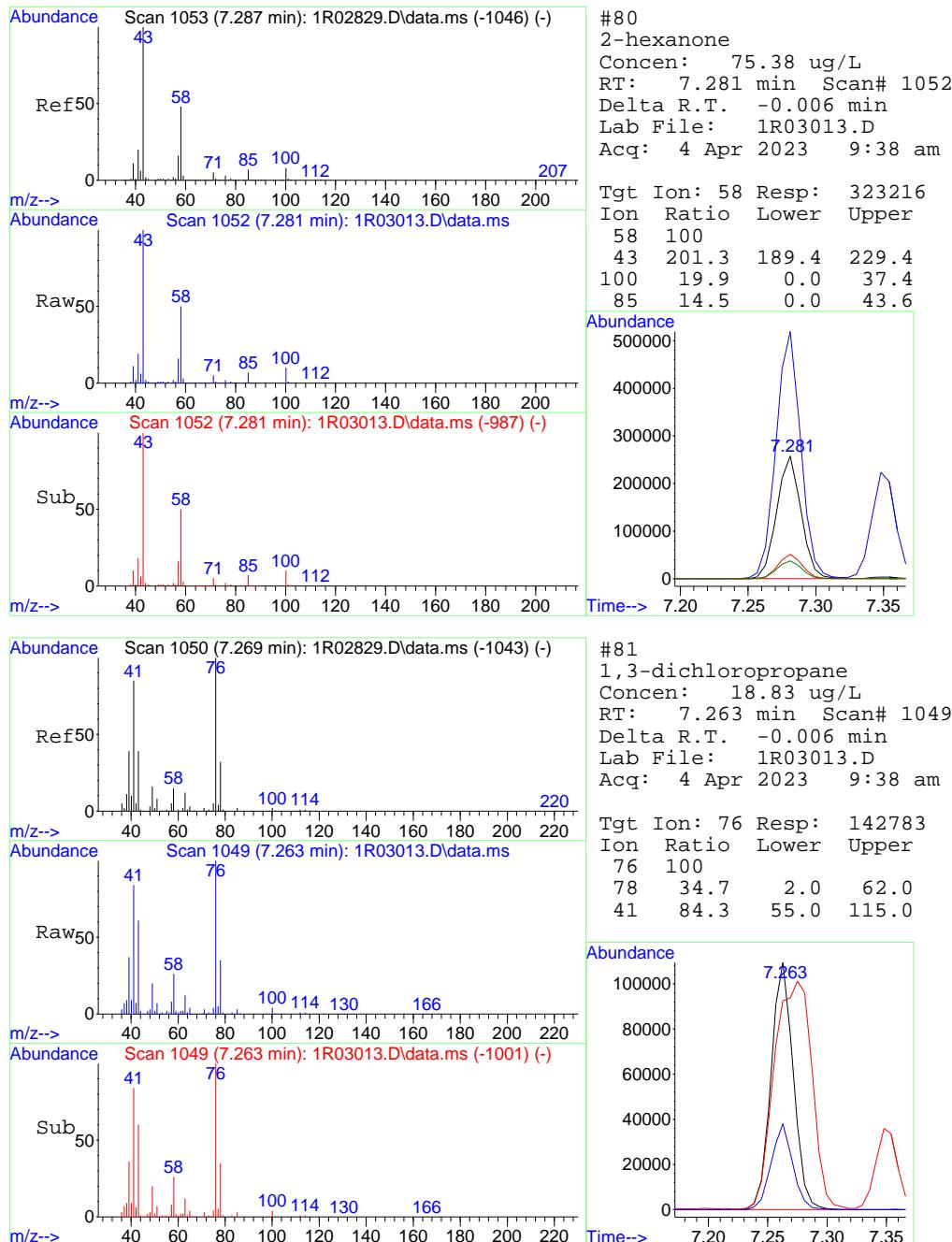


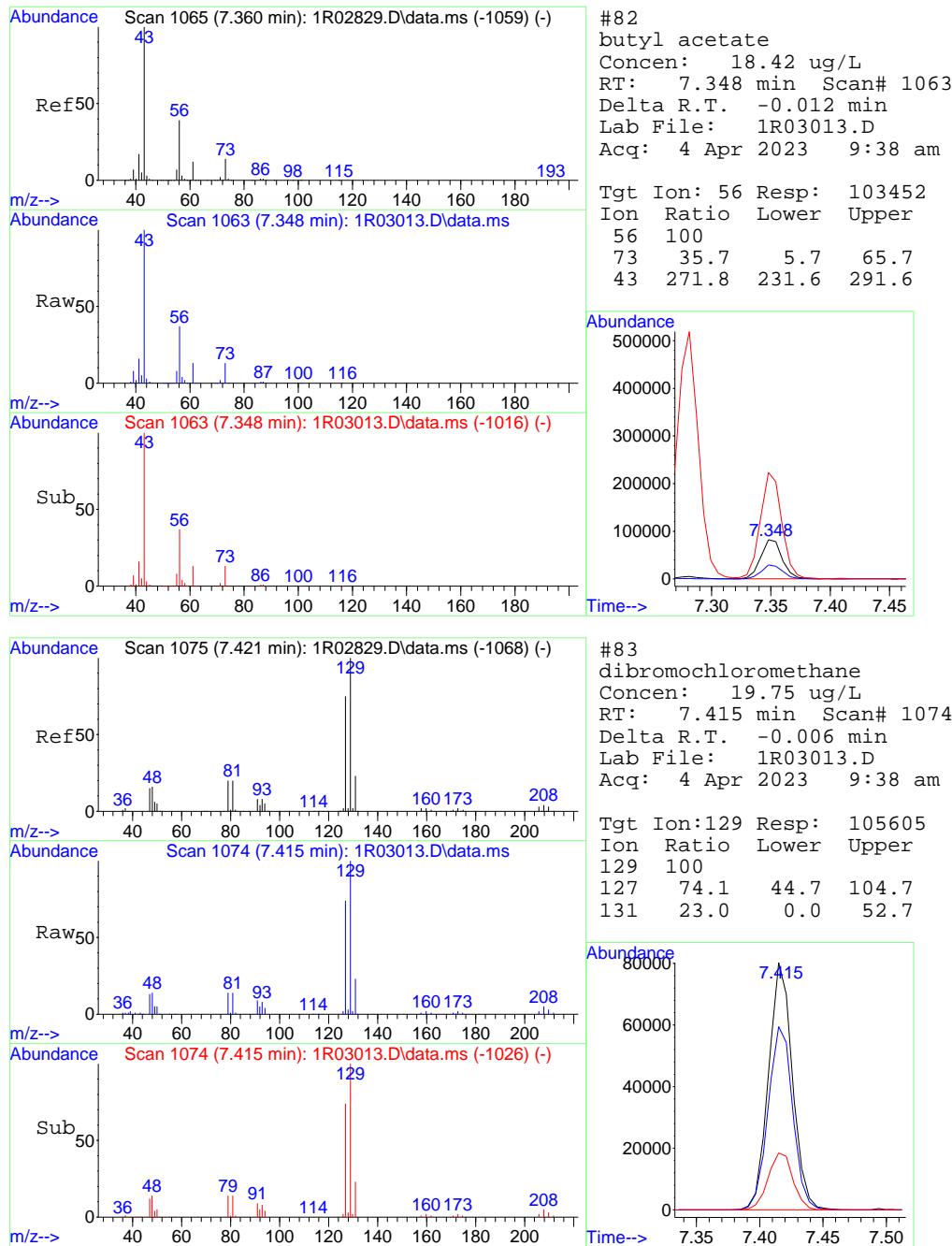


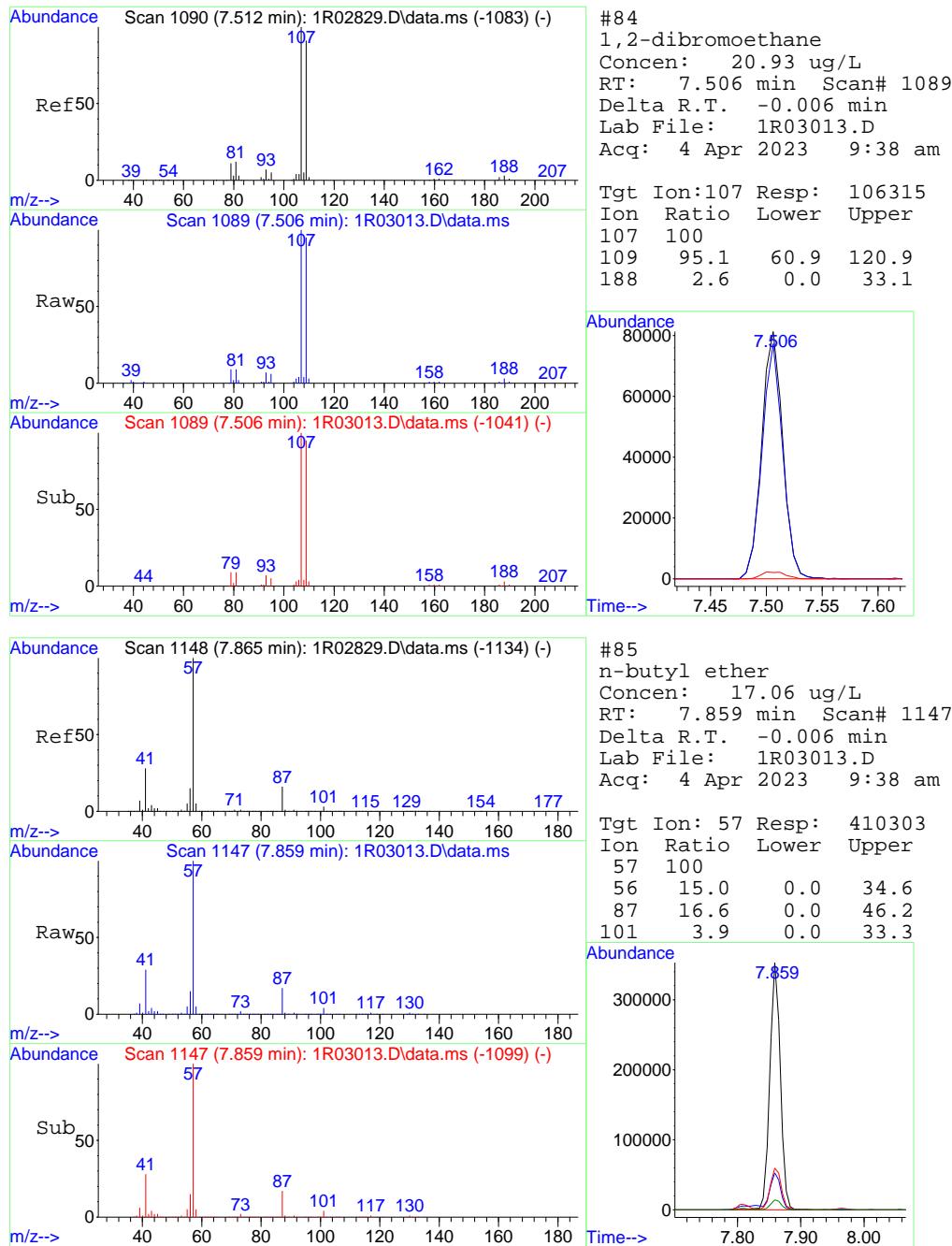


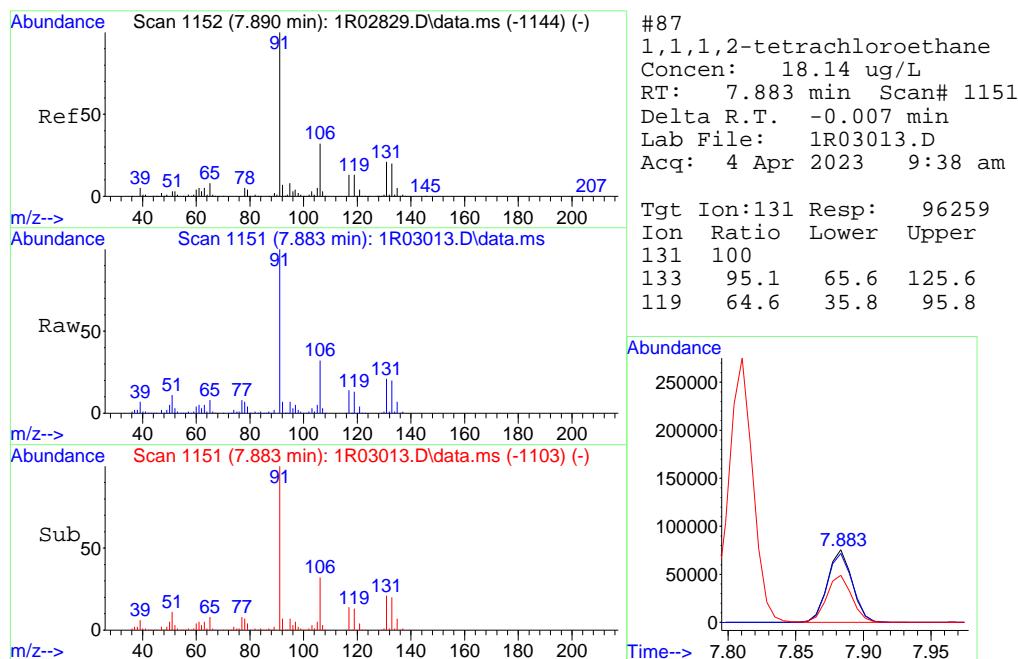
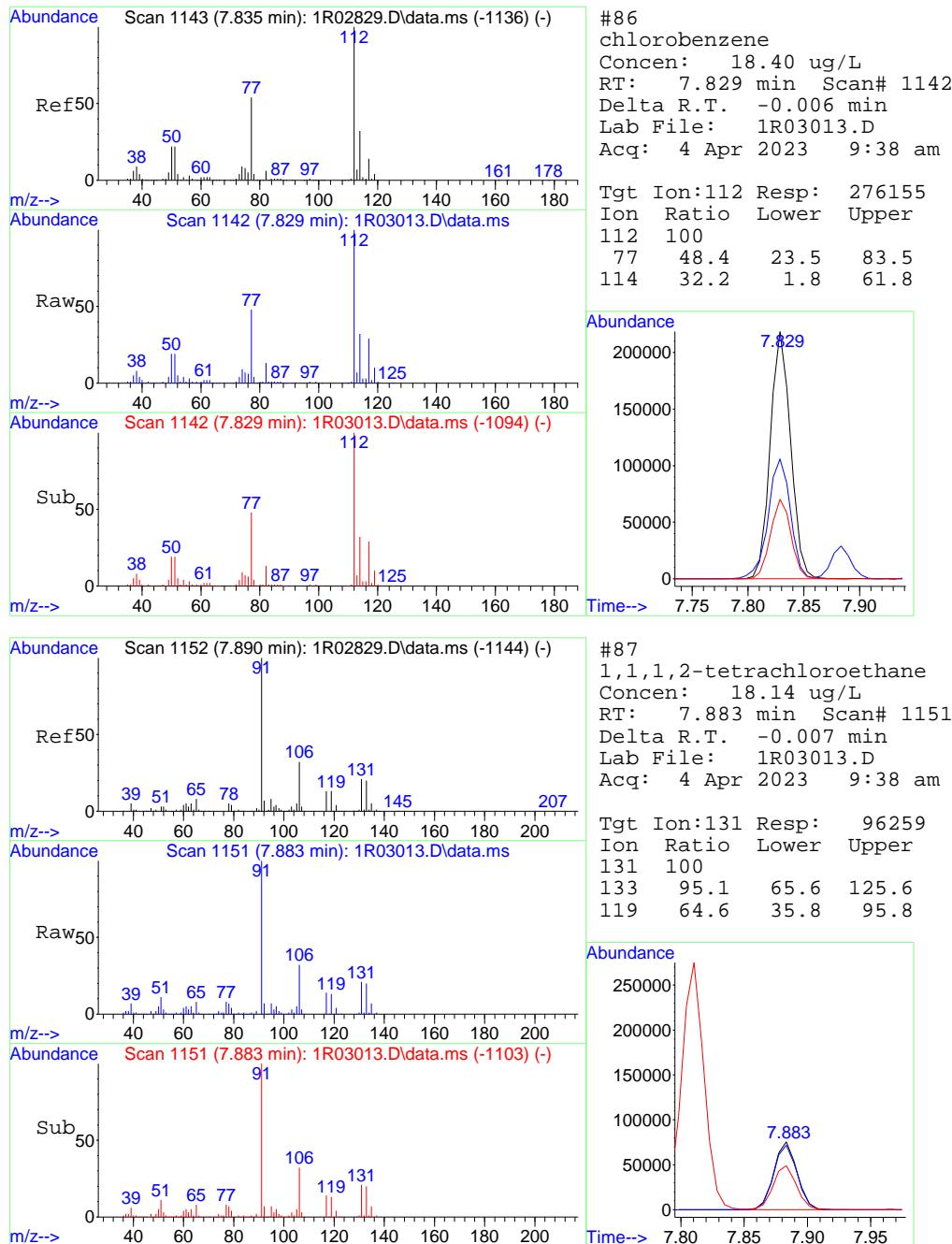


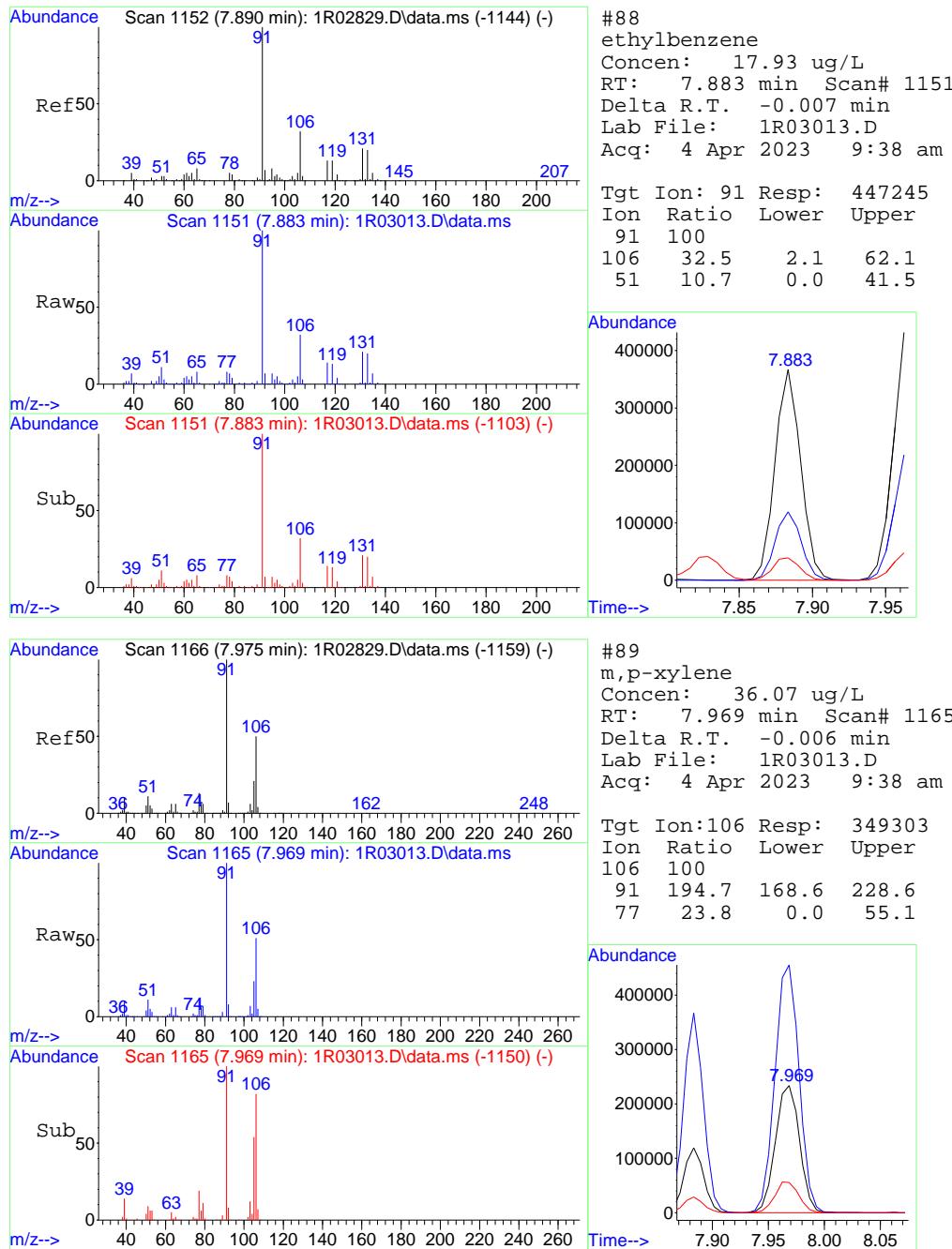


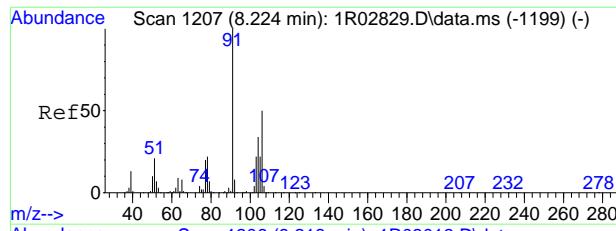






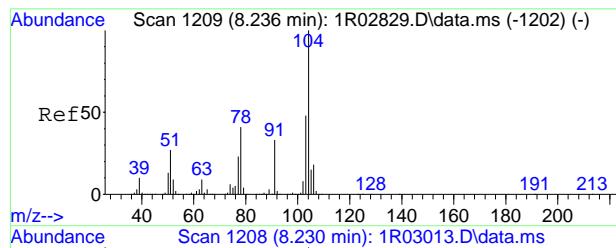
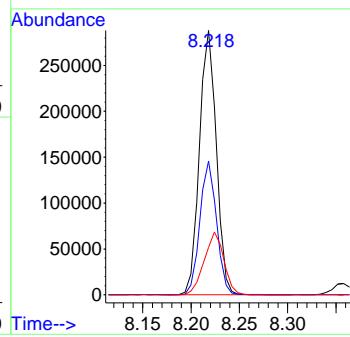
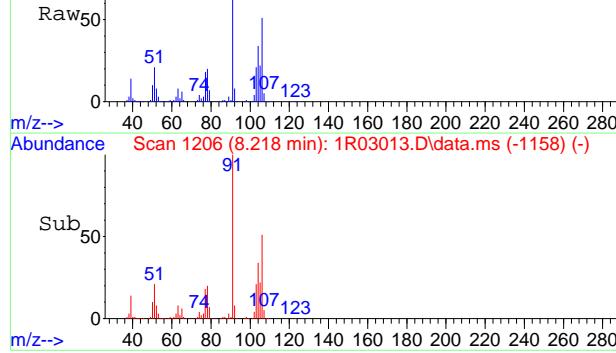
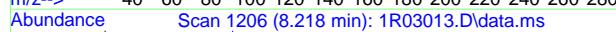






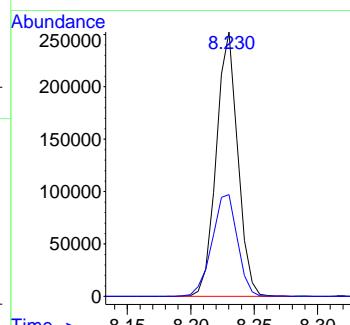
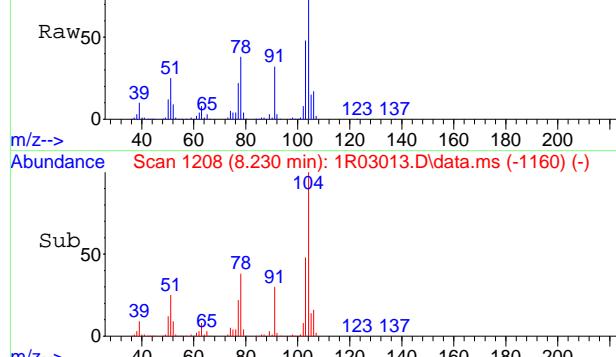
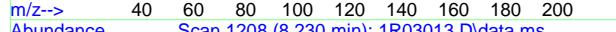
#90  
o-xylene  
Concen: 18.20 ug/L  
RT: 8.218 min Scan# 1206  
Delta R.T. -0.006 min  
Lab File: 1R03013.D  
Acq: 4 Apr 2023 9:38 am

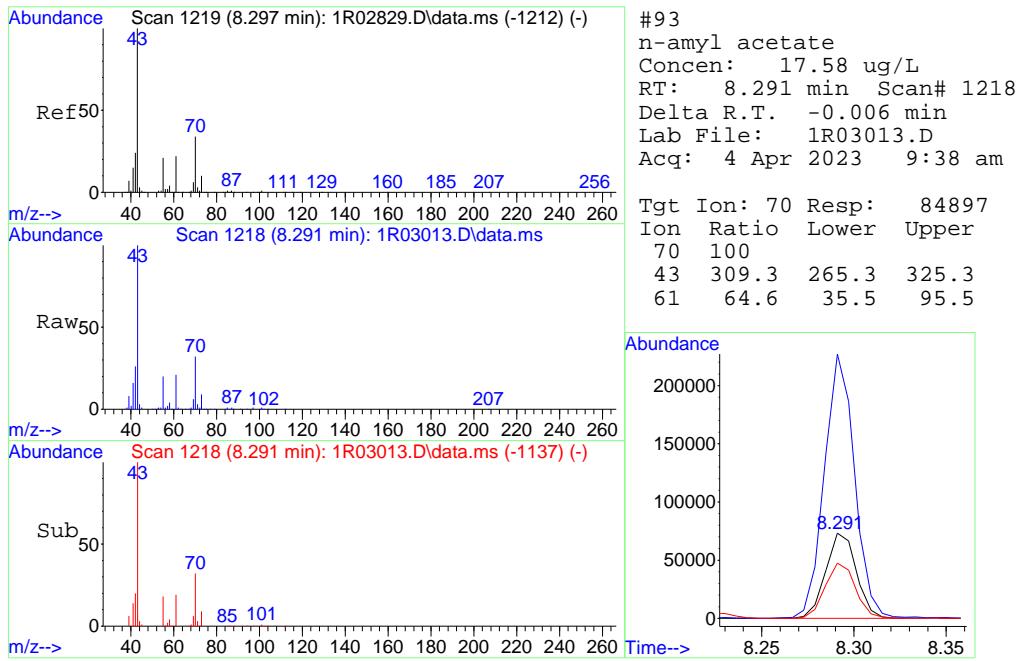
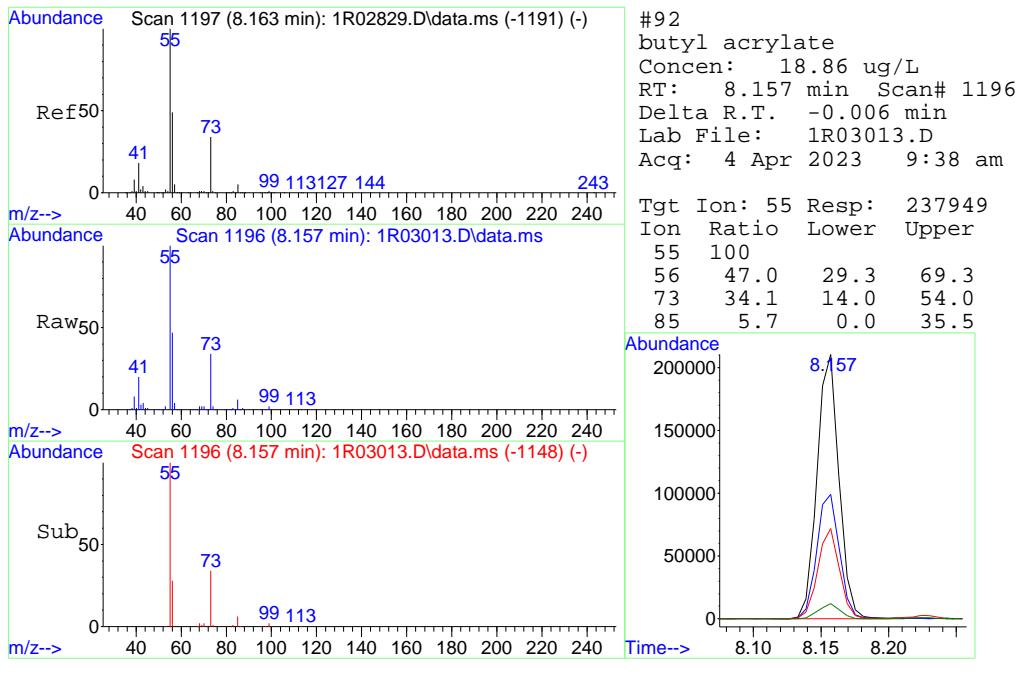
Tgt Ion: 91 Resp: 350219  
Ion Ratio Lower Upper  
91 100  
106 50.6 20.2 80.2  
77 18.0 0.0 50.1

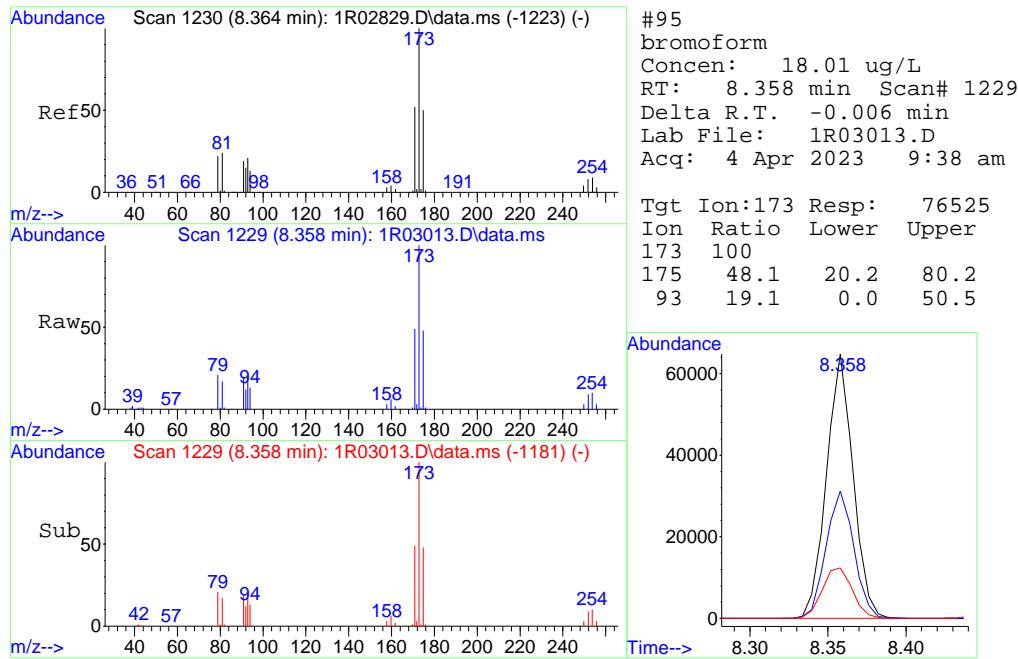
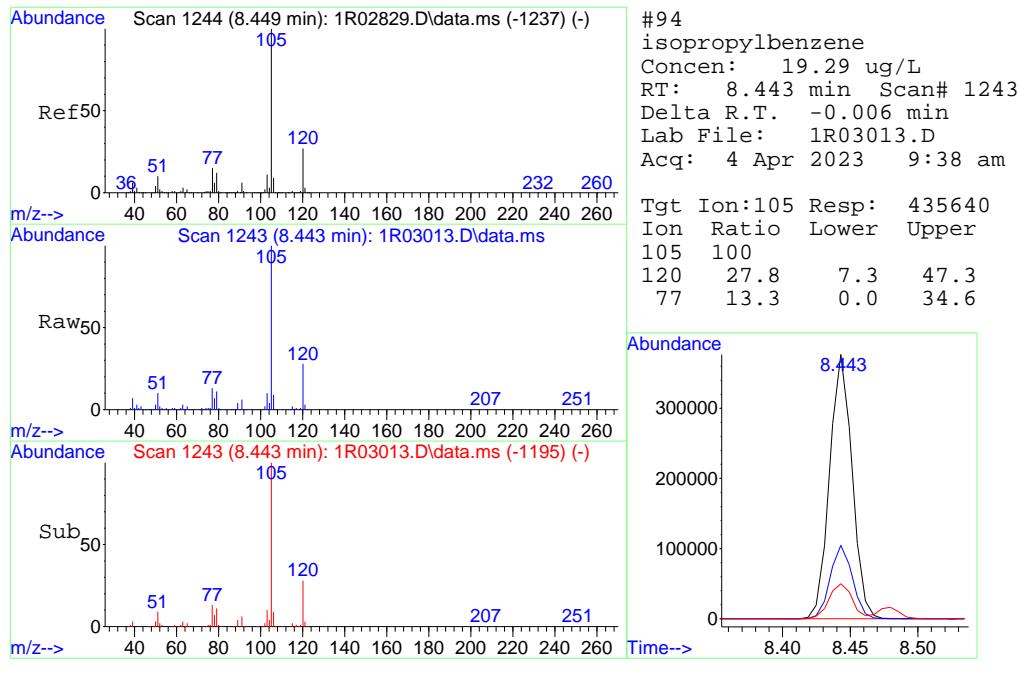


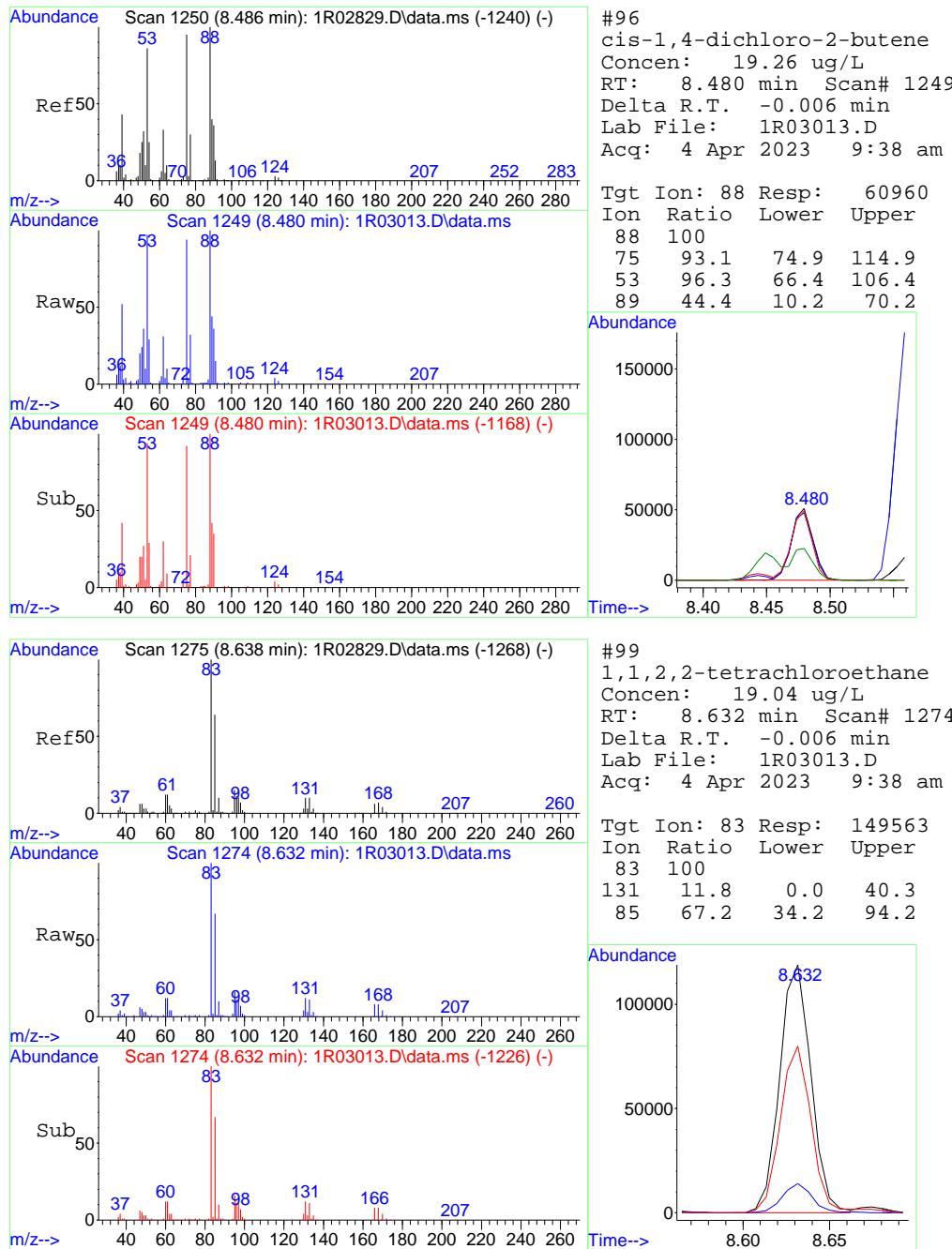
#91  
styrene  
Concen: 18.42 ug/L  
RT: 8.230 min Scan# 1208  
Delta R.T. -0.006 min  
Lab File: 1R03013.D  
Acq: 4 Apr 2023 9:38 am

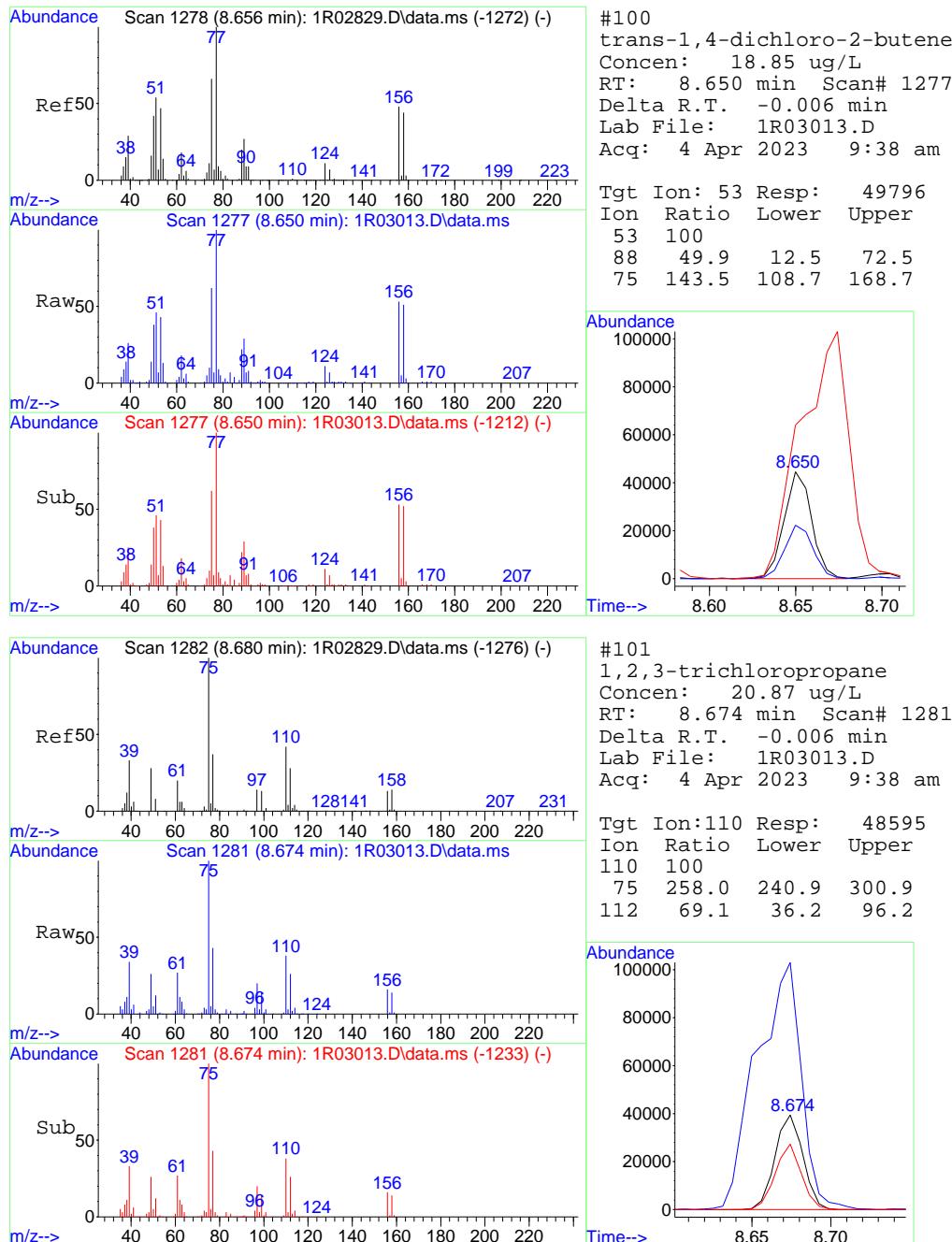
Tgt Ion: 104 Resp: 297520  
Ion Ratio Lower Upper  
104 100  
78 38.5 11.4 71.4

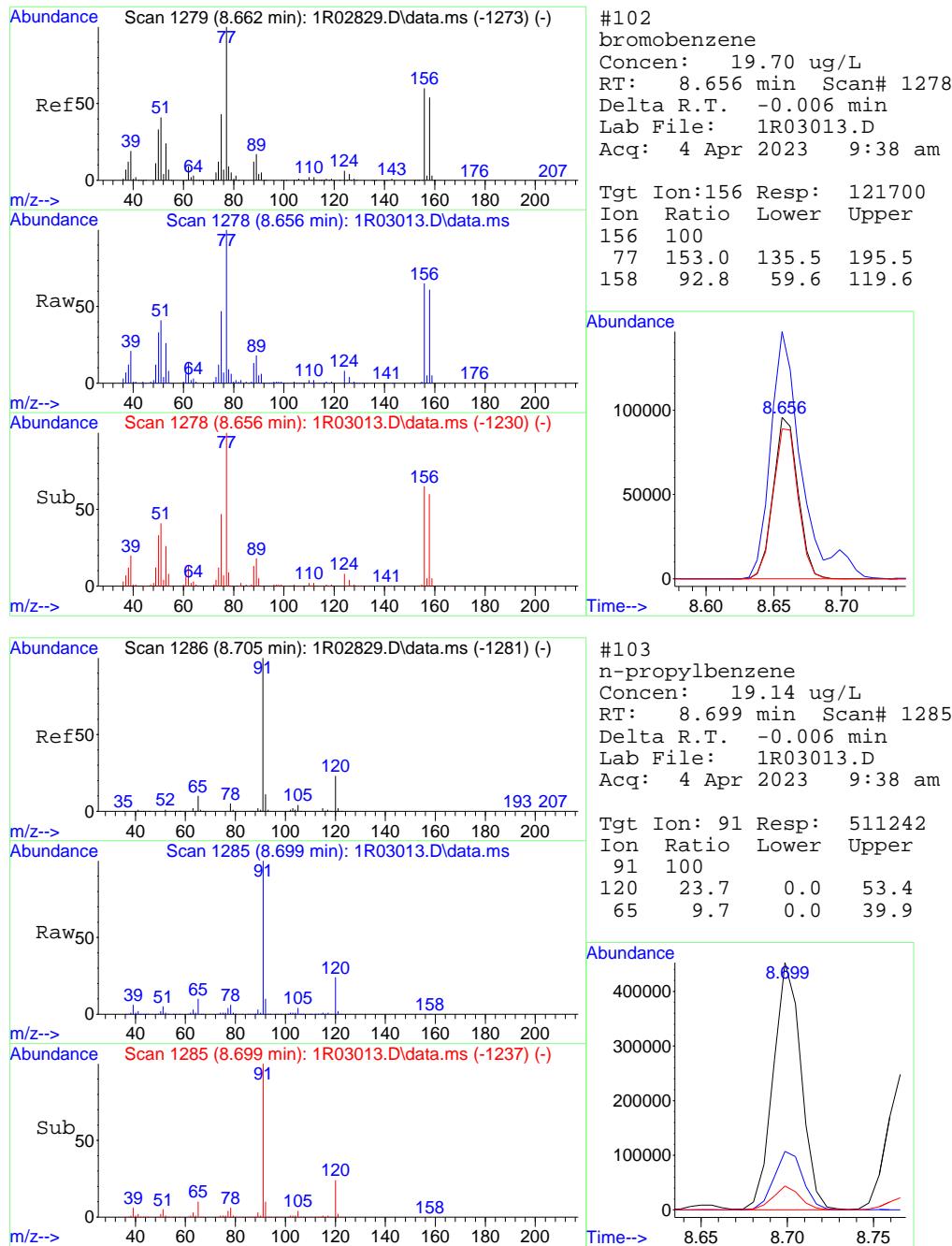


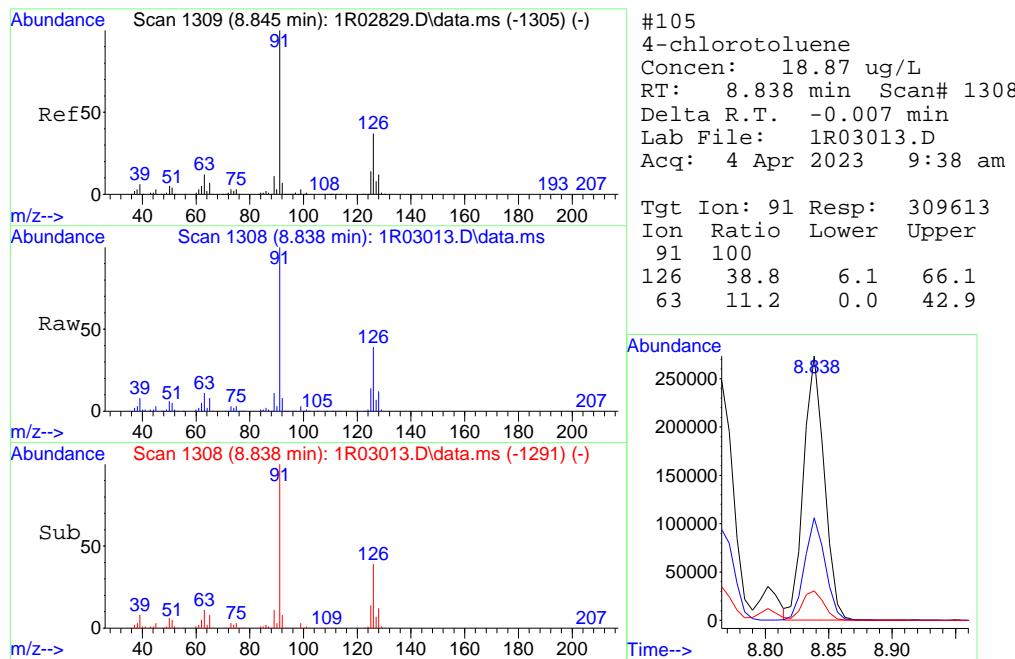
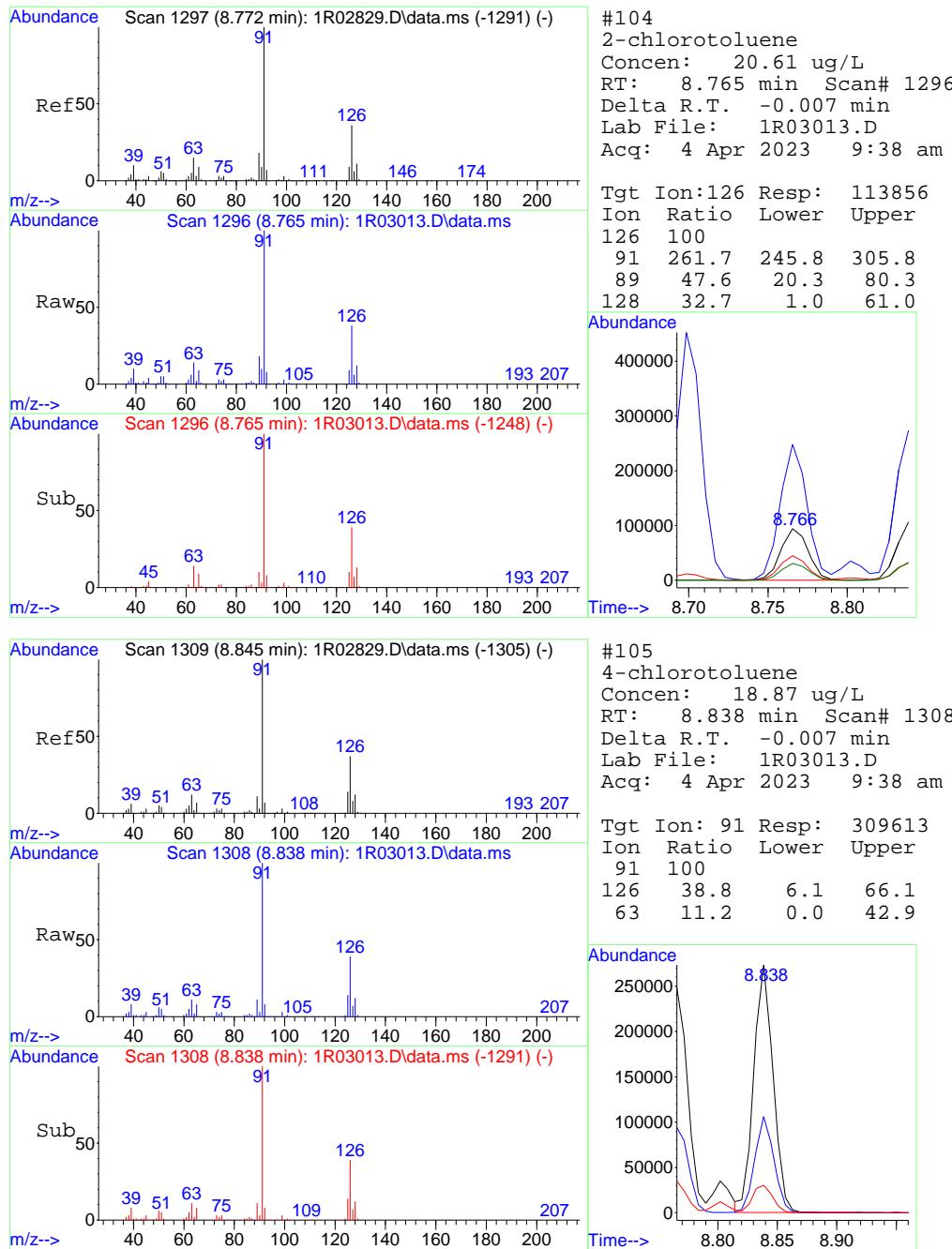


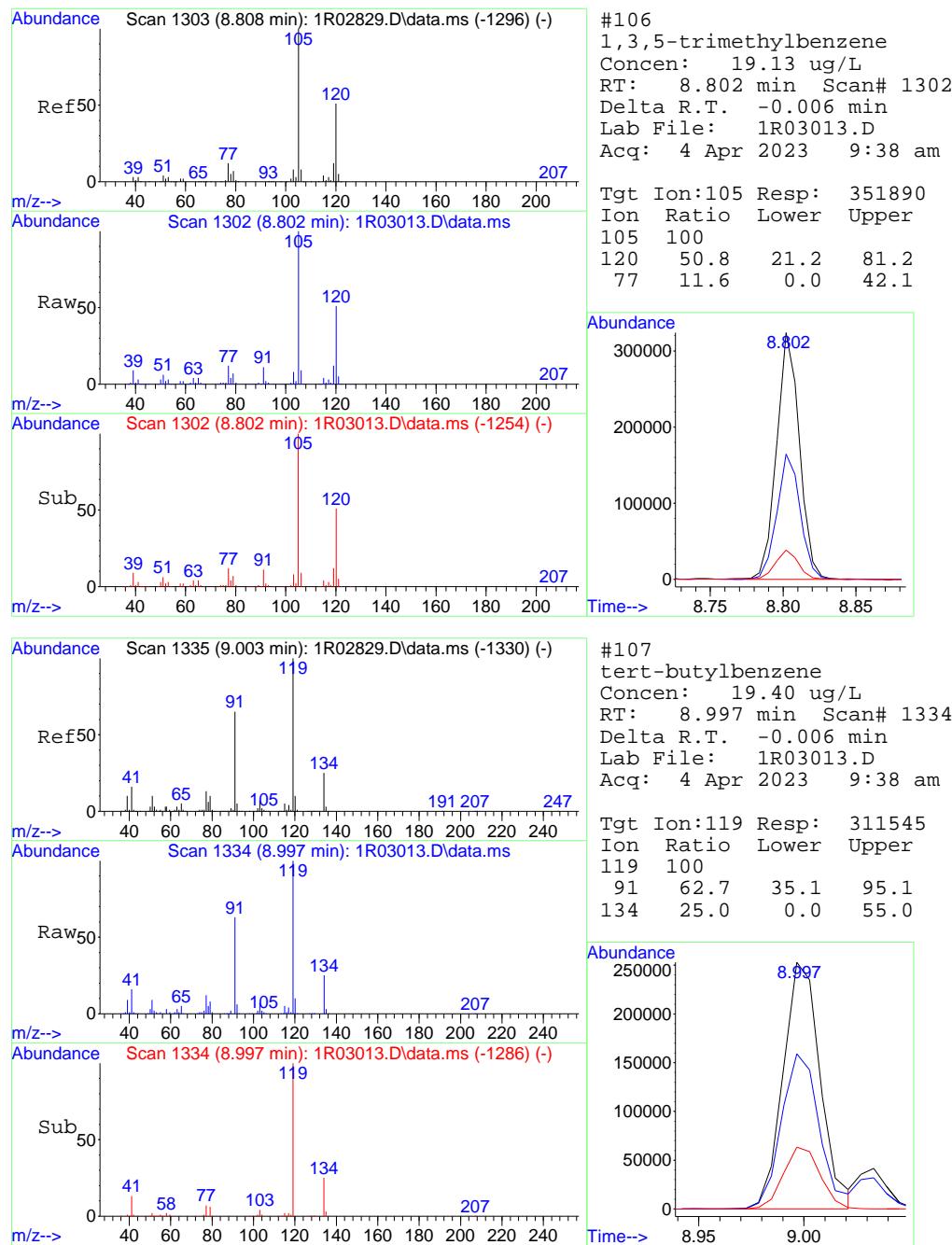


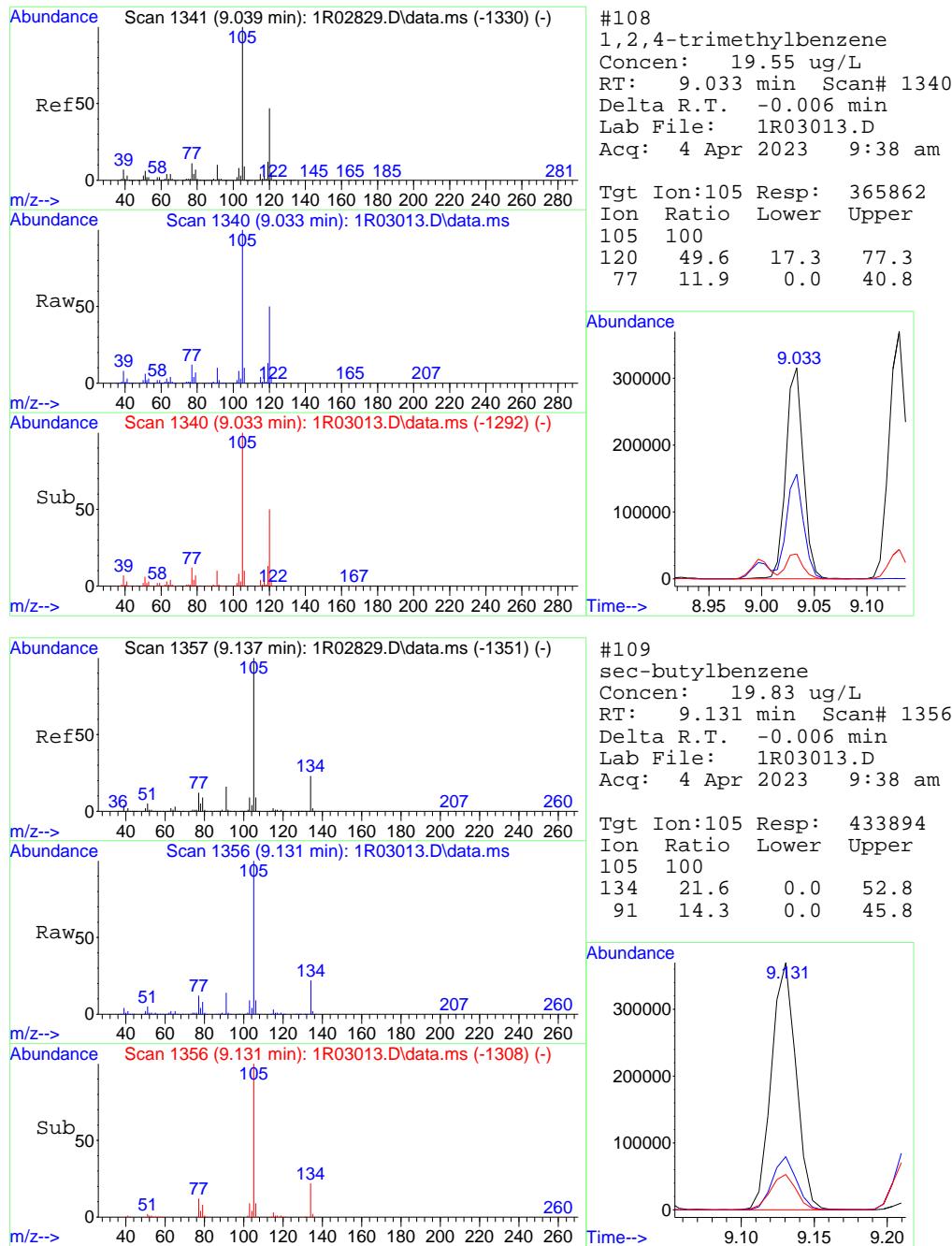


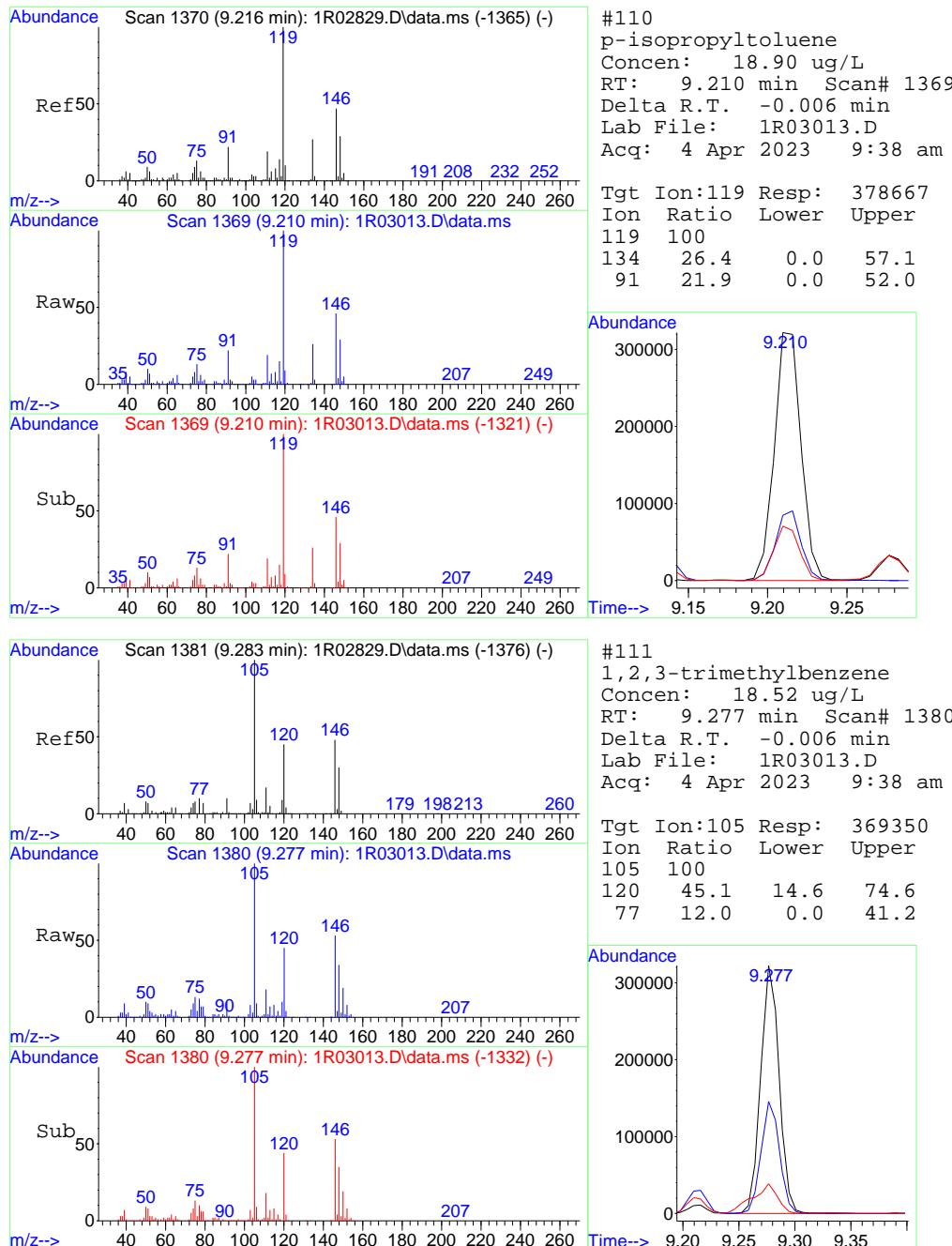


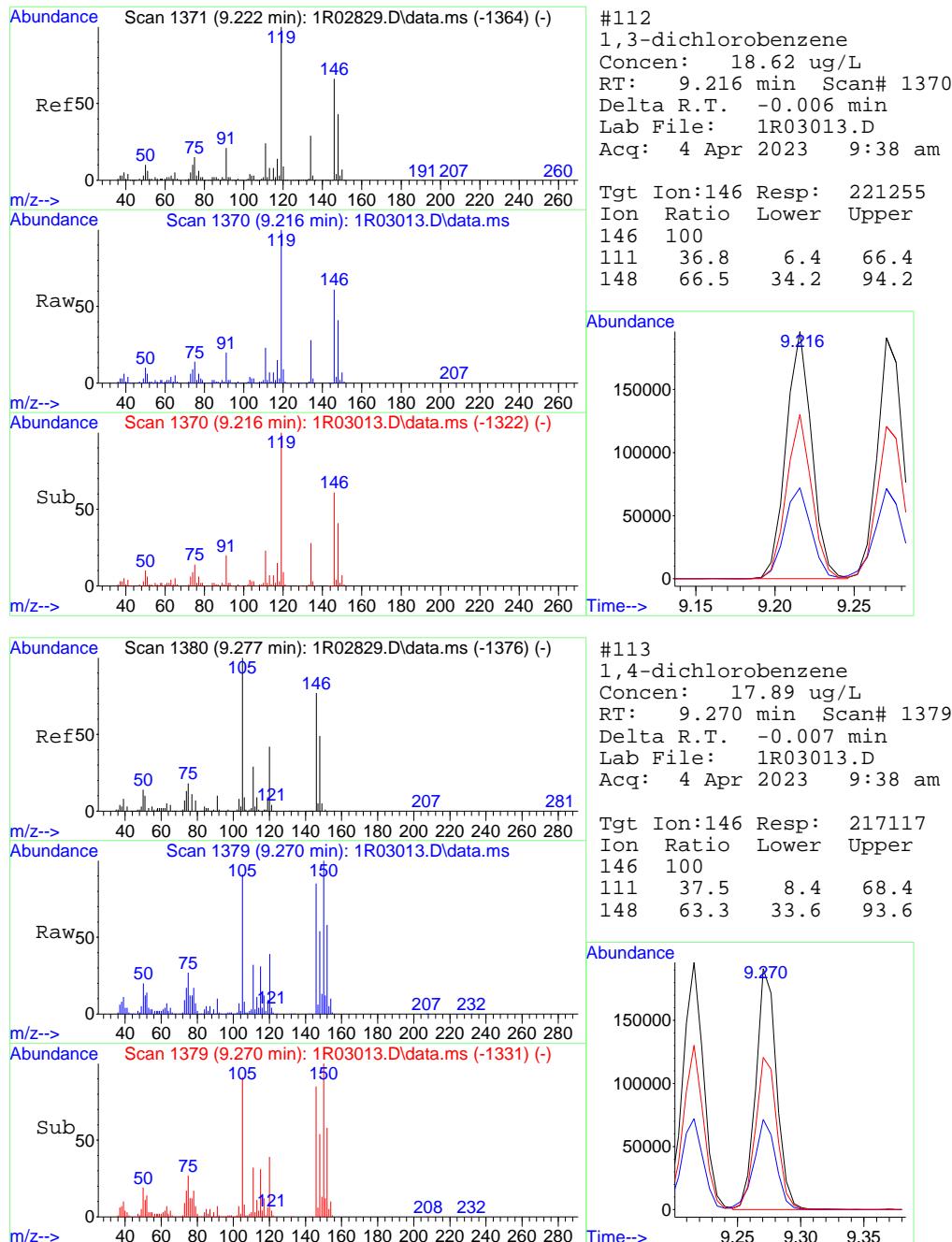


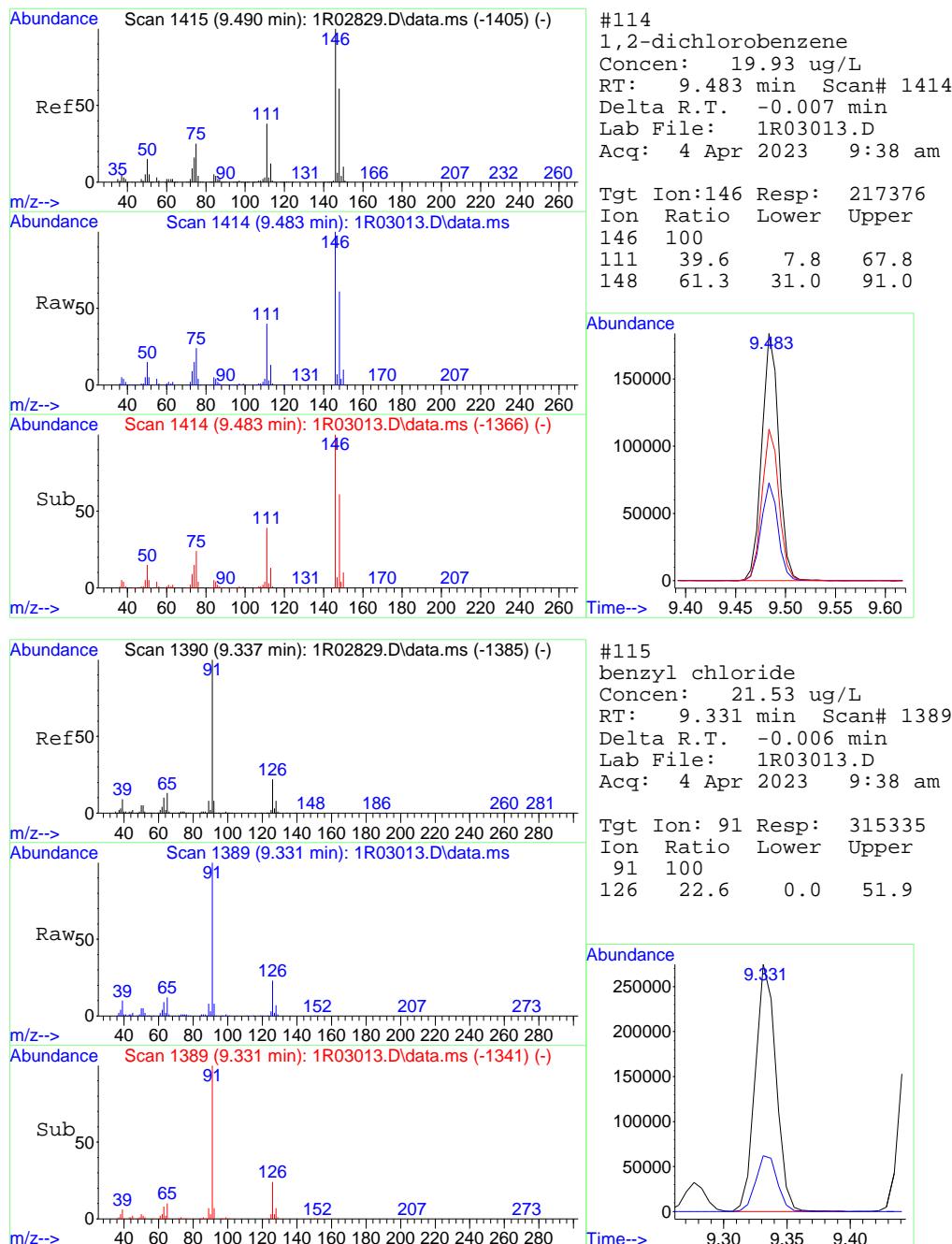


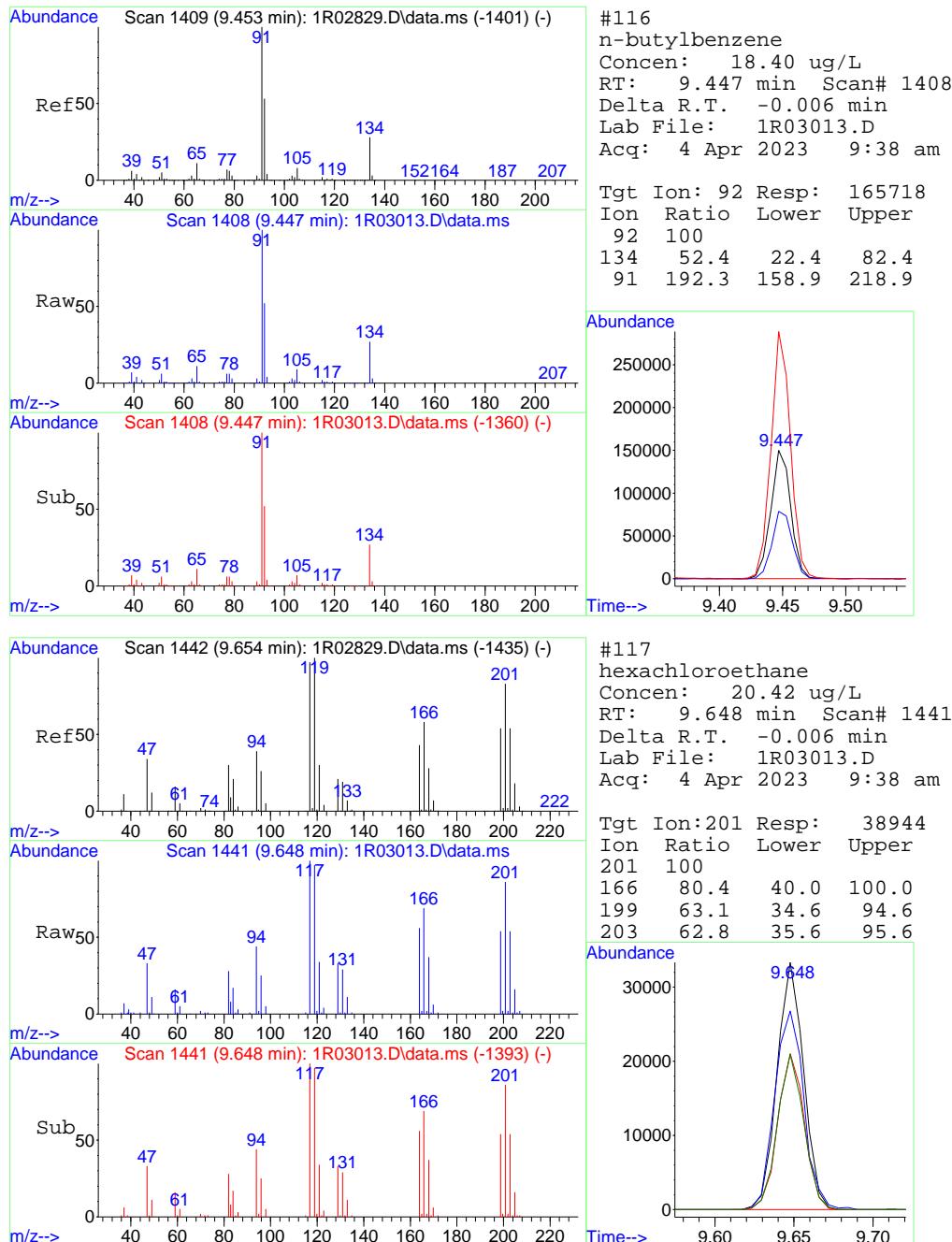


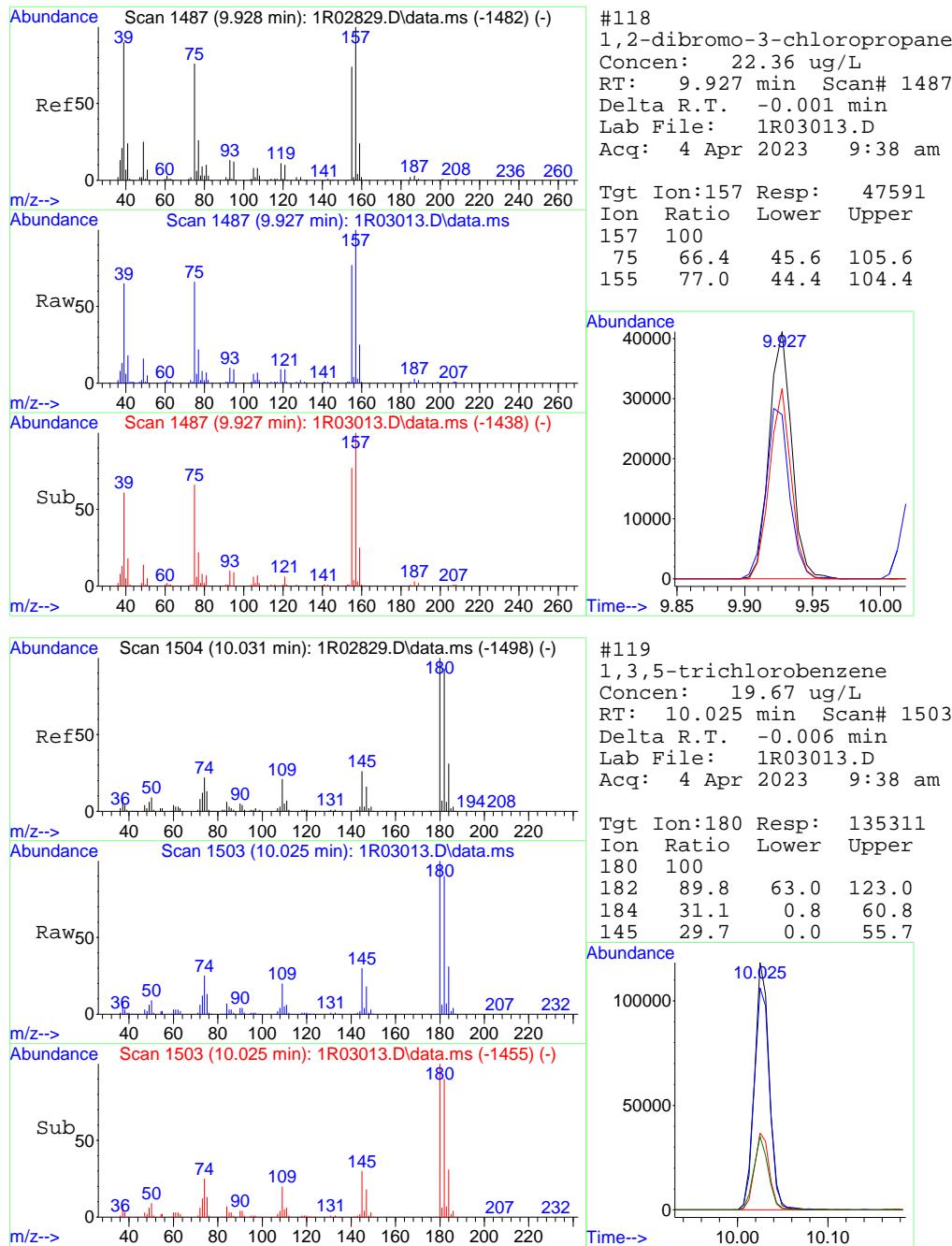


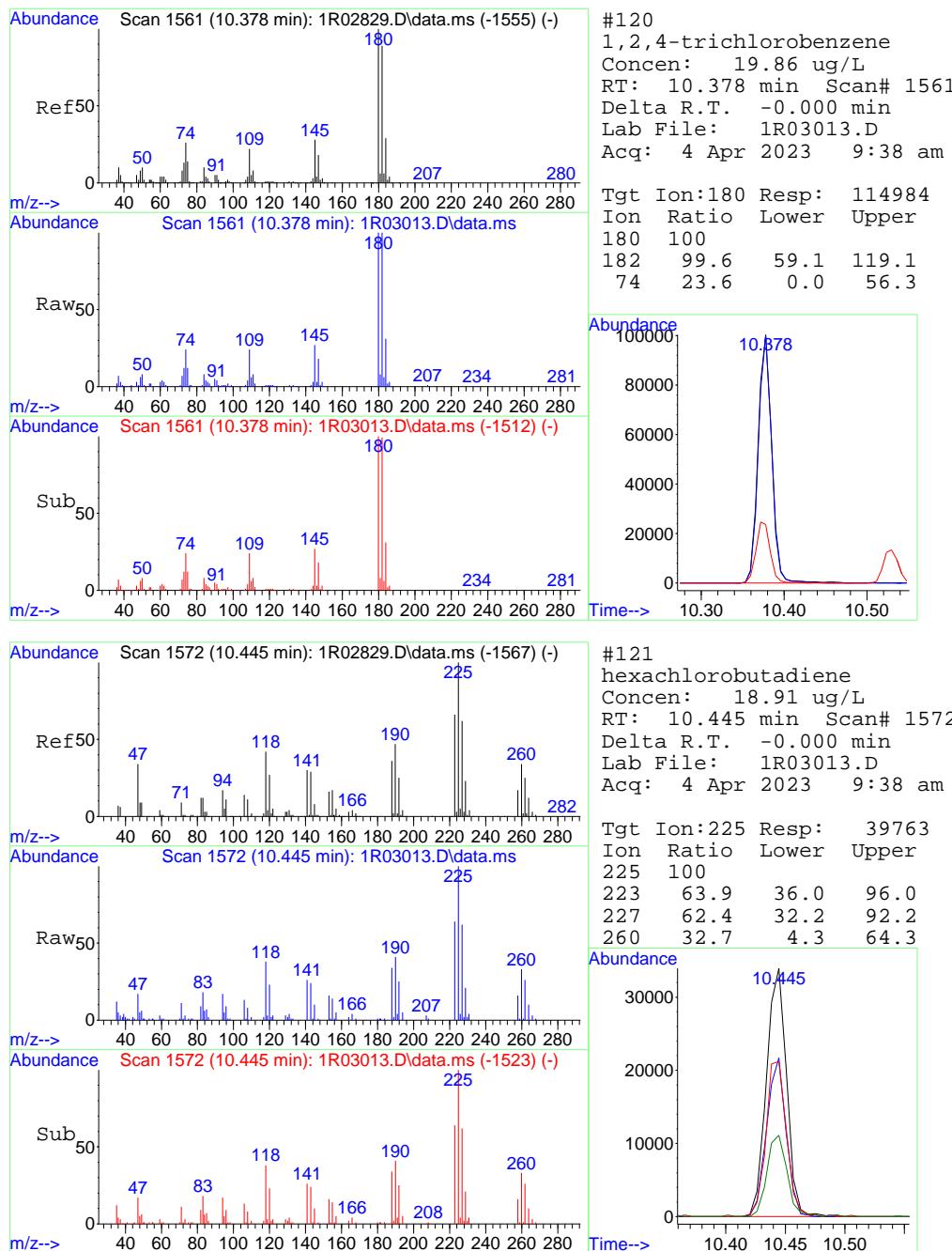


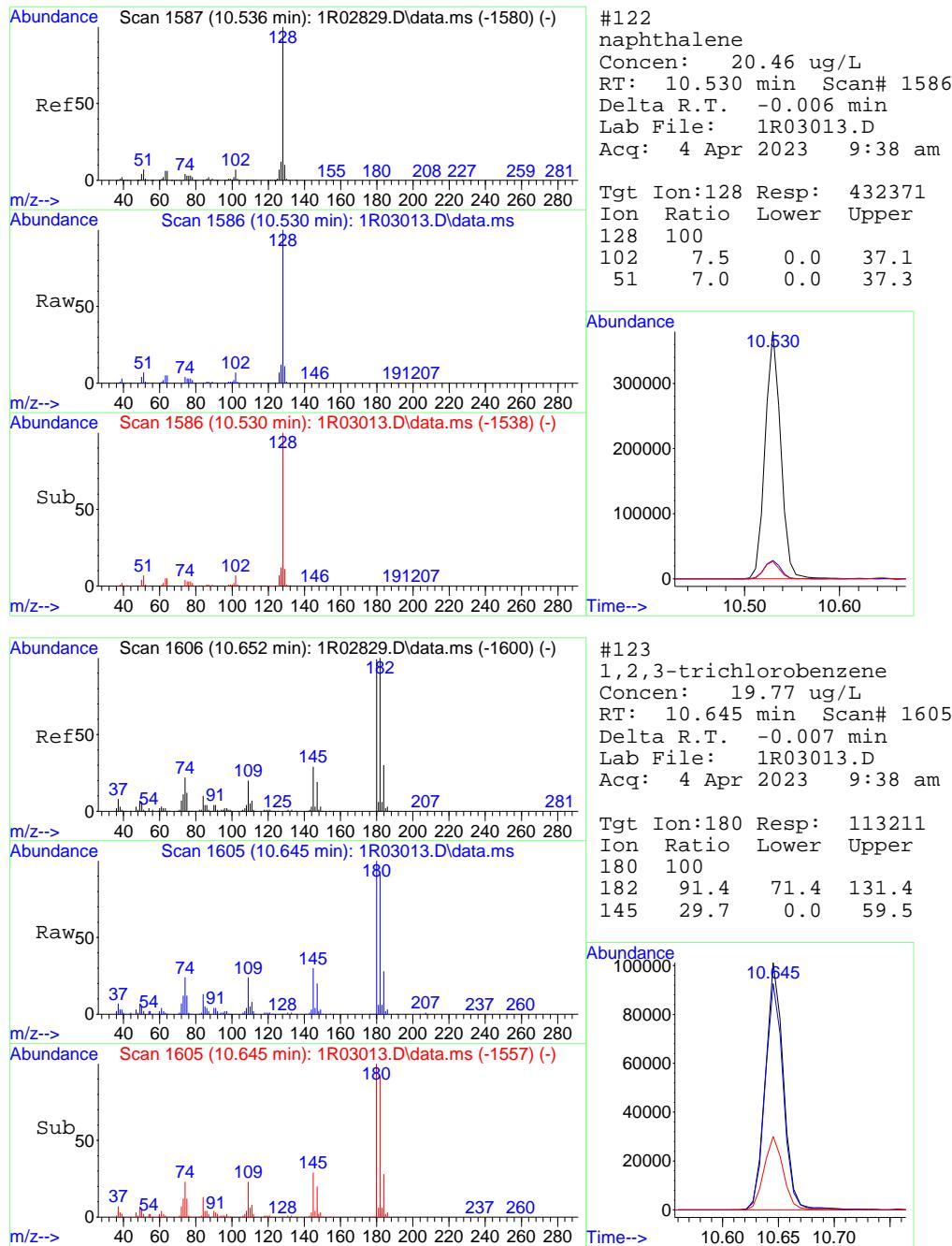


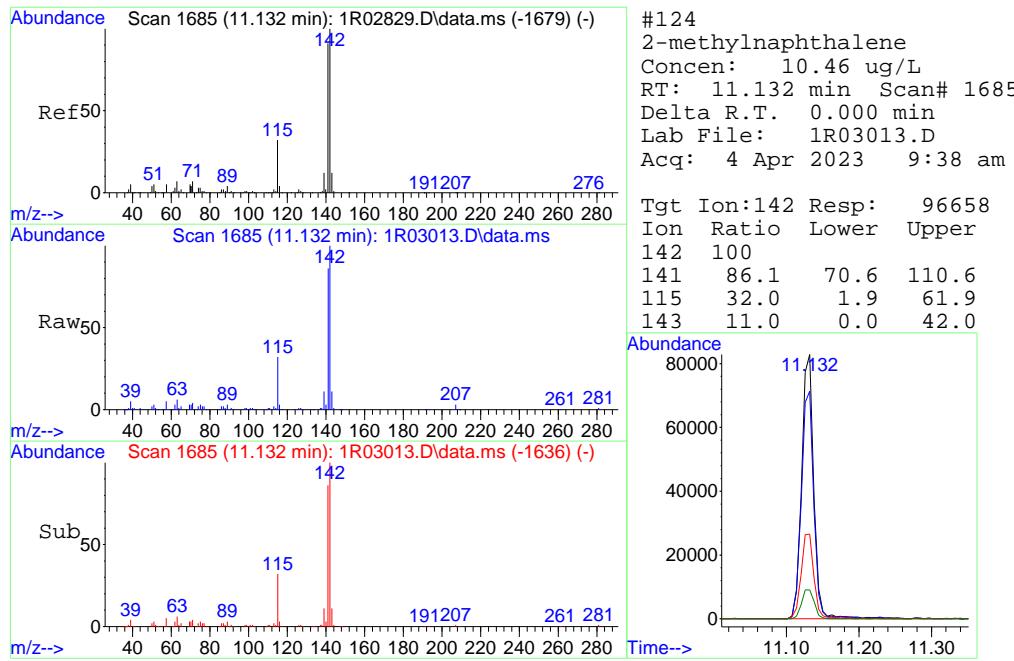














## VOLATILE by GCMS ANALYSIS LOG

Date: 04/04/23

Standard data

Lot#	Expiration Date	concentration
V023-2768-36	4/28/2023	100PPM
V023-2768-9	4/10/2023	100PPM
V023-2768-4	4/24/2023	100-10000PPM
		Internal/Surrogate data

Lot#	Expiration Date	concentration
V023-2768-5	5/3/2023	50-500PPM
paper:2222	8/1/2024	

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of SGS SOP EQA044.

Supervisor Signature: \_\_\_\_\_

Date: \_\_\_\_\_

Date Archived: \_\_\_\_\_

Date: \_\_\_\_\_

ALS	Sample ID	Method	File ID	MS Batch, Run ID, Pv,,, Dil.	Vial #	Purge Vol	IS/SURR	pH	Status	Comments
1	ib	MR0091	1R03012	MS67909,V1R0100.5,,,1		5			ok	
2	bfbcc91-20	MR0091	1R03013	MS67909,V1R0100.5,,,1		5			ok	9:38am.; 20uL ABK, AA, EC / 100mL
3	cc91-1	MR0091	1R03014	MS67909,V1R0100.5,,,1		5			ok	1uL ABK, AA, EC / 100mL
4	bs	MR0091	1R03015	MS67909,V1R0100.5,,,1		5			ok	50uL ABK, AA, EC / 100mL
5	ib	MR0091	1R03016	MS67909,V1R0100.5,,,1		5			ok	
6	mb	MR0091	1R03017	MS67909,V1R0100.5,,,1		5			ok	
7	j462679-6	MR0091	1R03018	MS67909,V1R0100.5,,,1	2	5		1	ok	
8	j462679-2	MR0091	1R03019	MS67909,V1R0100.5,,,5	4	10.0/50		1	ok	
9	j462679-5	MR0091	1R03020	MS67909,V1R0100.5,,100	2	0.5/50		1	ok	combine with 1r02953
10	j462679-4	MR0091	1R03021	MS67909,V1R0100.5,,,4	3	12.5/50		1	ok	combine with 1r02946
11	j462679-2	MR0091	1R03022	MS67909,V1R0100.5,,50	4	1.0/50		1	ok	
12	j463115-1	MR0091	1R03023	MS67909,V1R0100.5,,,1	4	5		1	ok	
13	j463115-2	MR0091	1R03024	MS67909,V1R0100.5,,,1	4	5		1	ok	
14	j463115-3	MR0091	1R03025	MS67909,V1R0100.5,,,1	4	5		1	ok	
15	j463115-4	MR0091	1R03026	MS67909,V1R0100.5,,,1	5	5		1	ok	
16	j462888-1	MR0091	1R03027	MS67909,V1R0100.5,,,1	8	5		1	ok	
17	j462853-1ms	MR0091	1R03028	MS67909,V1R0100.5,,,4	2	12.5/50		1	ok	25uL ABK, AA, EC / 50mL
18	j462853-1msd	MR0091	1R03029	MS67909,V1R0100.5,,,4	2	12.5/50		1	ok	25uL ABK, AA, EC / 50mL
19	ib	MR0091	1R03030	MS67909,V1R0100.5,,,1	5				ok	
20	j462888-4	MR0091	1R03031	MS67909,V1R0100.5,,,1	7	5		1	ok	
21	j462888-5	MR0091	1R03032	MS67909,V1R0100.5,,,1	1	5		1	ok	







VOLATILE BY GCMS ANALYSIS LOG

Date: 04/04/23

Standard date

Storage Details		
Lot#	Expiration Date	concentration
V023-2768-34	4/26/2023	(0)PPM
V023-2768-9	4/10/2023	100PPM
V023-2768-4	4/24/2023	100-10000PPM

internal/Surrogate data			
Lot#	Expiration Date	concentration	
W023-2768-	5/3/2023	50-500PPM	
paper:2222	8/1/2024		

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified in conomy with the criteria of SGS SOP FD044

Customer Signature \_\_\_\_\_ Date \_\_\_\_\_

110





VOLATILE BY GCMS ANALYSIS LOG

Date: 04/04/23

- 11 -

Storage Conditions			
Lot#	Expiration Date	concentration	
V023-2768-8	4/28/2023	100PPM	
V023-2768-9	4/10/2023	100PPM	
V023-2768-1	4/24/2023	100-1000PPM	

Internal/Surrogate data			
Lot#	Expiration Date	concentration	
V0023-2768-	5/3/2023	5G-500PPM	
paper:2222	8/1/2024		

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of SGS SOP FOA.

Consumer Signature

8

Date: \_\_\_\_\_

Date Archived:

Method: v8260d

Print Analyst Name: nickw

Batch ID: v1r100





VOLATILE by GCMS ANALYSIS LOG

Date: 04/04/23

## Standard data

Lot#	Expiration Date	concentration
V023-2768-94	4/28/2023	100PPM
V023-2768-93	4/10/2023	100PPM
V023-2768-93	4/24/2023	100-1000PPM

internal/surrogate data		
Lot#	Expiration Date	concentration
W0023-2768-	5/3/2023	50-500PPM
paper2222	8/1/2024	

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of SGS SOP EQA044.

Supervisor Signature:

Date:

Date Archived:

Method: v8260d  
Cal. Method: m1R91

Analyst Signature: Meichen

Batch ID: v1r100





VOLATILE BY GCMS ANALYSIS LOG

Date: 04/04/23

Standards

Storage Details		
Lot#	Expiration Date	concentration
V023-2768-34	4/26/2023	(0)PPM
V023-2768-9	4/10/2023	100PPM
V023-2768-4	4/24/2023	100-10000PPM

Internal/Surrogate data

Lot#	Expiration Date	concentration
V023-2768-1	5/3/2023	50-500PPM
paper-2222	8/1/2024	

ANALYTICAL INVESTIGATIONS AND OBSERVATIONS ON THE FOLLOWING ELEMENTS WHICH HAVE BEEN EXAMINED AND CERTIFIED BY ANALYSTS OF THE FORENSIC LABORATORY.

Constitutive Signatures

Date-

Date Archived:

Print Analyst Name: nickw

Print Analyst Name: nickw

Analyst Signature: MeiCh

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Initial Cal Method: m1B91

Date Archived:

Data





Date: 3/29/23

## Standard data

Standard data			
Lot#	Expiration Date	Date	concentration
ABK v023-2768-81.12	4/24/2023	100-10000ppm	
EC v023-2768-82.6	4/3/2023	100 ppm	
AA v023-2768-85.20	4/28/2023	100 ppm	
EXT EC v023-2768-89.8	4/5/2023	100ppm	
Internal/Surrogate data			
Lot#	Expiration Date	Date	concentration
EXT Chlorodifluoromethane v023-2768-64	4/14/2023	100ppm	
EXT ABK v023-2768-80.15	4/24/2023	100-10000ppm	
EXT EC v023-2768-89.6	4/5/2023	100ppm	
EXT Acro v023-2768-86.30	4/28/2023	100ppm	
EXT PA v023-2768-73	4/18/2023	100ppm	
IS v023-2768-55	4/3/2023	50/500ppm	

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of SGS SOP EQA044.

Signature: Blalley

Supervisor Signature-

ALS	Sample ID	Method	File ID	MS Batch, Run ID, P#,,,	Di	Vial #	Purge Vol	IS/SURR	pH	Status	Comments
	ib		1R2811				5ml			ok	
	ib		2R2812				5ml			ok	
bfb			1R2813				5ml			ok	
bfb			2R2814				5ml			ok	
ic91-0,2			1R2815				5ml			ok	2 uL ABK EC AA / 1000 mL
ic91-0,2			2R2816				5ml			ok	2 uL ABK EC AA / 1000 mL
ic91-0,5			1R2817				5ml			ok	5 uL ABK EC AA / 1000 mL
ic91-0,5			2R2818				5ml			ok	5 uL ABK EC AA / 1000 mL
ic91-1			1R2819				5ml			ok	1 uL ABK EC AA / 100 mL
ic91-1			2R2820				5ml			ok	1 uL ABK EC AA / 100 mL
ic91-2			1R2821				5ml			ok	2 uL ABK EC AA / 100 mL
ic91-2			2R2822				5ml			ok	2 uL ABK EC AA / 100 mL
ic91-4			1R2823				5ml			ok	4 uL ABK EC AA / 100 mL
ic91-4			2R2824				5ml			ok	4 uL ABK EC AA / 100 mL
ic91-8			1R2825				5ml			ok	8 uL ABK EC AA / 100 mL
ic91-8			2R2826				5ml			ok	8 uL ABK EC AA / 100 mL



## VOLATILE by GCMS ANALYSIS LOG

Date: 3/29/23

## Standard data

Lot#		Expiration Date	concentration
ABK v023-2767-81.12		4/24/2023	100-10000ppm
EC v023-2768-82.6		4/3/2023	100 ppm
AA v023-2768-85.20		4/28/2023	100 ppm
EXT EC v023-2768-89.8		4/5/2023	100ppm
	internal/Surrogate data		
Lot#		Expiration Date	concentration
EXT Chlорodifluoromethane v023-2768-64		4/14/2023	100ppm
EXT ABK v023-2768-80.15		4/24/2023	100-10000ppm
EXT EC v023-2768-89.6		4/5/2023	100ppm
EXT Acro v023-2768-86.30		4/28/2023	100ppm
EXT PA v023-2768-73		4/18/2023	100ppm
IS v023-2768-55		4/3/2023	50/50ppm

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of SGS SOP EQA044.

Supervisor Signature: \_\_\_\_\_

Date: \_\_\_\_\_

ALS	Sample ID	Method	File ID	MS Patch, Run ID, Pv,,, Dil	Vial #	Purge Vol	IS/SURR	pH	Status	Comments
ic91-20			1R2827			5ml			ok	20 ul ABK EC AA / 100 mL
ic91-20			2R2828			5ml			ok	20 ul ABK EC AA / 100 mL
ic91-50			1R2829			5ml			ok	50 ul ABK EC AA / 100 mL
ic91-50			2R2830			5ml			ok	50 ul ABK EC AA / 100 mL
ic91-100			1R2831			5ml			ok	100 ul ABK EC AA / 100 mL
ic91-100			2R2832			5ml			ok	100 ul ABK EC AA / 100 mL
ic91-200			1R2833			5ml			ok	200 ul ABK EC AA / 100 mL
ic91-200			2R2834			5ml			ok	200 ul ABK EC AA / 100 mL
ib			1R2835			5ml			ok	
ib			2R2836			5ml			ok	
ib			1R2837			5ml			ok	
ib			2R2838			5ml			ok	
icv91-50			1R2839			5ml			ok	50 ul EXT ABK EC Acro / 100 mL
icv91-50			2R2840			5ml			ok	50 ul EXT ABK EC Acro / 100 mL
icv91-50			1R2841			5ml			ok	500 ul EXT PA Chlorodifluoromethane / 100 mL
icv91-50			2R2842			5ml			ok	500 ul EXT PA Chlorodifluoromethane / 100 mL



VOLATILE BY GC/MS ANALYSIS LOG

Date: 3/29/23

### Standard data

Lot#		Expiration Date	concentration
Internal/Surrogate data			
ABK v023-2768-81.12		4/24/2023	100-10000ppm
EC v023-2768-82.6		4/3/2023	100 ppm
AA v023-2768-83.20		4/28/2023	100 ppm
EXT EC v023-2768-89.8		4/5/2023	100ppm
Lot#		Expiration Date	concentration
Internal/Surrogate data			
EXT Chlorodifluoromethane v023-2768-64		4/14/2023	100ppm
EXT TABK v023-2768-80.15		4/24/2023	100-10000ppm
EXT EC v023-2768-89.6		4/5/2023	100ppm
EXT Acro v023-2768-96.30		4/28/2023	100ppm
EXT PA v023-2768-73		4/18/2023	100ppm
IS v023-2768-55		4/3/2023	50/5000ppm

Manually integrated chromatographic peaks in the following recordable files have been reviewed and verified as

Supervisor Signature:

Date-

Batch ID: V1R91

Print Analyst Name: Brashant S

Processed By: Mohini H

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Mathcad: 18360D

Method: Mand

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**General Chemistry****QC Data Summaries**

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**Includes the following where applicable:**

- Method Blank and Blank Spike Summaries
- Duplicate Summaries
- Matrix Spike Summaries
- Instrument Runlogs/QC

METHOD BLANK AND SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRYLogin Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chloride	GP45907/GN39825	2.0	0.0	mg/l	80	83.7	104.6	90-110%
Dissolved Organic Carbon	GP45881/GN39916	1.0	0.0	mg/l	10	9.59	95.9	90-110%
Fluoride	GP45907/GN39825	0.20	0.0	mg/l	2	2.07	103.5	90-110%
Iron, Ferrous	GN39908	0.20	0.0	mg/l	3.0	2.9	96.7	85-115%
Nitrogen, Nitrate + Nitrite	GP45893/GN39816	0.10	0.0	mg/l	2	2.08	104.0	90-110%
Nitrogen, Nitrite	GN39776	0.010	0.0	mg/l	0.04	0.039	97.5	90-110%
Sulfate	GP45907/GN39825	2.0	0.0	mg/l	80	87.0	108.8	90-110%
Total Organic Carbon	GP45861/GN39817	1.0	0.0	mg/l	10	9.21	92.1	90-110%

## Associated Samples:

Batch GN39776: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GN39908: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45861: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45881: JD62888-1F, JD62888-2F, JD62888-3F, JD62888-4F  
 Batch GP45893: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45907: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 (\*) Outside of QC limits

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chloride	GP45907/GN39825	JD62888-3	mg/l	183	182	0.5	0-20%
Fluoride	GP45907/GN39825	JD62888-3	mg/l	0.32	0.32	0.0	0-20%
Nitrogen, Nitrate + Nitrite	GP45893/GN39816	JD62498-5	mg/l	2.4	2.4	0.0	0-20%
Sulfate	GP45907/GN39825	JD62888-3	mg/l	96.8	99.2	2.4	0-20%

Associated Samples:

Batch GP45893: JD62888-1, JD62888-2, JD62888-3, JD62888-4

Batch GP45907: JD62888-1, JD62888-2, JD62888-3, JD62888-4

(\*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chloride	GP45907/GN39825	JD62888-3	mg/l	183	80	262	98.8	80-120%
Dissolved Organic Carbon	GP45881/GN39916	JD62888-2F	mg/l	0.62	10	9.8	91.8	66-128%
Fluoride	GP45907/GN39825	JD62888-3	mg/l	0.32	2	2.6	114.0	80-120%
Iron, Ferrous	GN39908	JD62888-1	mg/l	0.0	3.0	2.9	96.7	92-113%
Nitrogen, Nitrate + Nitrite	GP45893/GN39816	JD62498-5	mg/l	2.4	1	3.3	45.0N(a)	90-110%
Nitrogen, Nitrite	GN39776	JD62888-1	mg/l	0.0	0.04	0.037	92.5	22-140%
Sulfate	GP45907/GN39825	JD62888-3	mg/l	96.8	80	181	105.3	80-120%
Total Organic Carbon	GP45861/GN39817	JD62808-1	mg/l	5.0	10	14.7	97.0	71-132%

Associated Samples:

Batch GN39776: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GN39908: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45861: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45881: JD62888-1F, JD62888-2F, JD62888-3F, JD62888-4F  
 Batch GP45893: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45907: JD62888-1, JD62888-2, JD62888-3, JD62888-4

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

(a) Spike recovery indicates possible matrix interference.

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MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Dissolved Organic Carbon	GP45881/GN39916	JD62888-2F	mg/l	0.62	10	9.7	1.0	20%
Iron, Ferrous	GN39908	JD62888-1	mg/l	0.0	3.0	2.9	0.0	20%
Nitrogen, Nitrite	GN39776	JD62888-1	mg/l	0.0	0.04	0.037	0.0	20%
Total Organic Carbon	GP45861/GN39817	JD62808-1	mg/l	5.0	10	14.4	2.1	10%

Associated Samples:

Batch GN39776: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GN39908: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45861: JD62888-1, JD62888-2, JD62888-3, JD62888-4  
 Batch GP45881: JD62888-1F, JD62888-2F, JD62888-3F, JD62888-4F

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

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SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: E033023W1.NO32 Date Analyzed: 03/30/23 Methods: EPA 353.2/LACHAT  
Analyst: MM Run ID: GN39816  
Parameters: Nitrogen, Nitrate + Nitrite

Time	Sample Description	Dilution Factor	PS Recov	Comments
14:09	GN39816-STD1	1		STDA
14:10	GN39816-STD2	1		STDB
14:11	GN39816-STD3	1		STDC
14:12	GN39816-STD4	1		STDD
14:13	GN39816-STD5	1		STDE
14:14	GN39816-STD6	1		STDF
14:15	GN39816-STD7	1		STDG
14:16	GN39816-STD8	1		STDH
14:18	GN39816-ICV1	1		
14:20	GN39816-ICB1	1		
14:21	GN39816-CCV1	1		
14:22	GN39816-CCB1	1		
14:29	GN39816-CCV2	1		
14:30	GN39816-CCB2	1		
14:31	GP45893-MB1	1		
14:32	GP45893-B1	1		
14:33	GP45893-S1	1		
14:34	GP45893-S2	1		
14:35	GP45893-D1	1		
14:36	JD62498-5	1		(sample used for QC only; not part of login JD62888)
14:37	ZZZZZZ	1		
14:38	ZZZZZZ	1		
14:39	ZZZZZZ	1		
14:40	ZZZZZZ	1		
14:41	GN39816-CCV3	1		
14:43	GN39816-CCB3	1		
14:44	ZZZZZZ	1		
14:45	ZZZZZZ	1		
14:46	ZZZZZZ	1		
14:47	ZZZZZZ	1		
14:48	ZZZZZZ	1		
14:49	JD62808-1	1		(sample used for QC only; not part of login JD62888)
14:50	ZZZZZZ	1		

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: E033023W1.NO32 Date Analyzed: 03/30/23 Methods: EPA 353.2/LACHAT  
Analyst: MM Run ID: GN39816  
Parameters: Nitrogen, Nitrate + Nitrite

Time	Sample Description	Dilution Factor	PS Recov	Comments
14:51	JD62888-1	1		
14:52	JD62888-2	1		
14:53	JD62888-3	1		
14:54	GN39816-CCV4	1		
14:55	GN39816-CCB4	1		
14:57	JD62888-4	1		
14:58	ZZZZZ	1		
14:59	ZZZZZ	1		
15:00	GP45894-MB1	1		
15:01	GP45894-B1	1		
15:02	GP45894-S1	1		
15:03	GP45894-S2	1		
15:04	GP45894-D1	1		
15:05	JD62596-14	1		(sample used for QC only; not part of login JD62888)
15:06	ZZZZZ	1		
15:07	GN39816-CCV5	1		
15:08	GN39816-CCB5	1		
15:09	ZZZZZ	1		
15:11	JD62515-1	1		(sample used for QC only; not part of login JD62888)
15:12	ZZZZZ	1		
15:13	ZZZZZ	1		
15:14	ZZZZZ	1		
15:15	ZZZZZ	1		
15:16	ZZZZZ	1		
15:17	ZZZZZ	1		
15:18	ZZZZZ	1		
15:19	ZZZZZ	1		
15:20	GN39816-CCV6	1		
15:21	GN39816-CCB6	1		
15:22	ZZZZZ	1		
15:23	ZZZZZ	1		
15:25	ZZZZZ	1		
15:26	ZZZZZ	1		

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: E033023W1.NO32 Date Analyzed: 03/30/23 Methods: EPA 353.2/LACHAT  
Analyst: MM Run ID: GN39816  
Parameters: Nitrogen, Nitrate + Nitrite

Time	Sample Description	Dilution Factor	PS Recov	Comments
15:27	ZZZZZZ	1		
15:28	ZZZZZZ	1		
15:29	JD62888-1	4		
15:30	JD62888-2	2		
15:31	JD62888-3	3		
15:32	ZZZZZZ	3		
15:33	GN39816-CCV7	1		
15:34	GN39816-CCB7	1		
15:35	ZZZZZZ	100		
15:36	ZZZZZZ	200		
15:37	ZZZZZZ	2		
15:39	GP45900-MB1	1		
15:40	GP45900-B1	1		
15:41	GP45900-S1	1		
15:42	GP45900-D1	1		
15:43	JD62928-1	1		(sample used for QC only; not part of login JD62888)
15:44	ZZZZZZ	1		
15:45	ZZZZZZ	1		
15:46	GN39816-CCV8	1		
15:51	GN39816-CCV9	1		
15:53	GN39816-CCB8	1		
15:54	ZZZZZZ	1		
15:55	ZZZZZZ	1		
15:56	ZZZZZZ	1		
15:57	ZZZZZZ	1		
15:58	ZZZZZZ	1		
15:59	ZZZZZZ	1		
16:00	ZZZZZZ	1		
16:03	GN39816-CCV10	1		
16:04	GN39816-CCB9	1		
16:08	GN39816-CCV11	1		
16:09	GN39816-CCB10	1		
16:10	ZZZZZZ	8		

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SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888

Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: E033023W1.NO32

Date Analyzed: 03/30/23

Methods: EPA 353.2/LACHAT

Analyst: MM

Run ID: GN39816

Parameters: Nitrogen, Nitrate + Nitrite

Time	Sample Description	Dilution Factor	PS Recov	Comments
16:13	GN39816-CCV12	1		
16:14	GN39816-CCB11	1		

Refer to raw data for calibration curve and standards.

Instrument QC Summary  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: E033023W1.NO32      Date Analyzed: 03/30/23      Methods: EPA 353.2/LACHAT  
Run ID: GN39816      Units: mg/l

Sample Number	Parameter	Result	RL	IDL/MDL	True Value	% Recov.	QC Limits
GN39816-ICV1	Nitrogen, Nitrate + Nitrite	2.13	0.10	0.090	2	106.5	90-110
GN39816-ICB1	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV1	Nitrogen, Nitrate + Nitrite	2.50	0.10	0.090	2.5	100.0	90-110
GN39816-CCB1	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV2	Nitrogen, Nitrate + Nitrite	2.51	0.10	0.090	2.5	100.4	90-110
GN39816-CCB2	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV3	Nitrogen, Nitrate + Nitrite	2.49	0.10	0.090	2.5	99.6	90-110
GN39816-CCB3	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV4	Nitrogen, Nitrate + Nitrite	2.39	0.10	0.090	2.5	95.6	90-110
GN39816-CCB4	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV5	Nitrogen, Nitrate + Nitrite	2.43	0.10	0.090	2.5	97.2	90-110
GN39816-CCB5	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV6	Nitrogen, Nitrate + Nitrite	2.40	0.10	0.090	2.5	96.0	90-110
GN39816-CCB6	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV7	Nitrogen, Nitrate + Nitrite	2.41	0.10	0.090	2.5	96.4	90-110
GN39816-CCB7	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV8	Nitrogen, Nitrate + Nitrite	2.42	0.10	0.090	2.5	96.8	90-110
GN39816-CCV9	Nitrogen, Nitrate + Nitrite	2.42	0.10	0.090	2.5	96.8	90-110
GN39816-CCB8	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV10	Nitrogen, Nitrate + Nitrite	2.40	0.10	0.090	2.5	96.0	90-110
GN39816-CCB9	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV11	Nitrogen, Nitrate + Nitrite	2.42	0.10	0.090	2.5	96.8	90-110
GN39816-CCB10	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			
GN39816-CCV12	Nitrogen, Nitrate + Nitrite	2.41	0.10	0.090	2.5	96.4	90-110
GN39816-CCB11	Nitrogen, Nitrate + Nitrite	0.090 U	0.10	0.090			

( ! ) Outside of QC limits

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D230330W1.TXT      Date Analyzed: 03/01/23      Methods: SM5310 B-11/14  
Analyst: MB      Run ID: GN39817  
Parameters: Total Organic Carbon

Time	Sample Description	Dilution Factor	PS Recov	Comments
11:50	GN39817-STD1	1		STDA
12:01	GN39817-STD2	1		STDB
12:18	GN39817-STD3	1		STDC
12:33	GN39817-STD4	1		STDD
12:45	GN39817-STD5	1		STDE
12:58	GN39817-STD6	1		STDF
09:28	ZZZZZ	1		
09:44	GN39817-CRI1	1		
09:59	GN39817-HSTD1	1		
10:30	GN39817-ICV1	1		
10:43	GN39817-ICB1	1		
12:03	GN39817-CCV1	1		
12:16	GN39817-CCB1	1		
12:32	ZZZZZ	1		
12:43	GP45861-MB1	1		
12:55	GP45861-B1	1		
14:04	JD62808-1	1		(sample used for QC only; not part of login JD62888)
14:13	GP45861-S1	1		
14:24	GP45861-MSD1	1		
14:35	JD62888-4	1		
14:46	ZZZZZ	1		
14:57	ZZZZZ	1		
15:08	ZZZZZ	1		
15:20	GN39817-CCVA1	1		
15:33	GN39817-CCB2	1		
15:42	ZZZZZ	1		
15:53	JD62888-1	1		
16:04	JD62888-2	1		
16:15	JD62888-3	1		
16:27	GN39817-CCV2	1		
16:40	GN39817-CCB3	1		
16:51	ZZZZZ	1		

Refer to raw data for calibration curve and standards.

Instrument QC Summary  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D230330W1.TXT

Date Analyzed: 03/01/23

Run ID: GN39817

Methods: SM5310 B-11/14

Units: mg/l

Sample Number	Parameter	Result	RL	IDL/MDL	True Value	% Recov.	QC Limits
GN39817-CRI1	Total Organic Carbon	1.10	1.0	0.72	1	110.0	70-130
GN39817-HSTD1	Total Organic Carbon	47.1	1.0	0.72	50	94.2	90-110
GN39817-ICV1	Total Organic Carbon	19.9	1.0	0.72	20	99.5	90-110
GN39817-ICB1	Total Organic Carbon	0.72 U	1.0	0.72			
GN39817-CCV1	Total Organic Carbon	24.1	1.0	0.72	25	96.4	90-110
GN39817-CCB1	Total Organic Carbon	0.809	1.0	0.72			
GN39817-CCVA1	Total Organic Carbon	46.4	1.0	0.72	50	92.8	
GN39817-CCB2	Total Organic Carbon	0.72 U	1.0	0.72			
GN39817-CCV2	Total Organic Carbon	24.1	1.0	0.72	25	96.4	90-110
GN39817-CCB3	Total Organic Carbon	0.72 U	1.0	0.72			

( ! ) Outside of QC limits

8.6

8

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D23033001.TXT  
Analyst: SS  
Parameters: Sulfate

Date Analyzed: 03/20/23  
Run ID: GN39825  
Methods: EPA 300/SW846 9056A

Time	Sample Description	Dilution Factor	PS Recov	Comments
13:03	GN39825-STD1	1		STDA
13:16	GN39825-STD2	1		STDC
13:29	GN39825-STD3	1		STDD
13:42	GN39825-STD4	1		STDE
13:55	GN39825-STD5	1		STDF
14:08	GN39825-STD6	1		STDG
14:21	GN39825-ICV1	1		
14:36	ZZZZZ	1		
14:49	ZZZZZ	1		
15:02	ZZZZZ	1		
15:15	GN39825-CCV1	1		
15:28	GN39825-CCB1	1		
15:41	ZZZZZ	1		
15:54	ZZZZZ	1		
16:07	GP45876-S1	15		
16:20	GP45876-D1	15		
16:33	JD62756-1	15		(sample used for QC only; not part of login JD62888)
16:46	ZZZZZ	5		
16:59	ZZZZZ	3		
17:12	ZZZZZ	10		
17:25	ZZZZZ	16		
17:38	ZZZZZ	30		
17:51	GN39825-CCV2	1		
18:04	GN39825-CCB2	1		
18:17	ZZZZZ	34		
18:31	ZZZZZ	20		
18:44	ZZZZZ	3		
18:57	ZZZZZ	5		
19:10	ZZZZZ	3		
19:23	ZZZZZ	7		
19:36	ZZZZZ	14		
19:49	GP45907-MB1	1		
20:02	GP45907-B1	1		

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D23033001.TXT  
Analyst: SS  
Parameters: Sulfate

Date Analyzed: 03/20/23  
Run ID: GN39825  
Methods: EPA 300/SW846 9056A

Time	Sample Description	Dilution Factor	PS Recov	Comments
20:15	GP45907-S1	1		
20:28	GN39825-CCV3	1		
20:41	GN39825-CCB3	1		
20:54	GP45907-D1	1		
21:07	JD62888-3	1		
21:20	ZZZZZ	1		
21:33	ZZZZZ	1		
21:46	ZZZZZ	1		
21:59	ZZZZZ	1		
22:12	ZZZZZ	1		
22:25	JD62888-1	1		
22:39	JD62888-2	1		
22:51	ZZZZZ	1		
23:04	GN39825-CCV4	1		
23:17	GN39825-CCB4	1		
23:30	JD62888-4	1		
23:43	ZZZZZ	1		
13:46	ZZZZZ	1		
13:59	ZZZZZ	1		
14:12	ZZZZZ	1		
14:25	ZZZZZ	1		
14:38	ZZZZZ	1		
14:51	ZZZZZ	1		
15:04	ZZZZZ	1		
15:17	ZZZZZ	1		
15:30	GN39825-CCV5	1		
15:43	GN39825-CCB5	1		
15:56	ZZZZZ	1		
16:09	GP45909-MB1	1		
16:22	GP45909-B1	1		
16:35	GP45909-S1	1		
16:48	GP45909-D1	1		
17:01	JD62596-14	1		(sample used for QC only; not part of login JD62888)

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D23033001.TXT  
Analyst: SS  
Parameters: Sulfate

Date Analyzed: 03/20/23  
Run ID: GN39825  
Methods: EPA 300/SW846 9056A

Time	Sample Description	Dilution Factor	PS Recov	Comments
17:14	ZZZZZZ	1		
17:27	ZZZZZZ	1		
17:40	ZZZZZZ	1		
17:53	ZZZZZZ	1		
18:06	GN39825-CCV6	1		
18:06	GN39825-CCV7	1		
18:19	GN39825-CCB6	1		
18:19	GN39825-CCB7	1		
18:32	ZZZZZZ	1		
18:45	ZZZZZZ	1		
18:58	ZZZZZZ	1		
19:11	ZZZZZZ	1		
19:24	ZZZZZZ	1		
19:37	ZZZZZZ	1		
19:50	ZZZZZZ	1		
20:04	ZZZZZZ	1		
20:17	ZZZZZZ	1		
20:30	ZZZZZZ	1		
20:43	GN39825-CCV8	1		
20:56	GN39825-CCB8	1		
21:09	ZZZZZZ	1		
21:22	ZZZZZZ	1		
21:35	ZZZZZZ	1		
21:48	ZZZZZZ	1		
22:01	ZZZZZZ	1		
22:14	ZZZZZZ	32		
22:27	GN39825-CCV9	1		
22:40	GN39825-CCB9	1		

Refer to raw data for calibration curve and standards.

Instrument QC Summary  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D23033001.TXT Date Analyzed: 03/20/23 Methods: EPA 300/SW846 9056A  
Run ID: GN39825 Units: mg/l

Sample Number	Parameter	Result	RL	IDL/MDL	True Value	% Recov.	QC Limits
GN39825-ICV1	Sulfate	100	2.0	0.89	100	100.0	90-110
GN39825-CCV1	Sulfate	209	2.0	0.89	200	104.5	90-110
GN39825-CCB1	Sulfate	0.89 U	2.0	0.89			
GN39825-CCV2	Sulfate	210	2.0	0.89	200	105.0	90-110
GN39825-CCB2	Sulfate	0.89 U	2.0	0.89			
GN39825-CCV3	Sulfate	208	2.0	0.89	200	104.0	90-110
GN39825-CCB3	Sulfate	0.89 U	2.0	0.89			
GN39825-CCV4	Sulfate	212	2.0	0.89	200	106.0	90-110
GN39825-CCB4	Sulfate	1.31	2.0	0.89			
GN39825-CCV5	Sulfate	211	2.0	0.89	200	105.5	90-110
GN39825-CCB5	Sulfate	0.89 U	2.0	0.89			
GN39825-CCV6	Sulfate	210	2.0	0.89	200	105.0	90-110
GN39825-CCV7	Sulfate	210	2.0	0.89	200	105.0	90-110
GN39825-CCB6	Sulfate	0.89 U	2.0	0.89			
GN39825-CCB7	Sulfate	0.89 U	2.0	0.89			
GN39825-CCV8	Sulfate	210	2.0	0.89	200	105.0	90-110
GN39825-CCB8	Sulfate	0.89 U	2.0	0.89			
GN39825-CCV9	Sulfate	210	2.0	0.89	200	105.0	90-110
GN39825-CCB9	Sulfate	0.89 U	2.0	0.89			

( ! ) Outside of QC limits

8.7  
8

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D230331W1.TXT Date Analyzed: 03/01/23 Methods: SM5310 B-11, SM5310 B-11/14  
Analyst: MB Run ID: GN39916  
Parameters: Dissolved Organic Carbon

Time	Sample Description	Dilution Factor	PS Recov	Comments
11:50	GN39916-STD1	1		STDA
12:01	GN39916-STD2	1		STDB
12:18	GN39916-STD3	1		STDC
12:33	GN39916-STD4	1		STDD
12:45	GN39916-STD5	1		STDE
12:58	GN39916-STD6	1		STDF
09:28	ZZZZZ	1		
09:44	GN39916-CRI1	1		
09:59	GN39916-HSTD1	1		
10:30	GN39916-ICV1	1		
10:43	GN39916-ICB1	1		
12:08	GN39916-CCV1	1		
12:22	GN39916-CCB1	1		
12:35	ZZZZZ	1		
12:51	GP45856-MB1	1		
15:43	GP45856-B1	1		
17:12	JD62800-1	1		(sample used for QC only; not part of login JD62888)
17:23	GP45856-S1	1		
17:35	GP45856-MSD1	1		
17:46	ZZZZZ	1		
17:57	ZZZZZ	1		
18:08	ZZZZZ	1		
18:20	ZZZZZ	1		
18:31	GN39916-CCVA1	1		
18:43	GN39916-CCB2	1		
18:54	ZZZZZ	1		
19:06	ZZZZZ	1		
19:17	ZZZZZ	1		
19:26	ZZZZZ	1		
19:37	GP45881-MB1	1		
19:49	GP45881-B1	1		
19:59	JD62888-2F	1		
20:11	GP45881-S1	1		

SGS Instrument Runlog  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D230331W1.TXT Date Analyzed: 03/01/23 Methods: SM5310 B-11, SM5310 B-11/14  
Analyst: MB Run ID: GN39916  
Parameters: Dissolved Organic Carbon

Time	Sample Description	Dilution Factor	PS Recov	Comments
20:22	GP45881-MSD1	1		
20:35	JD62888-4F	1		
20:45	GN39916-CCV2	1		
21:06	GN39916-CCB3	1		
21:17	JD62888-1F	1		
21:28	JD62888-3F	1		
21:37	GP45915-MB1	1		
21:49	GP45915-B1	1		
22:02	JD62928-1	1		(sample used for QC only; not part of login JD62888)
22:11	GP45915-S1	1		
22:22	GP45915-MSD1	1		
22:33	ZZZZZ	1		
22:44	ZZZZZ	1		
22:55	ZZZZZ	1		
23:07	GN39916-CCVA2	1		
23:20	GN39916-CCB4	1		
23:29	ZZZZZ	1		
23:40	ZZZZZ	1		
23:51	ZZZZZ	1		
00:03	GN39916-CCV3	1		
00:14	GN39916-CCB5	1		
00:25	ZZZZZ	1		
09:51	GN39916-CCVA3	1		
10:03	GN39916-CCB6	1		
10:13	ZZZZZ	1		
10:24	ZZZZZ	5		
11:16	ZZZZZ	5		
11:27	ZZZZZ	20		
11:39	GN39916-CCV4	1		
12:00	GN39916-CCB7	1		
12:09	ZZZZZ	1		

Refer to raw data for calibration curve and standards.

Instrument QC Summary  
Inorganics Analyses

Login Number: JD62888  
Account: FLSNYNY - Fleming-Lee Shue, Inc.  
Project: 388 Bridge Street, Brooklyn, NY

File ID: D230331W1.TXT Date Analyzed: 03/01/23 Methods: SM5310 B-11, SM5310 B-11/14  
Run ID: GN39916 Units: mg/l

Sample Number	Parameter	Result	RL	IDL/MDL	True Value	% Recov.	QC Limits
GN39916-CRI1	Total Organic Carbon	1.10	1.0	0.72	1	110.0	70-130
GN39916-HSTD1	Total Organic Carbon	47.1	1.0	0.72	50	94.2	90-110
GN39916-ICV1	Total Organic Carbon	19.9	1.0	0.72	20	99.5	90-110
GN39916-ICB1	Total Organic Carbon	0.72 U	1.0	0.72			
GN39916-CCV1	Total Organic Carbon	23.0	1.0	0.72	25	92.0	90-110
GN39916-CCB1	Total Organic Carbon	0.72 U	1.0	0.72			
GN39916-CCVA1	Total Organic Carbon	46.4	1.0	0.72	50	92.8	
GN39916-CCB2	Total Organic Carbon	0.72 U	1.0	0.72			
GN39916-CCV2	Total Organic Carbon	22.9	1.0	0.72	25	91.6	90-110
GN39916-CCB3	Total Organic Carbon	0.72 U	1.0	0.72			
GN39916-CCVA2	Total Organic Carbon	46.5	1.0	0.72	50	93.0	
GN39916-CCB4	Total Organic Carbon	0.72 U	1.0	0.72			
GN39916-CCV3	Total Organic Carbon	23.0	1.0	0.72	25	92.0	90-110
GN39916-CCB5	Total Organic Carbon	0.825	1.0	0.72			
GN39916-CCVA3	Total Organic Carbon	48.0	1.0	0.72	50	96.0	
GN39916-CCB6	Total Organic Carbon	0.72 U	1.0	0.72			
GN39916-CCV4	Total Organic Carbon	24.8	1.0	0.72	25	99.2	90-110
GN39916-CCB7	Total Organic Carbon	0.72 U	1.0	0.72			

( ! ) Outside of QC limits

**General Chemistry****Raw Data**

Test: Nitrogen, Nitrite  
 Product: NO<sub>2</sub>  
 Method: SM4500NC2 B-11 (Aqueous)  Units: mg/l  
 SM4500NO2 B-11 M (Solids)  mg/kg

Analyst: MAHENDRP  
 GN Batch ID: GN39776  
 GP Batch ID:  
 Date: 3/30/2023  
 Instrument ID: L

Pipet II GLASS PIPETTE CLASS A

Original Calibration Information													
Calibration Date: 1/14/2023													
Known:	Blank	Std 1	Std 2	Std 3	Std 4	Std 5	Std 6	Std 7					
Absorbance:	0.000	0.010	0.026	0.050	0.075	0.100	0.200						
Actual Value:	0.000	0.033	0.085	0.170	0.257	0.345	0.694						
% RE:	#DIV/0!	0.010	0.025	0.050	0.075	0.100	0.200						
	-0.83	-0.20	0.97	0.59	0.12	-0.17							
Continuing Calibration Check Standards Data:													
Calibration Date: 3/30/2023													
Known:	Blank	Std 1	Std 6			Correlation Coeff. =	0.99998						
Absorbance:	0.000	0.010	0.200			Slope =	0.28782						
Recovery:	0.000	0.032	0.688			Intercept =	0.00058						
Actual Value:	0.000	97.9%	99.3%										
0.010	0.199												
Bottle #	Sample ID	Time Analyzed	Initial Wt (g) or Vol (ml)	Final Vol (ml)	Dilution	Sample Abs	Background Abs.	Result From Curve (mg/L)	Final Result	DL	Units	Factor	pH between 5 and 9 (Y or N)
	ICV	8:10	50	50	1	0.326	NA	0.094	0.094	NA	mg/l	NA	
	CCV	8:36	50	50	1	0.328	NA	0.095	0.095	NA	mg/l	NA	
	CCB	8:36	50	50	1	0.000	NA	0.001	0.001	NA	mg/l	NA	
	GN39776-MB1	8:36	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	
	GN39776-B1	8:36	50	50	1	0.132	0.000	0.039	0.039	0.010	mg/l	1	
4	GN39776-S1	8:36	50	50	1	0.128	0.000	0.037	0.037	0.010	mg/l	1	
4	GN39776-MSD1	8:36	50	50	1	0.128	0.000	0.037	0.037	0.010	mg/l	1	
4	JD62888-1	8:36	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	
4	JD62888-2	8:36	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
4	JD62888-3	8:36	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
4	JD62888-4	8:36	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
6	JD62896-1	8:36	50	50	1	0.050	0.048	0.001	0.001	0.010	mg/l	1	Y
6	JD62897-1	8:36	50	50	1	0.088	0.035	0.016	0.016	0.010	mg/l	1	Y
	CCVA	8:36	50	50	1	0.685	0.000	0.198	0.198	NA	mg/l	NA	
	CCB	8:36	50	50	1	0.000	NA	0.001	0.001	NA	mg/l	NA	
8	JD62926-1	12:56	50	50	1	0.068	0.000	0.020	0.020	0.010	mg/l	1	Y
9	JD62928-2	13:10	50	50	1	0.005	0.000	0.002	0.002	0.010	mg/l	1	Y
9	JD62928-3	13:10	50	50	1	0.006	0.000	0.002	0.002	0.010	mg/l	1	Y
9	JD62928-4	13:10	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
9	JD62929-1	13:10	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
9	JD62929-2	13:10	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
9	JD62929-3	13:10	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
11	JD62934-1	14:15	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
3	JD62934-1	14:17	50	50	1	0.000	0.000	0.001	0.001	0.010	mg/l	1	Y
		14:20	50	50	1	0.000	0.000	0.001	0.001				
	CCV	14:20	50	50	1	0.330	NA	0.066	0.096	NA	mg/l	NA	
	CCB	14:20	50	50	1	0.000	NA	0.001	0.001	NA	mg/l	NA	
	CCVA		50	50	1		NA			NA	mg/l	NA	
	CCB		50	50	1		NA			NA	mg/l	NA	

Analyst: MAR Date: 3/30/23 QC Reviewer: \_\_\_\_\_ Date: \_\_\_\_\_3/30/23



Test: Nitrogen, Nitrite

Product: NO2

Method: SM18 4500 NO2B (aqueous)  
SM18 4500 NO2B M (solids)Units: mg/l  
mg/kgAnalyst: MPW  
GNBatch ID: 121397+6  
GPBatch ID:  
Date: 3/30/23**Preparation Batch QC Summary**

Units = \_\_\_\_\_

Method Blank ID: (GN39776-MB) Date: 3/30/23 Result: <DL DL: 0.01 <DL: 0.01  
 Spike Blank ID: -01 Date: 1 Result: 0.039 Spike: 0.04 %Rec: 97.5  
 MS ID: +51 Samp. Result: 0.032 MS Result: 0.037 Spike: 0.04 %Rec: 92.5  
 MSDUP ID: -MSD1 Samp. Result: 0.030 MSDUP Result: 0.032 Spike: 0.04 %Rec: 92.5  
 MSDUP ID: JD62885-1 MSDUP Result: 0.032 MS Result: 0.037 %RPD: 0.00

Method Blank ID: \_\_\_\_\_ Date: \_\_\_\_\_ Result: \_\_\_\_\_ DL: \_\_\_\_\_ <DL: \_\_\_\_\_  
 Spike Blank ID: \_\_\_\_\_ Date: \_\_\_\_\_ Result: \_\_\_\_\_ Spike: 0.04 %Rec: 97.5  
 MS ID: \_\_\_\_\_ Samp. Result: \_\_\_\_\_ MS Result: \_\_\_\_\_ Spike: 0.04 %Rec: 92.5  
 MSDUP ID: \_\_\_\_\_ Samp. Result: \_\_\_\_\_ MSDUP Result: \_\_\_\_\_ Spike: 0.04 %Rec: 92.5  
 MSDUP ID: \_\_\_\_\_ MSDUP Result: \_\_\_\_\_ MS Result: \_\_\_\_\_ %RPD: 0.00

**Analysis Batch QC Summary**

Units = \_\_\_\_\_

ICV (Ext): 3/30/23 Result: 0.094 TV: 0.1 %Rec: 94

CCV: 3/30/23 Result: 0.095 TV: 0.1 %Rec: 95  
 CCVA: \_\_\_\_\_ Result: 0.198 TV: 0.2 %Rec: 99.5  
 CCV: \_\_\_\_\_ Result: 0.096 TV: 0.1 %Rec: 96  
 CCVA: \_\_\_\_\_ Result: 0.196 TV: 0.2 %Rec: 98  
 CCV: \_\_\_\_\_ Result: 0.097 TV: 0.1 %Rec: 97  
 CCVA: \_\_\_\_\_ Result: 0.197 TV: 0.2 %Rec: 99

CCB: \_\_\_\_\_ Result: \_\_\_\_\_ DL: \_\_\_\_\_ <DL: \_\_\_\_\_  
 CCB: \_\_\_\_\_ Result: \_\_\_\_\_ DL: \_\_\_\_\_ <DL: \_\_\_\_\_

**Reagent Reference Numbers:**

QC ---- BS/MS/MSD ---- 2 ML OF 1 PPM NO2 STD -- 50 ML DI/H2O/SAMPLE ---- TV 0.04 MG/L

MS/MSD SAMPLE NO --- QC1 = JD62885-1

QC2 = \_\_\_\_\_

Analyst: MPW Date: 3/30/23

Comments: \_\_\_\_\_

on 39778



## Reagent Information Log - Nitrite as Nitrogen

All standards and stocks were made as described in the SOP for this method (circle one): Y or N  
If no (N), see attached page for standards prep.

Form: GN087A-42  
Rev. Date: 3/30/2023



# GENERAL CHEMISTRY STANDARD PREPARATION LOG

Product: 1402  
GN or GP Number: Opn 3977

Standard	Manufacturer/Lot Number	Concentration	Expiration Date
1000 mg/l Nitrite Standard	ERA LOT# 261221M	1000 PPM	12/14/2023
1000 mg/l Nitrite Standard (ICV)	ERA LOT# 120421M	1000 PPM	4/7/2023
Standard Description	Lot or Tracking #	Initial Volume	Autopipet ID
10 PPM NO2 STANDARD	GN-3/30/2023	1.0 ml	1000 mg/l Nitrite Standard A
1 PPM NO2 STANDARD	GN-3/30/2023	10.0 ml	10 PPM NO2 STANDARD A
10 PPM NO2 EXTERNAL	GN-3/30/2023	1.0 ml	Nitrite Standard (ICV) A
1PPM NO2 EXTERNAL	GN-3/30/2023	10.0 ml	10 PPM NO2 EXTERNAL A
Standard Description	Intermediate or Stock used to prepare standard	Initial Volume	Autopipet ID
0.2 mg/L (high)	GN-3/30/2023	20.0 ml	1.0 mg/L
0.1 mg/L		10.0 ml	1.0 mg/L
0.075 mg/L		7.5 ml	1.0 mg/L
0.05 mg/L		5.0 ml	1.0 mg/L
0.025 mg/L		2.5 ml	1.0 mg/L
0.01 mg/L (low)	GN-3/30/2023	1.0 ml	1.0 mg/L
0 mg/L	NA	NA	NA
CCV	GN-3/30/2023	10.0 ml	1.0 mg/L
CCVA	GN-3/30/2023	20.0 ml	1.0 mg/L
ICV	GN-3/30/2023	10.0 ml	1.0 mg/L

\* If Class A glass pipets are used, enter an A. For balances or autopipets, then enter the appropriate Accutest ID number.

Form: GN121-01  
Rev Date: 1/13/09

LABORATORY REVIEW SIGNATURE FORM  
(To be stored with the raw data)File ID: E033023W1.NO32  
Analyst: MMDate Analyzed: 03/30/23  
Run ID: GN39816  
Methods: EPA 353.2/LACHAT

The following analyst(s) have reviewed this run and attest that, to the best of their knowledge, this documentation is complete and correct:

Analyst: MM Date 03/30/23

Analyst: \_\_\_\_\_ Date \_\_\_\_\_

The following supervisor or their designee has reviewed this run and attests that, to the best of their knowledge, this documentation is complete and correct:

Supervisor (or designee): C Date 3/31/239.2  
9

GN39816

Author: Chemistry

Date : 3/30/2023

Original Run Filename: OM\_3-30-2023\_02-08-21PM.OMN Created: 3/30/2023 2:08:21 PM

Original Run Author's Signature: [Chemistry]

Current Run Filename: OM\_3-30-2023\_02-08-21PM.OMN Last Modified: 3/30/2023 2:21:19 PM

Current Run Author's Signature: [Chemistry]

Description: Default New Run

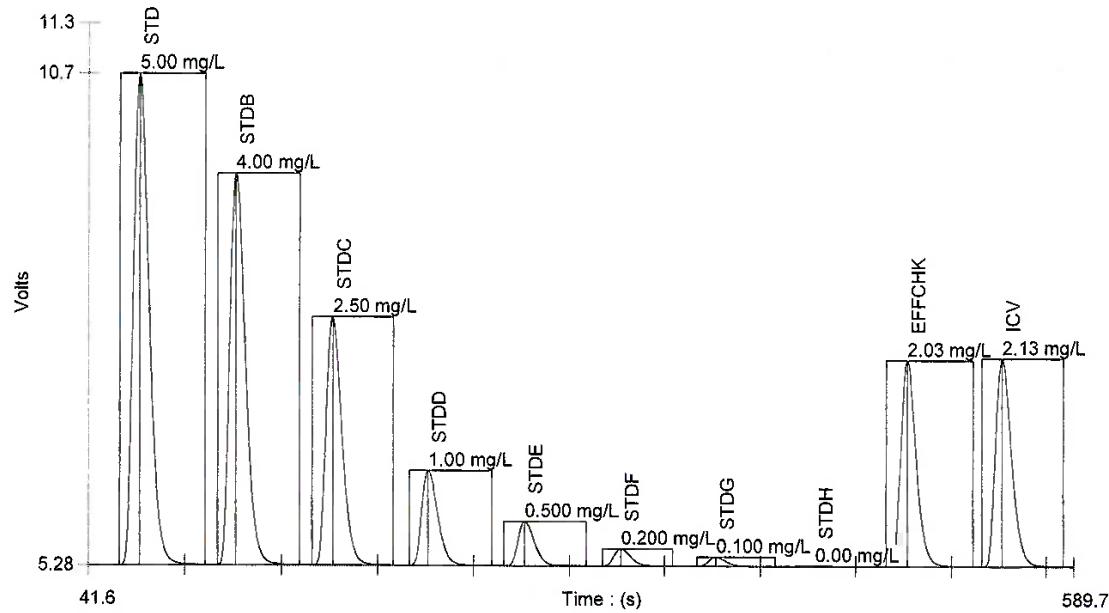
e033023w1.no32

Sample	Rep.	Cup No.	Channel 1	Detection Time	MDF
			NO32 (mg/L)		
STDA	1	1	5.00	3/30/2023@2:09:14 PM	
STDB	1	2	4.00	3/30/2023@2:10:19 PM	
STDC	1	3	2.50	3/30/2023@2:11:25 PM	
STDH	1	4	1.00	3/30/2023@2:12:30 PM	
STDE	1	5	0.500	3/30/2023@2:13:35 PM	
STDF	1	6	0.200	3/30/2023@2:14:40 PM	
STDG	1	7	0.100	3/30/2023@2:15:44 PM	
STDH	1	8	0.00	3/30/2023@2:16:49 PM	
EFFCHK	1	9	2.03	3/30/2023@2:17:53 PM	
Known Conc:			2.00		
Calibration: Table/Fig. : 1					
ICV	1	10	2.13	3/30/2023@2:18:57 PM	
Known Conc:			2.00		
ICB	1	11	0.0194	3/30/2023@2:20:01 PM	
Known Conc:			0.00		
CCV	1	12	2.50	3/30/2023@2:21:06 PM	
Known Conc:			2.50		
CCB	1	13	0.0206	3/30/2023@2:22:10 PM	
Known Conc:			0.00		

106.5 %

100%

Channel 1 - Set: 1 / 2



Date : 3/30/2023

Author: Chemistry

Channel 1 - Set: 2 / 2

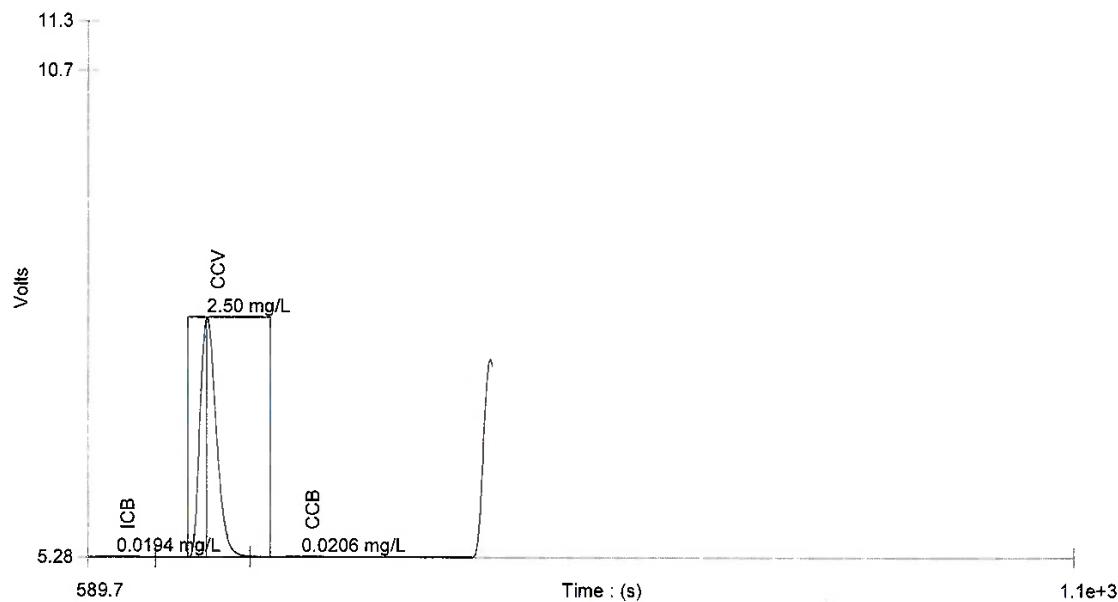
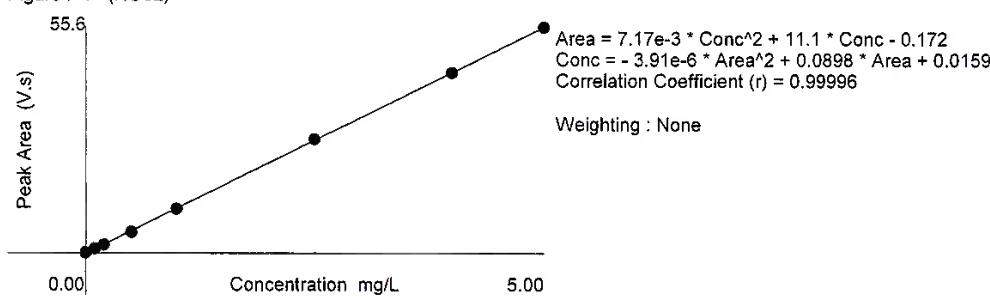


Table : 1 (NO32)

	Known Conc. (mg/L)	Rep.	Peak Area (V.s)	Peak Height (V)	% RSD	% Residual	Det. Conc (mg/L)	Detection Date	Detection Time
1	5.00	1	55.6	5.42	0.0	7.6e-3	5.00	3/30/2023	2:09:14 PM
2	4.00	1	44.3	4.33	0.0	0.2	3.99	3/30/2023	2:10:19 PM
3	2.50	1	27.9	2.73	0.0	-1.0	2.52	3/30/2023	2:11:25 PM
4	1.00	1	10.8	1.05	0.0	1.6	0.984	3/30/2023	2:12:30 PM
5	0.500	1	5.10	0.489	0.0	5.4	0.474	3/30/2023	2:13:35 PM
6	0.200	1	1.99	0.188	0.0	3.2	0.194	3/30/2023	2:14:40 PM
7	0.100	1	1.07	0.0997	0.0	-13.6	0.112	3/30/2023	2:15:44 PM
8	0.00	1	0.0461	4.82e-3			0.0200	3/30/2023	2:16:49 PM

0.8%  
12.0%

Figure : 1 (NO32)



Author: Chemistry

Date : 3/30/2023

Original Run Filename: OM\_3-30-2023\_02-28-02PM.OMN Created: 3/30/2023 2:28:02 PM

Original Run Author's Signature: [Chemistry]

Current Run Filename: OM\_3-30-2023\_02-28-02PM.OMN Last Modified: 3/30/2023 3:44:53 PM

Current Run Author's Signature: [Chemistry]

Description: Default New Run

e033023w2. no32

Sample	Rep.	Cup No.	Channel 1 NO32 (mg/L)	Detection Time	MDF
CCV	1	12	2.51	3/30/2023@2:29:03 PM	
		Known Conc:	2.50		→ 100.4%
CCB	1	13	0.0199	3/30/2023@2:30:07 PM	
		Known Conc:	0.00		
GP45893-MB1	1	14	0.0190	3/30/2023@2:31:11 PM	
GP45893-B1	1	15	2.08	3/30/2023@2:32:15 PM	
GP45893-S1	1	16	3.29	3/30/2023@2:33:20 PM	
GP45893-S2	1	17	1.34	3/30/2023@2:34:25 PM	
GP45893-D1	1	18	2.43	3/30/2023@2:35:31 PM	
JD62498-5	1	19	2.37	3/30/2023@2:36:36 PM	
JD62498-1	1	20	3.41	3/30/2023@2:37:41 PM	
JD62498-2	1	21	0.0935	3/30/2023@2:38:46 PM	
JD62498-4	1	22	0.991	3/30/2023@2:39:51 PM	
JD62498-6	1	23	0.678	3/30/2023@2:40:55 PM	
CCV	1	12	2.49	3/30/2023@2:41:59 PM	
		Known Conc:	2.50		→ 99.6%
CCB	1	13	0.0205	3/30/2023@2:43:03 PM	
		Known Conc:	0.00		
JD62498-7	1	24	1.81	3/30/2023@2:44:08 PM	
JD62498-8	1	25	0.137	3/30/2023@2:45:13 PM	
JD62498-9	1	26	0.781	3/30/2023@2:46:17 PM	
JD62498-10	1	27	0.812	3/30/2023@2:47:21 PM	
JD62498-11	1	28	1.09	3/30/2023@2:48:25 PM	
JD62808-1	1	29	0.447	3/30/2023@2:49:29 PM	
JD62808-2	1	30	0.226	3/30/2023@2:50:33 PM	
JD62888-1	1	31	8.22	3/30/2023@2:51:39 PM	
JD62888-2	1	32	5.44	3/30/2023@2:52:44 PM	
JD62888-3	1	33	7.04	3/30/2023@2:53:49 PM	
CCV	1	12	2.39	3/30/2023@2:54:53 PM	
		Known Conc:	2.50		→ 95.6
CCB	1	13	0.0205	3/30/2023@2:55:57 PM	
		Known Conc:	0.00		
JD62888-4	1	34	0.0233	3/30/2023@2:57:03 PM	
JD62896-1	1	35	4.50	3/30/2023@2:58:08 PM	
JD62897-1	1	36	6.58	3/30/2023@2:59:13 PM	
GP45894-MB1	1	37	0.0195	3/30/2023@3:00:17 PM	
GP45894-B1	1	38	2.00	3/30/2023@3:01:22 PM	
GP45894-S1	1	39	1.23	3/30/2023@3:02:27 PM	
GP45894-S2	1	40	2.66	3/30/2023@3:03:31 PM	
GP45894-D1	1	41	0.324	3/30/2023@3:04:35 PM	
JD62596-14	1	42	0.312	3/30/2023@3:05:40 PM	
JD62596-7	1	43	2.92	3/30/2023@3:06:44 PM	
CCV	1	12	2.43	3/30/2023@3:07:48 PM	
		Known Conc:	2.50		→ 97.2%
CCB	1	13	0.0206	3/30/2023@3:08:52 PM	
		Known Conc:	0.00		
JD62596-8	1	44	2.94	3/30/2023@3:09:56 PM	
JD62515-1	1	45	1.73	3/30/2023@3:11:00 PM	
JD62515-2	1	46	1.43	3/30/2023@3:12:06 PM	
JD62515-3	1	47	0.996	3/30/2023@3:13:12 PM	
JD62515-4	1	48	1.72	3/30/2023@3:14:17 PM	
JD62758-1	1	49	0.0850	3/30/2023@3:15:22 PM	
JD62758-4	1	50	1.63	3/30/2023@3:16:27 PM	
JD62758-5	1	51	1.61	3/30/2023@3:17:32 PM	
JD62758-7	1	52	1.60	3/30/2023@3:18:36 PM	
JD62758-8	1	53	0.437	3/30/2023@3:19:41 PM	
CCV	1	12	2.40	3/30/2023@3:20:45 PM	
		Known Conc:	2.50		→ 96%
CCB	1	13	0.0178	3/30/2023@3:21:49 PM	
		Known Conc:	0.00		

- 1 -

9.2  
9

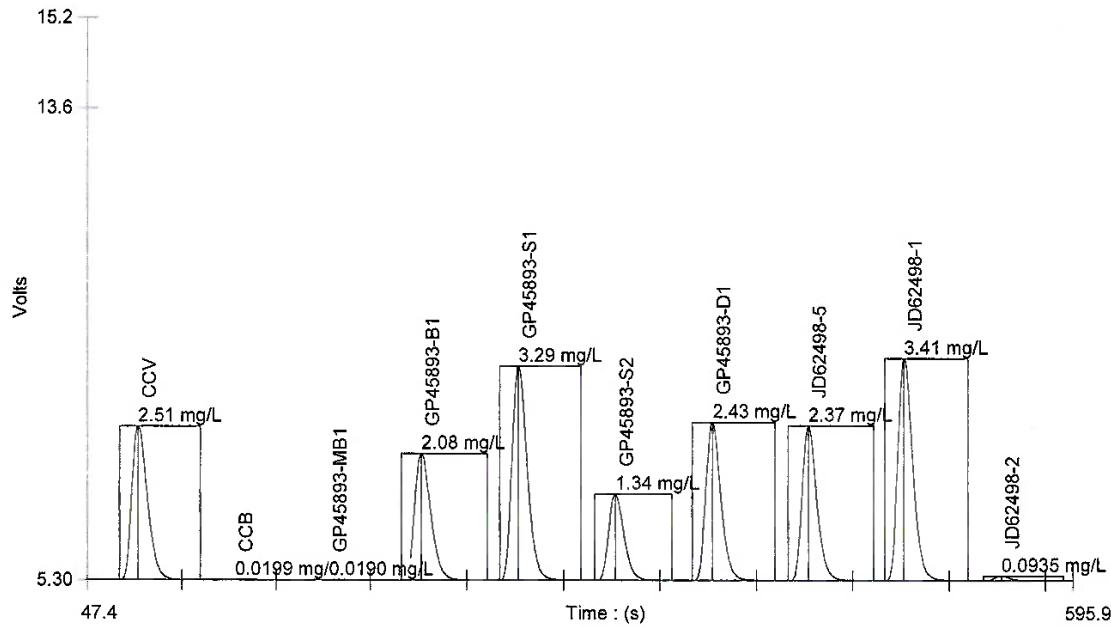
Author: Chemistry

Date : 3/30/2023

JD62758-9	1	54	0.0298	3/30/2023@3:22:54 PM
JD62758-10	1	55	0.0240	3/30/2023@3:23:58 PM
JD62758-11	1	56	0.0285	3/30/2023@3:25:02 PM
JD62758-12	1	57	0.281	3/30/2023@3:26:06 PM
JD62789-1	1	58	17.0	3/30/2023@3:27:10 PM
LA89062-1	1	63	4.83	3/30/2023@3:28:16 PM
JD62888-1	1	59	2.29	3/30/2023@3:29:20 PM
JD62888-2	1	60	2.75	3/30/2023@3:30:24 PM
JD62888-3	1	61	2.49	3/30/2023@3:31:29 PM
JD62897-1	1	62	2.37	3/30/2023@3:32:34 PM
CCV	1	12	2.41	3/30/2023@3:33:38 PM
Known Conc:		2.50		
Calibration:		Table/Fig. : 1		
CCB	1	13	0.0173	3/30/2023@3:34:42 PM
Known Conc:		0.00		
JD62789-1	1	78	3.25	3/30/2023@3:35:47 PM
JD62789-1	1	79	1.74	3/30/2023@3:36:52 PM
LA89062-1	1	80	1.44	3/30/2023@3:37:57 PM
GP45900-MB1	1	64	0.0200	3/30/2023@3:39:01 PM
GP45900-B1	1	65	2.00	3/30/2023@3:40:07 PM
GP45900-S1	1	66	1.95	3/30/2023@3:41:11 PM
GP45900-D1	1	67	0.997	3/30/2023@3:42:16 PM
JD62928-1	1	68	0.978	3/30/2023@3:43:21 PM
JD62928-2	1	69	0.432	3/30/2023@3:44:26 PM
JD62928-3	1	70	0.0653	3/30/2023@3:45:30 PM
CCV	1	12	2.42	3/30/2023@3:46:34 PM
Known Conc:		2.50		

- 2 ml Sample + 6 mL DI  
 - 4 mL Sample + 4 mL DI  
 - 3 mL Sample + 6 mL DI  
 - 96.4%  
 - not needed = 1 mL + 99 mL DI  
 - 4 mL (1:100) + 4 mL DI  
 - 3 mL Sample + 3 mL DI  
 - 100%  
 - 96.8%

Channel 1 - Set: 1 / 8

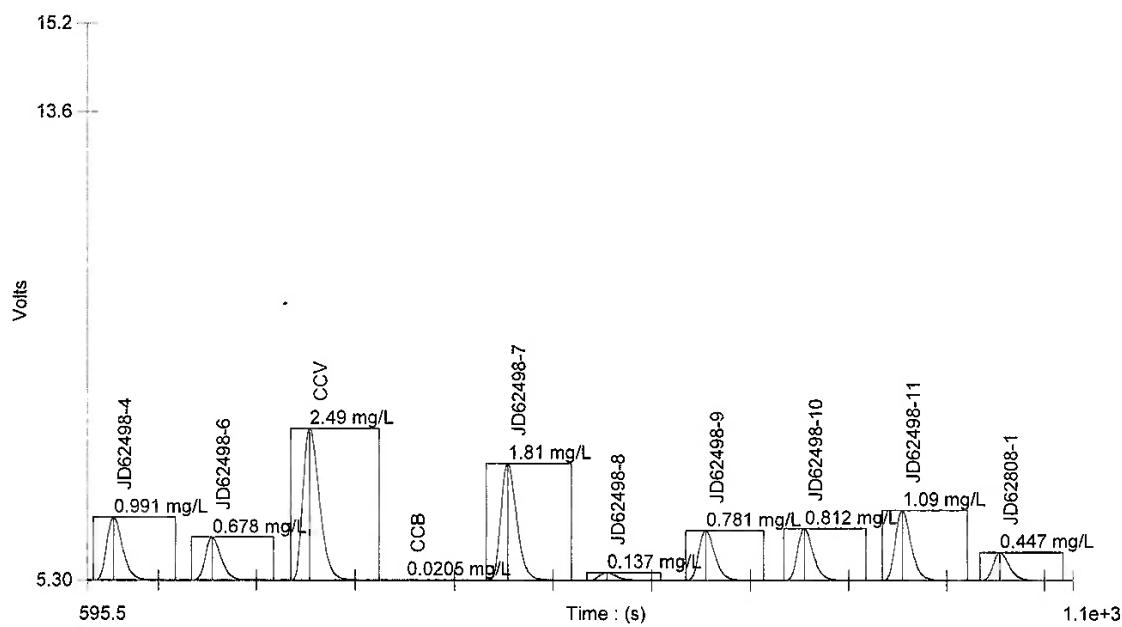


Date : 3/30/2023

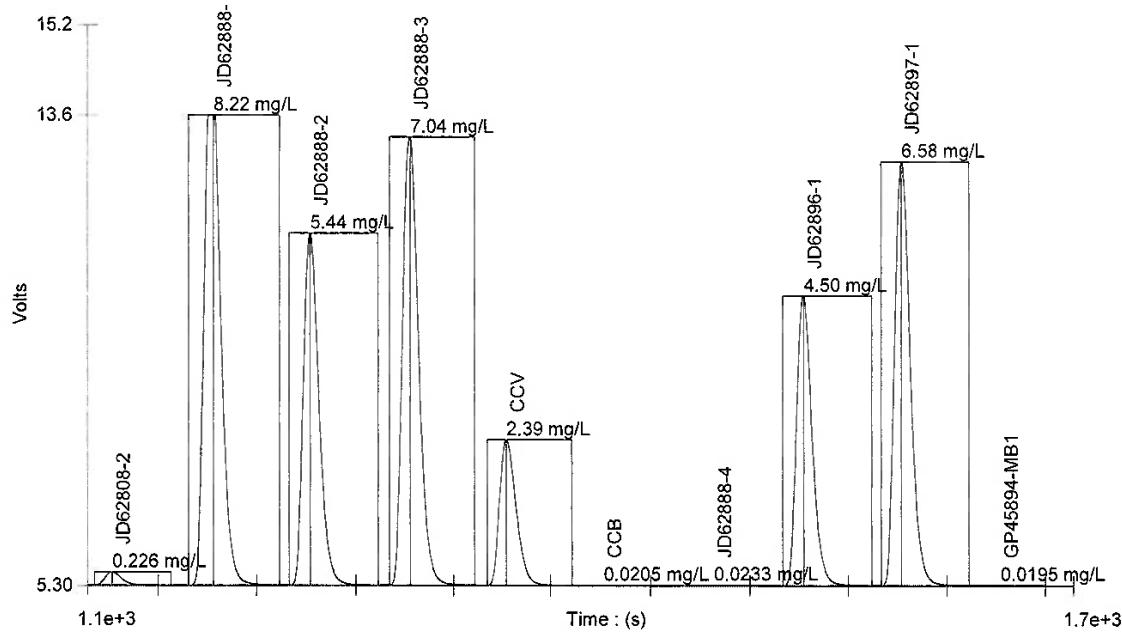
9  
2  
9

Author: Chemistry

Channel 1 - Set: 2 / 8



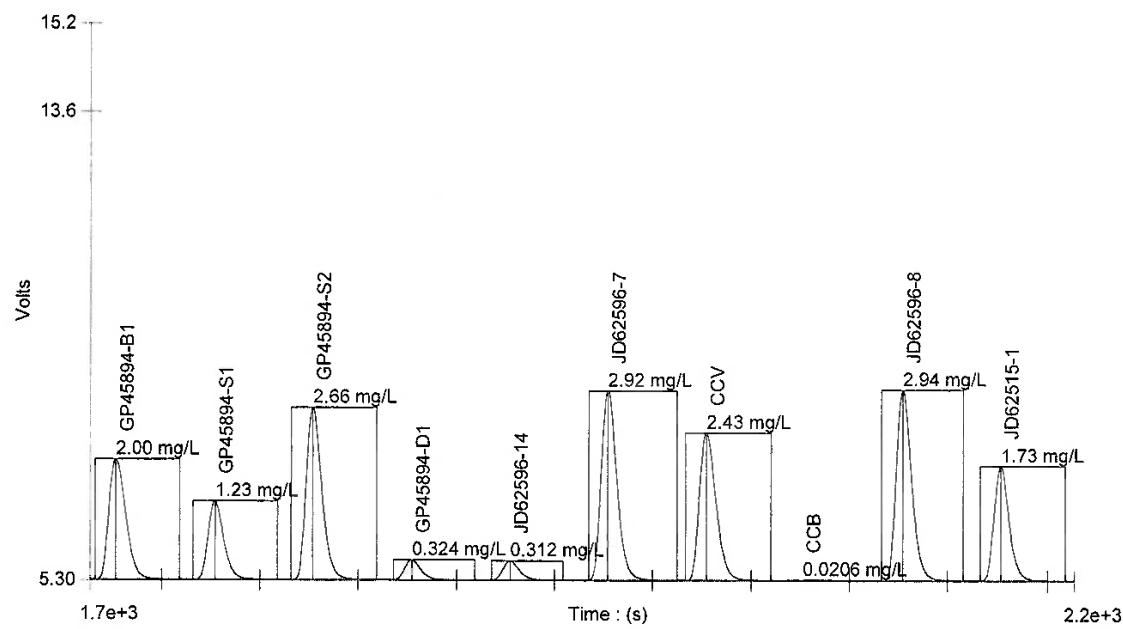
Channel 1 - Set: 3 / 8



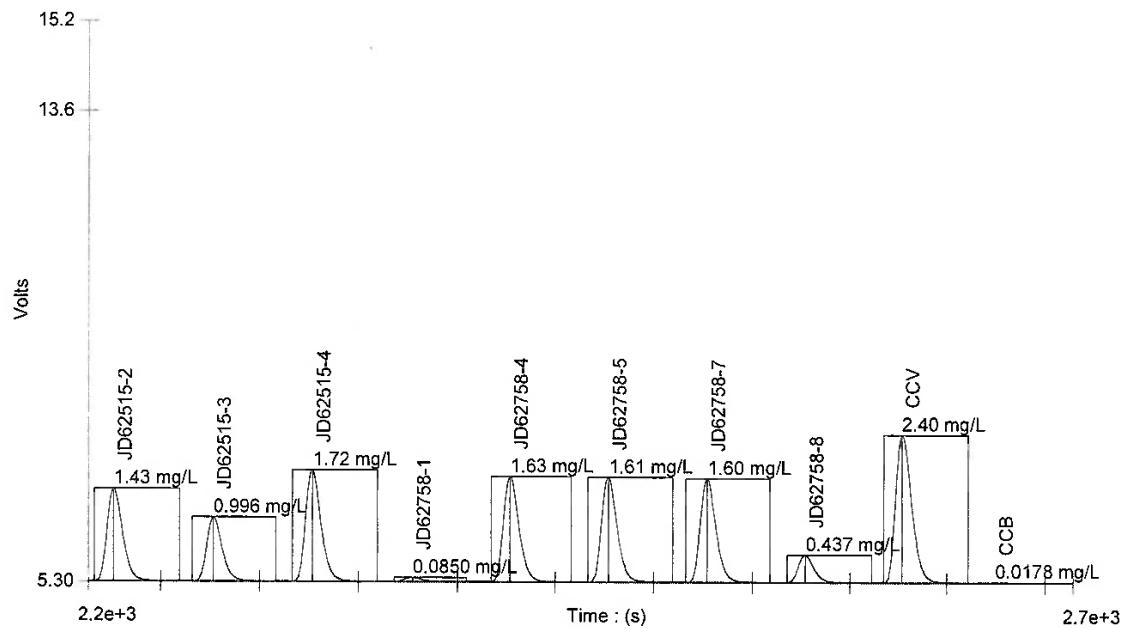
Date : 3/30/2023

Author: Chemistry

Channel 1 - Set: 4 / 8



Channel 1 - Set: 5 / 8

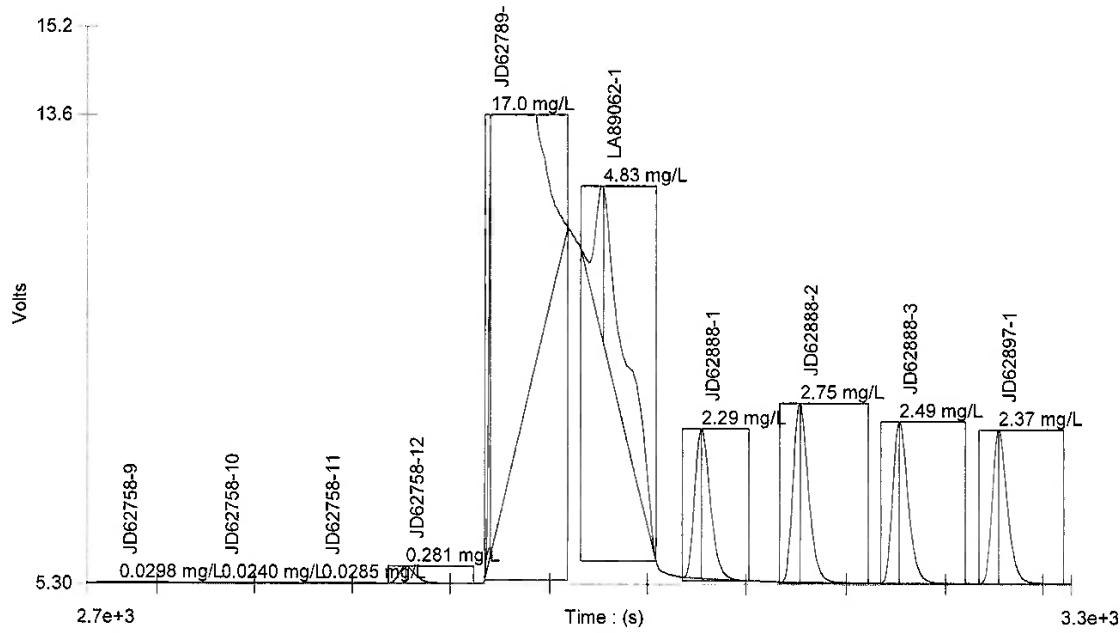


Date : 3/30/2023

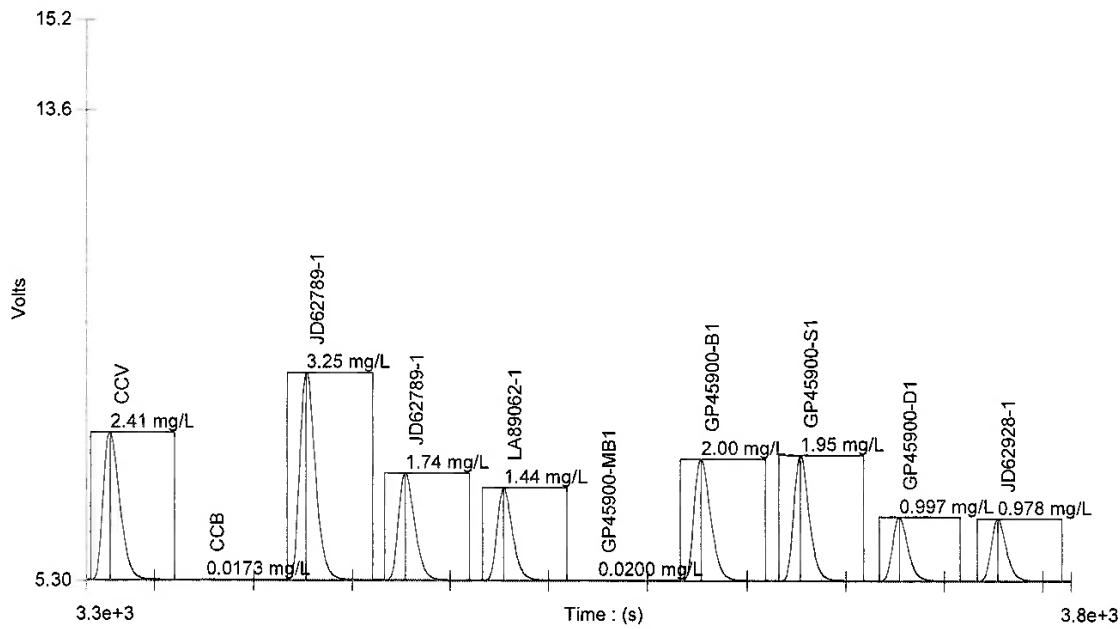
929

Author: Chemistry

Channel 1 - Set: 6 / 8



Channel 1 - Set: 7 / 8



Date : 3/30/2023

Author: Chemistry

Channel 1 - Set: 8 / 8

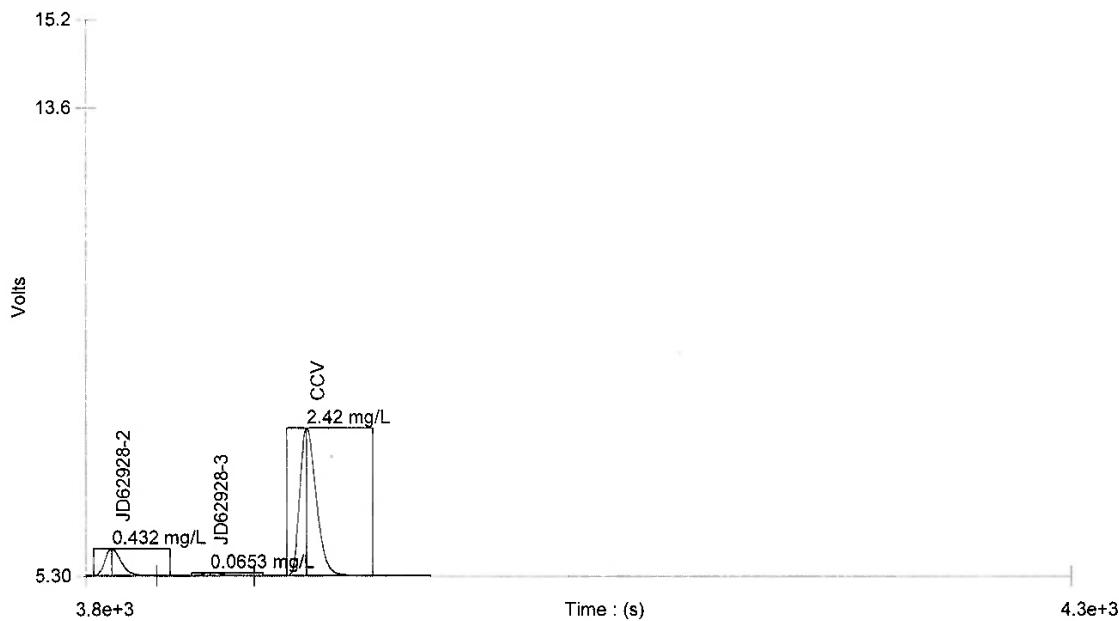
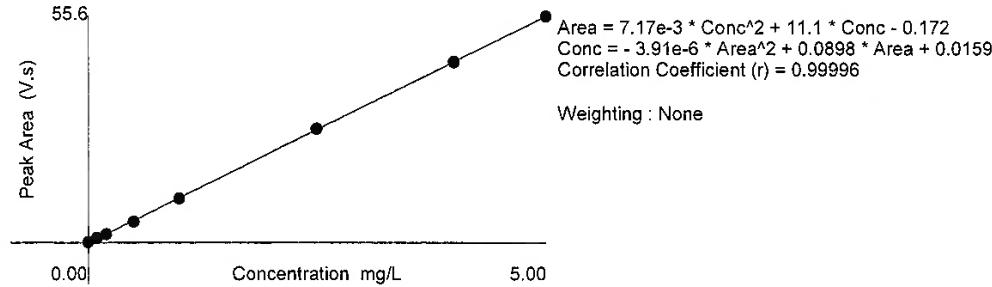


Table : 1 (NO32)

	Known Conc. (mg/L)	Rep.	Peak Area (V.s)	Peak Height (V)	% RSD	% Residual	Det. Conc (mg/L)	Detection Date	Detection Time
1	5.00	1	55.6	5.42	0.0	7.6e-3	5.00	3/30/2023	2:09:14 PM
2	4.00	1	44.3	4.33	0.0	0.2	3.99	3/30/2023	2:10:19 PM
3	2.50	1	27.9	2.73	0.0	-1.0	2.52	3/30/2023	2:11:25 PM
4	1.00	1	10.8	1.05	0.0	1.6	0.984	3/30/2023	2:12:30 PM
5	0.500	1	5.10	0.489	0.0	5.4	0.474	3/30/2023	2:13:35 PM
6	0.200	1	1.99	0.188	0.0	3.2	0.194	3/30/2023	2:14:40 PM
7	0.100	1	1.07	0.0997	0.0	-13.6	0.112	3/30/2023	2:15:44 PM
8	0.00	1	0.0461	4.82e-3			0.0200	3/30/2023	2:16:49 PM

Figure : 1 (NO32)



Date : 3/30/2023

Author: Chemistry

Original Run Filename: OM\_3-30-2023\_03-51-04PM.OMN Created: 3/30/2023 3:51:04 PM

Original Run Author's Signature: [Chemistry]

Current Run Filename: OM\_3-30-2023\_03-51-04PM.OMN Last Modified: 3/30/2023 4:12:23 PM

Current Run Author's Signature: [Chemistry]

Description: Default New Run

e033023w3.r032

96.8%

Sample	Rep.	Cup No.	Channel 1	Detection Time	MDF
			NO32 (mg/L)		
CCV	1	12	2.42	3/30/2023@3:51:57 PM	
		Known Conc:	2.50		
CCB	1	13	0.0257	3/30/2023@3:53:01 PM	
		Known Conc:	0.00		
		Calibration:	Table/Fig. : 1		
JD62928-4	1	71	0.0243	3/30/2023@3:54:05 PM	
JD62929-1	1	72	0.115	3/30/2023@3:55:09 PM	
JD62929-2	1	73	0.0466	3/30/2023@3:56:14 PM	
JD62929-3	1	74	1.21	3/30/2023@3:57:18 PM	
LA89039-1	1	75	2.95	3/30/2023@3:58:22 PM	
LA89039-2	1	76	12.1	3/30/2023@3:59:27 PM	
LA89039-3	1	77	4.15	3/30/2023@4:00:33 PM	
CCV	1	12	2.40	3/30/2023@4:03:10 PM	
		Known Conc:	2.50		
CCB	1	13	0.0226	3/30/2023@4:04:14 PM	
		Known Conc:	0.00		
CONF	1	1	4.85	3/30/2023@4:05:52 PM	
CCV	1	12	2.42	3/30/2023@4:08:29 PM	
		Known Conc:	2.50		
CCB	1	13	0.0226	3/30/2023@4:09:33 PM	
		Known Conc:	0.00		
LA89039-2	1	81	2.93	3/30/2023@4:10:38 PM	8.00
CCV	1	12	2.41	3/30/2023@4:13:16 PM	
		Known Conc:	2.50		
CCB	1	13	0.0242	3/30/2023@4:14:20 PM	
		Known Conc:	0.00		

96.1%

96.8%

96.8%

1 ml Sample + 7 mL DI

96.4%

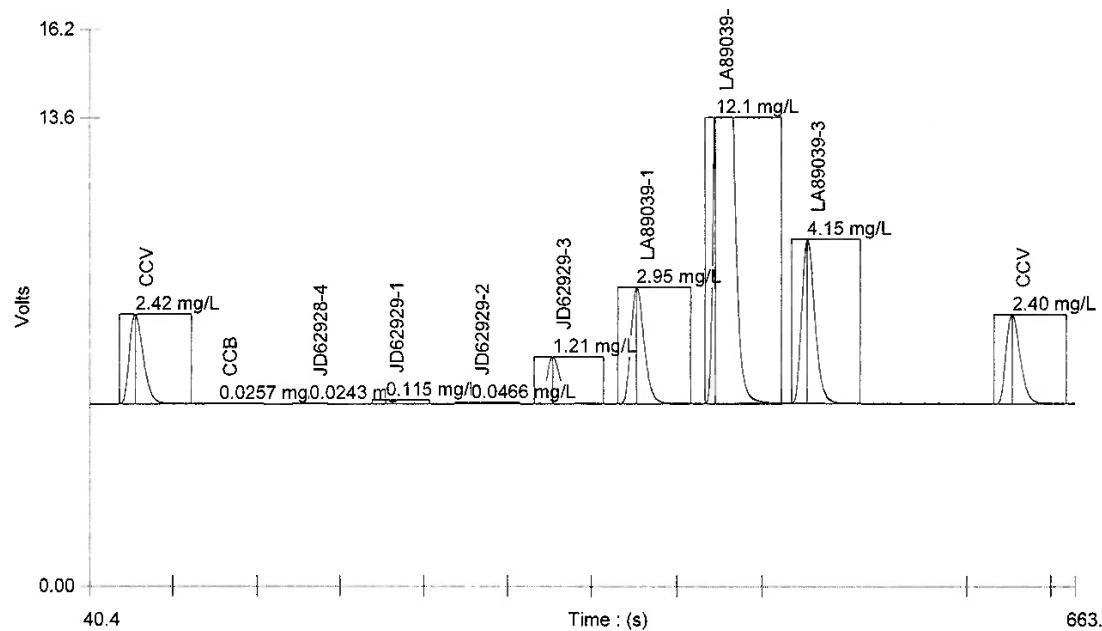
9.2

9

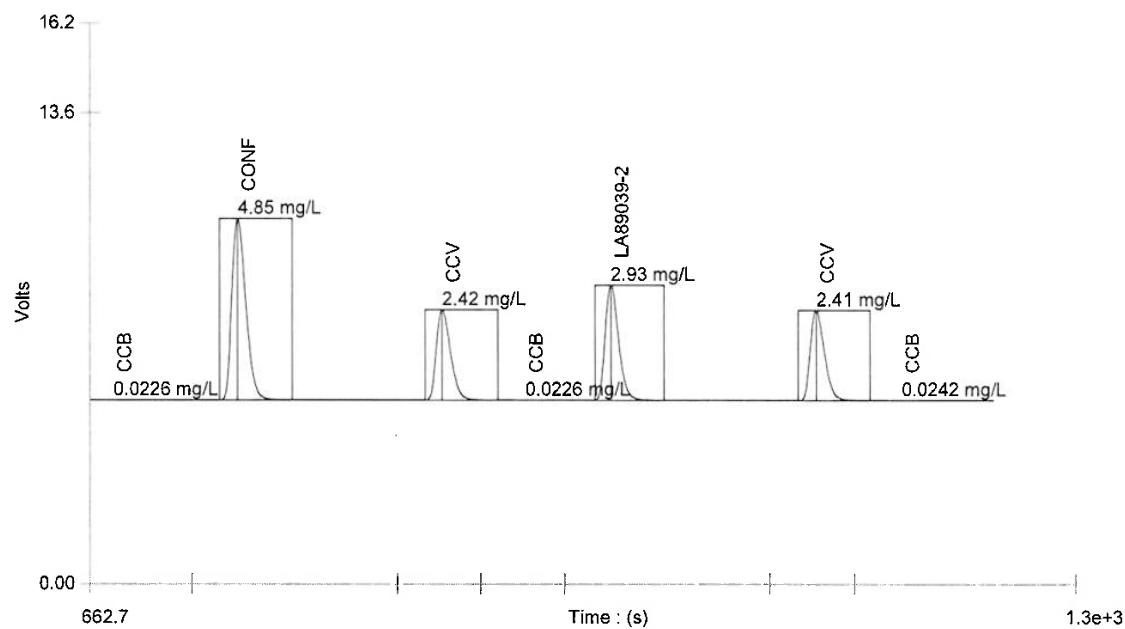
Date : 3/30/2023

Author: Chemistry

Channel 1 - Set: 1 / 2



Channel 1 - Set: 2 / 2



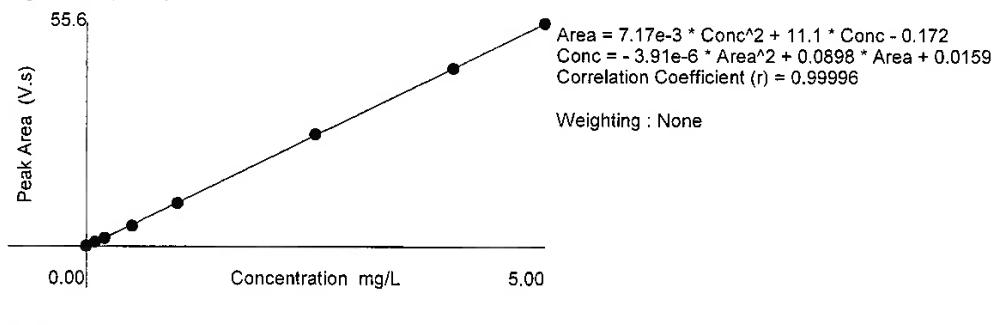
Author: Chemistry

Date : 3/30/2023

Table : 1 (NO32)

	Known Conc. (mg/L)	Rep.	Peak Area (V.s)	Peak Height (V)	% RSD	% Residual	Det. Conc (mg/L)	Detection Date	Detection Time
1	5.00	1	55.6	5.42	0.0	7.6e-3	5.00	3/30/2023	2:09:14 PM
2	4.00	1	44.3	4.33	0.0	0.2	3.99	3/30/2023	2:10:19 PM
3	2.50	1	27.9	2.73	0.0	-1.0	2.52	3/30/2023	2:11:25 PM
4	1.00	1	10.8	1.05	0.0	1.6	0.984	3/30/2023	2:12:30 PM
5	0.500	1	5.10	0.489	0.0	5.4	0.474	3/30/2023	2:13:35 PM
6	0.200	1	1.99	0.188	0.0	3.2	0.194	3/30/2023	2:14:40 PM
7	0.100	1	1.07	0.0997	0.0	-13.6	0.112	3/30/2023	2:15:44 PM
8	0.00	1	0.0461	4.82e-3			0.0200	3/30/2023	2:16:49 PM

Figure : 1 (NO32)



SGS Dayton

Analyst: \_\_\_\_\_ MM \_\_\_\_\_

Prep Date: 03/30/2023

Pipette ID: 52 and 57

Batch ID: GP45893

pH paper Lot: 206722 XP 03/01/2025

Nitrate-Nitrite (NO<sub>3</sub>) AQUEOUS QC Prep Log / Sample pH Log

Sample ID	Standard Prep	Std Final Volume		SAMPLE pH 5 - 9 ( Y / N )	Chlorine Present ( Y / N ) If Y, add sodium thiosulfate
EDTA BUFFER SOLN.	N/A	N/A		9.1	
MB				Y	
BS	1.0ML OF 100PPM	50MLS	carrier		
MS1	0.5ML OF 100PPM	50MLS	JD62498-5		
MS2	0.5ML OF 100PPM	50MLS	JD62808-1		
DUP			JD62498-5		
JD62498-1					
JD62498-2					
JD62498-4					
JD62498-5					
JD62498-6					
JD62498-7					
JD62498-8					
JD62498-9					
JD62498-10					
JD62498-11					
JD62808-1					
JD62808-2					
JD62888-1					
JD62888-2					
JD62888-3					
JD62888-4					
JD62896-1					
JD62897-1					

Form GN319-04 Aqueous

Reviewer:

Rev date: 03/06/2023

Date:

SGS Dayton

Analyst: \_\_\_\_\_ MM \_\_\_\_\_

Prep Date: 03/30/2023

Pipette ID: 52 and 57

Batch ID: GP45894

pH paper Lot: 206722 XP 03/01/2025

Nitrate-Nitrite (NO<sub>3</sub>) AQUEOUS QC Prep Log / Sample pH Log

Sample ID	Standard Prep	Std Final Volume		SAMPLE pH 5 - 9 (Y / N)	Chlorine Present (Y / N) If Y, add sodium thiosulfate
EDTA BUFFER SOLN.	N/A	N/A		9.1	
MB				Y	
BS	1.0ML OF 100PPM	50MLS	carrier		
MS1	0.5ML OF 100PPM	50MLS	JD62596-14		
MS2	0.5ML OF 100PPM	50MLS	JD62515-1		
DUP			JD62596-14		
JD62596-14					
JD62596-7					
JD62596-8					
JD62515-1					
JD62515-2					
JD62515-3					
JD62515-4					
JD62758-1					
JD62758-4					
JD62758-5					
JD62758-7					
JD62758-8					
JD62758-9					
JD62758-10					
JD62758-11					
JD62758-12					
JD62789-1					
LA89062-1					

Form GN319-04 Aqueous

**Reviewer:**

Rev date: 03/06/2023

Date:

SGS Dayton

Analyst: MM

Prep Date: 03/30/2023

Pipette ID: 52 and 57

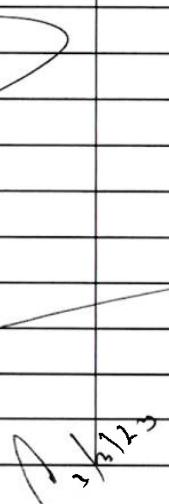
Batch ID: GP45900

pH paper Lot: 206722 XP 03/01/2025

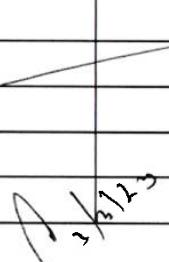
## Nitrate-Nitrite (NO32) AQUEOUS QC Prep Log / Sample pH Log

Sample ID	Standard Prep	Std Final Volume		SAMPLE pH 5 - 9 ( Y / N )	Chlorine Present (Y / N) If Y, add sodium thiosulfate
EDTA BUFFER SOLN.	N/A	N/A		9.1	
MB				Y	
BS	1.0ML OF 100PPM	50MLS	carrier		
MS1	0.5ML OF 100PPM	50MLS	JD62928-1		
DUP			JD62928-1		
JD62928-1					
JD62928-2					
JD62928-3					
JD62928-4					
JD62929-1					
JD62929-2					
JD62929-3					
LA89039-1					
LA89039-2					
LA89039-3					

Form GN319-04 Aqueous

Reviewer: 

Rev date: 03/06/2023

Date: 9.2  
9



GENERAL CHEMISTRY STANDARD PREPARATION LOG

Product:  
Analyst:  
Date:

NO32 MM 3/30/2023

**GN39816**

\*\*If Class A glass pipettes are used, enter an A.

Form: GN205-03  
Rev Date: 03/22/2022



## Reagent Information Log - Nitrate Lachat Autoanalyzer

<u>Reagent</u>	<u>Reagent # or Manufacturer/Lot</u>	<u>Expiration Date</u>
Nitrate Stock Solution	GNE12-72667-NO32	6/10/2023
Ammonium Chloride Buffer Solution	GNE11-72430-NO3	5/23/2023
Sulfanilamide Color Reagent	GNE3-73551-NO32	4/2/2023
1:1 NH4OH	GNE11-72312-NO3	6/15/2023
Carrier Solution	GNE1-73258-NO32	8/2/2023
1000 ppm Nitrite Solution	GNE12-2665-NO32	6/10/2023
Nitrate External Stock Solution	GNE12-72666-NO32	6/10/2023

Reason codes for data corrections: 1-reviewer error correction; 2-transcription error; 3-computer error; 4-analyst error

Form: GN087A-43  
Rev. Date: 7/19/06

LABORATORY REVIEW SIGNATURE FORM  
(To be stored with the raw data)

File ID: D230330W1.TXT      Date Analyzed: 03/01/23      Methods: SM5310 B-11/14  
Analyst: MB      Run ID: GN39817

The following analyst(s) have reviewed this run and attest that, to the best of their knowledge, this documentation is complete and correct:

Analyst: MB      Date 3-30-23

Analyst: \_\_\_\_\_ Date \_\_\_\_\_

The following supervisor or their designee has reviewed this run and attests that, to the best of their knowledge, this documentation is complete and correct:

Supervisor (or designee):         Date 3/11/23

9.3  
6

C:\TOC-L\Data\d230330w1.toc.tlx

	Sample Name	Sample ID	Origin	Manual Dilution	Result	Comment
1	WASHCONF		TOCAQSW846.met	1.000	NPOC:0.5131mg/L	
2	CRI		TOCAQSW846.met	1.000	NPOC:1.097mg/L	
3	HSTD		TOCAQSW846.met	1.000	NPOC:47.05mg/L	
4	ICV		TOCAQSW846.met	1.000	NPOC:19.89mg/L	
5	ICB		TOCAQSW846.met	1.000	NPOC:0.3547mg/L	
6	CCV		TOCAQSW846.met	1.000	NPOC:24.13mg/L	
7	CCB		TOCAQSW846.met	1.000	NPOC:0.8087mg/L	
8	SPARGERCHK		TOCAQSW846.met	1.000	NPOC:0.6197mg/L	
9	GP45861-MB1	[toc]	TOCAQ.met	1.000	NPOC:0.4633mg/L	
10	GP45861-B1		TOCAQ.met	1.000	NPOC:9.211mg/L	
11	JD62808-1		TOCAQ.met	1.000	NPOC:5.029mg/L	
12	GP45861-S1	JD62808-1	TOCAQ.met	1.000	NPOC:14.65mg/L	
13	GP45861-MSD1	JD62808-1	TOCAQ.met	1.000	NPOC:14.41mg/L	
14	JD62888-4	field blank	TOCAQ.met	1.000	NPOC:0.3454mg/L	
15	JD62808-2		TOCAQ.met	1.000	NPOC:5.617mg/L	
16	JD62808-3		TOCAQ.met	1.000	NPOC:5.159mg/L	
17	JD62808-4		TOCAQ.met	1.000	NPOC:5.154mg/L	
18	CCVA		TOCAQ.met	1.000	NPOC:46.44mg/L	
19	CCB		TOCAQ.met	1.000	NPOC:0.6588mg/L	
20	JD62808-5		TOCAQ.met	1.000	NPOC:4.828mg/L	
21	JD62888-1		TOCAQ.met	1.000	NPOC:1.012mg/L	
22	JD62888-2		TOCAQ.met	1.000	NPOC:0.8136mg/L	
23	JD62888-3		TOCAQ.met	1.000	NPOC:1.073mg/L	
24	CCV		TOCAQ.met	1.000	NPOC:24.09mg/L	
25	CCB		TOCAQ.met	1.000	NPOC:0.6822mg/L	
26	SPARGERCHK		TOCAQ.met	1.000	NPOC:0.3946mg/L	

d230330 w1 toc

GN39817

MB

3-30-23

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GN Batch ID: GN39817  
Date: 3-30-23

Test: Total Organic Carbon

d230330wl.tocProduct: TOC or DOCMethod: SM5310 B, C, or D-11, SW846 9060M

Note: Refer to raw data and LIMS for information not shown below.

Autosampler Position #	Sample ID	pH	Dilution Factor	Bottle #	Comments
1	WASHCONF	<2			
2	CRI				
3	HSTD				
4	ICV				
5	ICB				
6	CCV				
7	CCB				
8	SPARGERCHK				
9	GP45861-MB1			-	[toc]
10	GP45861-B1			-	
11	JD62808-1			12	
12	GP45861-S1			12	JD62808-1
13	GP45861-MSD1			12	JD62808-1
14	JD62888-4			5	Field Blank
15	JD62808-2			12	
16	JD62808-3			1	
17	JD62808-4			1	
18	CCVA				
19	CCB				
20	JD62808-5			5	
21	JD62888-1			5	
22	JD62888-2			5	
23	JD62888-3			5	
24	CCV				
25	CCB				
26	SPARGERCHK				

Analyst: MB Date: 3-30-23 QCReviewer: \_\_\_\_\_ Date: \_\_\_\_\_

Comments: BSP: 500ul of 1000ppm KHP up to 50mL with DI H2O TV= 10 mg/L

MS/MSD: 200ul of 1000ppm KHP up to 20mL with sample TV=10 mg/L

ICV: 5mL of 100 ppm Sucrose up to 25mL with DI H2O TV=20 mg/L

Form: GN054-04

Rev. Date: 2/27/18

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## Reagent Information Log - TOC/DOC

GN39817

<u>Reagent</u>	<u>Expiration Date</u>	<u>Reagent # or Manufacturer/Lot</u>
Potassium Hydrogen Phthalate (KHP), Stock Solution 1000 mg/L	4/13/2023	GNE1-72979-TOC
Carbonate/Bicarbonate Stock Solution	4/13/2023	GNE1-72982-TOC
Sparger Check Solution	4/13/2023	GNE3-70727-TOC
CCV Solution	4/14/2023	GNE3-70678-TOC
CCVA Solution (50 ppm)	4/14/2023	GNE3-70679-TOC
CRI Check	4/13/2023	GNE3-70729-TOC
HCl	4/10/2024	VWR CHEMICALS LOT #2022041418
pH Hydron Paper	8/1/2024	FISHER LOT #222221
Sucrose Stock Solution	4/14/2023	GNE3-70676-TOC

All standards and stocks were made as described in the SOP for this method (circle one):  Y or N  
If no (N), see attached page for standards prep.

Form: GN087A67-043  
Rev Date: 3/17/2023

9.3  
9



## GENERAL CHEMISTRY STANDARD PREPARATION LOG FOR TOC (AQ)

Product:

TOC

Analyst:

Mβ

Date:

3-30-23

GN/GP number:

Intermediate Standard Description	Stock used to prepare standard	Stock concentration (mg/L)	Stock volume used (mL or g)	Balance/Pipette ID*	Diluent(a)	Final Volume	Final Conc. of Intermediate (mg/L)	Expiration Date	Analyst	Date
GNE1-72979-TOC	Fisher Lot: 178879	KHP	2.1528 g	B-39	DI H <sub>2</sub> O	1000 mL	1000 ppm	4/13/2023	MB	1/12/2023
GNE3-70728-TOC	GNE1-72979-TOC	1000 ppm	20 mL	Class A Pipet	DI H <sub>2</sub> O	200 mL	100 ppm	4/13/2023	MB	3/28/2023
GNE3-70676-TOC	Fisher Lot: 210554	Sucrose	0.0486 g	B-39	DI H <sub>2</sub> O	200 mL	100 ppm	4/14/2023	MB	3/17/2023
KHP STDs										
GNE3-70729-TOC	GNE3-70728-TOC	100 ppm	1.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	10	4/13/2023	MB	3/28/2023
GNE3-70730-TOC	GNE3-70728-TOC	100 ppm	2.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	20	4/13/2023	MB	3/28/2023
GNE3-70731-TOC	GNE3-70728-TOC	100 ppm	5.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	50	4/13/2023	MB	3/28/2023
GNE3-70732-TOC	GNE3-70728-TOC	100 ppm	10.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	10.0	4/13/2023	MB	3/28/2023
GNE3-70733-TOC	GNE3-70728-TOC	100 ppm	20.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	20.0	4/13/2023	MB	3/28/2023
GNE3-70734-TOC	GNE3-70728-TOC	100 ppm	30.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	30.0	4/13/2023	MB	3/28/2023
GNE3-70735-TOC	GNE3-70728-TOC	100 ppm	50.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	50.0	4/13/2023	MB	3/28/2023
KHP STDs										
GNE3-70677-TOC	Acros Lot: A0383564	KHP	0.2307 g	B-39	DI H <sub>2</sub> O	100 mL	100 ppm	4/14/2023	MB	3/17/2023
GNE3-70678-TOC	GNE3-70677-TOC	100 ppm	50 mL	Class A Pipet	DI H <sub>2</sub> O	100 mL	25 ppm	4/14/2023	MB	3/17/2023
GNE3-70679-TOC	GNE3-70677-TOC	100 ppm	100 mL	Class A Pipet	DI H <sub>2</sub> O	100 mL	50 ppm	4/14/2023	MB	3/17/2023

\*If Class A glass pipettes are used, enter an A.

(a) Diluent reagent reference number.

All strikeouts must be initialed, dated, and reason applied if not transcription error

Expiration date:

# TOC-Control L Report

Chem36  
d230201wl toc tx

**Instr.Information**

Instrument Options  
Catalyst

TOC/ASI/Sparge Kit/  
Regular Sensitivity

**Cal. Curve**

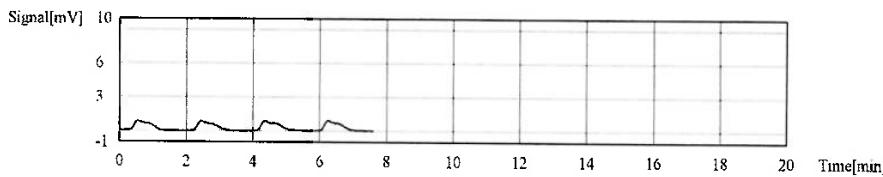
Sample Name: Untitled  
 Sample ID: Untitled  
 Cal. Curve: d230201wl.2023\_02\_01\_11\_45\_27.cal  
 Status: Completed

Type	Anal.
Standard	NPOC

Conc: 0.000mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	2.806	100uL	1.000	*****		2/1/2023 11:50:08 AM
2	2.648	100uL	1.000	*****		2/1/2023 11:52:23 AM
3	2.915	100uL	1.000	*****		2/1/2023 11:54:38 AM
4	2.824	100uL	1.000	*****		2/1/2023 11:56:53 AM

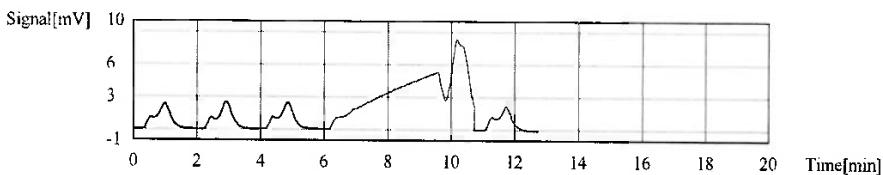
Acid Add. 0.000%  
 Sp. Time 600.0sec  
 Mean Area 2.798



Conc: 1.000mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	7.402	100uL	1.000	*****		2/1/2023 12:01:30 PM
2	7.958	100uL	1.000	*****		2/1/2023 12:03:47 PM
3	7.934	100uL	1.000	*****		2/1/2023 12:06:08 PM
4	57.64	100uL	1.000	T*****	E	2/1/2023 12:11:23 PM
5	7.582	100uL	1.000	*****		2/1/2023 12:14:09 PM

Acid Add. 0.000%  
 Sp. Time 600.0sec  
 Mean Area 7.719



Conc: 2.000mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	16.98	100uL	1.000	*****		2/1/2023 12:18:47 PM
2	17.18	100uL	1.000	*****		2/1/2023 12:21:13 PM
3	17.02	100uL	1.000	*****		2/1/2023 12:23:41 PM
4	16.83	100uL	1.000	*****		2/1/2023 12:26:00 PM

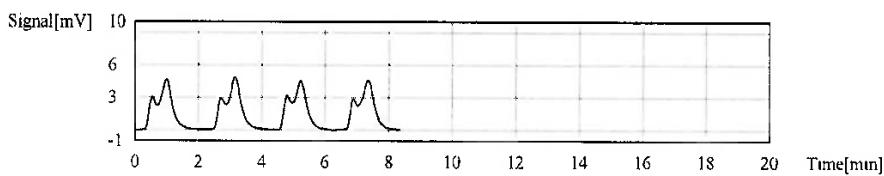
9.3

6

# TOC-Control L Report

Chem36  
d230201wl.toc.tlx

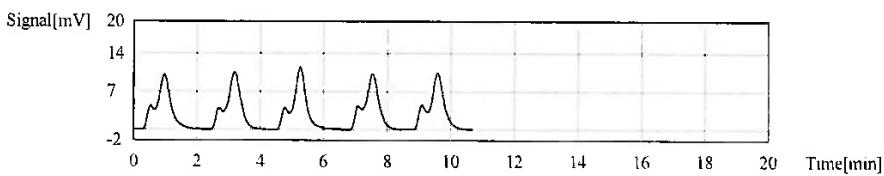
Acid Add. 0.000%  
Sp. Time 600.0sec  
Mean Area 17.00



Conc: 5.000mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	32.71	100µL	1.000	*****	E	2/1/2023 12:30:53 PM
2	33.99	100µL	1.000	*****		2/1/2023 12:33:18 PM
3	34.26	100µL	1.000	*****		2/1/2023 12:35:58 PM
4	33.02	100µL	1.000	*****		2/1/2023 12:38:19 PM
5	33.81	100µL	1.000	*****		2/1/2023 12:40:48 PM

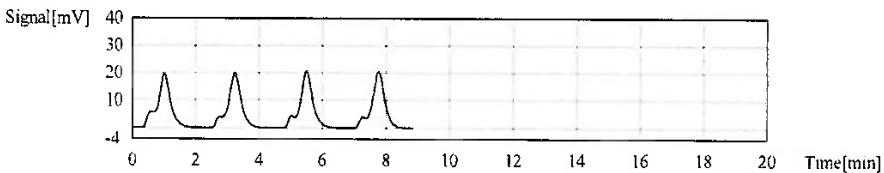
Acid Add. 0.000%  
Sp. Time 600.0sec  
Mean Area 33.77



Conc: 10.00mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	58.33	100µL	1.000	*****		2/1/2023 12:45:28 PM
2	58.49	100µL	1.000	*****		2/1/2023 12:48:06 PM
3	58.80	100µL	1.000	*****		2/1/2023 12:50:42 PM
4	58.79	100µL	1.000	*****		2/1/2023 12:53:08 PM

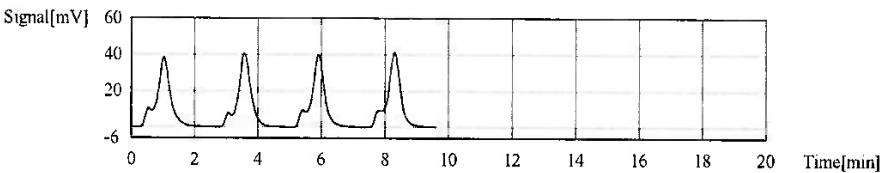
Acid Add. 0.000%  
Sp. Time 600.0sec  
Mean Area 58.60



Conc: 20.00mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	116.5	100µL	1.000	*****		2/1/2023 12:58:24 PM
2	117.3	100µL	1.000	*****		2/1/2023 1:01:04 PM
3	116.9	100µL	1.000	*****		2/1/2023 1:03:48 PM
4	117.7	100µL	1.000	*****		2/1/2023 1:06:32 PM

Acid Add. 0.000%  
Sp. Time 600.0sec  
Mean Area 117.1



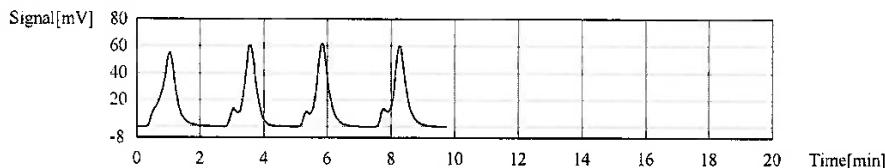
Conc: 30.00mg/L

# TOC-Control L Report

Chem365  
d230201wl.toc.tlx

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	174.8	100µL	1.000	*****		2/1/2023 1:11:44 PM
2	174.5	100µL	1.000	*****		2/1/2023 1:14:24 PM
3	176.4	100µL	1.000	*****		2/1/2023 1:17:10 PM
4	176.3	100µL	1.000	*****		2/1/2023 1:20:00 PM

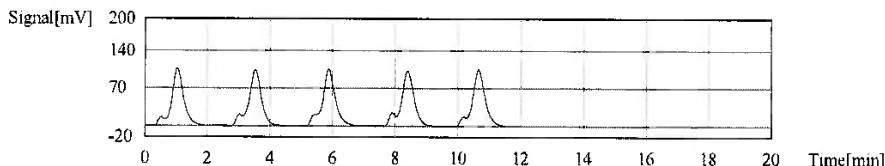
Acid Add. 0.000%  
Sp. Time 600.0sec  
Mean Area 175.5



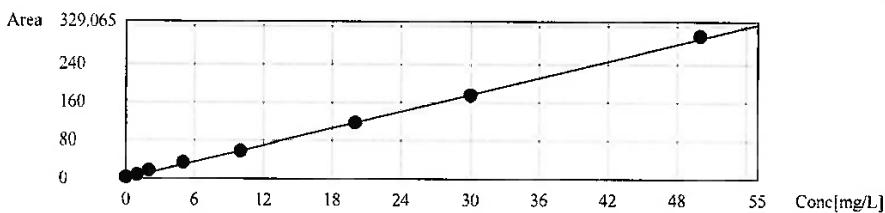
Conc: 50.00mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	292.2	100µL	1.000	*****	E	2/1/2023 1:25:12 PM
2	299.6	100µL	1.000	*****		2/1/2023 1:27:56 PM
3	301.2	100µL	1.000	*****		2/1/2023 1:30:48 PM
4	295.9	100µL	1.000	*****		2/1/2023 1:33:24 PM
5	299.9	100µL	1.000	*****		2/1/2023 1:36:10 PM

Acid Add. 0.000%  
Sp. Time 600.0sec  
Mean Area 299.2



Slope: 5.875  
Intercept 0.000  
 $r^2$  0.9994  
 $r$  0.9997  
Zero Shift Yes



9.3  
6

# TOC-Control L Report

Chem56  
d230330wl.toc.tbx

**Instr.Information**

Instrument Options  
Catalyst

TOC/ASI/Sparge Kit/  
Regular Sensitivity

**Sample**

Sample Name: WASHCONF  
Sample ID:  
Origin:  
Status Completed  
Chk. Result

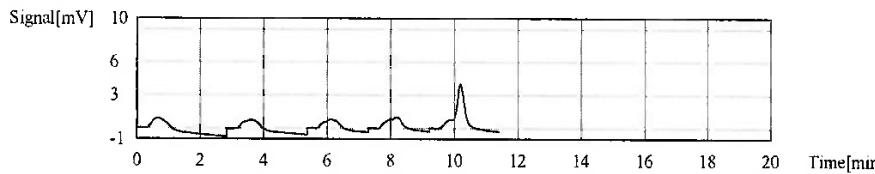
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.5131mg/L

## 1. Det

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2 179	0.3709mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:28:51 AM	
2	3 358	0.5715mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:31:38 AM	
3	2 918	0.4966mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:33:55 AM	
4	3 605	0.6136mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:36:05 AM	
5	8 560	1.457mg/L	100uL	1 000	E d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:38:31 AM	

Mean Conc. 0.5131mg/L  
CV Conc 20.75%

**Sample**

Sample Name: CRI  
Sample ID:  
Origin:  
Status Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 1.097mg/L

## 1. Det

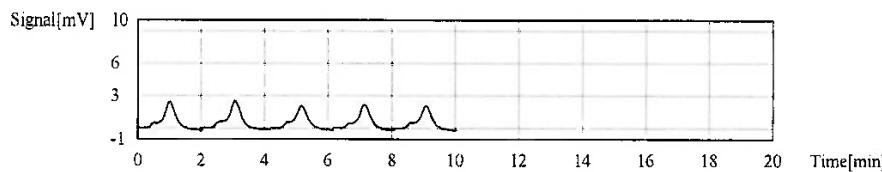
Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	6 943	1.182mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:44:02 AM	
2	7 321	1.246mg/L	100uL	1 000	E d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:46:26 AM	
3	6 342	1.079mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:48:39 AM	
4	6 418	1.092mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:50:52 AM	
5	6 077	1.034mg/L	100uL	1 000	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:53:04 AM	

# TOC-Control L Report

Chem365  
d230330wl.toc.tlk

Mean Conc. 1.097mg/L  
CV Conc 5.63%



## Sample

Sample Name: HSTD  
 Sample ID:  
 Origin: TOCAQSW846 met  
 Status Completed  
 Chk. Result

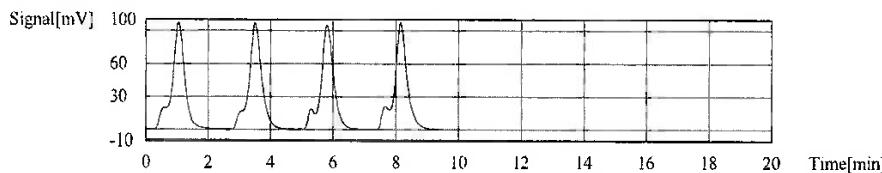
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 47.05mg/L

### 1. Det

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	277.7	47.26mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 9:59:00 AM
2	278.4	47.38mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:01:34 AM
3	271.5	46.21mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:04:11 AM
4	278.2	47.35mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:06:52 AM

Mean Conc. 47.05mg/L  
CV Conc 1.20%



## Sample

Sample Name: ICV  
 Sample ID:  
 Origin: TOCAQSW846 met  
 Status Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 19.89mg/L

### 1. Det

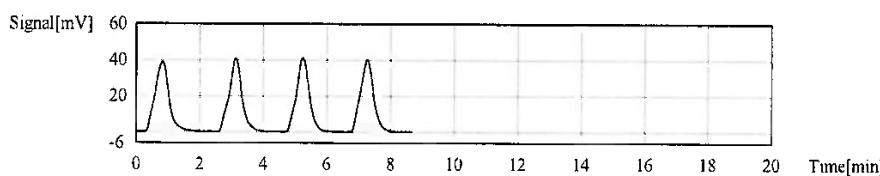
Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	117.1	19.93mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:30:58 AM
2	116.8	19.88mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:33:18 AM
3	116.8	19.88mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:35:32 AM
4	116.8	19.88mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:37:54 AM

# TOC-Control L Report

Chem56  
d230330wl.toc.tlx

Mean Conc. 19.89mg/L  
CV Conc 0.13%



## Sample

Sample Name: ICB  
 Sample ID:  
 Origin: TOCAQSW846.met  
 Status: Completed  
 Chk. Result

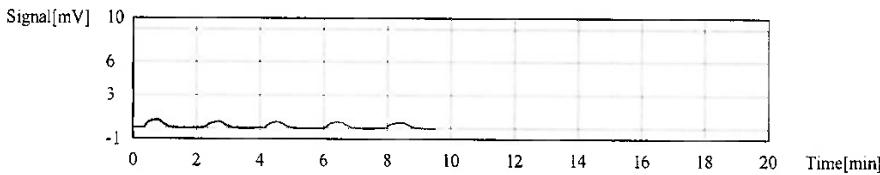
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.3547mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2.269	0.3862mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:45:21 AM
2	1.999	0.3402mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:45:26 AM
3	1.770	0.3013mg/L	100uL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:47:31 AM
4	2.034	0.3462mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.est	3/1/2023 10:49:36 AM
5	2.035	0.3464mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/1/2023 10:51:41 AM

Mean Conc. 0.3547mg/L  
CV Conc 5.96%



## Sample

Sample Name: CCV  
 Sample ID:  
 Origin: TOCAQSW846.met  
 Status: Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 24.13mg/L

### 1. Det

Anal.: NPOC

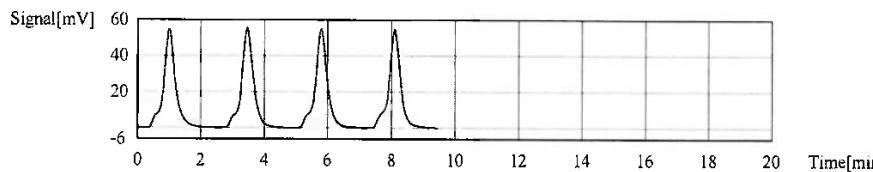
9.3  
6

# TOC-Control L Report

Chem365  
d230330wl.toc.tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	140.6	23.93mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:03:20 PM
2	143.3	24.39mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:05:54 PM
3	143.4	24.41mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:08:23 PM
4	139.7	23.78mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:11:01 PM

Mean Conc. 24.13mg/L  
CV Conc 1.33%



## Sample

Sample Name: CCB  
 Sample ID:  
 Origin: TOCAQSW846.met  
 Status Completed  
 Chk. Result

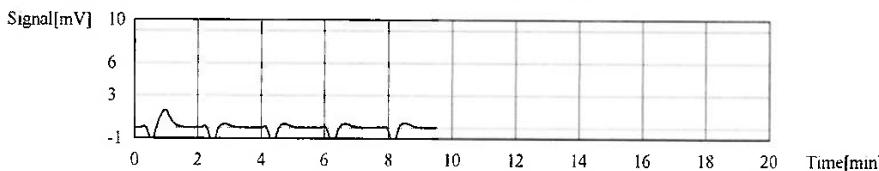
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC: 0.8087mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	6.871	1.169mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:16:39 PM
2	0.000	0.000mg/L	100µL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 12:18:44 PM
3	3.512	0.5977mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:20:49 PM
4	4.075	0.6936mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:22:54 PM
5	4.547	0.7739mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 12:24:59 PM

Mean Conc. 0.8087mg/L  
CV Conc 31.05%



## Sample

Sample Name: SPARGERCHK  
 Sample ID:  
 Origin: TOCAQSW846.met  
 Status Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC: 0.6197mg/L

### 1. Det

9.3

6

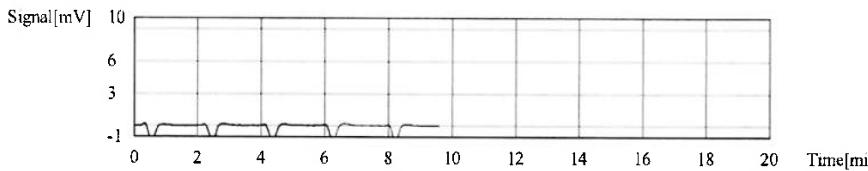
# TOC-Control L Report

Chem36  
d230330w1.toc.tlx

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	0.000	0.000mg/L	100uL	1 000	E	d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:30:28 PM
2	3.310	0.5634mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:32:33 PM
3	3.653	0.6217mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:34:38 PM
4	4.972	0.8462mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:36:48 PM
5	2.629	0.4475mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:38:53 PM

Mean Conc. 0.6197mg/L  
CV Conc 27.03%



## Sample

Sample Name: GP45861-MB1  
Sample ID: [toc]  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

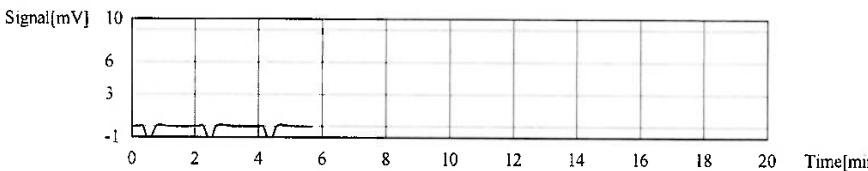
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.4633mg/L

1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2.767	0.4709mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:43:37 PM
2	3.067	0.5220mg/L	100uL	1 000	E	d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:45:46 PM
3	2.677	0.4556mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/30/2023 12:47:51 PM

Mean Conc. 0.4633mg/L  
CV Conc 2.34%



## Sample

Sample Name: GP45861-B1  
Sample ID: [toc]  
Origin: TOCAQ.met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.211mg/L

1. Det

9.3  
6

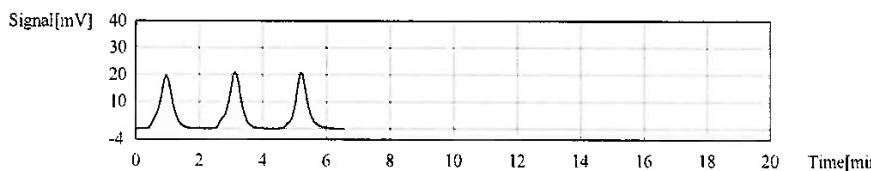
# TOC-Control L Report

Chem36  
d230330wl.toc.tlx

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	54.07	9.203mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/30/2023 12:55:02 PM
2	58.13	9.894mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/30/2023 12:57:18 PM
3	54.17	9.220mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/30/2023 12:59:49 PM

Mean Conc. 9.211mg/L  
CV Conc 0.13%



## Sample

Sample Name: JD62808-1  
Sample ID:  
Origin:  
Status TOCAQ.met  
Chk. Result Completed

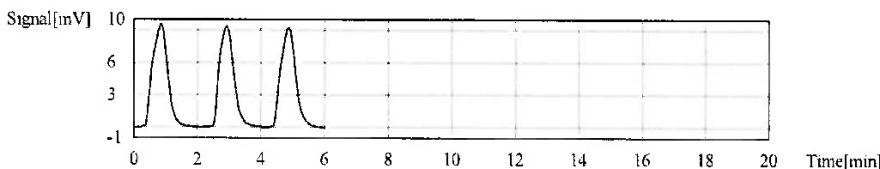
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 5.029mg/L

## 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	31.52	5.365mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/30/2023 2:02:00 PM
2	29.29	4.985mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/30/2023 2:04:05 PM
3	29.81	5.074mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/30/2023 2:06:13 PM

Mean Conc. 5.029mg/L  
CV Conc 1.24%



## Sample

Sample Name: GP45861-S1  
Sample ID: JD62808-1  
Origin:  
Status TOCAQ.met  
Chk. Result Completed

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 14.65mg/L

## 1. Det

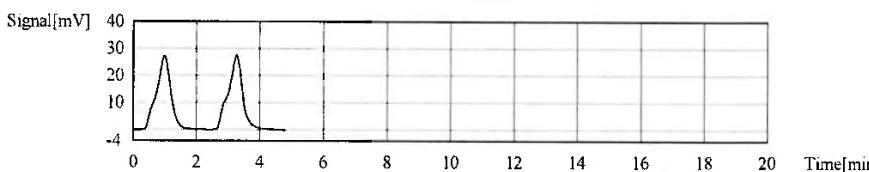
Anal.: NPOC

# TOC-Control L Report

Chem36  
d230330wl.toc.tbx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	84.92	14.45mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 2:13:18 PM
2	87.28	14.85mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 2:16:04 PM

Mean Conc. 14.65mg/L  
CV Conc 1.94%



## Sample

Sample Name: GP45861-MSD1  
 Sample ID: JD62808-1  
 Origin: TOCAQ met  
 Status: Completed  
 Chk. Result

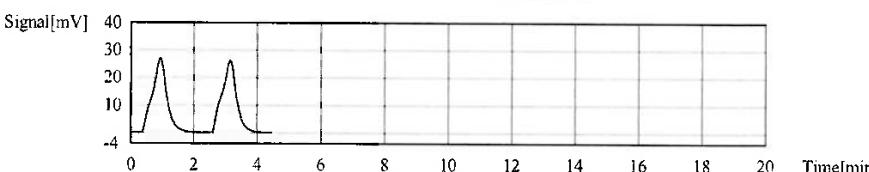
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 14.41mg/L

1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	85.50	14.55mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 2:24:24 PM
2	83.81	14.26mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 2:26:52 PM

Mean Conc. 14.41mg/L  
CV Conc 1.41%



## Sample

Sample Name: JD62888-4  
 Sample ID: field blank  
 Origin: TOCAQ met  
 Status: Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.3454mg/L

1. Det

Anal.: NPOC

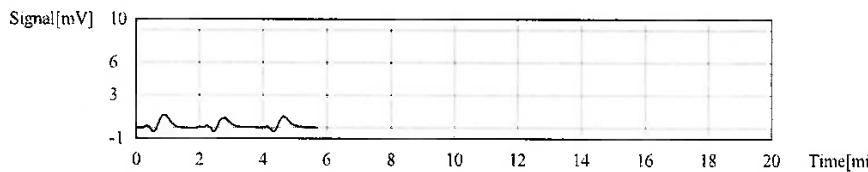
# TOC-Control L Report

Chem36  
d230330w1 toc.tbx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	2.345	0.3991mg/L	100uL	1.000		d230201w1.2023_02_01_11_45_27.cal	3/30/2023 2:35:14 PM
2	1.714	0.2917mg/L	100uL	1.000		d230201w1.2023_02_01_11_45_27.cal	3/30/2023 2:37:20 PM
3	3.441	0.5857mg/L	100uL	1.000	E	d230201w1.2023_02_01_11_45_27.cal	3/30/2023 2:39:24 PM

Mean Conc.  
CV Conc

0.3454mg/L  
21.98%



## Sample

Sample Name: JD62808-2  
 Sample ID:  
 Origin: TOCAQ.met  
 Status Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 5.617mg/L

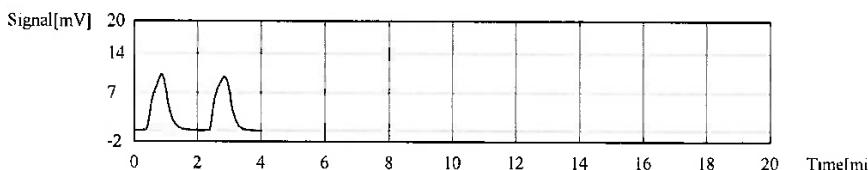
### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	32.75	5.574mg/L	100uL	1.000		d230201w1.2023_02_01_11_45_27.cal	3/30/2023 2:46:30 PM
2	33.25	5.659mg/L	100uL	1.000		d230201w1.2023_02_01_11_45_27.cal	3/30/2023 2:48:41 PM

Mean Conc.  
CV Conc

5.617mg/L  
1.07%



## Sample

Sample Name: JD62808-3  
 Sample ID:  
 Origin: TOCAQ.met  
 Status Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 5.159mg/L

### 1. Det

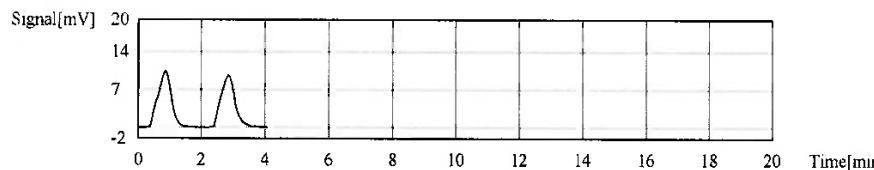
Anal.: NPOC

# TOC-Control L Report

Chem36  
d230330wl.toc.tbx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	31.21	5.312mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 2:57:40 PM
2	29.41	5.006mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 2:59:55 PM

Mean Conc. 5.159mg/L  
CV Conc 4.20%



## Sample

Sample Name: JD62808-4  
Sample ID:  
Origin:  
Status TOCAQ met  
Chk. Result Completed

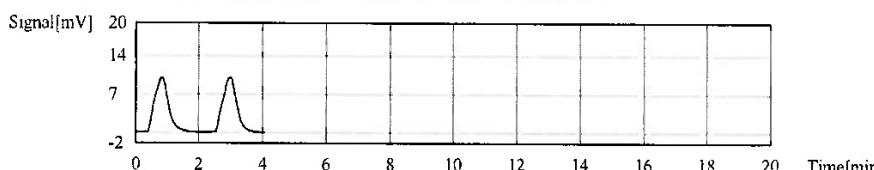
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 5.154mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	30.29	5.155mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 3:08:57 PM
2	30.27	5.152mg/L	100µL	1.000	d230201wl.2023_02_01_11_45_27.cal		3/30/2023 3:11:03 PM

Mean Conc. 5.154mg/L  
CV Conc 0.05%



## Sample

Sample Name: CCVA  
Sample ID:  
Origin:  
Status TOCAQ met  
Chk. Result Completed

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 46.44mg/L

### 1. Det

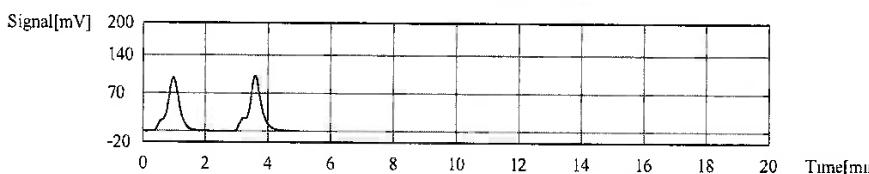
Anal.: NPOC

# TOC-Control L Report

Chem36  
d230330wl toc.tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	267.8	45.58mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27.cal		3/30/2023 3:20:33 PM
2	277.9	47.30mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27.cal		3/30/2023 3:23:15 PM

Mean Conc. 46.44mg/L  
CV Conc 2.62%



## Sample

Sample Name: CCB  
Sample ID.  
Origin: TOCAQ.met  
Status Completed  
Chk. Result

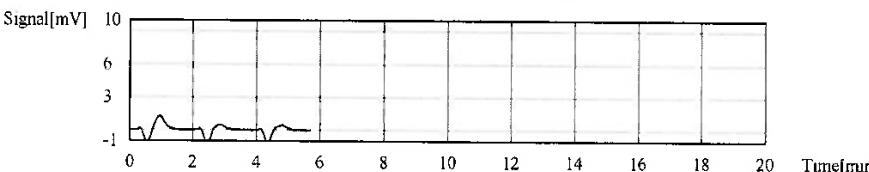
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.6588mg/L

1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	1459	0.2483mg/L	100uL	1.000	E	d230201wl2023_02_01_11_45_27.cal	3/30/2023 3:31:02 PM
2	3756	0.6393mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27.cal		3/30/2023 3:33:11 PM
3	3986	0.6784mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27.cal		3/30/2023 3:35:16 PM

Mean Conc. 0.6588mg/L  
CV Conc 4.20%



9.3

6

## Sample

Sample Name: JD62808-5  
Sample ID.  
Origin: TOCAQ.met  
Status Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 4.828mg/L

1. Det

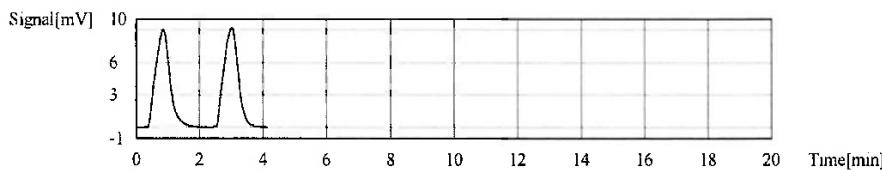
Anal.: NPOC

# TOC-Control L Report

Chem56  
d230330wl.toc.tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	28.00	4.766mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/30/2023 3:42:26 PM
2	28.73	4.890mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/30/2023 3:44:33 PM

Mean Conc. 4.828mg/L  
CV Conc 1.82%



## Sample

Sample Name: JD62888-1  
Sample ID:  
Origin: TOCAQ.net  
Status: Completed  
Chk. Result:

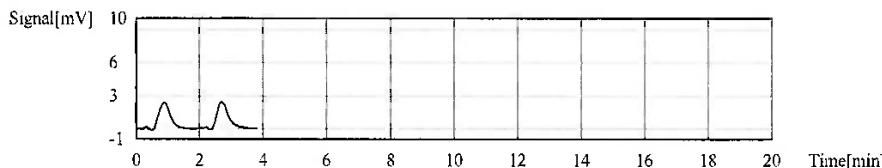
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 1.012mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	5.984	1.018mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/30/2023 3:53:19 PM
2	5.912	1.006mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/30/2023 3:55:25 PM

Mean Conc. 1.012mg/L  
CV Conc 0.86%



## Sample

Sample Name: JD62888-2  
Sample ID:  
Origin: TOCAQ.net  
Status: Completed  
Chk. Result:

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.8136mg/L

### 1. Det

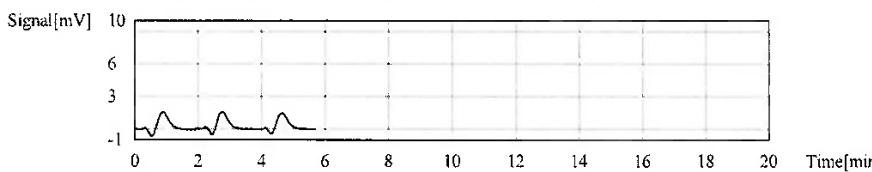
Anal.: NPOC

# TOC-Control L Report

Cham56  
d230330wl.toc.rlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	5 167	0.8794mg/L	100µL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:04:29 PM
2	3 (95	0.5268mg/L	100µL	1 000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:06:35 PM
3	4 394	0.7479mg/L	100µL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:08:39 PM

Mean Conc. 0.8136mg/L  
CV Conc 11.43%



## Sample

Sample Name: JD62888-3  
Sample ID:  
Origin:  
Status Completed  
Chk. Result

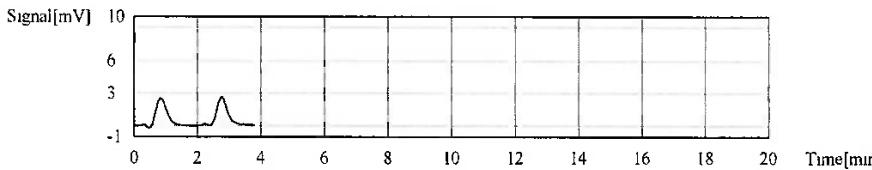
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.00X	NPOC 1.073mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	6 437	1.096mg/L	100µL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:15:39 PM
2	6 168	1.050mg/L	100µL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:17:43 PM

Mean Conc. 1.073mg/L  
CV Conc 3.02%



## Sample

Sample Name: CCV  
Sample ID:  
Origin:  
Status Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 24.09mg/L

### 1. Det

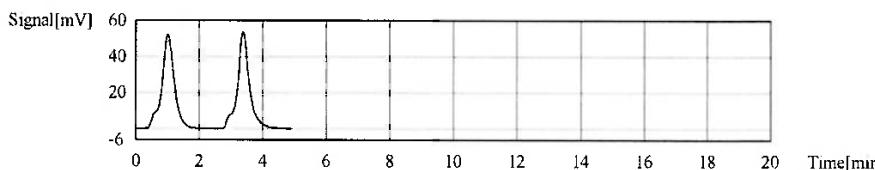
Anal.: NPOC

# TOC-Control L Report

Chem365  
d230350wl.toc.lbx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	141.8	24.13mg/L	100uL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:27:16 PM
2	141.3	24.05mg/L	100uL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:30:02 PM

Mean Conc. 24.09mg/L  
CV Conc 0.25%



## Sample

Sample Name: CCB  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

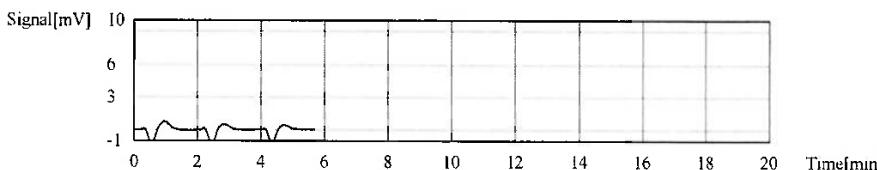
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.6822mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	0.1383	0.02354mg/L	100uL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:37:59 PM
2	4.364	0.7427mg/L	100uL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:40:09 PM
3	3.653	0.6217mg/L	100uL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:42:14 PM

Mean Conc. 0.6822mg/L  
CV Conc 12.54%



## Sample

Sample Name: SPARGERCHK  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.3946mg/L

### 1. Det

Anal.: NPOC

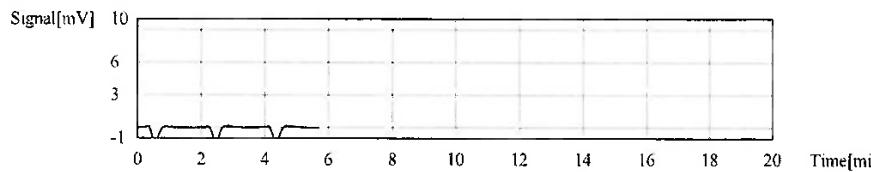
# TOC-Control L Report

Chem36  
d230330wl.toc.flx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	0.000	0.000mg/L	100uL	1.000	E	d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:49:07 PM
2	2.194	0.3734mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:51:14 PM
3	2.443	0.4158mg/L	100uL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/30/2023 4:53:20 PM

Mean Conc.  
CV Conc

0.3946mg/L  
7.59%



9.3

9

LABORATORY REVIEW SIGNATURE FORM  
(To be stored with the raw data)File ID: D23033001.TXT  
Analyst: SSDate Analyzed: 03/20/23  
Run ID: GN39825  
Methods: EPA 300/SW846 9056A

The following analyst(s) have reviewed this run and attest that, to the best of their knowledge, this documentation is complete and correct:

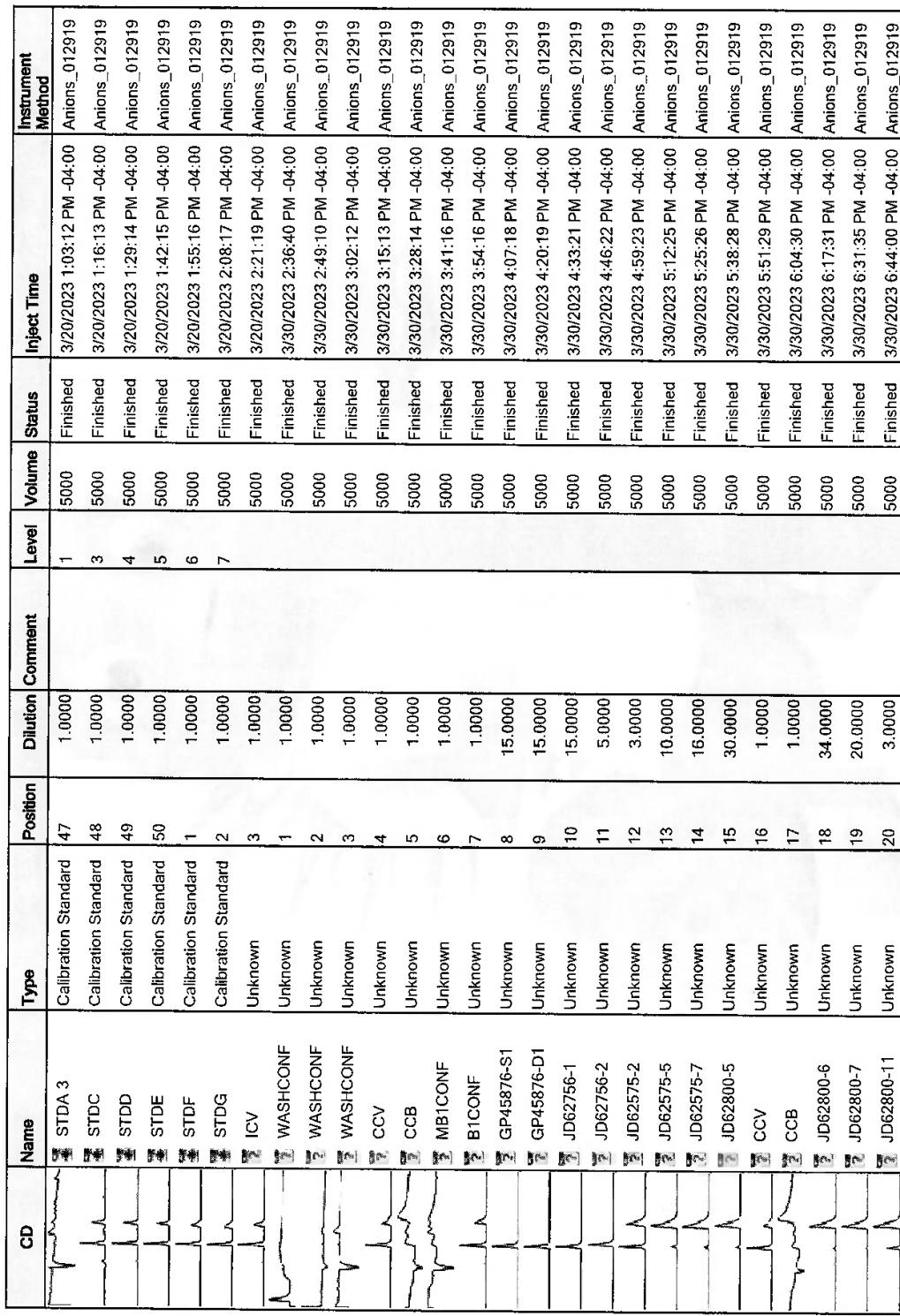
Analyst: JD Date 4/3/23

Analyst: \_\_\_\_\_ Date \_\_\_\_\_

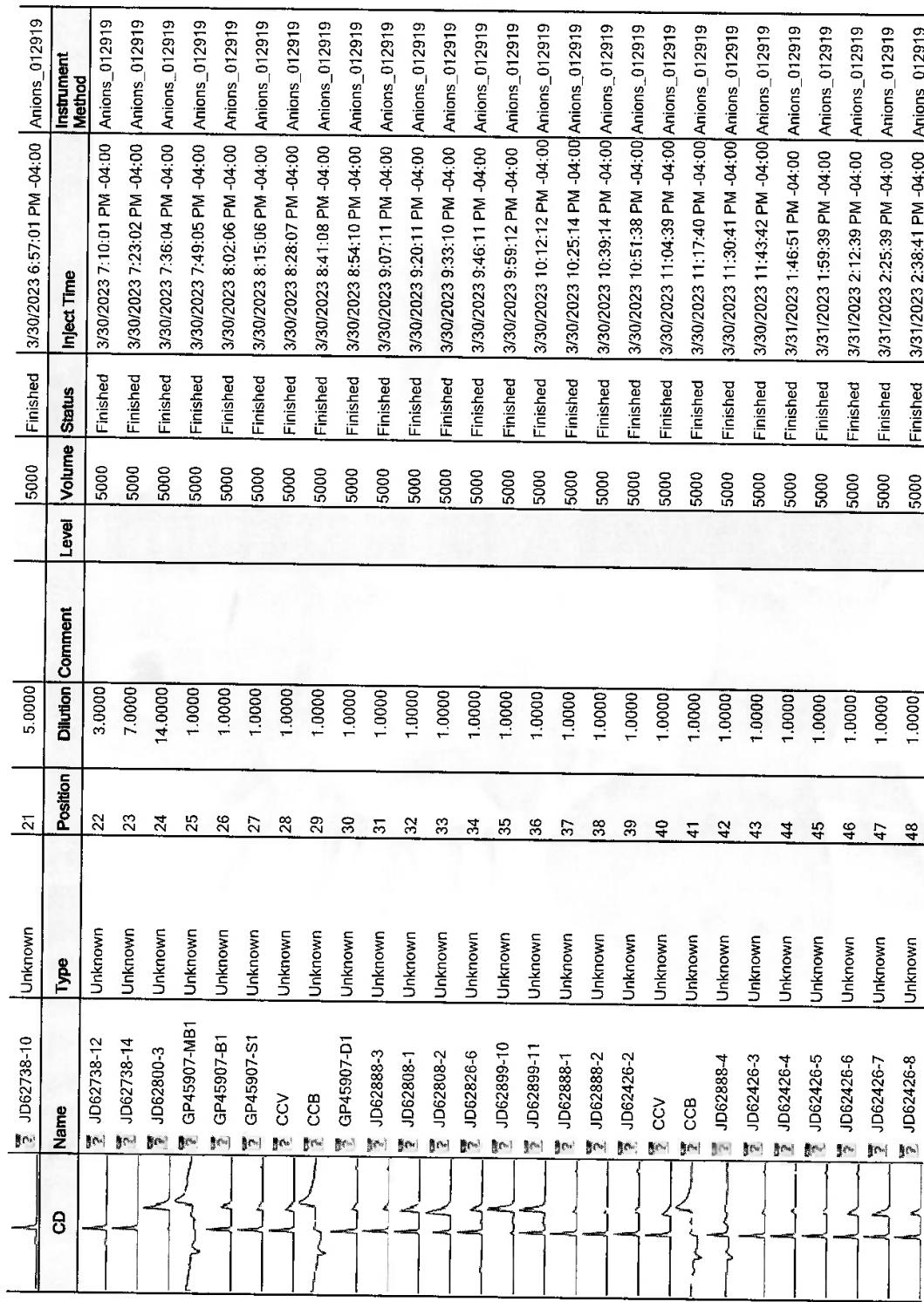
The following supervisor or their designee has reviewed this run and attests that, to the best of their knowledge, this documentation is complete and correct:

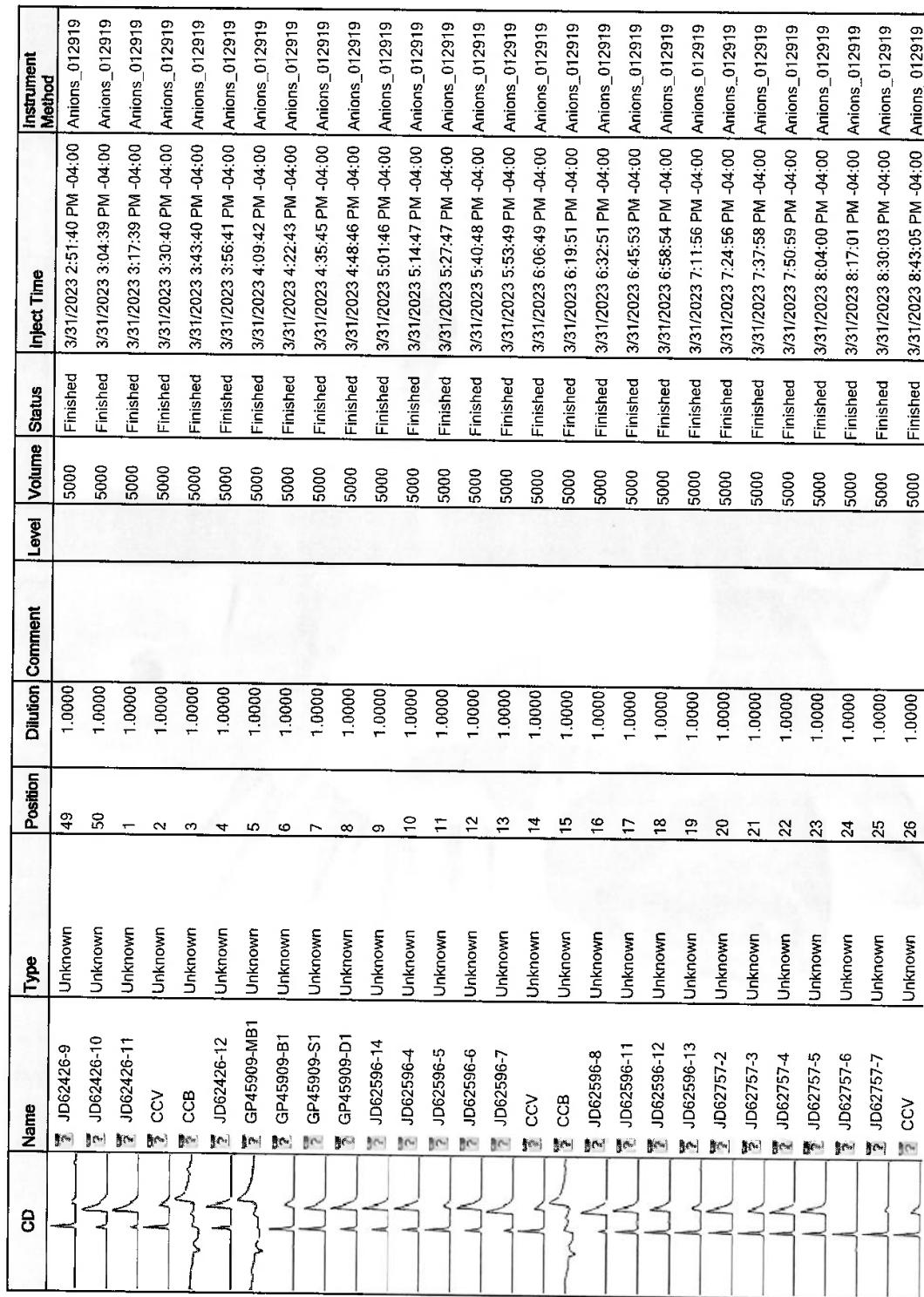
Supervisor (or designee): C Date 4/3/23

9.4 9



Sequence: D20333001  
Last Update Operator: Chemistry





Sequence: D22033001  
 Last Update Operator: Chemistry

CD	Name	Type	Position	Dilution	Comment	Level	Volume	Status	Inject Time	Instrument	Method
	CCB	Unknown	27	1.0000			5000	Finished	3/31/2023 8:56:06 PM -04:00	Anions_012919	
	JD62757-8	Unknown	28	1.0000			5000	Finished	3/31/2023 9:09:08 PM -04:00	Anions_012919	
	JD62757-9	Unknown	29	1.0000			5000	Finished	3/31/2023 9:22:10 PM -04:00	Anions_012919	
	JD62757-10	Unknown	30	1.0000			5000	Finished	3/31/2023 9:35:11 PM -04:00	Anions_012919	
	JD62757-11	Unknown	31	1.0000			5000	Finished	3/31/2023 9:48:13 PM -04:00	Anions_012919	
	JD62757-12	Unknown	32	1.0000			5000	Finished	3/31/2023 10:01:15 PM -04:00	Anions_012919	
	JD62575-7	Unknown	33	32.0000			5000	Finished	3/31/2023 10:14:16 PM -04:00	Anions_012919	
	CCV	Unknown	34	1.0000			5000	Finished	3/31/2023 10:27:17 PM -04:00	Anions_012919	
	CCB	Unknown	35	1.0000			5000	Finished	3/31/2023 10:40:18 PM -04:00	Anions_012919	

Chromeleon 7, Version 7.2.8.10783, Thermo Fisher Scientific

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Printed by Chemistry  
 03/04/23 10:06



## Reagent Information Log - IONC IC<sup>D</sup> - Water /Soil

GN Batch ID#: GN39825

<u>Reagent</u>	<u>Reagent # or Man./Lot</u>	<u>Exp. Date</u>
Standard Stocks (for Calibration Curve) (4000 PPM CHL/SO4)	SCP S220726024	9/30/2023
Standard Stocks (for Calibration Curve) (100 PPM BRO)	SCP S220726023	9/30/2023
Standard Stocks (for Calibration Curve) (100 PPM F)	SCP S220726025	9/30/2023
ICV MIX I (200 PPM CHL/SO4)	SCP S220726030	9/30/2023
ICV MIX II (100 PPM F/BRO)	SCP S221213009	12/31/2023
ICV	GNE3-73533-IC	4/1/2023
Eluent	220950716314	9/30/2024
CCV	GNE3-73534-IC	4/1/2023
Spiking Solution	SCP S220726027	9/30/2023
Filter lot numbers	TISCH 290064158	NA

Form : GN087A-76  
Rev. Date: 4/7/09

4.9

**Anions Standard Preparation Log**

Class A pipet used :  
Autopipet used : 53

		Date	GN Number	3/30/23	C-N39825						
Intermediate Standard Description	Stock Standard used for preparation	Stock concentration	Anion	Stock volume or weight used with units	Diluent	Final Volume	Intermediate (mg/l)	Final Conc. of Standard (mg/l)	Expiration Date	Analyst	Date
GNE3-73526-IC	SPEC S220726024	4000ppm	CH <sub>3</sub> SO <sub>4</sub>	2.5 mL	DI	100 mL	100ppm	4/1/2023	SS	3/1/2023	
GNE3-73525-IC	SPEC S220726025	100ppm	F	20mL	DI	100 mL	20ppm	4/1/2023	SS	3/1/2023	
GNE3-73527-IC	SPEC S220726023	100ppm	BRO	20mL	DI	100 mL	20ppm	4/1/2023	SS	3/1/2023	
ICV	SPEC S220726030	200 ppm	CH <sub>3</sub> SO <sub>4</sub>	50 mL	DI	100 mL	100 ppm	4/1/2023	SS	3/1/2023	
GNE3-73533-IC	SPEC S221213009	100 ppm	F/BRO	3 mL	DI	3 ppm					
Standard Description	Intermediate or Stock used to prepare standard	Intermediate or Stock concentration	Anion	Intermediate or Stock volume used in ml	Diluent	Final Volume	Final Conc. of Standard (mg/l)	Expiration Date	Analyst	Date	
STD C	GNE3-73526-IC	100ppm	CH <sub>3</sub> SO <sub>4</sub>	2.0ml	DI	100 mL	0.5ppm	4/1/2023	SS	3/1/2023	
STD D	GNE3-73528-IC	100ppm	F	0.5mL	DI	100 mL	0.2ppm	4/1/2023	SS	3/1/2023	
STD E	GNE3-73529-IC	100ppm	CH <sub>3</sub> SO <sub>4</sub>	10mL	DI	100 mL	1ppm	4/1/2023	SS	3/1/2023	
STD F	GNE3-73530-IC	100ppm	BRO	1mL	DI	100 mL	0.4ppm	4/1/2023	SS	3/1/2023	
STD G	GNE3-73531-IC	100ppm	CH <sub>3</sub> SO <sub>4</sub>	50mL	DI	100 mL	50ppm	4/1/2023	SS	3/1/2023	
GNE3-73532-IC	GNE3-73525-IC	20ppm	F	5mL	DI	100 mL	1ppm	4/1/2023	SS	3/1/2023	
CCV	SPEC S220726024	4000ppm	CH <sub>3</sub> SO <sub>4</sub>	10mL	DI	200 mL	200ppm	4/1/2023	SS	3/1/2023	
GNE3-73534-IC	SPEC S220726023	100ppm	BRO	6mL	DI	200 mL	300ppm	4/1/2023	SS	3/1/2023	

\* If Class A glass pipets are used, enter an A. For balances or autopipets, then enter the appropriate SGS ID number.

NO2/NC3/P4 included in curve but not reported

Form: GN318-01

Rev Date: 6/23/2020



Analyst: SS

Product: ANIONS

Pipette ID: \_\_\_\_\_ 48 \_\_\_\_\_

Prep Date: 3/31/2023

Balance ID: \_\_\_\_\_

Batch ID: GN39825

## Sample Prep Log

**Reviewer:**

Date:

Form: GN166-04  
Revised: 10/27/16



Analyst JI

Analyst JD Product Anions Autopipette # 43

Date 3/20/23

Product Anons

Autopipette # 43

Date 3/20/23

### Class A Vol. Flask

## Sample Dilution Prep Log

9.4

QC Reviewer: \_\_\_\_\_ Date: \_\_\_\_\_

Date:

Form: GN165-01  
Rev. Date: 2/25/03

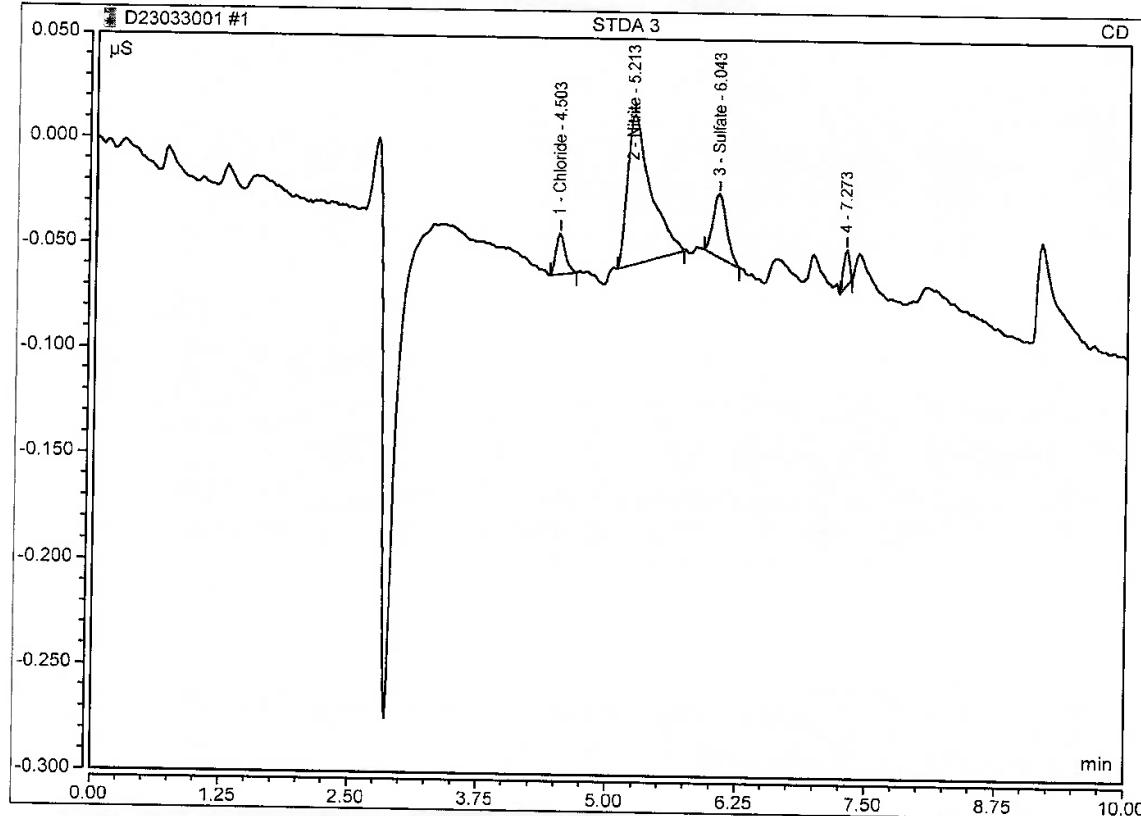
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

Page 1 of 92  
 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	STDA 3	Inj. Vol.:	5000.00
Injection Type:	Calibration Standard	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	20-Mar-2023 / 13:03	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	4.50	Chloride	BMB	0.002	0.020	0.2462
2	5.21	Nitrite	BMB	0.019	0.072	0.0374
3	6.04	Sulfate	BMB	0.005	0.030	0.1273
<b>TOTAL:</b>				0.03	0.12	0.41



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

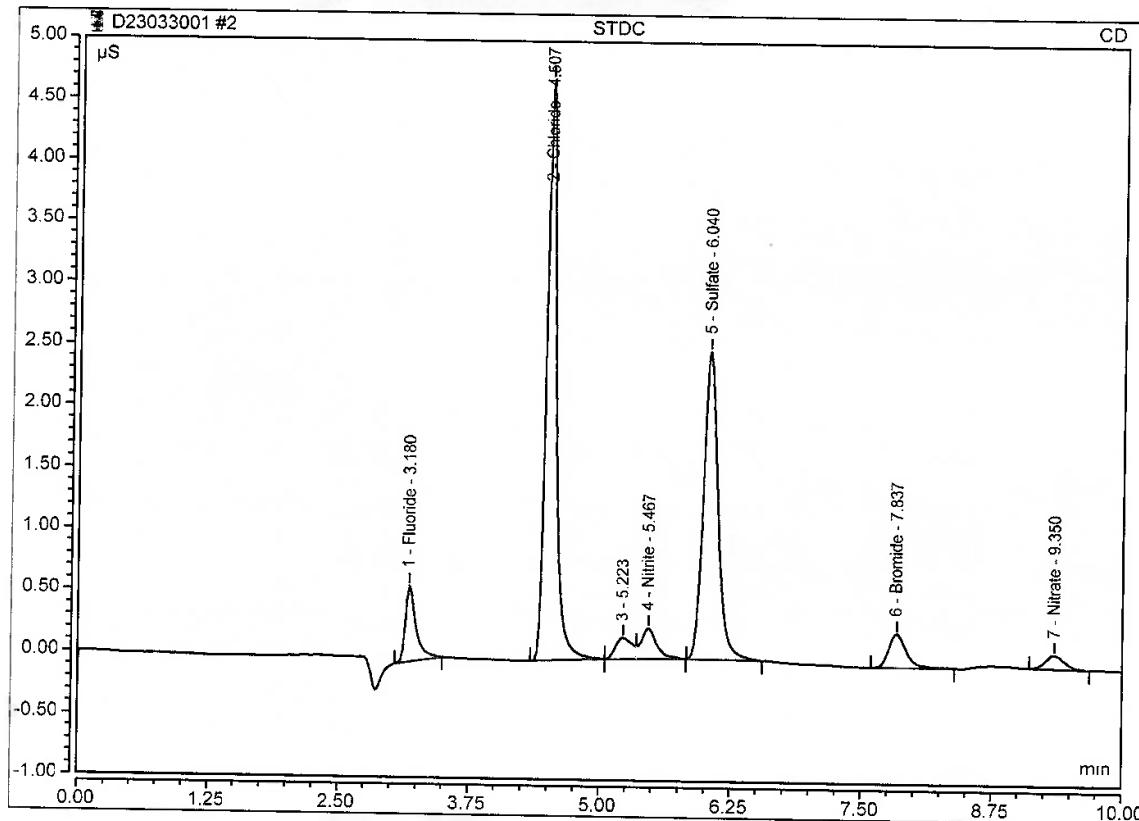
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

Page 2 of 92  
 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	STDC	Inj. Vol.:	5000.00
Injection Type:	Calibration Standard	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	20-Mar-2023 / 13:16	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.18	Fluoride	BMB	0.072	0.614	0.2117
2	4.51	Chloride	BMB	0.531	4.724	1.9116
4	5.47	Nitrite	MB	0.043	0.249	0.2009
5	6.04	Sulfate	BMB	0.407	2.497	1.9051
6	7.84	Bromide	BMB	0.054	0.273	0.5404
7	9.35	Nitrate	BMB	0.023	0.110	0.1311
<b>TOTAL:</b>				1.13	8.47	4.90



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

9.4

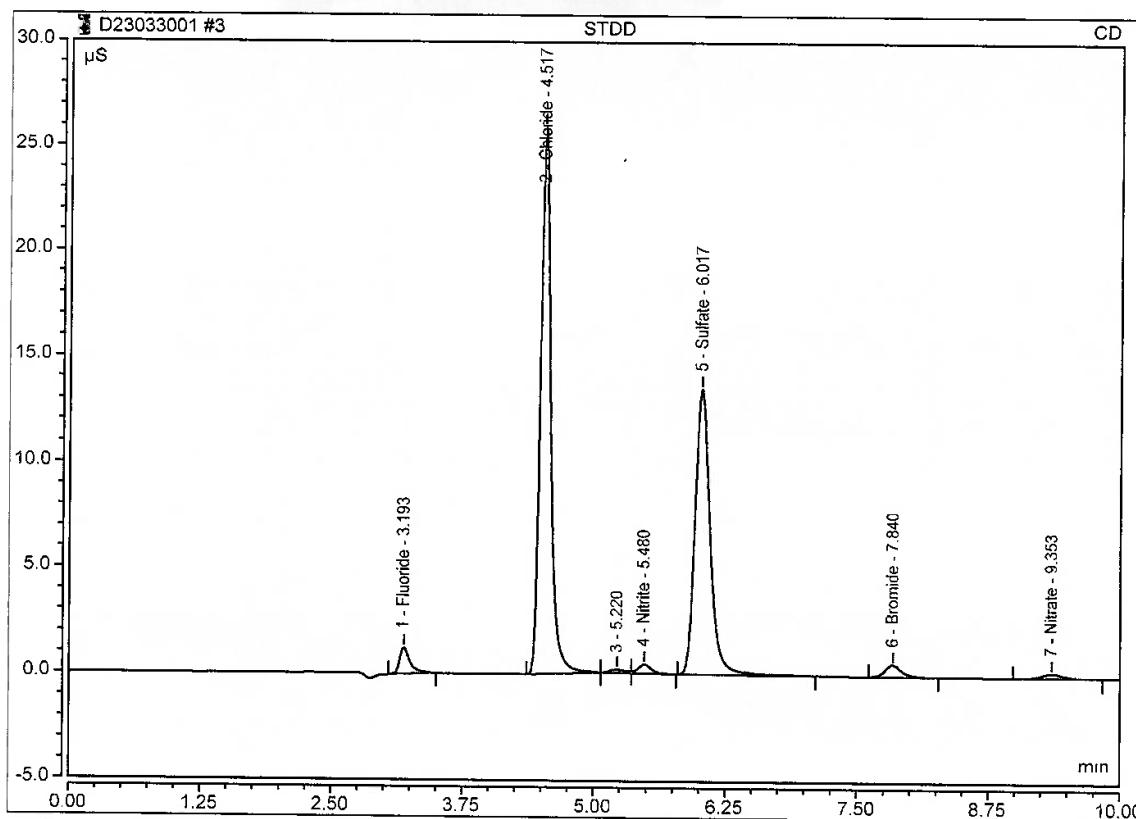
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 Instrument: Integron\_1  
 Sequence: D23033001

Page 3 of 92  
 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	STDD	Inj. Vol.:	5000.00
Injection Type:	Calibration Standard	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	20-Mar-2023 / 13:29	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB	0.142	1.247	0.3807
2	4.52	Chloride	BMB	2.952	25.959	9.5331
4	5.48	Nitrite	MB	0.063	0.436	0.3441
5	6.02	Sulfate	BMB	2.229	13.488	9.9420
6	7.84	Bromide	BMB	0.102	0.561	0.9520
7	9.35	Nitrate	BMB	0.053	0.222	0.3856
<b>TOTAL:</b>				5.54	41.91	21.54



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

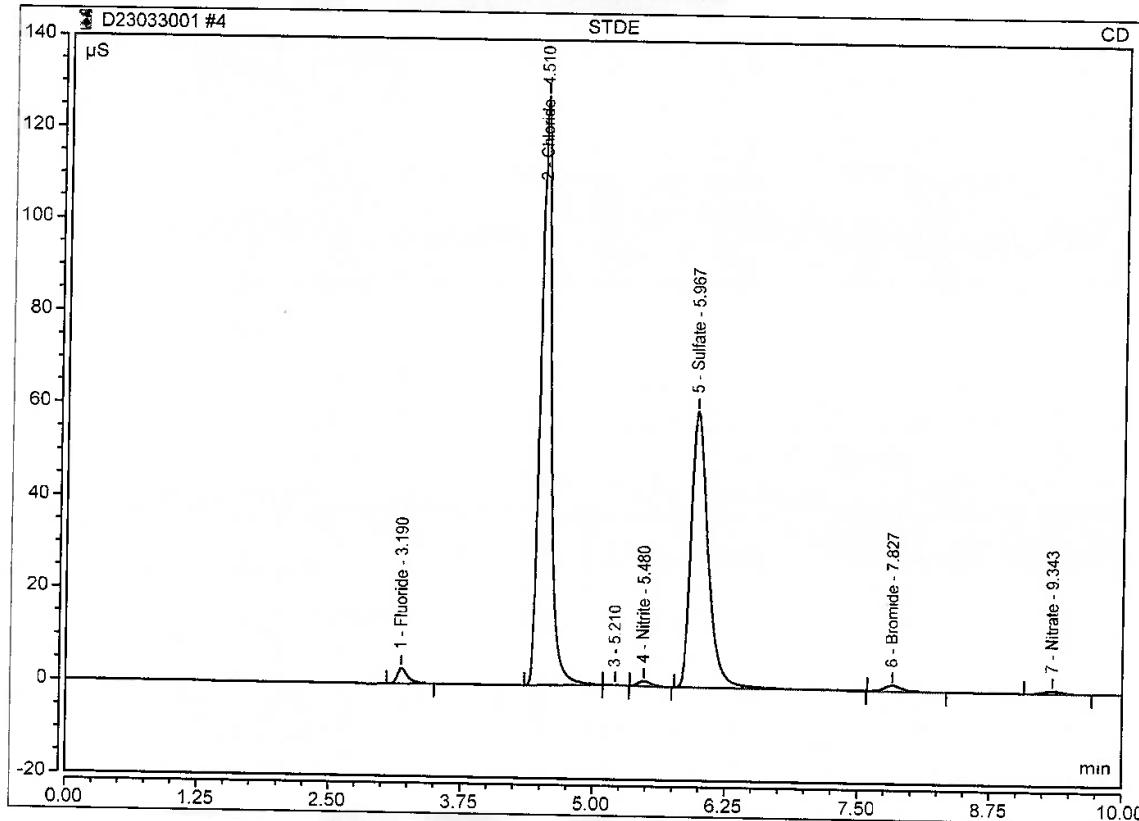
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

Page 4 of 92  
 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	STDE	Inj. Vol.:	5000.00
Injection Type:	Calibration Standard	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	20-Mar-2023 / 13:42	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB	0.367	3.280	0.9240
2	4.51	Chloride	BMb	15.177	127.825	48.0197
4	5.48	Nitrite	BMB	0.137	1.065	0.8460
5	5.97	Sulfate	BMB	11.089	59.673	49.0404
6	7.83	Bromide	BMB	0.215	1.178	1.9155
7	9.34	Nitrate	BMB	0.113	0.522	0.8930
<b>TOTAL:</b>				27.10	193.54	101.64



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

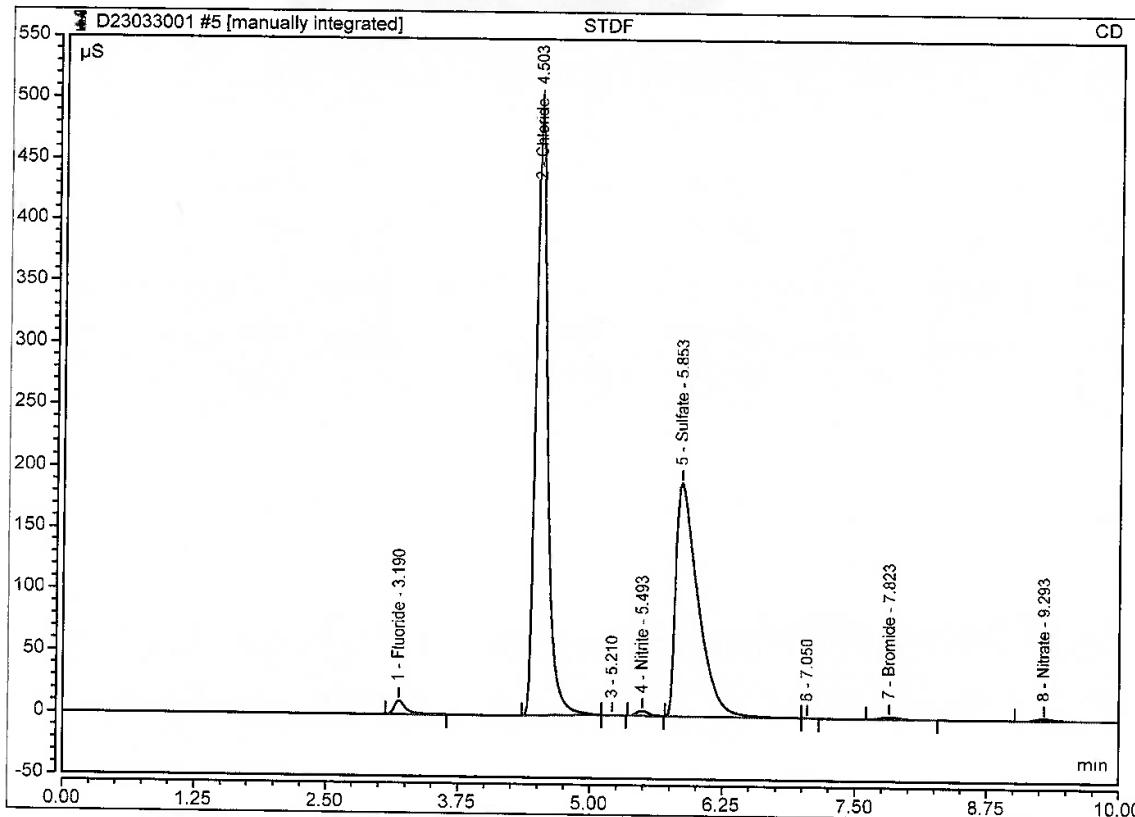
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 Sequence: D23033001

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 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	STDF	Inj. Vol.:	5000.00
Injection Type:	Calibration Standard	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Ini. Date / Time:	20-Mar-2023 / 13:55	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB*	1.353	11.060	3.3103
2	4.50	Chloride	BMb	63.586	494.275	200.4161
4	5.49	Nitrite	BMB	0.512	4.013	3.4034
5	5.85	Sulfate	BMb	45.842	189.675	202.3956
7	7.82	Bromide	BMB	0.333	1.853	2.9320
8	9.29	Nitrate	BMB	0.432	1.928	3.5834
<b>TOTAL:</b>				112.06	702.80	416.04



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

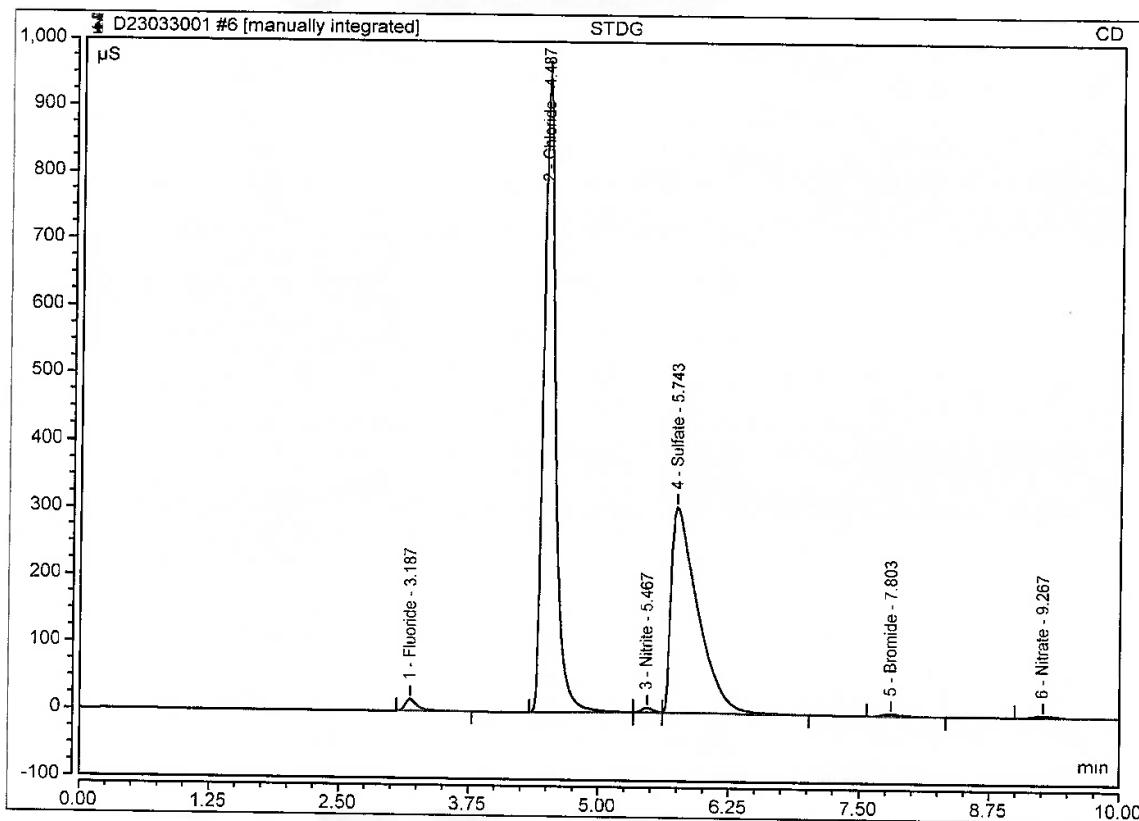
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 Sequence: D23033001

Page 6 of 92  
 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	STDG	Inj. Vol.:	5000.00
Injection Type:	Calibration Standard	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	20-Mar-2023 / 14:08	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB*	2.372	17.718	5.7733
2	4.49	Chloride	BMb	127.656	951.102	402.1195
3	5.47	Nitrite	bM	0.863	7.083	5.8057
4	5.74	Sulfate	MB	90.332	306.943	398.7169
5	7.80	Bromide	BMB	0.594	3.245	5.1600
6	9.27	Nitrate	BMB	0.778	3.419	6.5070
<b>TOTAL:</b>				222.59	1289.51	824.08



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

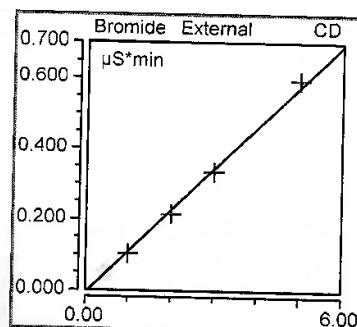
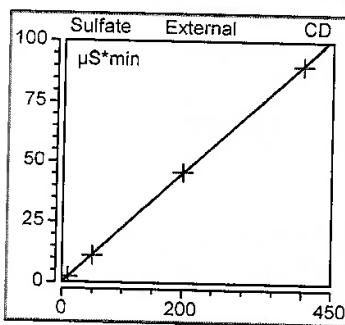
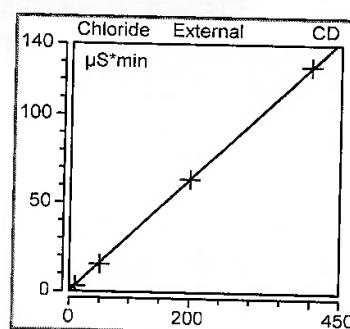
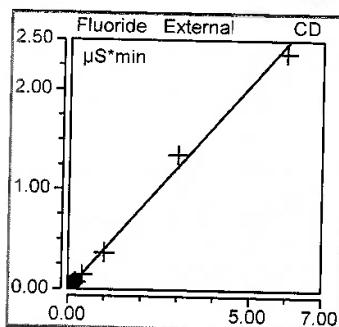
9.4  
9

## Calibration Batch Report

Sequence: D23033001	Injection Volu 5,000.00
Instrument M Anions_012 919	Operator: Chemistry
Inj. Date / Tin 20/03/2023 / 13:16	Run Time: 10

## ibration Summary

Peak Name	Eval.Type	Cal.Type	Points	Offset (C0)	Slope (C1)	Curve (C2)	Coeff.Det. %
Fluoride	Area	, WithOffset,	5.000	-0.016	0.413	0.000	99.3896
Chloride	Area	, WithOffset,	6.000	-0.076	0.318	0.000	99.9771
Nitrite	Area	, WithOffset,	6.000	0.013	0.146	0.000	98.8884
Sulfate	Area	, WithOffset,	6.000	-0.024	0.227	0.000	99.9900
Bromide	Area	, WithOffset,	5.000	-0.009	0.117	0.000	99.7028
Nitrate	Area	, WithOffset,	5.000	0.008	0.118	0.000	92.5272
AVERAGE:				-0.0174	0.2232	0.0000	98.4125

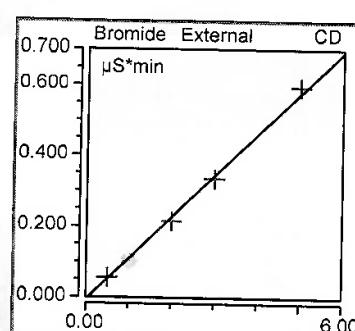
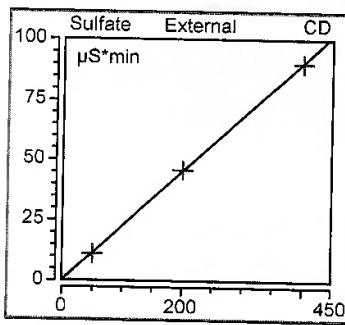
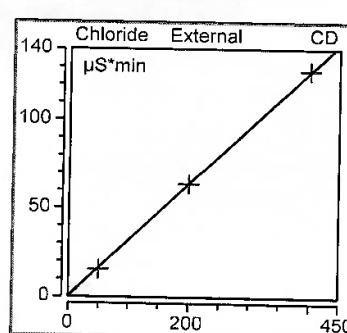
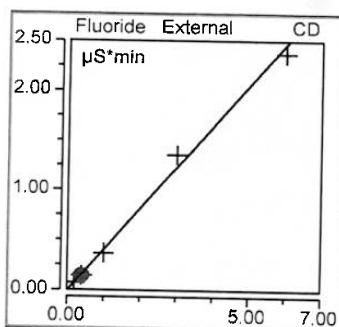


## Calibration Batch Report

Sequence: D23033001	Injection Volu 5,000.00
Instrument M Anlons_012 919	Operator: Chemistry
Inj. Date / Tin 20/03/2023 / 13:29	Run Time: 10

## bration Summary

Peak Name	Eval.Type	Cal.Type	Points	Offset (C0)	Slope (C1)	Curve (C2)	Coeff.Det. %
Fluoride	Area	, WithOffset,	5.000	-0.016	0.413	0.000	99.3896
Chloride	Area	, WithOffset,	6.000	-0.076	0.318	0.000	99.9771
Nitrite	Area	, WithOffset,	6.000	0.013	0.146	0.000	98.8884
Sulfate	Area	, WithOffset,	6.000	-0.024	0.227	0.000	99.9900
Bromide	Area	, WithOffset,	5.000	-0.009	0.117	0.000	99.7028
Nitrate	Area	, WithOffset,	5.000	0.008	0.118	0.000	92.5272
AVERAGE:				-0.0174	0.2232	0.0000	98.4125

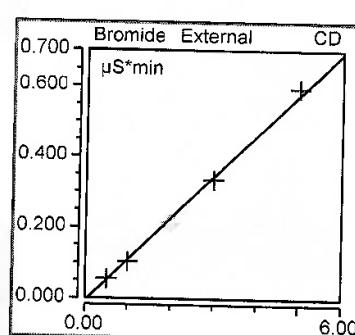
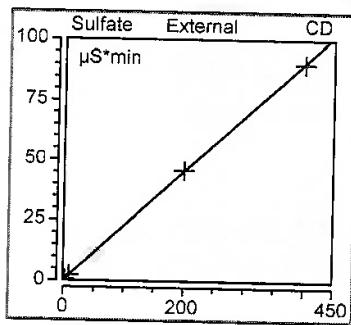
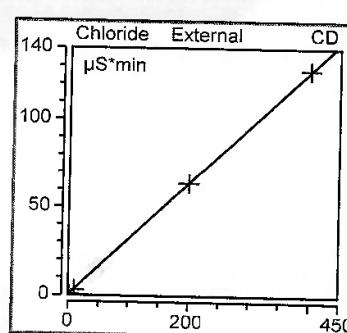
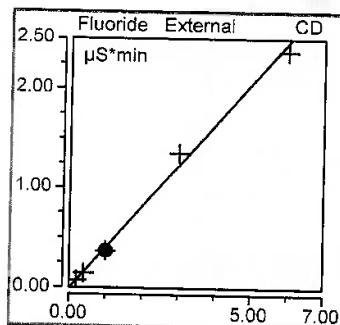


## Calibration Batch Report

Sequence: D23033001	Injection Volu 5,000.00
Instrument M Anions_012 919	Operator: Chemistry
Inj. Date / Tin 20/03/2023 / 13:42	Run Time: 10

## ibration Summary

Peak Name	Eval.Type	Cal.Type	Points	Offset (C0)	Slope (C1)	Curve (C2)	Coeff.Det. %
Fluoride	Area	, WithOffset,	5.000	-0.016	0.413	0.000	99.3096
Chloride	Area	, WithOffset,	6.000	-0.076	0.318	0.000	99.9771
Nitrite	Area	, WithOffset,	6.000	0.013	0.146	0.000	98.8884
Sulfate	Area	, WithOffset,	6.000	-0.024	0.227	0.000	99.9900
Bromide	Area	, WithOffset,	5.000	-0.009	0.117	0.000	99.7028
Nitrate	Area	, WithOffset,	5.000	0.008	0.118	0.000	92.5272
<b>AVERAGE:</b>				-0.0174	0.2232	0.0000	98.4125

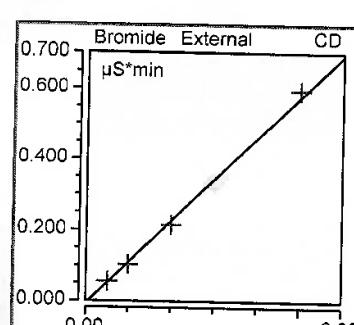
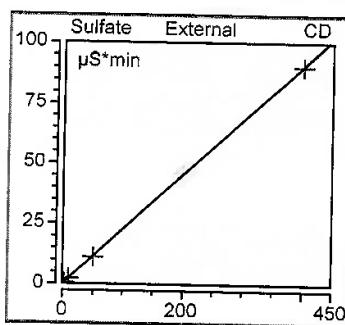
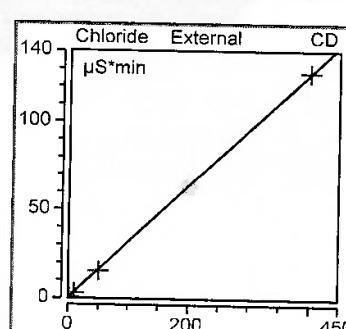
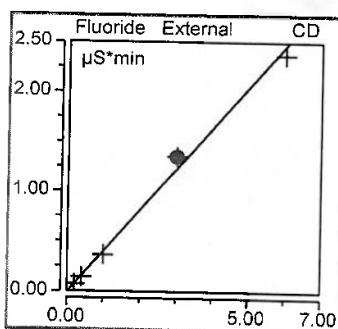


## Calibration Batch Report

Sequence: D23033001	Injection Volu 5,000.00
Instrument M Anions_012 919	Operator: Chemistry
Inj. Date / Tin 20/03/2023 / 13:55	Run Time: 10

## ibration Summary

Peak Name	Eval.Type	Cal.Type	Points	Offset (C0)	Slope (C1)	Curve (C2)	Coeff.Det. %
Fluoride	Area	, WithOffset,	5.000	-0.016	0.413	0.000	99.3896
Chloride	Area	, WithOffset,	6.000	-0.076	0.318	0.000	99.9771
Nitrite	Area	, WithOffset,	6.000	0.013	0.146	0.000	98.8884
Sulfate	Area	, WithOffset,	6.000	-0.024	0.227	0.000	99.9900
Bromide	Area	, WithOffset,	5.000	-0.009	0.117	0.000	99.7028
Nitrate	Area	, WithOffset,	5.000	0.008	0.118	0.000	92.5272
<b>AVERAGE:</b>				-0.0174	0.2232	0.0000	98.4125

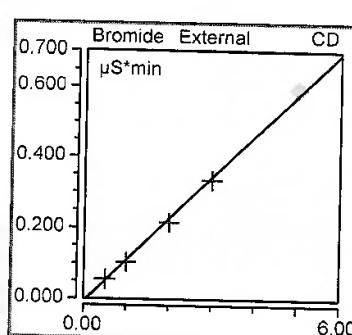
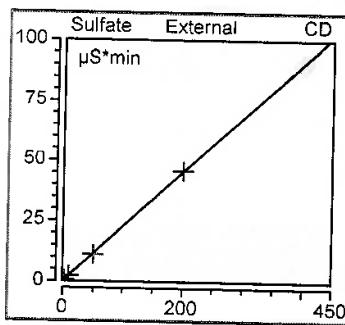
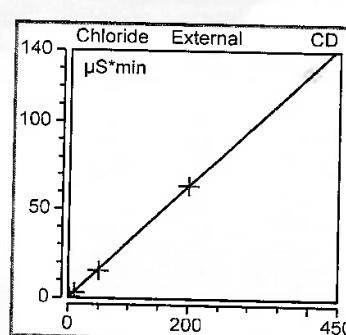
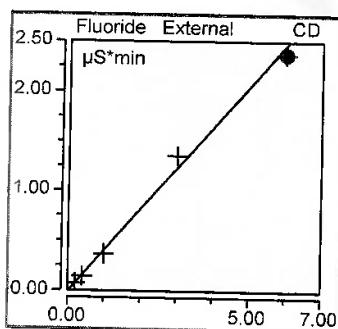


## Calibration Batch Report

Sequence: D23033001	Injection Volu 5,000.00
Instrument M Anions_012 919	Operator: Chemistry
Inj. Date / Tin 20/03/2023 / 14:08	Run Time: 10

## ibration Summary

Peak Name	Eval.Type	Cal.Type	Points	Offset (C0)	Slope (C1)	Curve (C2)	Coeff.Det. %
Fluoride	Area	, WithOffset,	5.000	-0.016	0.413	0.000	99.3896
Chloride	Area	, WithOffset,	6.000	-0.076	0.318	0.000	99.9771
Nitrite	Area	, WithOffset,	6.000	0.013	0.146	0.000	98.8884
Sulfate	Area	, WithOffset,	6.000	-0.024	0.227	0.000	99.9900
Bromide	Area	, WithOffset,	5.000	-0.009	0.117	0.000	99.7028
Nitrate	Area	, WithOffset,	5.000	0.008	0.118	0.000	92.5272
<b>AVERAGE:</b>				-0.0174	0.2232	0.0000	98.4125



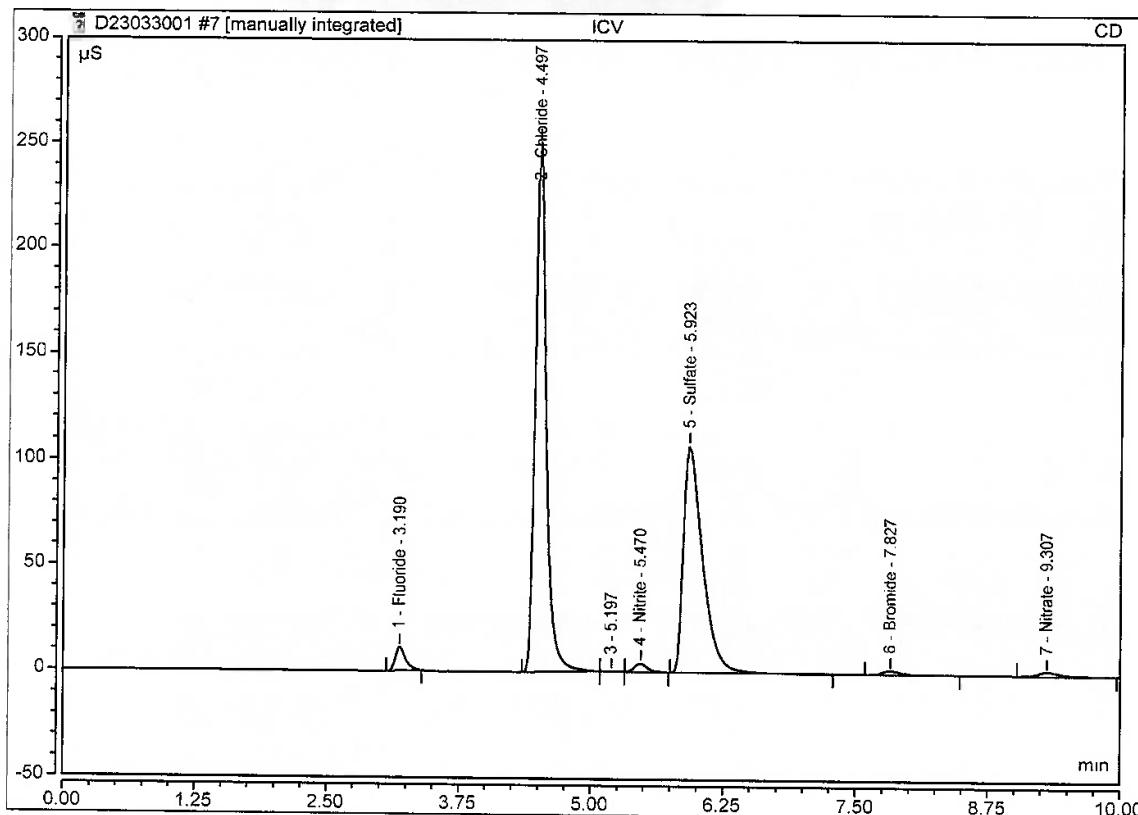
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	ICV	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	20-Mar-2023 / 14:21	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\cdot}\text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB*	1.233	11.224	3.0206
2	4.50	Chloride	BMB	30.784	251.670	97.1516
4	5.47	Nitrite	BMB	0.541	4.014	3.6015
5	5.92	Sulfate	BMB	22.642	107.490	100.0221
6	7.83	Bromide	BMB	0.351	1.906	3.0819
7	9.31	Nitrate	BMB	0.440	1.984	3.6532
<b>TOTAL:</b>				55.99	378.29	210.53



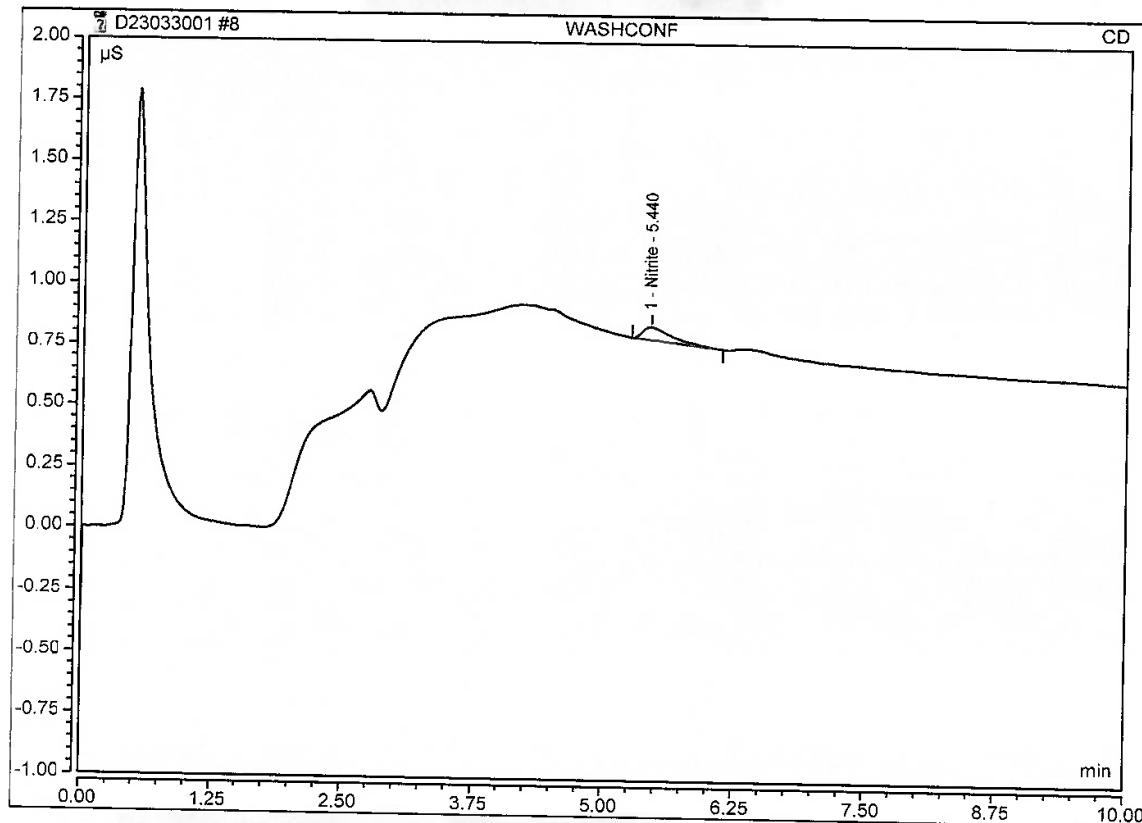
Logged on User: Chemistry  
Instrument: Integriion\_1  
Sequence: D23033001

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## Peak Integration Report

Sample Name:	WASHCONF	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 14:36	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	5.44	Nitrite	BMB	0.017	0.053	0.0270
		TOTAL:		0.02	0.05	0.03

9.4  
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Anion/Integration

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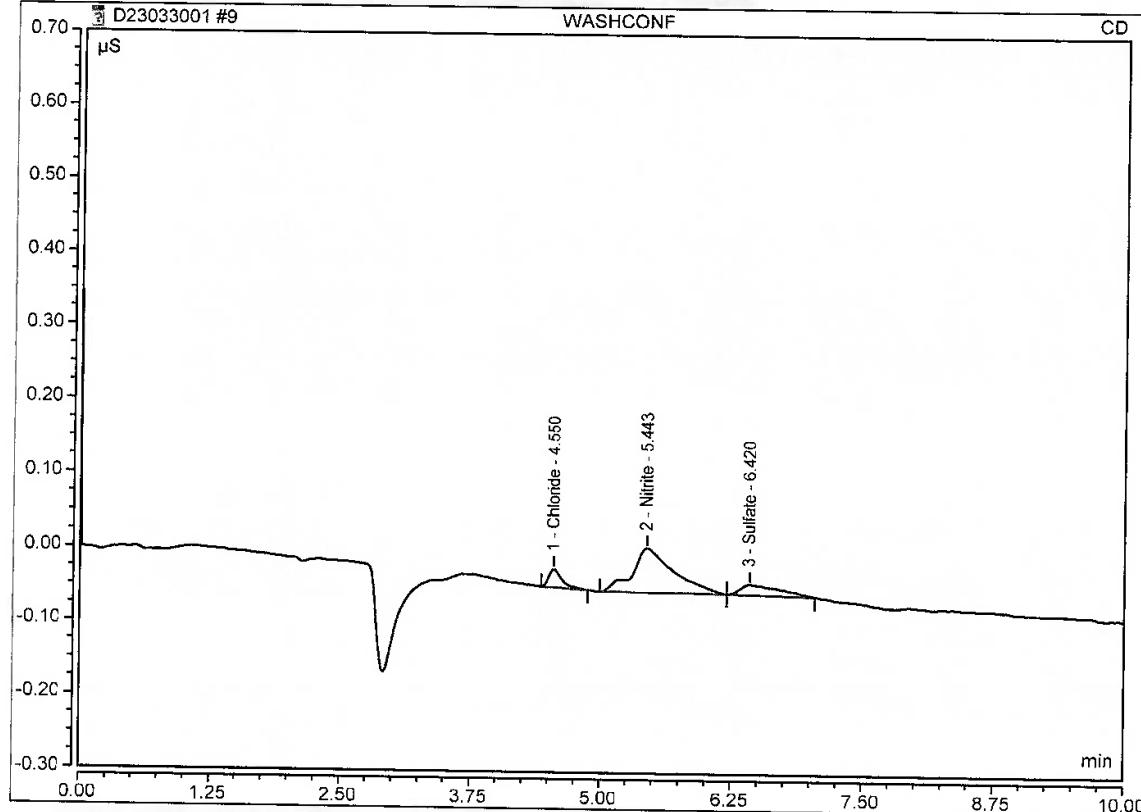
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 Instrument: Integriion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	WASHCONF	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 14:49	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	4.55	Chloride	BMB	0.004	0.025	
2	5.44	Nitrite	BMb	0.028	0.061	
3	6.42	Sulfate	bMB	0.006	0.014	
<b>TOTAL:</b>				0.04	0.10	0.49



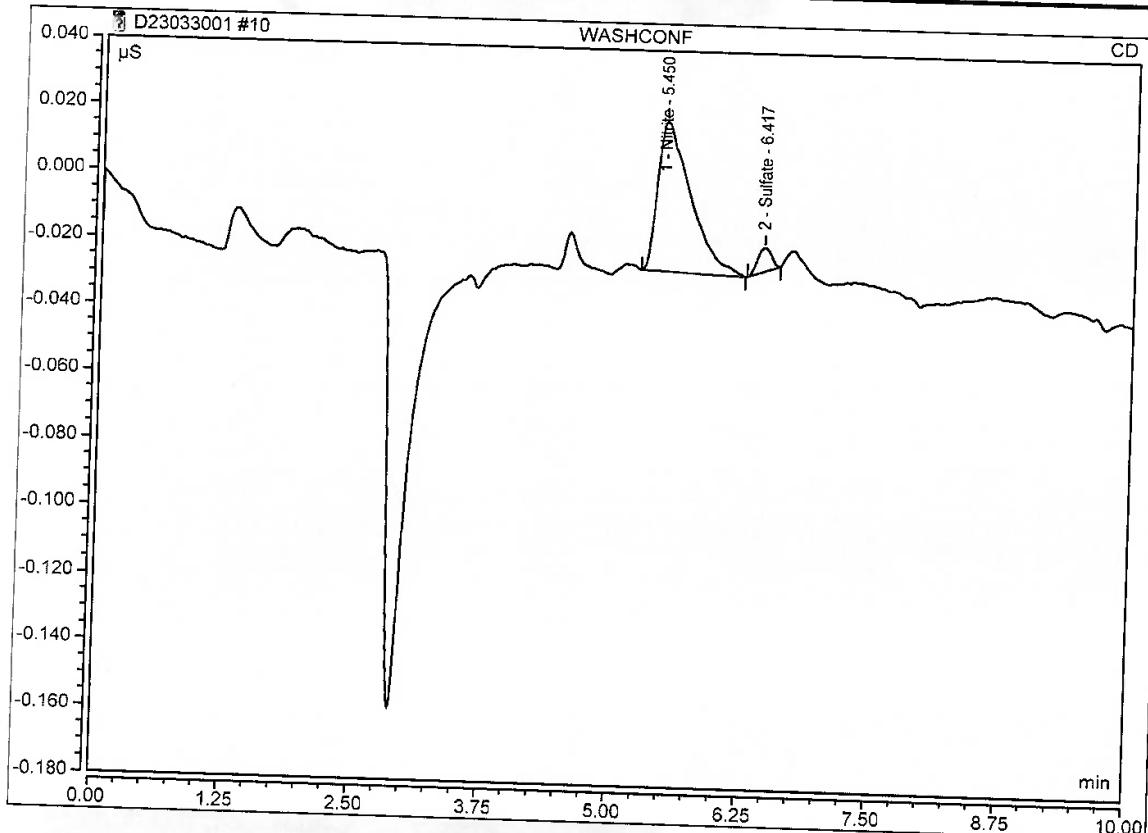
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	WASHCONF	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 15:02	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}/\text{min}$	Height $\mu\text{S}$	Amount
1	5.45	Nitrite	BMB	0.017	0.045	
2	6.42	Sulfate	BMB	0.001	0.007	
		TOTAL:		0.02	0.05	0.14



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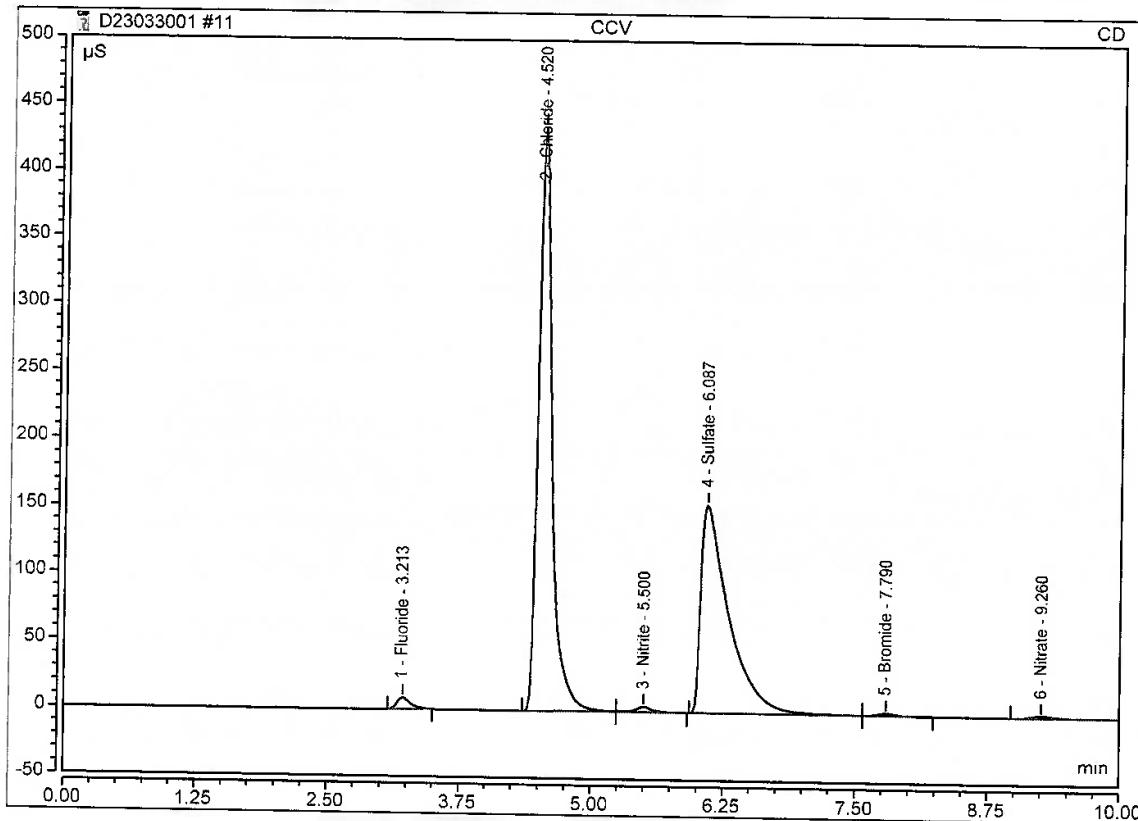
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCV	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 15:15	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	1.255	8.554	3.0728
2	4.52	Chloride	BMb	65.236	435.198	205.6109
3	5.50	Nitrite	bMB	0.626	3.784	4.1825
4	6.09	Sulfate	BMb	47.356	154.580	209.0746
5	7.79	Bromide	bMB	0.323	1.658	2.8433
6	9.26	Nitrate	BMB	0.425	1.700	3.5284
<b>TOTAL:</b>				115.22	605.47	428.31



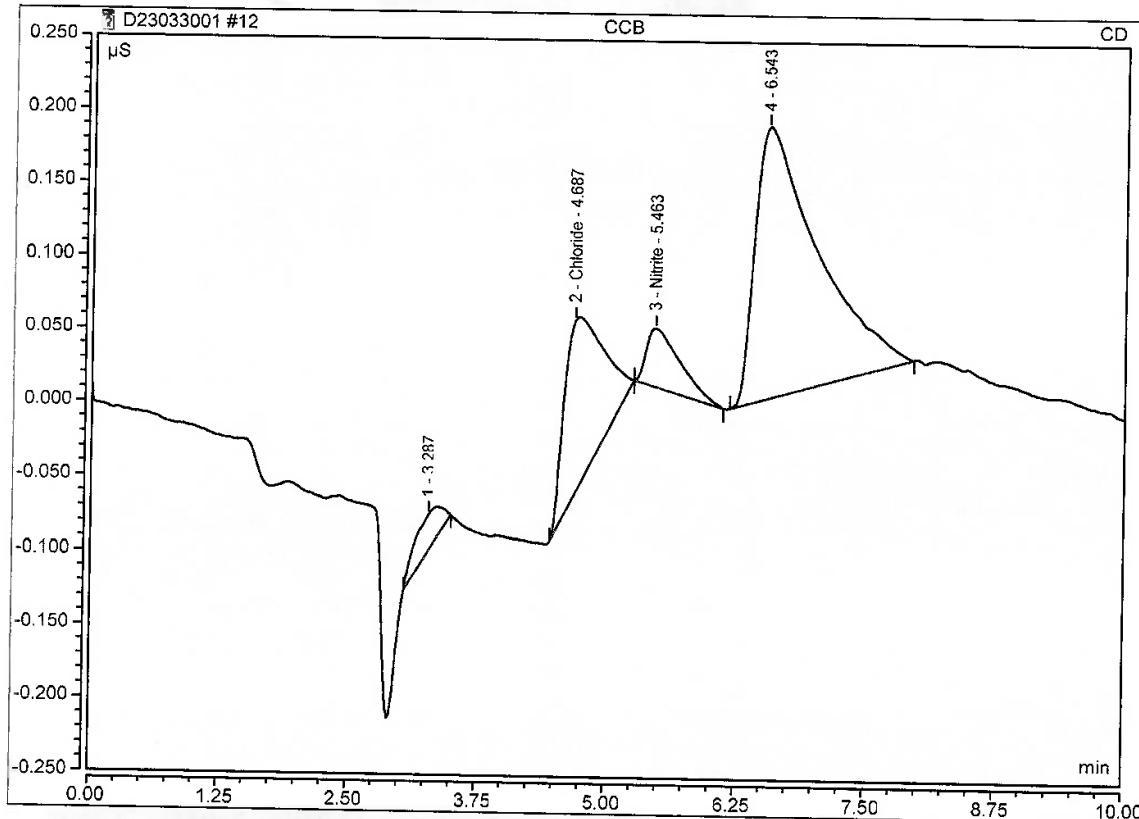
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 15:28	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.69	Chloride	BMb	0.049	0.118	0.3926
3	5.46	Nitrite	bMB	0.015	0.040	0.0149
		TOTAL:		0.06	0.16	0.41



Anion/Integration

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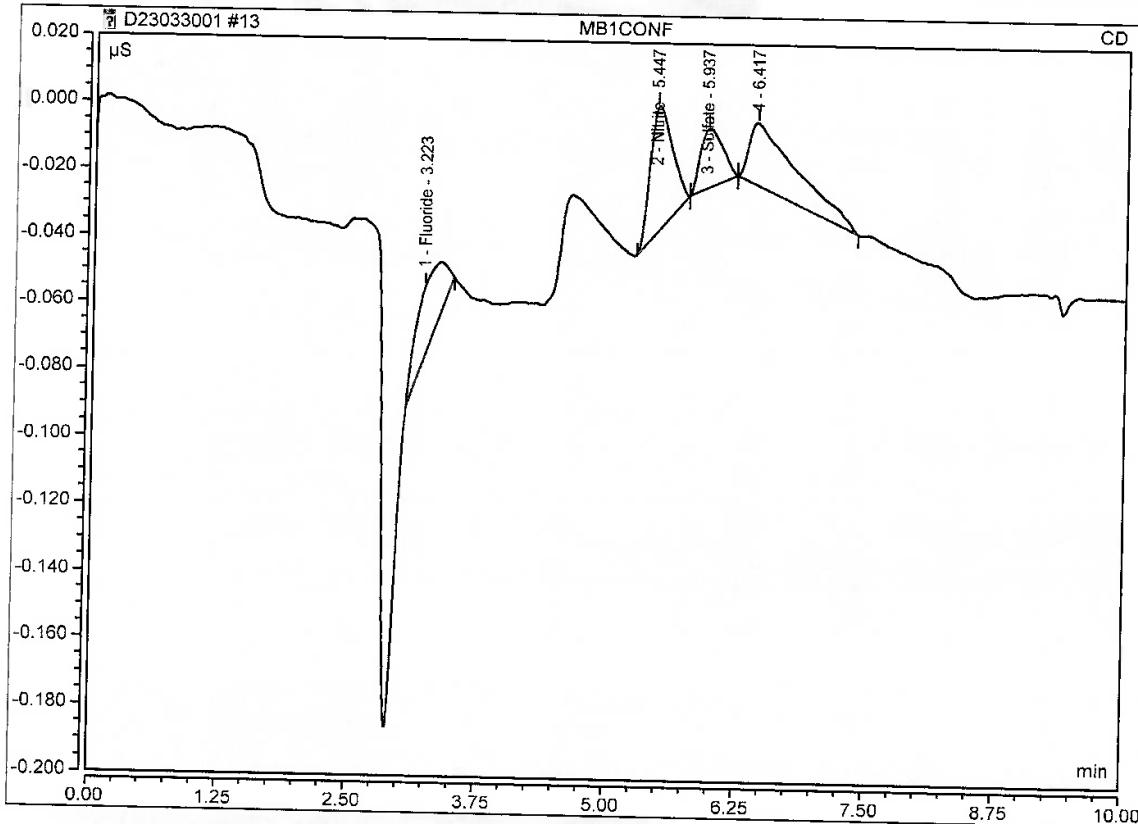
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	MB1CONF	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Ini. Date / Time:	30-Mar-2023 / 15:41	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.006	0.021	
2	5.45	Nitrite	BMb	0.009	0.038	
3	5.94	Sulfate	bMb	0.004	0.018	
<b>TOTAL:</b>				0.02	0.08	0.18



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Anion/Integration

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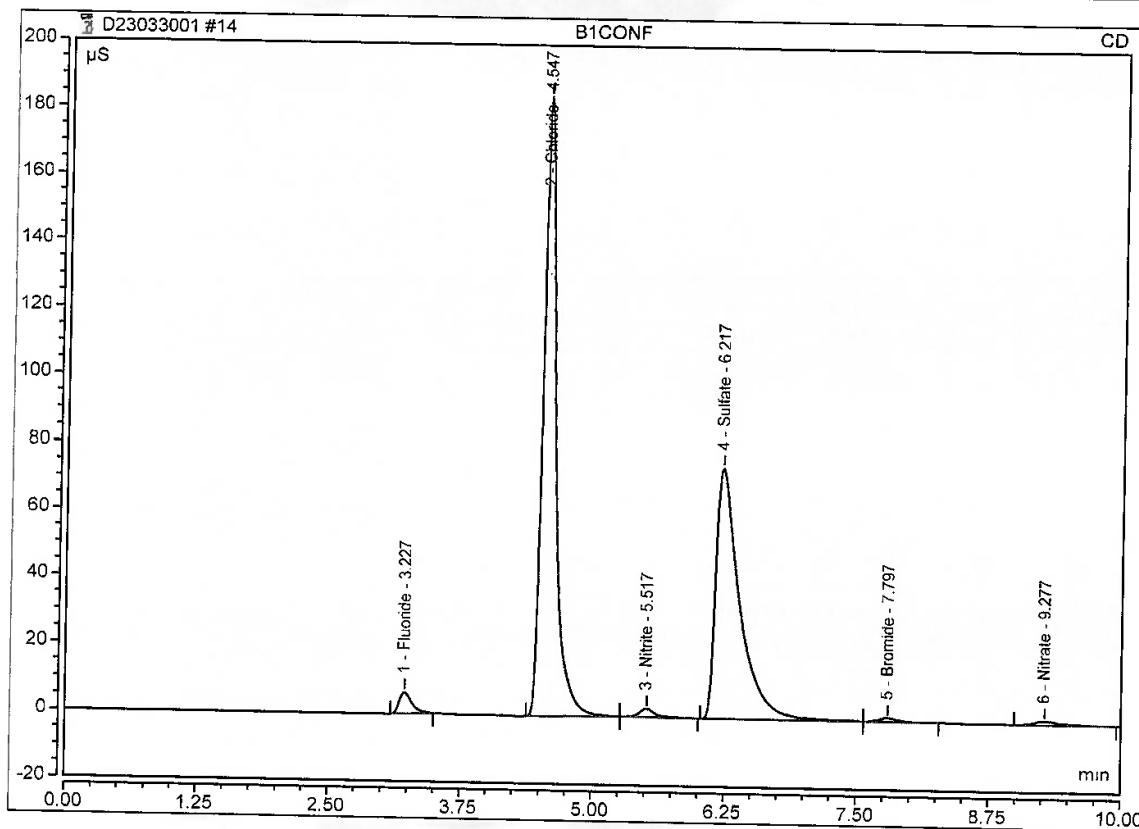
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	B1CONF	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Ini. Date / Time:	30-Mar-2023 / 15:54	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.23	Fluoride	BMB	0.865	6.296	
2	4.55	Chloride	BMb	26.519	184.429	
3	5.52	Nitrite	bMB	0.424	2.501	
4	6.22	Sulfate	BMb	19.333	74.783	
5	7.80	Bromide	bMB	0.211	1.096	
6	9.28	Nitrate	BMB	0.280	1.111	
<b>TOTAL:</b>				47.63	270.22	178.26



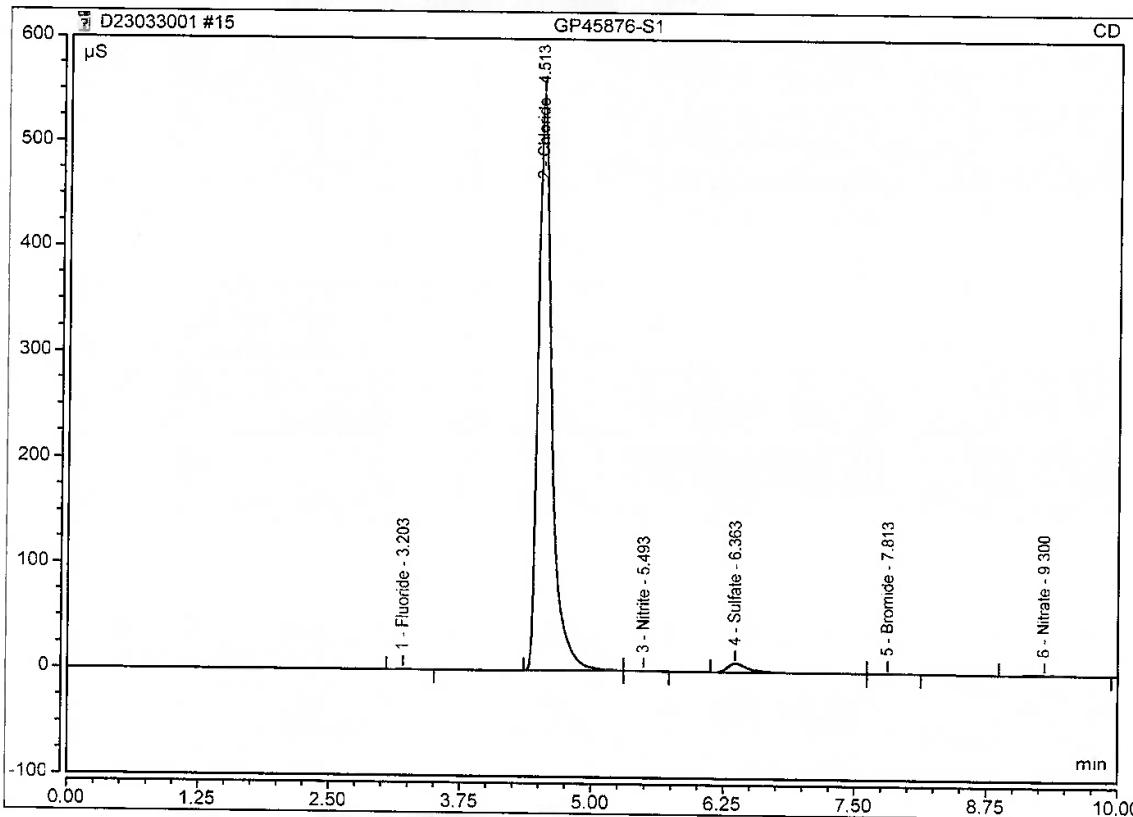
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45876-S1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	15.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 16:07	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.20	Fluoride	BMB	0.067	0.441	2.9956
2	4.51	Chloride	BMb	83.090	547.294	3927.2653
3	5.49	Nitrite	bMB	0.049	0.299	3.6841
4	6.36	Sulfate	BMb	2.063	8.600	138.1509
5	7.81	Bromide	bMB	0.017	0.091	3.3934
6	9.30	Nitrate	BMB	0.122	0.477	14.4725
<b>TOTAL:</b>				85.41	557.20	4089.96



Anion/Integration

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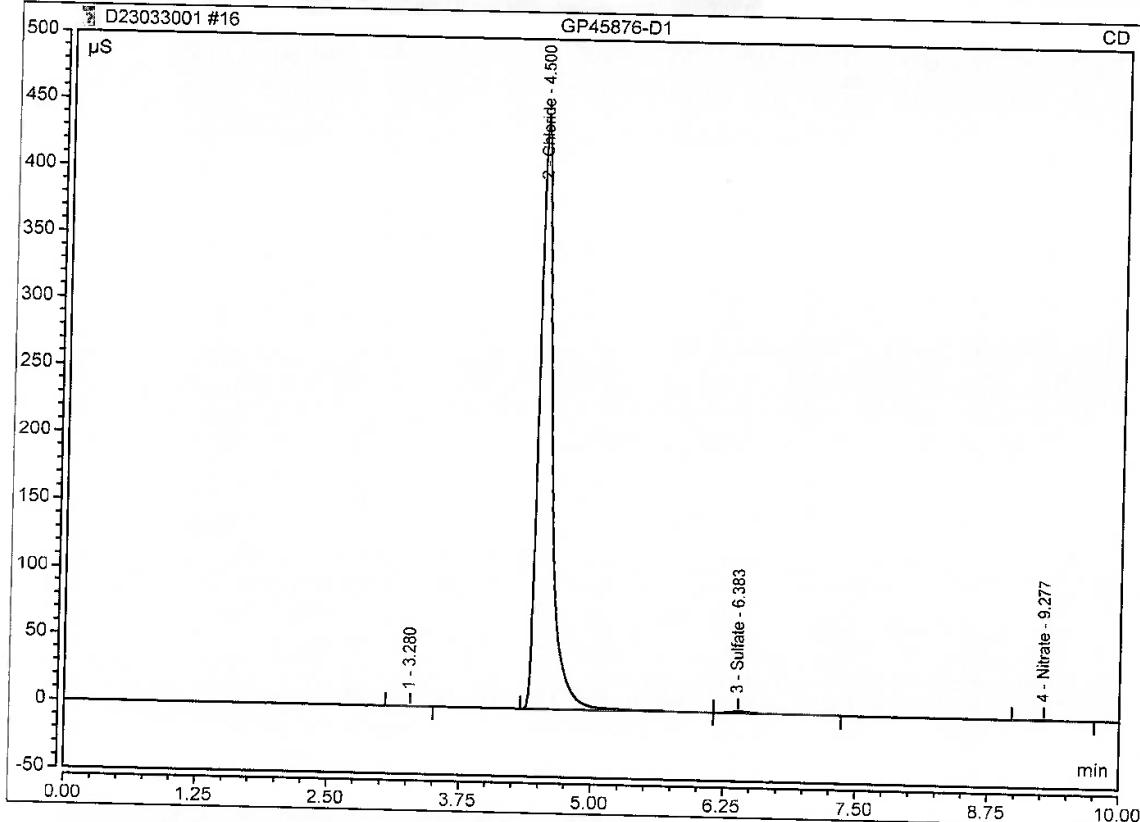
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45876-D1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	15.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 16:20	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
2	4.50	Chloride	BMB	67.153	444.627	
3	6.38	Sulfate	bMB	0.299	1.278	
4	9.28	Nitrate	BMB	0.078	0.314	
<b>TOTAL:</b>				67.53	446.22	3205.02



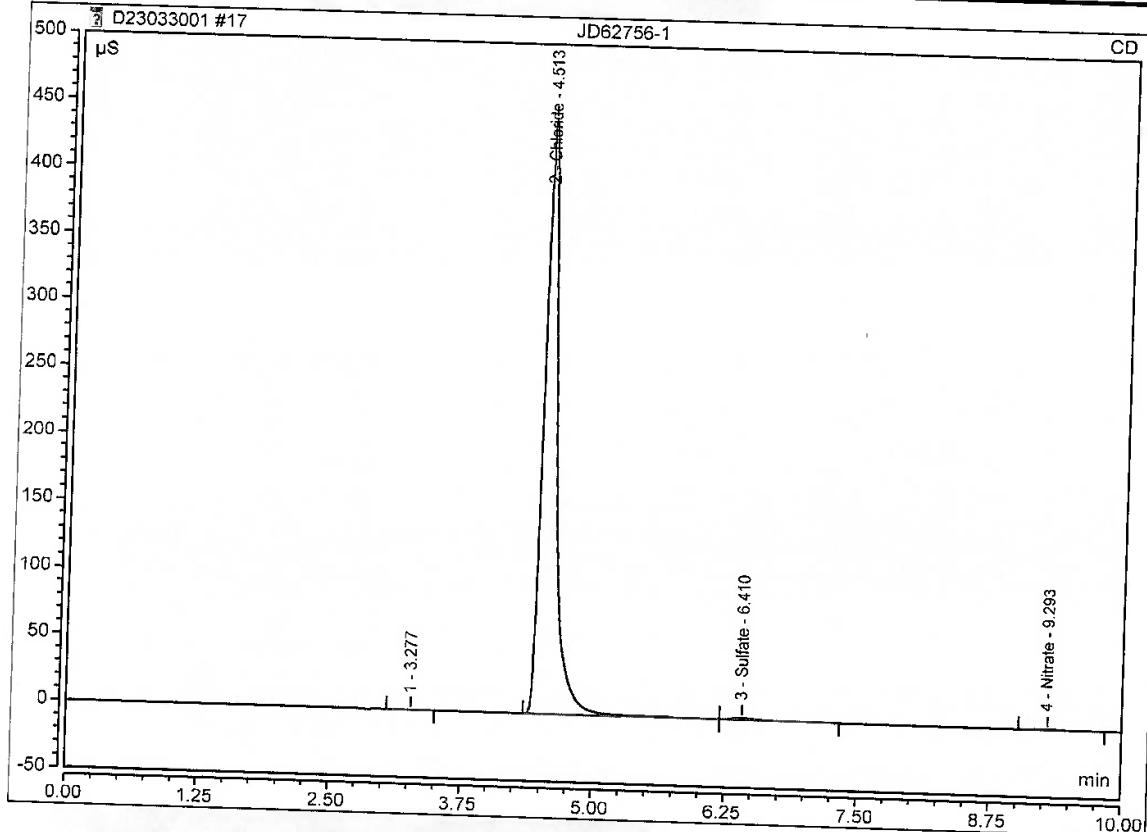
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62756-1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	15.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 16:33	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.51	Chloride	BMb	67.492	443.865	
3	6.41	Sulfate	bMB	0.290	1.275	3190.6989
4	9.29	Nitrate	BMB	0.086	0.344	20.8306
<b>TOTAL:</b>				67.87	445.48	9.9248
						3221.45



Anion/Integration

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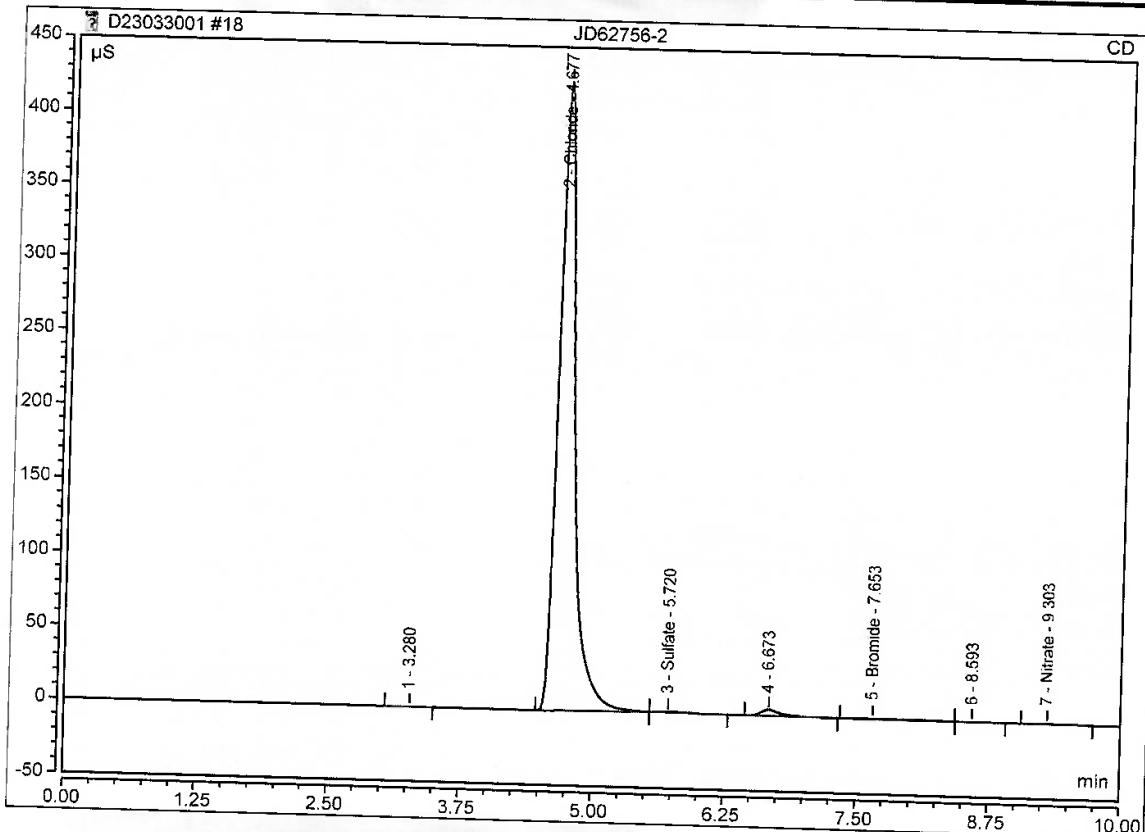
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62756-2	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	5.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 16:46	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
2	4.68	Chloride	BMb	77.707	426.492	
3	5.72	Sulfate	bMB	0.135	0.458	1224.3612
5	7.65	Bromide	BMb	0.239	0.370	3.5156
7	9.30	Nitrate	BMB	0.092	0.399	10.6313
<b>TOTAL:</b>				78.17	427.72	3.5839
						1242.09



Anion/Integration

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 Version 7.2.8.10783

9.4  
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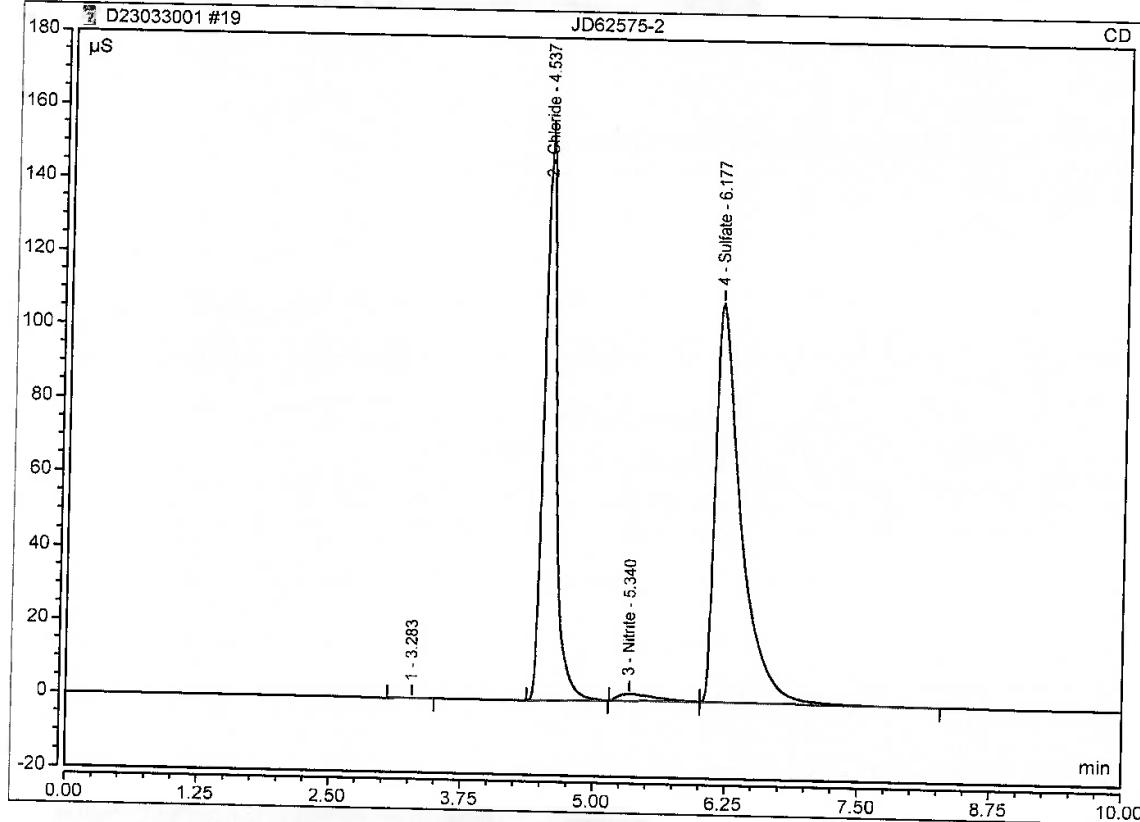
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62575-2	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	3.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 16:59	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
2	4.54	Chloride	BMB	21.299	153.578	
3	5.34	Nitrite	BMB	0.755	1.908	
4	6.18	Sulfate	BMB	30.178	107.953	
<b>TOTAL:</b>				52.23	263.44	616.89



Anion/Integration

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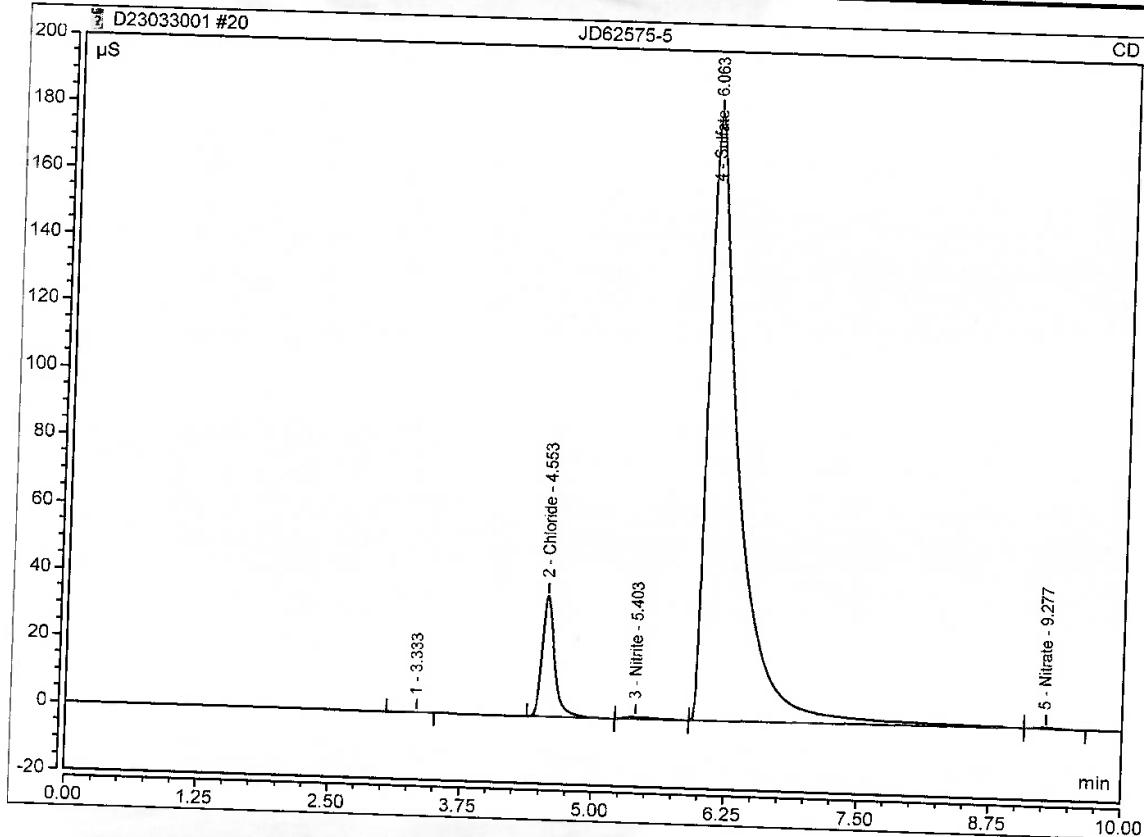
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 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62575-5	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	10.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 17:12	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}/\text{min}$	Height $\mu\text{S}$	Amount
2	4.55	Chloride	BMB	4.920	35.932	
3	5.40	Nitrite	BMB	0.198	0.630	157.2780
4	6.06	Sulfate	BMb	59.974	181.436	12.6339
5	9.28	Nitrate	bMB	0.078	0.342	2647.5641
<b>TOTAL:</b>				65.17	218.34	5.9495
						2823.43



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

9.4  
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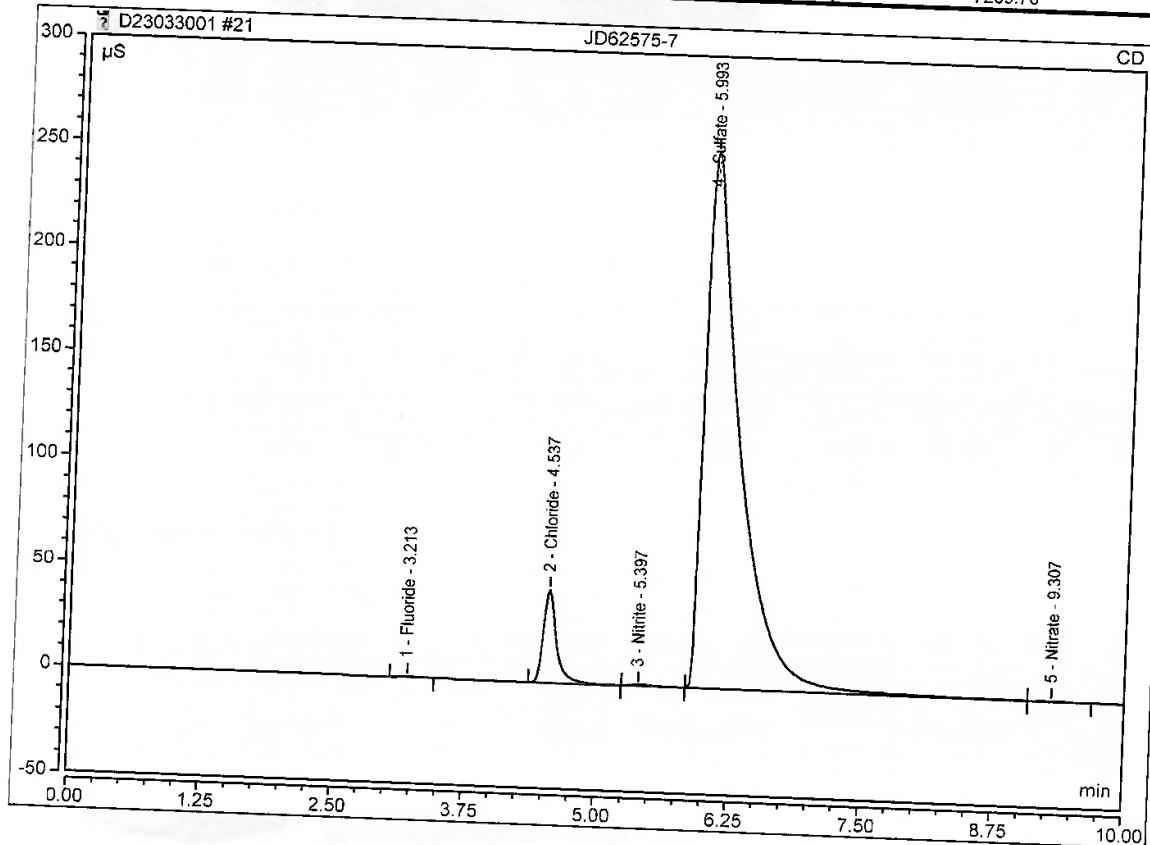
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 Sequence: D23033001

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## Peak Integration Report

Sample Name:	JD62575-7	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	16.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 17:25	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.063	0.335	
2	4.54	Chloride	BMB	6.328	44.109	
3	5.40	Nitrite	BMB	0.163	0.609	
4	5.99	Sulfate	BMb	98.050	254.846	
5	9.31	Nitrate	bMB	0.032	0.139	
<b>TOTAL:</b>				104.64	300.04	7269.76



Anion/Integration

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 Version 7.2.8.10783

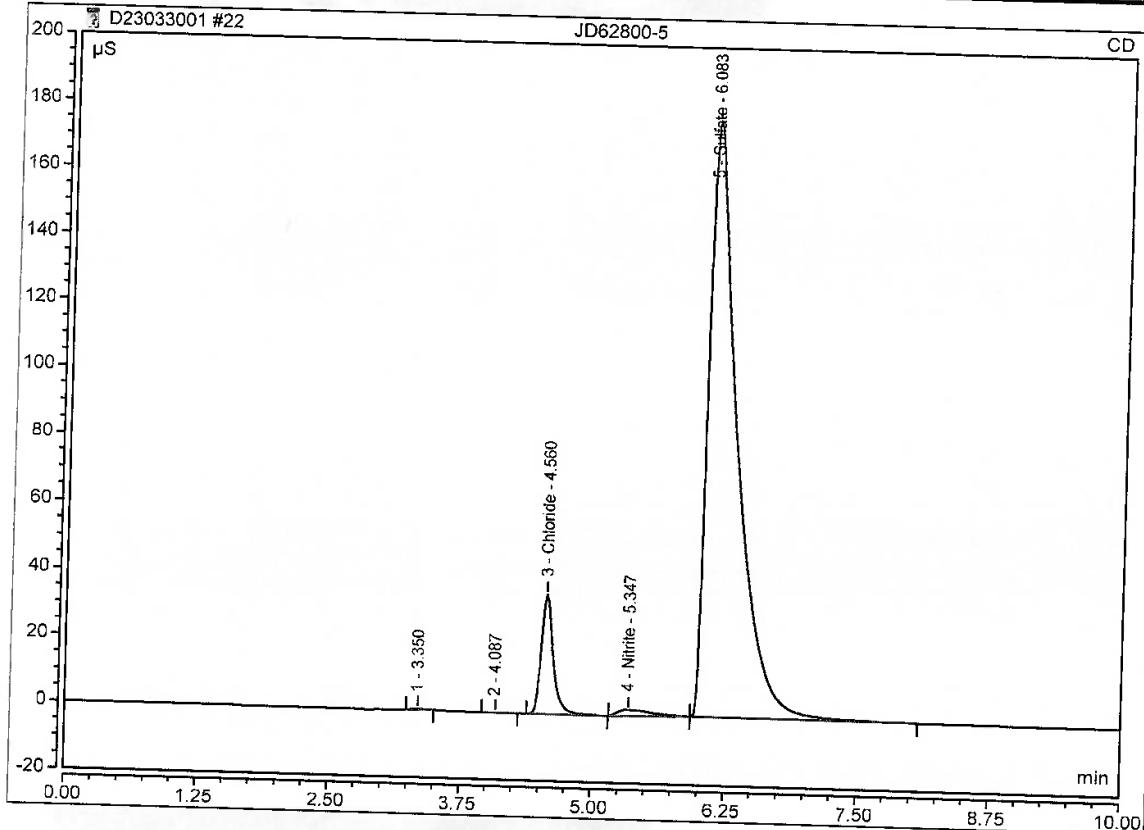
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62800-5	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	30.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 17:38	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\cdot}\text{min}$	Height $\mu\text{S}$	Amount
3	4.56	Chloride	BMB	4.709	35.607	
4	5.35	Nitrite	BMB	0.684	1.854	
5	6.08	Sulfate	BMB	56.497	177.984	
<b>TOTAL:</b>				61.89	215.44	7482.4050
						8071.82



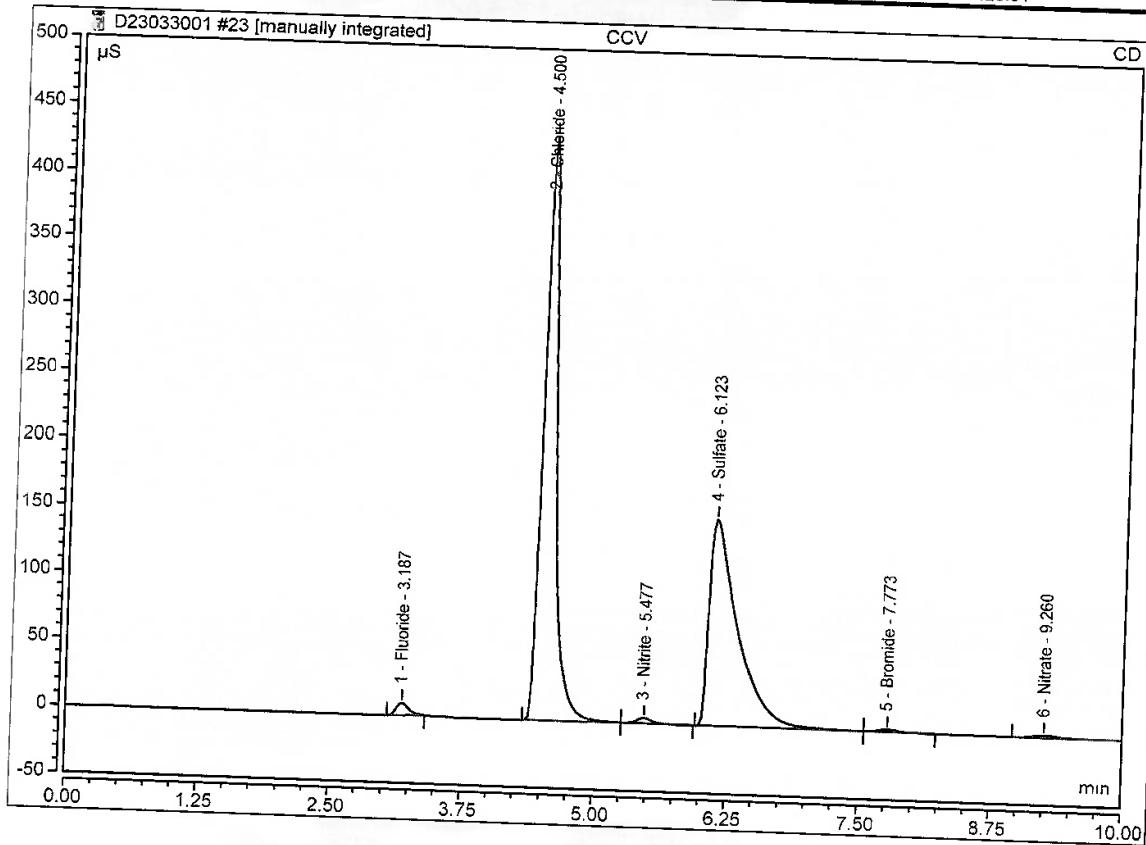
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCV	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 17:51	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB*	1.256	9.204	
2	4.50	Chloride	BMb	64.847	435.403	3.0744
3	5.48	Nitrite	bMB	0.665	3.952	204.3879
4	6.12	Sulfate	BMb	47.595	153.703	4.4491
5	7.77	Bromide	bMB	0.348	1.776	210.1317
6	9.26	Nitrate	BMB	0.458	1.828	3.0556
<b>TOTAL:</b>				115.17	605.87	3.8066
						428.91



Anion/Integration

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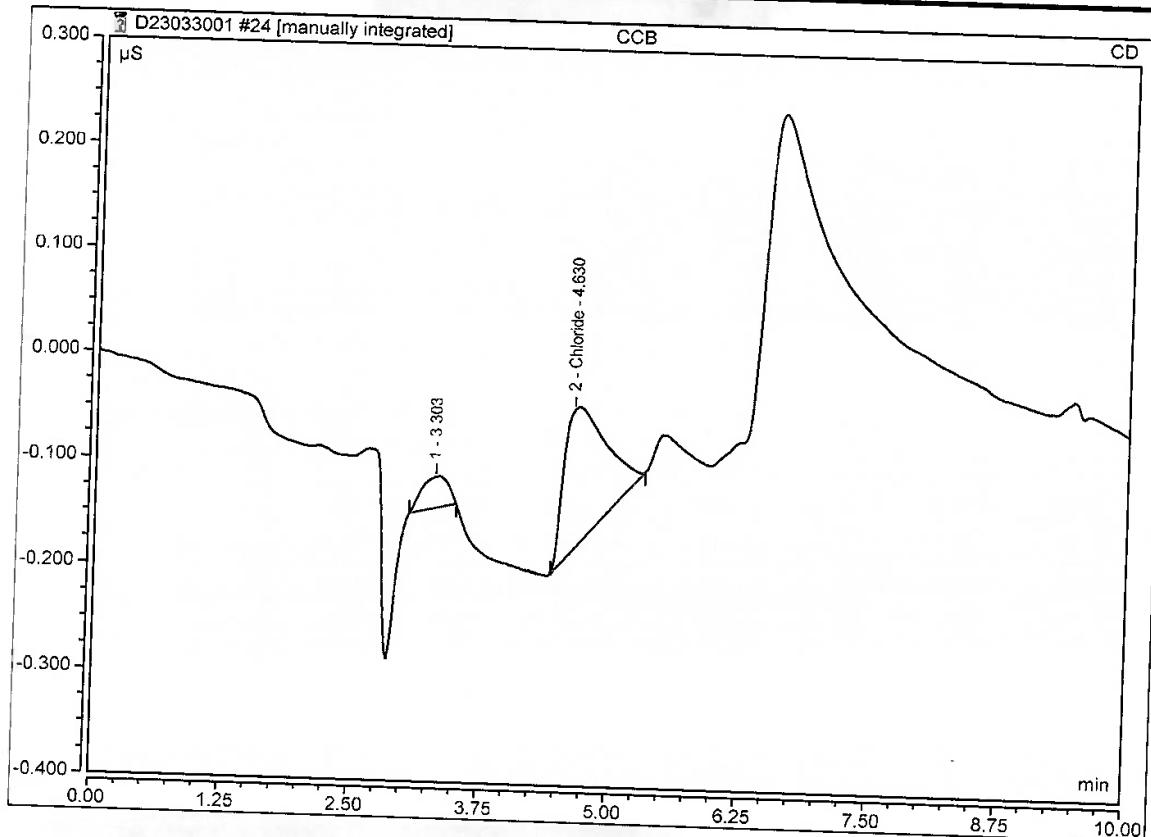
Logged on User: Chemistry  
 Instrument: Integri<sup>n</sup>\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 18:04	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.63	Chloride	BMB	0.059	0.133	0.4260
		TOTAL:		0.06	0.13	0.43



Anion/Integration

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9.4  
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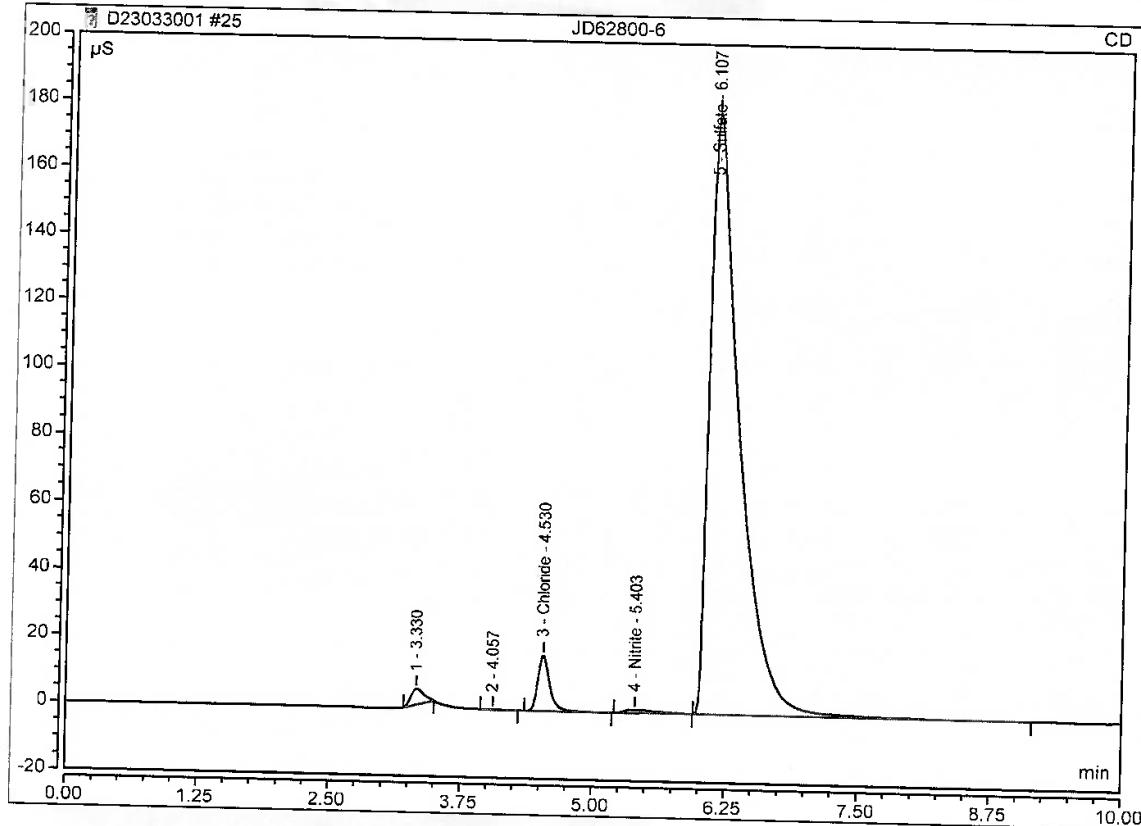
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62800-6	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	34.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 18:17	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
3	4.53	Chloride	BMB	2.160	16.639	
4	5.40	Nitrite	BMB	0.343	1.030	
5	6.11	Sulfate	BMB	58.348	179.602	
<b>TOTAL:</b>				60.85	197.27	9073.67



Anion/Integration

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9.4  
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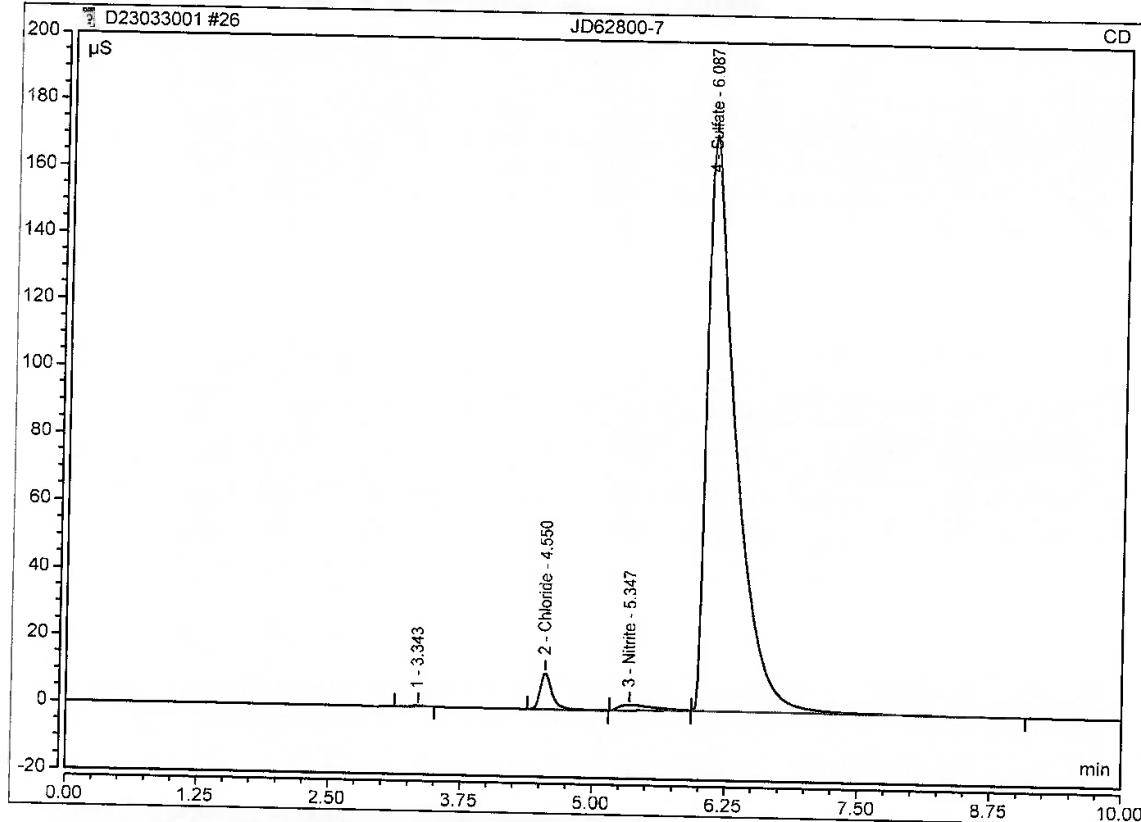
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 Instrument: Integriion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62800-7	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	20.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 18:31	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
2	4.55	Chloride	BMB	1.405	10.795	
3	5.35	Nitrite	BMB	0.612	1.678	
4	6.09	Sulfate	BMB	53.897	171.631	
<b>TOTAL:</b>				55.91	184.10	4933.93



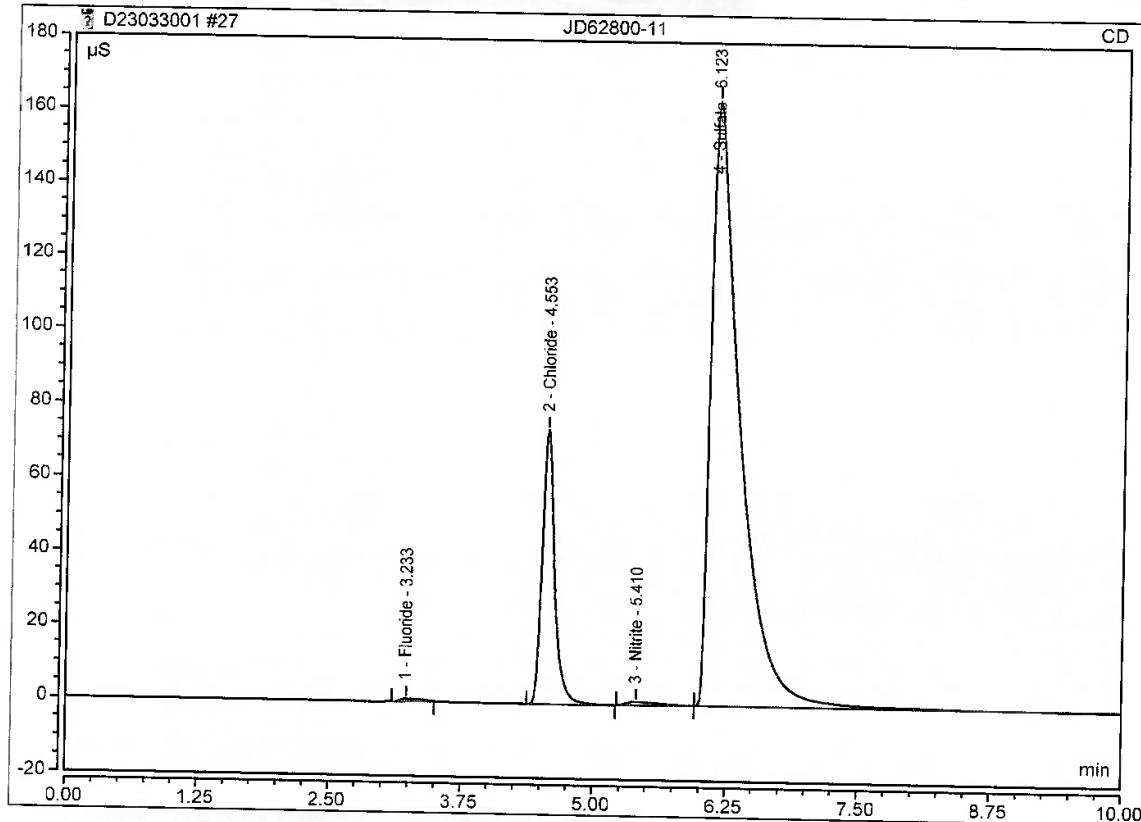
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62800-11	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	3.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 18:44	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.23	Fluoride	BMB	0.147	0.677	
2	4.55	Chloride	BMB	10.215	74.169	
3	5.41	Nitrite	BMB	0.311	0.927	
4	6.12	Sulfate	BMB	53.461	164.136	
<b>TOTAL:</b>				64.13	239.91	812.52



Anion/Integration

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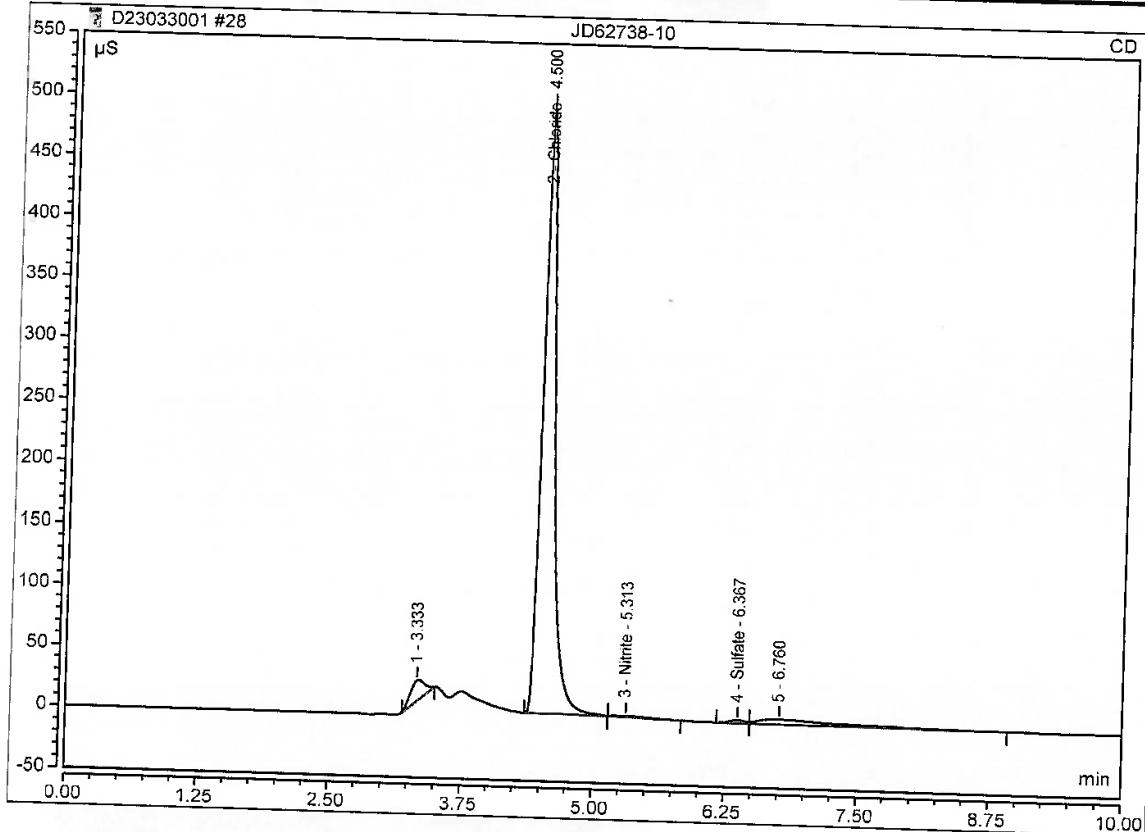
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62738-10	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	5.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 18:57	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
2	4.50	Chloride	BMB	70.091	492.606	
3	5.31	Nitrite	BMB	0.367	1.111	1104.4768
4	6.37	Sulfate	BM	0.365	2.429	12.0867
<b>TOTAL:</b>				<b>70.82</b>	<b>496.15</b>	<b>8.5847</b>
						<b>1125.15</b>



Anion/Integration

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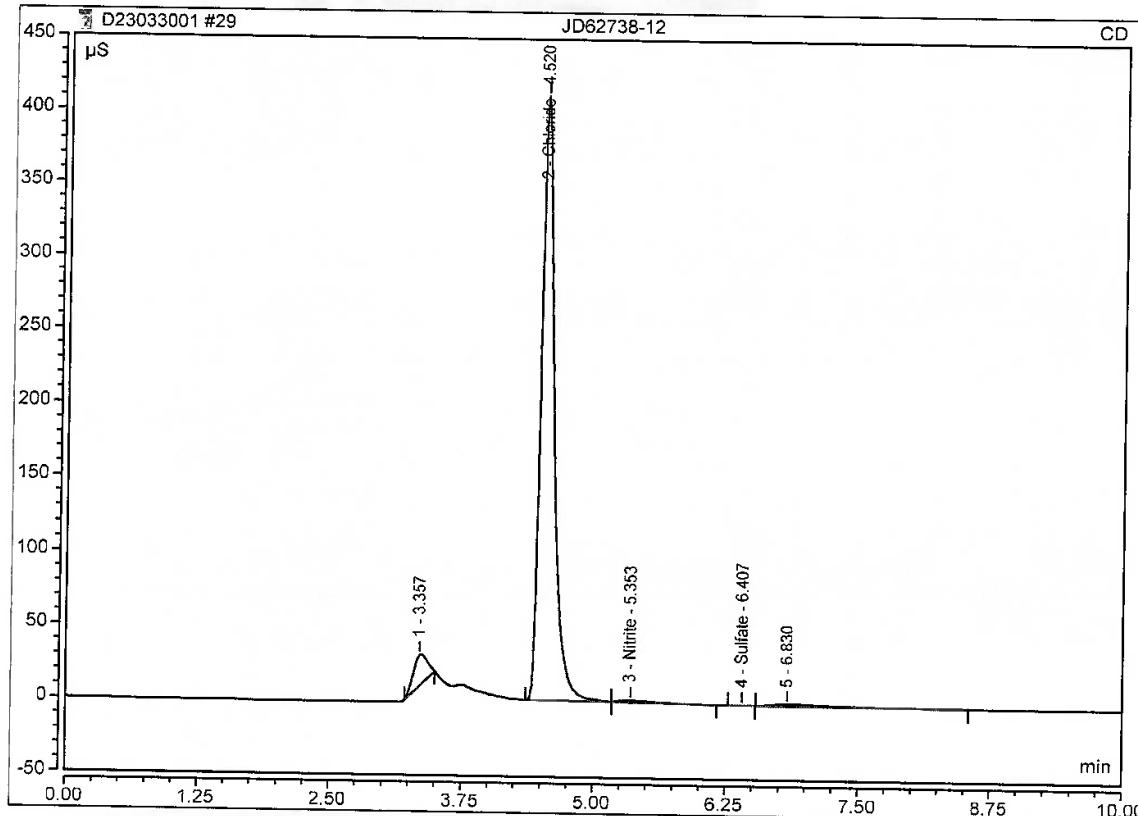
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62738-12	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	3.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 19:10	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
2	4.52	Chloride	BMB	58.784	410.315	555.8958
3	5.35	Nitrite	BMB	0.469	1.258	9.3311
4	6.41	Sulfate	BMB	0.007	0.054	0.4159
<b>TOTAL:</b>				59.26	411.63	565.64



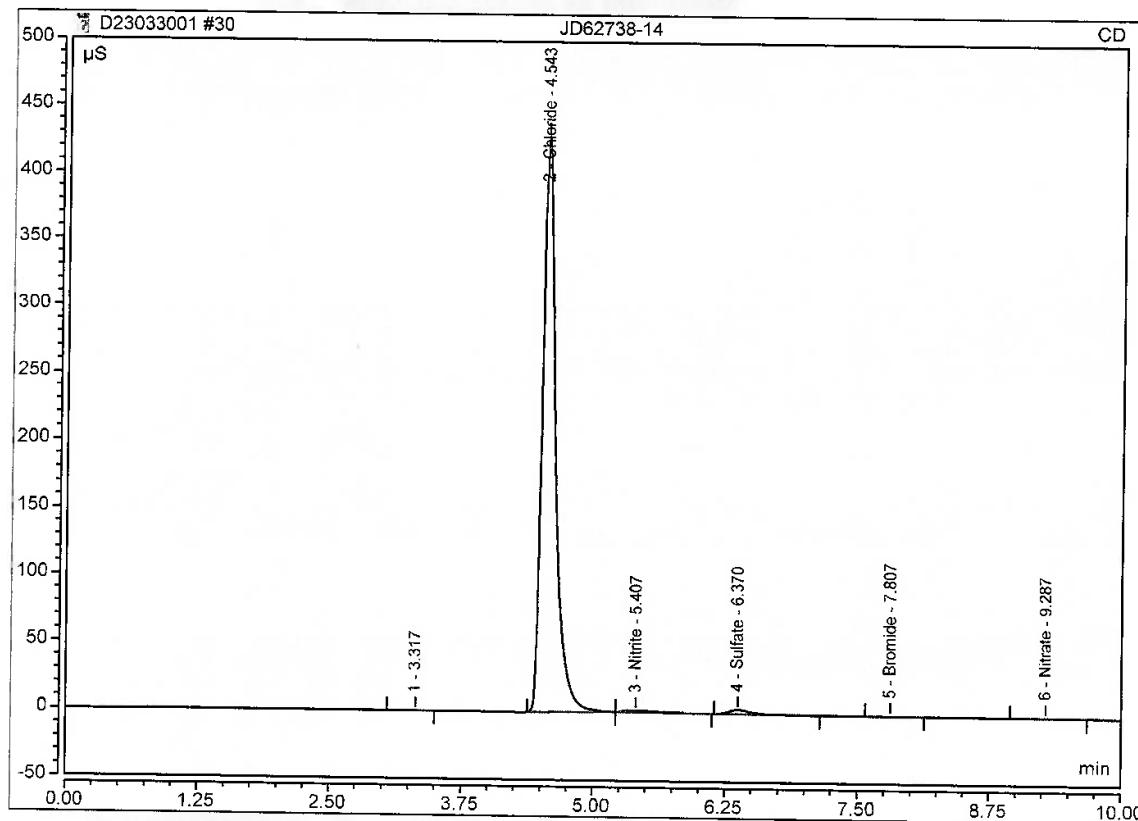
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62738-14	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	7.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 19:23	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.54	Chloride	BMB	62.572	429.260	1380.5725
3	5.41	Nitrite	bMB	0.431	1.149	19.9968
4	6.37	Sulfate	BMB	0.702	3.239	22.4417
5	7.81	Bromide	BMB	0.020	0.108	1.7760
6	9.29	Nitrate	BMB	0.044	0.185	2.1475
<b>TOTAL:</b>				<b>63.77</b>	<b>433.94</b>	<b>1426.93</b>



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Anion/Integration

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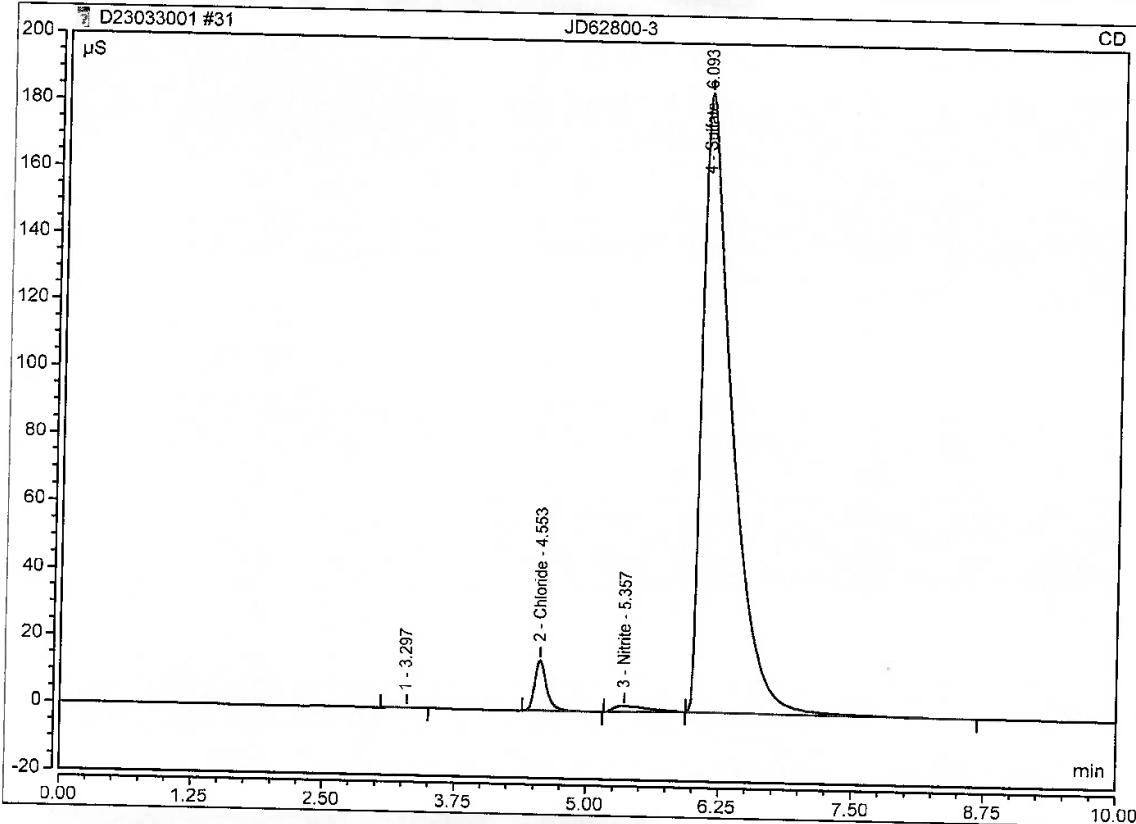
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62800-3	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	14.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 19:36	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
2	4.55	Chloride	BMB	1.984	15.071	
3	5.36	Nitrite	BMB	0.655	1.775	
4	6.09	Sulfate	BMB	59.342	184.952	
<b>TOTAL:</b>				61.98	201.80	3819.74



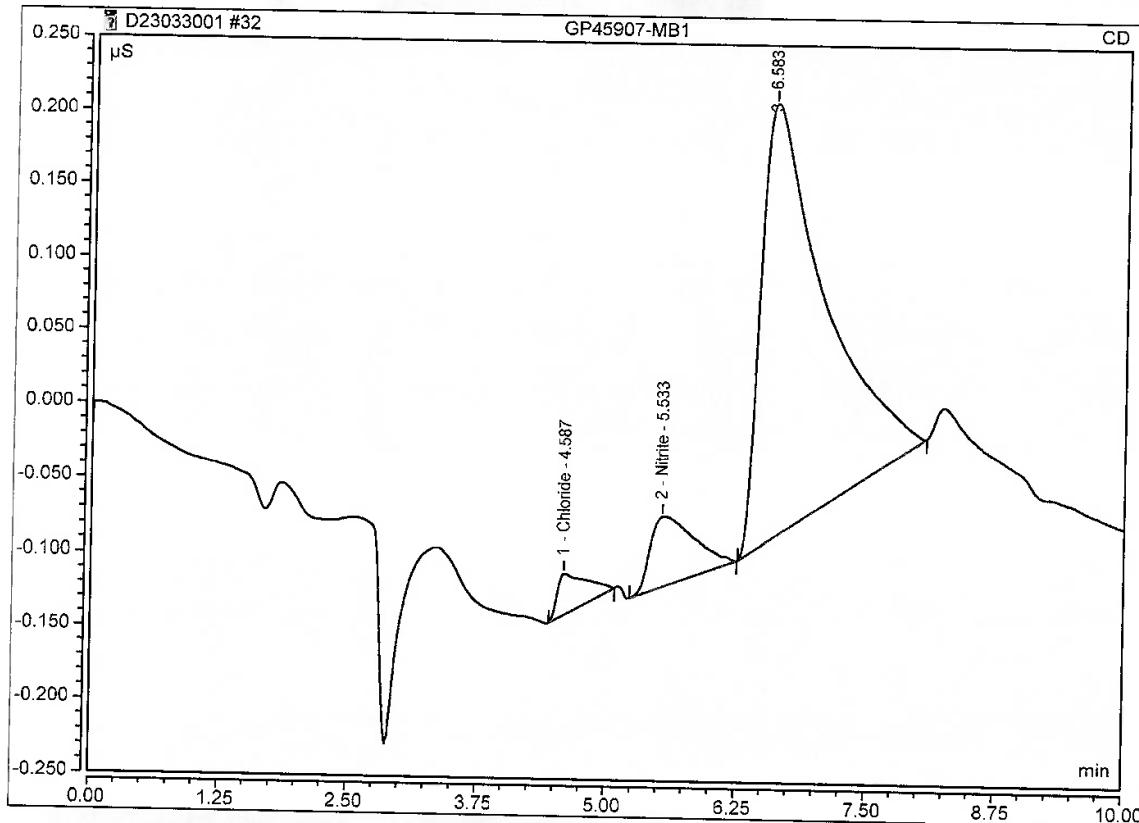
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45907-MB1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 19:49	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	4.59	Chloride	BMB	0.009	0.028	
2	5.53	Nitrite	BMB	0.023	0.047	
		TOTAL:		0.03	0.08	0.33



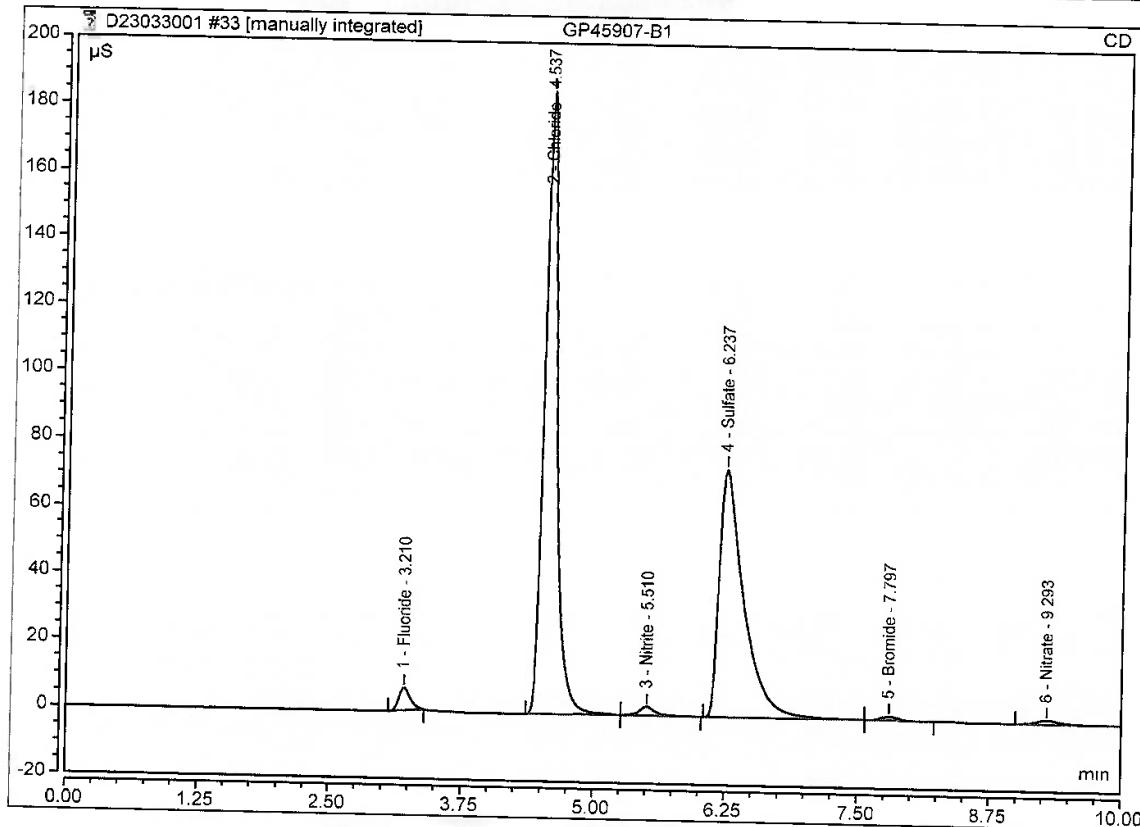
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45907-B1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 20:02	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB*	0.839	6.681	2.0674
2	4.54	Chloride	BMb	26.518	185.575	83.7222
3	5.51	Nitrite	bMB	0.443	2.584	2.9362
4	6.24	Sulfate	BMb	19.687	73.882	86.9828
5	7.80	Bromide	bMB	0.222	1.152	1.9772
6	9.29	Nitrate	BMB	0.287	1.158	2.3569
<b>TOTAL:</b>				48.00	271.03	180.04



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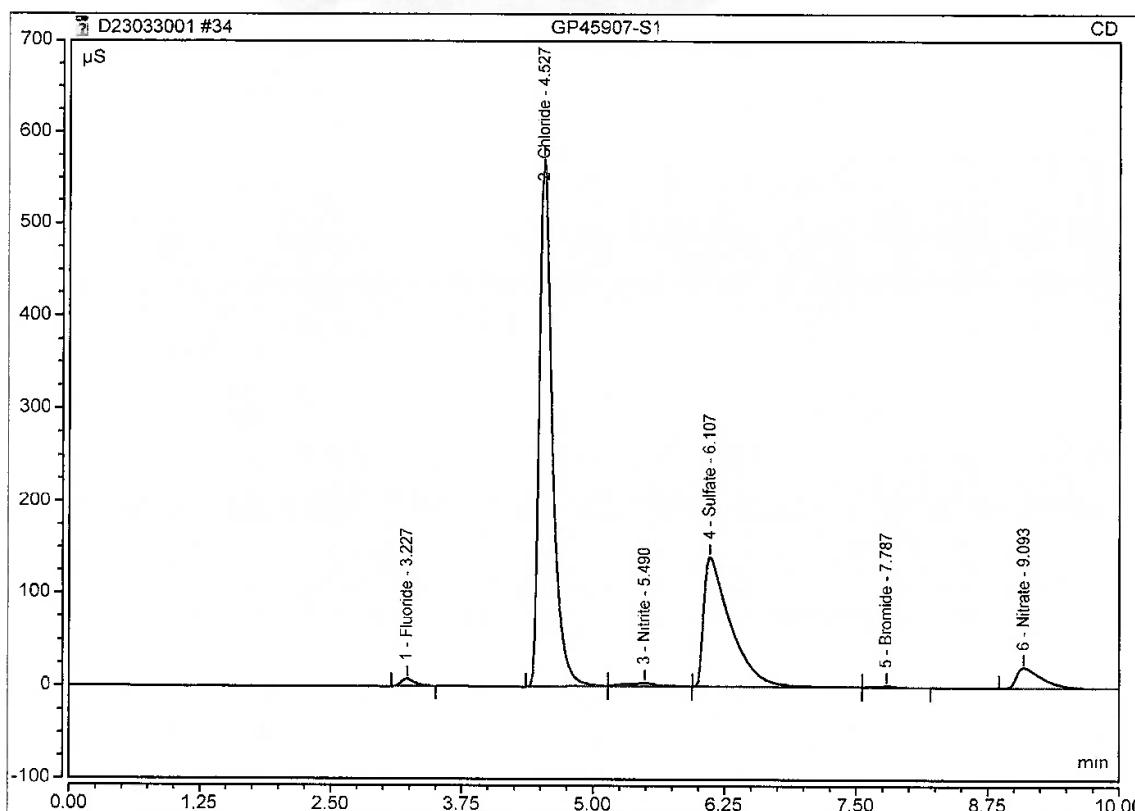
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### Peak Integration Report

<b>Sample Name:</b>	GP45907-S1	<b>Inj. Vol.:</b>	5000.00
<b>Injection Type:</b>	Unknown	<b>Dilution Factor:</b>	1.0000
<b>Instrument Method:</b>	Anions_012919	<b>Operator:</b>	Chemistry
<b>Inj. Date / Time:</b>	30-Mar-2023 / 20:15	<b>Run Time:</b>	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.23	Fluoride	BMB	1.056	7.466	2.5904
2	4.53	Chloride	BMb	83.001	571.998	261.5382
3	5.49	Nitrite	bMB	0.999	3.260	6.7292
4	6.11	Sulfate	BMb	40.945	139.656	180.7860
5	7.79	Bromide	bMB	0.249	1.309	2.2106
6	9.09	Nitrate	BMB	5.895	21.513	49.7493
<b>TOTAL:</b>				132.14	745.20	503.60



Anion/Integration

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9.4  
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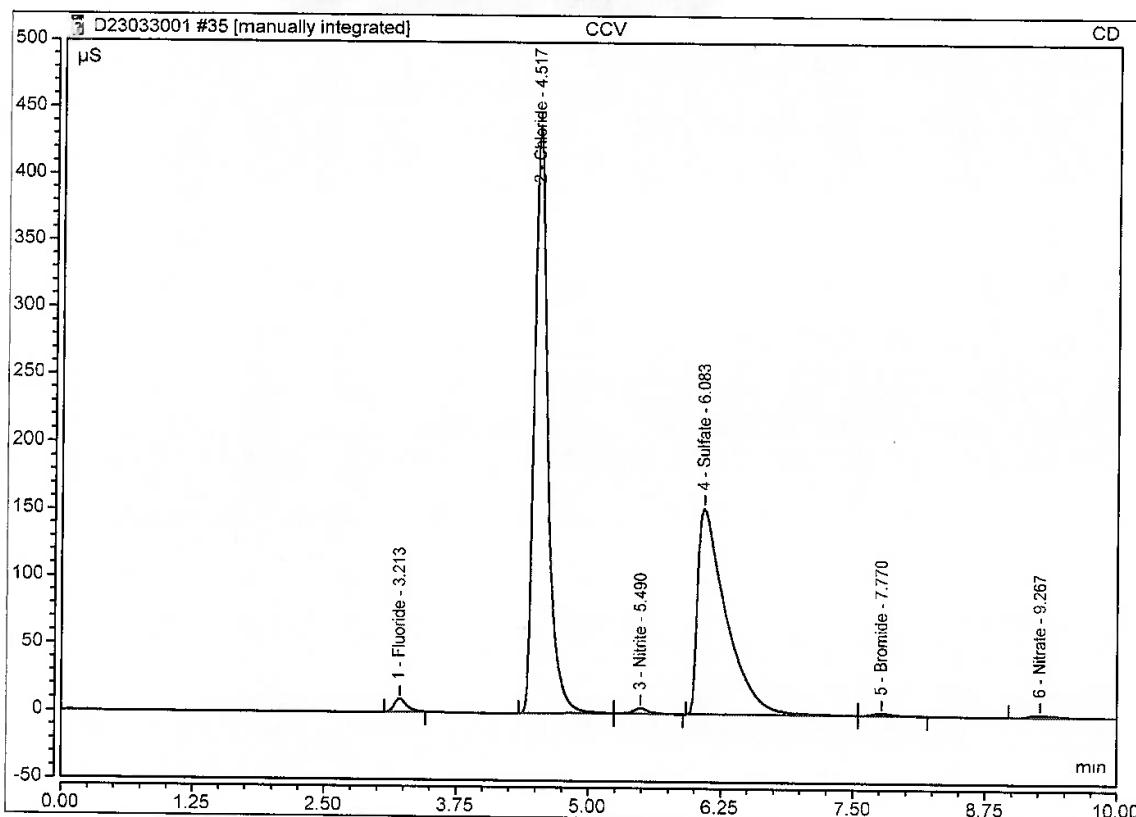
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 Sequence: D23033001

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### Peak Integration Report

<b>Sample Name:</b>	CCV	<b>Inj. Vol.:</b>	5000.00
<b>Injection Type:</b>	Unknown	<b>Dilution Factor:</b>	1.0000
<b>Instrument Method:</b>	Anions_012919	<b>Operator:</b>	Chemistry
<b>Inj. Date / Time:</b>	30-Mar-2023 / 20:28	<b>Run Time:</b>	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB*	1.341	9.828	3.2801
2	4.52	Chloride	BMb	64.814	438.481	204.2850
3	5.49	Nitrite	bMB	0.628	3.951	4.1958
4	6.08	Sulfate	BMb	47.215	153.438	208.4553
5	7.77	Bromide	bMB	0.346	1.799	3.0420
6	9.27	Nitrate	BMB	0.461	1.813	3.8343
<b>TOTAL:</b>				114.81	609.31	427.09



Anion/Integration

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 Version 7.2.8.10783

9.4  
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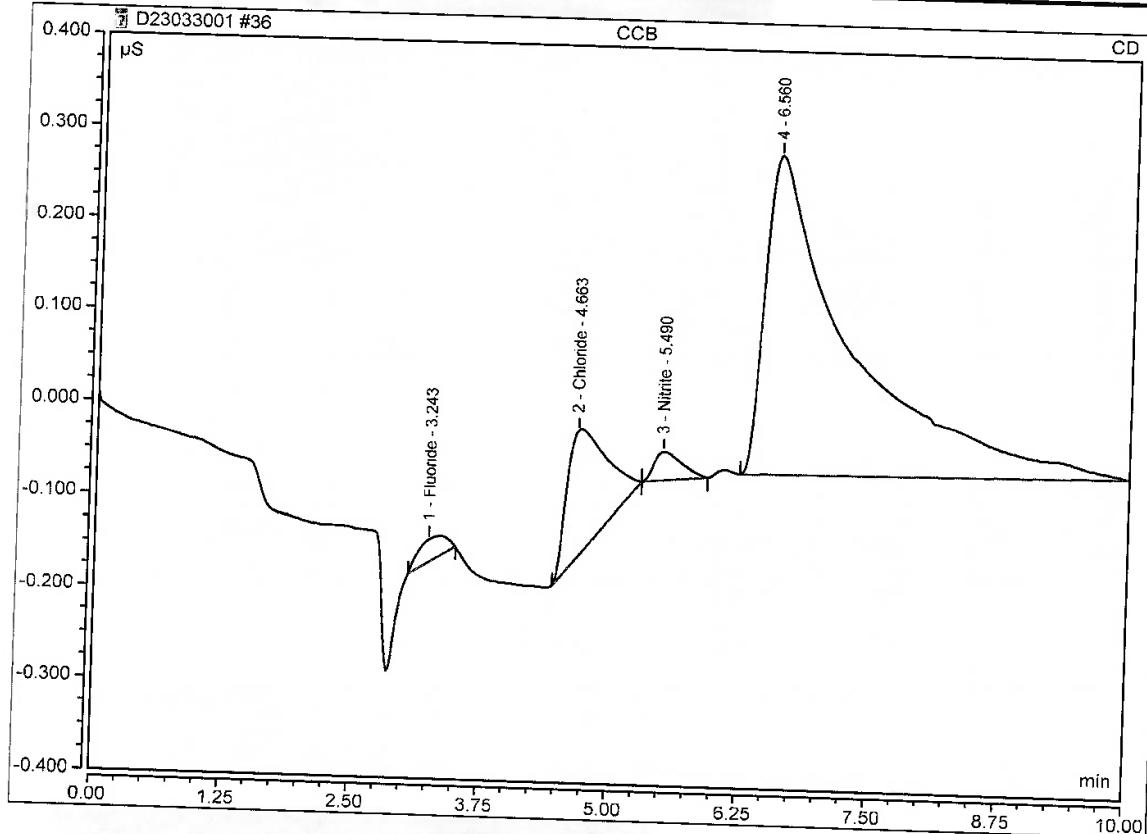
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 20:41	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}/\text{min}$	Height $\mu\text{S}$	Amount
1	3.24	Fluoride	BMB	0.008	0.025	
2	4.66	Chloride	BMb	0.058	0.139	0.0563
3	5.49	Nitrite	bMB	0.009	0.031	0.4214
<b>TOTAL:</b>				0.07	0.19	n.a. 0.48



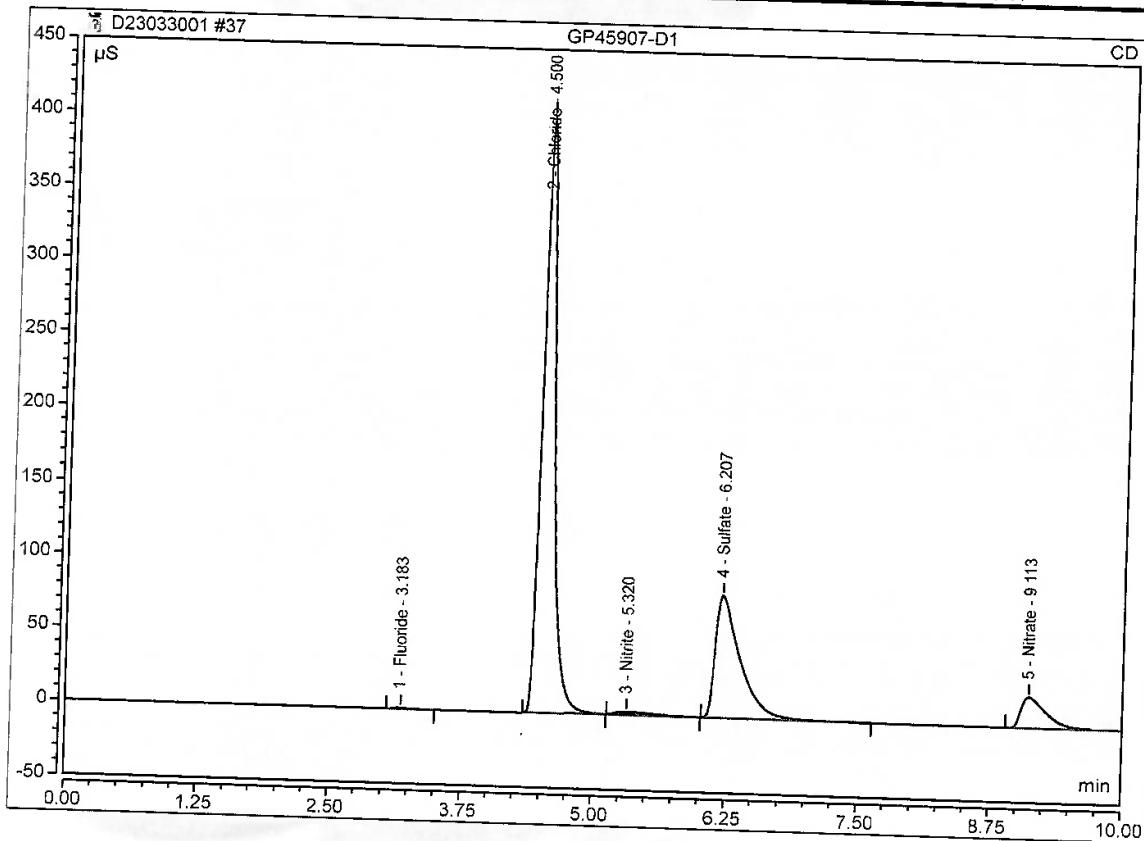
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45907-D1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 20:54	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.18	Fluoride	BMB	0.119	0.801	
2	4.50	Chloride	BMB	57.823	404.908	0.3242
3	5.32	Nitrite	BMB	0.760	1.910	182.2758
4	6.21	Sulfate	BMB	22.460	82.733	5.1014
5	9.11	Nitrate	BMB	5.710	20.735	99.2178
<b>TOTAL:</b>				86.87	511.09	48.1870
						335.11



Anion/Integration

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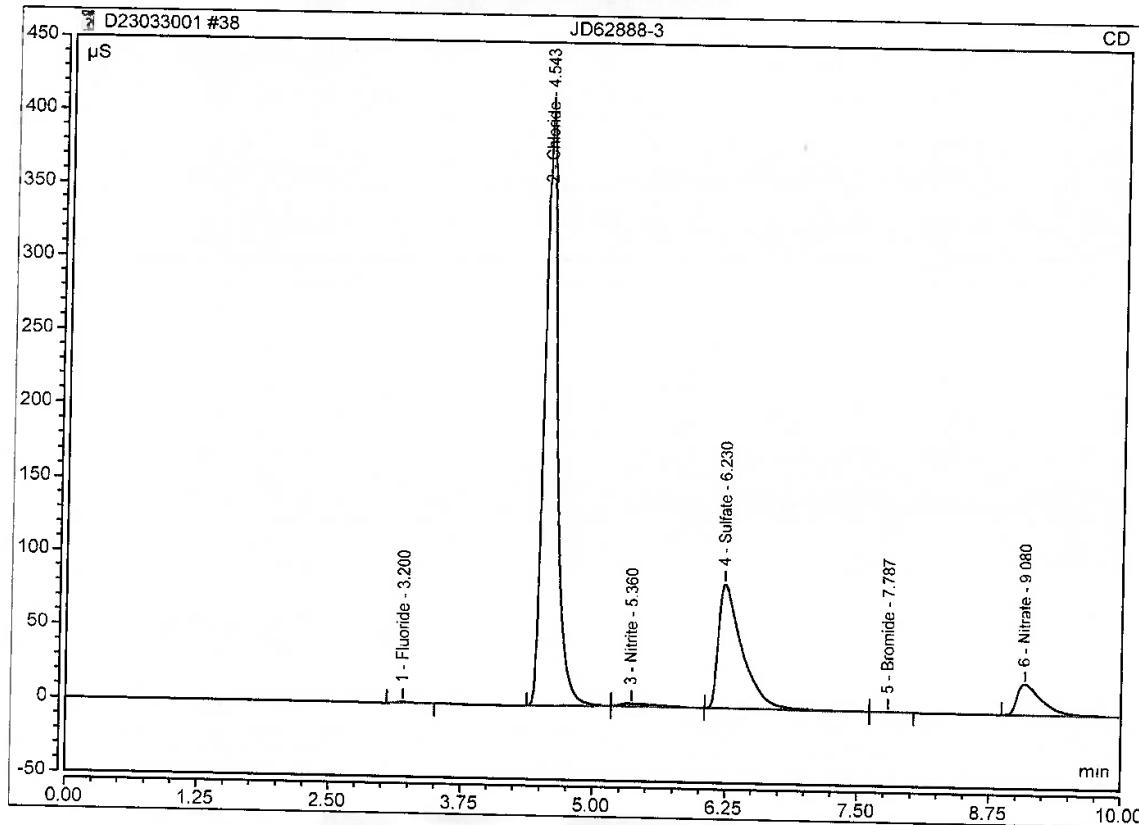
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62888-3	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 21:07	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.20	Fluoride	BMB	0.115	0.794	0.3162
2	4.54	Chloride	BMb	58.088	403.323	183.1095
3	5.36	Nitrite	bMB	0.753	1.896	5.0522
4	6.23	Sulfate	BMb	21.914	84.015	96.8068
5	7.79	Bromide	bMB	0.014	0.079	0.1966
6	9.08	Nitrate	BMB	5.601	20.603	47.2642
<b>TOTAL:</b>				86.48	510.71	332.75



Anion/Integration

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 Version 7.2.8.10783

9.4  
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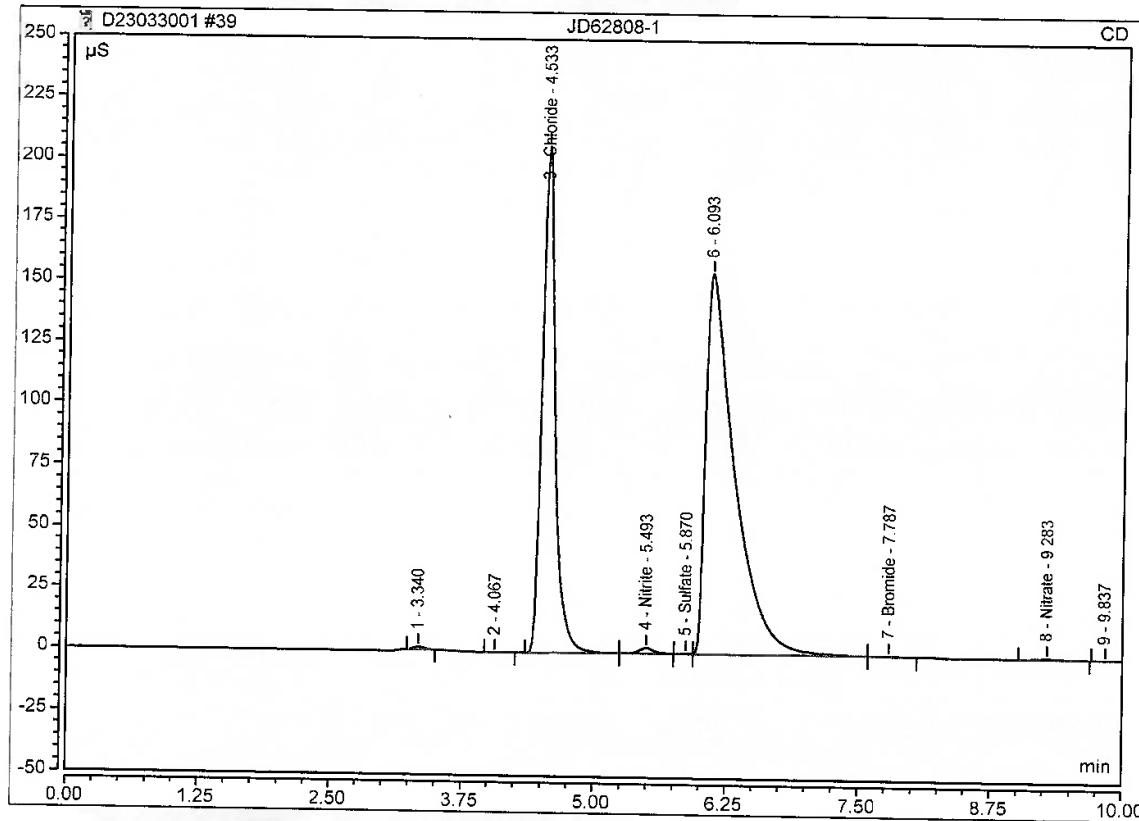
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62808-1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Ini. Date / Time:	30-Mar-2023 / 21:20	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
3	4.53	Chloride	BMb	29.685	206.534	93.6922
4	5.49	Nitrite	bMB	0.358	2.300	2.3579
5	5.87	Sulfate	BM	0.025	0.238	0.2167
7	7.79	Bromide	bMB	0.029	0.161	0.3268
8	9.28	Nitrate	BMB	0.094	0.370	0.7278
<b>TOTAL:</b>				30.19	209.60	97.32



Anion/Integration

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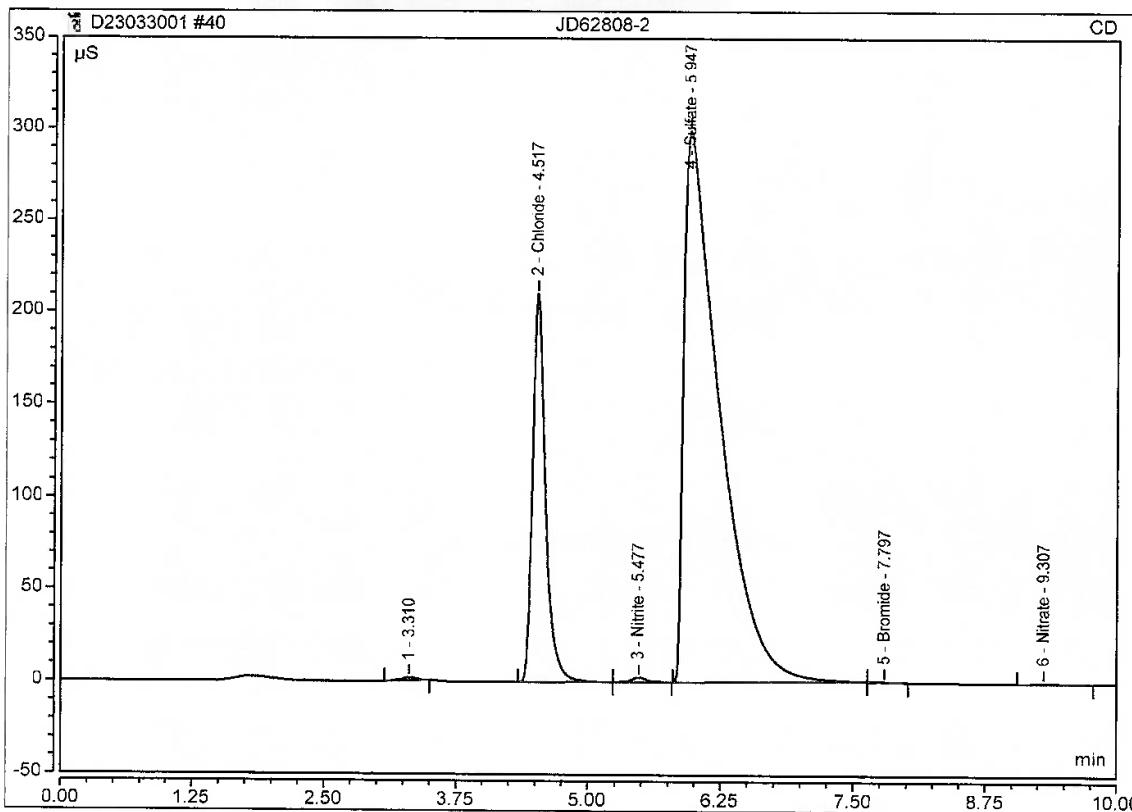
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62808-2	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 21:33	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\cdot}\text{min}$	Height $\mu\text{S}$	Amount
2	4.52	Chloride	BMB	30.199	210.482	95.3118
3	5.48	Nitrite	bMB	0.400	2.529	2.6415
4	5.95	Sulfate	BMB	117.252	299.371	517.5084
5	7.80	Bromide	bMB	0.025	0.148	0.2907
6	9.31	Nitrate	BMB	0.098	0.402	0.7652
<b>TOTAL:</b>				147.97	512.93	616.52



Anion/Integration

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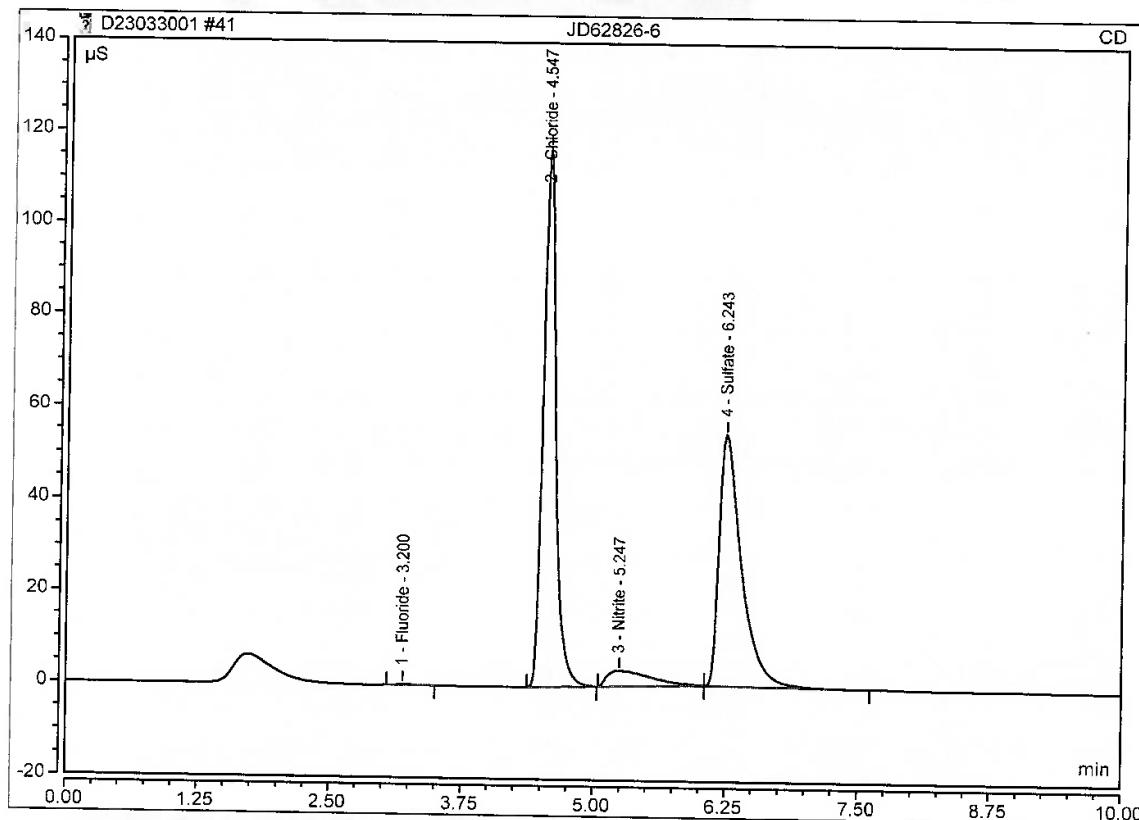
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62826-6	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 21:46	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.20	Fluoride	BMB	0.033	0.173	0.1174
2	4.55	Chloride	BMB	16.220	117.268	51.3018
3	5.25	Nitrite	BMb	1.638	3.442	11.0953
4	6.24	Sulfate	bMB	13.315	54.439	58.8646
TOTAL:				31.21	175.32	121.38



Anion/Integration

Chromleon (c) Dionex 1996-2009  
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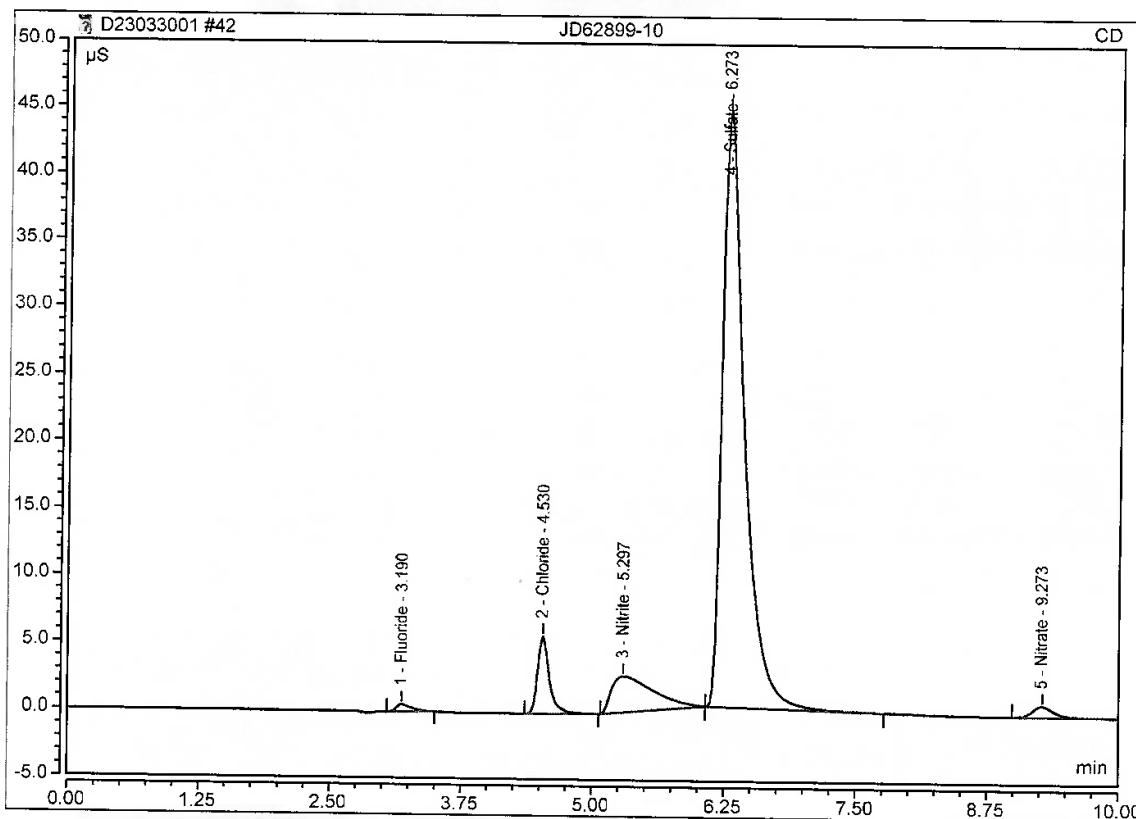
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62899-10	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 21:59	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB	0.083	0.514	0.2373
2	4.53	Chloride	BMB	0.739	5.702	2.5677
3	5.30	Nitrite	BMb	1.214	2.660	8.2021
4	6.27	Sulfate	bMB	10.800	44.474	47.7663
5	9.27	Nitrate	BMB	0.185	0.798	1.4960
<b>TOTAL:</b>				13.02	54.15	60.27



Anion/Integration

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9.4  
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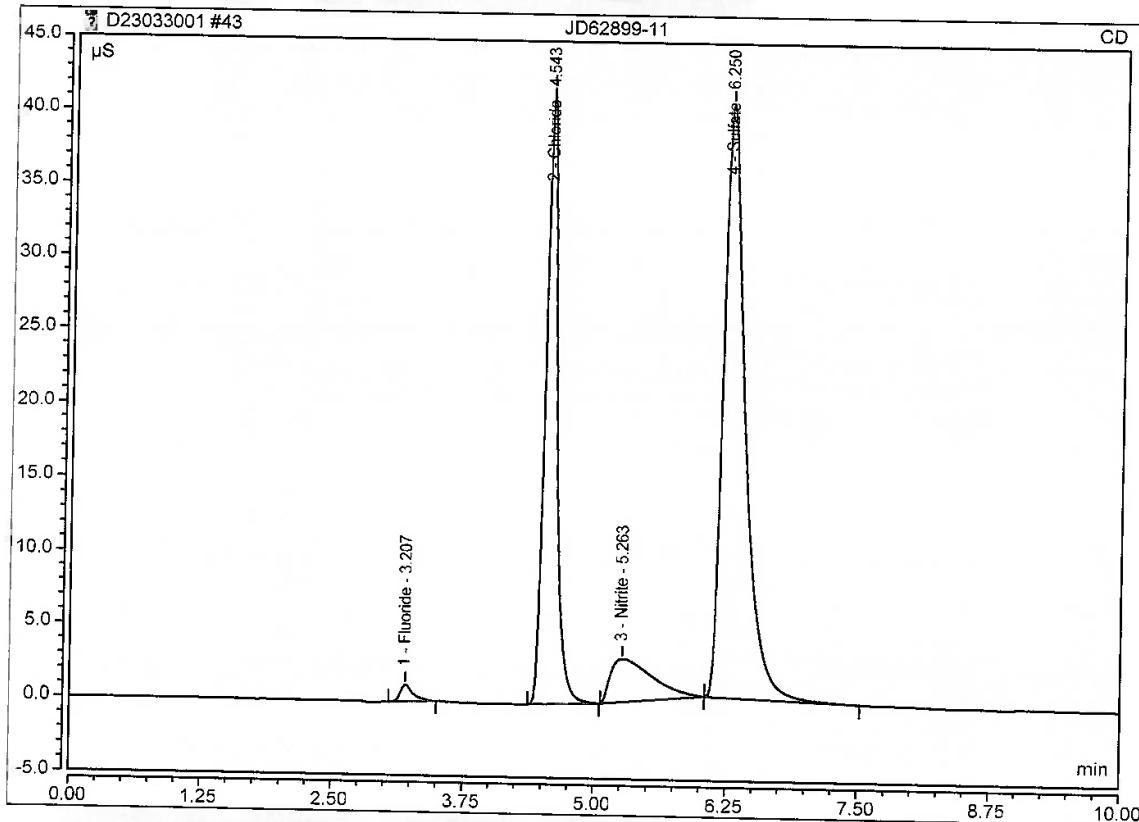
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62899-11	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 22:12	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.158	1.123	
2	4.54	Chloride	BMB	5.504	41.875	
3	5.26	Nitrite	BMb	1.332	2.931	
4	6.25	Sulfate	bMB	9.594	40.347	
TOTAL:				16.59	86.28	69.43



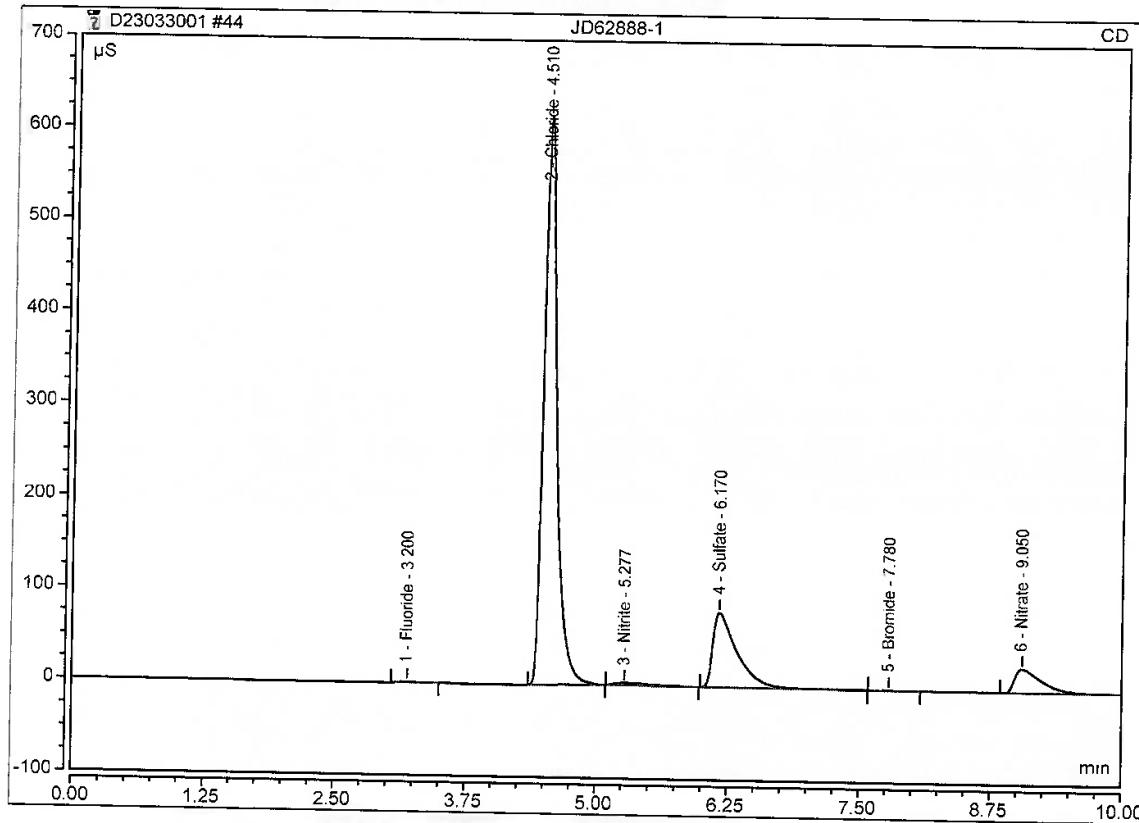
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62888-1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 22:25	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
1	3.20	Fluoride	BMB	0.152	1.062	0.4042
2	4.51	Chloride	BMB	89.886	617.599	283.2152
3	5.28	Nitrite	BMB	0.857	2.154	5.7599
4	6.17	Sulfate	BMB	21.167	81.095	93.5124
5	7.78	Bromide	BMB	0.041	0.231	0.4328
6	9.05	Nitrate	BMB	7.371	25.542	62.2191
<b>TOTAL:</b>				119.47	727.68	445.54



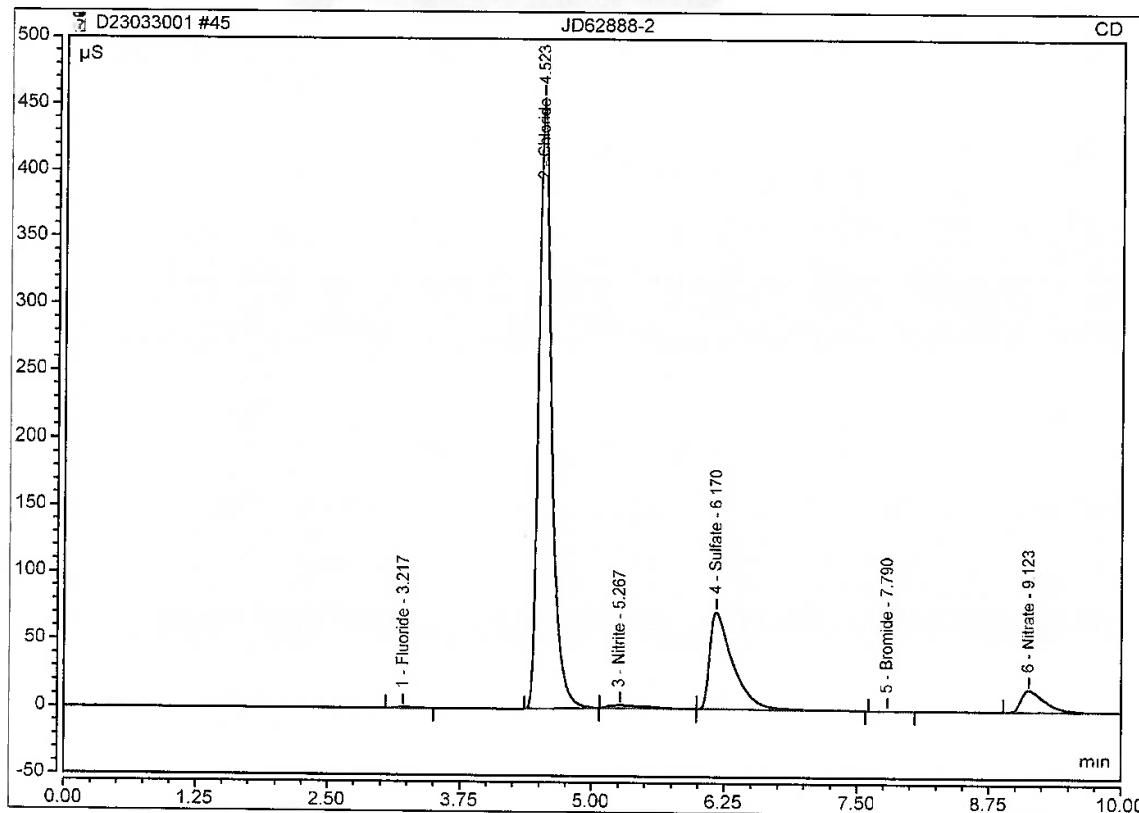
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62888-2	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 22:39	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.102	0.716	0.2851
2	4.52	Chloride	BMB	64.042	455.509	201.8523
3	5.27	Nitrite	BMB	0.994	2.383	6.6976
4	6.17	Sulfate	BMB	18.084	72.087	79.9065
5	7.79	Bromide	BMB	0.018	0.106	0.2365
6	9.12	Nitrate	BMB	4.288	16.415	36.1686
<b>TOTAL:</b>				87.53	547.22	325.15



Anion/Integration

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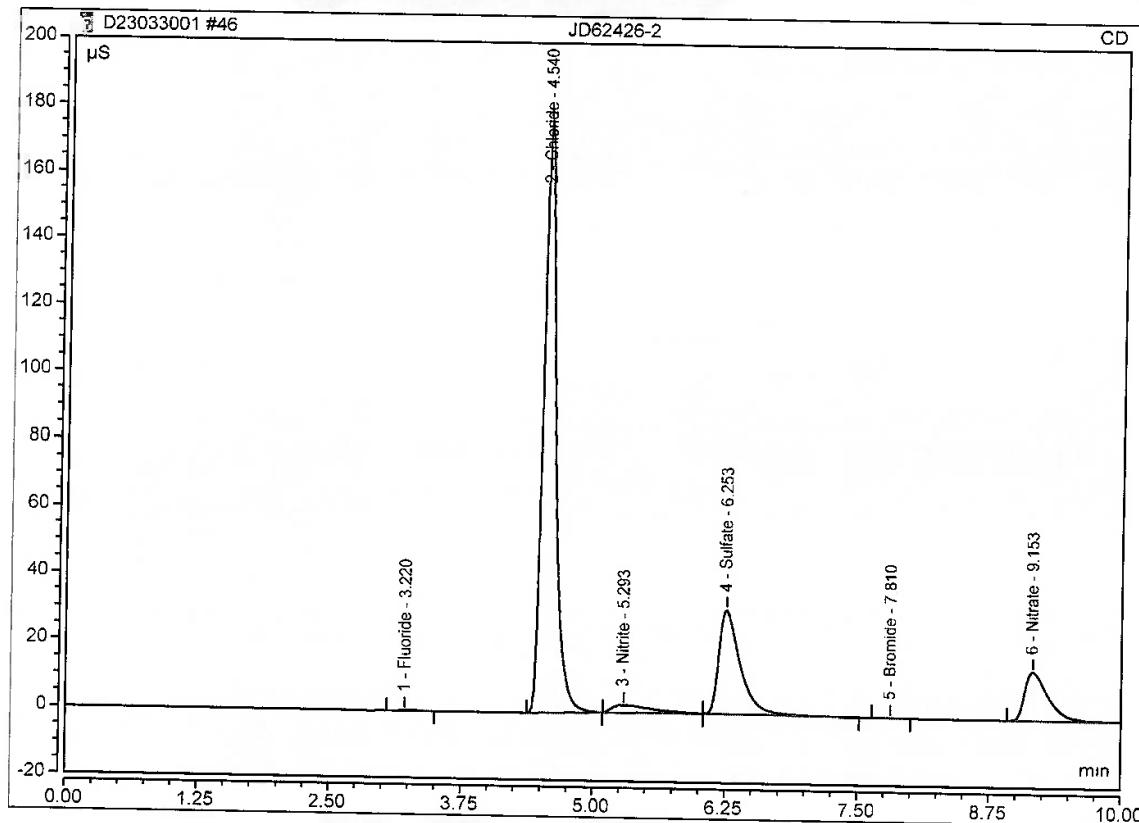
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62426-2	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 22:51	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.032	0.181	0.1147
2	4.54	Chloride	BMB	24.362	176.958	76.9359
3	5.29	Nitrite	BMB	0.970	2.261	6.5314
4	6.25	Sulfate	bMB	7.167	30.972	31.7343
5	7.81	Bromide	BMB	0.003	0.021	0.1074
6	9.15	Nitrate	BMB	3.654	14.342	30.8104
<b>TOTAL:</b>				36.19	224.73	146.23



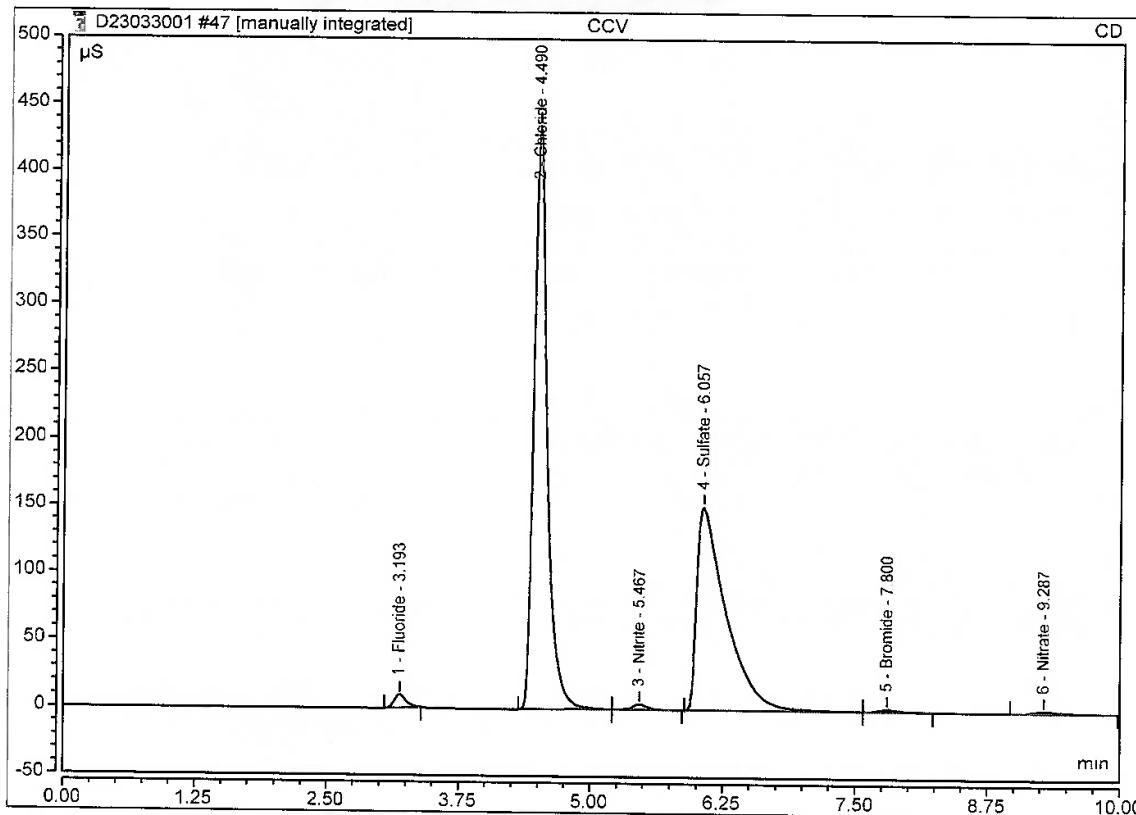
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### Peak Integration Report

Sample Name:	CCV	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 23:04	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB*	1.258	9.686	3.0794
2	4.49	Chloride	BMb	64.034	435.398	201.8292
3	5.47	Nitrite	bMB	0.623	3.923	4.1645
4	6.06	Sulfate	BMb	48.098	151.344	212.3488
5	7.80	Bromide	bMB	0.343	1.757	3.0167
6	9.29	Nitrate	BMB	0.459	1.831	3.8118
<b>TOTAL:</b>				114.81	603.94	428.25



Anion/Integration

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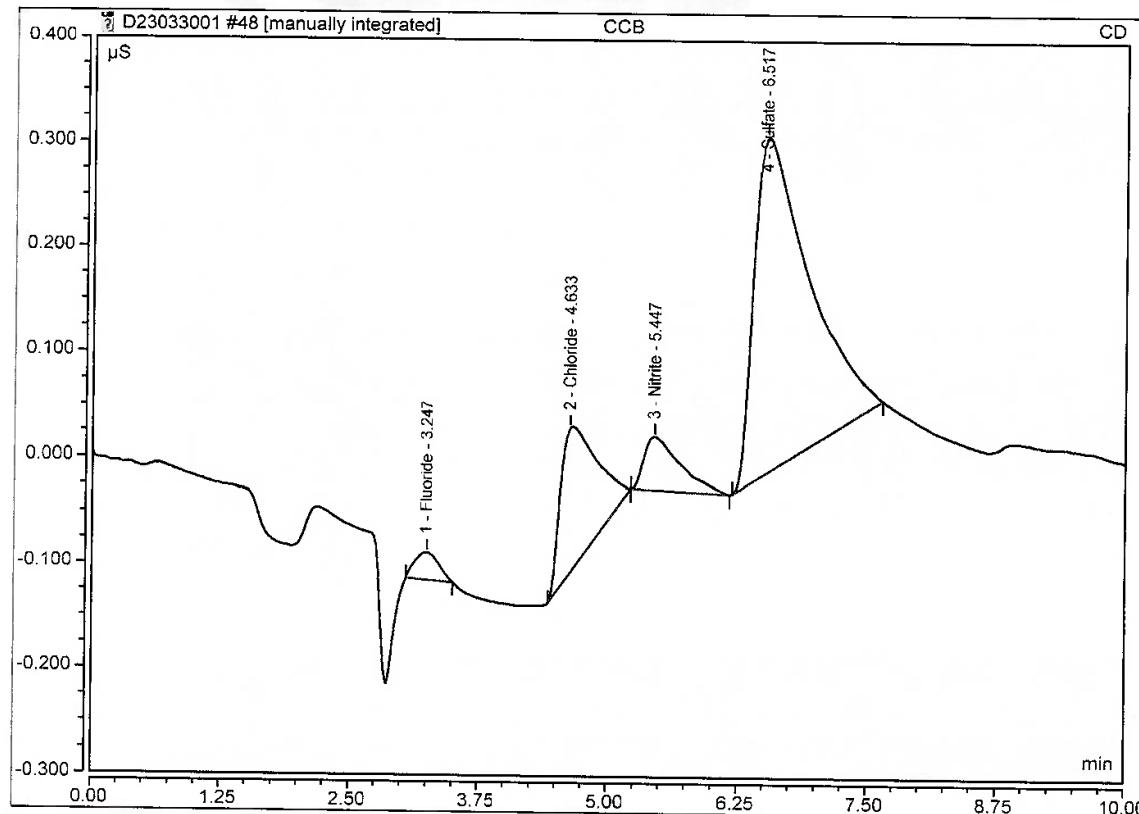
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 23:17	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.25	Fluoride	BMB	0.007	0.026	0.0548
2	4.63	Chloride	BMb	0.054	0.137	0.4090
3	5.45	Nitrite	bMB	0.020	0.050	0.0486
4	6.52	Sulfate	BMB*	0.196	0.319	0.9722
<b>TOTAL:</b>				0.28	0.53	1.48



Anion/Integration

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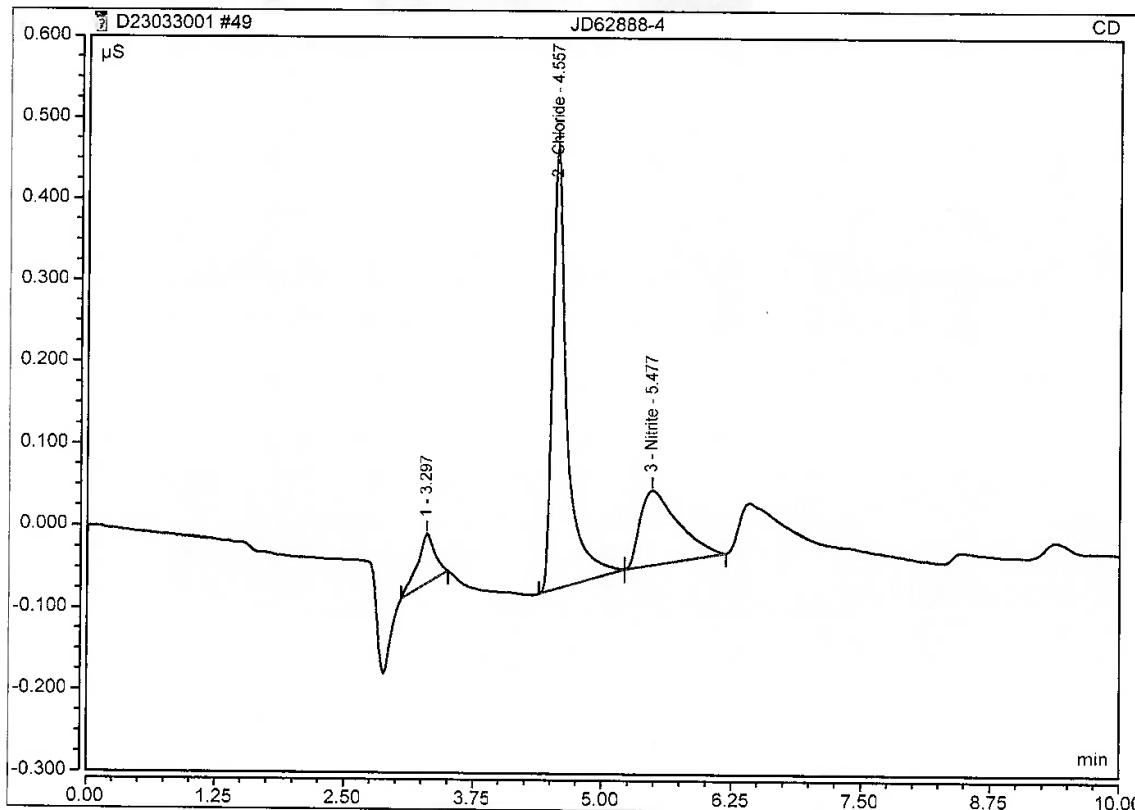
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### Peak Integration Report

Sample Name:	JD62888-4	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	30-Mar-2023 / 23:30	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.56	Chloride	BMb	0.087	0.544	0.5142
3	5.48	Nitrite	bMB	0.038	0.090	0.1709
<b>TOTAL:</b>				0.13	0.63	0.69



Anion/Integration

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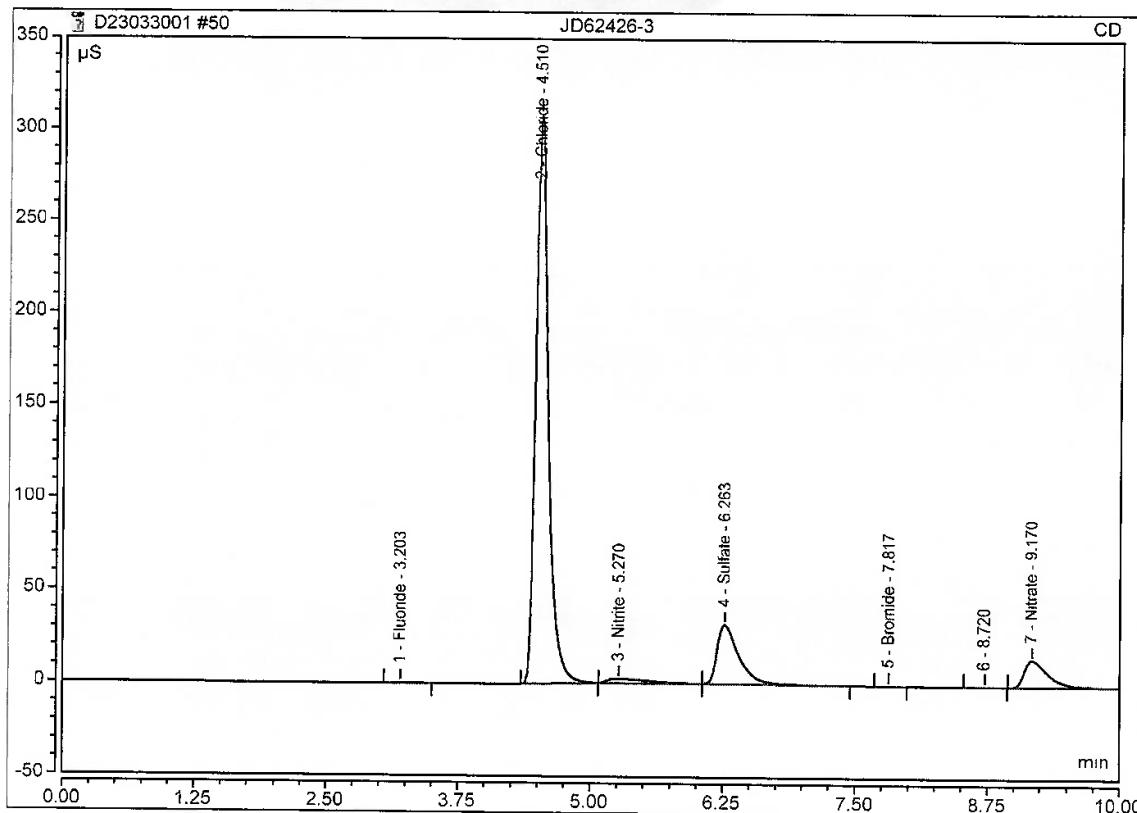
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 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

<b>Sample Name:</b>	JD62426-3	<b>Inj. Vol.:</b>	5000.00
<b>Injection Type:</b>	Unknown	<b>Dilution Factor:</b>	1.0000
<b>Instrument Method:</b>	Anions_012919	<b>Operator:</b>	Chemistry
<b>Inj. Date / Time:</b>	30-Mar-2023 / 23:43	<b>Run Time:</b>	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.20	Fluoride	BMB	0.027	0.150	0.1023
2	4.51	Chloride	BMB	42.725	308.972	134.7448
3	5.27	Nitrite	BMB	1.013	2.312	6.8259
4	6.26	Sulfate	BMB	7.498	32.048	33.1919
5	7.82	Bromide	BMB	0.002	0.016	0.0992
7	9.17	Nitrate	BMB	3.623	14.298	30.5518
<b>TOTAL:</b>				54.89	357.80	205.52



Anion/Integration

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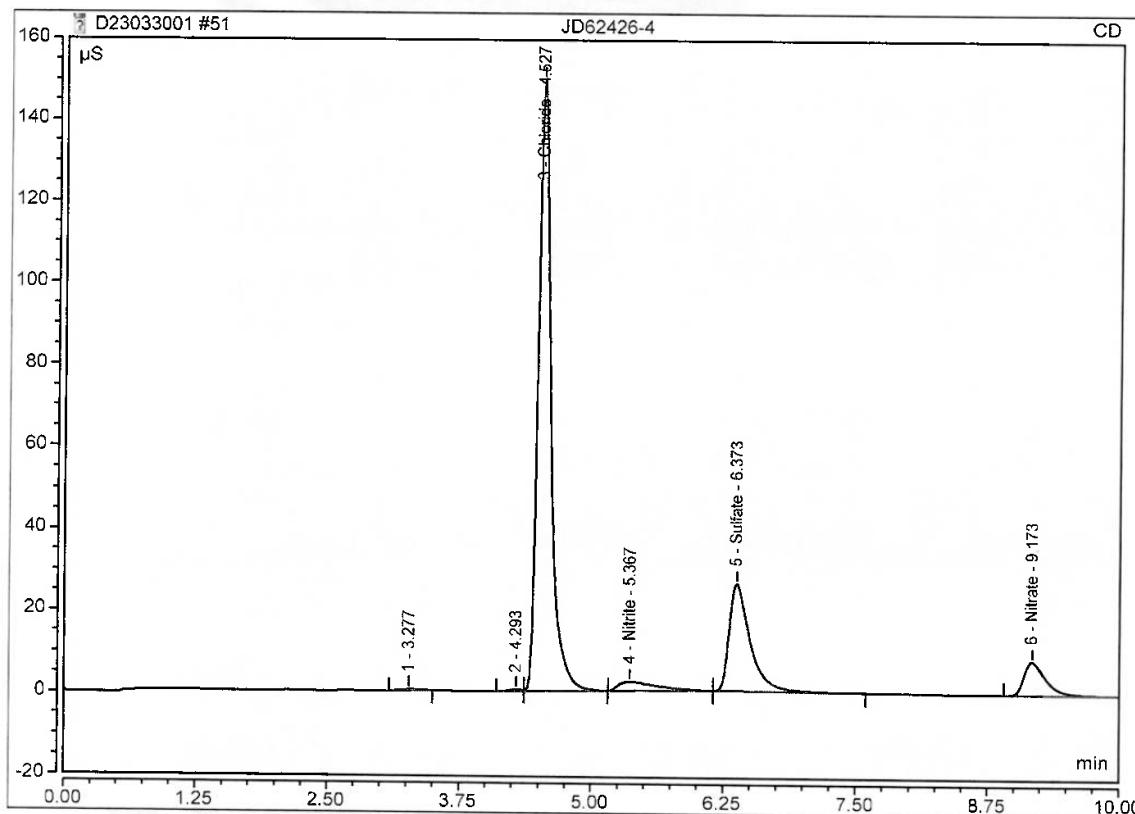
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62426-4	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 13:46	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
3	4.53	Chloride	MB	20.933	149.473	66.1383
4	5.37	Nitrite	BMB	0.940	2.137	6.3261
5	6.37	Sulfate	bMB	6.017	26.139	26.6584
6	9.17	Nitrate	BMB	1.961	8.032	16.5025
<b>TOTAL:</b>				29.85	185.78	115.63



Anion/Integration

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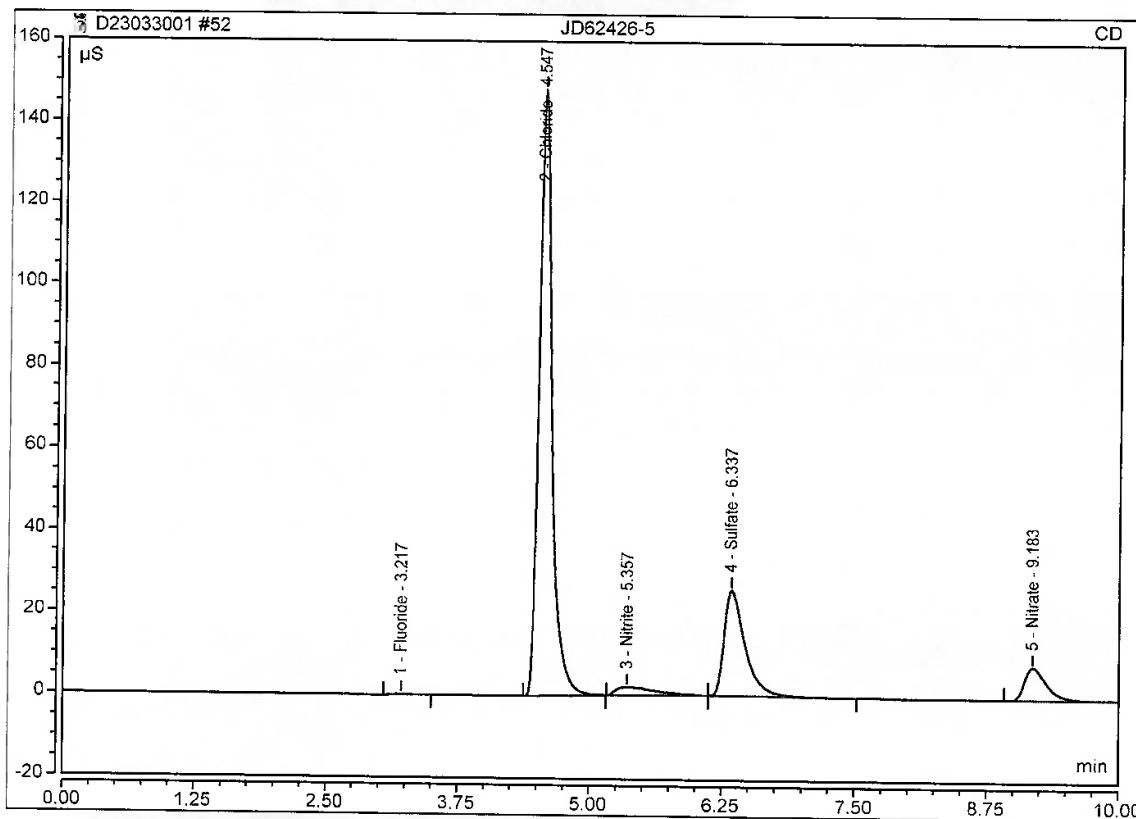
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62426-5	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 13:59	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.035	0.196	0.1223
2	4.55	Chloride	BMB	20.944	148.185	66.1731
3	5.36	Nitrite	BMb	0.877	2.050	5.9004
4	6.34	Sulfate	bMB	5.877	25.761	26.0409
5	9.18	Nitrate	BMB	1.958	8.001	16.4836
<b>TOTAL:</b>				29.69	184.19	114.72



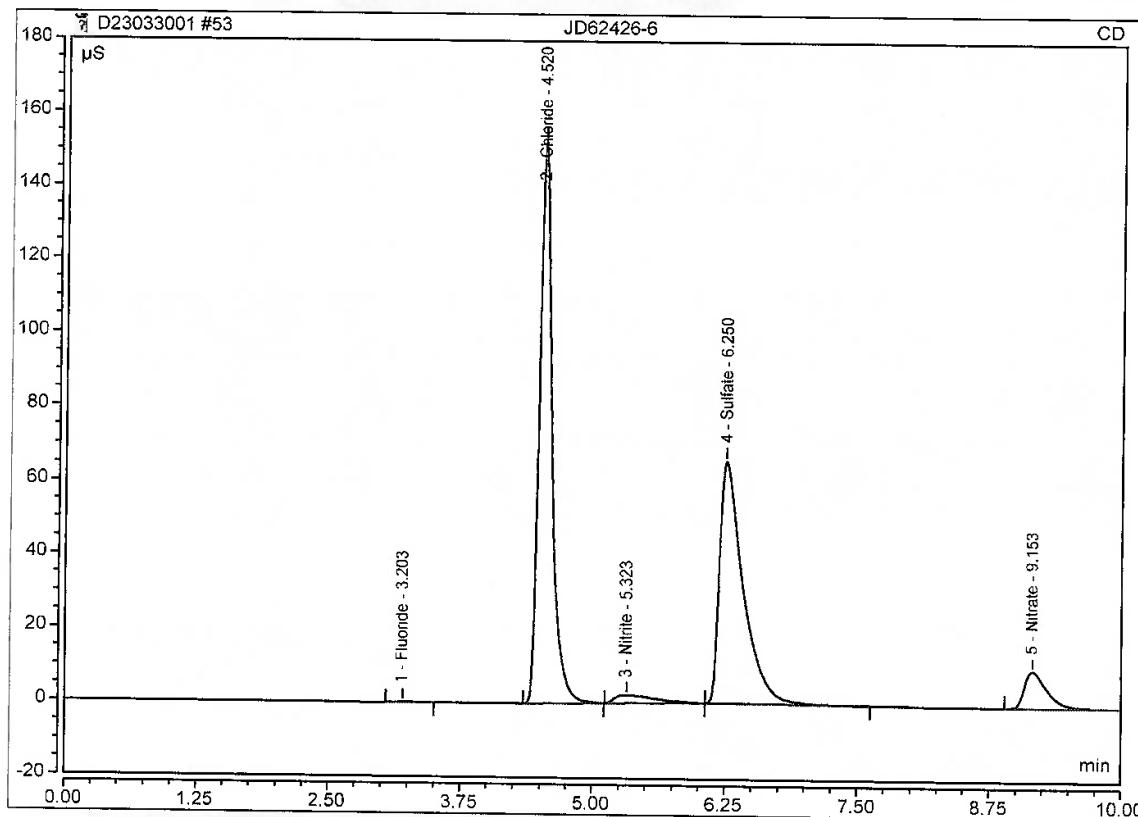
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62426-6	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 14:12	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.20	Fluoride	BMB	0.030	0.181	0.1101
2	4.52	Chloride	BMB	21.599	152.627	68.2371
3	5.32	Nitrite	BMb	0.962	2.231	6.4786
4	6.25	Sulfate	bMB	16.966	65.826	74.9741
5	9.15	Nitrate	BMB	2.500	9.981	21.0568
<b>TOTAL:</b>				42.06	230.85	170.86



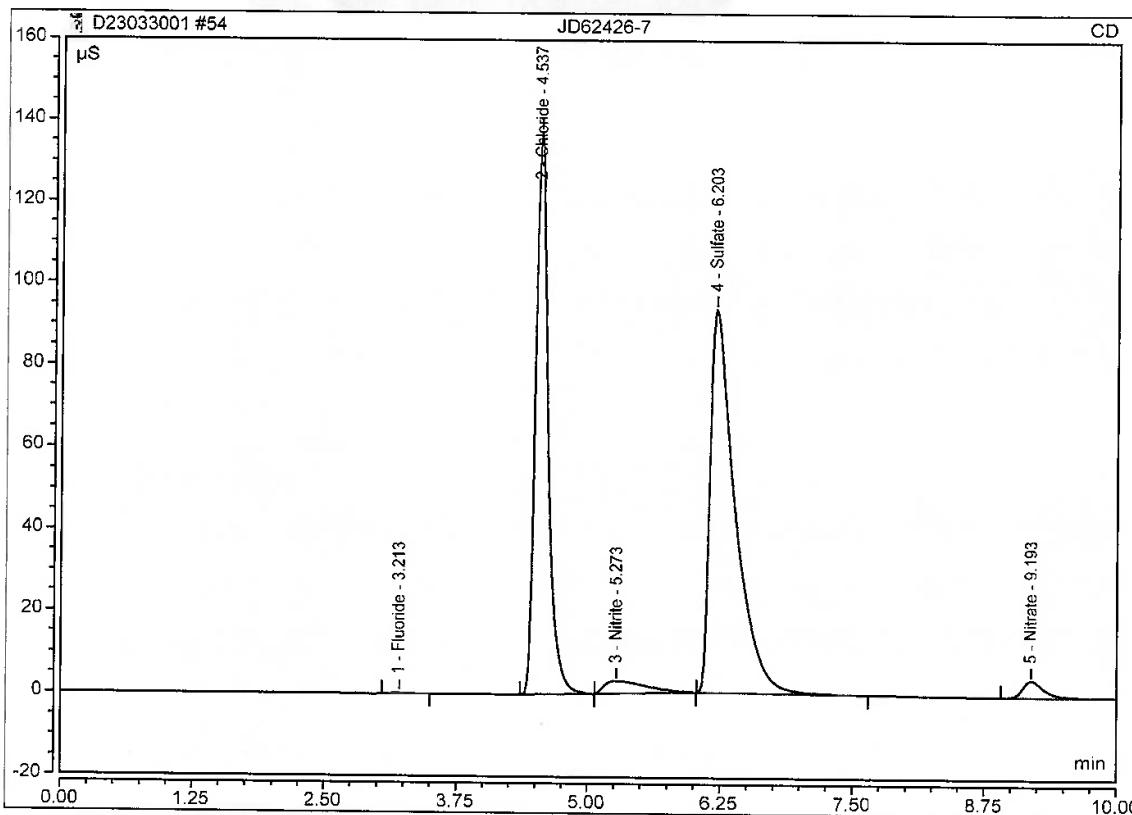
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### Peak Integration Report

Sample Name:	JD62426-7	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 14:25	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.038	0.218	0.1306
2	4.54	Chloride	BMB	19.115	137.514	60.4161
3	5.27	Nitrite	BMB	1.387	3.019	9.3826
4	6.20	Sulfate	BMB	25.424	93.359	112.2960
5	9.19	Nitrate	BMB	0.950	4.063	7.9657
<b>TOTAL:</b>				46.91	238.17	190.19



Anion/Integration

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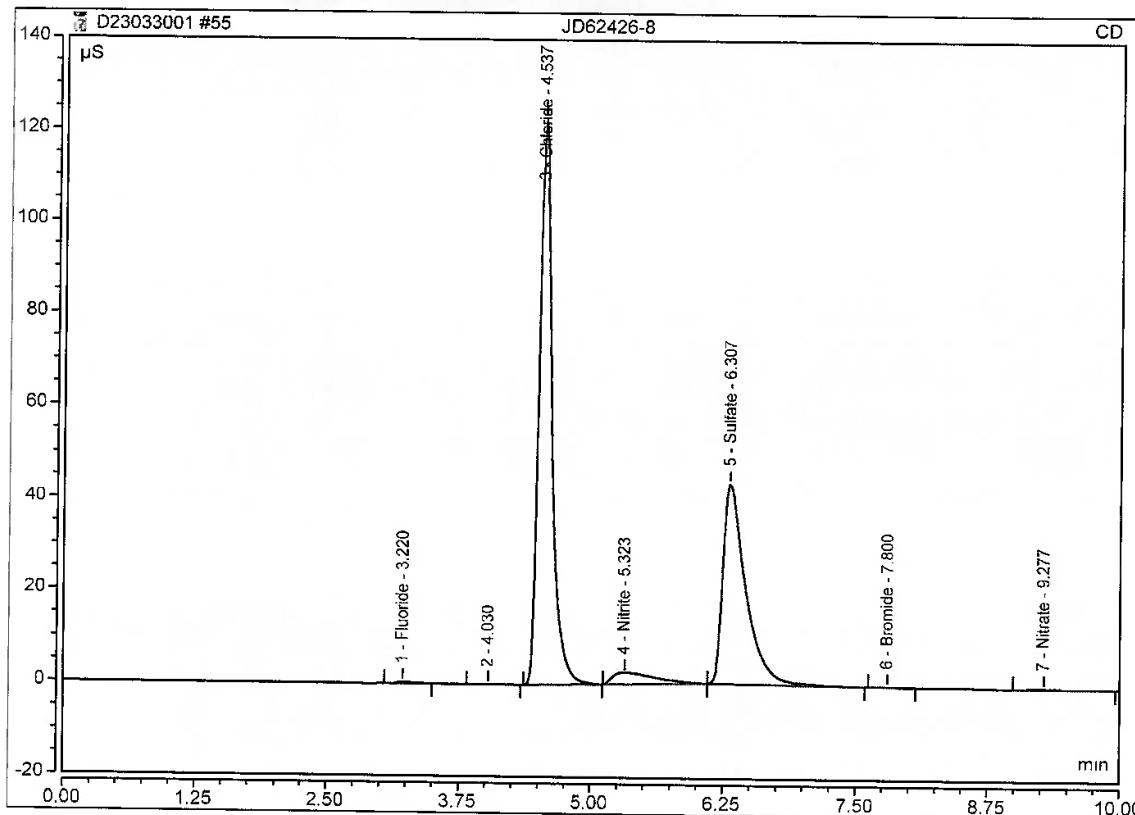
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62426-8	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 14:38	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.047	0.293	0.1522
3	4.54	Chloride	BMB	17.288	122.822	54.6638
4	5.32	Nitrite	BMb	1.098	2.448	7.4110
5	6.31	Sulfate	bMB	10.708	43.460	47.3575
6	7.80	Bromide	BMB	0.007	0.039	0.1394
7	9.28	Nitrate	BMB	0.036	0.149	0.2414
<b>TOTAL:</b>				29.18	169.21	109.97



Anion/Integration

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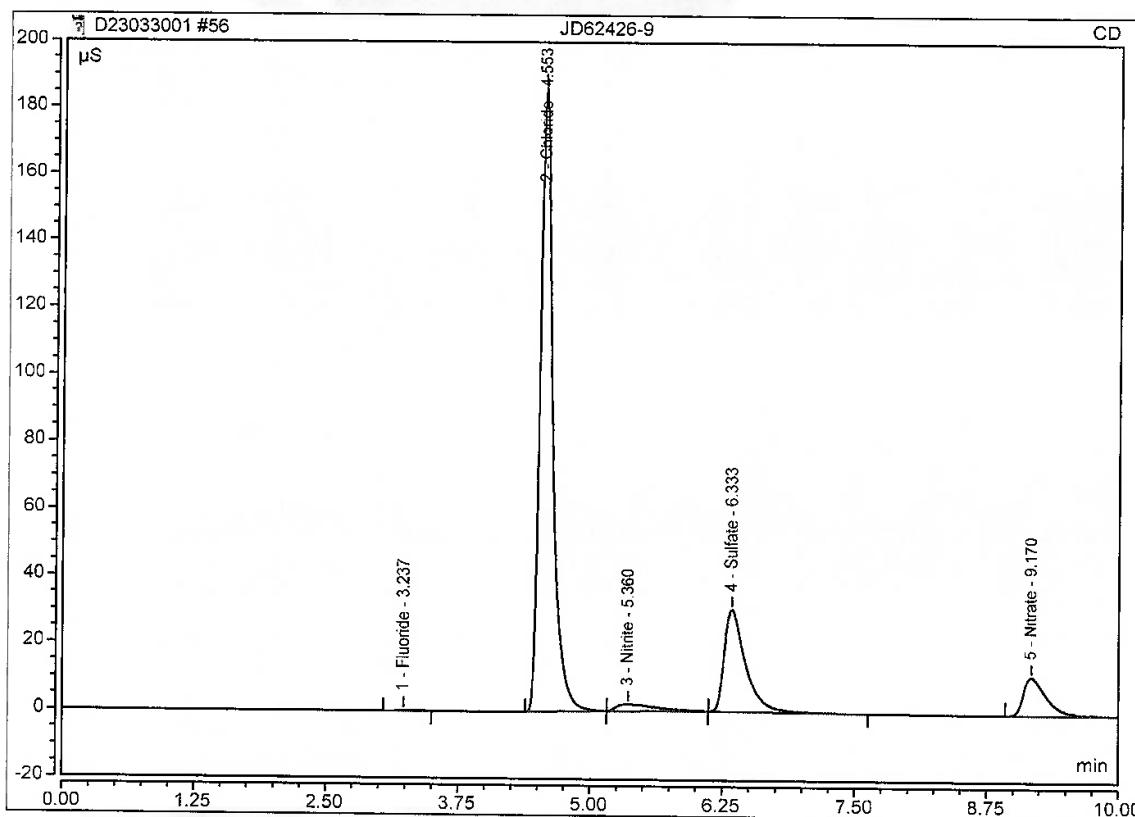
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### Peak Integration Report

Sample Name:	JD62426-9	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 14:51	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.24	Fluoride	BMB	0.039	0.208	0.1317
2	4.55	Chloride	BMB	26.355	186.523	83.2086
3	5.36	Nitrite	BMb	0.860	2.017	5.7808
4	6.33	Sulfate	bMB	7.022	30.359	31.0919
5	9.17	Nitrate	BMB	2.784	11.220	23.4608
TOTAL:				37.06	230.33	143.67



Anion/Integration

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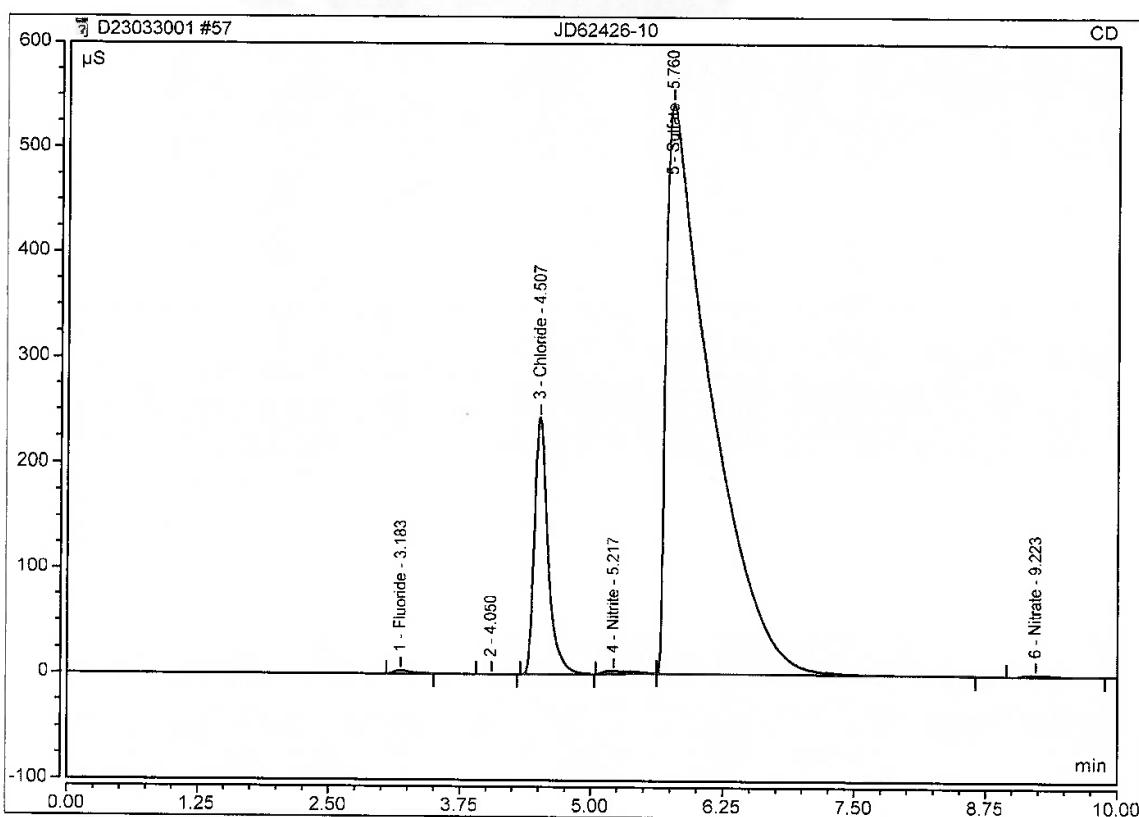
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62426-10	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 15:04	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.18	Fluoride	BMB	0.506	3.202	1.2613
3	4.51	Chloride	BMB	36.238	243.851	114.3219
4	5.22	Nitrite	BMB	0.939	2.859	6.3201
5	5.76	Sulfate	BMB	265.796	542.364	1172.9944
6	9.22	Nitrate	BMB	0.400	1.661	3.3120
<b>TOTAL:</b>				303.88	793.94	1298.21



Anion/Integration

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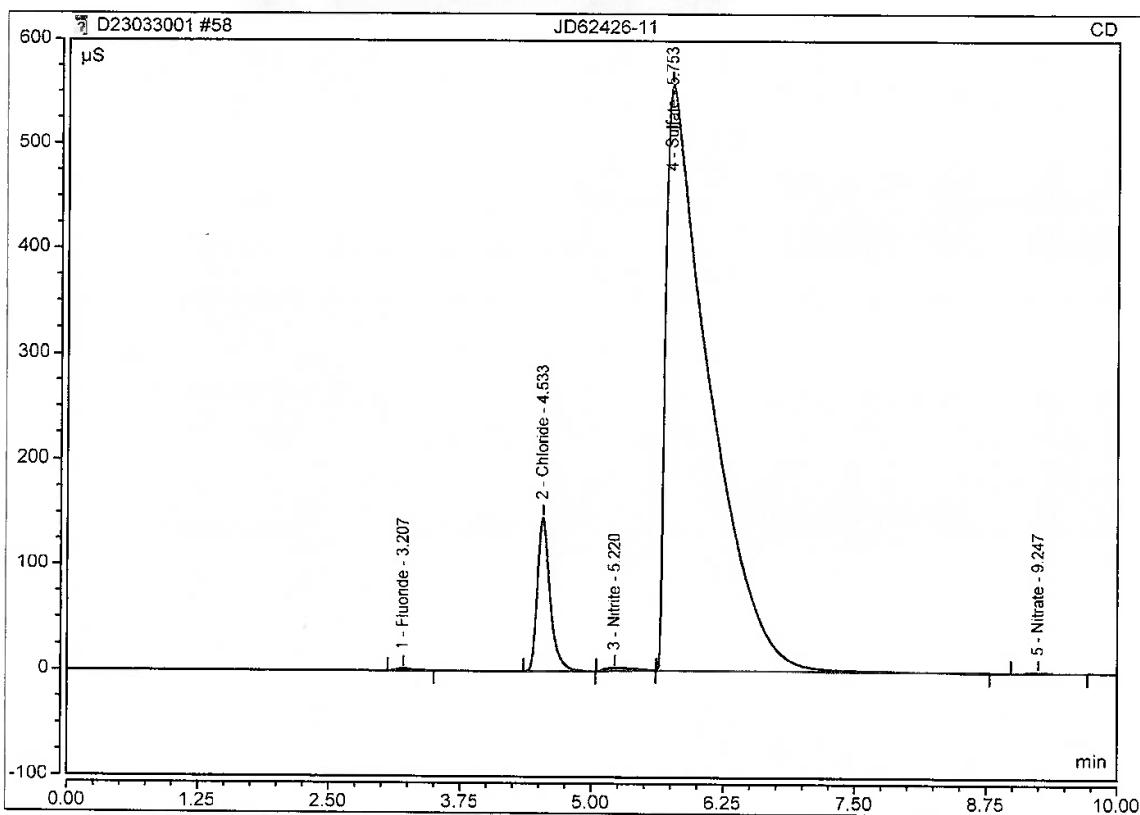
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### Peak Integration Report

Sample Name:	JD62426-11	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 15:17	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.324	2.050	0.8222
2	4.53	Chloride	BMB	20.913	146.179	66.0772
3	5.22	Nitrite	BMB	0.964	2.962	6.4961
4	5.75	Sulfate	BMB	264.571	556.527	1167.5879
5	9.25	Nitrate	BMB	0.131	0.560	1.0389
<b>TOTAL:</b>				286.90	708.28	1242.02



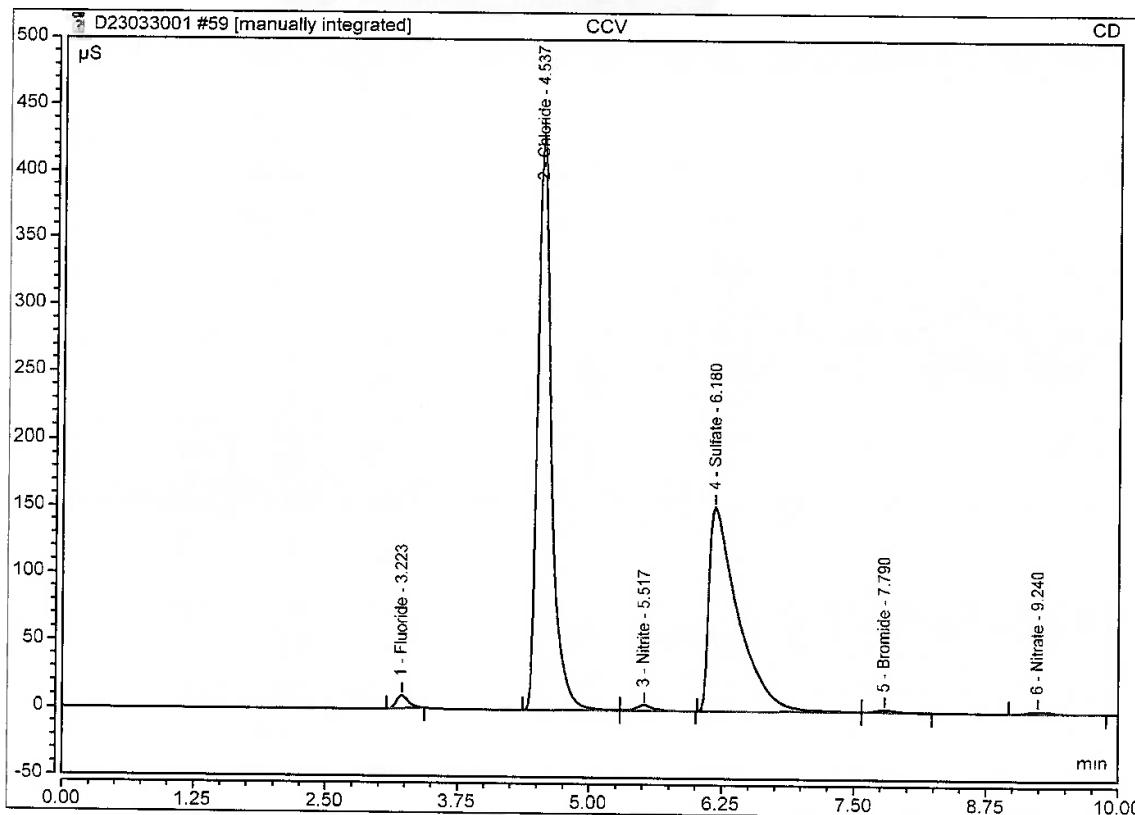
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 Sequence: D23033001

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### Peak Integration Report

<b>Sample Name:</b>	CCV	<b>Inj. Vol.:</b>	5000.00
<b>Injection Type:</b>	Unknown	<b>Dilution Factor:</b>	1.0000
<b>Instrument Method:</b>	Anions_012919	<b>Operator:</b>	Chemistry
<b>Inj. Date / Time:</b>	31-Mar-2023 / 15:30	<b>Run Time:</b>	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB*	1.270	9.430	3.1100
2	4.54	Chloride	BMb	64.659	430.604	203.7963
3	5.52	Nitrite	bMB	0.698	3.931	4.6755
4	6.18	Sulfate	8Mb	47.704	152.247	210.6119
5	7.79	Bromide	bMB	0.342	1.770	3.0088
6	9.24	Nitrate	BMB	0.440	1.797	3.6548
<b>TOTAL:</b>				115.11	599.78	428.86



Anion/Integration

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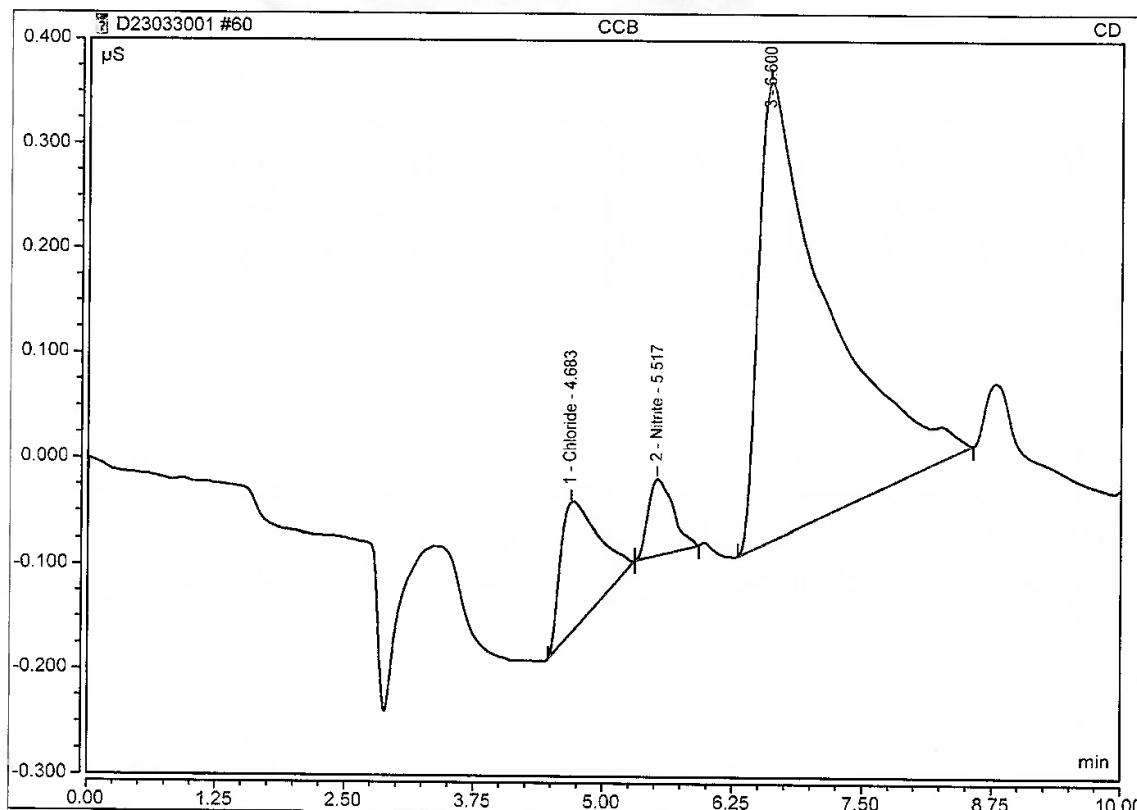
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### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 15:43	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	4.68	Chloride	BMb	0.051	0.124	0.3987
2	5.52	Nitrite	bMB	0.020	0.072	0.0463
TOTAL:				0.07	0.20	0.44



Anion/Integration

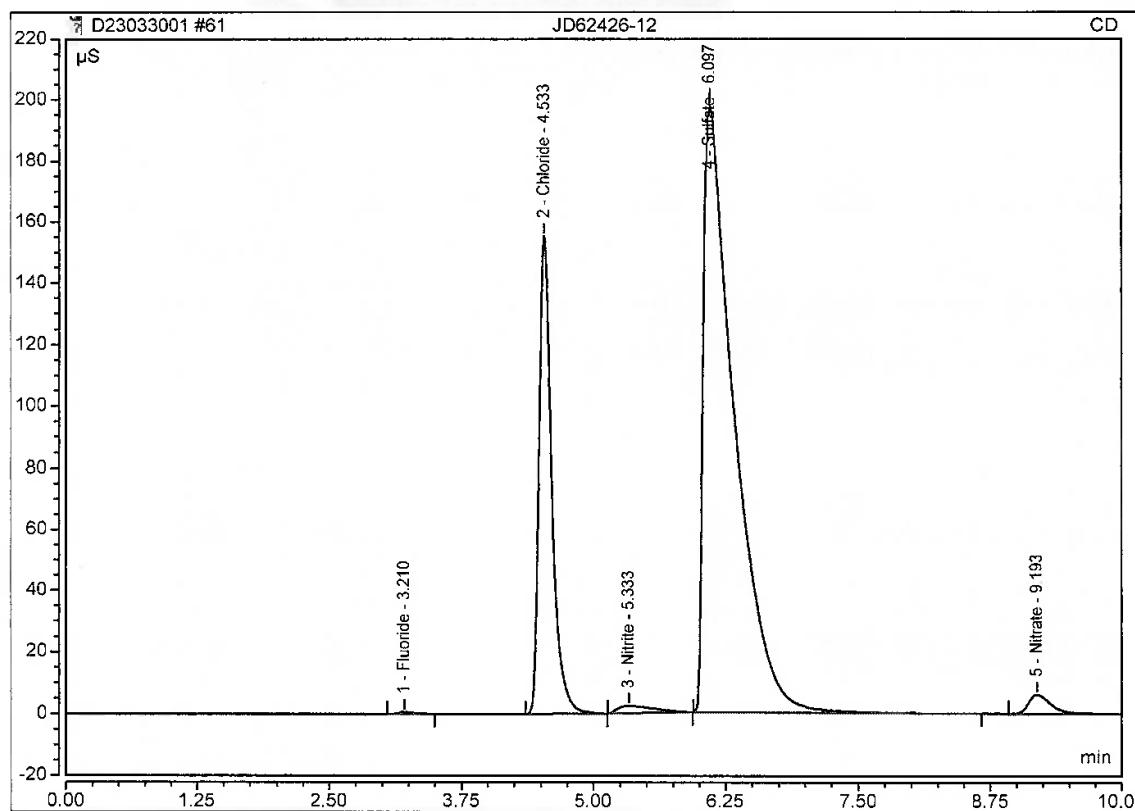
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Logged on User: Chemistry  
Instrument: Integron\_1  
Sequence: D23033001

## Peak Integration Report

Sample Name:	JD62426-12	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 15:56	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.072	0.483	0.2122
2	4.53	Chloride	BMB	21.466	155.368	67.8184
3	5.33	Nitrite	BMB	0.884	2.256	5.9459
4	6.10	Sulfate	BMB	66.669	197.625	294.3010
5	9.19	Nitrate	BMB	1.462	6.079	12.2885
<b>TOTAL:</b>				90.55	361.81	380.57



Anion/Integration

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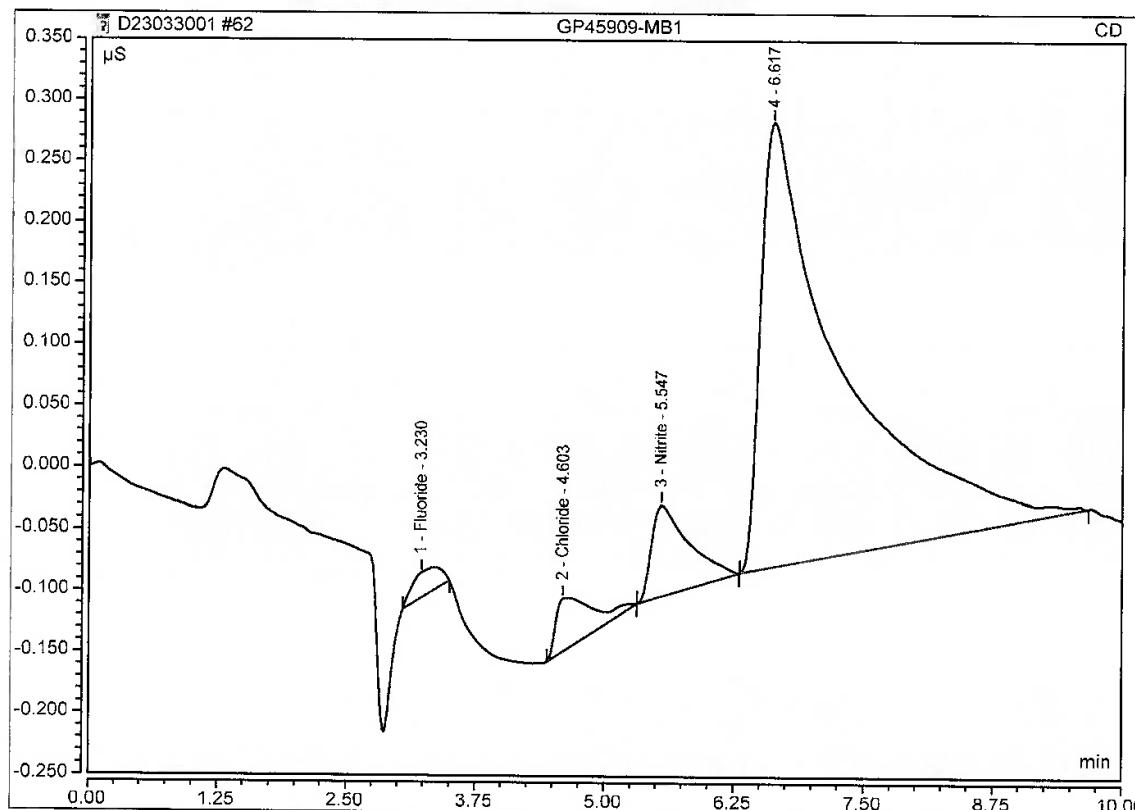
Logged on User: Chemistry  
 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45909-MB1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 16:09	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.23	Fluoride	BMB	0.007	0.021	0.0533
2	4.60	Chloride	BMb	0.016	0.043	0.2910
3	5.55	Nitrite	bMB	0.031	0.074	0.1221
<b>TOTAL:</b>				0.05	0.14	0.47



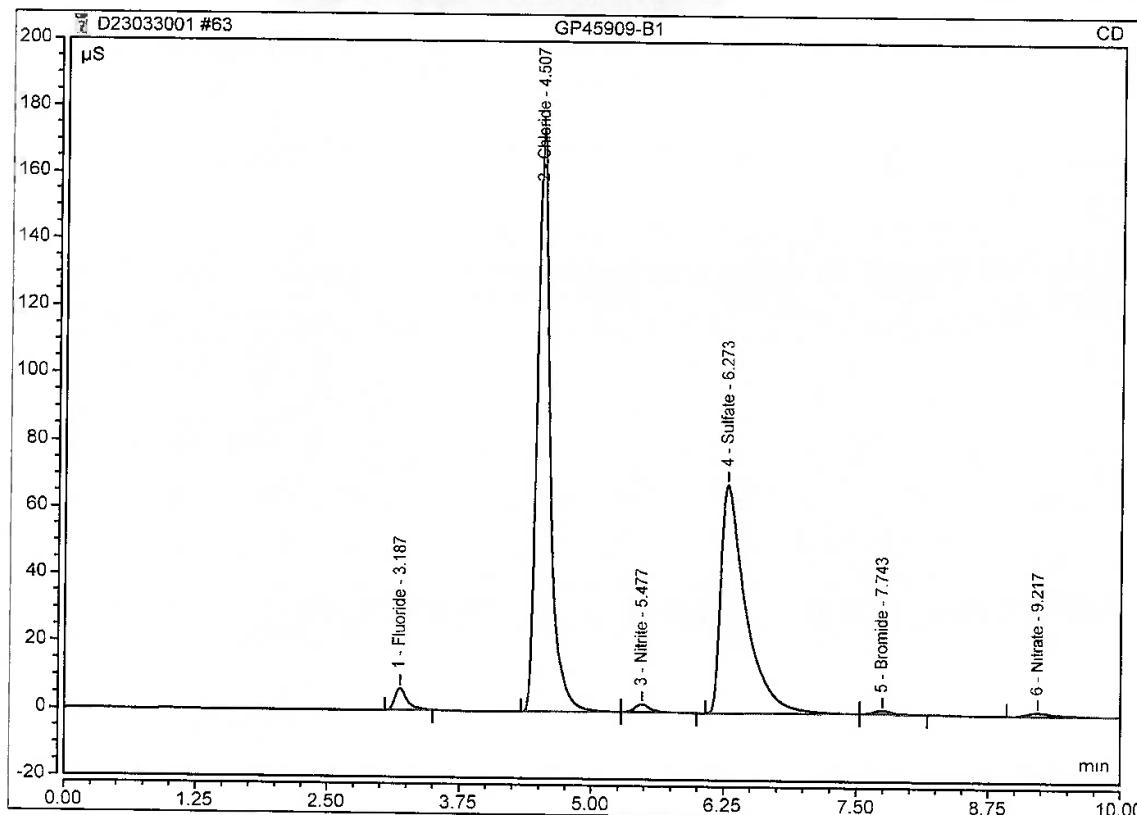
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45909-B1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Ini. Date / Time:	31-Mar-2023 / 16:22	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.19	Fluoride	BMB	0.860	6.549	2.1170
2	4.51	Chloride	BMb	24.968	173.842	78.8415
3	5.48	Nitrite	bMB	0.379	2.366	2.5016
4	6.27	Sulfate	BMb	18.465	68.212	81.5869
5	7.74	Bromide	bMB	0.202	1.064	1.8068
6	9.22	Nitrate	BMB	0.266	1.047	2.1852
<b>TOTAL:</b>				45.14	253.08	169.04



Anion/Integration

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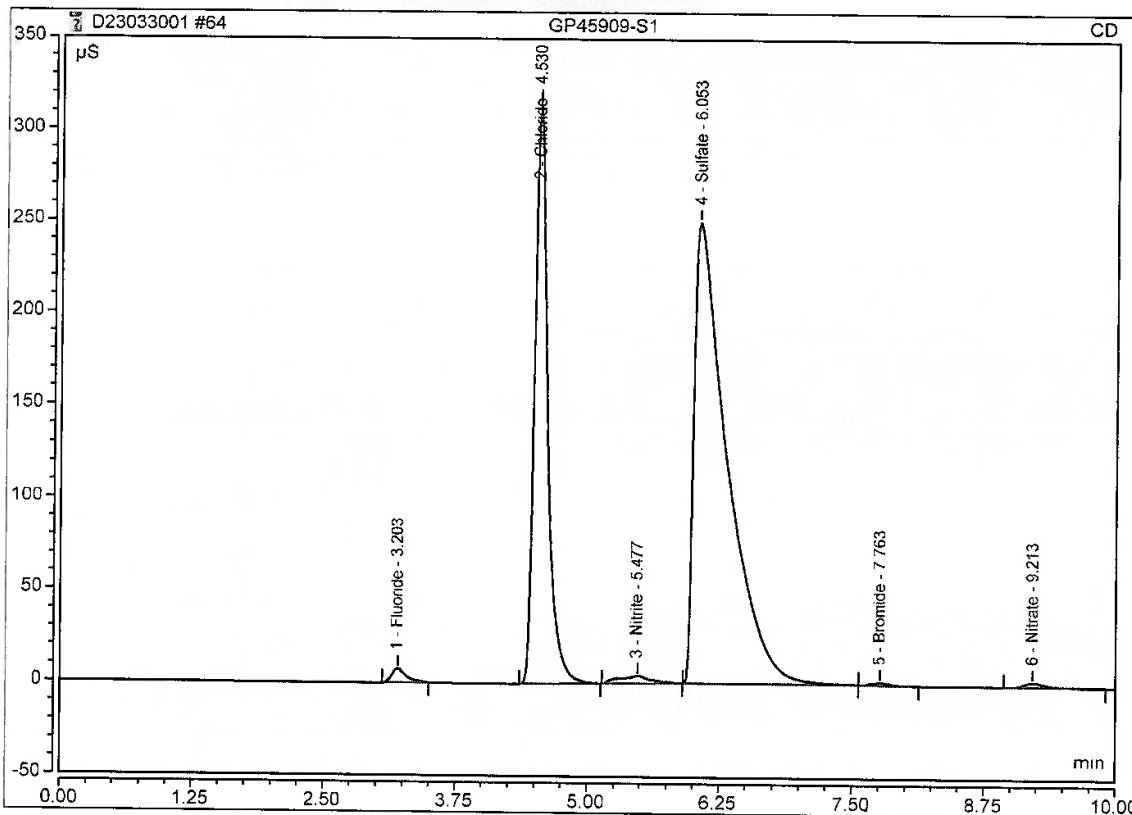
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 Instrument: Integriion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45909-S1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 16:35	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.20	Fluoride	BMB	1.075	7.365	2.6364
2	4.53	Chloride	BMB	44.892	314.147	141.5670
3	5.48	Nitrite	BMB	1.280	3.739	8.6505
4	6.05	Sulfate	BMB	90.189	249.601	398.0876
5	7.76	Bromide	bMB	0.225	1.200	2.0067
6	9.21	Nitrate	BMB	0.513	2.183	4.2739
<b>TOTAL:</b>				138.17	578.24	557.22



Anion/Integration

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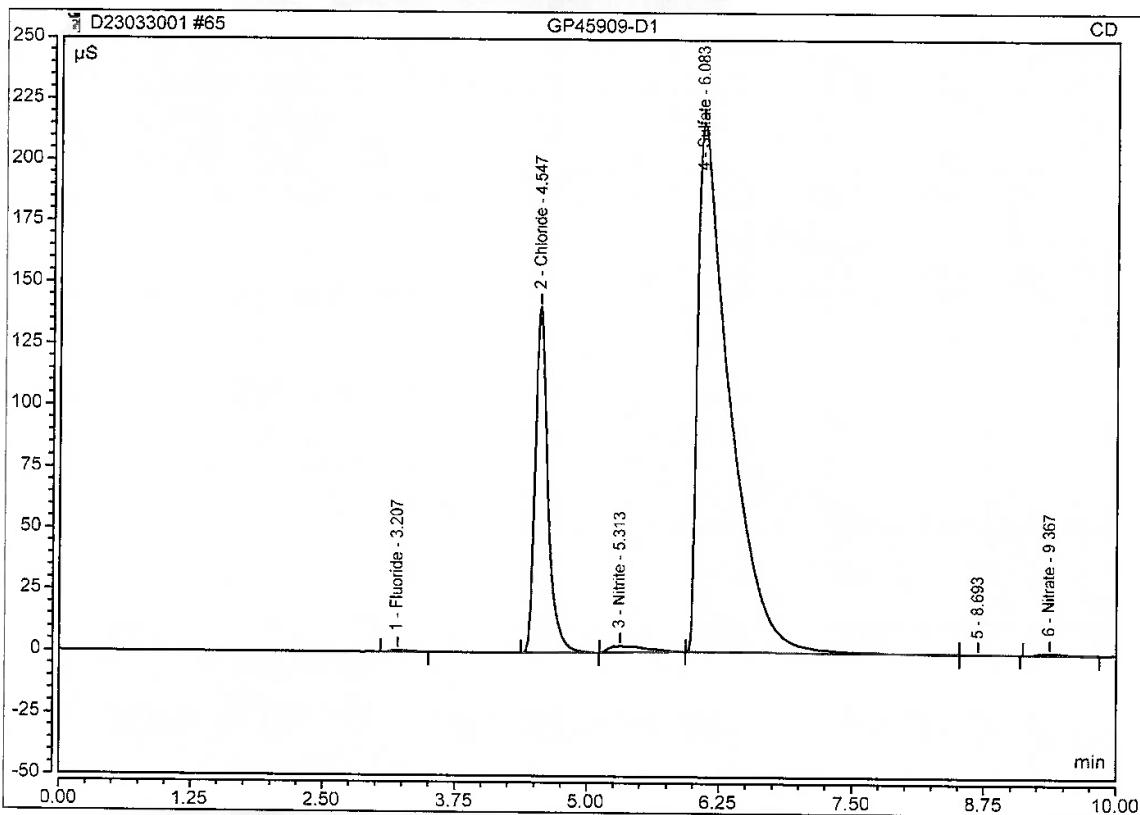
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 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	GP45909-D1	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 16:48	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.112	0.767	0.3073
2	4.55	Chloride	BMB	19.513	140.487	61.6694
3	5.31	Nitrite	BMB	1.068	2.604	7.2013
4	6.08	Sulfate	BMb	75.549	215.960	333.4830
6	9.37	Nitrate	BMB	0.203	0.889	1.6521
<b>TOTAL:</b>				96.44	360.71	404.31



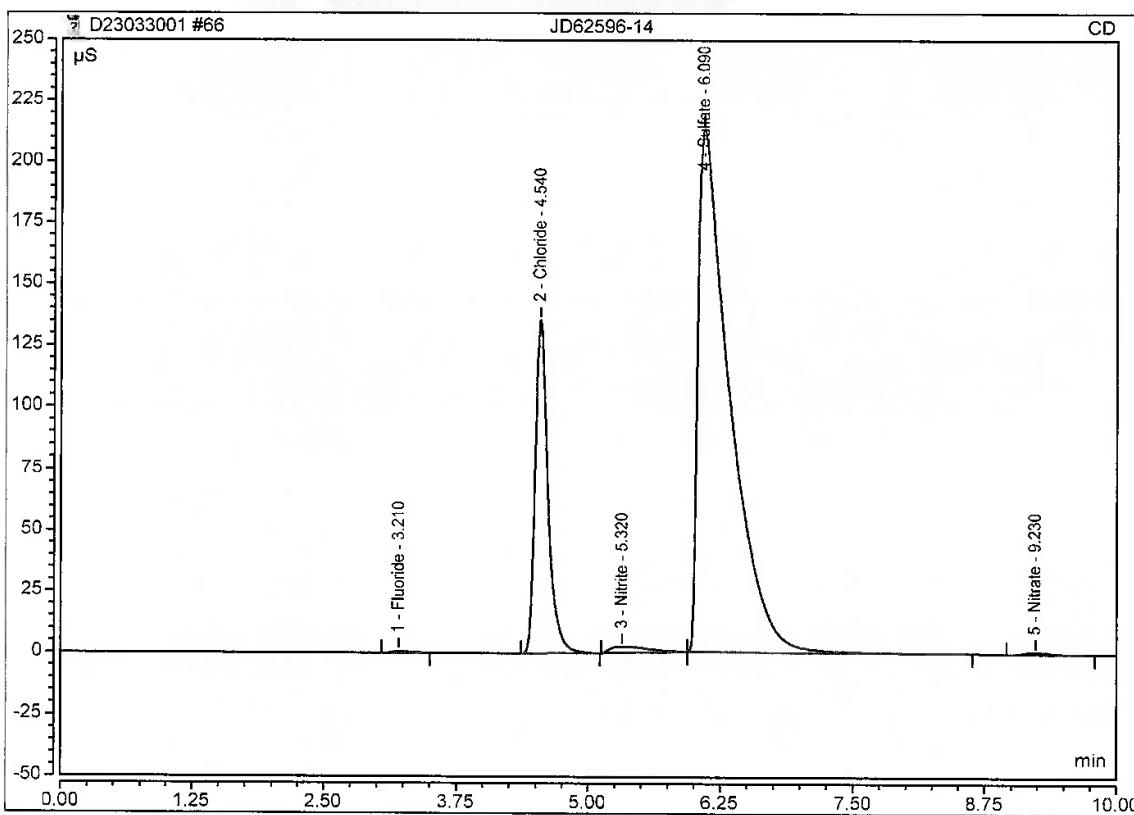
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 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-14	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 17:01	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\cdot}\text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.110	0.734	0.3038
2	4.54	Chloride	BMB	19.179	135.527	60.6188
3	5.32	Nitrite	BMB	1.062	2.604	7.1639
4	6.09	Sulfate	BMB	73.797	213.028	325.7548
5	9.23	Nitrate	BMB	0.199	0.857	1.6150
<b>TOTAL:</b>				94.35	352.75	395.46



Anion/Integration

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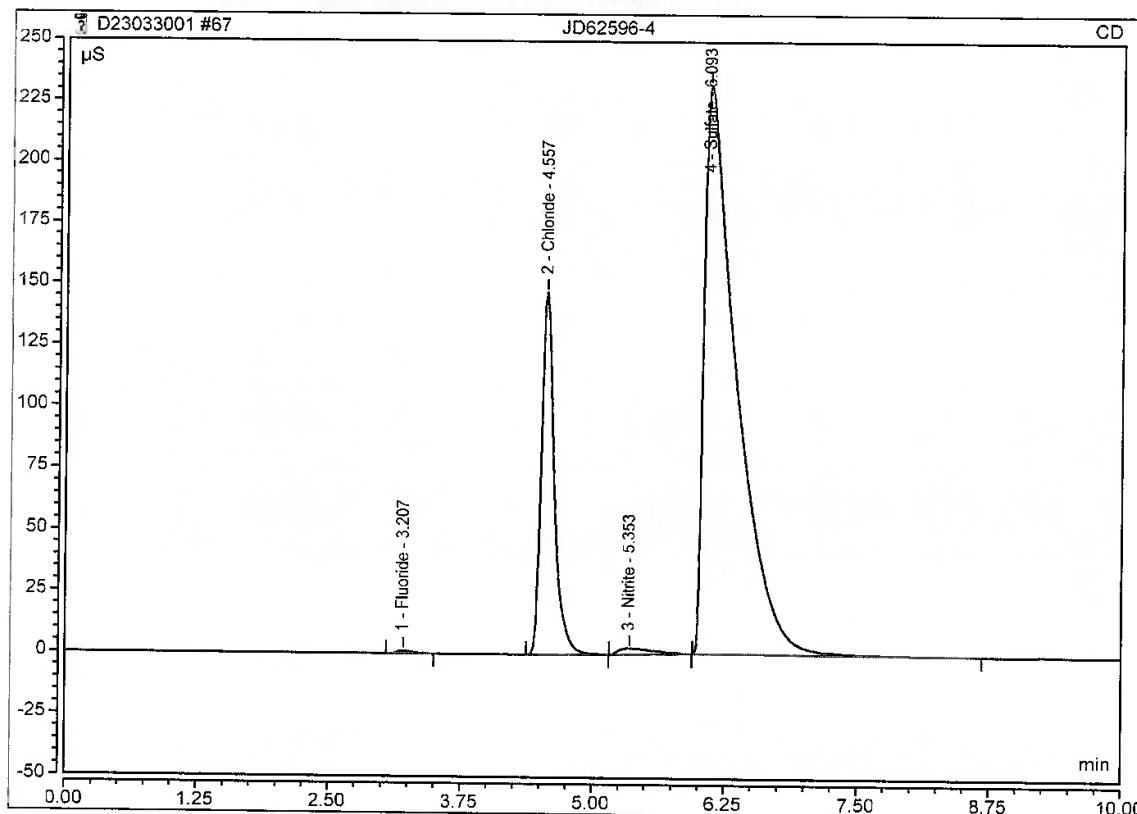
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-4	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 17:14	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.174	1.202	0.4593
2	4.56	Chloride	BMB	21.106	147.209	66.6857
3	5.35	Nitrite	BMB	0.965	2.462	6.4968
4	6.09	Sulfate	BMB	82.344	231.832	363.4708
TOTAL:				104.59	382.70	437.11



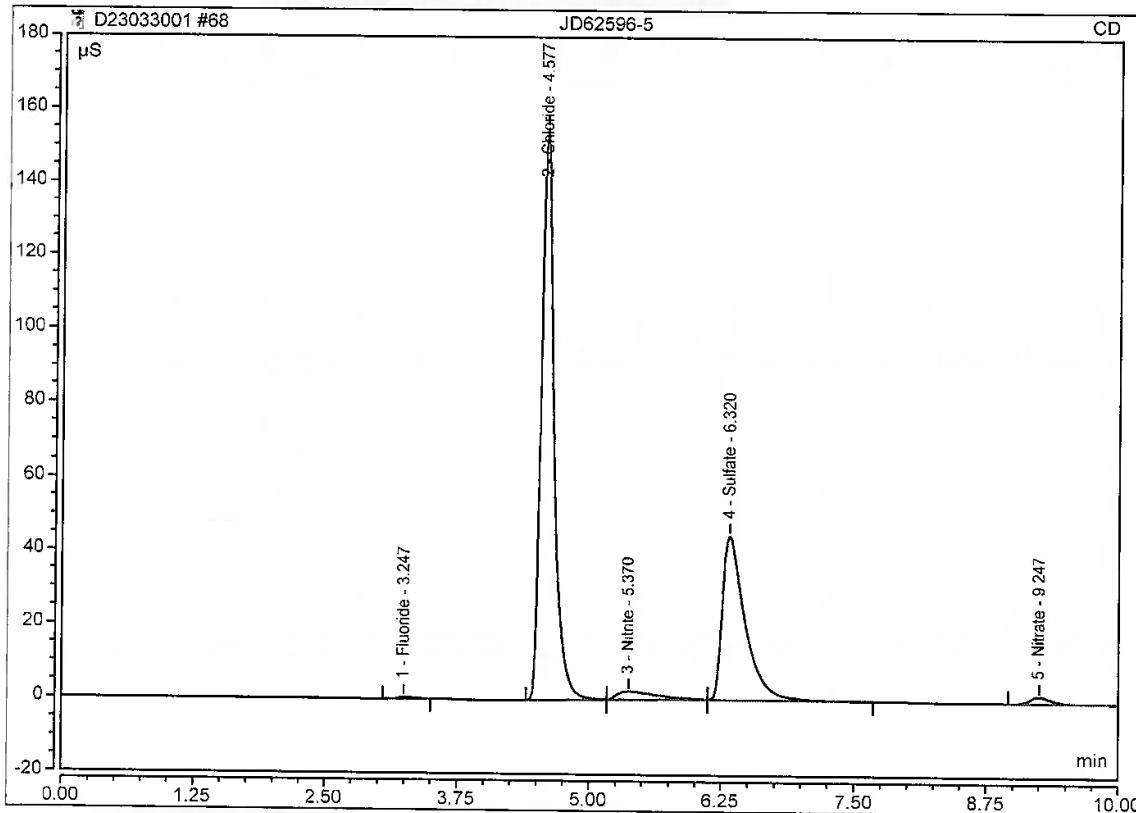
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-5	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 17:27	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.25	Fluoride	BMB	0.064	0.395	0.1922
2	4.58	Chloride	BMB	21.793	155.336	68.8473
3	5.37	Nitrite	BMb	0.937	2.168	6.3086
4	6.32	Sulfate	bMB	10.992	44.474	48.6104
5	9.25	Nitrate	BMB	0.428	1.818	3.5481
<b>TOTAL:</b>				34.21	204.19	127.51



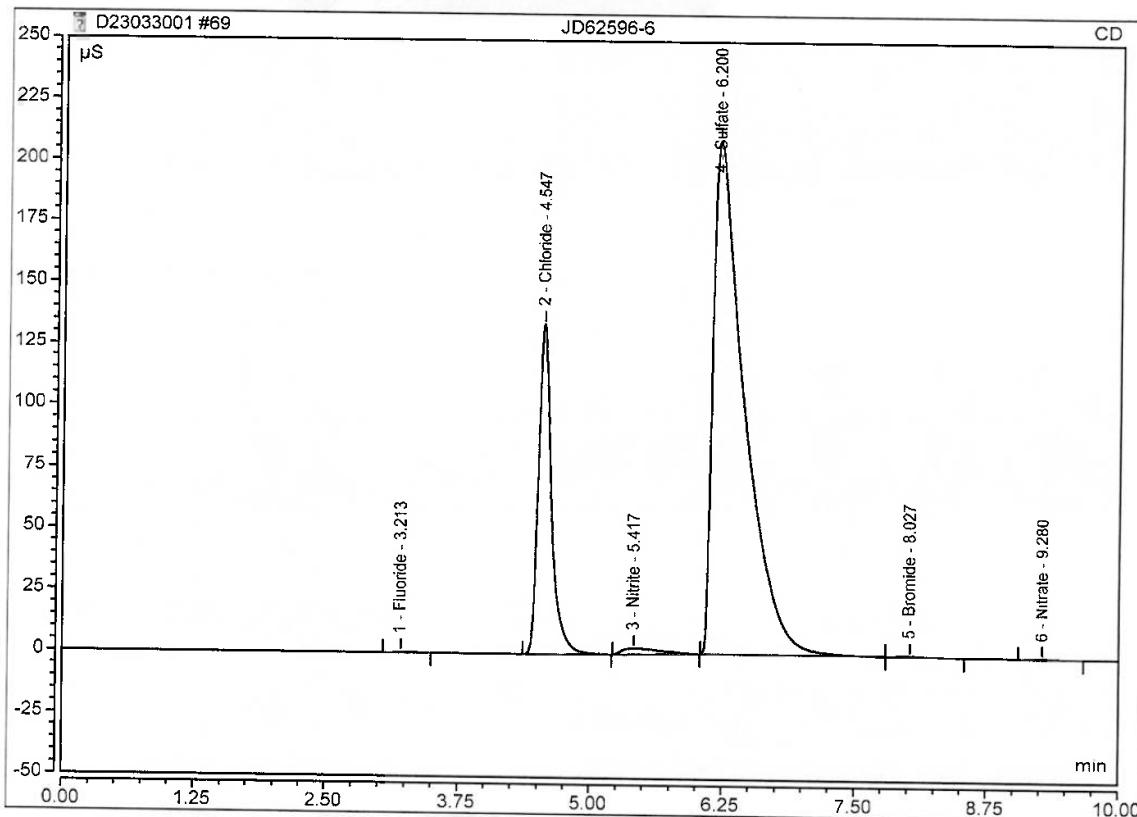
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-6	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 17:40	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.054	0.294	0.1687
2	4.55	Chloride	BMB	19.319	134.323	61.0596
3	5.42	Nitrite	BMB	1.044	2.514	7.0382
4	6.20	Sulfate	BMb	71.493	209.126	315.5875
5	8.03	Bromide	bMB	0.090	0.247	0.8500
6	9.28	Nitrate	BMB	0.049	0.221	0.3529
<b>TOTAL:</b>				92.05	346.72	385.06



Anion/Integration

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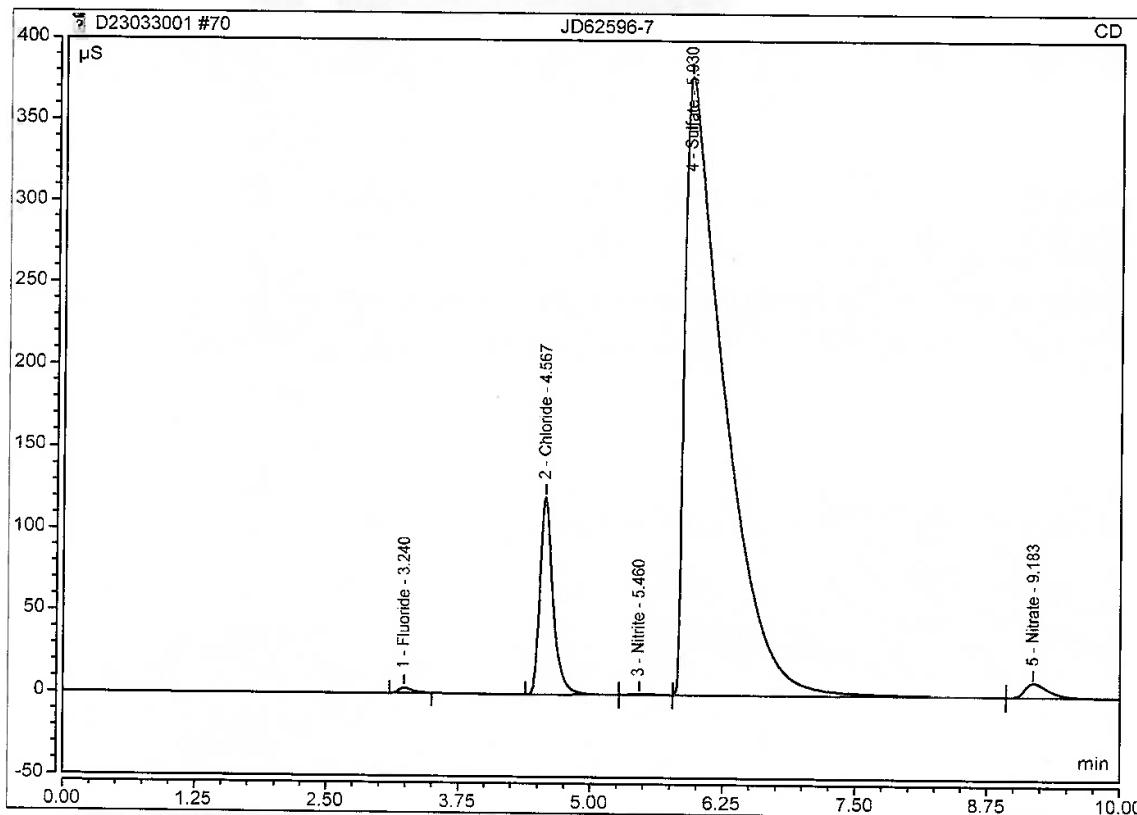
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-7	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 17:53	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
1	3.24	Fluoride	BMB	0.487	3.349	1.2147
2	4.57	Chloride	BMb	17.363	120.813	54.8995
3	5.46	Nitrite	bMB	0.125	0.523	0.7647
4	5.93	Sulfate	BMb	162.778	377.926	718.4042
5	9.18	Nitrate	bMB	2.305	8.832	19.4138
<b>TOTAL:</b>				183.06	511.44	794.70



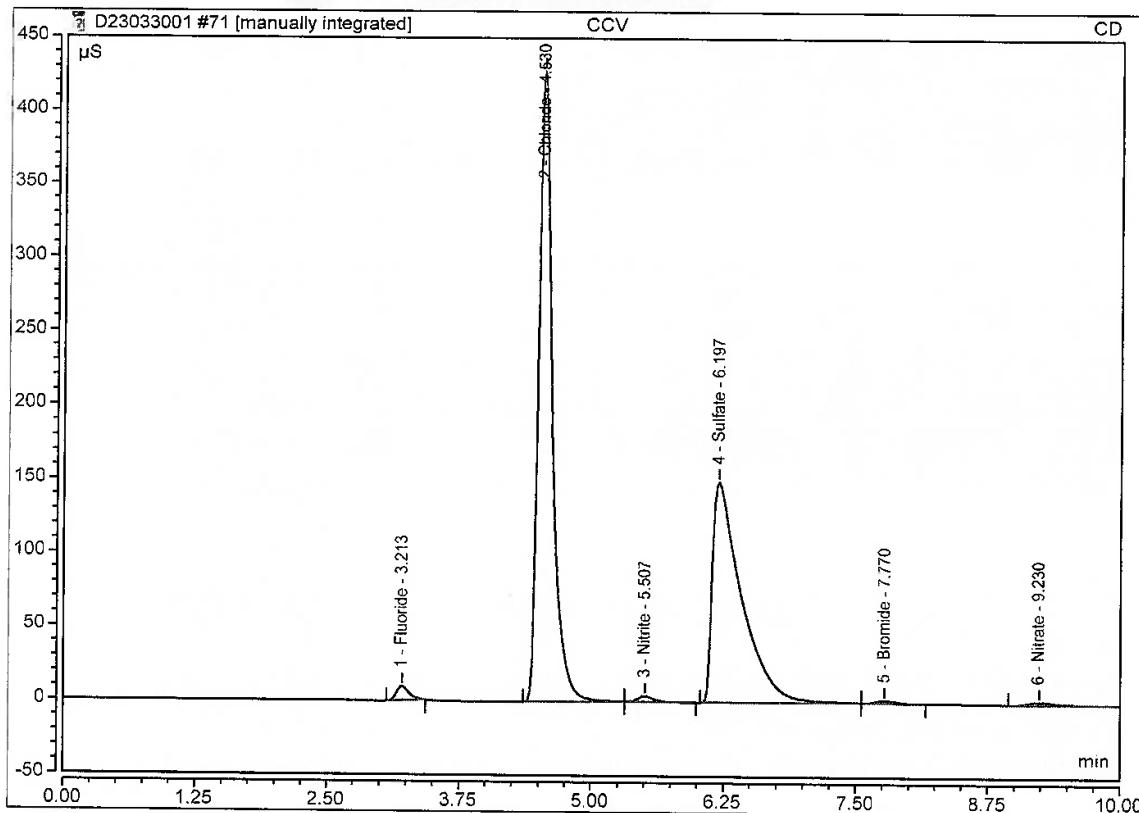
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 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCV	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 18:06	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB*	1.283	9.409	3.1394
2	4.53	Chloride	BMb	64.676	427.563	203.8504
3	5.51	Nitrite	bMB	0.638	3.871	4.2692
4	6.20	Sulfate	BMb	47.535	149.889	209.8686
5	7.77	Bromide	bMB	0.342	1.765	3.0100
6	9.23	Nitrate	BMB	0.452	1.811	3.7566
<b>TOTAL:</b>				114.93	594.31	427.89



Anion/Integration

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9.4  
9

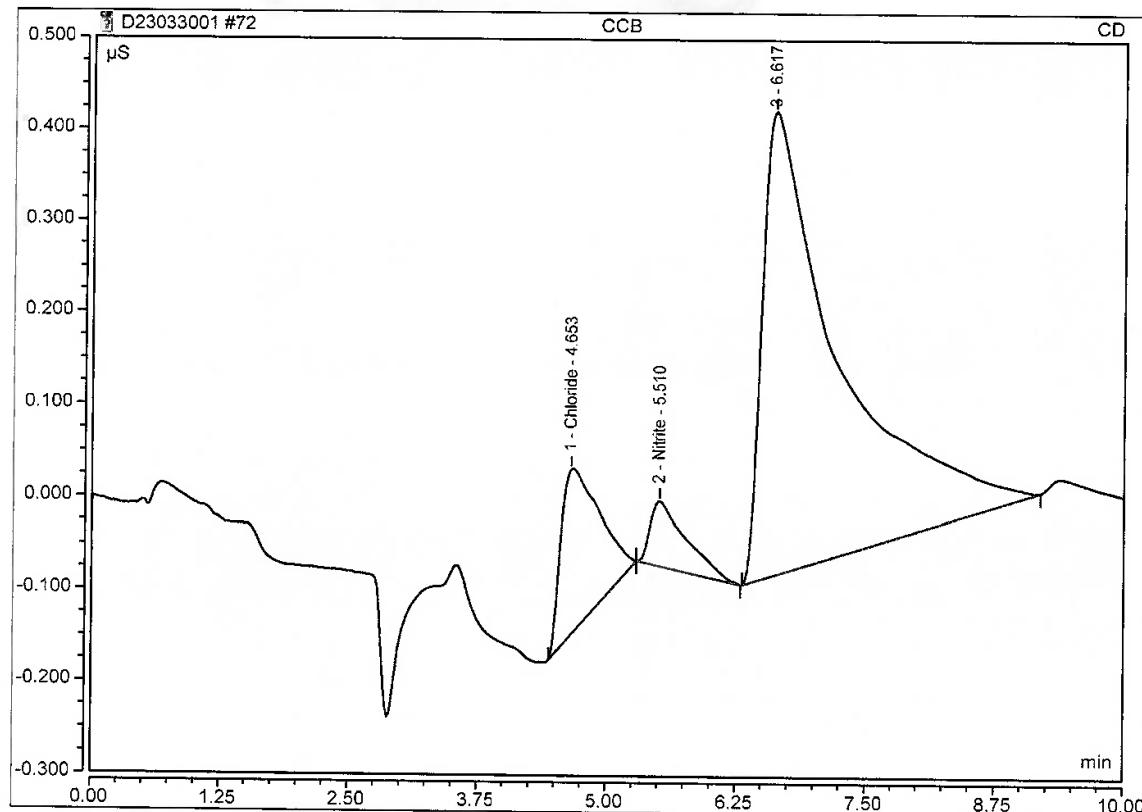
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 18:19	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	4.65	Chloride	BMb	0.076	0.180	0.4784
2	5.51	Nitrite	bMB	0.029	0.070	0.1063
<b>TOTAL:</b>				0.10	0.25	0.58



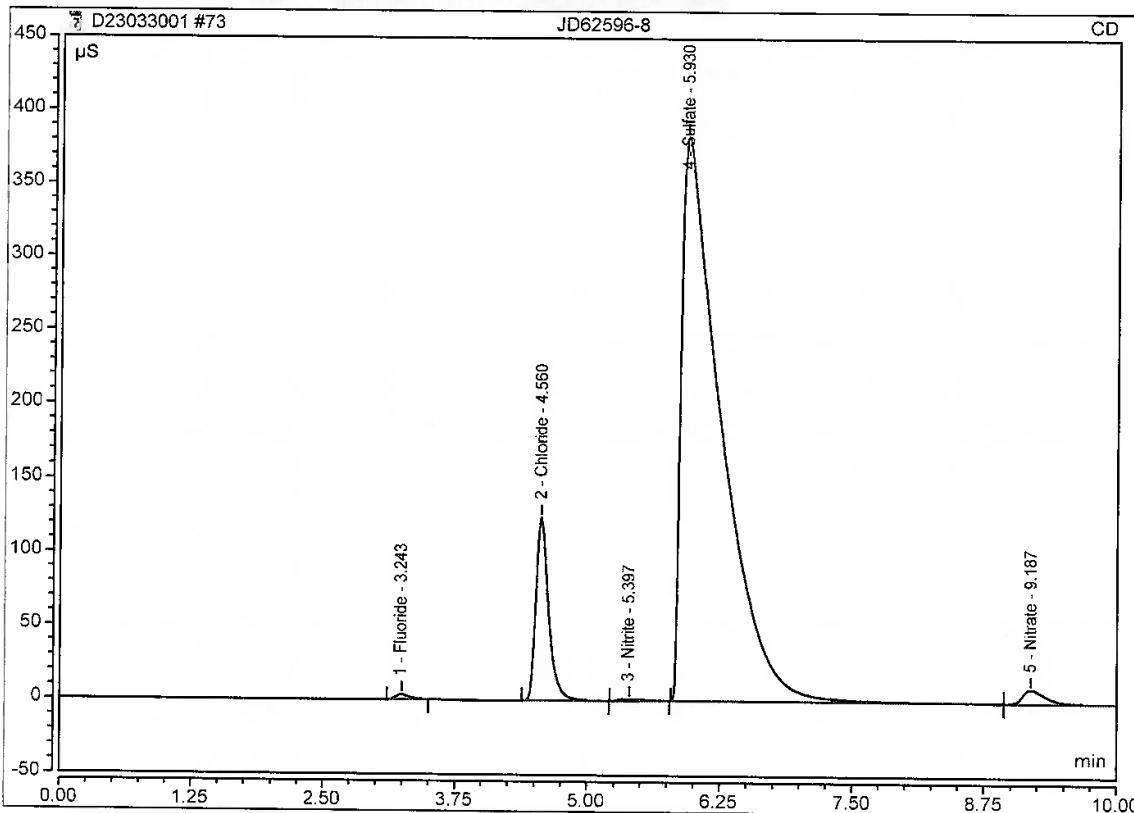
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 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-8	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 18:32	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.24	Fluoride	BMB	0.499	3.456	1.2437
2	4.56	Chloride	BMB	17.365	124.458	54.9083
3	5.40	Nitrite	BMB	0.345	1.185	2.2686
4	5.93	Sulfate	BMb	165.953	382.906	732.4135
5	9.19	Nitrate	bMB	2.313	9.267	19.4775
<b>TOTAL:</b>				<b>186.48</b>	<b>521.27</b>	<b>810.31</b>



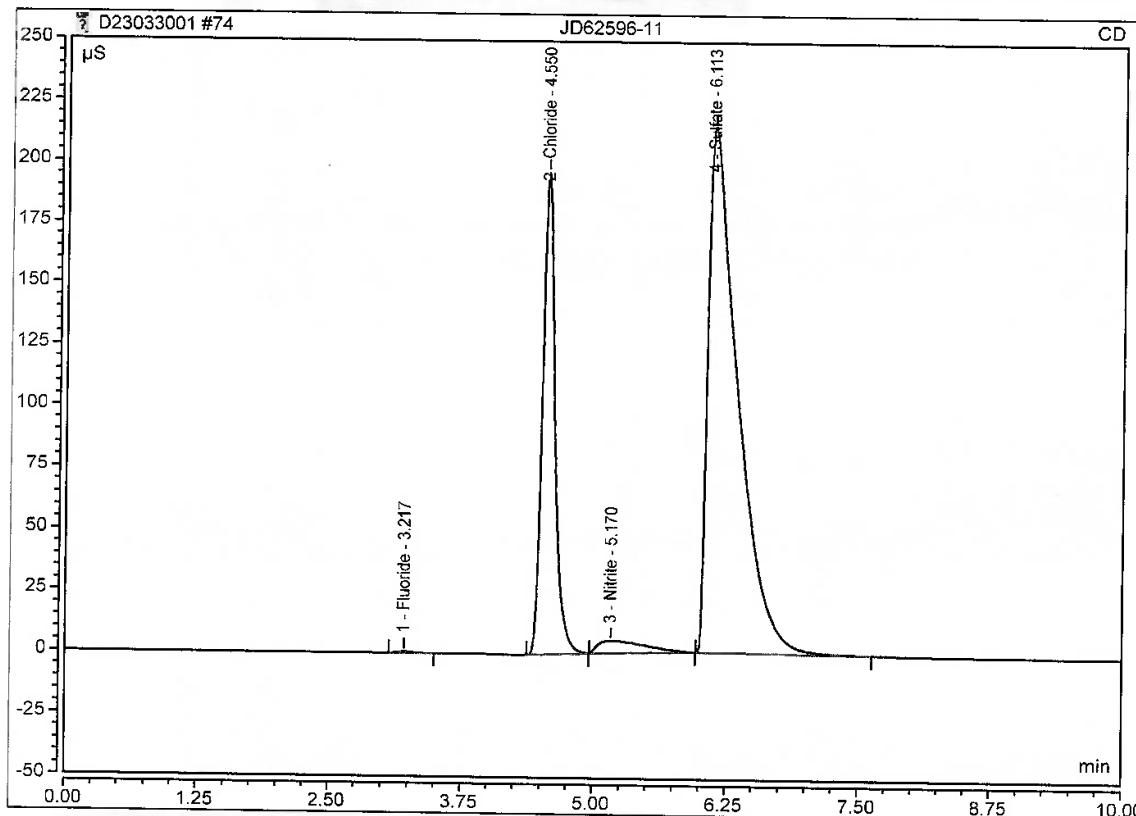
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-11	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 18:45	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.095	0.571	0.2678
2	4.55	Chloride	BMB	27.095	196.337	85.5372
3	5.17	Nitrite	BMB	2.605	5.030	17.6982
4	6.11	Sulfate	BMB	69.556	214.302	307.0401
<b>TOTAL:</b>				99.35	416.24	410.54



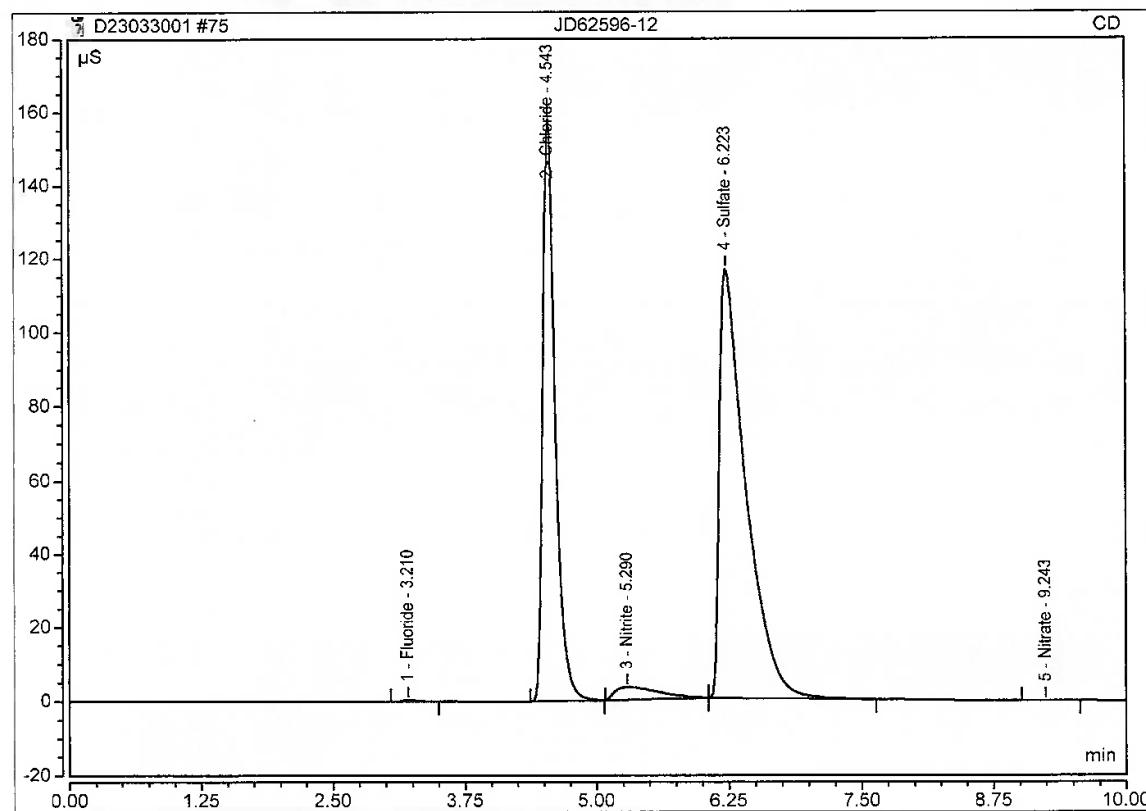
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-12	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 18:58	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.062	0.390	0.1870
2	4.54	Chloride	BMB	21.726	158.311	68.6357
3	5.29	Nitrite	BMB	1.644	3.454	11.1354
4	6.22	Sulfate	BMB	33.392	116.347	147.4558
5	9.24	Nitrate	BMB	0.008	0.036	0.0003
<b>TOTAL:</b>				56.83	278.54	227.41



Anion/Integration

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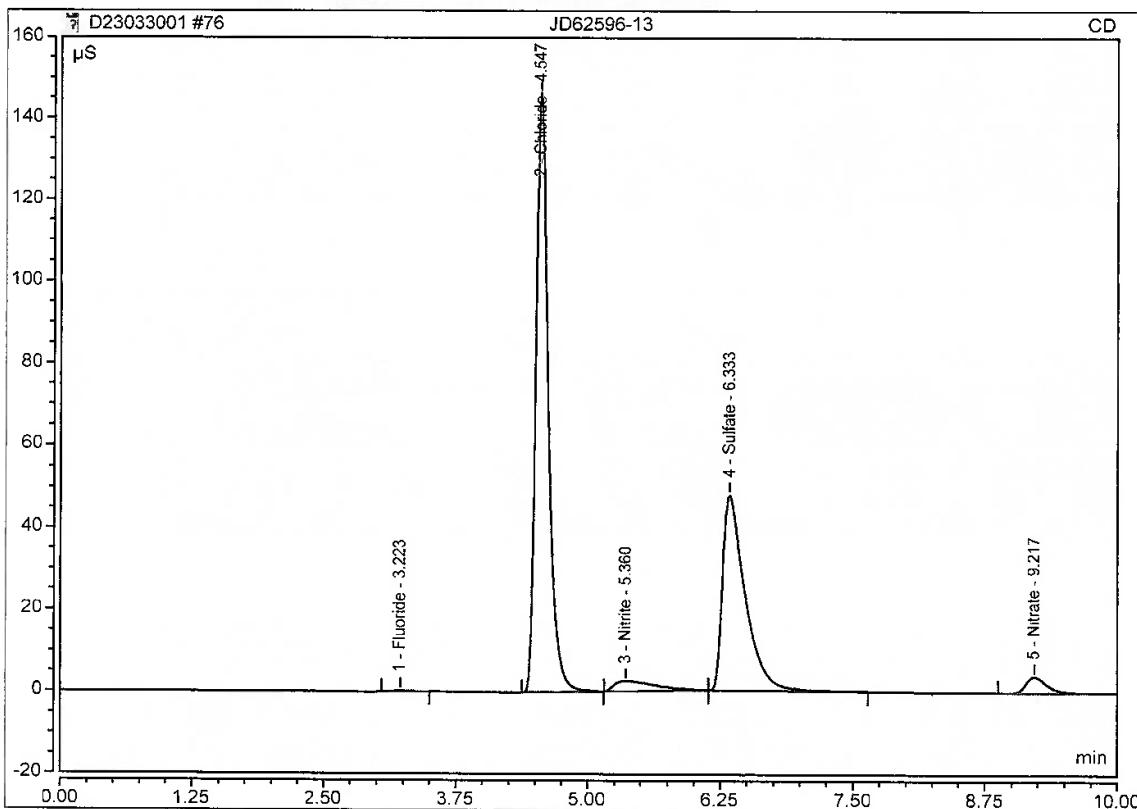
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 Instrument: Integrion\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62596-13	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 19:11	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.044	0.285	0.1442
2	4.55	Chloride	BMB	20.027	146.587	63.2887
3	5.36	Nitrite	BMb	1.106	2.469	7.4617
4	6.33	Sulfate	bMB	11.848	47.911	52.3883
5	9.22	Nitrate	BMB	0.921	3.978	7.7187
<b>TOTAL:</b>				33.95	201.23	131.00



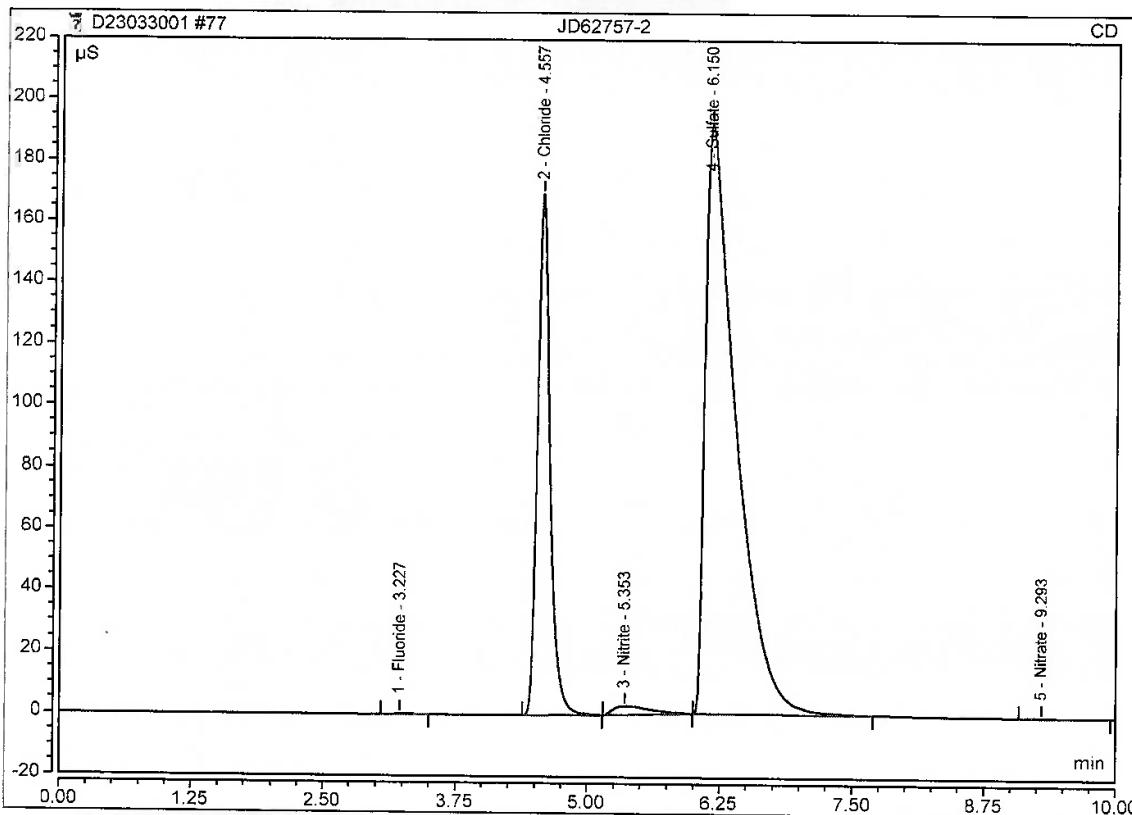
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62757-2	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 19:24	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.23	Fluoride	BMB	0.037	0.188	0.1262
2	4.56	Chloride	BMB	23.451	169.639	74.0674
3	5.35	Nitrite	BMB	1.103	2.643	7.4427
4	6.15	Sulfate	BMB	64.944	191.971	286.6868
5	9.29	Nitrate	BMB	0.028	0.115	0.1684
<b>TOTAL:</b>				89.56	364.56	368.49



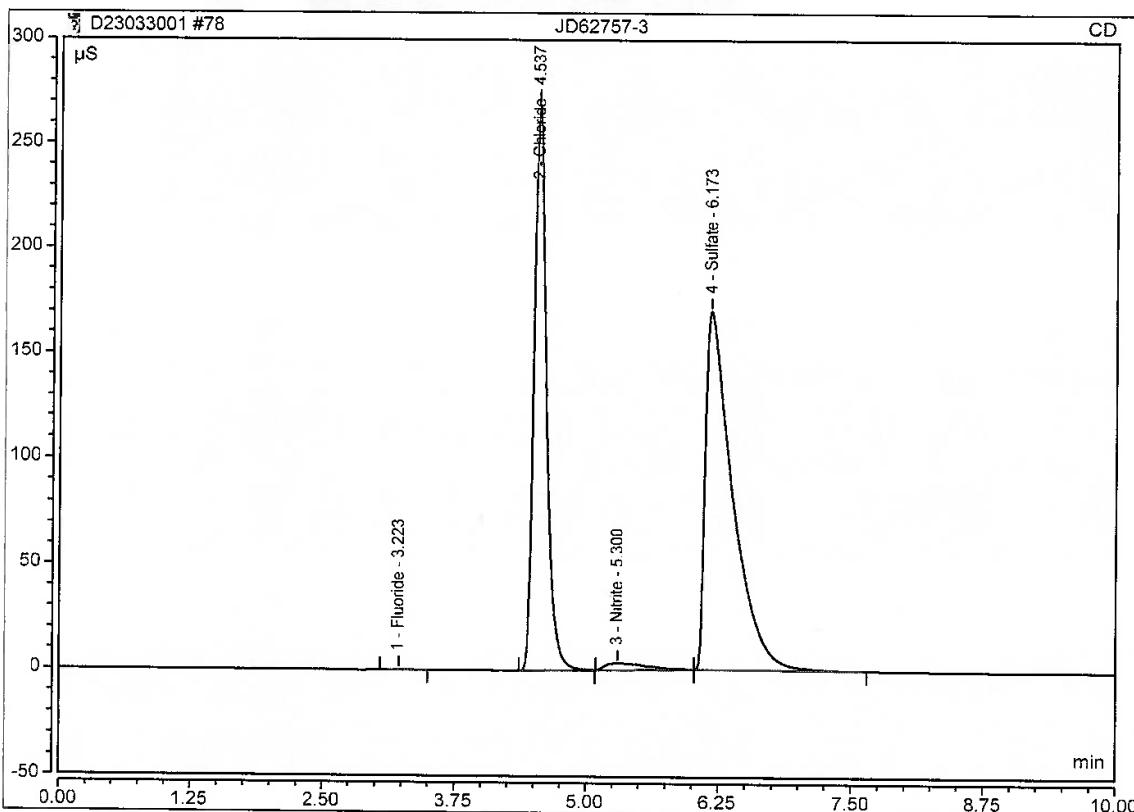
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62757-3	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 19:37	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.049	0.237	0.1553
2	4.54	Chloride	BMB	36.853	268.798	116.2583
3	5.30	Nitrite	BMB	1.531	3.336	10.3616
4	6.17	Sulfate	BMB	50.604	169.927	223.4089
TOTAL:				89.04	442.30	350.18



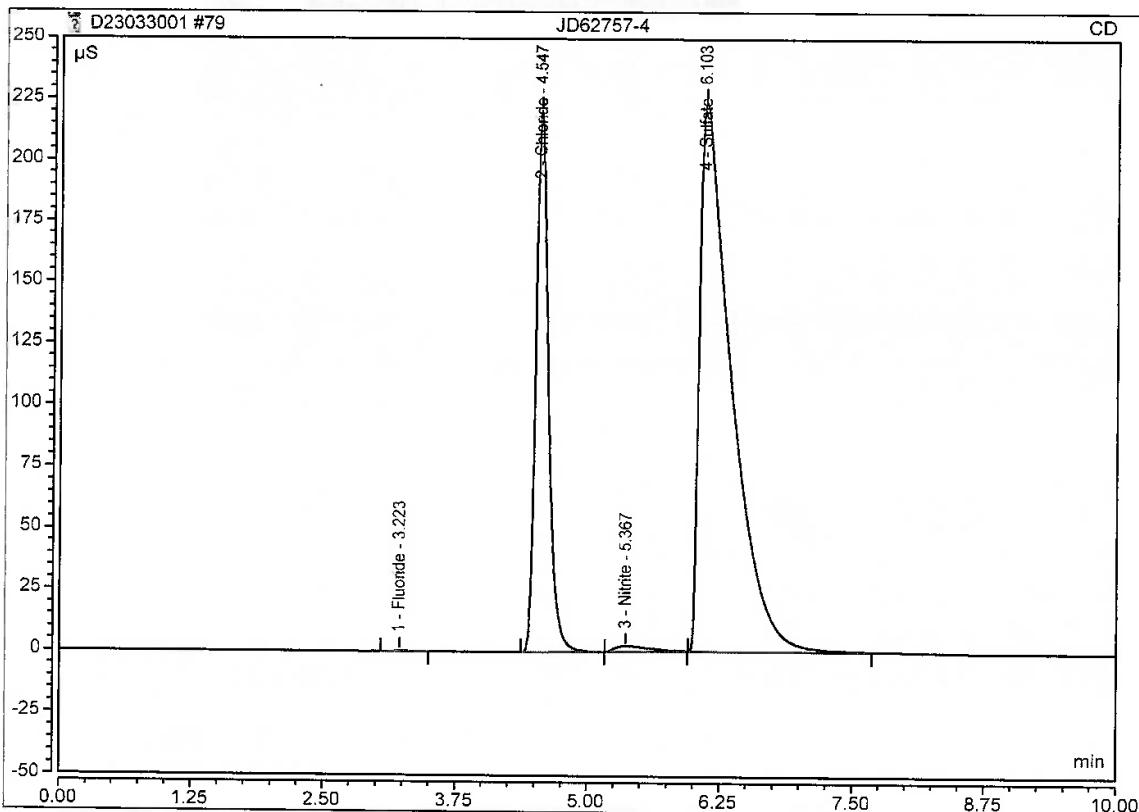
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 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62757-4	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 19:50	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.035	0.201	0.1227
2	4.55	Chloride	BMB	30.269	220.959	95.5307
3	5.37	Nitrite	BMB	0.883	2.284	5.9420
4	6.10	Sulfate	BMB	77.423	223.814	341.7546
TOTAL:				108.61	447.26	443.35



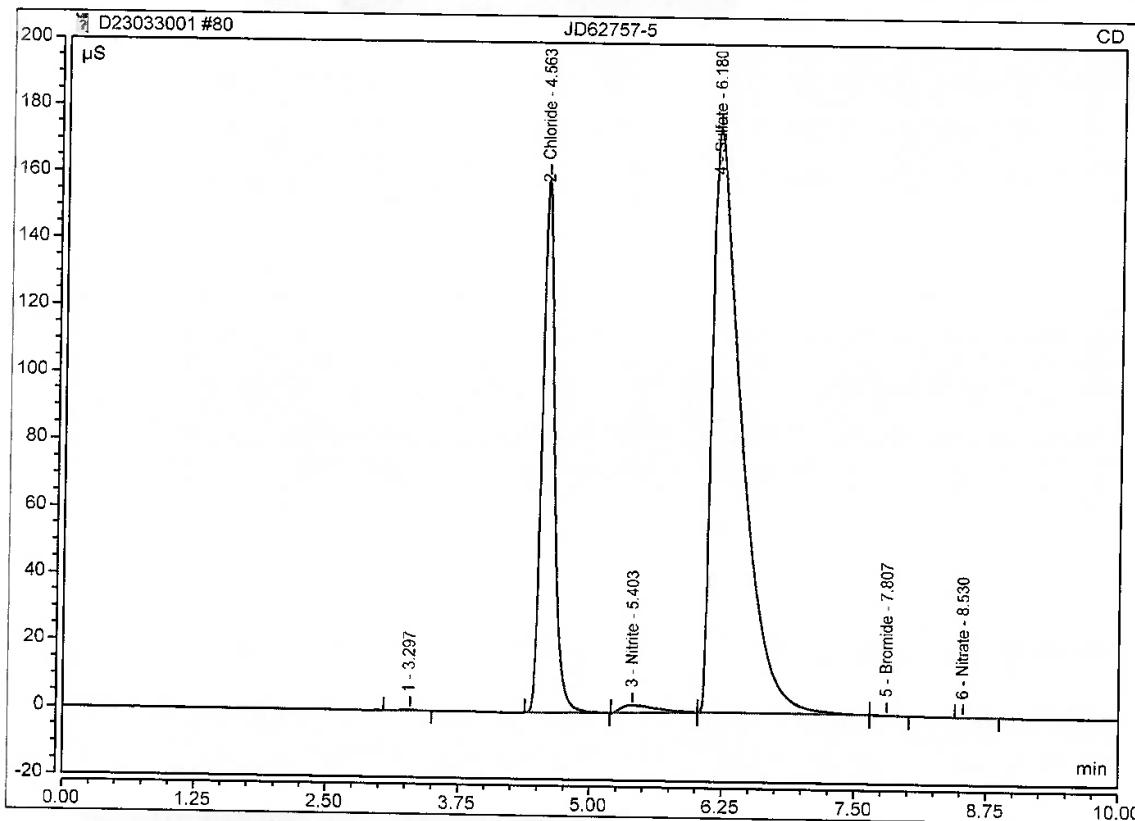
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	JD62757-5	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 20:04	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.56	Chloride	BMB	21.592	159.601	68.2135
3	5.40	Nitrite	BMB	0.835	2.120	5.6139
4	6.18	Sulfate	BMb	55.482	174.553	244.9341
5	7.81	Bromide	bMB	0.011	0.065	0.1702
6	8.53	Nitrate	BMB	0.005	0.031	n.a.
<b>TOTAL:</b>				77.92	336.37	318.93



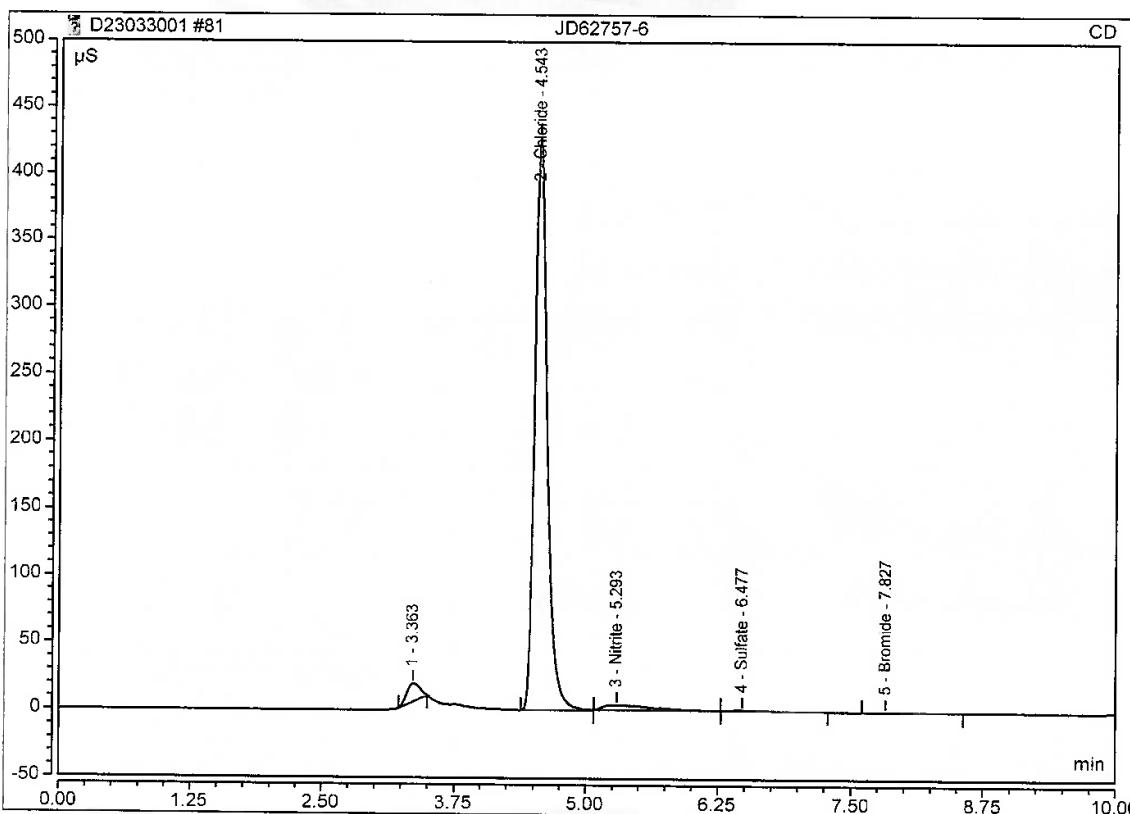
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	JD62757-6	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 20:17	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.54	Chloride	BMB	58.914	428.061	185.7088
3	5.29	Nitrite	BMb	2.038	3.833	13.8294
4	6.48	Sulfate	bMB	0.151	0.567	0.7726
5	7.83	Bromide	BMB	0.022	0.067	0.2675
TOTAL:				61.13	432.53	200.58



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

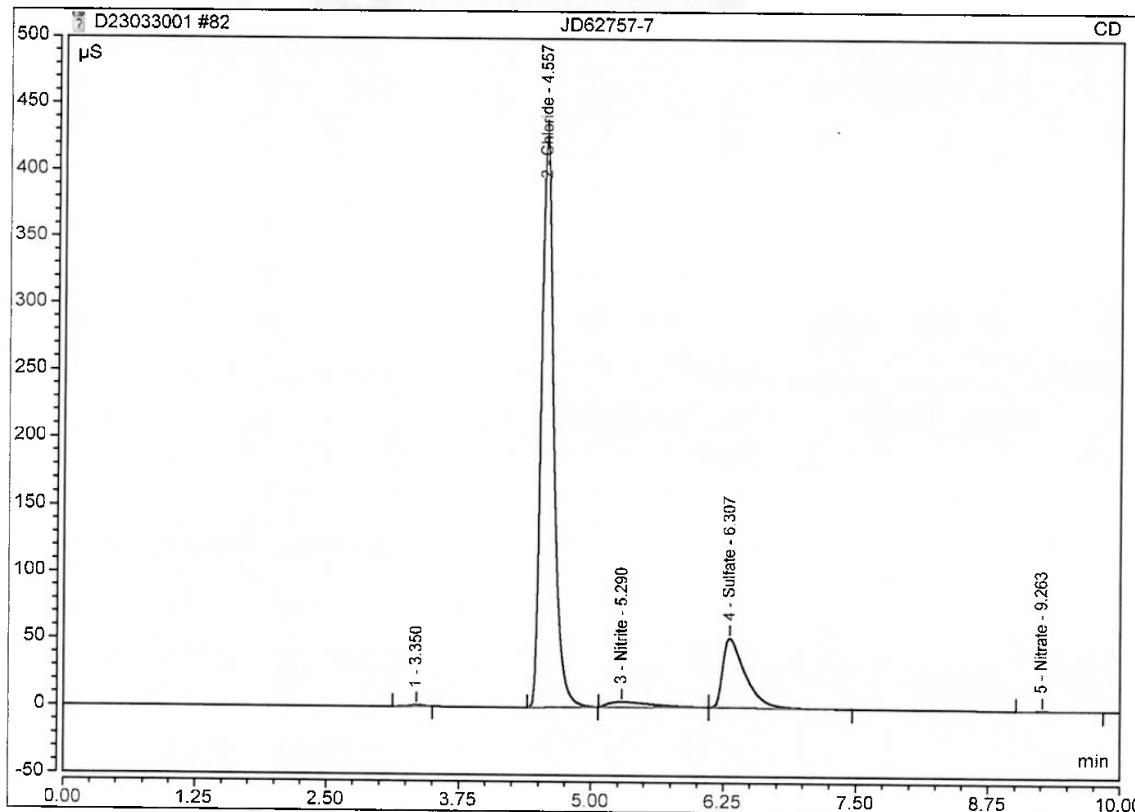
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	JD62757-7	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 20:30	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^{\star}\text{min}$	Height $\mu\text{S}$	Amount
2	4.56	Chloride	BMB	58.925	429.318	185.7425
3	5.29	Nitrite	BMb	1.917	3.953	12.9993
4	6.31	Sulfate	bMB	12.673	51.749	56.0277
5	9.26	Nitrate	BMB	0.021	0.087	0.1092
<b>TOTAL:</b>				73.53	485.11	254.88



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

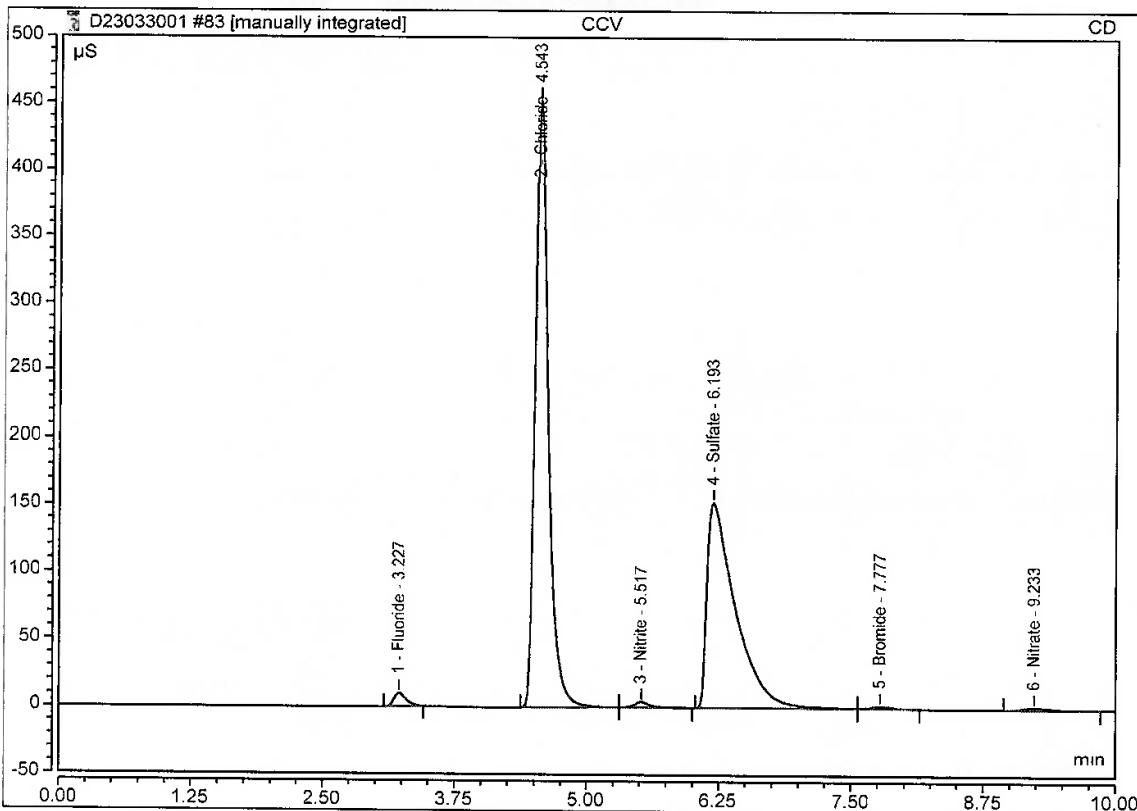
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 Instrument: Integron\_1  
 Sequence: D23033001

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 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	CCV	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 20:43	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.23	Fluoride	BMB*	1.342	10.067	3.2836
2	4.54	Chloride	BMb	65.022	451.957	204.9382
3	5.52	Nitrite	bMB	0.656	4.079	4.3931
4	6.19	Sulfate	BMb	47.572	153.039	210.0313
5	7.78	Bromide	bMB	0.339	1.824	2.9821
6	9.23	Nitrate	BMB	0.449	1.862	3.7322
<b>TOTAL:</b>				115.38	622.83	429.36



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

9.4  
9

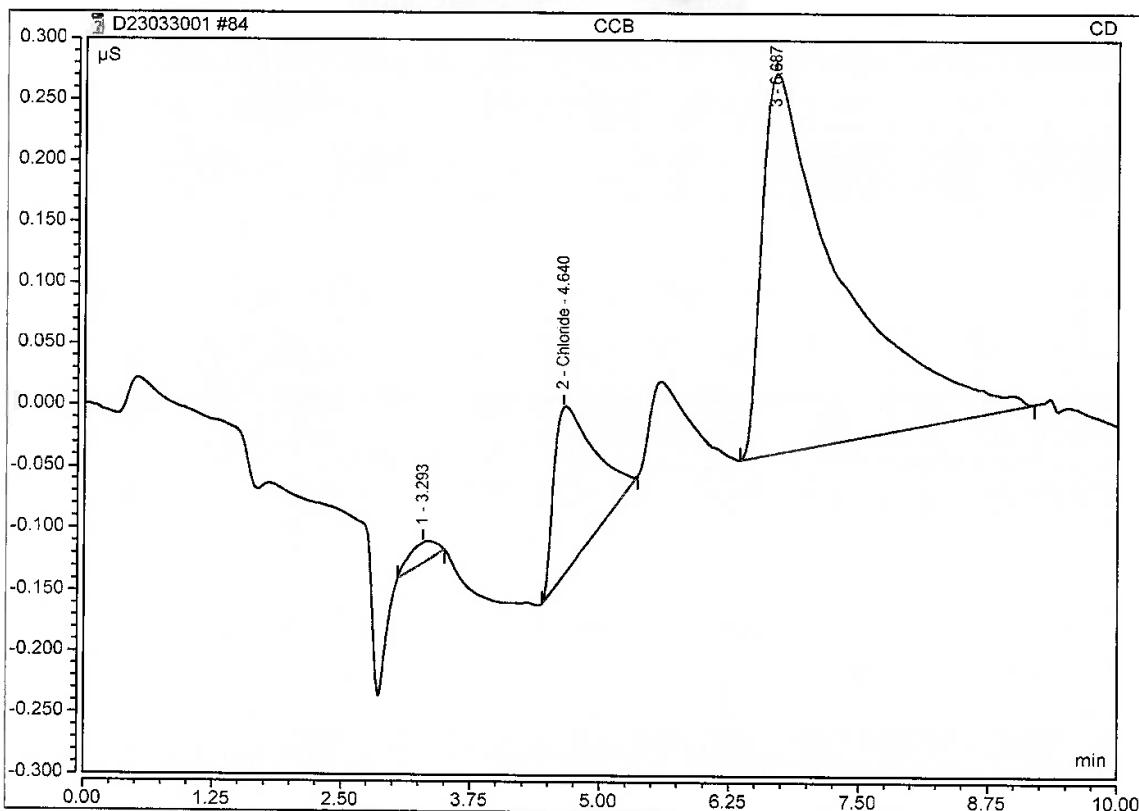
Logged on User: Chemistry  
Instrument: Integriion\_1  
Sequence: D23033001

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### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 20:56	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.64	Chloride	BMB	0.059	0.136	0.4250
		TOTAL:		0.06	0.14	0.42



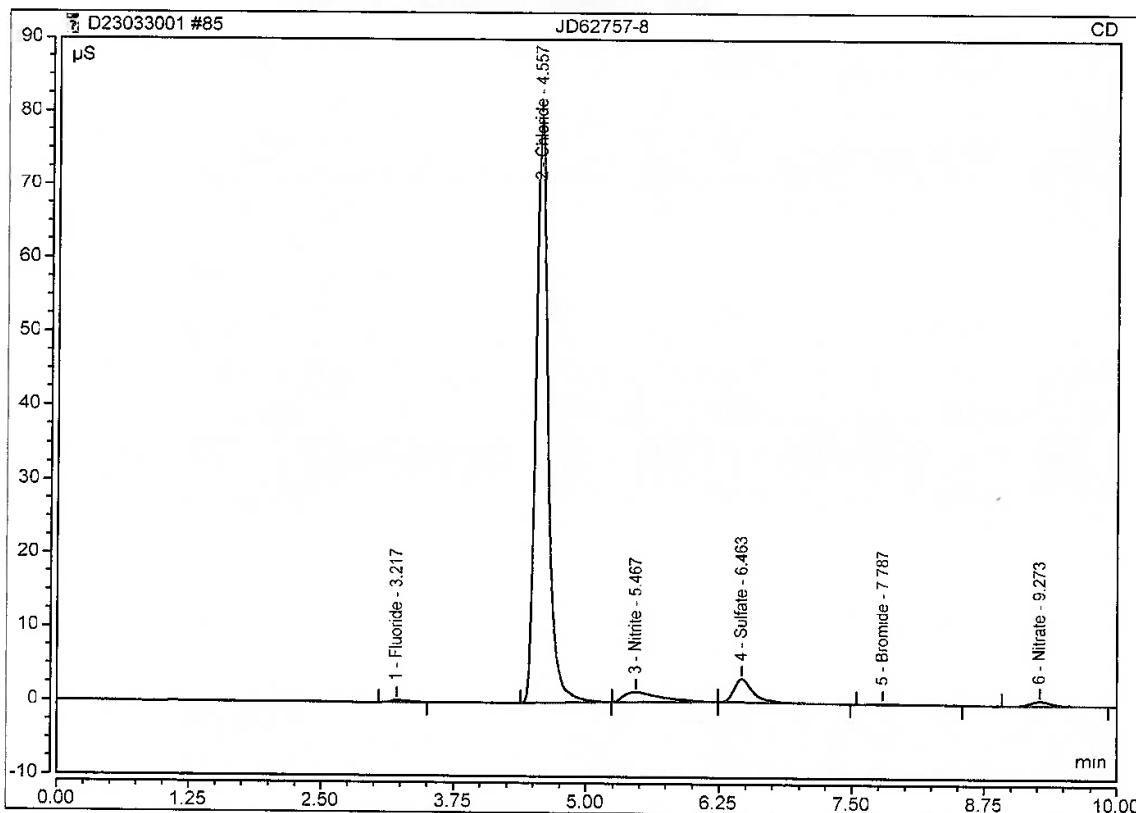
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	JD62757-8	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 21:09	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.048	0.317	0.1535
2	4.56	Chloride	BMB	10.689	79.760	33.8905
3	5.47	Nitrite	BMb	0.547	1.335	3.6445
4	6.46	Sulfate	bMB	0.672	3.124	3.0749
5	7.79	Bromide	BMB	0.030	0.098	0.3321
6	9.27	Nitrate	BMB	0.150	0.644	1.2012
<b>TOTAL:</b>				12.14	85.28	42.30



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

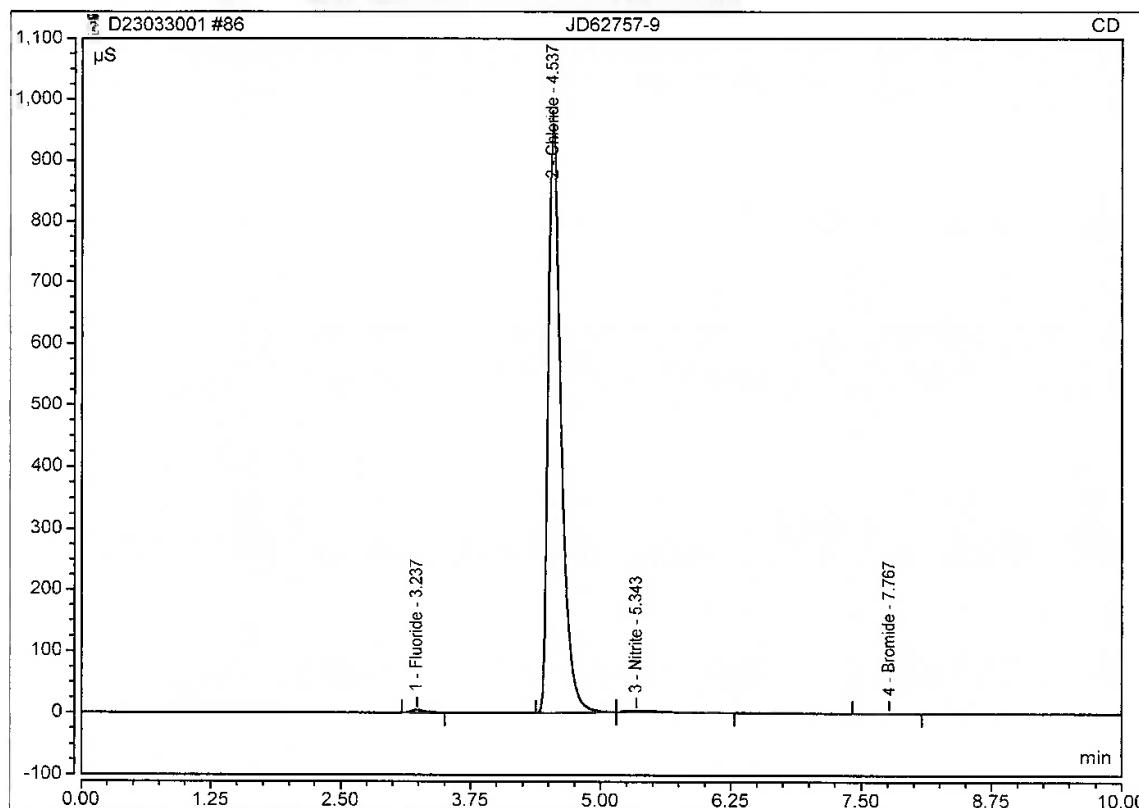
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62757-9	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 21:22	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.24	Fluoride	BMB	0.649	4.556	1.6063
2	4.54	Chloride	BMB	138.281	960.587	435.5677
3	5.34	Nitrite	BMB	1.141	2.608	7.7002
4	7.77	Bromide	BMB	0.041	0.205	0.4283
<b>TOTAL:</b>				140.11	967.96	445.30



Anion/Integration

Chromleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

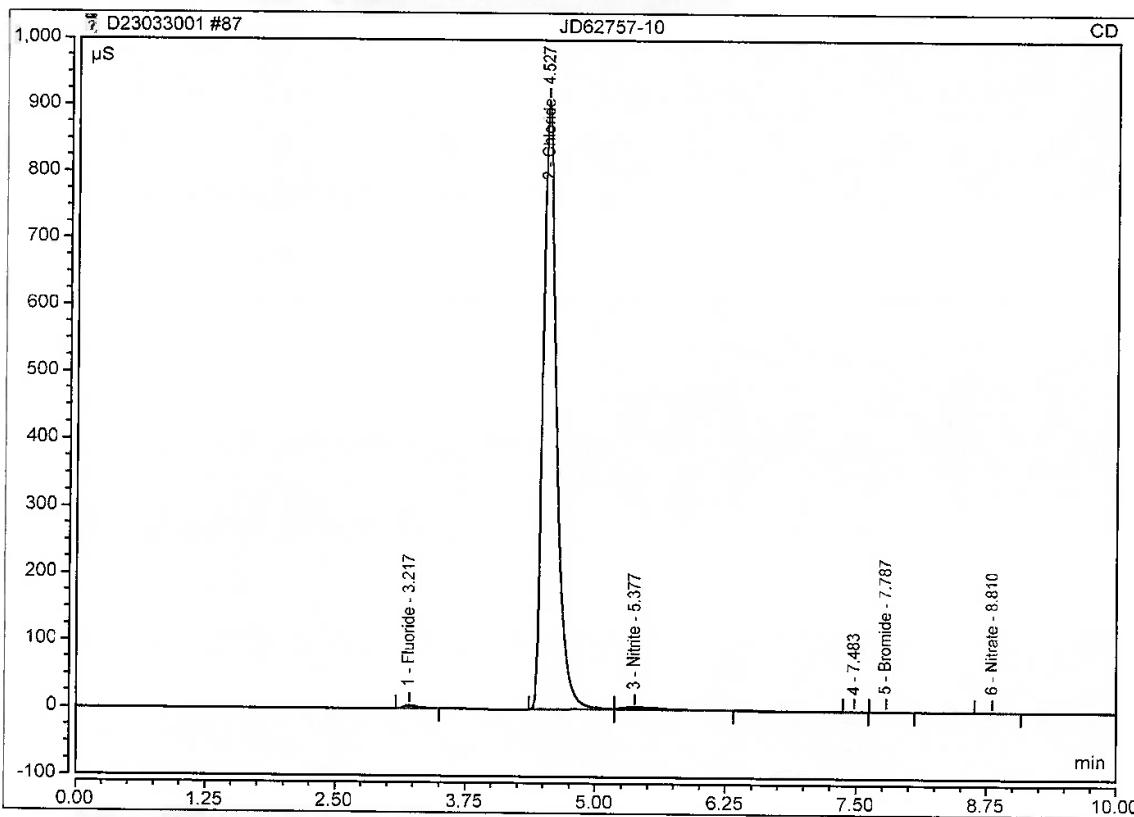
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 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62757-10	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 21:35	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}^*\text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB	0.627	4.414	1.5541
2	4.53	Chloride	BMb	135.465	908.501	426.7031
3	5.38	Nitrite	bMB	1.068	2.392	7.2040
5	7.79	Bromide	BMB	0.017	0.095	0.2233
6	8.81	Nitrate	BMB	0.005	0.026	n.a.
<b>TOTAL:</b>				137.18	915.43	435.68



Anion/Integration

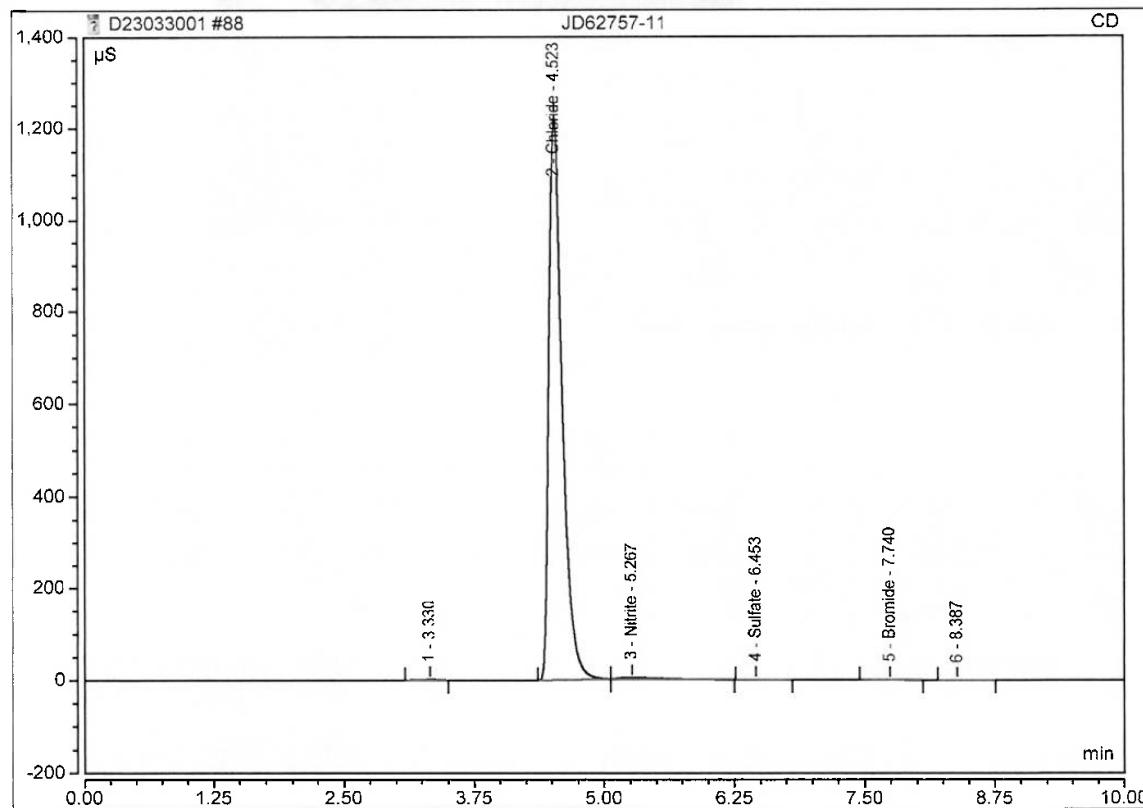
Chromeleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

Logged on User: Chemistry  
Instrument: Integron\_1  
Sequence: D23033001

## Peak Integration Report

Sample Name:	JD62757-11	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 21:48	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}'\text{min}$	Height $\mu\text{S}$	Amount
2	4.52	Chloride	BMB	179.115	1227.770	564.1182
3	5.27	Nitrite	BMB	1.716	3.467	11.6261
4	6.45	Sulfate	BMB	0.087	0.450	0.4901
5	7.74	Bromide	BMB	0.028	0.115	0.3197
<b>TOTAL:</b>				180.95	1231.80	576.55



Anion/Integration

Chromleon (c) Dionex 1996-2009  
Version 7.2.8.10783

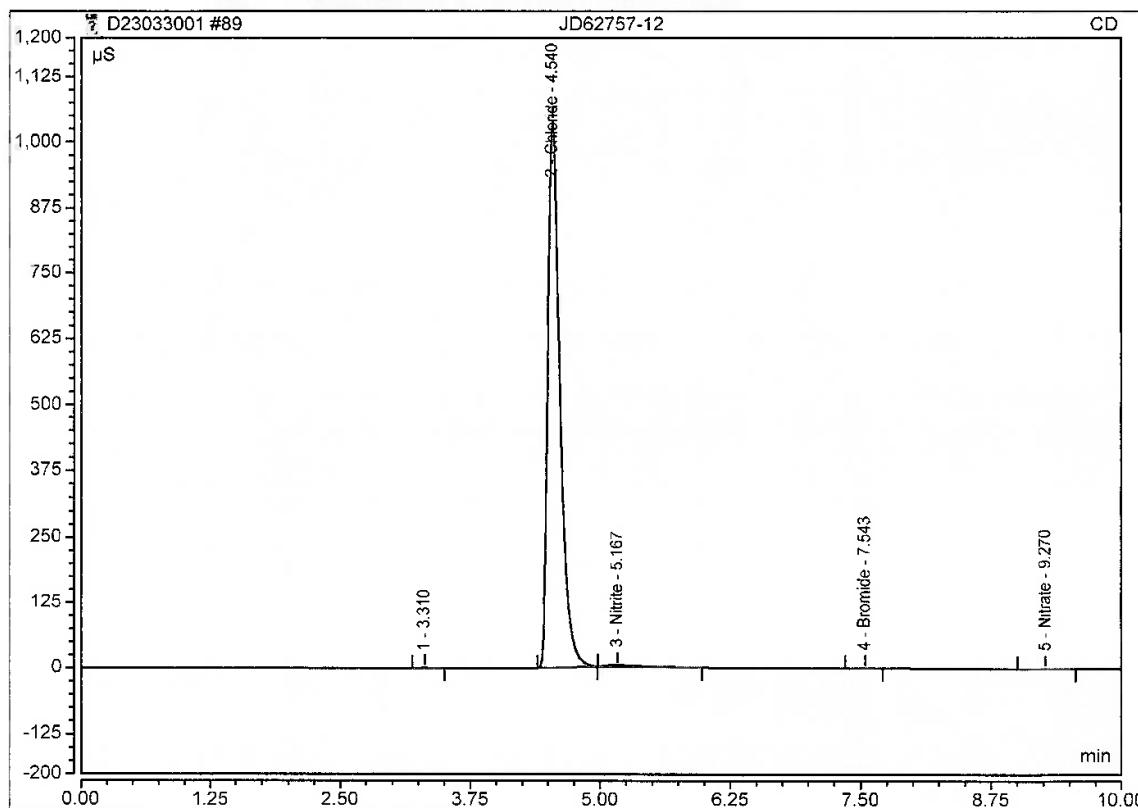
Logged on User: Chemistry  
 Instrument: Integron\_1  
 Sequence: D23033001

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### Peak Integration Report

Sample Name:	JD62757-12	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 22:01	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.54	Chloride	BMB	141.315	1040.070	445.1193
3	5.17	Nitrite	BMB	1.328	2.830	8.9787
4	7.54	Bromide	BMB	0.010	0.057	0.1675
5	9.27	Nitrate	BMB	0.004	0.021	n.a.
<b>TOTAL:</b>				142.66	1042.98	454.27



Anion/Integration

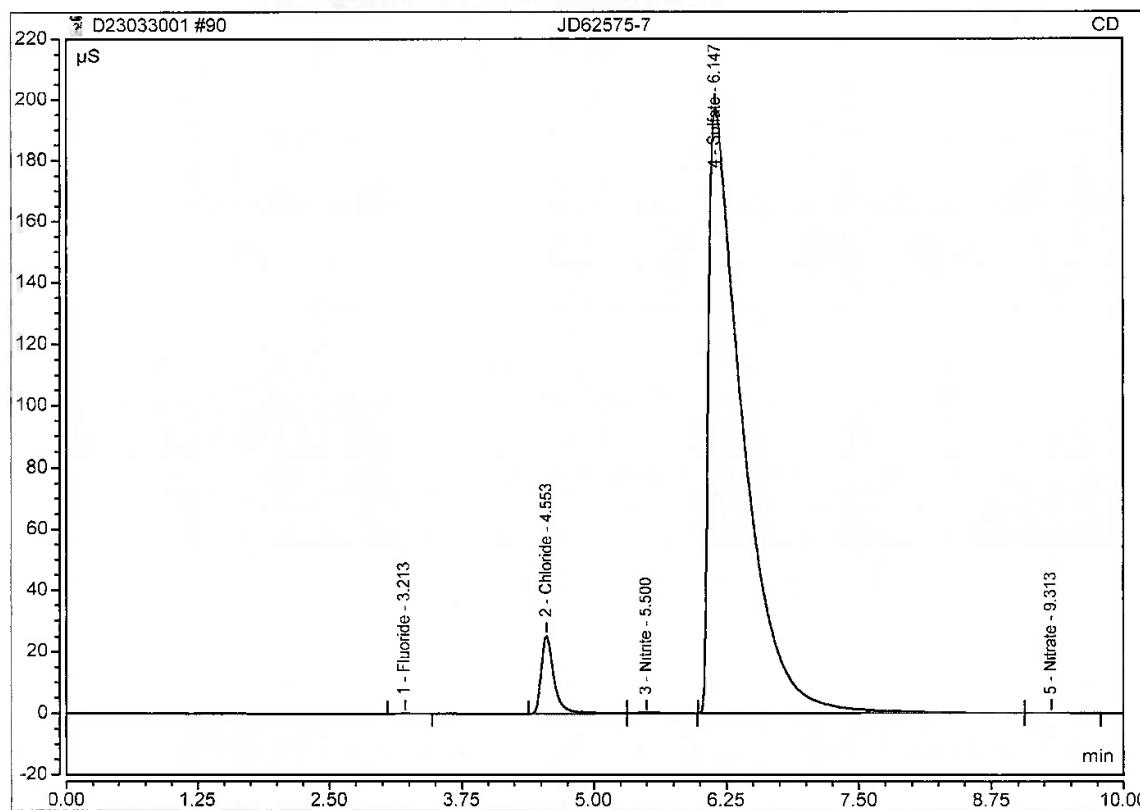
Chromeleon (c) Dionex 1996-2009  
 Version 7.2.8.10783

Logged on User: Chemistry  
Instrument: Integron\_1  
Sequence: D23033001

## Peak Integration Report

Sample Name:	JD62575-7	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	32.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 22:14	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
1	3.21	Fluoride	BMB	0.041	0.221	4.3973
2	4.55	Chloride	BMb	3.545	25.476	364.7702
3	5.50	Nitrite	bMB	0.047	0.158	7.3961
4	6.15	Sulfate	BMb	72.823	197.534	10286.6382
5	9.31	Nitrate	bMB	0.022	0.092	3.9766
TOTAL:				76.48	223.48	10667.18



Anion/Integration

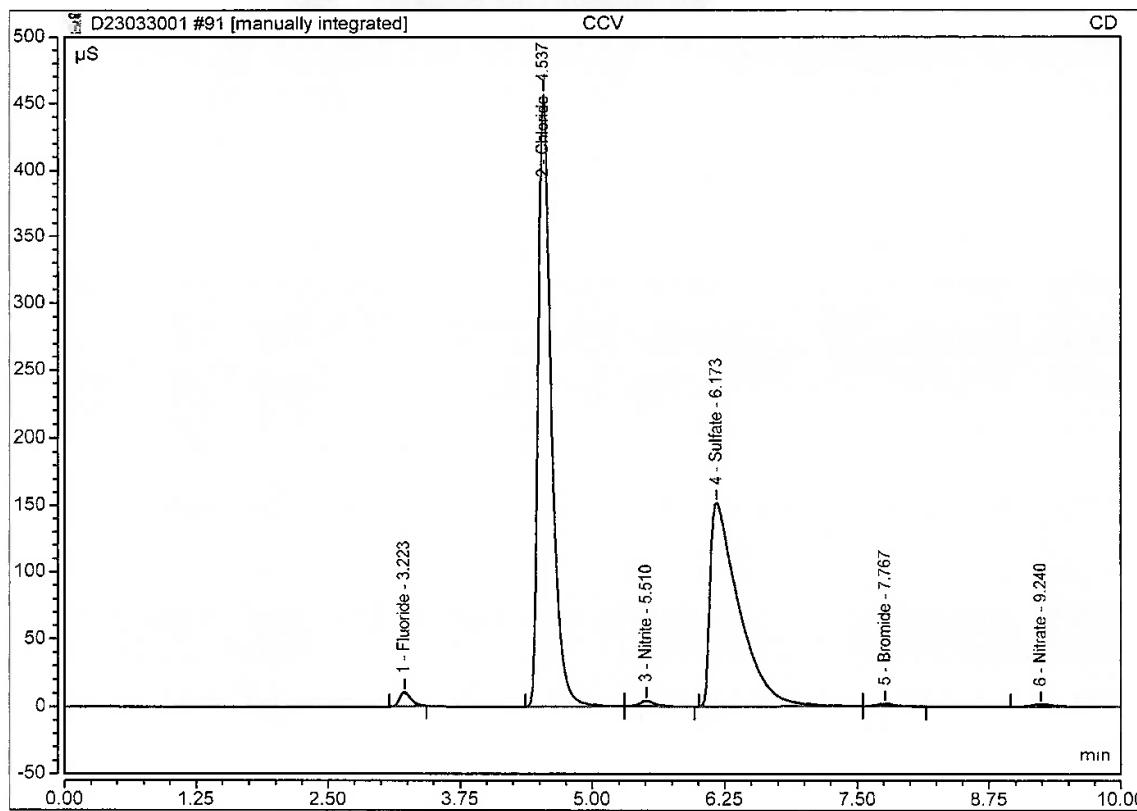
Chromleon (c) Dionex 1996-2009  
Version 7.2.8.10783

Logged on User: Chemistry  
Instrument: Integron\_1  
Sequence: D23033001

## Peak Integration Report

Sample Name:	CCV	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 22:27	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S}/\text{min}$	Height $\mu\text{S}$	Amount
1	3.22	Fluoride	BMB*	1.328	10.529	3.2483
2	4.54	Chloride	BMb	64.965	456.859	204.7582
3	5.51	Nitrite	bMB	0.643	4.136	4.3028
4	6.17	Sulfate	BMb	47.562	152.199	209.9867
5	7.77	Bromide	bMB	0.347	1.856	3.0457
6	9.24	Nitrate	BMB	0.460	1.894	3.8245
TOTAL:				115.30	627.47	429.17



Anion/Integration

Chromleon (c) Dionex 1996-2009  
Version 7.2.8.10783

9.4

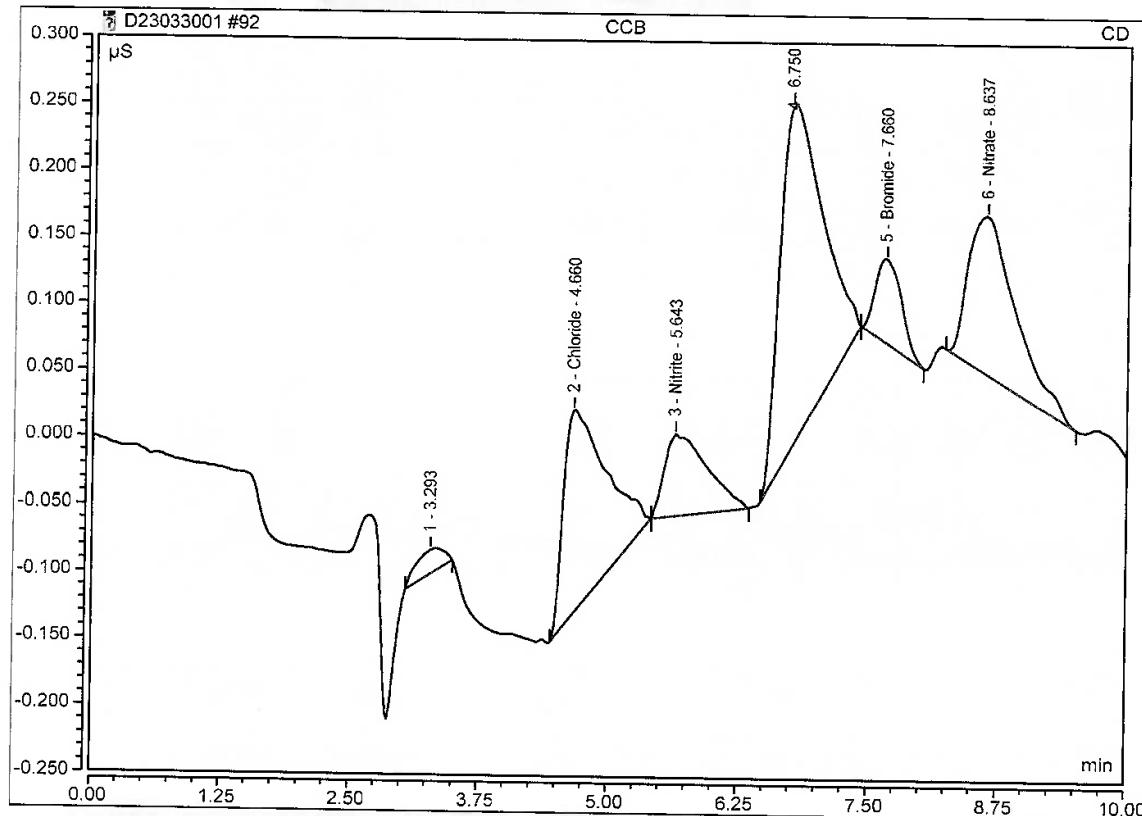
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 Sequence: D23033001

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 4/3/2023 10:39 AM

### Peak Integration Report

Sample Name:	CCB	Inj. Vol.:	5000.00
Injection Type:	Unknown	Dilution Factor:	1.0000
Instrument Method:	Anions_012919	Operator:	Chemistry
Inj. Date / Time:	31-Mar-2023 / 22:40	Run Time:	10.00

No.	Time min	Peak Name	Peak Type	Area $\mu\text{S} \cdot \text{min}$	Height $\mu\text{S}$	Amount
2	4.66	Chloride	BMB	0.070	0.152	0.4589
3	5.64	Nitrite	bMB	0.029	0.060	0.1061
5	7.66	Bromide	bMB	0.018	0.063	0.2367
6	8.64	Nitrate	BMB	0.063	0.118	0.4679
TOTAL:				0.18	0.39	1.27



Test: Iron, Ferrous  
 Product: FE2  
 Method: SM3500FE-B-11

Analyst: MARCINK  
 GN Batch ID: GN39908  
 GP Batch ID: N/A  
 Date: 4/2/2023  
 Instrument ID: M

Original Calibration Information		Calibration Date: 1/19/2023							
		Blank	Std 1	Std 2	Std 3	Std 4	Std 5	Std 6	Std 7
Known:	0.000	0.100	0.200	0.500	1.000	2.000	4.000	6.000	
Absorbance:	0.000	0.012	0.024	0.056	0.120	0.233	0.474	0.711	
Actual Value:	0.003	0.104	0.206	0.493	1.016	1.970	4.004	6.005	
% RE:	#D/Vol	-4.26	-2.78	1.49	-1.59	-1.51	-0.10	-0.08	
Continuing Calibration Check Standards Data:		Correlation Coeff. = 0.99998							
Known:	0.000	0.100	0.600	Std 6					
Absorbance:	0.000	0.011	0.097						
Recovery:	0.0%	95.8%	98.1%						
Actual Value:	0.003	0.096	0.587						
Bottle #	Sample ID	Time Analyzed	Initial Wt (g) or Vol (ml)	Final Vol (ml)	Dilution	Sample Abs.	Background Abs.	Result From Curve (mg/L)	RL (Normal RL = 0.200 mg/l)
ICV	13:15	50	50	1	0.113	NA	0.957	0.957	NA
CCV	13:49	50	50	1	0.110	NA	0.932	0.932	NA
GN39908-MB1	13:49	50	50	1	0.000	0.000	0.003	0.003	0.200
GN39908-B1	13:49	50	50	1	0.342	0.000	2.890	2.890	0.200
GN39908-S1	13:49	50	50	1	0.349	0.000	2.949	2.949	0.200
GN39908-MSD1	13:49	50	50	1	0.348	0.000	2.941	2.941	0.200
1 JD62888-1	13:49	50	50	1	0.002	0.000	0.020	0.020	0.200
1 JD62888-2	13:49	50	50	1	0.002	0.000	0.020	0.020	0.200
1 JD62888-3	13:49	50	50	1	0.001	0.000	0.011	0.011	0.200
1 JD62888-4	13:49	50	50	1	0.000	0.000	0.003	0.003	0.200
CCVA	13:49	50	50	1	0.699	NA	5.903	5.903	NA

Analyst: \_\_\_\_\_

Date: \_\_\_\_\_ QC Reviewer: \_\_\_\_\_ Date: \_\_\_\_\_

6 56



Test: IRON, FERROUS

Product: FE2

Method: SM18 3500FE B-2011 (aqueous) Units: (mg/l)

Analyst: MK  
 GNBatch ID: GN39908  
 GPBatch ID:  
 Date: 9/2/23

**Preparation Batch QC Summary**Units = mg/lMethod Blank ID: GN39908 4101 Date: 4/2/2023 Result: <OL DL: \_\_\_\_\_ <DL: \_\_\_\_\_Spike Blank ID: -81 Date: ↓ Result: 2.810 Spike: 3.00 %Rec.: 96.33%MS Blank ID: -51 Date: ↓ Results: 2.149 Samp. Result: 0.020 Spike: 3.0 %Rec.: 96.33%MS Duplicate ID: -MS01 Date: ↓ Results: 2.941 Samp. Result: 0.020 Spike: 3.0 %Rec.: 97.37%

Duplicate ID: \_\_\_\_\_ Date: \_\_\_\_\_ Dup. Result: \_\_\_\_\_ Samp. Result: \_\_\_\_\_ %RPD: \_\_\_\_\_

Matrix spike Duplicate Result: \_\_\_\_\_ Matrix spike Result: \_\_\_\_\_ %RPD: \_\_\_\_\_

**Analysis Batch QC Summary**Units = \_\_\_\_\_  
ICV (Ext): 4/2/23 Result: 0.957 TV: 1.0 %Rec: 95.77%CCV: \_\_\_\_\_ Result: 0.932 TV: 1.00 %Rec.: 93.2%CCVA: \_\_\_\_\_ Result: 5.703 TV: 6.00 %Rec.: 78.47%CCV: \_\_\_\_\_ Result: 1.000 TV: 1.00 %Rec.: 100.0%CCVA: \_\_\_\_\_ Result: 6.000 TV: 6.00 %Rec.: 100.0%CCV: \_\_\_\_\_ Result: 1.000 TV: 1.00 %Rec.: 100.0%CCVA: \_\_\_\_\_ Result: 6.000 TV: 6.00 %Rec.: 100.0%

CCB: \_\_\_\_\_ Result: \_\_\_\_\_ DL: \_\_\_\_\_ &lt;DL: \_\_\_\_\_

**Reagent Reference Numbers:**

BSP/MS---1.5 ML OF 200PPM STOCK--100ML SAMPLE AND DI (WATER)


Analyst: \_\_\_\_\_ Date: \_\_\_\_\_

Comments: \_\_\_\_\_

Form: GN032-01

Rev. Date: 9/4/09]



## Reagent Information Log

**Test Name: FERROUS IRON (FE2)**

All standards and stocks were made as described in the SOP for this method (circle one):  Y or  N  
If no (N), see attached page for standards prep.

Form: GN087-01  
Rev. Date: 4/2/2023



## GENERAL CHEMISTRY STANDARD PREPARATION LOG

Product: Fe2

GN or GP Number: \_\_\_\_\_

Reagent Standards		Tracking #		Expiration					
200 PPM FE2 STOCK		GNE3-3/22/2023		5/22/2023					
200 PPM FE2 ICV		GNE3-3/22/2023		5/22/2023					
Intermediate Standard Description	Stock used to prepare standard	Stock concentration	volume or weight used with units	Balance or Autopipet ID (*)	Additional Reagents	Diluent	Final Volume	Final Conc. of Intermediate (mg/l)	Expiration Date
10 PPM FE2 STD	GNE3-3/22/2023	200 mg/L	5.00 mL	A		DI water	100mL	10.0	04/02/23
gn-4/2/2023									
10 PPM FE2 ICV	GNE3-3/22/2023	200 mg/L	5.00 mL	A		DI water	100mL	10.0	04/02/23
gn-4/2/2023									
Intermediate Standard Description	Intermediate or Stock used to prepare standard	Intermediate or Stock concentration	Balance or Stock volume used in mL	Autopipet ID (*)	Additional Reagents	Diluent	Final Volume	Final Conc. of Standard (mg/l)	Expiration Date
BLANK	NA	0 mg/L	0.00	A	2.00 mL of concentrated	DI water	100 mL	0.00	04/02/23
0.10 mg/L	gn-4/2/2023	10 mg/L	1.00	A	HCl and 1.00 mL of	DI water	100 mL	0.10	04/02/23
0.20 mg/L	gn-4/2/2023	10 mg/L	2.00	A	hydroxylamine hydrochloride added to each standard.	DI water	100 mL	0.20	
0.50 mg/L	gn-4/2/2023	10 mg/L	5.00	A		DI water	100 mL	0.50	
1.00 mg/L	gn-4/2/2023	10 mg/L	10.00	A		DI water	100 mL	1.00	
2.00 mg/L	GNE3-3/22/2023	200 mg/L	1.00	A		DI water	100 mL	2.00	
4.00 mg/L	GNE3-3/22/2023	200 mg/L	2.00	A		DI water	100 mL	4.00	
6.00 mg/L	GNE3-3/22/2023	200 mg/L	3.00	A		DI water	100 mL	6.00	05/22/23
ICV	gn-4/2/2023	10 mg/L	10.00	A		DI water	100 mL	1.00	04/02/23
CCV	gn-4/2/2023	10 mg/L	10.00	A		DI water	100 mL	1.00	04/02/23
CCVA	GNE3-3/22/2023	200 mg/L	3.00			DI water	100 mL	6.00	05/22/23

\* If Class A glass pipets are used, enter an A. For balances or autopipets, then enter the appropriate Accutest ID number.

Form: GN121-01  
Rev. Date: 1/13/09

LABORATORY REVIEW SIGNATURE FORM  
(To be stored with the raw data)

File ID: 0230331W1.TXT      Date Analyzed: 03/01/23      Methods: SM5310 B-11, SM5310 B-11/14  
Analyst: MB      Run ID: GN39916

The following analyst(s) have reviewed this run and attest that, to the best of their knowledge, this documentation is complete and correct:

Analyst: MB      Date 4-3-23

Analyst: \_\_\_\_\_ Date \_\_\_\_\_

The following supervisor or their designee has reviewed this run and attests that, to the best of their knowledge, this documentation is complete and correct:

Supervisor (or designee):  Date 4/3/23

9.6  
9

C:\TOC-L\Data\d230331w1.toc.tlx

9  
6

	Sample Name	Sample ID	Origin	Manual Dilution	Result	Comment
1	WASHCONF		TOCAQSW846.met	1.000	NPOC:0.5131mg/L	
2	CRI		TOCAQSW846.met	1.000	NPOC:1.097mg/L	
3	HSTD		TOCAQSW846.met	1.000	NPOC:47.05mg/L	
4	ICV		TOCAQSW846.met	1.000	NPOC:19.89mg/L	
5	ICB		TOCAQSW846.met	1.000	NPOC:0.3547mg/L	
6	CCV		TOCAQSW846.met	1.000	NPOC:22.99mg/L	
7	CCB		TOCAQSW846.met	1.000	NPOC:0.3092mg/L	
8	SPARGERCHK		TOCAQSW846.met	1.000	NPOC:0.3068mg/L	
9	GP45856-MB1	[toc]	TOCAQ.met	1.000	NPOC:0.8149mg/L	
10	GP45856-B1		TOCAQ.met	1.000	NPOC:9.343mg/L	
11	JD62800-1		TOCAQ.met	1.000	NPOC:1.196mg/L	
12	GP45856-S1	JD62800-1	TOCAQ.met	1.000	NPOC:9.196mg/L	
13	GP45856-MSD1	JD62800-1	TOCAQ.met	1.000	NPOC:9.487mg/L	
14	JD62800-2		TOCAQ.met	1.000	NPOC:3.895mg/L	
15	JD62800-3		TOCAQ.met	1.000	NPOC:6.549mg/L	
16	JD62800-4		TOCAQ.met	1.000	NPOC:0.6508mg/L	
17	JD62800-5	(2)	TOCAQ.met	1.000	NPOC:189.0mg/L	non-acidified, see rerun
18	CCVA		TOCAQ.met	1.000	NPOC:46.43mg/L	
19	CCB		TOCAQ.met	1.000	NPOC:0.7008mg/L	
20	JD62800-6	(2)	TOCAQ.met	1.000	NPOC:358.1mg/L	non-acidified, see rerun
21	JD62800-7		TOCAQ.met	1.000	NPOC:14.96mg/L	
22	JD62800-8		TOCAQ.met	1.000	NPOC:0.7392mg/L	
23	JD62800-9	field blank	TOCAQ.met	1.000	NPOC:0.4505mg/L	
24	GP45881-MB1	[doc]	TOCAQ.met	1.000	NPOC:0.1368mg/L	
25	GP45881-B1		TOCAQ.met	1.000	NPOC:9.592mg/L	
26	JD62888-2F		TOCAQ.met	1.000	NPOC:0.6218mg/L	
27	GP45881-S1	JD62888-2F	TOCAQ.met	1.000	NPOC:9.807mg/L	
28	GP45881-MSD1	JD62888-2F	TOCAQ.met	1.000	NPOC:9.703mg/L	
29	JD62888-4F	field blank	TOCAQ.met	1.000	NPOC:0.2537mg/L	
30	CCV		TOCAQ.met	1.000	NPOC:22.91mg/L	
31	CCB		TOCAQ.met	1.000	NPOC:0.5875mg/L	
32	JD62888-1F		TOCAQ.met	1.000	NPOC:1.017mg/L	
33	JD62888-3F		TOCAQ.met	1.000	NPOC:0.7451mg/L	
34	GP45915-MB1	[toc]	TOCAQ.met	1.000	NPOC:0.3734mg/L	
35	GP45915-B1		TOCAQ.met	1.000	NPOC:9.660mg/L	
36	JD62928-1		TOCAQ.met	1.000	NPOC:1.266mg/L	
37	GP45915-S1	JD62928-1	TOCAQ.met	1.000	NPOC:9.891mg/L	
38	GP45915-MSD1	JD62928-1	TOCAQ.met	1.000	NPOC:10.51mg/L	
39	JD62928-4	field blank	TOCAQ.met	1.000	NPOC:0.5833mg/L	
40	JD62928-2		TOCAQ.met	1.000	NPOC:2.618mg/L	
41	JD62928-3		TOCAQ.met	1.000	NPOC:1.072mg/L	
42	CCVA		TOCAQ.met	1.000	NPOC:46.47mg/L	
43	CCB		TOCAQ.met	1.000	NPOC:0.4340mg/L	
44	JD62929-1		TOCAQ.met	1.000	NPOC:2.236mg/L	
45	JD62929-2		TOCAQ.met	1.000	NPOC:5.310mg/L	
46	JD62929-3		TOCAQ.met	1.000	NPOC:0.8721mg/L	
47	CCV		TOCAQ.met	1.000	NPOC:22.98mg/L	
48	CCB		TOCAQ.met	1.000	NPOC:0.8250mg/L	
49	SPARGERCHK		TOCAQ.met	1.000	NPOC:0.4635mg/L	
50	CCVA		TOCAQ.met	1.000	NPOC:47.95mg/L	
51	CCB		TOCAQ.met	1.000	NPOC:0.6175mg/L	
52	JD62800-5	(2)	TOCAQ.met	1.000	NPOC:58.68mg/L	overrange, see rerun
53	JD62800-6	↓	TOCAQ.met	5.000	NPOC:277.5mg/L	overrange, see rerun

4/3/2023 12:14:34 PM

GN39916

MB

4-3-23

1/4

J230331wl.toc

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	Sample Name	Sample ID	Origin	Manual Dilution	Result	Comment
54	JD62800-5		TOCAQ.met	5.000	NPOC:58.56mg/L	
55	JD62800-6		TOCAQ.met	20.00	NPOC:280.6mg/L	
56	CCV		TOCAQ.met	1.000	NPOC:24.82mg/L	
57	CCB		TOCAQ.met	1.000	NPOC:0.4846mg/L	
58	SPARGERCHK		TOCAQ.met	1.000	NPOC:0.3626mg/L	

jd230331wl.toc

GN39916

MB

4-3-23



GN Batch ID: GN39916  
Date: 3-31-23

Test: Total Organic Carbon

Product: TOC or DOC

J230331 wl.toc

Method: **SM5310 B, C, or D-11, SW846 9060M**

Note: Refer to raw data and LIMS for information not shown below.

Autosampler Position #	Sample ID	pH	Dilution Factor	Bottle #	Comments
1	WASHCONF	<2			
2	CRI				
3	HSTD				
4	ICV				
5	ICB				
6	CCV				
7	CCB				
8	SPARGERCHK				
9	GP45856-MB1			-	[toc]
10	GP45856-B1			-	
11	JD62800-1			31	
12	GP45856-S1			32	JD62800-1
13	GP45856-MSD1			32	JD62800-1
14	JD62800-2			13	
15	JD62800-3			13	
16	JD62800-4			13	
17	JD62800-5			13	
18	CCVA				
19	CCB				
20	JD62800-6			13	
21	JD62800-7			13	
22	JD62800-8			13	
23	JD62800-9			13	Field Blank
24	GP45881-MB1			-	[doc]
25	GP45881-B1			-	
26	JD62888-2F			6	
27	GP45881-S1			6	JD62888-2F

Analyst: M3 Date: 4-3-23 QC Reviewer: \_\_\_\_\_ Date: \_\_\_\_\_

Comments: BSP: 500ul of 1000ppm KHP up to 50mL with DI H2O TV= 10 mg/L

MS/MSD: 200ul of 1000ppm KHP up to 20mL with sample TV=10 mg/L

ICV: 5mL of 100 ppm Sucrose up to 25mL with DI H2O TV=20 mg/L

Form: GN054-04

Rev. Date: 2/27/18

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GN Batch ID: GN39916  
Date: 3-31-23

## Test: Total Organic Carbon

Product:  or DOC

Method: **SM5310 B, C, or D-11, SW846 9060M**

Note: Refer to raw data and LIMS for information not shown below

Autosampler Position #	Sample ID	pH	Dilution Factor	Bottle #	Comments
28	GP45881-MSD1	<2		6	JD62888-2F
29	JD62888-4F			6	Field Blank
30	CCV				
31	CCB				
32	JD62888-1F			6	
33	JD62888-3F			6	
34	GP45915-MB1			-	[toc]
35	GP45915-B1			-	
36	JD62928-1			14	
37	GP45915-S1			14	JD62928-1
38	GP45915-MSD1			14	JD62928-1
39	JD62928-4			14	Field Blank
40	JD62928-2			14	
41	JD62928-3			14	
42	CCVA				
43	CCB				
44	JD62929-1			12	
45	JD62929-2			12	
46	JD62929-3			13	
47	CCV				
48	CCB				
49	SPARGERCHK				

Analyst: MB Date: 4-3-23 QC Reviewer:    Date:     
Comments:

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**Form: GN054-04**  
**Rev. Date: 2/27/18**



Analyst MB      Product Toc / DOC      Autopipette # 53  
Date 4-3-23      Batch ID GN39916      Class A Vol. Flask

## Sample Dilution Prep Log

OC Reviewer

Dat

Form: GN165-01  
Rev. Date: 2/25/03

9.6



## Reagent Information Log - TOC/DOC

GN 39916

<u>Reagent</u>	<u>Expiration Date</u>	<u>Reagent # or Manufacturer/Lot</u>
Potassium Hydrogen Phthalate (KHP), Stock Solution 1000 mg/L	4/13/2023	GNE1-72979-TOC
Carbonate/Bicarbonate Stock Solution	4/13/2023	GNE1-72982-TOC
Sparger Check Solution	4/13/2023	GNE3-70727-TOC
CCV Solution	4/14/2023	GNE3-70678-TOC
CCVA Solution (50 ppm)	4/14/2023	GNE3-70679-TOC
CRF Check	4/13/2023	GNE3-70729-TOC
HCl	4/10/2024	VWR CHEMICALS LOT #2022041418
pH Hydron Paper	8/1/2024	FISHER LOT #222221
Sucrose Stock Solution	4/14/2023	GNE3-70676-TOC

All standards and stocks were made as described in the SOP for this method (circle one):  Y or N  
 If no (N), see attached page for standards prep.

Form: GN087A67-043  
 Rev Date: 3/17/2023

9.6



## GENERAL CHEMISTRY STANDARD PREPARATION LOG FOR TOC (AQ)

Product: TOC  
 Analyst: Mß  
 Date: 4-3-25

Standard Description	Stock used to prepare standard	Stock concentration (mg/L)	Stock volume used (mL or g)	Balance/Pipette ID*	Diluent(a)	Final Volume	Final Conc. of Intermediate (mg/L)	Expiration Date	Analyst	Date
GNE1-72979-TOC	Fisher Lot: 178879	KHP	2.1528 g	B-39	DI H <sub>2</sub> O	1000 mL	1000 ppm	4/13/2023	MB	1/12/2023
GNE3-70728-TOC	GNE1-72979-TOC	1000 ppm	20 mL	Class A Pipet	DI H <sub>2</sub> O	200 mL	100 ppm	4/13/2023	MB	3/28/2023
GNE3-70676-TOC	Fisher Lot: 210554	Sucrose	0.0486 g	B-39	DI H <sub>2</sub> O	200 mL	100 ppm	4/14/2023	MB	3/17/2023
<b>KHP STDS</b>		Intermediate or Stock used to prepare standard	Intermediate or Stock concentration (mg/L)	Intermediate or Stock volume used (mL)	Balance/Pipette ID*	Diluent	Final Volume	Final Conc. of Standard (mg/L)	Expiration Date	Analyst
GNE3-70729-TOC	GNE3-70728-TOC	100 ppm	1.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	1.0	4/13/2023	MB	3/28/2023
GNE3-70730-TOC	GNE3-70728-TOC	100 ppm	2.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	2.0	4/13/2023	MB	3/28/2023
GNE3-70731-TOC	GNE3-70728-TOC	100 ppm	5.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	5.0	4/13/2023	MB	3/28/2023
GNE3-70732-TOC	GNE3-70728-TOC	100 ppm	10.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	10.0	4/13/2023	MB	3/28/2023
GNE3-70733-TOC	GNE3-70728-TOC	100 ppm	20.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	20.0	4/13/2023	MB	3/28/2023
GNE3-70734-TOC	GNE3-70728-TOC	100 ppm	30.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	30.0	4/13/2023	MB	3/28/2023
GNE3-70735-TOC	GNE3-70728-TOC	100 ppm	50.0	Class A Pipet	DI H <sub>2</sub> O	100 mL	50.0	4/13/2023	MB	3/28/2023
<b>KHP STDS</b>										
GNE3-70677-TOC	Acros Lot: A0383584	KHP	0.2307 g	B-39	DI H <sub>2</sub> O	100 mL	100 ppm	4/14/2023	MB	3/17/2023
GNE3-70678-TOC	GNE3-70677-TOC	100 ppm	50 mL	Class A Pipet	DI H <sub>2</sub> O	100 mL	25 ppm	4/14/2023	MB	3/17/2023
GNE3-70679-TOC	GNE3-70677-TOC	100 ppm	100 mL	Class A Pipet	DI H <sub>2</sub> O	100 mL	50 ppm	4/14/2023	MB	3/17/2023

\*If Class A glass pipettes are used, enter an A.

(a) Diluent reagent reference number:

All strikeouts must be initialed, dated, and reason applied if not transcription error

Expiration date:

Form: GN199-02  
 Rev. Date: 3/17/2023

# TOC-Control L Report

Chem36  
d230201wl.toc.tlx

**Instr. Information**

Instrument Options  
Catalyst

TOC/ASI/Sparge Kit/  
Regular Sensitivity

**Cal. Curve**

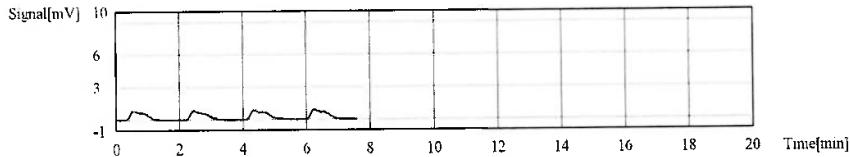
Sample Name: Untitled  
 Sample ID: Untitled  
 Cal. Curve: d230201wl.2023\_02\_01\_11\_45\_27.cal  
 Status: Completed

Type	Anal.
Standard	NPOC

Conc: 0.000mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	2 806	100uL	1 000	*****		2/1/2023 11:50:08 AM
2	2 648	100uL	1 000	*****		2/1/2023 11:52:23 AM
3	2 915	100uL	1 000	*****		2/1/2023 11:54:38 AM
4	2 824	100uL	1 000	*****		2/1/2023 11:56:53 AM

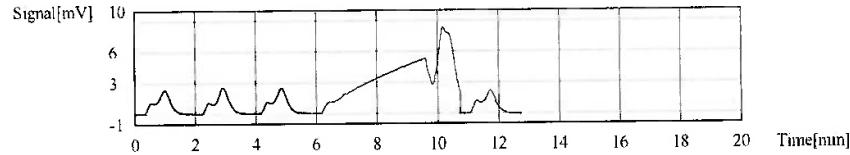
Acid Add. 0.000%  
 Sp. Time 600.0usec  
 Mean Area 2.798



Conc: 1.000mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	7 402	100uL	1 000	*****		2/1/2023 12:01:30 PM
2	7 958	100uL	1 000	*****		2/1/2023 12:03:47 PM
3	7 934	100uL	1 000	*****		2/1/2023 12:06:08 PM
4	57 74	100uL	1 000	T*****	E	2/1/2023 12:11:23 PM
5	7 582	100uL	1 000	*****		2/1/2023 12:14:09 PM

Acid Add. 0.000%  
 Sp. Time 600.0usec  
 Mean Area 7.719



Conc: 2.000mg/L

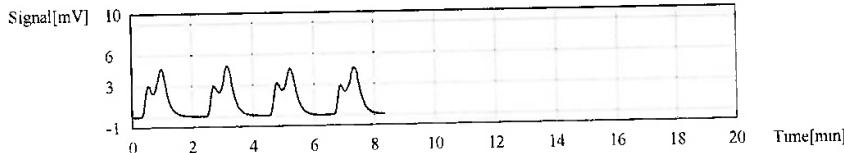
No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	16 98	100uL	1 050	*****		2/1/2023 12:18:47 PM
2	17 18	100uL	1 000	*****		2/1/2023 12:21:13 PM
3	17 02	100uL	1 050	*****		2/1/2023 12:23:41 PM
4	16 83	100uL	1 000	*****		2/1/2023 12:26:00 PM

# TOC-Control L Report

Chem36  
d230201wl.toc fix

Acid Add.  
Sp. Time  
Mean Area

0.000%  
600.0sec  
17.00

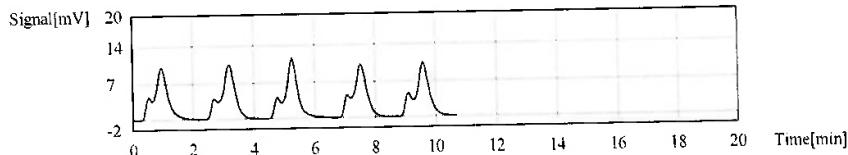


Conc. 5.000mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	32.71	100uL	1.000	*****	E	2/1/2023 12:30:53 PM
2	33.99	100uL	1.000	*****		2/1/2023 12:33:18 PM
3	34.26	100uL	1.000	*****		2/1/2023 12:35:58 PM
4	33.02	100uL	1.000	*****		2/1/2023 12:38:19 PM
5	33.81	100uL	1.000	*****		2/1/2023 12:40:48 PM

Acid Add.  
Sp. Time  
Mean Area

0.000%  
600.0sec  
33.77

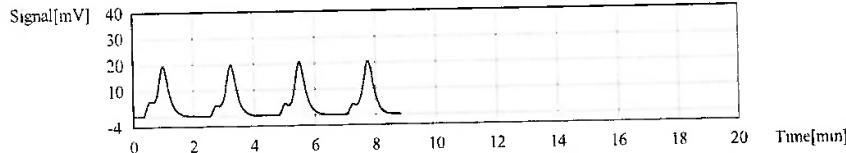


Conc: 10.00mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	58.33	100uL	1.000	*****		2/1/2023 12:45:28 PM
2	58.49	100uL	1.000	*****		2/1/2023 12:48:06 PM
3	58.80	100uL	1.000	*****		2/1/2023 12:50:42 PM
4	58.79	100uL	1.000	*****		2/1/2023 12:53:08 PM

Acid Add.  
Sp. Time  
Mean Area

0.000%  
600.0sec  
58.60

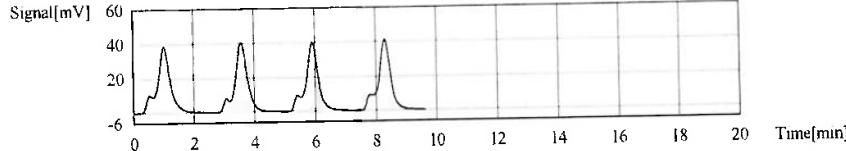


Conc: 20.00mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	116.5	100uL	1.000	*****		2/1/2023 12:58:24 PM
2	117.3	100uL	1.000	*****		2/1/2023 1:01:04 PM
3	116.9	100uL	1.000	*****		2/1/2023 1:03:48 PM
4	117.7	100uL	1.000	*****		2/1/2023 1:06:32 PM

Acid Add.  
Sp. Time  
Mean Area

0.000%  
600.0sec  
117.1

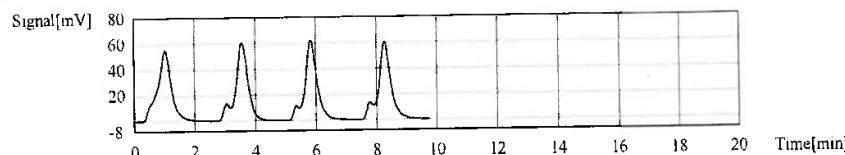


Conc. 30.00mg/L

## TOC-Control L Report

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	174.8	100 $\mu$ L	1.000	*****		2/1/2023 1:11:44 PM
2	174.5	100 $\mu$ L	1.000	*****		2/1/2023 1:14:24 PM
3	176.4	100 $\mu$ L	1.000	*****		2/1/2023 1:17:10 PM
4	176.3	100 $\mu$ L	1.000	*****		2/1/2023 1:20:00 PM

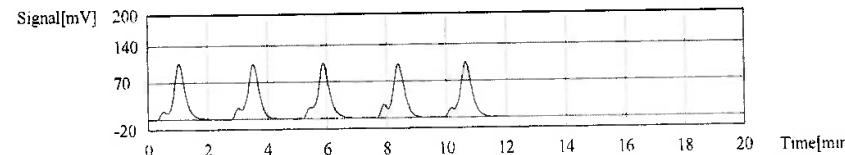
Acid Add. 0.000%  
 Sp. Time 600.0sec  
 Mean Area 175.5



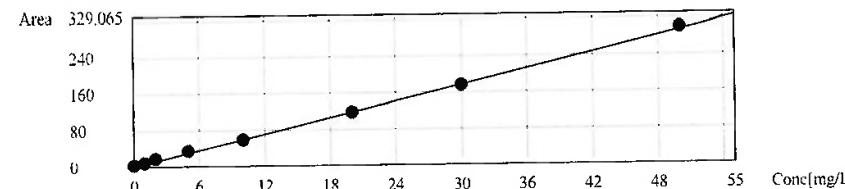
Conc: 50.00mg/L

No.	Area	Inj. Vol.	Aut. Dil.	Rem.	Ex.	Date / Time
1	292.2	100 $\mu$ L	1.000	*****	E	2/1/2023 1:25:12 PM
2	299.6	100 $\mu$ L	1.000	*****		2/1/2023 1:27:56 PM
3	301.2	100 $\mu$ L	1.000	*****		2/1/2023 1:30:48 PM
4	295.9	100 $\mu$ L	1.000	*****		2/1/2023 1:33:24 PM
5	299.9	100 $\mu$ L	1.000	*****		2/1/2023 1:36:10 PM

Acid Add. 0.000%  
 Sp. Time 600.0sec  
 Mean Area 299.2



Slope: 5.875  
 Intercept 0.000  
 r^2 0.9994  
 r 0.9997  
 Zero Shift Yes



# TOC-Control L Report

Chem365  
d230331w1 toc tlx

**Instr.Information**

Instrument Options  
Catalyst

TOC/ASI/Sparge Kit/  
Regular Sensitivity

**Sample**

Sample Name: WASHCONF  
 Sample ID:  
 Origin:  
 Status: Completed  
 Chk. Result

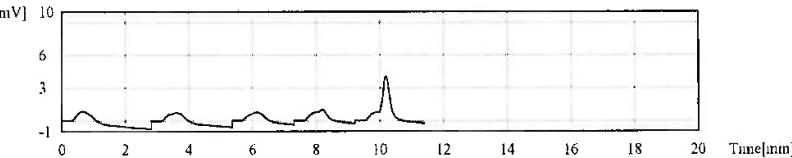
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.5131mg/L

1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2 179	0.3709mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:28:51 AM
2	3 358	0.5715mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:31:38 AM
3	2 918	0.4966mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:33:55 AM
4	3 605	0.6136mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:36:05 AM
5	8 560	1.457mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:38:31 AM

Mean Conc. 0.5131mg/L  
 CV Conc. 20.75%

**Sample**

Sample Name: CRI  
 Sample ID:  
 Origin:  
 Status: Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 1.097mg/L

1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	6 943	1.182mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:44:02 AM
2	7 321	1.246mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:46:26 AM
3	6 342	1.079mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:48:39 AM
4	6 418	1.022mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:50:52 AM
5	6 077	1.034mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/1/2023 9:53:04 AM

# TOC-Control L Report

**Sample**

Sample Name: HSTD  
 Sample ID:  
 Origin: TOCAQSW846 met  
 Status Completed  
 Chk. Result

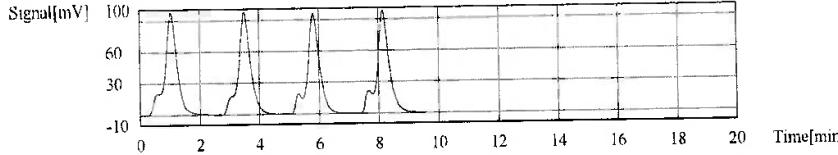
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 47.05mg/L

## 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	277.7	47.26mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 9:59:00 AM	
2	278.4	47.38mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:01:34 AM	
3	271.5	46.21mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:04:11 AM	
4	278.2	47.35mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:06:52 AM	

Mean Conc. 47.05mg/L  
 CV Conc. 1.20%

**Sample**

Sample Name: ICV  
 Sample ID:  
 Origin: TOCAQSW846 met  
 Status Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 19.89mg/L

## 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	117.1	19.93mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:30:58 AM	
2	116.8	19.88mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:33:18 AM	
3	116.8	19.88mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:35:32 AM	
4	116.8	19.88mg/L	100μL	1.000	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:37:54 AM	

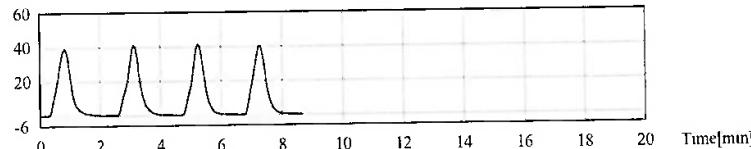
# TOC-Control L Report

Chem36  
d230331wl toc tlx

Mean Conc.  
CV Conc.

19.89mg/L  
0.13%

Signal[mV]



## Sample

Sample Name: ICB  
 Sample ID: TOCAQSW846 met  
 Origin:  
 Status: Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.3547mg/L

## 1. Det

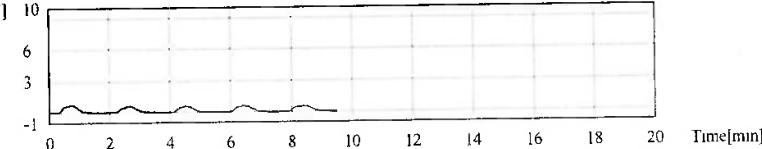
Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Auf. Dil.	Ex.	Cal. Curve	Date / Time
1	2.269	0.3862mg/L	100uL	1.000		d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:43:21 AM
2	1.999	0.3402mg/L	100uL	1.000		d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:45:26 AM
3	1.770	0.3013mg/L	100uL	1.000	E	d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:47:31 AM
4	2.034	0.3462mg/L	100uL	1.000		d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:49:36 AM
5	2.035	0.3464mg/L	100uL	1.000		d230201wl_2023_02_01_11_45_27.cal	3/1/2023 10:51:41 AM

Mean Conc.  
CV Conc.

0.3547mg L  
5.96%

Signal[mV]



## Sample

Sample Name: CCV  
 Sample ID: TOCAQSW846 met  
 Origin:  
 Status: Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 22.99mg/L

## 1. Det

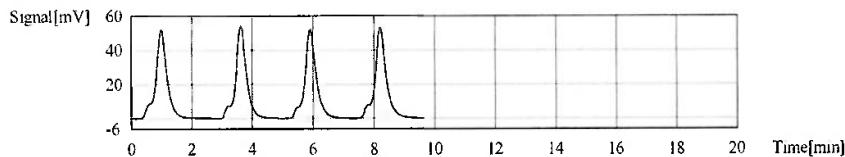
Anal.. NPOC

# TOC-Control L Report

Chem360  
d230331wl.toc.tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	134.2	22.84mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:08:49 PM	
2	136.0	23.15mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:11:22 PM	
3	135.1	22.99mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:13:55 PM	
4	135.1	22.99mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:16:52 PM	

Mean Conc. 22.99mg/L  
CV Conc 0.54%



## Sample

Sample Name: CCB  
Sample ID:  
Origin:  
Status: Completed  
Chk. Result

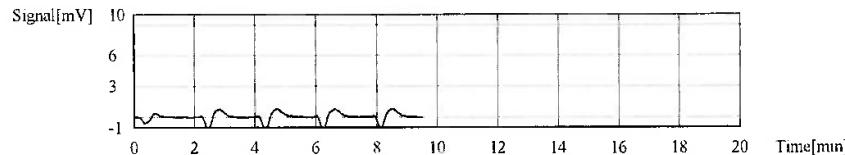
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.3092mg/L

## 1. Det

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	1.491	0.2538mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:22:07 PM	
2	4.765	0.8110mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:24:12 PM	
3	5.607	0.9543mg/L	100uL	1.000	E d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:26:17 PM	
4	0.4005	0.06816mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:28:22 PM	
5	0.6100	0.1038mg/L	100uL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 12:30:27 PM	

Mean Conc. 0.3092mg/L  
CV Conc 111.28%



## Sample

Sample Name: SPARGERCHK  
Sample ID:  
Origin:  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.3068mg/L

## 1. Det

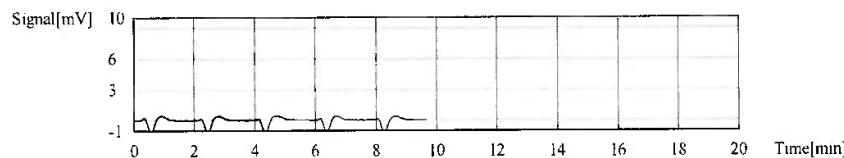
# TOC-Control L Report

Chem36  
d230331wl toc tlx

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	0.000	0.000mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:35 58 PM
2	3.632	0.6182mg/L	100uL	1 000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:38 03 PM
3	5.422	0.9228mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:40 16 PM
4	0.000	0.000mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:42 21 PM
5	3.579	0.6091mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:44 26 PM

Mean Conc. 0.3068mg/L  
CV Conc 115.48%

**Sample**

Sample Name: GP45856-MB1  
Sample ID: [toc]  
Origin: TOCAQ.net  
Status: Completed  
Chk. Result

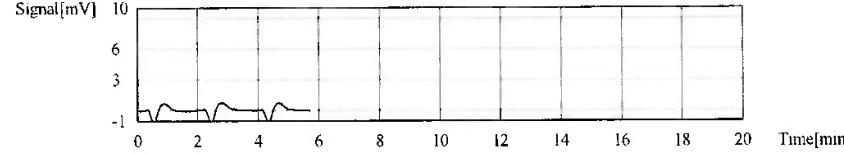
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1 000	NPOC 0.8149mg/L

## 1. Det

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	0.3410	0.05804mg/L	100uL	1 000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:49 10 PM
2	4.968	0.8455mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:51 15 PM
3	4.608	0.7843mg/L	100uL	1 000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 12:53 20 PM

Mean Conc. 0.8149mg/L  
CV Conc 5.32%

**Sample**

Sample Name: GP45856-B1  
Sample ID:  
Origin: TOCAQ.net  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1 000	NPOC 0.343mg/L

## 1. Det

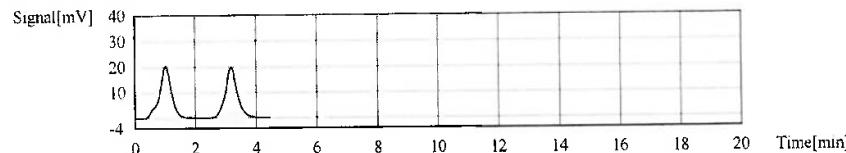
5:30

4/3/2023 12:14:47 PM

# TOC-Control L Report

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	55.52	9.449mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 3:43:42 PM
2	54.27	9.237mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 3:46:05 PM

Mean Conc. 9.343mg/L  
CV Conc 1.61%Sample

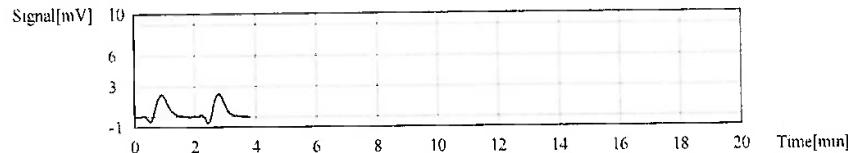
Sample Name: JD62800-1  
 Sample ID:  
 Origin: TOCAQ met  
 Status: Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 1.196mg/L

## 1. Det

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	7.086	1.206mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:12:28 PM
2	6.974	1.187mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:14:34 PM

Mean Conc. 1.196mg/L  
CV Conc 1.13%Sample

Sample Name: GP45856-S1  
 Sample ID: JD62800-1  
 Origin: TOCAQ met  
 Status: Completed  
 Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 9.196mg/L

## 1. Det

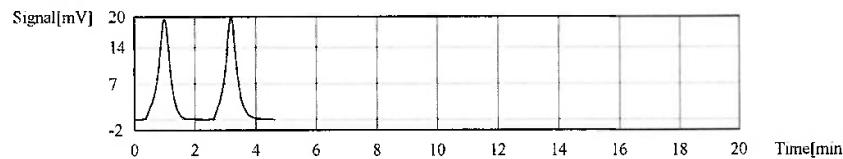
Anal.. NPOC

# TOC-Control L Report

Chem365  
d230331wl toc.ttx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	54.20	9.225mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:23:56 PM
2	53.86	9.167mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:26:30 PM

Mean Conc. 9.196mg/L  
CV Conc 0.44%



## Sample

Sample Name: GP45856-MSD1  
Sample ID: JD62800-1  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

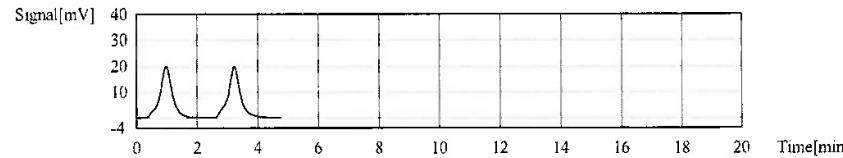
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 9.487mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	55.60	9.463mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:35:07 PM
2	55.88	9.511mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:37:48 PM

Mean Conc. 9.487mg/L  
CV Conc 0.36%



## Sample

Sample Name: JD62800-2  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 3.895mg/L

### 1. Det

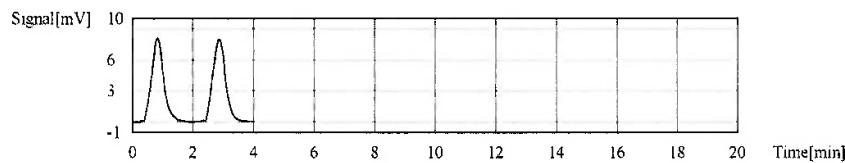
Anal.: NPOC

# TOC-Control L Report

Chem56  
d230331wl.toc.ltx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	23 15	3.940mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:46:04 PM
2	22 62	3.850mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:48:14 PM

Mean Conc. 3.895mg/L  
CV Conc 1.64%



## Sample

Sample Name: JD62800-3  
Sample ID:  
Origin: TOCAQ met  
Status Completed  
Chk. Result

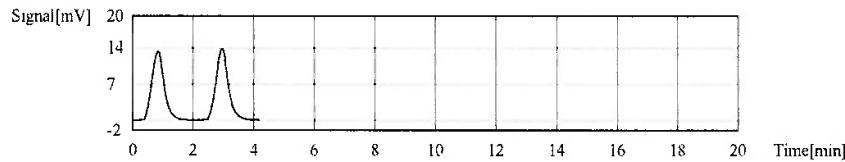
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 6.549mg/L

### 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	38 44	6.542mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:57:17 PM
2	38 52	6.556mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 5:59:32 PM

Mean Conc. 6.549mg/L  
CV Conc 0.15%



## Sample

Sample Name: JD62800-4  
Sample ID:  
Origin: TOCAQ.met  
Status Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 6.508mg/L

### 1. Det

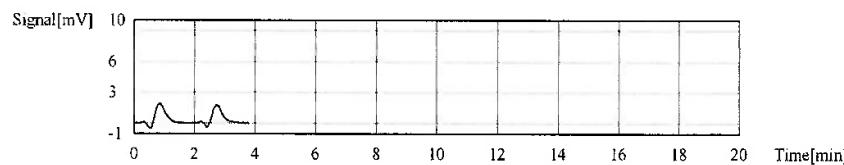
Anal. NPOC

# TOC-Control L Report

Chem36  
d230331w1 toc.ttx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	3.923	0.6677mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27 cal		3/31/2023 6:08:15 PM
2	3.725	0.6340mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27 cal		3/31/2023 6:10:20 PM

Mean Conc. 0.6508mg/L  
CV Conc 3.66%



## Sample

Sample Name: JD62800-5  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

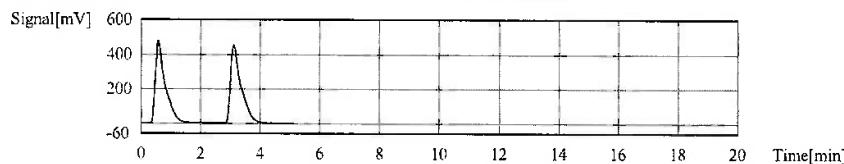
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 189.0mg/L

### 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	1089	185.3mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27 cal		3/31/2023 6:20:01 PM
2	1132	192.7mg/L	100uL	1.000	d230201wl2023_02_01_11_45_27 cal		3/31/2023 6:22:56 PM

Mean Conc. 189.0mg/L  
CV Conc 2.74%



## Sample

Sample Name: CCVA  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 46.43mg/L

### 1. Det

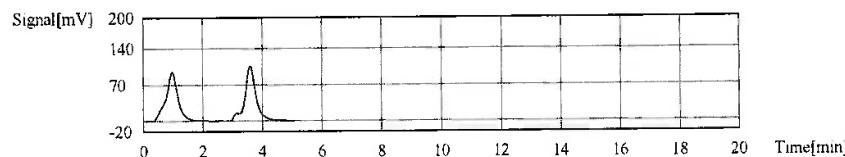
Anal. NPOC

# TOC-Control L Report

Chem36  
d230331wl toc tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	270.4	46.02mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 6:31:14 PM
2	275.2	46.84mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 6:34:00 PM

Mean Conc. 46.43mg/L  
CV Conc 1.24%



## Sample

Sample Name: CCB  
Sample ID:  
Origin: TOCAQ.met  
Status: Completed  
Chk. Result

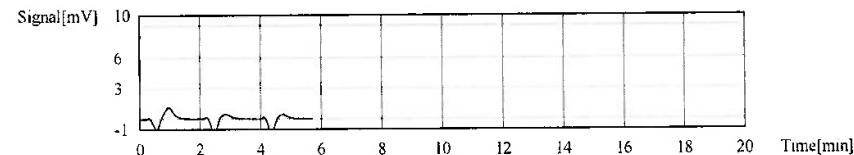
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.7008mg/L

## 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	5.271	0.8971mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 6:41:43 PM
2	4.463	0.7596mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 6:43:48 PM
3	3.772	0.6420mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 6:45:53 PM

Mean Conc. 0.7008mg/L  
CV Conc 11.87%



## Sample

Sample Name: JD62800-6  
Sample ID:  
Origin: TOCAQ.met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 358.1mg/L

## 1. Det

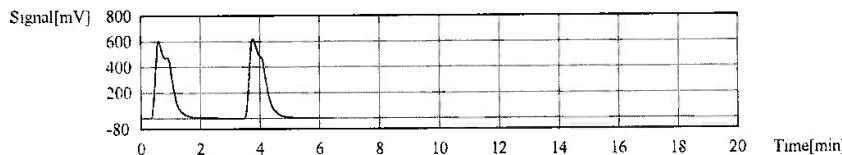
Anal. NPOC

# TOC-Control L Report

Chem360  
d230331wl toc.tb

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2080	354.0mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 6:54:07 PM
2	2128	362.2mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 6:58:01 PM

Mean Conc. 358.1mg/L  
CV Conc 1.61%



## Sample

Sample Name: JD62800-7  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

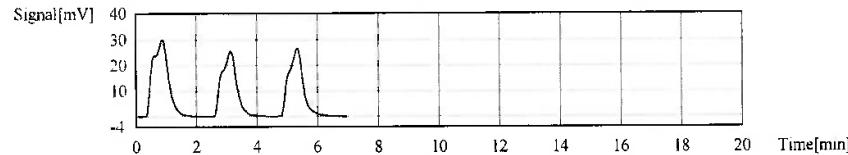
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 14.96mg/L

## 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	108.0	18.38mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7:04:23 PM
2	87.02	14.81mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7:06:46 PM
3	88.78	15.11mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7:08:27 PM

Mean Conc. 14.96mg/L  
CV Conc 1.42%



## Sample

Sample Name: JD62800-8  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.7392mg/L

## 1. Det

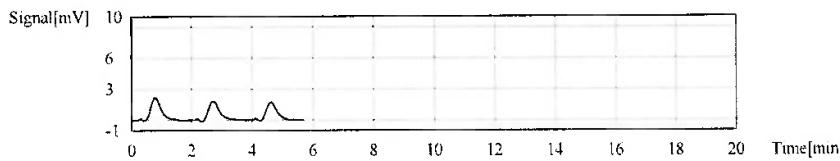
Anal. NPOC

# TOC-Control L Report

Chem36  
d230231wl.toc.ttx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	5.344	0.9095mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7 15 11 PM
2	4.416	0.7516mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7 17 20 PM
3	4.270	0.7267mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7 19 25 PM

Mean Conc. 0.7392mg/L  
CV Conc 2.38%



## Sample

Sample Name: JD62800-9  
Sample ID: field blank  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

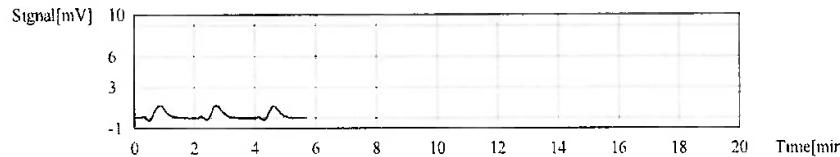
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.4505mg/L

### 1. Det

Anal.. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	2.760	0.4697mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7 26 27 PM
2	3.086	0.5252mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7 28 32 PM
3	2.514	0.4313mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 7 30 37 PM

Mean Conc. 0.4505mg/L  
CV Conc 6.04%



## Sample

Sample Name: GP45881-MB1  
Sample ID: [doc]  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.1368mg/L

### 1. Det

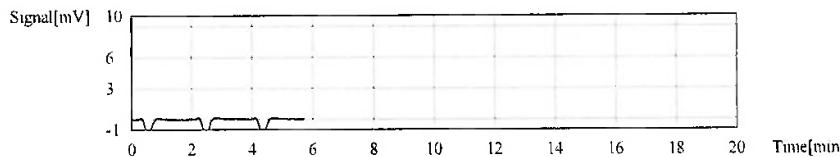
Anal.. NPOC

# TOC-Control L Report

Chem365  
d230331w1 toc.ttx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2.411	0.4103mg/L	100uL	1.000		d230201w1 2023_02_01_11_45_27.cal	3/31/2023 7:37:29 PM
2	0.000	0.000mg/L	100uL	1.000		d230201w1 2023_02_01_11_45_27.cal	3/31/2023 7:39:39 PM
3	0.000	0.000mg/L	100uL	1.000		d230201w1 2023_02_01_11_45_27.cal	3/1/2023 7:41:44 PM

Mean Conc. 0.1368mg/L  
CV Conc 173.2%



## Sample

Sample Name: GP45881-B1  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

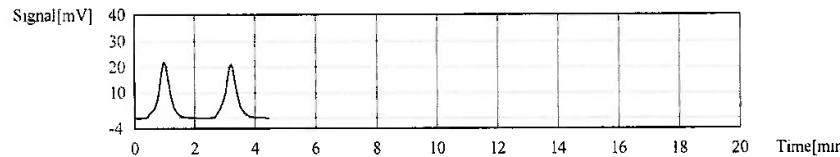
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.592mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	56.16	9.558mg/L	100uL	1.000		d230201w1 2023_02_01_11_45_27.cal	3/31/2023 7:49:00 PM
2	56.56	9.626mg/L	100uL	1.000		d230201w1 2023_02_01_11_45_27.cal	3/31/2023 7:51:22 PM

Mean Conc. 9.592mg/L  
CV Conc 0.50%



## Sample

Sample Name: JD62888-2F  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.6218mg/L

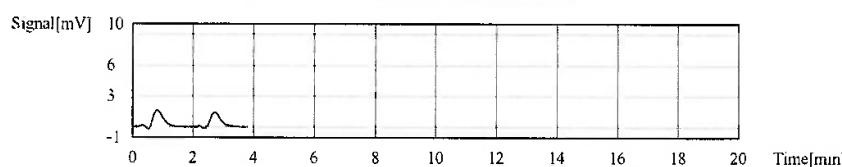
### 1. Det

Anal.: NPOC

# TOC-Control L Report

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	3 619	0.6159mg/L	100ul	1 000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 7:59:48 PM
2	3 688	0.6277mg/L	100ul	1 000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 8:01:56 PM

Mean Conc. 0.6218mg/L  
CV Conc 1.34%



## Sample

Sample Name: GP45881-S1  
Sample ID: JD62888-2F  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

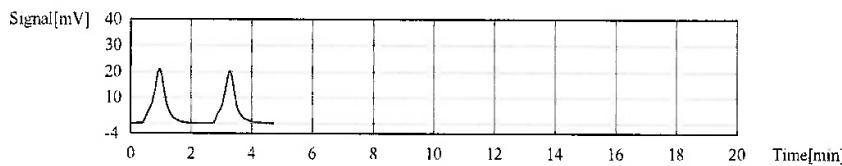
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1 000	NPOC @ 807mg/L

1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	58 95	10.03mg/L	100ul	1 000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 8:11:24 PM
2	56 29	9.580mg/L	100ul	1 000	d230201wl 2023_02_01_11_45_27 cal		3/31/2023 8:13:56 PM

Mean Conc. 9.807mg/L  
CV Conc 3.26%



## Sample

Sample Name: GP45881-MSD1  
Sample ID: JD62888-2F  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1 000	NPOC @ 703mg/L

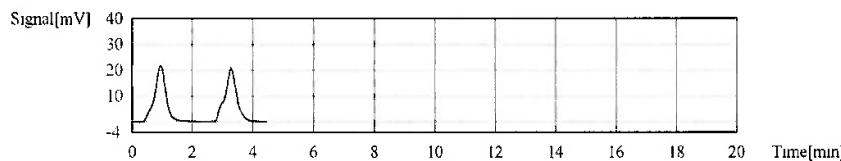
1. Det

Anal.: NPOC

# TOC-Control L Report

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	57 58	0.800mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/31/2023 8:22:36 PM
2	56 44	9.606mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/31/2023 8:24:51 PM

Mean Conc. 0.703mg/L  
CV Conc 1.41%

**Sample**

Sample Name: JD62888-4F  
Sample ID: field blank  
Origin: TOCA/Q met  
Status: Completed  
Chk. Result

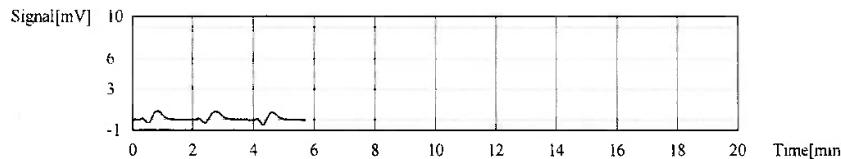
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.2537mg/L

**1. Det**

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	2706	0.4606mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/31/2023 8:33:16 PM
2	1671	0.2844mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/31/2023 8:35:26 PM
3	1310	0.2230mg/L	100uL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	3/31/2023 8:37:31 PM

Mean Conc. 0.2537mg/L  
CV Conc 17.13%

**Sample**

Sample Name: CCV  
Sample ID:  
Origin: TOCA/Q met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 22.91mg/L

**1. Det**

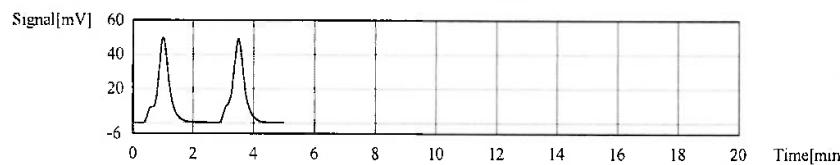
Anal.: NPOC

# TOC-Control L Report

Chem36  
d230201w1.toc.ttx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	132.2	22.50mg/L	100uL	1.000	E	d230201w1.2023_02_01_11_45_27.cal	3/31/2023 8:45:12 PM
2	137.0	23.32mg/L	100uL	1.000	E	d230201w1.2023_02_01_11_45_27.cal	3/31/2023 8:47:57 PM

Mean Conc. 22.91mg/L  
CV Conc 2.52%



## Sample

Sample Name: CCB  
 Sample ID:  
 Origin: TOCAQ met  
 Status: Completed  
 Chk. Result: 9.6

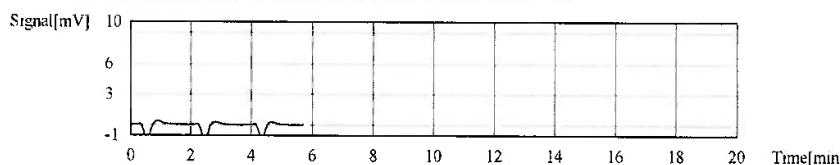
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.5875mg/L

## 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	4913	0.8362mg/L	100uL	1.000	E	d230201w1.2023_02_01_11_45_27.cal	3/31/2023 9:04:09 PM
2	3768	0.6413mg/L	100uL	1.000	E	d230201w1.2023_02_01_11_45_27.cal	3/31/2023 9:06:16 PM
3	3136	0.5337mg/L	100uL	1.000	E	d230201w1.2023_02_01_11_45_27.cal	3/31/2023 9:08:21 PM

Mean Conc. 0.5875mg/L  
CV Conc 12.95%



## Sample

Sample Name: JD62888-1F  
 Sample ID:  
 Origin: TOCAQ met  
 Status: Completed  
 Chk. Result:

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 1.017mg/L

## 1. Det

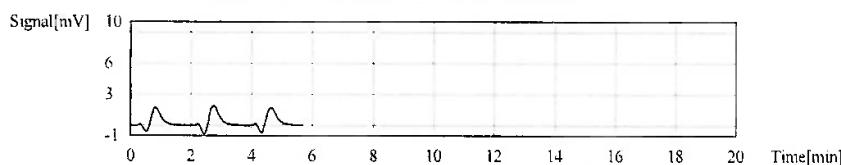
Anal.: NPOC

# TOC-Control L Report

Chem365  
d230331wl.toc.ltx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	5 296	0.9014mg/L	100uL	1 000	E	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 9:15:19 PM
2	6 043	1.029mg/L	100uL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 9:17:28 PM
3	5 907	1.005mg/L	100uL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 9:19:33 PM

Mean Conc. 1.017mg/L  
CV Conc 1.61%



## Sample

Sample Name: JD62888-3F  
Sample ID.  
Origin:  
Status Completed  
Chk. Result

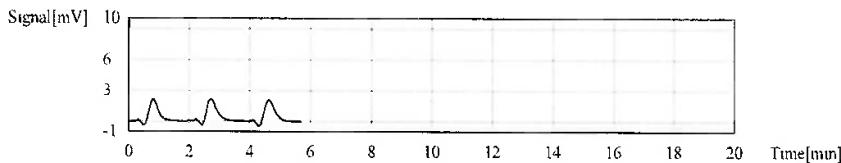
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1 (xx)	NPOC 0.7451mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	4 063	0.6915mg/L	100uL	1 000	E	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 9:26:29 PM
2	4 390	0.7472mg/L	100uL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 9:28:38 PM
3	4 366	0.7431mg/L	100uL	1 000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 9:30:43 PM

Mean Conc. 0.7451mg/L  
CV Conc 0.39%



## Sample

Sample Name: GP45915-MB1  
Sample ID. [toc]  
Origin: TOCAQ.met  
Status Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1 (xx)	NPOC 0.3734mg/L

### 1. Det

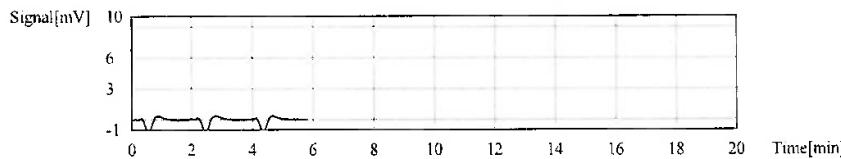
Anal.: NPOC

# TOC-Control L Report

Chem360  
d230331wl toc.ttx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	3 132	0.5331mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27.cal	3/31/2023 9:37:39 PM
2	3 450	0.5872mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27.cal	3/31/2023 9:39:44 PM
3	0.000	0.0000mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27.cal	3/31/2023 9:41:56 PM

Mean Conc. 0.3734mg/L  
CV Conc 86.91%



## Sample

Sample Name: GP-15915-B1  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

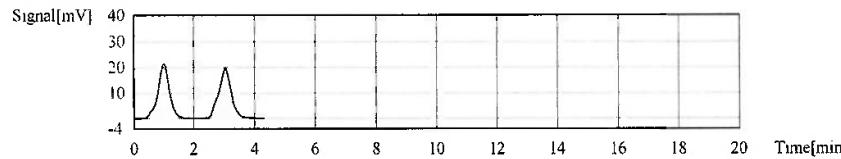
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.660mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	57.21	9.737mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27.cal	3/31/2023 9:49:02 PM
2	56.30	9.582mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27.cal	3/31/2023 9:51:25 PM

Mean Conc. 9.660mg/L  
CV Conc 1.13%



## Sample

Sample Name: JD62928-1  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 1.266mg/L

### 1. Det

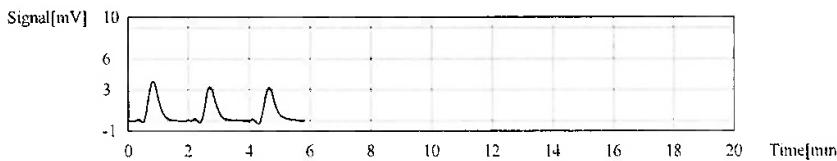
Anal.: NPOC

# TOC-Control L Report

Chem360  
d230331wl toc ltx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	8943	1.522mg/L	100uL	1.000	B	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:00:09 PM
2	7095	1.208mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:02:05 PM
3	7779	1.324mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:04:17 PM

Mean Conc. 1.266mg/L  
CV Conc 6.50%



## Sample

Sample Name: GP45915-S1  
Sample ID: JD62928-1  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

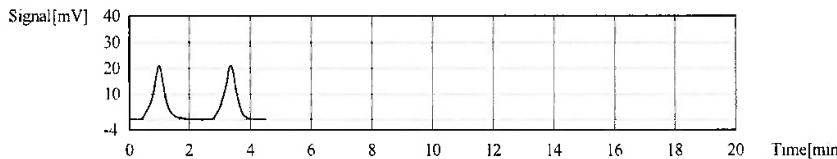
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 9.891mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	5822	9.909mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:11:38 PM
2	5801	9.873mg/L	100uL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:13:56 PM

Mean Conc. 9.891mg/L  
CV Conc 0.26%



## Sample

Sample Name: GP45915-MSD1  
Sample ID: JD62928-1  
Origin: TOCAQ.met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 10.51mg/L

### 1. Det

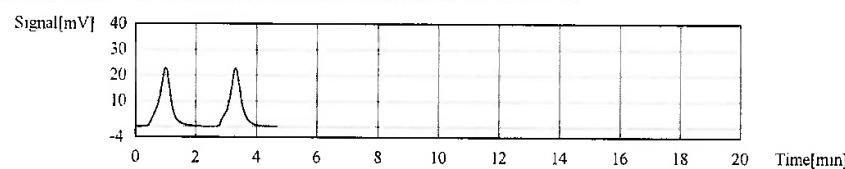
Anal.: NPOC

# TOC-Control L Report

Chem36  
d230331wl.toc.tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	62.14	10.58mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 10:22:46 PM
2	61.36	10.44mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 10:25:16 PM

Mean Conc. 10.51mg/L  
CV Conc 0.89%



## Sample

Sample Name: JD62928-4  
Sample ID: field blank  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

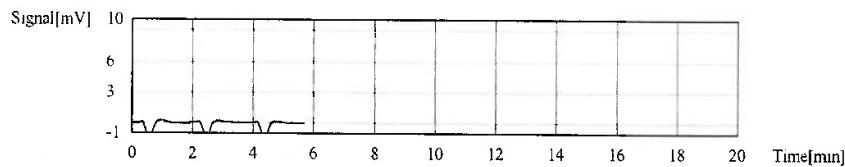
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.5833mg/L

### 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	3.439	0.5853mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 10:33:30 PM
2	3.708	0.6311mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 10:35:39 PM
3	3.134	0.5334mg/L	100µL	1.000		d230201wl.2023_02_01_11_45_27.cal	3/31/2023 10:37:44 PM

Mean Conc. 0.5833mg/L  
CV Conc 8.38%



## Sample

Sample Name: JD62928-2  
Sample ID: field blank  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 2.618mg/L

### 1. Det

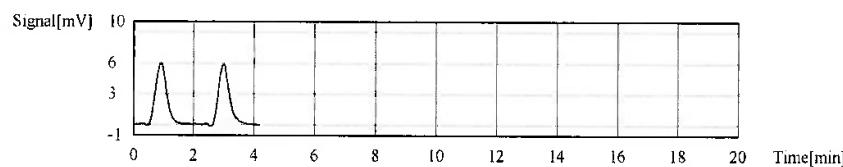
Anal. NPOC

# TOC-Control L Report

Chem36  
d230331wl toc ttx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	15.59	2.653mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:44:52 PM
2	15.17	2.582mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:47:08 PM

Mean Conc. 2.618mg/L  
CV Conc 1.93%



## Sample

Sample Name: JD62928-3  
Sample ID:  
Origin:  
Status Completed  
Chk. Result

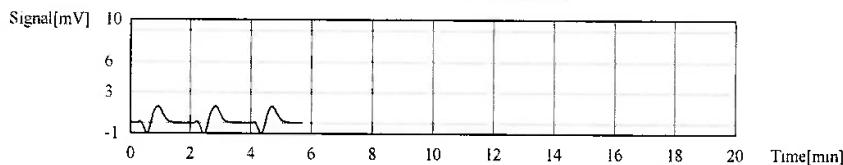
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 1.072mg/L

### 1. Det

Anal. NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	6.230	1.060mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:55:50 PM
2	2.342	0.3986mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 10:58:00 PM
3	6.370	1.084mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27 cal	3/31/2023 11:00:05 PM

Mean Conc. 1.072mg/L  
CV Conc 1.57%



## Sample

Sample Name: CCVA  
Sample ID:  
Origin:  
Status Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 46.47mg/L

### 1. Det

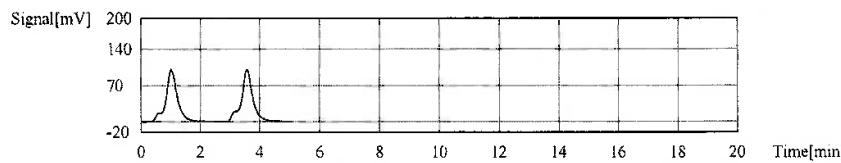
Anal. NPOC

# TOC-Control L Report

Chem56  
d230331wl toc thx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	270.1	45.97mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 11:07:30 PM
2	276.0	46.97mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 11:10:28 PM

Mean Conc. 46.47mg/L  
CV Conc 1.53%



## Sample

Sample Name: CCB  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

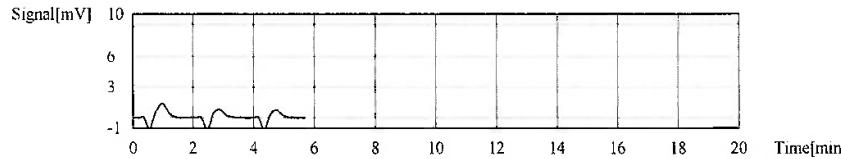
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.4340mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	6.382	1.086mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 11:18:11 PM
2	4.886	0.8316mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 11:20:20 PM
3	0.2143	0.03647mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	3/31/2023 11:22:30 PM

Mean Conc. 0.4340mg/L  
CV Conc 129.54%



## Sample

Sample Name: JD62929-1  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 2.236mg/L

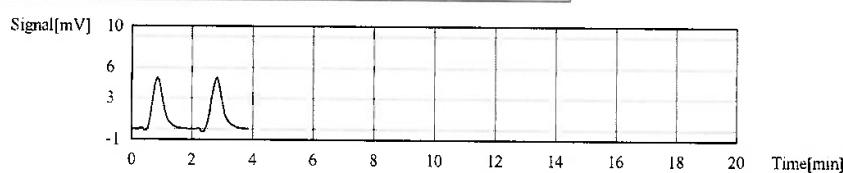
### 1. Det

Anal.: NPOC

# TOC-Control L Report

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	12.71	2.163mg/L	100μL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 11:29:30 PM	
2	13.56	2.308mg/L	100μL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 11:31:38 PM	

Mean Conc. 2.236mg/L  
CV Conc 4.58%



## Sample

Sample Name: JD62929-2  
Sample ID:  
Origin: TOCAQ met  
Status Completed  
Chk. Result

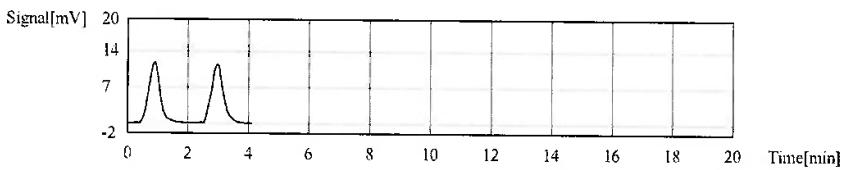
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 5.310mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	30.87	5.254mg/L	100μL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 11:40:44 PM	
2	31.53	5.366mg/L	100μL	1.000	d230201wl.2023_02_01_11_45_27.cal	3/31/2023 11:42:53 PM	

Mean Conc. 5.310mg/L  
CV Conc 1.50%



## Sample

Sample Name: JD62929-3  
Sample ID:  
Origin: TOCAQ met  
Status Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.8721mg/L

### 1. Det

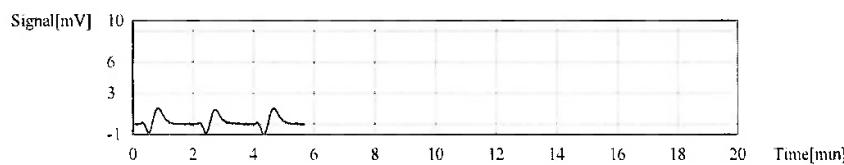
Anal. NPOC

# TOC-Control L Report

Chem36  
d230331w1 toc tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	4.878	0.8302mg/L	100uL	1.000	E	d230201w1 2023_02_01_11_45_27 cal	3/31/2023 11:51:41 PM
2	2.610	0.3421mg/L	100uL	1.000	E	d230201w1 2023_02_01_11_45_27 cal	3/31/2023 11:53:49 PM
3	5.570	0.9140mg/L	100uL	1.000	E	d230201w1 2023_02_01_11_45_27 cal	3/31/2023 11:55:54 PM

Mean Conc. 0.8721mg/L  
CV Conc 6.79%



## Sample

Sample Name: CCV  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

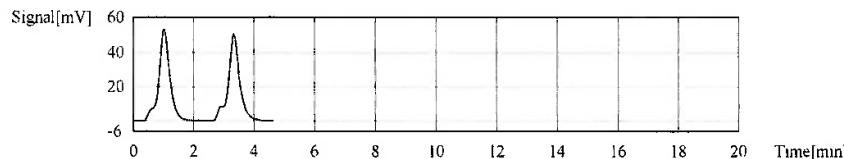
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 22.98mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil	Ex.	Cal. Curve	Date / Time
1	135.7	23.10mg/L	100uL	1.000	E	d230201w1 2023_02_01_11_45_27 cal	4/1/2023 12:03:15 AM
2	134.3	22.86mg/L	100uL	1.000	E	d230201w1 2023_02_01_11_45_27 cal	4/1/2023 12:05:49 AM

Mean Conc. 22.98mg/L  
CV Conc 0.73%



## Sample

Sample Name: CCB  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.8250mg/L

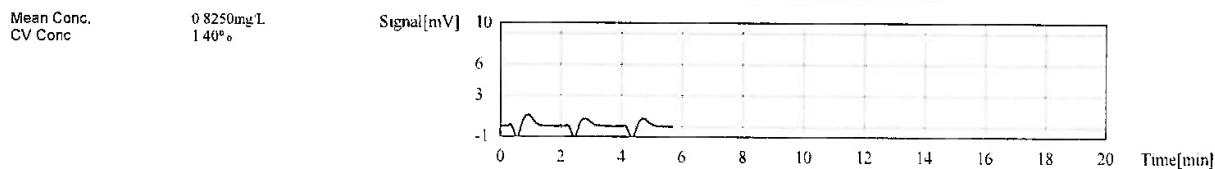
### 1. Det

Anal.: NPOC

# TOC-Control L Report

Chem36  
d230331wl toc lfx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	4.895	0.8331mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/1/2023 12:14:01 AM
2	4.183	0.7119mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/1/2023 12:16:10 AM
3	4.799	0.8168mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/1/2023 12:18:15 AM

**Sample**

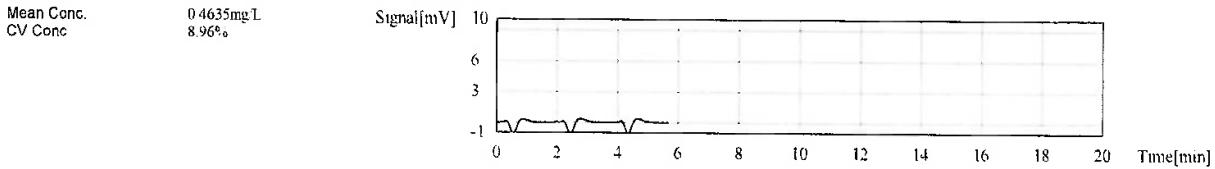
Sample Name: SPARGERCHK  
 Sample ID:  
 Origin:  
 Status TOCAQ.met  
 Chk. Result Completed

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.4635mg/L

## 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2.551	0.4342mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/1/2023 12:25:12 AM
2	2.896	0.4929mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/1/2023 12:27:21 AM
3	0.000	0.0000mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/1/2023 12:29:26 AM

**Sample**

Sample Name: CCVA  
 Sample ID:  
 Origin:  
 Status TOCAQ.met  
 Chk. Result Completed

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 47.95mg/L

## 1. Det

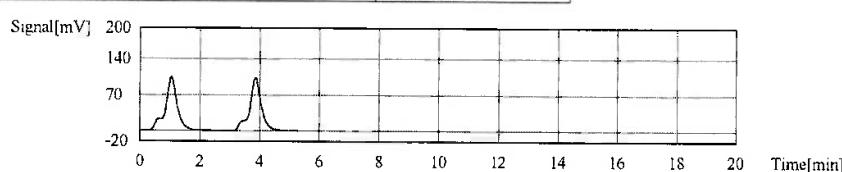
Anal.: NPOC

# TOC-Control L Report

Chem56  
d230331wl toc ltx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	284.9	48.40mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/3/2023 9:51:33 AM
2	278.6	47.42mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/3/2023 9:54:17 AM

Mean Conc. 47.95mg/L  
CV Conc 1.58%



## Sample

Sample Name: CCB  
Sample ID:  
Origin:  
Status: TOCAQ met  
Chk. Result: Completed

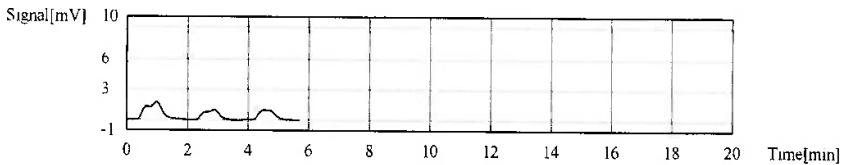
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.6175mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	6.082	1.035mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/3/2023 10:01:48 AM
2	3.638	0.6192mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/3/2023 10:03:53 AM
3	3.618	0.6158mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27 cal	4/3/2023 10:05:58 AM

Mean Conc. 0.6175mg/L  
CV Conc 0.39%



## Sample

Sample Name: JD62800-5  
Sample ID:  
Origin:  
Status: TOCAQ met  
Chk. Result: Completed

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 58.68mg/L

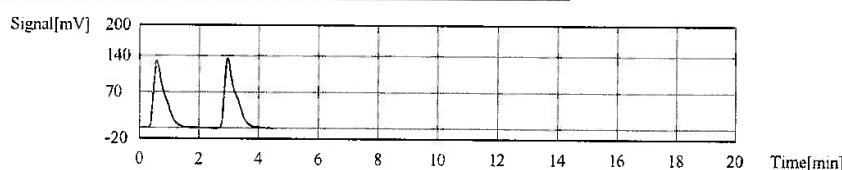
### 1. Det

Anal. NPOC

# TOC-Control L Report

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	336.9	57.34mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 10:13:26 AM
2	352.6	60.01mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 10:15:51 AM

Mean Conc. 58.68mg/L  
CV Conc 3.22%

**Sample**

Sample Name: JD62800-6  
Sample ID:  
Origin:  
Status TOCAQ met  
Chk. Result Completed

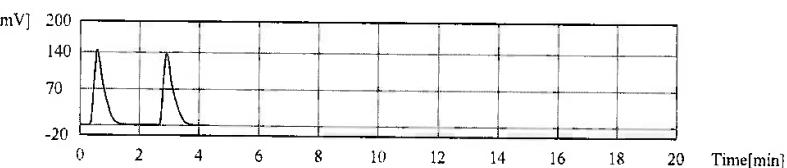
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	5.000	NPOC 277.5mg/L

## 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	331.6	282.2mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 10:24:33 AM
2	320.5	272.7mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 10:26:49 AM

Mean Conc. 277.5mg/L  
CV Conc 2.41%

**Sample**

Sample Name: JD62800-5  
Sample ID:  
Origin:  
Status TOCAQ met  
Chk. Result Completed

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	5.000	NPOC 58.56mg/L

## 1. Det

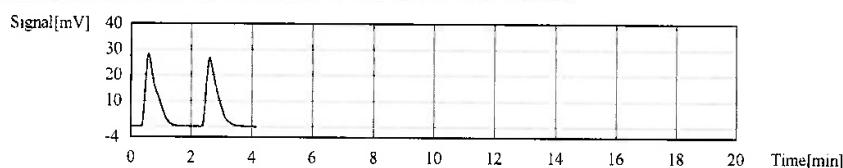
Anal. NPOC

# TOC-Control L Report

Chem36  
d230331wl toc tlx

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	70.16	59.71mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 11:16:33 AM
2	67.47	57.42mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 11:18:51 AM

Mean Conc. 58.56mg/L  
CV Conc 2.76%



## Sample

Sample Name: JD62800-6  
Sample ID:  
Origin:  
Status: TOCAQ met  
Chk. Result: Completed

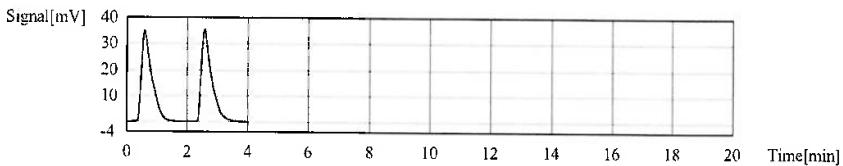
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	20.00	NPOC 280.6mg/L

### 1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	82.41	280.5mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 11:27:41 AM
2	82.48	280.8mg/L	100uL	1.000	d230201wl 2023_02_01_11_45_27 cal		4/3/2023 11:29:54 AM

Mean Conc. 280.6mg/L  
CV Conc 0.06%



## Sample

Sample Name: CCV  
Sample ID:  
Origin:  
Status: TOCAQ met  
Chk. Result: Completed

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 24.82mg/L

### 1. Det

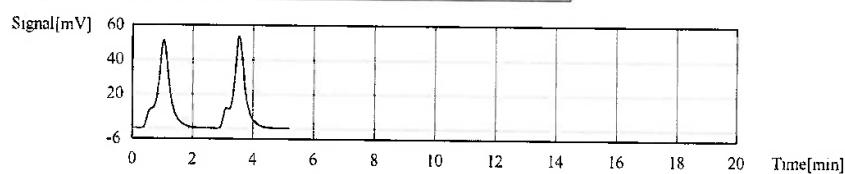
Anal.: NPOC

# TOC-Control L Report

Chem36  
d230331wl toc tlk

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	146.2	24.88mg/L	100µL	1.000	d230201wl 2023_02_01_11_45_27.cal		4/3/2023 11:39:24 AM
2	145.5	24.76mg/L	100µL	1.000	d230201wl 2023_02_01_11_45_27.cal		4/3/2023 11:42:18 AM

Mean Conc. 24.82mg/L  
CV Conc 0.34%



## Sample

Sample Name: CCB  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

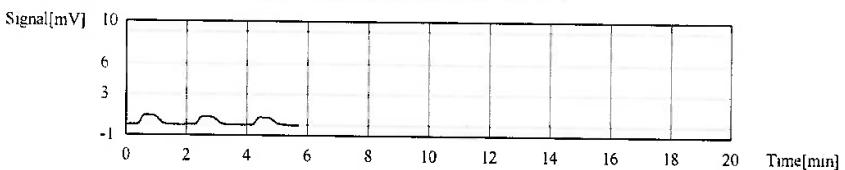
Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.4846mg/L

1. Det

Anal.: NPOC

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	3.499	0.5955mg/L	100µL	1.000	E	d230201wl 2023_02_01_11_45_27.cal	4/3/2023 11:58:32 AM
2	2.928	0.4983mg/L	100µL	1.000	d230201wl 2023_02_01_11_45_27.cal		4/3/2023 12:00:36 PM
3	2.767	0.4709mg/L	100µL	1.000	d230201wl 2023_02_01_11_45_27.cal		4/3/2023 12:02:41 PM

Mean Conc. 0.4846mg/L  
CV Conc 4.00%



## Sample

Sample Name: SPARGERCHK  
Sample ID:  
Origin: TOCAQ met  
Status: Completed  
Chk. Result

Type	Anal.	Manual Dilution	Result
Unknown	NPOC	1.000	NPOC 0.3626mg/L

1. Det

Anal. NPOC

# TOC-Control L Report

No.	Area	Conc.	Inj. Vol.	Aut. Dil.	Ex.	Cal. Curve	Date / Time
1	2.087	0.3552mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27.cal	4/3/2023 12:09:42 PM
2	2.174	0.3700mg/L	100µL	1.000		d230201wl 2023_02_01_11_45_27.cal	4/3/2023 12:11:47 PM

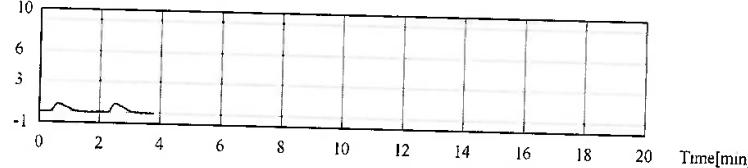
Mean Conc.

0.3626mg/L

CV Conc

2.89%

Signal[mV]

9.6  
9

30/30

4/3/2023 12:14:47 PM