

**TECHNICAL MEMORANDUM NO. 3A
FOR
REMEDIAL INVESTIGATION/FEASIBILITY STUDY
CITY DISPOSAL CORPORATION LANDFILL
(DUNN LANDFILL)**

(PELA Reference No. 495207)

Appendix G

**Results of Analyses
Soil and Ground-Water Samples**

Volume 2 of 16

Technical Reports,
Special Sampling - September 1990

SAMPLES: CA4753-CA4756, CA4757-4762, CA4763

April 1991



OHM Corporation

ETC Environmental Testing
and Certification Corp.

Technical Report

for

WASTE MANAGEMENT, INC.

| <i>Chain of Custody Data Required for ETC Data Management Summary Reports</i> | | | | | | |
|---|------------------------|-----------------|---------------------|-------------|-------------|----------------------|
| <i>ETC Sample No.</i> | <i>Company</i> | <i>Facility</i> | <i>Sample Point</i> | <i>Date</i> | <i>Time</i> | <i>Elapsed Hours</i> |
| CA4757-CA4762 | WASTE MANAGEMENT, INC. | 405 | | | | |

Richard P. Albert
Vice President, General Manager

This Technical Report is an INSITESM service generated by LODESTARSM Data Management Software.

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| Semivolatile Data | NA |
| QC Summary Data | NA |
| Sample Data | NA |
| Standard Data | NA |
| Raw QC Data | NA |
| Pesticide/PCB Data | NA |
| QC Summary Data | NA |
| Sample Data | NA |
| Standard Data | NA |
| Raw QC Data | NA |
| Chain of Custody | 205 |

ETC

CASE NARRATIVE

SDG NARRATIVE

This technical report submitted by ETC Corporation contains the analytical results and required deliverables for WASTE MANAGEMENT, INC. Site 405 samples as identified below:


| <u>ETC ID</u> | <u>CLIENT ID</u> | <u>ETC ID</u> | <u>CLIENT ID</u> |
|---------------|------------------|---------------|------------------|
| CA4757 | 2GWPZ10 | CA4760 | 2GWPZ115 |
| CA4758 | 2GWPZ5 | CA4761 | 2GWPZ11I |
| CA4759 | 2GWPZ9 | CA4762 | 2GWPZ11D |

During the preparation and analysis of these samples, the following was observed:

VOLATILES:

(OV70357): Surrogate recoveries outside of control limits for samples CA4760, CA4761 and CA4762 may be attributed to sample matrix. The results have been confirmed by a replicate analysis.

Release of the data contained in this hardcopy data package has been authorized by the following signature.



Gregory G. Morrison
Laboratory Manager

10-22-98
Date

ETC

METHODOLOGY

INTRODUCTION

Environmental Testing and Certification Corp.
284 Raritan Center Parkway, CN 7808
Edison, New Jersey 08818-7808
(201) 225-6700

This report contains analytical results in tabular form for your sample. It includes comprehensive data for each analytical process. Associated quality control data is also presented, including QC batch results for the laboratory blank, spiked blank, matrix spike and a replicated sample spike, as well as results from surrogate compound analyses. Quality control data for instrument performance is also included. Other appendices may include data system printouts and chain of custody records.

Analytical Methodology

ETC utilizes a wide variety of methodologies and EPA approved procedures which are listed below. Analytical results and Quality Assurance protocols are based upon the appropriate guideline dependent upon the level of deliverable requested.

- "Methods of Organic Chemical Analysis of Municipal and Industrial Wastewater", Federal Register Vol. 49, No. 209, October 26, 1984;
- "Test Methods for Evaluating Solid Waste", SW-846 Third Edition, September 1986, USEPA;
- "Standard Methods for the Examination of Water and Wastewater" 1985, 14th, 15th and 16th Edition;
- "Methods for Chemical Analysis of Water and Wastes" March 1983, EMSL, EPA 600 4-79-020;
- Organic Analysis: Multi-media, Multi-Concentration-IFB-CLP, SOW February 1988;
- Inorganic Analysis: Multi-media, Multi-Concentration IFB-CLP, SOW July 1987;
- Dioxin Analysis: Soil/Sediment/Water Matrix; Multi-Concentration, Selected Ion Monitoring with Jar Extraction Procedure; IFB WA86-K357.
- "Methods for the Determination of Organic Compounds in Drinking Water" December 1988, EPA-600/4-88/039;
- "Handbook for Analytical Quality Control in Water and Wastewater Laboratories", EPA-600/4-79-019, March 1979;
- "National Enforcement Investigation Center Policies and Procedures Manual, EPA-330/9/78/001-R, Revised May 1986.

ETC

ETC Network Laboratories are certified to perform analytical analyses upon samples collected throughout the United States. It may be required that this environmental data be reported with reference to a certified laboratory. The Laboratory Identification Numbers for ETC Corp., Edison NJ Laboratory are summarized below for your information and reference. Please contact your Program Manager should you require certification verification for an ETC Network Laboratory.

Certification Summary, January 1990

| <u>State Agency</u> | <u>Certification Number</u> |
|---|------------------------------------|
| Alabama Department of Environmental Management | 40280 |
| Arizona Department of Environmental Quality | 322 |
| California Department of Health Services | 162 |
| Connecticut Department of Health Services | 0511 |
| Florida Department of Health and Rehabilitative Services | E87074 & 87262 |
| Illinois Environmental Protection Agency | 100224 |
| Kansas Department of Health and Environment | E-148 & E1122 |
| New Hampshire Department of Environmental Services | 198948-B |
| New Jersey Department of Environmental Protection | 12257 |
| New York Department of Public Health | 10586 |
| Oklahoma Water Resources Board | 8703 |
| Pennsylvania Department of Environmental Regulation | 68-323 |
| South Carolina Department of Health and Environmental Control | 94002 |
| Tennessee Department of Health and Environmental Laboratory Services | 02915 |
| Utah Department of Health | E-91 |
| Virginia Department of General Services | 00113 |
| Wisconsin Department of Natural Resources | 999464070 |

ETC

VOLATILE DATA

ETC

QC SUMMARY

2A
WATER VOLATILE SURROGATE RECOVERY

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

| | EPA SAMPLE NO. | S1 (TOL)# | S2 (BFB)# | S3 (DCE)# | OTHER | TOT OUT |
|----|-------------------|--------------|--------------|--------------|-------|------------|
| 01 | UCLK01 | 101 | 106 | 95 | | 0 |
| 02 | A4753 | 102 | 107 | 99 | | 0 |
| 03 | A4755 | 101 | 107 | 101 | | 0 |
| 04 | A4763 | 100 | 107 | 103 | | 0 |
| 05 | A4756 | 100 | 107 | 102 | | 0 |
| 06 | A4758 | 100 | 110 | 103 | | 0 |
| 07 | UCLK2 | 99 | 103 | 108 | | 0 |
| 08 | A4754 | 103 | 105 | 112 | | 0 |
| 09 | A4757 | 101 | 106 | 114 | | 0 |
| 10 | A4759 | 101 | 104 | 113 | | 0 |
| 11 | A4760 | 111 | 95 | 111 | | 1 |
| 12 | UCLK4 | 101 | 111 | 105 | | 0 |
| 13 | A4762 | 104 | 127 * | 129 * | | 2 |
| 14 | A4761 | 101 | 126 * | 130 * | | 2 |
| 15 | UCLK5 | 98 | 112 | 97 | | 0 |
| 16 | A4755MS | 105 | 112 | 109 | | 0 |
| 17 | A4755MSD | 98 | 112 | 106 | | 0 |
| 18 | | | | | | |
| 19 | A4762 DL | 100 | 121 * | 119 * | | 2 |
| 20 | A4761 DL | 100 | 121 * | 122 * | | 2 |
| 21 | A4760 | 98 | 118 * | 116 * | | 2 |
| 22 | A4760 DL | 104 | 128 * | 128 * | | 2 |
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CA 10-16-90

CA 10-16-90

QC LIMITS

S1 (TOL) = Toluene-d8 (88-110)
 S2 (BFB) = Bromofluorobenzene (86-115)
 S3 (DCE) = 1,2-Dichloroethane-d4 (76-114)

Column to be used to flag recovery values
 * Values outside QC limits
 D Surrogates diluted out

3A
WATER VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix Spike - EPA Sample No.: A4755

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC # | QC LIMITS REC. |
|--------------------|--------------------------|-----------------------------------|-------------------------------|------------------|----------------------|
| 1,1-Dichloroethene | 50.000 | 0.000 | 54.109 | 108 | 161-145 |
| Trichloroethene | 50.000 | 0.000 | 48.018 | 96 | 171-120 |
| Benzene | 50.000 | 0.000 | 41.577 | 83 | 176-127 |
| Toluene | 50.000 | 0.000 | 49.472 | 99 | 176-125 |
| Chlorobenzene | 50.000 | 0.000 | 47.831 | 96 | 175-130 |

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC # | % RPD # | QC LIMITS RPD REC. |
|--------------------|--------------------------|--------------------------------|-------------------|------------|-------------------------|
| 1,1-Dichloroethene | 50.000 | 51.446 | 103 | 5 | 14 161-145 |
| Trichloroethene | 50.000 | 49.846 | 100 | 4 | 14 171-120 |
| Benzene | 50.000 | 42.915 | 86 | 3 | 11 176-127 |
| Toluene | 50.000 | 45.757 | 92 | 8 | 13 176-125 |
| Chlorobenzene | 50.000 | 49.015 | 98 | 2 | 13 175-130 |

Column to be used to flag recovery and RPD values with an asterisk

Values outside of limits

RPD: 0 out of 5 outside limits

Matrix Spike Recovery: 0 out of 10 outside limits

Comments:

VOLATILE METHOD ^{4A} BLANK SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0918 Lab Sample ID: VBLK01
 Date Analyzed 09/26/90 Time Analyzed: 1725
 Matrix: (soil/water) WATER Level: (low/med) LOW
 Instrument ID: GC/MS C

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | TIME ANALYZED |
|----|-------------------|------------------|----------------|------------------|
| 01 | A4753 | CA4753U | >C0920 | 1904 |
| 02 | A4755 | CA4755U | >C0922 | 2043 |
| 03 | A4763 | CA4763U | >C0923 | 2132 |
| 04 | A4756 | CA4756U | >C0924 | 2221 |
| 05 | A4758 | CA4758U | >C0925 | 2310 |
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Comments: _____

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0929 Lab Sample ID: VBLK2
 Date Analyzed 09/27/90 Time Analyzed: 1814
 Matrix: (soil/water) WATER Level:(low/med) LOW
 Instrument ID: GC/MS C

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | TIME ANALYZED |
|----|-------------------|------------------|----------------|------------------|
| 01 | A4754 | CA4754U2 | >C0931 | 1952 |
| 02 | A4757 | CA4757U2 | >C0932 | 2041 |
| 03 | A4759 | CA4759U2 | >C0933 | 2130 |
| 04 | | | | |
| 05 | A4760 | CA4760U2 | >C0934 | 2220 |
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Comments: _____

Page 1 of 1

5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name:ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0910 BFB Injection Date:09/26/90
 Instrument ID: GC/MS C BFB Injection Time:0854
 Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap)PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 25.0 |
| 75 | 30.0 - 60.0% of mass 95 | 58.3 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 7.0 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 81.4 |
| 175 | 5.0 - 9.0% of mass 174 | 6.5 (8.0)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 78.9 (96.9)1 |
| 177 | 5.0 - 9.0% of mass 176 | 5.9 (7.4)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------|---------------|-------------|---------------|---------------|
| 01 | VOA | QC70357US | >C0911 | 09/26/90 | 1124 |
| 02 | USTD150 | QC70357US | >C0912 | 09/26/90 | 1213 |
| 03 | USTD100 | QC70357US | >C0913 | 09/26/90 | 1303 |
| 04 | USTD50 | QC70357US | >C0914 | 09/26/90 | 1352 |
| 05 | USTD20 | QC70357US | >C0915 | 09/26/90 | 1441 |
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5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0917 BFB Injection Date: 09/26/90
 Instrument ID: GC/MS C BFB Injection Time: 1625
 Matrix: (soil/water) WATER Level: (low/med) LOW Column: (pack/cap) PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 23.3 |
| 75 | 30.0 - 60.0% of mass 95 | 59.4 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 6.9 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 84.0 |
| 175 | 5.0 - 9.0% of mass 174 | 7.0 (8.4)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 81.3 (96.7)1 |
| 177 | 5.0 - 9.0% of mass 176 | 6.0 (7.4)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------|---------------|-------------|---------------|---------------|
| 01 | UCLK01 | QC70357U | >C0918 | 09/26/90 | 1725 |
| 02 | A4753 | CA4753U | >C0920 | 09/26/90 | 1904 |
| 03 | A4755 | CA4755U | >C0922 | 09/26/90 | 2043 |
| 04 | A4763 | CA4763U | >C0923 | 09/26/90 | 2132 |
| 05 | A4756 | CA4756U | >C0924 | 09/26/90 | 2221 |
| 06 | A4758 | CA4758U | >C0925 | 09/26/90 | 2310 |
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5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0926 BFB Injection Date: 09/27/90
 Instrument ID: GC/MS C BFB Injection Time: 0858
 Matrix: (soil/water) WATER Level: (low/med) LOW Column: (pack/cap) PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 27.2 |
| 75 | 30.0 - 60.0% of mass 95 | 59.1 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 7.1 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 77.4 |
| 175 | 5.0 - 9.0% of mass 174 | 6.0 (7.7)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 73.7 (95.2)1 |
| 177 | 5.0 - 9.0% of mass 176 | 5.9 (8.0)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|------------------|----------------------|----------------------|---------------------|-----------------|
| 01 | U5TD50 | QC70357US | >C0927 | 09/27/90 | 1656 |
| 02 | UBLK2 | QC70357U2 | >C0927 | 09/27/90 | 1814 |
| 03 | A4754 | CA4754U2 | >C0931 | 09/27/90 | 1952 |
| 04 | | | | | |
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10-19-90

5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name:ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Lab File ID: >C0928

BFB Injection Date:09/27/90

Instrument ID: GC/MS C

BFB Injection Time:1741

Matrix:(soil/water) WATER

Level:(low/med) LOW

Column:(pack/cap)PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 23.5 |
| 75 | 30.0 - 60.0% of mass 95 | 58.6 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 6.7 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 87.9 |
| 175 | 5.0 - 9.0% of mass 174 | 7.8 (8.9)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 87.8 (99.8)1 |
| 177 | 5.0 - 9.0% of mass 176 | 6.8 (7.7)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|-------------------|------------------|----------------|------------------|------------------|
| 01 | VBLK2 | QC70357U2 | >C0929 | 09/27/90 | 1814 |
| 02 | A4754 | CA4754U2 | >C0931 | 09/27/90 | 1952 |
| 03 | A4757 | CA4757U2 | >C0932 | 09/27/90 | 2041 |
| 04 | A4759 | CA4759U2 | >C0933 | 09/27/90 | 2130 |
| 05 | A4760 | CA4760U2 | >C0934 | 09/27/90 | 2220 |
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5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name:ETCND

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Lab File ID: >C0957

BFB Injection Date:10/02/90

Instrument ID: GC/MS C

BFB Injection Time:1148

Matrix:(soil/water) WATER

Level:(low/med) LOW

Column:(pack/cap)PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 20.2 |
| 75 | 30.0 - 60.0% of mass 95 | 45.2 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 6.1 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 92.0 |
| 175 | 5.0 - 9.0% of mass 174 | 8.0 (8.7)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 89.6 (97.3)1 |
| 177 | 5.0 - 9.0% of mass 176 | 6.5 (7.2)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|-------------------|------------------|----------------|------------------|------------------|
| 01 | USTD200 | QC70357US | >C0958 | 10/02/90 | 1215 |
| 02 | USTD150 | QC70357US | >C0959 | 10/02/90 | 1304 |
| 03 | USTD100 | QC70357US | >C0960 | 10/02/90 | 1353 |
| 04 | USTD50 | QC70357US | >C0961 | 10/02/90 | 1443 |
| 05 | USTD20 | QC70357US | >C0962 | 10/02/90 | 1532 |
| 06 | VBLK4 | QC70357U4 | >C0963 | 10/02/90 | 1621 |
| 07 | A4762 | CA4762U4 | >C0969 | 10/02/90 | 2137 |
| 08 | A4761 | CA4761U4 | >C0970 | 10/02/90 | 2225 |
| 09 | | | | | |
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5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name:ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0971 BFB Injection Date:10/03/90
 Instrument ID: GC/MS C BFB Injection Time:0930
 Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap)PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 18.1 |
| 75 | 30.0 - 60.0% of mass 95 | 49.8 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 8.0 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 94.9 |
| 175 | 5.0 - 9.0% of mass 174 | 7.9 (8.3)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 93.5 (98.5)1 |
| 177 | 5.0 - 9.0% of mass 176 | 7.6 (8.1)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------|---------------|-------------|---------------|---------------|
| 01 | USTD50 | QC70357US | >C0972 | 10/03/90 | 1001 |
| 02 | USTD200 | QC70357US | >C0975 | 10/03/90 | 1305 |
| 03 | USTD150 | QC70357US | >C0976 | 10/03/90 | 1355 |
| 04 | USTD100 | QC70357US | >C0977 | 10/03/90 | 1445 |
| 05 | USTD20 | QC70357US | >C0978 | 10/03/90 | 1535 |
| 06 | UCLK5 | QC70357U5 | >C0979 | 10/03/90 | 1625 |
| 07 | A4755MS | CA4755US | >C0981 | 10/03/90 | 1819 |
| 08 | A4755MSD | CA4755UR | >C0982 | 10/03/90 | 1907 |
| 09 | | | | | |
| 10 | | | | | |
| 11 | | | | | |
| 12 | | | | | |
| 13 | | | | | |
| 14 | | | | | |
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| 22 | | | | | |

ETC

SAMPLE DATA

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4757

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4757U2

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0932

Level: (low/med) LOW

Date Received: 09/25/90

% Moisture: not dec.

Date Analyzed: 09/27/90

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | | Q |
|------------|----------------------------|----------------------|------|----|
| | | (ug/L or ug/Kg) | UG/L | |
| 74-87-3 | Chloromethane | | 10 | 1U |
| 74-83-9 | Bromomethane | | 10 | 1U |
| 75-01-4 | Vinyl Chloride | | 10 | 1U |
| 75-00-3 | Chloroethane | | 10 | 1U |
| 75-09-2 | Methylene Chloride | | 15 | 1U |
| 67-64-1 | Acetone | | 10 | 1U |
| 75-15-0 | Carbon Disulfide | | 15 | 1U |
| 75-35-4 | 1,1-Dichloroethane | | 15 | 1U |
| 75-34-3 | 1,1-Dichloroethane | | 15 | 1U |
| 540-59-0 | 1,2-Dichloroethene (total) | | 15 | 1U |
| 67-66-3 | Chloroform | | 15 | 1U |
| 107-06-2 | 1,2-Dichloroethane | | 15 | 1U |
| 78-93-3 | 2-Butanone | | 10 | 1U |
| 71-55-6 | 1,1,1-Trichloroethane | | 15 | 1U |
| 56-23-5 | Carbon Tetrachloride | | 15 | 1U |
| 108-05-4 | Vinyl Acetate | | 10 | 1U |
| 75-27-4 | Bromodichloromethane | | 15 | 1U |
| 78-87-5 | 1,2-Dichloropropane | | 15 | 1U |
| 10061-01-5 | cis-1,3-Dichloropropene | | 15 | 1U |
| 79-01-6 | Trichloroethene | | 15 | 1U |
| 124-48-1 | Dibromochloromethane | | 15 | 1U |
| 79-00-5 | 1,1,2-Trichloroethane | | 15 | 1U |
| 71-43-2 | Benzene | | 15 | 1U |
| 10061-02-6 | trans-1,3-Dichloropropene | | 15 | 1U |
| 75-25-2 | Bromoform | | 15 | 1U |
| 108-10-1 | 4-Methyl-2-Pentanone | | 10 | 1U |
| 591-78-6 | 2-Hexanone | | 10 | 1U |
| 127-18-4 | Tetrachloroethene | | 15 | 1U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 15 | 1U |
| 108-88-3 | Toluene | | 15 | 1U |
| 108-90-7 | Chlorobenzene | | 15 | 1U |
| 100-41-4 | Ethylbenzene | | 15 | 1U |
| 100-42-5 | Styrene | | 15 | 1U |
| 1330-20-7 | Xylene (total) | | 15 | 1U |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETCNS

Contract: _____

A4757

Lab Code: _____ Case No.: _____

SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER

Lab Sample ID: CA4757

Sample wt/vol: 5 (g/mL) HL

Lab File ID: >C0932

Level: (low/med) LOW

Date Received: 09/25/90

Moisture: not dec. _____

Date Analyzed: 09/27/90

Column: (pack/cap) PACK

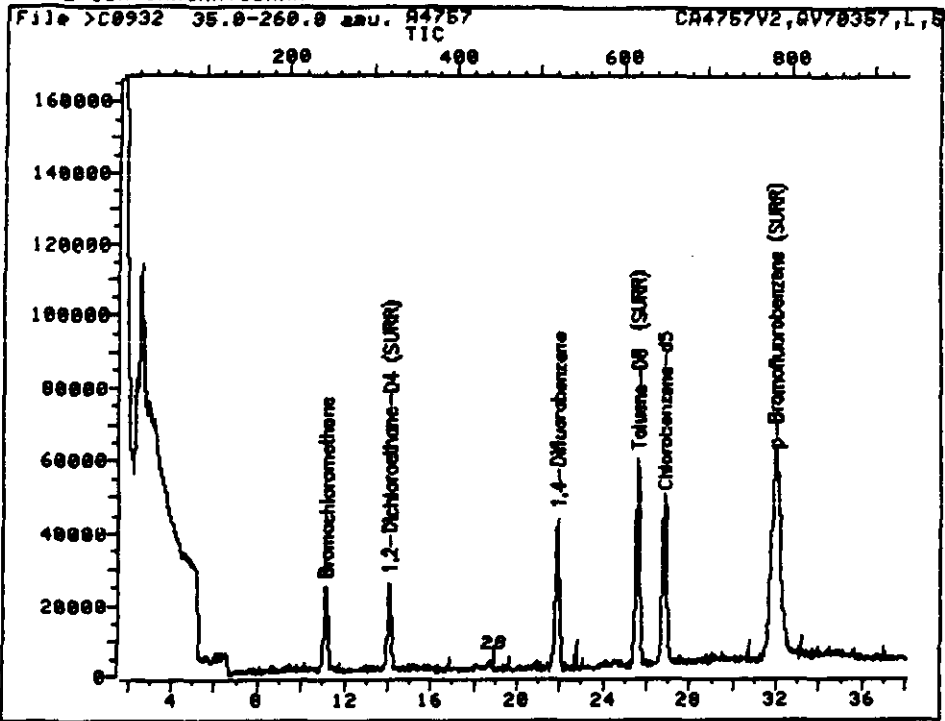
Dilution Factor: 1

CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

Number TICs found: 0

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
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TOTAL ION CHROMATOGRAM



Data File: >C0932
Name: A4757
Misc: CA4757V2,QU70357,L,5,,

Quant Output File: ^C0932::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900927 17:41

Operator ID: MGRMS
Quant Time: 900927 21:19
Injected at: 900927 20:41

QUANT REPORT

Page 1

Operator ID: MGRMS
 Output File: ^C0932::AQ
 Data File: >C0932::U4
 Name: A4757
 Misc: CA4757U2,QU70357,L,5,,

Quant Rev: 7 Quant Time: 900927 21:19
 Injected at: 900927 20:41
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900927 17:41

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|---------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.10 | 244 | 39142 | 250.00 | NG | 96 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.04 | 320 | 144677M | 285.97 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.84 | 521 | 183849 | 250.00 | NG | 98 |
| 28) Trichloroethylene | 18.74 | 441 | 3412 | 18.49 | NG | 88 |
| 36) *Chlorobenzene-d5 | 26.78 | 648 | 173277 | 250.00 | NG | 78 |
| 41) Toluene-D8 (SURR) | 25.56 | 617 | 255796 | 252.76 | NG | 88 |
| 45) p-Bromofluorobenzene (SURR) | 31.90 | 780 | 206793 | 264.21 | NG | 86 |

* Compound is ISTD

CA
 10-16-90.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CA4758

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4758U

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0925

Level: (low/med) LOW

Date Received: ~~10/17/90~~ 09/25/90 (CR)

% Moisture: not dec.

Date Analyzed: 09/26/90 10-6-90

Column: (pack/cap) PACK

Dilution Factor: 1

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

| | | | |
|------------|----------------------------|-----|----|
| 74-87-3 | Chloromethane | 10 | 10 |
| 74-83-9 | Bromomethane | 10 | 10 |
| 75-01-4 | Vinyl Chloride | 10 | 10 |
| 75-00-3 | Chloroethane | 10 | 10 |
| 75-09-2 | Methylene Chloride | 15 | 10 |
| 67-64-1 | Acetone | 10 | 10 |
| 75-15-0 | Carbon Disulfide | 15 | 10 |
| 75-35-4 | 1,1-Dichloroethane | 15 | 10 |
| 75-34-3 | 1,1-Dichloroethane | 15 | 10 |
| 540-59-0 | 1,2-Dichloroethene (total) | 15 | 10 |
| 67-66-3 | Chloroform | 15 | 10 |
| 107-06-2 | 1,2-Dichloroethane | 15 | 10 |
| 78-93-3 | 2-Butanone | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | 15 | 10 |
| 56-23-5 | Carbon Tetrachloride | 15 | 10 |
| 108-05-4 | Vinyl Acetate | 10 | 10 |
| 75-27-4 | Bromodichloromethane | 15 | 10 |
| 78-87-5 | 1,2-Dichloropropane | 15 | 10 |
| 10061-01-5 | cis-1,3-Dichloropropene | 15 | 10 |
| 79-01-6 | Trichloroethene | 156 | 10 |
| 124-48-1 | Dibromochloromethane | 15 | 10 |
| 79-00-5 | 1,1,2-Trichloroethane | 15 | 10 |
| 71-43-2 | Benzene | 15 | 10 |
| 10061-02-6 | trans-1,3-Dichloropropene | 15 | 10 |
| 75-25-2 | Bromoform | 15 | 10 |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 10 |
| 591-78-6 | 2-Hexanone | 10 | 10 |
| 127-18-4 | Tetrachloroethene | 12 | 10 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 15 | 10 |
| 108-88-3 | Toluene | 15 | 10 |
| 108-90-7 | Chlorobenzene | 15 | 10 |
| 100-41-4 | Ethylbenzene | 15 | 10 |
| 100-42-5 | Styrene | 15 | 10 |
| 1330-20-7 | Xylene (total) | 15 | 10 |

18
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

A4758

Lab Name: ETCNS Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: CA4758

Sample wt/vol: 5 (g/mL) ML Lab File ID: 76925

Level: (low/med) LOW Date Received: 09/25/90

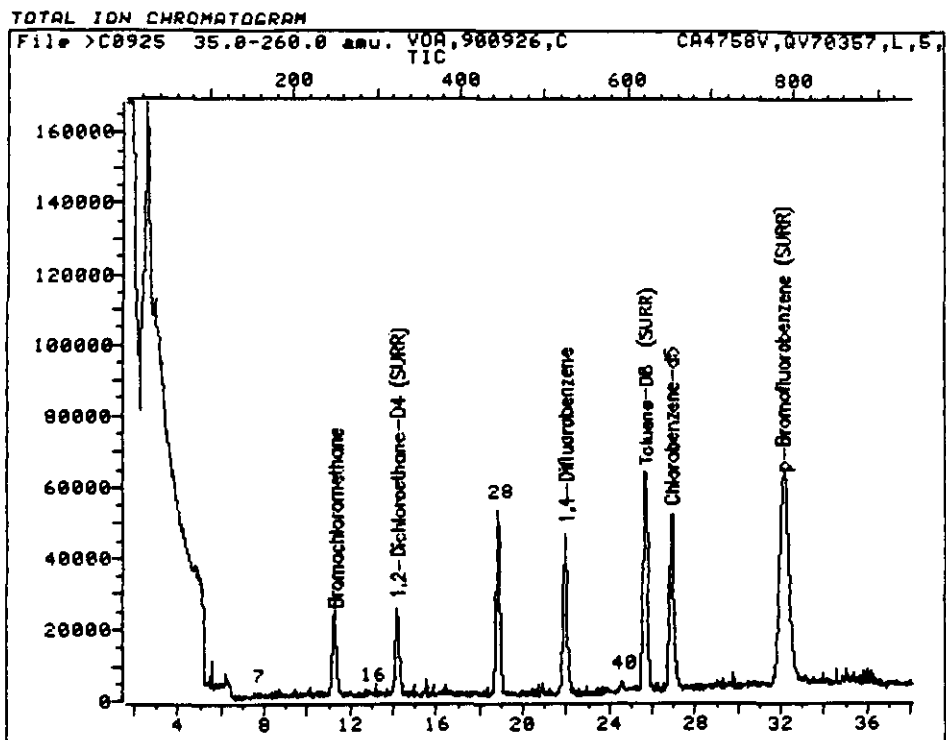
% Moisture: not des. _____ Date Analyzed: 09/26/90

Column: (pack/cap) PACK Dilution Factor: 1

CONCENTRATION UNITS:
 (ug/L or ug/kg) ug/L

Number TICs found: 0

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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Data File: >C0925::U0
Name: VOA,900926,C
Misc: CA4758V,QV70357,L,5,,

Quant Output File: ^C0925::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900926 16:16

Operator ID: JA8781
Quant Time: 900926 23:48
Injected at: 900926 23:10

QUANT REPORT

Operator ID: JA8781
 Output File: ^C0925::AQ
 Data File: >C0925::U0
 Name: VOA,900926,C
 Misc: CA4758U,QV70357,L,5,,

Quant Rev: 7 Quant Time: 900926 23:48
 Injected at: 900926 23:10
 Dilution Factor: 1.00000

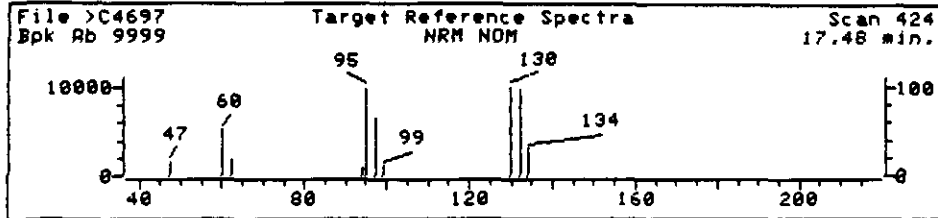
ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 16:16

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|-------|----|
| 1) *Bromochloromethane | 11.30 | 250 | 41198 | 250.00 | NG | 98 |
| 7) Methylene chloride | 7.65 | 156 | 2545 | 9.33 | NG | 89 |
| 16) 1,2-Trans-dichloroethylene | 12.93 | 292 | 2699 | 11.12 | NG | 89 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.21 | 325 | 148614 | 258.41 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.96 | 525 | 192371 | 250.00 | NG | 97 |
| 28) Trichloroethylene | 18.86 | 445 | 93350 | 277.87 | NG | 85 |
| 36) *Chlorobenzene-d5 | 26.90 | 652 | 182527 | 250.00 | NG | 74 |
| 40) Tetrachloroethylene | 24.56 | 592 | 3518 | 10.26 | NG | 87 |
| 41) Toluene-D8 (SURR) | 25.69 | 621 | 268674 | 249.93 | NG | 94 |
| 45) p-Bromofluorobenzene (SURR) | 32.10 | 786 | 218921 | 275.60 | NG | 84 |

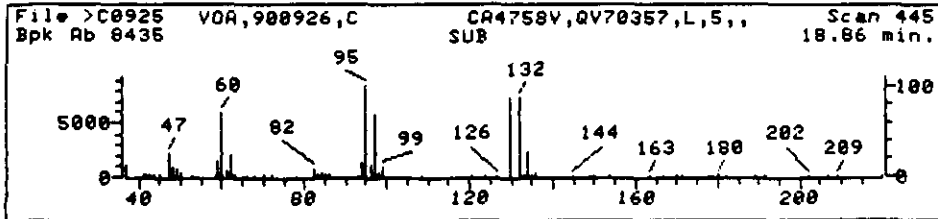
* Compound is ISTD

CA
10-17-90.

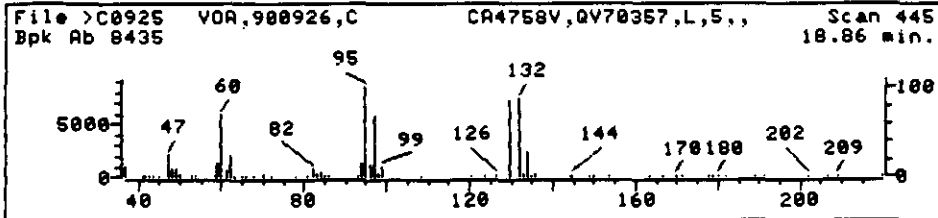
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



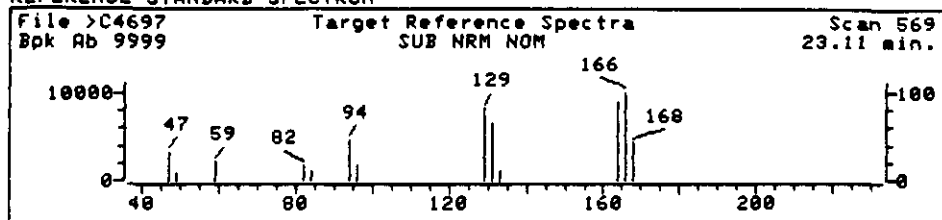
Data File: >C0925::U0
 Name: VOA,900926,C
 Misc: CA4758V,QV70357,L,5,,
 Quant Time: 900926 23:48
 Injected at: 900926 23:10

Quant Output File: ^C0925::AQ

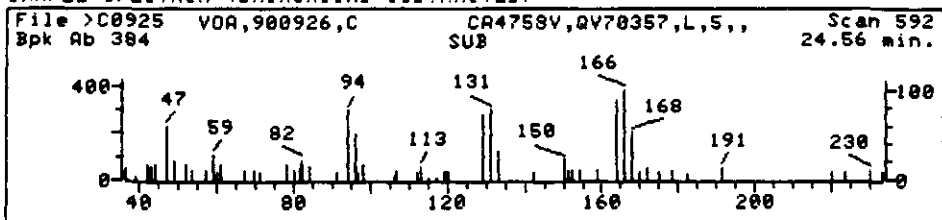
Quant ID File: IC1171::US
 Last Calibration: 900926 16:16

Compound No: 28
 Compound Name: Trichloroethylene
 Scan Number: 445
 Retention Time: 18.86 min.
 Quant Ion: 130.0
 Area: 93350
 Concentration: 277.87 NG
 q-value: 85

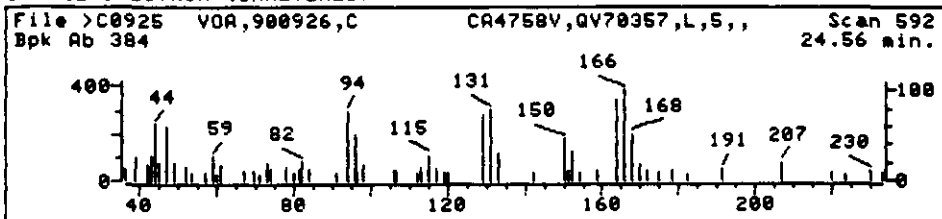
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >C0925::U0
Name: VOA,900926,C
Misc: CA4758V,QU70357,L,5,,
Quant Time: 900926 23:48
Injected at: 900926 23:10

Quant Output File: ^C0925::AQ
Quant ID File: IC1171::US
Last Calibration: 900926 16:16

Compound No: 40
Compound Name: Tetrachloroethylene
Scan Number: 592
Retention Time: 24.56 min.
Quant Ion: 164.0
Area: 3518
Concentration: 10.26 NG
q-value: 87

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4759

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4759U2

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0933

Level: (low/med) LOW

Date Received: 09/26/90

% Moisture: not dec.

Date Analyzed: 09/27/90

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | | Q |
|------------|----------------------------|----------------------|------|---|
| | | (ug/L or ug/Kg) | UG/L | |
| 74-87-3 | Chloromethane | 10 | | U |
| 74-83-9 | Bromomethane | 10 | | U |
| 75-01-4 | Vinyl Chloride | 10 | | U |
| 75-00-3 | Chloroethane | 10 | | U |
| 75-09-2 | Methylene Chloride | 5 | | U |
| 67-64-1 | Acetone | 10 | | U |
| 75-15-0 | Carbon Disulfide | 5 | | U |
| 75-35-4 | 1,1-Dichloroethene | 5 | | U |
| 75-34-3 | 1,1-Dichloroethane | 5 | | U |
| 540-59-0 | 1,2-Dichloroethene (total) | 14 | | |
| 67-66-3 | Chloroform | 5 | | U |
| 107-06-2 | 1,2-Dichloroethane | 5 | | U |
| 78-93-3 | 2-Butanone | 10 | | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | | U |
| 56-23-5 | Carbon Tetrachloride | 5 | | U |
| 108-05-4 | Vinyl Acetate | 10 | | U |
| 75-27-4 | Bromodichloromethane | 5 | | U |
| 78-87-5 | 1,2-Dichloropropane | 5 | | U |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | | U |
| 79-01-6 | Trichloroethene | 350 | | E |
| 124-48-1 | Dibromochloromethane | 5 | | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | | U |
| 71-43-2 | Benzene | 5 | | U |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | | U |
| 75-25-2 | Bromoform | 5 | | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | | U |
| 591-78-6 | 2-Hexanone | 10 | | U |
| 127-18-4 | Tetrachloroethene | 6 | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | | U |
| 108-88-3 | Toluene | 5 | | U |
| 108-90-7 | Chlorobenzene | 5 | | U |
| 100-41-4 | Ethylbenzene | 5 | | U |
| 100-42-5 | Styrene | 5 | | U |
| 1330-20-7 | Xylene (total) | 5 | | U |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETCNI Contract: _____

A4759

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: CA4759

Sample wt/vol: 5 (g/mL) HL Lab File ID: >C0933

Level: (low/med) LOW Date Received: 09/26/90

% Moisture: not dec. _____ Date Analyzed: 09/27/90

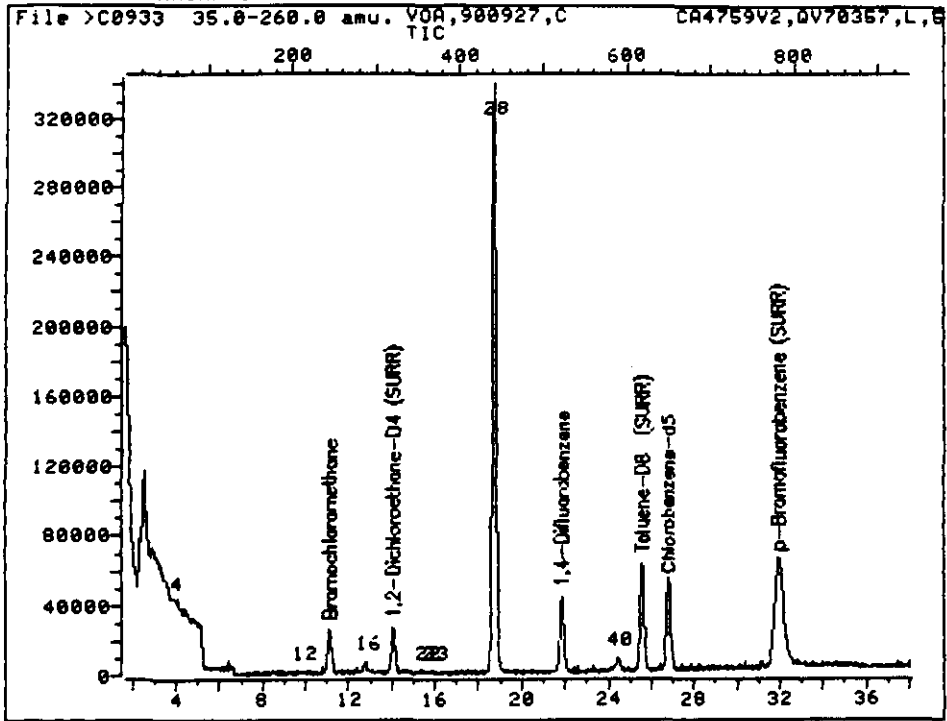
Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
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TOTAL ION CHROMATOGRAM



Data File: >C0933::U4
Name: VOA,900927,C
Misc: CA4759V2,QV70357,L,5,,

Quant Output File: ^C0933::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900927 17:41

Operator ID: MGRMS
Quant Time: 900927 22:09
Injected at: 900927 21:30

QUANT REPORT

Operator ID: MGRMS
 Output File: ^C0933::AQ
 Data File: >C0933::U4
 Name: UOA,900927,C
 Misc: CA4759U2,QU70357,L,5,,

Quant Rev: 7 Quant Time: 900927 22:09
 Injected at: 900927 21:30
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900927 17:41

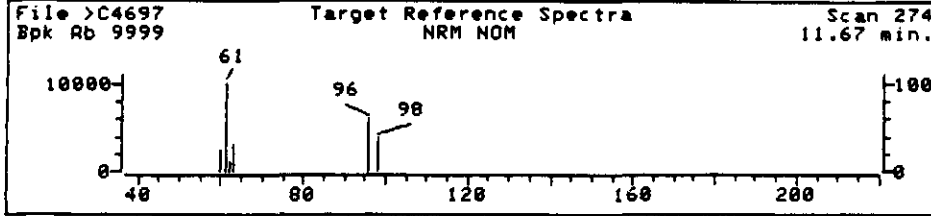
| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|-----------|
| 1) *Bromochloromethane | 11.11 | 244 | 42889 | 250.00 | NG | 92 |
| 4) Dichlorodifluoromethane | 3.97 | 60 | 10773 | 65.63 | NG | 97 |
| 12) Trichlorofluoromethane | 9.87 | 212 | 5307 | 6.91 | NG | 96 |
| 16) 1,2-Trans-dichloroethylene | 12.78 | 287 | 15613 | 68.19 | NG | 94 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.06 | 320 | 156685 | 282.65 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.85 | 521 | 192985 | 250.00 | NG | 96 |
| 22) 1,1,1-Trichloroethane | 15.53 | 358 | 3973 | 4.76 | NG | 90 |
| 23) Carbon tetrachloride | 15.92 | 368 | 2271 | 2.63 | NG | 83 |
| 28) Trichloroethylene | 18.71 | 440 | 599628 | 1756.73 | NG | 1834 * 92 |
| 36) *Chlorobenzene-d5 | 26.79 | 648 | 186569 | 250.00 | NG | 79 |
| 40) Tetrachloroethylene | 24.45 | 588 | 10502 | 28.65 | NG | 94 |
| 41) Toluene-D8 (SURR) | 25.58 | 617 | 273909 | 251.38 | NG | 89 |
| 45) p-Bromofluorobenzene (SURR) | 31.91 | 780 | 219231 | 260.15 | NG | 86 |

* Compound is ISTD

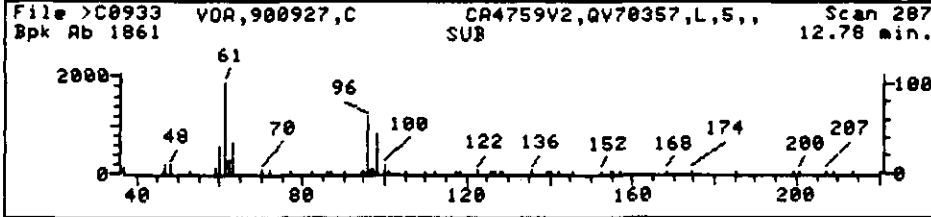
(Handwritten)
 10-16-90.

* value is from dilution run.

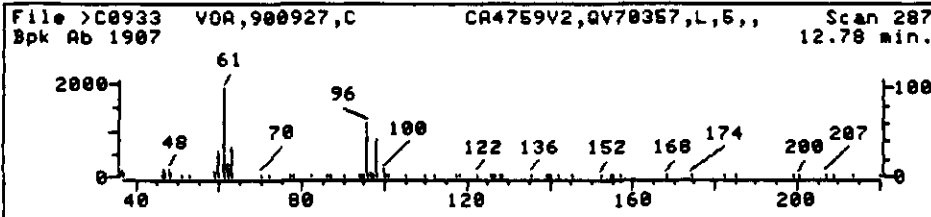
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

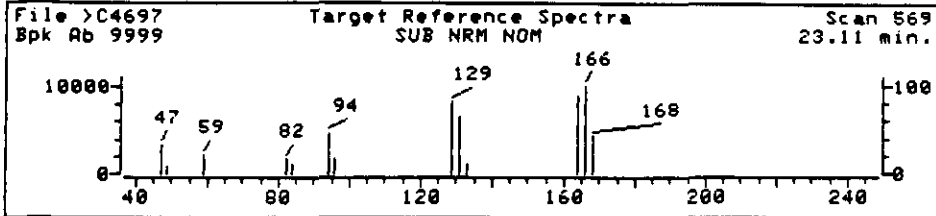


Data File: >C0933::U4
Name: VOA,900927,C
Misc: CA4759V2,QV70357,L,5,,
Quant Time: 900927 22:09
Injected at: 900927 21:30

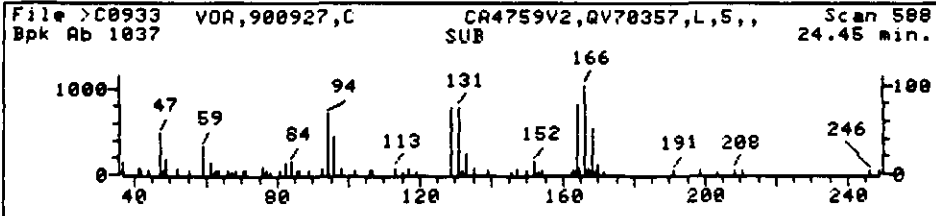
Quant Output File: ^C0933::AQ
Quant ID File: IC1171::US
Last Calibration: 900927 17:41

Compound No: 16
Compound Name: 1,2-Trans-dichloroethylene
Scan Number: 287
Retention Time: 12.78 min.
Quant Ion: 96.0
Area: 15613
Concentration: 68.19 NG
q-value: 94

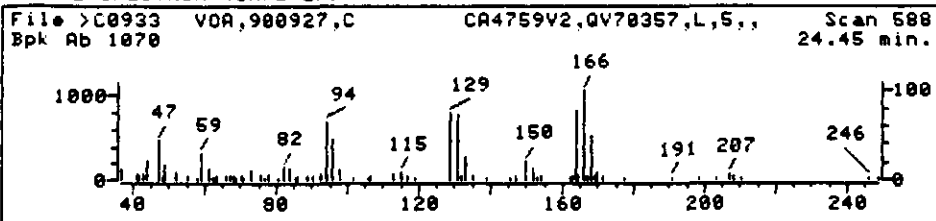
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >C0933::U4
 Name: VOA,900927,C
 Misc: CA4759V2,QV70357,L,5,,
 Quant Time: 900927 22:09
 Injected at: 900927 21:30

Quant Output File: ^C0933::AQ
 Quant ID File: IC1171::US
 Last Calibration: 900927 17:41

Compound No: 40
 Compound Name: Tetrachloroethylene
 Scan Number: 588
 Retention Time: 24.45 min.
 Quant Ion: 164.0
 Area: 10502
 Concentration: 28.65 NG
 q-value: 94

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4759DL

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4759U5

Sample wt/vol: 1.0 (g/mL) ML

Lab File ID: >C0983

Level: (low/med) LOW

Date Received: 09/ /90

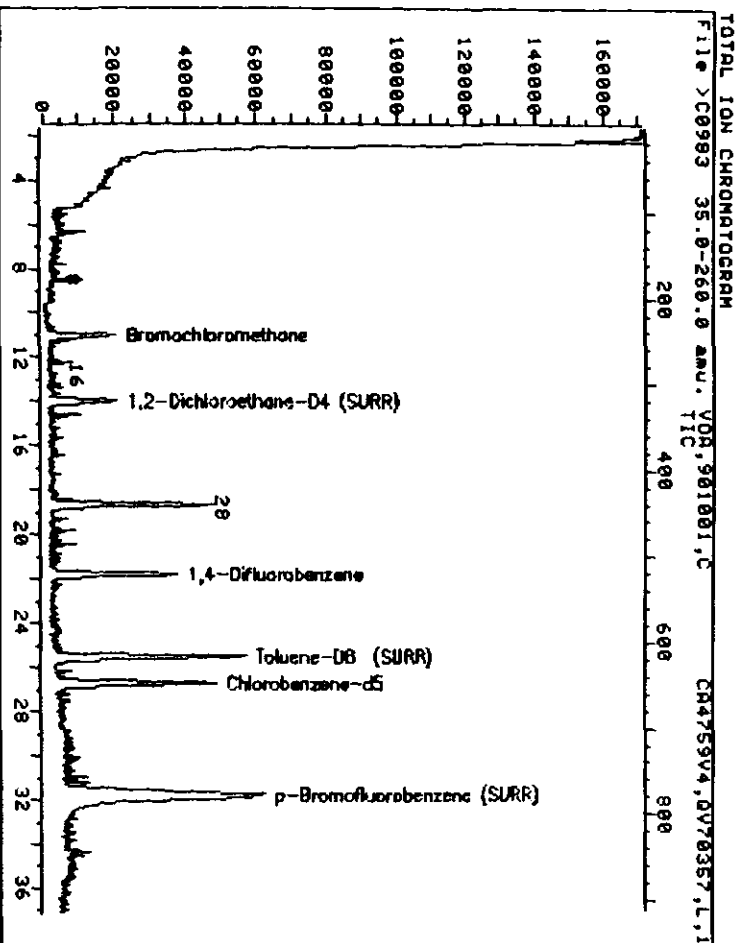
% Moisture: not dec.

Date Analyzed: 10/03/90

Column: (pack/cap) PACK

Dilution Factor: 5

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) | UG/L | Q |
|------------|----------------------------|---|------|-----|
| 74-87-3 | Chloromethane | | 150 | IUD |
| 74-83-9 | Bromomethane | | 150 | IUD |
| 75-01-4 | Vinyl Chloride | | 150 | IUD |
| 75-00-3 | Chloroethane | | 150 | IUD |
| 75-09-2 | Methylene Chloride | | 125 | IUD |
| 67-64-1 | Acetone | | 150 | IUD |
| 75-15-0 | Carbon Disulfide | | 125 | IUD |
| 75-35-4 | 1,1-Dichloroethane | | 125 | IUD |
| 75-34-3 | 1,1-Dichloroethane | | 125 | IUD |
| 540-59-0 | 1,2-Dichloroethane (total) | | 11 | IJD |
| 67-66-3 | Chloroform | | 125 | IUD |
| 107-06-2 | 1,2-Dichloroethane | | 125 | IUD |
| 78-93-3 | 2-Butanone | | 150 | IUD |
| 71-55-6 | 1,1,1-Trichloroethane | | 125 | IUD |
| 56-23-5 | Carbon Tetrachloride | | 125 | IUD |
| 108-05-4 | Vinyl Acetate | | 150 | IUD |
| 75-27-4 | Bromodichloromethane | | 125 | IUD |
| 78-87-5 | 1,2-Dichloropropane | | 125 | IUD |
| 10061-01-5 | cis-1,3-Dichloropropane | | 125 | IUD |
| 79-01-6 | Trichloroethene | | 1370 | ID |
| 124-48-1 | Dibromochloromethane | | 125 | IUD |
| 79-00-5 | 1,1,2-Trichloroethane | | 125 | IUD |
| 71-43-2 | Benzene | | 125 | IUD |
| 10061-02-6 | trans-1,3-Dichloropropane | | 125 | IUD |
| 75-25-2 | Bromoform | | 125 | IUD |
| 108-10-1 | 4-Methyl-2-Pentanone | | 150 | IUD |
| 591-78-6 | 2-Hexanone | | 150 | IUD |
| 127-18-4 | Tetrachloroethene | | 125 | IUD |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 125 | IUD |
| 108-88-3 | Toluene | | 125 | IUD |
| 108-90-7 | Chlorobenzene | | 125 | IUD |
| 100-41-4 | Ethylbenzene | | 125 | IUD |
| 100-42-5 | Styrene | | 125 | IUD |
| 1330-20-7 | Xylene (total) | | 125 | IUD |



Data File: >C0983:::U4

Quant Output File: >C0983:::AQ

Name: V08, 901001.C

Misc: C04759V4, QV70357.L, 1,,

ID File: IC1172:::US

Title: IFB, PP/V08, TCL, XUD013

Last Calibration: 901003 17:09

Operator ID: PT1575

Quant Time: 901003 20:45

Injected at: 901003 19:55

QUANT REPORT

Page 1

Operator ID: PT1575
 Output File: ^C0983::AQ
 Data File: >C0983::U4
 Name: UOA,901001,C
 Misc: CA4759U45QU70357,L,1,,

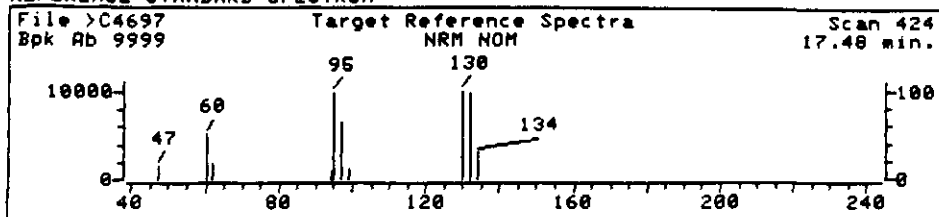
Quant Rev: 7 Quant Time: 901003 20:45
 Injected at: 901003 19:55
 Dilution Factor: 1.00000

ID File: IC1172::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901003 17:09

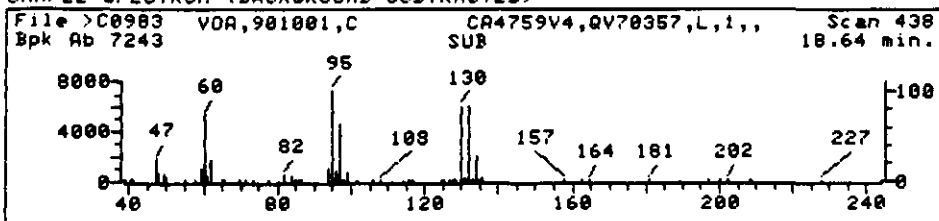
| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|-------------------|---------------|----|
| 1) *Bromochloromethane | 11.00 | 241 | 29878 | 250.00 | NG | 93 |
| 8) Acrolein | 8.40 | 174 | 2605 | 127.68 | NG | 98 |
| 16) 1,2-Trans-dichloroethylene | 12.70 | 285 | 2098 | 10.85 | NG | 95 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.98 | 318 | 109457 | 258.53 | NG | 95 |
| 20) *1,4-Difluorobenzene | 21.78 | 519 | 139647 | 250.00 | NG | 98 |
| 28) Trichloroethylene | 18.64 | 438 | 76309 | 366.89 | NG | 81 |
| 36) *Chlorobenzene-d5 | 26.73 | 646 | 162542 | 250.00 | NG | 69 |
| 41) Toluene-D8 (SURR) | 25.50 | 615 | 235191 | 244.68 | NG | 90 |
| 45) p-Bromofluorobenzene (SURR) | 31.78 | 776 | 219841 | 300.06 | NG | 88 |

* Compound is ISTD

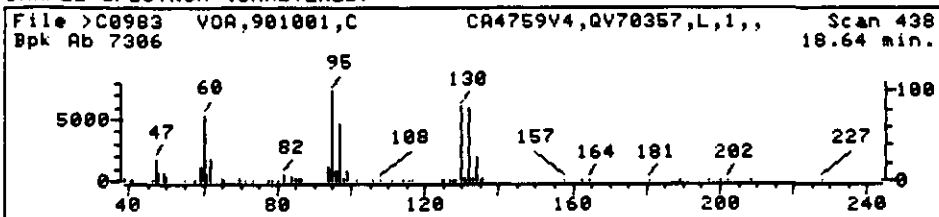
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >C0983::U4
 Name: VOA,901001,C
 Misc: CA4759V4,QV70357,L,1,,
 Quant Time: 901003 20:45
 Injected at: 901003 19:55

Quant Output File: ^C0983::AQ
 Quant ID File: IC1172::US
 Last Calibration: 901003 17:09

Compound No: 28
 Compound Name: Trichloroethylene
 Scan Number: 438
 Retention Time: 18.64 min.
 Quant Ion: 130.0
 Area: 76309
 Concentration: 366.89 NG
 q-value: 81

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

1A4760

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4760U5

Sample wt/vol: .025 (g/mL) ML

Lab File ID: >C0984

Level: (low/med) LOW MED

Date Received: 09/26/90

% Moisture: not dec.

Date Analyzed: 10/03/90

Column: (pack/cap) PACK

Dilution Factor: 200

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) | UG/L | Q |
|------------|----------------------------|---|-------|---|
| 74-87-3 | Chloromethane | | 2000 | U |
| 74-83-9 | Bromomethane | | 2000 | U |
| 75-01-4 | Vinyl Chloride | | 2000 | U |
| 75-00-3 | Chloroethane | | 2000 | U |
| 75-09-2 | Methylene Chloride | | 1000 | U |
| 67-64-1 | Acetone | | 14000 | |
| 75-15-0 | Carbon Disulfide | | 1000 | U |
| 75-35-4 | 1,1-Dichloroethene | | 1000 | U |
| 75-34-3 | 1,1-Dichloroethane | | 1000 | U |
| 540-59-0 | 1,2-Dichloroethene (total) | | 1000 | U |
| 67-66-3 | Chloroform | | 1000 | U |
| 107-06-2 | 1,2-Dichloroethane | | 1000 | U |
| 78-93-3 | 2-Butanone | | 9500 | |
| 71-55-6 | 1,1,1-Trichloroethane | | 1000 | U |
| 56-23-5 | Carbon Tetrachloride | | 1000 | U |
| 108-05-4 | Vinyl Acetate | | 2000 | U |
| 75-27-4 | Bromodichloromethane | | 1000 | U |
| 78-87-5 | 1,2-Dichloropropane | | 1000 | U |
| 10061-01-5 | cis-1,3-Dichloropropene | | 1000 | U |
| 79-01-6 | Trichloroethene | | 1000 | U |
| 124-48-1 | Dibromochloromethane | | 1000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 1000 | U |
| 71-43-2 | Benzene | | 1000 | U |
| 10061-02-6 | trans-1,3-Dichloropropene | | 1000 | U |
| 75-25-2 | Bromoform | | 1000 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | | 2000 | U |
| 591-78-6 | 2-Hexanone | | 2000 | U |
| 127-18-4 | Tetrachloroethene | | 1000 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 1000 | U |
| 108-88-3 | Toluene | | 2300 | |
| 108-90-7 | Chlorobenzene | | 1000 | U |
| 100-41-4 | Ethylbenzene | | 1000 | U |
| 100-42-5 | Styrene | | 1000 | U |
| 1330-20-7 | Xylene (total) | | 1000 | U |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETC NJ

Contract: _____

A4760

Lab Code: _____ Case No.: _____

SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER

Lab Sample ID: CA4760

Sample wt/vol: .025 (g/mL) HL

Lab File ID: >C0984

Level: (low/med) LOW MED (A)10-16-90

Date Received: 09/26/90

% Moisture: not dec. _____

Date Analyzed: 10/03/90

Column: (pack/cap) PALC

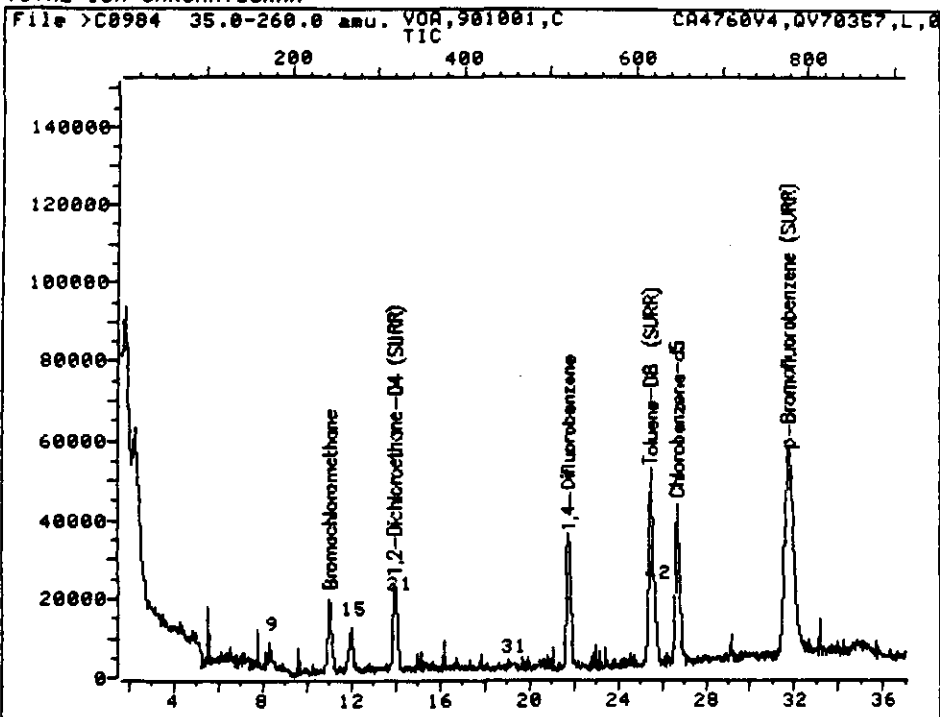
Dilution Factor: 100

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
| 3. | | | | |
| 4. | | | | |
| 5. | | | | |
| 6. | | | | |
| 7. | | | | |
| 8. | | | | |
| 9. | | | | |
| 10. | | | | |
| 11. | | | | |
| 12. | | | | |
| 13. | | | | |
| 14. | | | | |
| 15. | | | | |
| 16. | | | | |
| 17. | | | | |
| 18. | | | | |
| 19. | | | | |
| 20. | | | | |
| 21. | | | | |
| 22. | | | | |
| 23. | | | | |
| 24. | | | | |
| 25. | | | | |
| 26. | | | | |
| 27. | | | | |
| 28. | | | | |
| 29. | | | | |
| 30. | | | | |

TOTAL ION CHROMATOGRAM



Data File: >C0984::U4
Name: VOA,901001,C
Misc: CA4760V4,QV70357,L,0.025,,

Quant Output File: ^C0984::AQ

Id File: IC1172::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901003 17:09

Operator ID: PT1575
Quant Time: 901003 21:22
Injected at: 901003 20:44

QUANT REPORT

Page 1

Operator ID: PT1575
 Output File: ^C0984::AQ
 Data File: >C0984::U4
 Name: UOA,901001,C
 Misc: CA4760U~~4~~,QU70357,L,0.025,,

Quant Rev: 7 Quant Time: 901003 21:22
 Injected at: 901003 20:44
 Dilution Factor: 1.00000

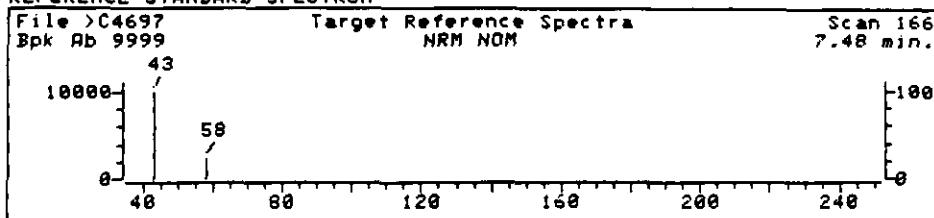
ID File: IC1172::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|-----------------|-------|-----|
| 1) *Bromochloromethane | 11.02 | 242 | 26590 | 250.00 | NG | 95 |
| 9) Acetone | 8.30 | 172 | 28010 | 361.08 | NG | 94 |
| 15) Tetrahydrofuran | 11.99 | 267 | 31422 | 1105.28 | NG | 100 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.96 | 318 | 108974 | 289.22 | NG | 98 |
| 20) *1,4-Difluorobenzene | 21.76 | 519 | 126921 | 250.00 | NG | 97 |
| 21) Methyl ethyl ketone | 14.04 | 320 | 5403 | 237.66 | NG | 93 |
| 31) Benzene | 19.20 | 453 | 2217 | 3.74 | NG | 50 |
| 36) *Chlorobenzene-d5 | 26.69 | 646 | 146632 | 250.00 | NG | 72 |
| 41) Toluene-D8 (SURR) | 25.48 | 615 | 211646 | 244.07 | NG | 89 |
| 42) Toluene | 25.68 | 620 | 28303 | 56.55 | NG | 98 |
| 45) p-Bromofluorobenzene (SURR) | 31.70 | 775 | 195826 | 296.28 | NG | 82 |

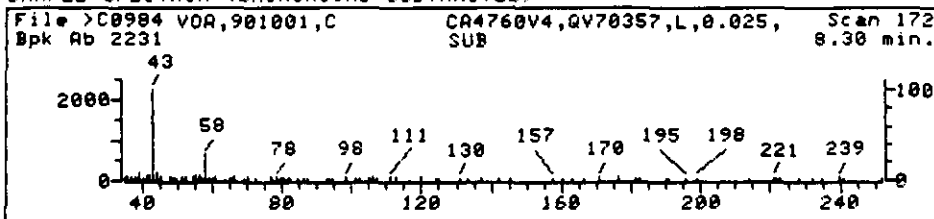
* Compound is ISTD

CV
10-16-90

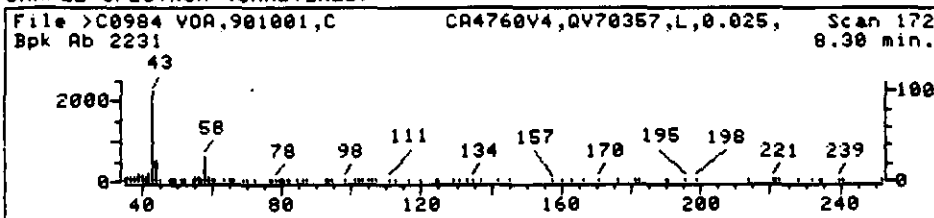
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



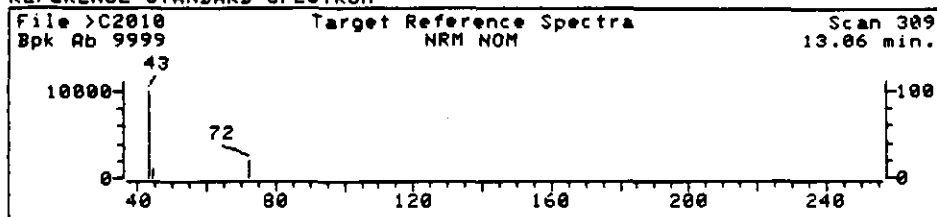
SAMPLE SPECTRUM (UNALTERED)



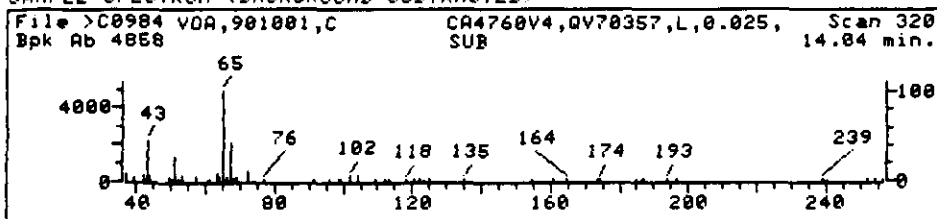
Data File: >C0984::U4 Quant Output File: ^C0984::AQ
 Name: VOA,901001,C
 Misc: CA4760V4,QV70357,L,0.025,,
 Quant Time: 901003 21:22 Quant ID File: IC1172::US
 Injected at: 901003 20:44 Last Calibration: 901003 17:09

Compound No: 9
 Compound Name: Acetone
 Scan Number: 172
 Retention Time: 8.30 min.
 Quant Ion: 43.0
 Area: 28010
 Concentration: 361.08 NG
 q-value: 94

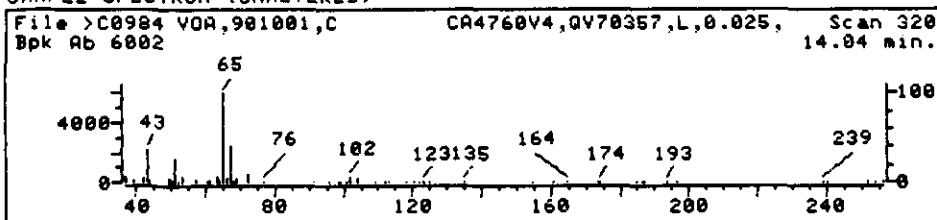
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



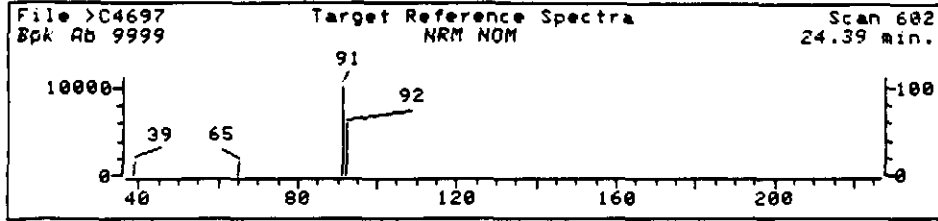
SAMPLE SPECTRUM (UNALTERED)



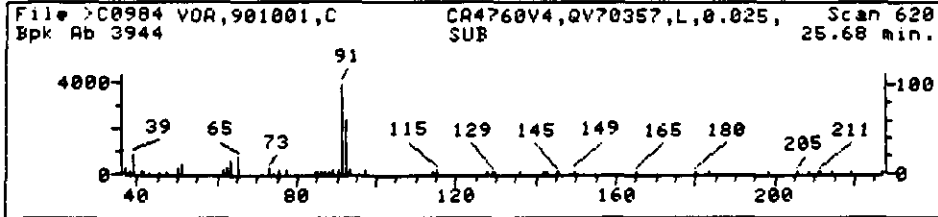
Data File: >C0984::U4 Quant Output File: ^C0984::AQ
 Name: VOA,901001,C
 Misc: CA4760V4,QV70357,L,0.025,,
 Quant Time: 901003 21:22 Quant ID File: IC1172::US
 Injected at: 901003 20:44 Last Calibration: 901003 17:09

Compound No: 21
 Compound Name: Methyl ethyl ketone
 Scan Number: 320
 Retention Time: 14.04 min.
 Quant Ion: 72.0
 Area: 5403
 Concentration: 237.66 NG
 q-value: 93

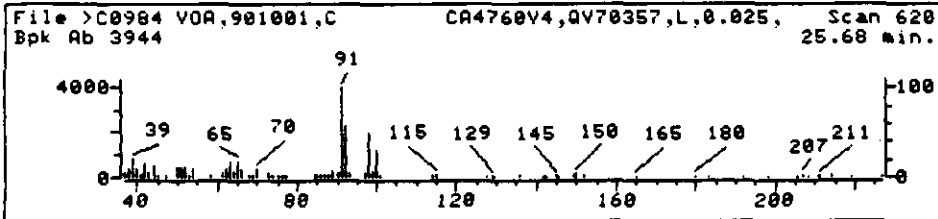
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



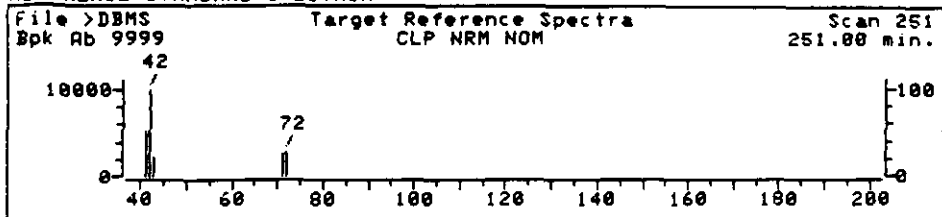
SAMPLE SPECTRUM (UNALTERED)



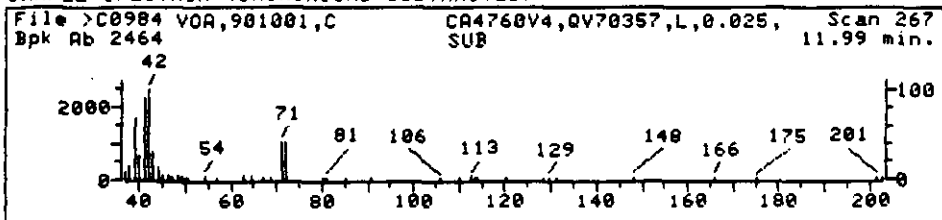
Data File: >C0984::U4 Quant Output File: ^C0984::AQ
 Name: VOA,901001,C
 Misc: CA4760V4,QV70357,L,0.025,,
 Quant Time: 901003 21:22 Quant ID File: IC1172::US
 Injected at: 901003 20:44 Last Calibration: 901003 17:09

Compound No: 42
 Compound Name: Toluene
 Scan Number: 620
 Retention Time: 25.68 min.
 Quant Ion: 92.0
 Area: 28303
 Concentration: 56.55 NG
 q-value: 98

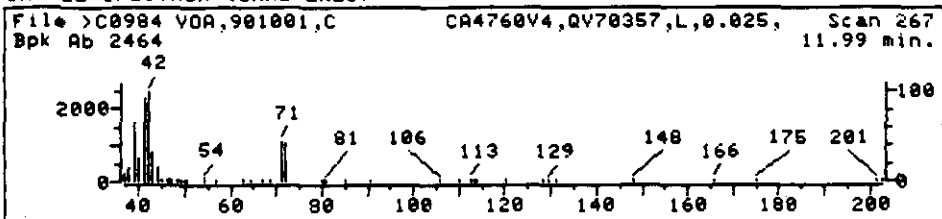
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



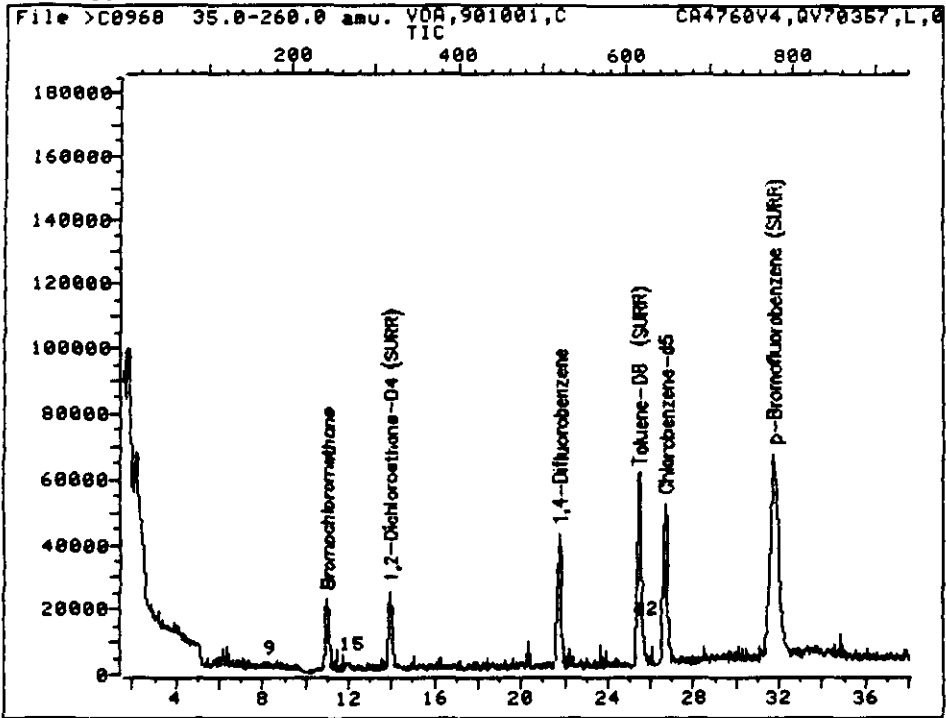
SAMPLE SPECTRUM (UNALTERED)



Data File: >C0984::U4 Quant Output File: ^C0984::AQ
 Name: VOA,901001,C
 Misc: CA4760V4,QV70357,L,0.025,,
 Quant Time: 901003 21:22 Quant ID File: IC1172::US
 Injected at: 901003 20:44 Last Calibration: 901003 17:09

Compound No: 15
 Compound Name: Tetrahydrofuran
 Scan Number: 267
 Retention Time: 11.99 min.
 Quant Ion: 42.0
 Area: 31422
 Concentration: 1105.28 NG
 q-value: 100

TOTAL ION CHROMATOGRAM



Data File: >C0968::U4
Name: VOA,901001,C
Misc: CA4760V4,QU70357,L,0.005,,

Quant Output File: ^C0968::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901002 16:53

Operator ID: JA8781
Quant Time: 901002 21:27
Injected at: 901002 20:48

QUANT REPORT

Page 1

Operator ID: JA8781
 Output File: ^C0968::AQ
 Data File: >C0968::U4
 Name: VOA,901001,C
 Misc: CA4760U4,QU70357,L,0.005,,

Quant Rev: 7 Quant Time: 901002 21:27
 Injected at: 901002 20:48
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|-----|
| 1) *Bromochloromethane | 11.00 | 243 | 30387 | 250.00 | NG | 92 |
| 9) Acetone | 8.29 | 173 | 6939 | 99.25 | NG | 84 |
| 15) Tetrahydrofuran | 11.97 | 268 | 4136 | 129.49 | NG | 100 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.95 | 319 | 125868 | 320.66 | NG | 91 |
| 20) *1,4-Difluorobenzene | 21.78 | 521 | 153691 | 250.00 | NG | 94 |
| 36) *Chlorobenzene-d5 | 26.71 | 648 | 178955 | 250.00 | NG | 67 |
| 41) Toluene-D8 (SURR) | 25.46 | 616 | 271240 | 258.82 | NG | 89 |
| 42) Toluene | 25.66 | 621 | 8329 | 14.02 | NG | 91 |
| 45) p-Bromofluorobenzene (SURR) | 31.76 | 778 | 237617 | 320.99 | NG | 78 |

* Compound is ISTD

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

| |
|-------|
| A4761 |
|-------|

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4761U4

Sample wt/vol: .02 (g/mL) ML

Lab File ID: >C0970

Level: (low/med) LOW

Date Received: 09/26/90

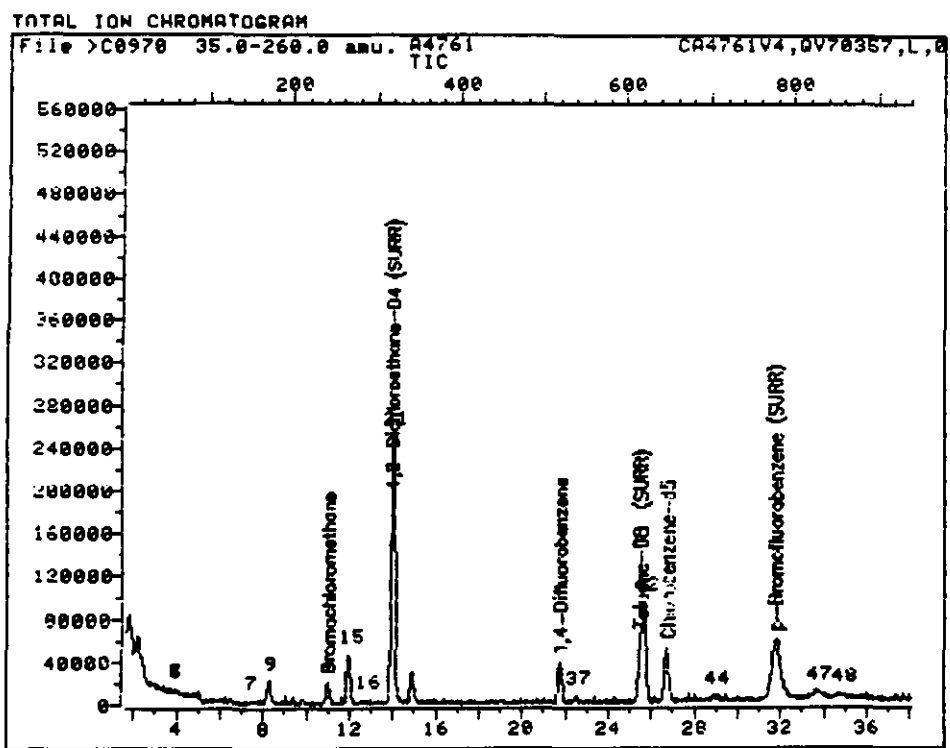
% Moisture: not dec.

Date Analyzed: 10/02/90

Column: (pack/cap) PACK

Dilution Factor: 250

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|----|
| 74-87-3 | Chloromethane | 12500 | IU |
| 74-83-9 | Bromomethane | 12500 | IU |
| 75-01-4 | Vinyl Chloride | 13300 | IU |
| 75-00-3 | Chloroethane | 12500 | IU |
| 75-09-2 | Methylene Chloride | 11300 | IU |
| 67-64-1 | Acetone | 140000 | IE |
| 75-15-0 | Carbon Disulfide | 11300 | IU |
| 75-35-4 | 1,1-Dichloroethene | 11300 | IU |
| 75-34-3 | 1,1-Dichloroethane | 11300 | IU |
| 540-59-0 | 1,2-Dichloroethene (total) | 11300 | IU |
| 67-66-3 | Chloroform | 11300 | IU |
| 107-06-2 | 1,2-Dichloroethane | 11300 | IU |
| 78-93-3 | 2-Butanone | 1830000 | IE |
| 71-55-6 | 1,1,1-Trichloroethane | 11300 | IU |
| 56-23-5 | Carbon Tetrachloride | 11300 | IU |
| 108-05-4 | Vinyl Acetate | 12500 | IU |
| 75-27-4 | Bromodichloromethane | 11300 | IU |
| 78-87-5 | 1,2-Dichloropropane | 11300 | IU |
| 10061-01-5 | cis-1,3-Dichloropropene | 11300 | IU |
| 79-01-6 | Trichloroethene | 11300 | IU |
| 124-48-1 | Dibromochloromethane | 11300 | IU |
| 79-00-5 | 1,1,2-Trichloroethane | 11300 | IU |
| 71-43-2 | Benzene | 11300 | IU |
| 10061-02-6 | trans-1,3-Dichloropropene | 11300 | IU |
| 75-25-2 | Bromoform | 11300 | IU |
| 108-10-1 | 4-Methyl-2-Pentanone | 13100 | IU |
| 591-78-6 | 2-Hexanone | 12500 | IU |
| 127-18-4 | Tetrachloroethene | 11300 | IU |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 11300 | IU |
| 108-88-3 | Toluene | 18000 | IU |
| 108-90-7 | Chlorobenzene | 11300 | IU |
| 100-41-4 | Ethylbenzene | 1700 | IU |
| 100-42-5 | Styrene | 11300 | IU |
| 1330-20-7 | Xylene (total) | 13300 | IU |



Data File: >C0970::U4 Quant Output File: ^C0970::AQ
 Name: A4761
 Misc: CA4761U4,QV70357,L,0.02,,

 Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

 Operator ID: JAB/81
 Quant Time: 901002 23:04
 Injected at: 901002 22:25

QUANT REPORT

Operator ID: JA8781
 Output File: ^C0970::AQ
 Data File: >C0970::U4
 Name: A4761
 Misc: CA4761U4,QU70357,L,0.02,,

Quant Rev: 7 Quant Time: 901002 23:04
 Injected at: 901002 22:25
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

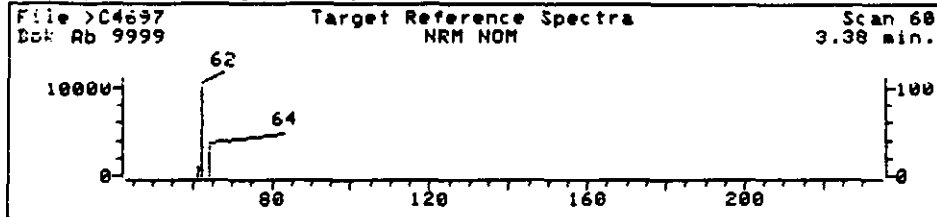
| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|-------|------------|
| 1) *Bromochloromethane | 11.01 | 243 | 28016 | 250.00 | NG | 96 |
| 4) Dichlorodifluoromethane | 3.95 | 61 | 3344 | 64.89 | NG | 90 |
| 5) Vinyl chloride | 3.95 | 61 | 4878 | 65.11 | NG | 86 |
| 7) Methylene chloride | 7.45 | 151 | 4490 | 36.17 | NG | 97 |
| 9) Acetone | 8.26 | 172 | 180294 | 2796.53 | NG | 3350* 97 |
| 15) Tetrahydrofuran | 11.94 | 267 | 132089 | 4485.60 | NG | 2200* 100 |
| 16) 1,2-Trans-dichloroethylene | 12.72 | 287 | 3092 | 16.85 | NG | 95 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.96 | 319 | 117337 | 324.22 | NG | 93 |
| 20) *1,4-Difluorobenzene | 21.75 | 520 | 145859 | 250.00 | NG | 96 |
| 21) Methyl ethyl ketone | 14.00 | 320 | 336682 | 16573.60 | NG | -10333* 95 |
| 36) *Chlorobenzene-d5 | 26.73 | 648 | 173819 | 250.00 | NG | 73 |
| 37) Methyl-iso-butyl ketone | 22.49 | 539 | 14691 | 61.76 | NG | 83 |
| 41) Toluene-D8 (SURR) | 25.48 | 616 | 255850 | 251.34 | NG | 90 |
| 42) Toluene | 25.67 | 621 | 213132 | 369.41 | NG | 93 |
| 44) Ethylbenzene | 28.98 | 706 | 6810 | 14.02 | NG | 82 |
| 45) p-Bromofluorobenzene (SURR) | 31.73 | 777 | 226610 | 315.17 | NG | 77 |
| 47) m-Xylene | 33.71 | 828 | 21377 | 33.59 | NG | 93 |
| 48) o+p-Xylenes | 34.87 | 858 | 21115 | 32.13 | NG | 85 |

* Compound is ISTD

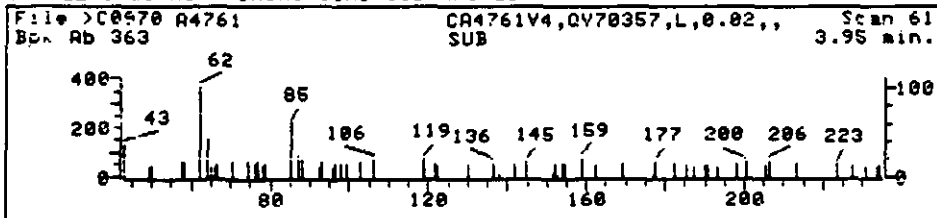
(21)
 10-16-90

* values are from dilution run.

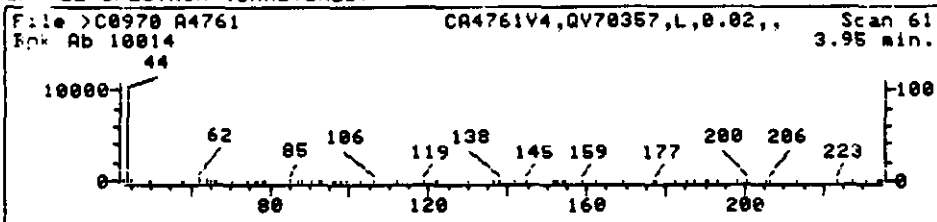
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

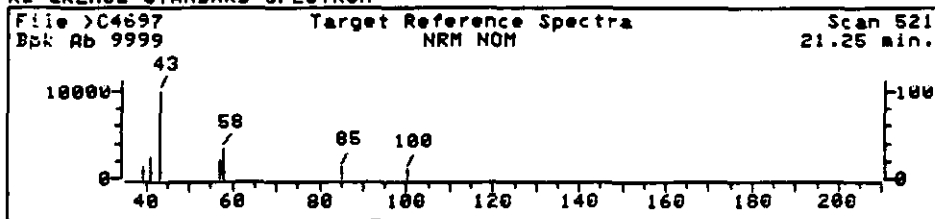


Data File: >C0970::U4
 Name: A4761
 Misc: CA4761V4,QV70357,L,0.02,,
 Quant Time: 901002 23:04
 Injected at: 901002 22:25

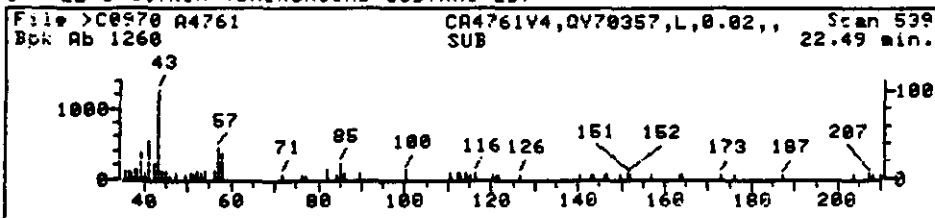
Quant Output File: >C0970::AQ
 Quant ID File: IC1171::US
 Last Calibration: 901002 16:53

Compound No: 5
 Compound Name: Vinyl chloride
 Scan Number: 61
 Retention Time: 3.95 min.
 Quant Ion: 62.0
 Area: 4878
 Concentration: 65.11 NG
 q-value: 86

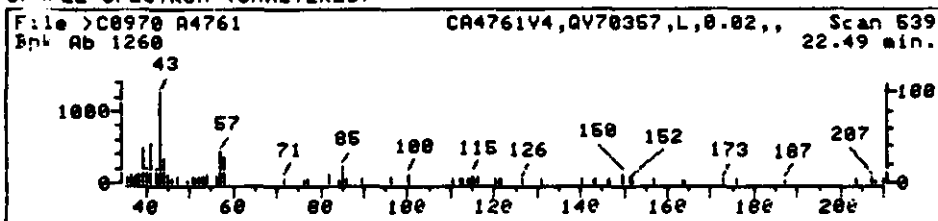
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



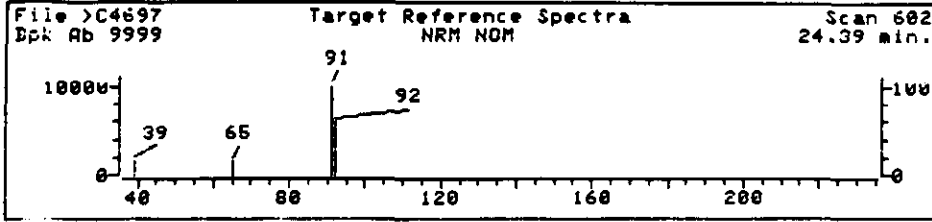
SAMPLE SPECTRUM (UNALTERED)



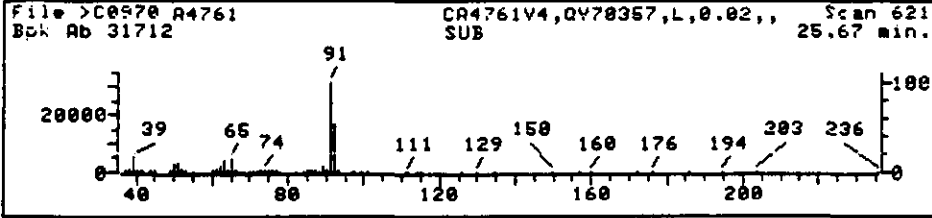
Data File: >C0970::U4 Quant Output File: ^C0970::AQ
 Name: A4761
 Misc: CA4761V4,QV70357,L,0.02,,
 Quant Time: 901002 23:04 Quant ID File: IC1171::US
 Injected at: 901002 22:25 Last Calibration: 901002 16:53

Compound No: 37
 Compound Name: Methyl-iso-butyl ketone
 Scan Number: 539
 Retention Time: 22.49 min.
 Quant Ion: 43.0
 Area: 14691
 Concentration: 61.76 NG
 q-value: 83

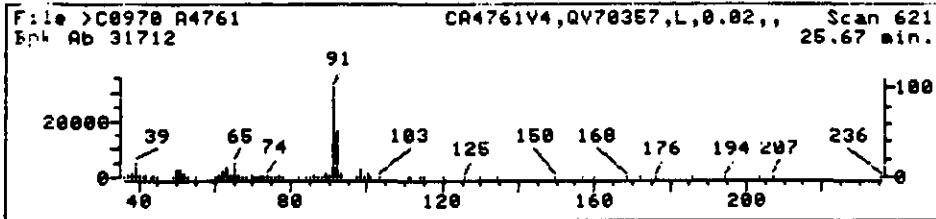
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

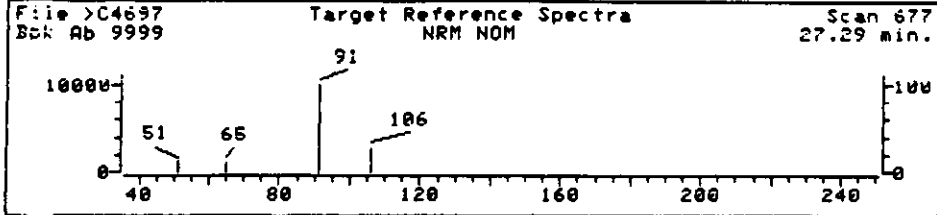


Data File: >C0970::U4
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 Misc: CA4761V4,QV70357,L,0.02,,
 Quant Time: 901002 23:04
 Injected at: 901002 22:25

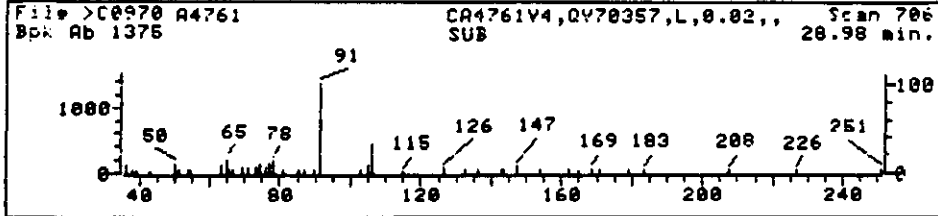
Quant Output File: ^C0970::AQ
 Quant ID File: IC1171::US
 Last Calibration: 901002 16:53

Compound No: 42
 Compound Name: Toluene
 Scan Number: 621
 Retention Time: 25.67 min.
 Quant Ion: 92.0
 Area: 213132
 Concentration: 369.41 NG
 q-value: 93

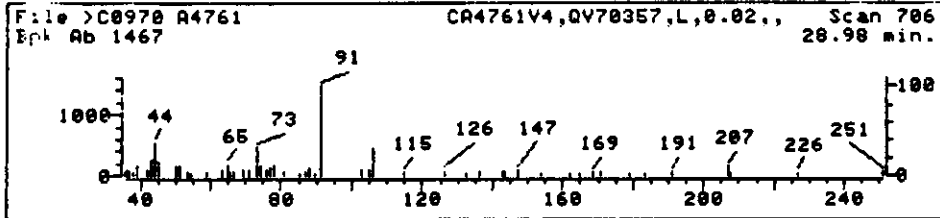
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >C0970::U4

Quant Output File: ^C0970::AQ

Name: A4761

Misc: CA4761V4,QV70357,L,0.02,,

Quant ID File: IC1171::US

Quant Time: 901002 23:04

Last Calibration: 901002 16:53

Injected at: 901002 22:25

Compound No: 44

Compound Name: Ethylbenzene

Scan Number: 706

Retention Time: 28.98 min.

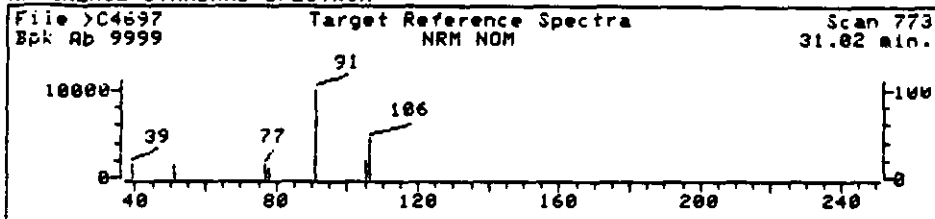
Quant Ion: 106.0

Area: 6810

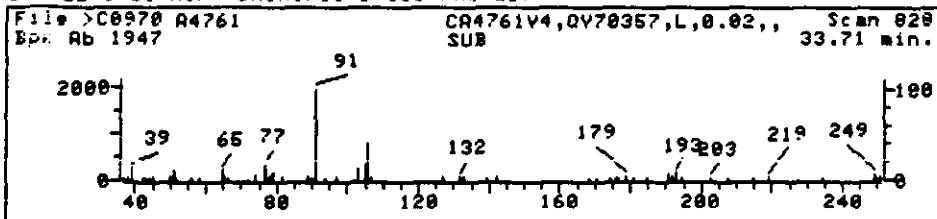
Concentration: 14.02 NG

q-value: 82

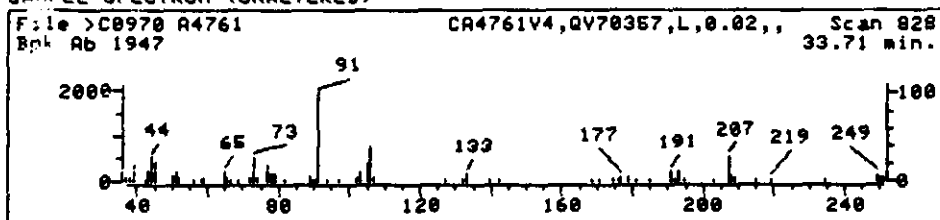
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

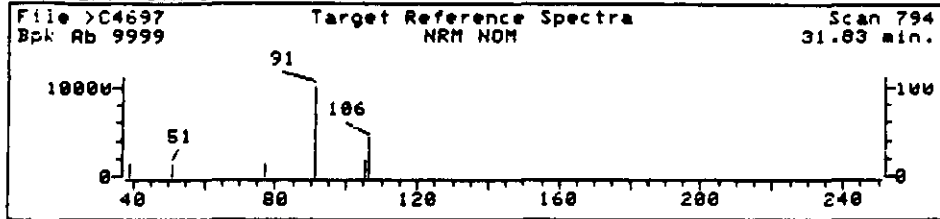


Data File: >C0970::U4
 Name: A4761
 Misc: CA4761V4,QV70357,L,0.02,,
 Quant Time: 901002 23:04
 Injected at: 901002 22:25

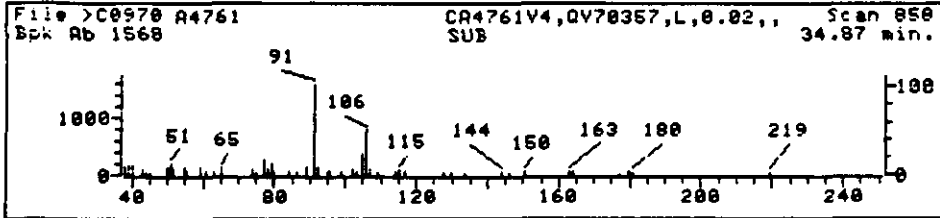
Quant Output File: ^C0970::AQ
 Quant ID File: IC1171::US
 Last Calibration: 901002 16:53

Compound No: 47
 Compound Name: m-Xylene
 Scan Number: 828
 Retention Time: 33.71 min.
 Quant Ion: 106.0
 Area: 21377
 Concentration: 33.59 NG
 q-value: 93

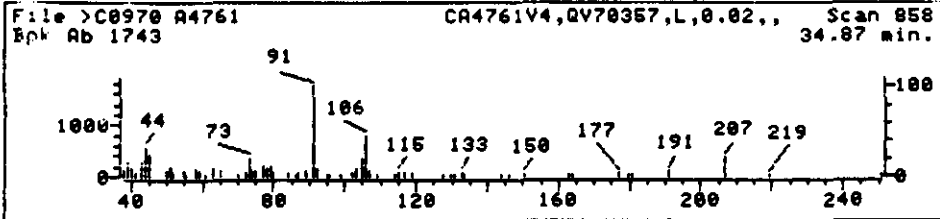
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



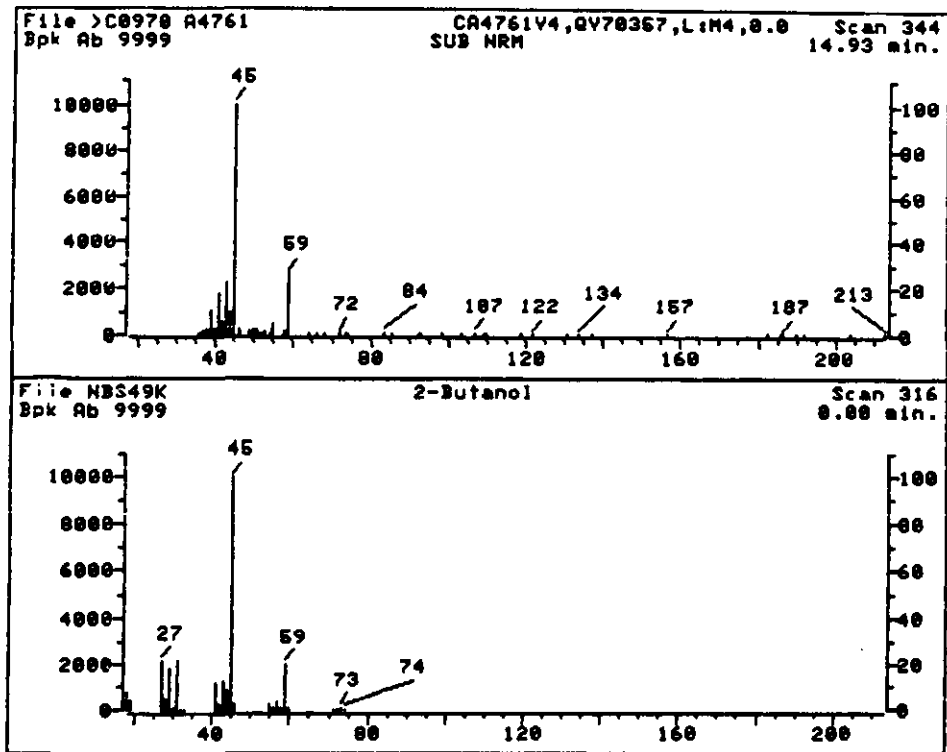
SAMPLE SPECTRUM (UNALTERED)



Data File: >C0970::U4
 Name: A4761
 Misc: CA4761V4,QV70357,L,0.02,,
 Quant Time: 901002 23:04
 Injected at: 901002 22:25

Quant Output File: ^C0970::AQ
 Quant ID File: IC1171::US
 Last Calibration: 901002 16:53

Compound No: 48
 Compound Name: o+p-Xylenes
 Scan Number: 858
 Retention Time: 34.87 min.
 Quant Ion: 106.0
 Area: 21115
 Concentration: 32.13 NG
 q-value: 85



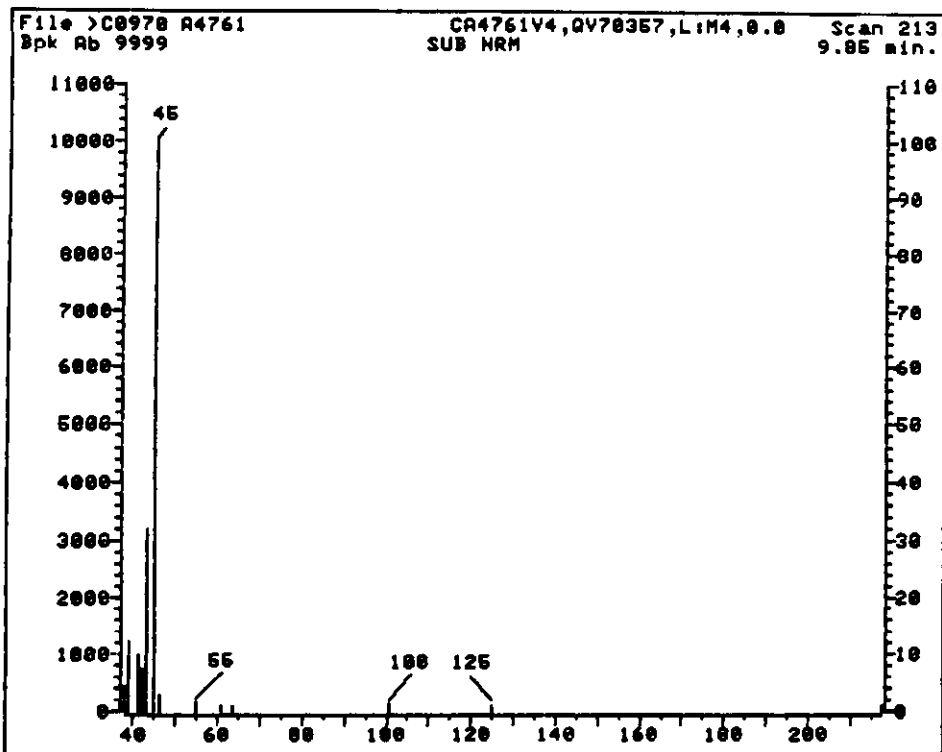
Data File: >C0970
Name: A4761
Misc Data: CA4761V4,QU70357,L:M4,0.02,, 1:250 DL
RT (min): 14.93
Scan: 344
Area: 347476 Rank: 6
Semi-quantitative Conc (uncorrected): 318.16 NG
Semi-quantitative Conc (corrected): 15907.80 ug/l
Calculated using Istd: Bromochloromethane @ 11.01 minutes

1. 2-Butanol

74 C4H10O

Sample file: >C0970 Spectrum #: 344
Search speed: 2 Tilting option: S No. of ion ranges searched: 45

| | Prob. | CAS # | CON # | ROOT | K | DK | #FLG | TILT | % | CON | C_1 | R_IV |
|----|-------|-------|-------|--------|----|----|------|------|-----|-----|-----|------|
| 1. | 60 | 78922 | 1715 | NBS49K | 35 | 48 | 2 | 0 | 100 | 14 | 30 | 12 |



Data File: >C0970
 Name: A4761
 Misc Data: CA4761V4,QU70357,L:M4,0.02,, 1:250 DL
 RT (min): 9.85
 Scan: 213
 Area: 57973 Rank: 8
 Semi-quantitative Conc (uncorrected): 53.08 NG
 Semi-quantitative Conc (corrected): 2654.06 ug/l
 Calculated using Istd: Bromochloromethane @ 11.01 minutes

No PBM hits for this scan.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4761DL

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4761U5

Sample wt/vol: .0 (g/mL) ML

Lab File ID: >C0986

Level: (low/med) LOW

Date Received: 09/26/90

% Moisture: not dec.

Date Analyzed: 10/03/90

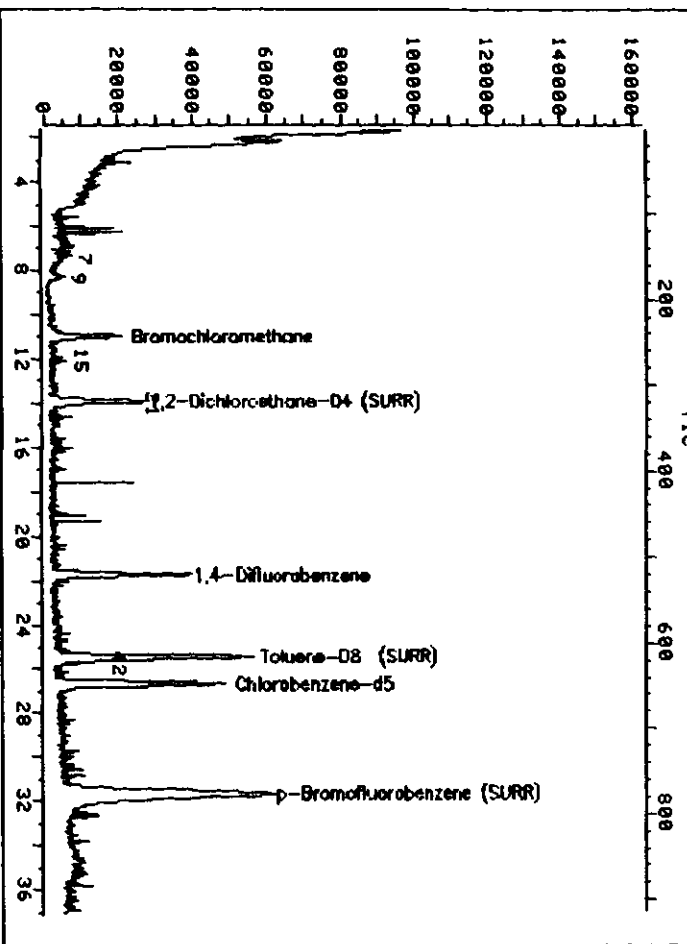
Column: (pack/cap) PACK

Dilution Factor: 5000

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|----|
| 74-87-3 | Chloromethane | 50000 | UD |
| 74-83-9 | Bromomethane | 50000 | UD |
| 75-01-4 | Vinyl Chloride | 50000 | UD |
| 75-00-3 | Chloroethane | 50000 | UD |
| 75-09-2 | Methylene Chloride | 25000 | UD |
| 67-64-1 | Acetone | 170000 | ID |
| 75-15-0 | Carbon Disulfide | 25000 | UD |
| 75-35-4 | 1,1-Dichloroethene | 25000 | UD |
| 75-34-3 | 1,1-Dichloroethane | 25000 | UD |
| 540-59-0 | 1,2-Dichloroethene (total) | 25000 | UD |
| 67-66-3 | Chloroform | 25000 | UD |
| 107-06-2 | 1,2-Dichloroethane | 25000 | UD |
| 78-93-3 | 2-Butanone | 520000 | ID |
| 71-55-6 | 1,1,1-Trichloroethane | 25000 | UD |
| 56-23-5 | Carbon Tetrachloride | 25000 | UD |
| 108-05-4 | Vinyl Acetate | 50000 | UD |
| 75-27-4 | Bromodichloromethane | 25000 | UD |
| 78-87-5 | 1,2-Dichloropropane | 25000 | UD |
| 10061-01-5 | cis-1,3-Dichloropropene | 25000 | UD |
| 79-01-6 | Trichloroethene | 25000 | UD |
| 124-48-1 | Dibromochloromethane | 25000 | UD |
| 79-00-5 | 1,1,2-Trichloroethane | 25000 | UD |
| 71-43-2 | Benzene | 25000 | UD |
| 10061-02-6 | trans-1,3-Dichloropropene | 25000 | UD |
| 75-25-2 | Bromoform | 25000 | UD |
| 108-10-1 | 4-Methyl-2-Pentanone | 50000 | UD |
| 591-78-6 | 2-Hexanone | 50000 | UD |
| 127-18-4 | Tetrachloroethene | 25000 | UD |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 25000 | UD |
| 108-88-3 | Toluene | 21000 | UD |
| 108-90-7 | Chlorobenzene | 25000 | UD |
| 100-41-4 | Ethylbenzene | 25000 | UD |
| 100-42-5 | Styrene | 25000 | UD |
| 1330-20-7 | Xylene (total) | 25000 | UD |

TOTAL ION CHROMATOGRAM

File: >C0986 35.0-260.0 amu. V09,901001.C CR4762V4,0V70357.L,0



Data File: >C0986::U4

Quant Output File: <C0986::AQ

Name: V09,901001.C

Misc: CA4762V4,0V70357.L,0.001,,
for 10/22/09

ID File: IC1172::US

Title: IFB, PP/V09, TCL, XV09A13

Last Calibration: 901003 17:09

Operator ID: PT1575

Quant Time: 901003 22:59

Injected at: 901003 22:21

QUANT REPORT

Operator ID: PT1575
 Output File: ^C0986::AQ
 Data File: >C0986::U4
 Name: VOA,901001,C
 Misc: CA4767U45QUV70357,L,0.001,,
 1

Quant Rev: 7 Quant Time: 901003 22:59
 Injected at: 901003 22:21
 Dilution Factor: 1.00000

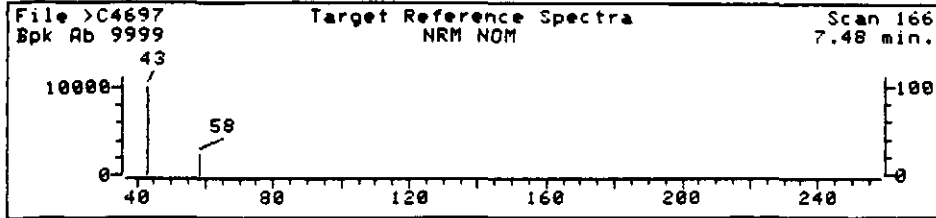
ID File: IC1172::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|-----|
| 1) *Bromochloromethane | 10.96 | 241 | 27061 | 250.00 | NG | 99 |
| 7) Methylene chloride | 7.43 | 150 | 6308 | 44.48 | NG | 92 |
| 9) Acetone | 8.29 | 172 | 13223 | 167.49 | NG ✓ | 89 |
| 15) Tetrahydrofuran | 11.97 | 267 | 4629 | 159.99 | NG ✓ | 100 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.91 | 317 | 116601 | 304.08 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.74 | 519 | 136448 | 250.00 | NG | 94 |
| 21) Methyl ethyl ketone | 13.98 | 319 | 12628 | 516.69 | NG ✓ | 97 |
| 36) *Chlorobenzene-d5 | 26.68 | 646 | 157154 | 250.00 | NG | 66 |
| 41) Toluene-D8 (SURR) | 25.43 | 614 | 231507 | 249.10 | NG | 83 |
| 42) Toluene | 25.62 | 619 | 11180 | 20.84 | NG ✓ | 95 |
| 45) p-Bromofluorobenzene (SURR) | 31.64 | 774 | 214732 | 303.13 | NG | 82 |

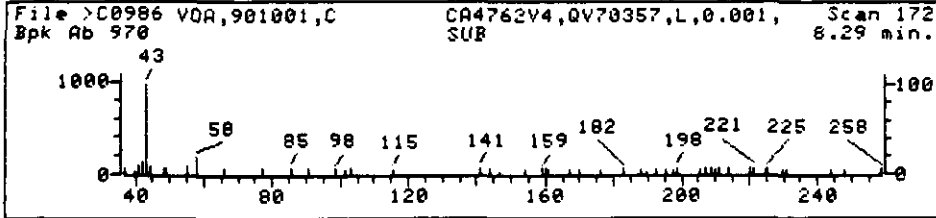
* Compound is ISTD

CA
 10-16-90.

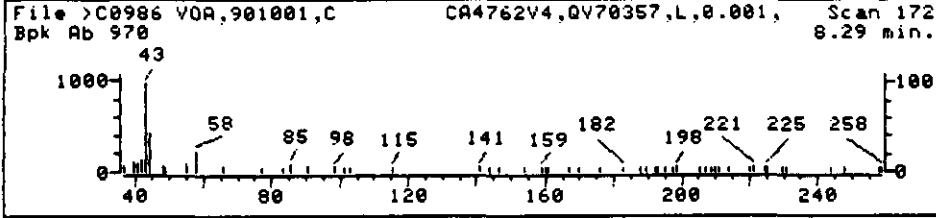
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



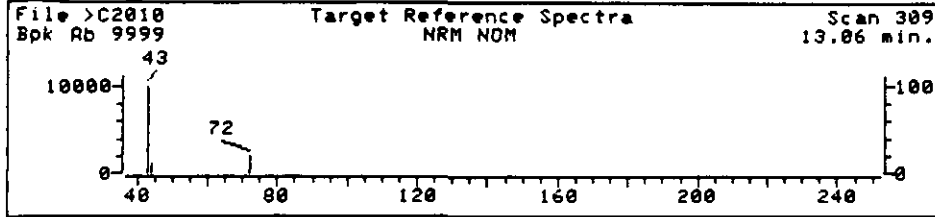
SAMPLE SPECTRUM (UNALTERED)



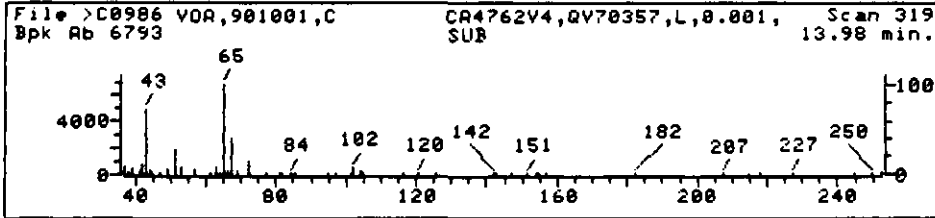
Data File: >C0986::U4 Quant Output File: ^C0986::AQ
 Name: UDA,901001,C *2/12/90*
 Misc: CA4762V4,QV70357,L,0.001,,
 Quant Time: 901003 22:59 Quant ID File: IC1172::US
 Injected at: 901003 22:21 Last Calibration: 901003 17:09

Compound No: 9
 Compound Name: Acetone
 Scan Number: 172
 Retention Time: 8.29 min.
 Quant Ion: 43.0
 Area: 13223
 Concentration: 167.49 NG
 q-value: 89

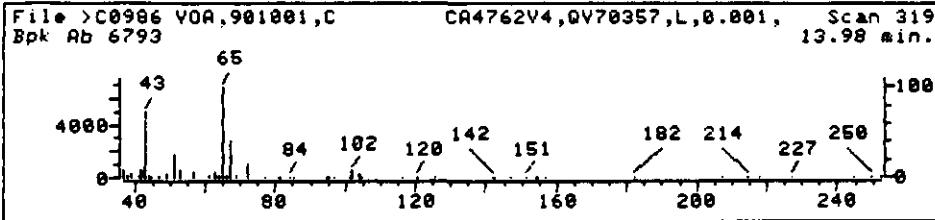
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



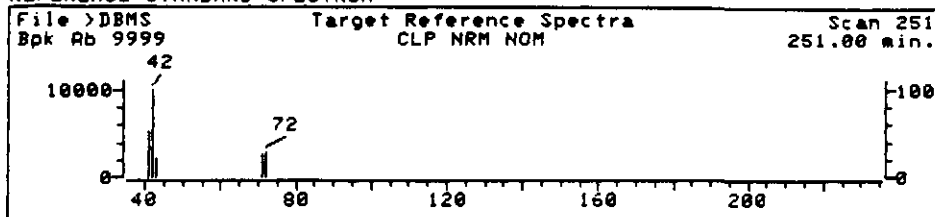
SAMPLE SPECTRUM (UNALTERED)



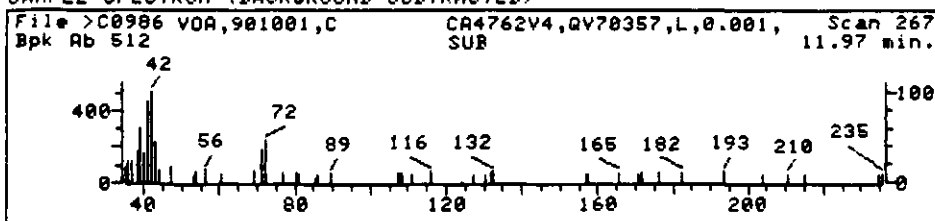
Data File: >C0986::U4 Quant Output File: ^C0986::AQ
 Name: VOA,901001,C *for 9/21/00*
 Misc: CA4762V4,QV70357,L,0.001,,
 Quant Time: 901003 22:59 Quant ID File: IC1172::US
 Injected at: 901003 22:21 Last Calibration: 901003 17:09

Compound No: 21
 Compound Name: Methyl ethyl ketone
 Scan Number: 319
 Retention Time: 13.98 min.
 Quant Ion: 72.0
 Area: 12628
 Concentration: 516.69 NG
 q-value: 97

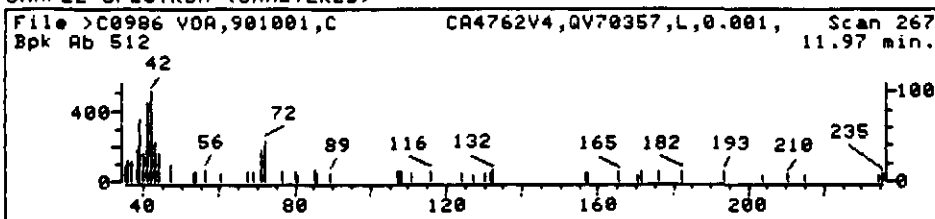
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >C0986::U4 Quant Output File: ^C0986::AQ
 Name: VOA,901001,C *on 1/22/90*
 Misc: CA4762V4,QV70357,L,0.001,,
 Quant Time: 901003 22:59 Quant ID File: IC1172::US
 Injected at: 901003 22:21 Last Calibration: 901003 17:09

Compound No: 15
 Compound Name: Tetrahydrofuran
 Scan Number: 267
 Retention Time: 11.97 min.
 Quant Ion: 42.0
 Area: 4629
 Concentration: 159.99 NG
 q-value: 100

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

| |
|-------|
| A4762 |
|-------|

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4762U4

Sample wt/vol: .5 (g/mL) ML

Lab File ID: >C0969

Level: (low/med) ~~LOW~~ MED ¹⁰⁻¹¹⁻⁹⁰

Date Received: 09/26/90

% Moisture: not dec.

Date Analyzed: 10/02/90

Column: (pack/cap) PACK

Dilution Factor: 10

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|----|
| 74-87-3 | Chloromethane | 100 | IU |
| 74-83-9 | Bromomethane | 100 | IU |
| 75-01-4 | Vinyl Chloride | 330 | I |
| 75-00-3 | Chloroethane | 100 | IU |
| 75-09-2 | Methylene Chloride | 50 | IU |
| 67-64-1 | Acetone | 5200 | IE |
| 75-15-0 | Carbon Disulfide | 50 | IU |
| 75-35-4 | 1,1-Dichloroethene | 50 | IU |
| 75-34-3 | 1,1-Dichloroethane | 50 | IU |
| 540-59-0 | 1,2-Dichloroethene (total) | 1400 | I |
| 67-66-3 | Chloroform | 50 | IU |
| 107-06-2 | 1,2-Dichloroethane | 50 | IU |
| 78-93-3 | 2-Butanone | 25000 | IE |
| 71-55-6 | 1,1,1-Trichloroethane | 50 | IU |
| 56-23-5 | Carbon Tetrachloride | 50 | IU |
| 108-05-4 | Vinyl Acetate | 100 | IU |
| 75-27-4 | Bromodichloromethane | 50 | IU |
| 78-87-5 | 1,2-Dichloropropane | 50 | IU |
| 10061-01-5 | cis-1,3-Dichloropropane | 50 | IU |
| 79-01-6 | Trichloroethene | 37 | IJ |
| 124-48-1 | Dibromochloromethane | 50 | IU |
| 79-00-5 | 1,1,2-Trichloroethane | 50 | IU |
| 71-43-2 | Benzene | 50 | IU |
| 10061-02-6 | trans-1,3-Dichloropropane | 50 | IU |
| 75-25-2 | Bromoform | 50 | IU |
| 108-10-1 | 4-Methyl-2-Pentanone | 110 | I |
| 591-78-6 | 2-Hexanone | 100 | IU |
| 127-18-4 | Tetrachloroethene | 50 | IU |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 50 | IU |
| 108-88-3 | Toluene | 570 | I |
| 108-90-7 | Chlorobenzene | 50 | IU |
| 100-41-4 | Ethylbenzene | 50 | IU |
| 100-42-5 | Styrene | 50 | IU |
| 1330-20-7 | Xylene (total) | 66 | I |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETC NJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: CA4762

Sample wt/vol: .5 (g/mL) HL Lab File ID: 7C0969

Level: (low/med) LOW Date Received: 09/26/90

Moisture: not dec. _____ Date Analyzed: 10/02/90

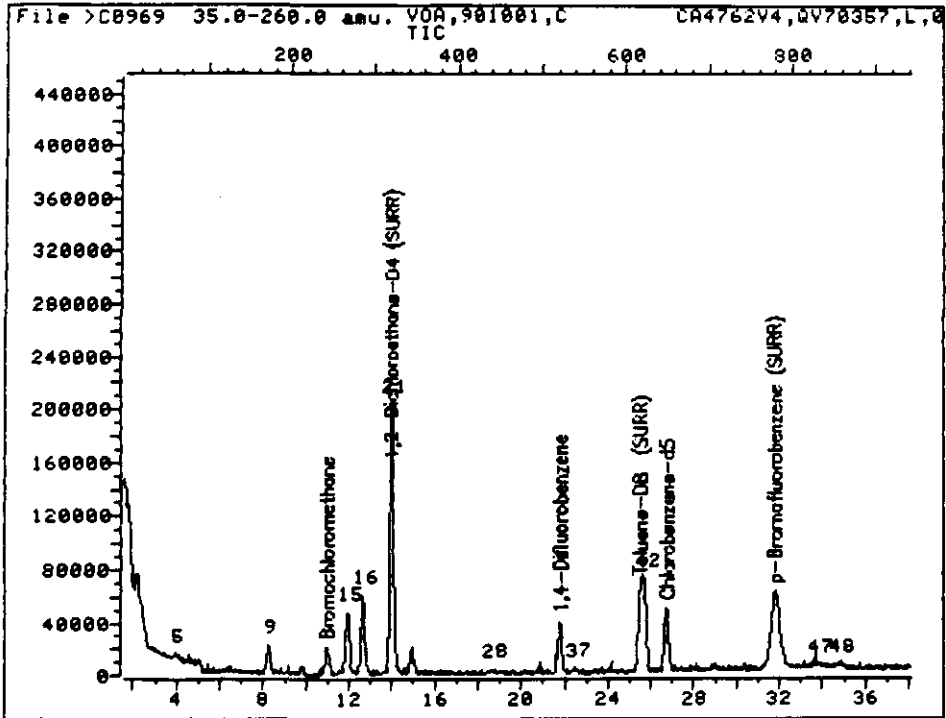
Column: (pack/cap) PALIK Dilution Factor: 10

Number TICs found: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|-------|------------|----|
| 1. | 2-Butanol | 14.92 | 440 | 11 |
| 2. | UNKNOWN | 9.89 | 160 | 11 |
| 3. | | | | |
| 4. | | | | |
| 5. | | | | |
| 6. | | | | |
| 7. | | | | |
| 8. | | | | |
| 9. | | | | |
| 10. | | | | |
| 11. | | | | |
| 12. | | | | |
| 13. | | | | |
| 14. | | | | |
| 15. | | | | |
| 16. | | | | |
| 17. | | | | |
| 18. | | | | |
| 19. | | | | |
| 20. | | | | |
| 21. | | | | |
| 22. | | | | |
| 23. | | | | |
| 24. | | | | |
| 25. | | | | |
| 26. | | | | |
| 27. | | | | |
| 28. | | | | |
| 29. | | | | |
| 30. | | | | |

TOTAL ION CHROMATOGRAM



Data File: >C0969::U4
Name: VOA,901001,C
Misc: CA4762V4,QU70357,L,0.5,,

Quant Output File: ^C0969::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901002 16:53

Operator ID: JA8781
Quant Time: 901002 22:15
Injected at: 901002 21:37

QUANT REPORT

Operator ID: JA8781
 Output File: ^C0969::AQ
 Data File: >C0969::U4
 Name: UOA,901001,C
 Misc: CA4762U4,QU70357,L,0.5,,

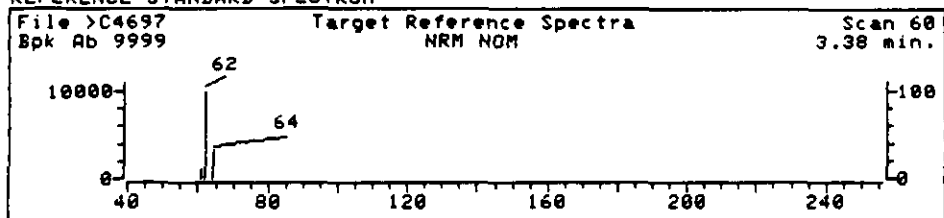
Quant Rev: 7 Quant Time: 901002 22:15
 Injected at: 901002 21:37
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901002 16:53

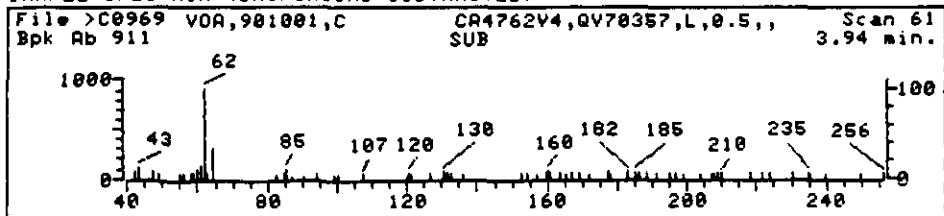
| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|----------|----------------------|-----|
| 1) *Bromochloromethane | 10.96 | 242 | 29250 | 250.00 | NG | 98 |
| 5) Vinyl chloride | 3.94 | 61 | 12789 | 163.49 | NG | 99 |
| 9) Acetone | 8.25 | 172 | 176238 | 2618.29 | NG ^{-2800*} | 98 |
| 15) Tetrahydrofuran | 11.93 | 267 | 137914 | 4485.83 | NG ^{-4980*} | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.67 | 286 | 134570 | 702.26 | NG | 93 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.95 | 319 | 122208 | 323.44 | NG | 93 |
| 20) *1,4-Difluorobenzene | 21.78 | 521 | 152364 | 250.00 | NG | 94 |
| 21) Methyl ethyl ketone | 13.99 | 320 | 266355 | 12551.88 | NG ^{-8075*} | 96 |
| 28) Trichloroethylene | 18.68 | 441 | 4193 | 18.63 | NG | 75 |
| 36) *Chlorobenzene-d5 | 26.75 | 649 | 179637 | 250.00 | NG | 70 |
| 37) Methyl-iso-butyl ketone | 22.52 | 540 | 13019 | 52.96 | NG | 89 |
| 41) Toluene-D8 (SURR) | 25.50 | 617 | 274359 | 260.80 | NG | 90 |
| 42) Toluene | 25.70 | 622 | 168759 | 283.03 | NG | 96 |
| 45) p-Bromofluorobenzene (SURR) | 31.76 | 778 | 235927 | 317.50 | NG | 80 |
| 47) m-Xylene | 33.81 | 831 | 10939 | 16.63 | NG | 95 |
| 48) o+p-Xylenes | 34.82 | 857 | 11268 | 16.59 | NG | 82 |

* Compound is ISTD

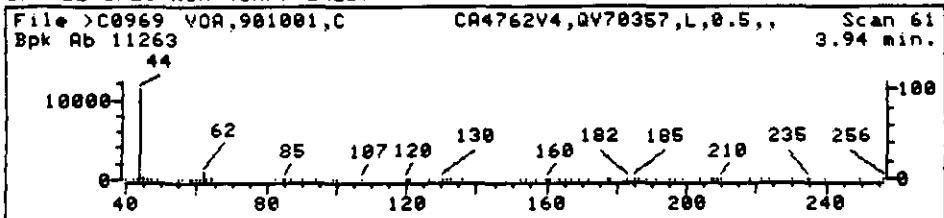
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

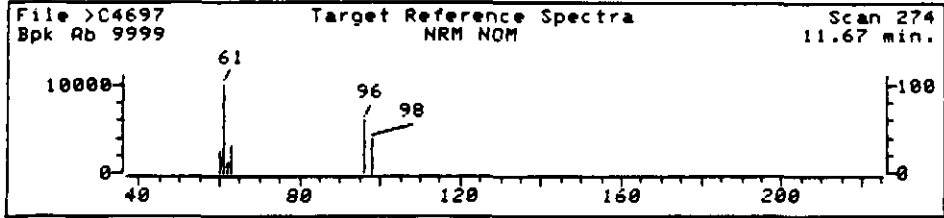


Data File: >C0969::U4
 Name: VOA,901001,C
 Misc: CA4762V4,QV70357,L,0.5,,
 Quant Time: 901002 22:15
 Injected at: 901002 21:37

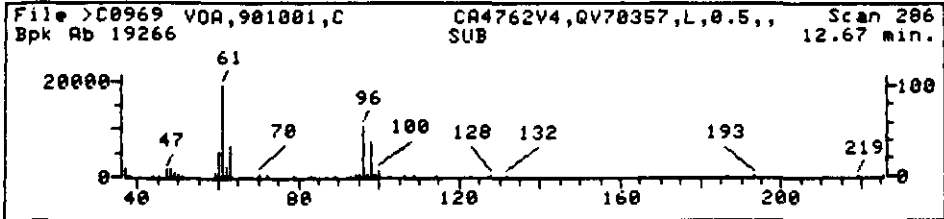
Quant Output File: ^C0969::AQ
 Quant ID File: IC1171::US
 Last Calibration: 901002 16:53

Compound No: 5
 Compound Name: Vinyl chloride
 Scan Number: 61
 Retention Time: 3.94 min.
 Quant Ion: 62.0
 Area: 12789
 Concentration: 163.49 NG
 q-value: 99

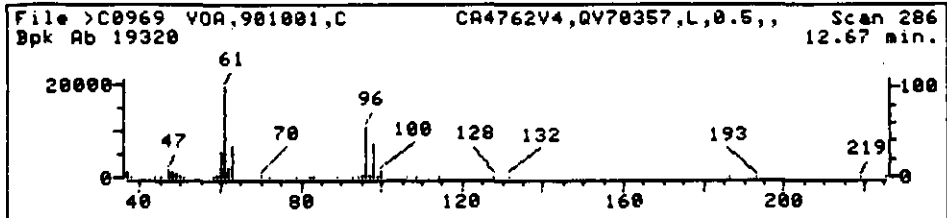
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

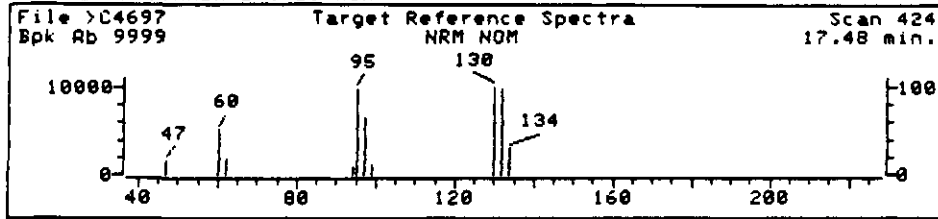


Data File: >C0969::U4
Name: VOA,901001,C
Misc: CA4762V4,QV70357,L,0.5,,
Quant Time: 901002 22:15
Injected at: 901002 21:37

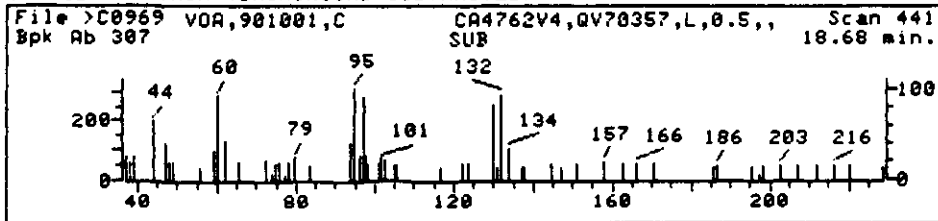
Quant Output File: ^C0969::AQ
Quant ID File: IC1171::US
Last Calibration: 901002 16:53

Compound No: 16
Compound Name: 1,2-Trans-dichloroethylene
Scan Number: 286
Retention Time: 12.67 min.
Quant Ion: 96.0
Area: 134570
Concentration: 702.26 NG
q-value: 93

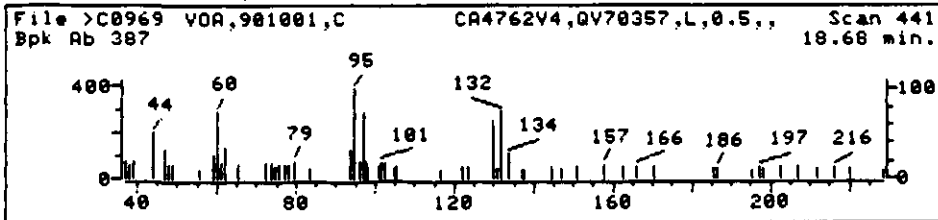
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



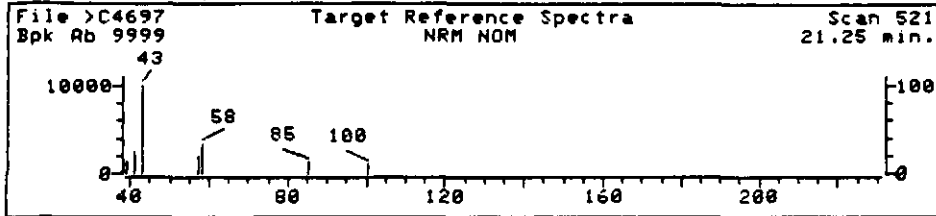
SAMPLE SPECTRUM (UNALTERED)



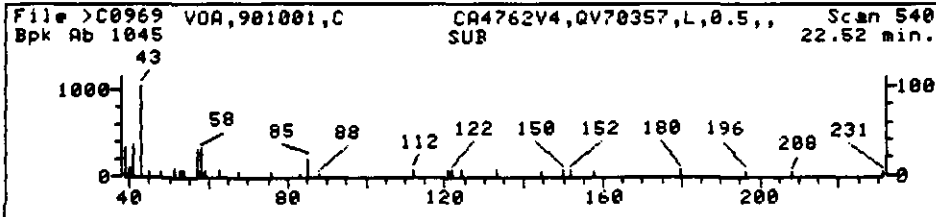
Data File: >C0969::U4 Quant Output File: ^C0969::AQ
 Name: VOA,901001,C
 Misc: CA4762V4,QV70357,L,0.5,,
 Quant Time: 901002 22:15 Quant ID File: IC1171::US
 Injected at: 901002 21:37 Last Calibration: 901002 16:53

Compound No: 28
 Compound Name: Trichloroethylene
 Scan Number: 441
 Retention Time: 18.68 min.
 Quant Ion: 130.0
 Area: 4193
 Concentration: 18.63 NG
 q-value: 75

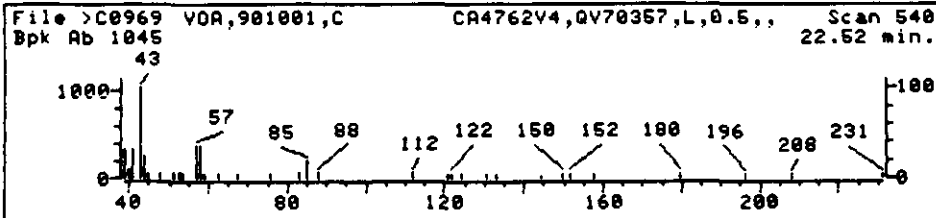
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



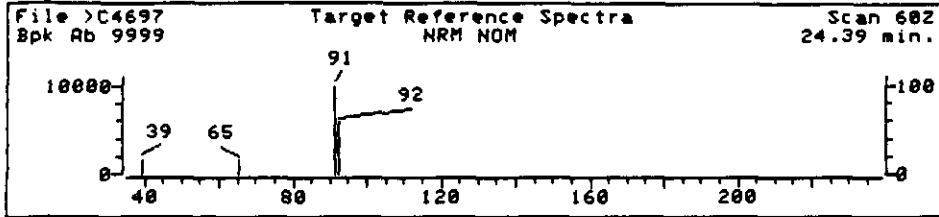
SAMPLE SPECTRUM (UNALTERED)



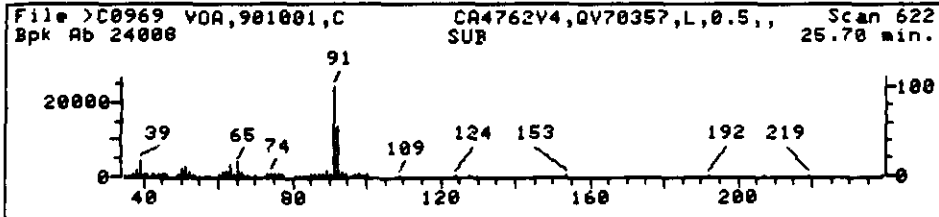
Data File: >C0969::U4 Quant Output File: ^C0969::AQ
 Name: VOA,901001,C
 Misc: CA4762V4,QV70357,L,0.5,,
 Quant Time: 901002 22:15 Quant ID File: IC1171::US
 Injected at: 901002 21:37 Last Calibration: 901002 16:53

Compound No: 37
 Compound Name: Methyl-iso-butyl ketone
 Scan Number: 540
 Retention Time: 22.52 min.
 Quant Ion: 43.0
 Area: 13019
 Concentration: 52.96 NG
 q-value: 89

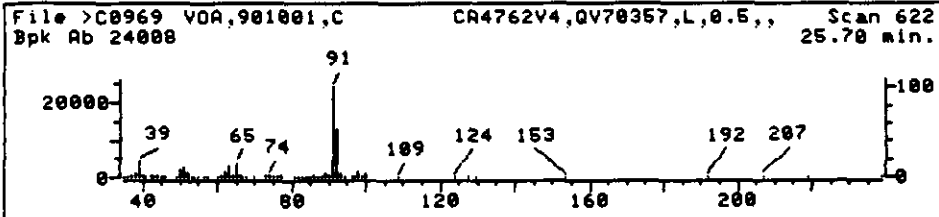
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

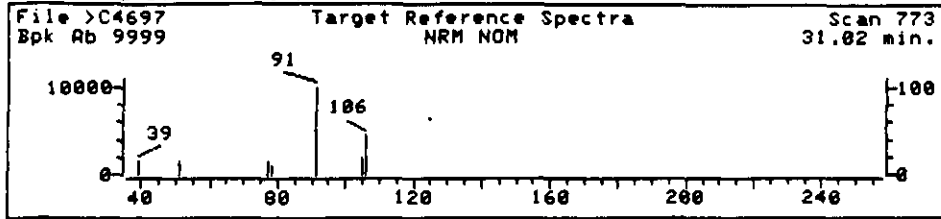


Data File: >C0969::U4
Name: VOA,901001,C
Misc: CA4762V4,QV70357,L,0.5,,
Quant Time: 901002 22:15
Injected at: 901002 21:37

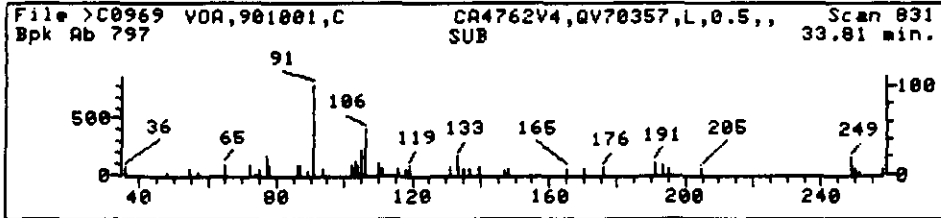
Quant Output File: ^C0969::AQ
Quant ID File: IC1171::US
Last Calibration: 901002 16:53

Compound No: 42
Compound Name: Toluene
Scan Number: 622
Retention Time: 25.70 min.
Quant Ion: 92.0
Area: 168759
Concentration: 283.03 NG
q-value: 96

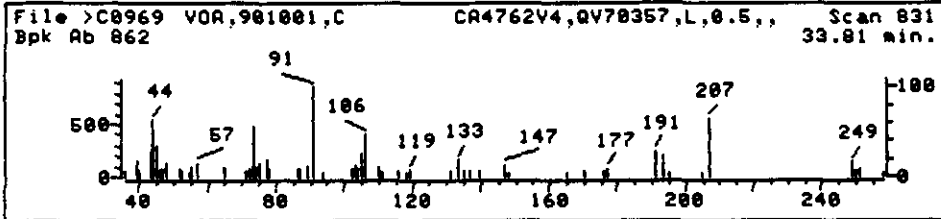
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

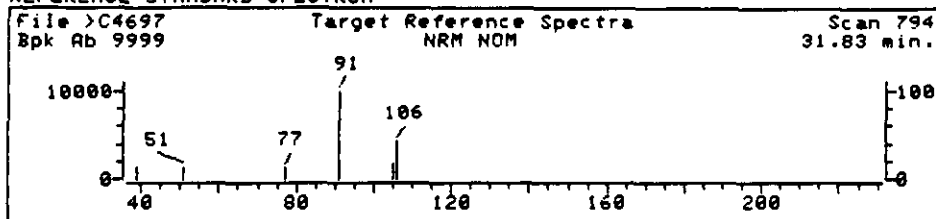


Data File: >C0969::U4
 Name: VOA,901001,C
 Misc: CA4762V4,QV70357,L,0.5,,
 Quant Time: 901002 22:15
 Injected at: 901002 21:37

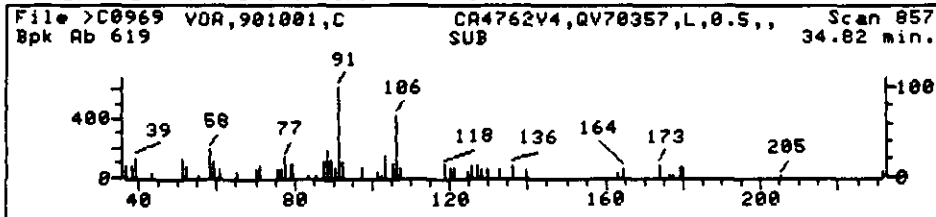
Quant Output File: ^C0969::AQ
 Quant ID File: IC1171::US
 Last Calibration: 901002 16:53

Compound No: 47
 Compound Name: m-Xylene
 Scan Number: 831
 Retention Time: 33.81 min.
 Quant Ion: 106.0
 Area: 10939
 Concentration: 16.63 NG
 q-value: 95

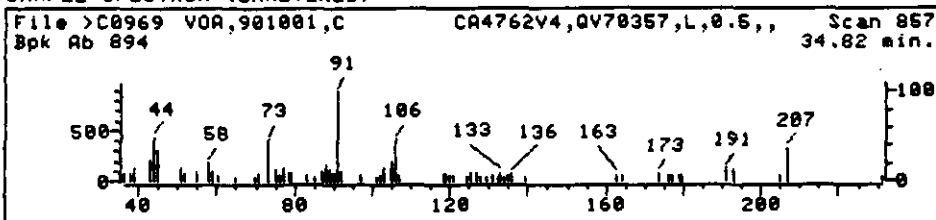
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



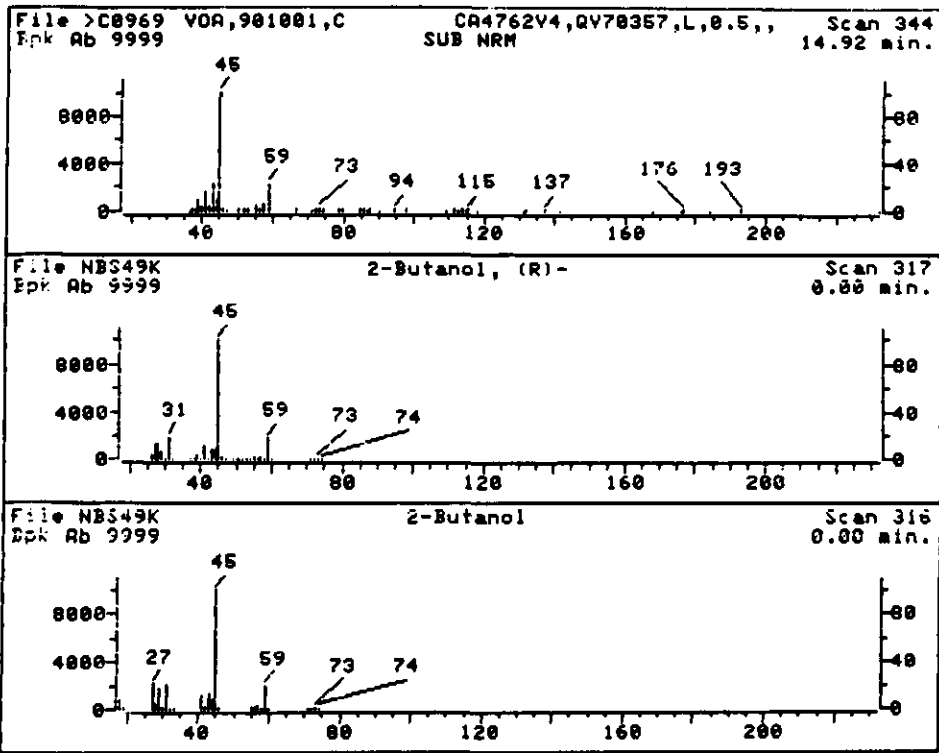
SAMPLE SPECTRUM (UNALTERED)



Data File: >C0969::U4
Name: VOA,901001,C
Misc: CA4762V4,QV70357,L,0.5,,
Quant Time: 901002 22:15
Injected at: 901002 21:37

Quant Output File: ^C0969::AQ
Quant ID File: IC1171::US
Last Calibration: 901002 16:53

Compound No: 48
Compound Name: o+p-Xylenes
Scan Number: 857
Retention Time: 34.82 min.
Quant Ion: 106.0
Area: 11268
Concentration: 16.59 NG
q-value: 82

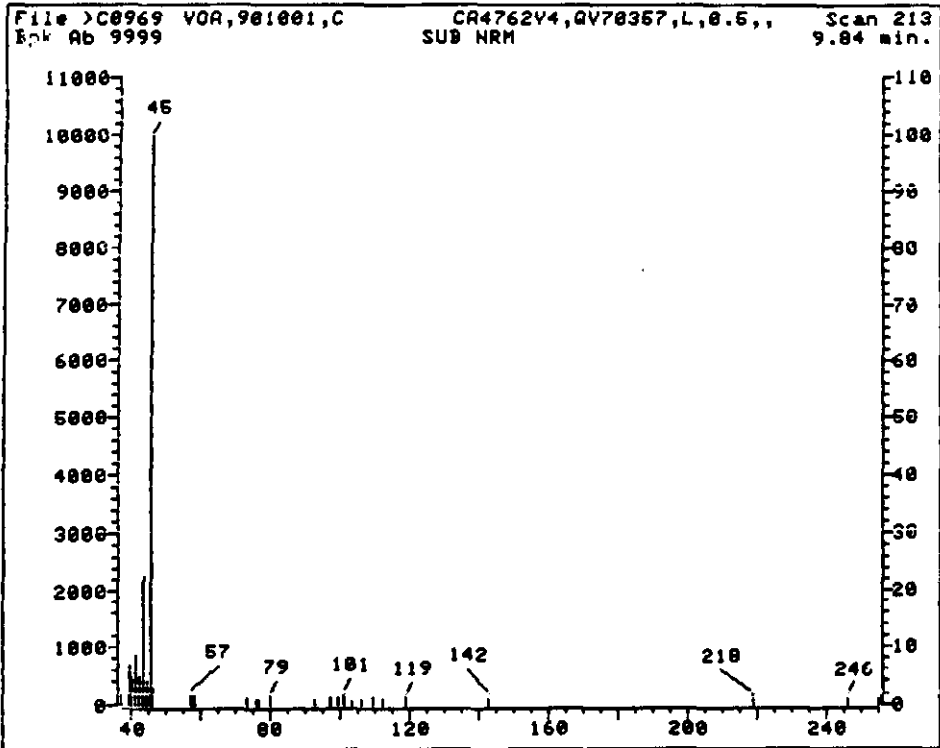


Data File: >C0969::U4
 Name: VOA,901001,C
 Misc Data: CA4762U4,QU70357,L,0.5,,
 Ri (min): 14.92
 Scan: 344
 Area: 245480 Rank: 7
 Semi-quantitative Conc (uncorrected): 219.74 NG
 Semi-quantitative Conc (corrected): 439.48 ug/l
 Calculated using Istd: Bromochloromethane @ 11.00 minutes

- 1. 2-Butanol, (R)- 74 C4H10O
- 2. 2-Butanol 74 C4H10O

Sample file: >C0969 Spectrum #: 344
 Search speed: 2 Tilting option: S No. of ion ranges searched: 45

| | Prob. | CAS # | CON # | ROUT | K | DK | #FLG | TILT | % | CON | C_I | R_IV |
|----|-------|----------|-------|--------|----|----|------|------|-----|-----|-----|------|
| 1. | 60 | 14898794 | 1716 | NBS49K | 36 | 44 | 0 | 0 | 84 | 14 | 30 | 19 |
| 2. | 52 | 78922 | 1715 | NBS49K | 43 | 40 | 2 | 0 | 100 | 16 | 20 | 14 |



Data File: >C0969::U4
 Name: VOA,901001,C
 Misc Data: CA4762V4,QV70357,L,0.5,,
 R: (min): 9.84
 Scan: 213
 Area: 88566 Rank: 8
 Semi-quantitative Conc (uncorrected): 79.28 NG
 Semi-quantitative Conc (corrected): 158.56 ug/l
 Calculated using Istd: Bromochloromethane @ 11.00 minutes

No PBM hits for this scan.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4762DL

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4762U5

Sample wt/vol: .0 (g/mL) ML

Lab File ID: >C0985

Level: (low/med) LOW

Date Received: 09/26/90

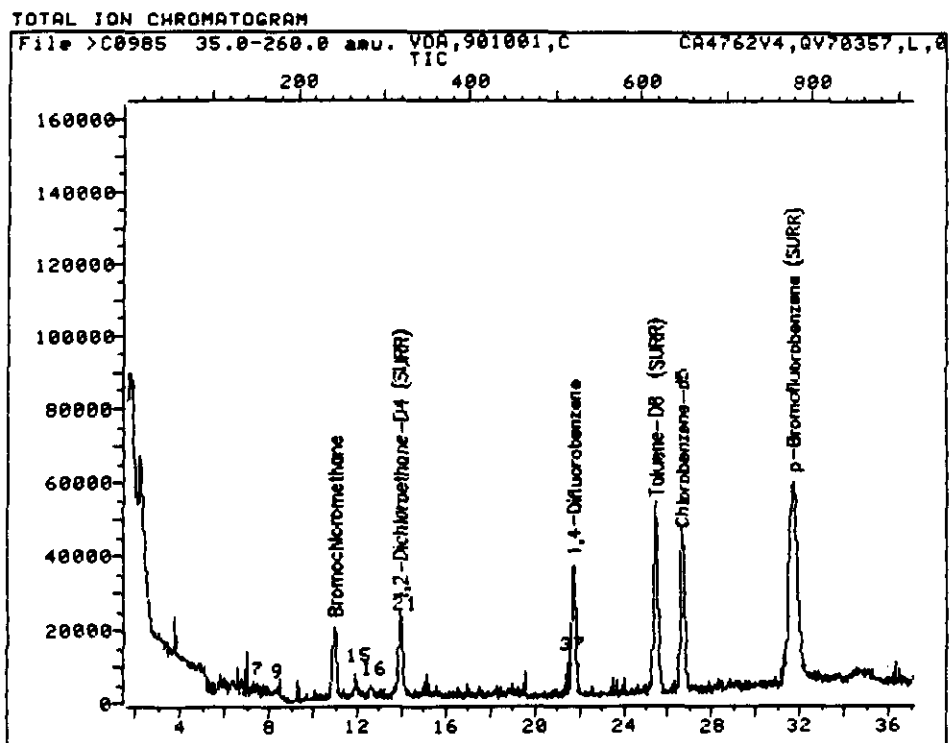
% Moisture: not dec.

Date Analyzed: 10/03/90

Column: (pack/cap) PACK

Dilution Factor: 200

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|----|
| 74-87-3 | Chloromethane | 2000 | UD |
| 74-83-9 | Bromomethane | 2000 | UD |
| 75-01-4 | Vinyl Chloride | 2000 | UD |
| 75-00-3 | Chloroethane | 2000 | UD |
| 75-09-2 | Methylene Chloride | 1000 | UD |
| 67-64-1 | Acetone | 5600 | ID |
| 75-15-0 | Carbon Disulfide | 1000 | UD |
| 75-35-4 | 1,1-Dichloroethene | 1000 | UD |
| 75-34-3 | 1,1-Dichloroethane | 1000 | UD |
| 540-59-0 | 1,2-Dichloroethene (total) | 1400 | 0 |
| 67-66-3 | Chloroform | 1000 | UD |
| 107-06-2 | 1,2-Dichloroethane | 1000 | UD |
| 78-93-3 | 2-Butanone | 16000 | ID |
| 71-55-6 | 1,1,1-Trichloroethane | 1000 | UD |
| 56-23-5 | Carbon Tetrachloride | 1000 | UD |
| 108-05-4 | Vinyl Acetate | 2000 | UD |
| 75-27-4 | Bromodichloromethane | 1000 | UD |
| 78-87-5 | 1,2-Dichloropropane | 1000 | UD |
| 10061-01-5 | cis-1,3-Dichloropropene | 1000 | UD |
| 79-01-6 | Trichloroethene | 1000 | UD |
| 124-48-1 | Dibromochloromethane | 1000 | UD |
| 79-00-5 | 1,1,2-Trichloroethane | 1000 | UD |
| 71-43-2 | Benzene | 1000 | UD |
| 10061-02-6 | trans-1,3-Dichloropropene | 1000 | UD |
| 75-25-2 | Bromoform | 1000 | UD |
| 108-10-1 | 4-Methyl-2-Pentanone | 2000 | UD |
| 591-78-6 | 2-Hexanone | 2000 | UD |
| 127-18-4 | Tetrachloroethene | 1000 | UD |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1000 | UD |
| 108-88-3 | Toluene | 630 | UD |
| 108-90-7 | Chlorobenzene | 1000 | UD |
| 100-41-4 | Ethylbenzene | 1000 | UD |
| 100-42-5 | Styrene | 1000 | UD |
| 1330-20-7 | Xylene (total) | 1000 | UD |



Data File: >C0985::U4
Name: VOA,901001,C
Misc: CA4762V4,QU70357,L,0.025,,

Quant Output File: ^C0985::AQ

Id File: IC1172::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901003 17:09

Operator ID: PT1575
Quant Time: 901003 22:11
Injected at: 901003 21:33

QUANT REPORT

Operator ID: PT1575 Quant Rev: 7 Quant Time: 901003 22:11
 Output File: ^C0985::AQ Injected at: 901003 21:33
 Data File: >C0985::U4 Dilution Factor: 1.00000
 Name: A4762DL
 Misc: CA4762U5,QU70357,L:M4,0.025,, 1:200 DL

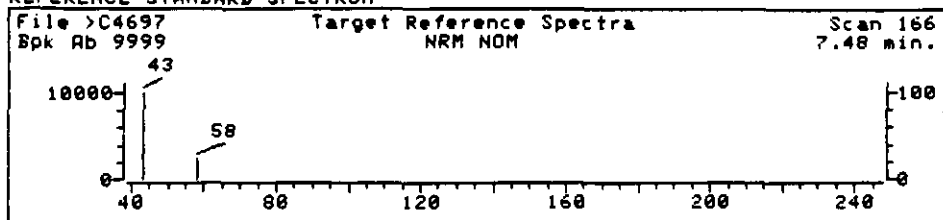
ID File: IC1172::US
 Title: IFB, PP/VOA, TCL, XUOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|-----|
| 1) *Bromochloromethane | 10.97 | 242 | 27092 | 250.00 | NG | 90 |
| 7) Methylene chloride | 7.40 | 150 | 3150 | 22.18 | NG | 98 |
| 9) Acetone | 8.29 | 173 | 11075 | 140.12 | NG | 80 |
| 15) Tetrahydrofuran | 11.94 | 267 | 7213 | 249.02 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.63 | 285 | 6113 | 34.88 | NG | 93 |
| 18) 1,2-Dichloroethane-04 (SURR) | 13.91 | 318 | 114687 | 298.74 | NG | 98 |
| 20) *1,4-Difluorobenzene | 21.75 | 520 | 132079 | 250.00 | NG | 93 |
| 21) Methyl ethyl ketone | 13.99 | 320 | 9552 | 403.76 | NG | 96 |
| 36) *Chlorobenzene-d5 | 26.64 | 646 | 149729 | 250.00 | NG | 74 |
| 37) Methyl-iso-butyl ketone | 21.48 | 513 | 2278 | 11.22 | NG | 72 |
| 41) Toluene-D8 (SURR) | 25.43 | 615 | 220625 | 249.16 | NG | 89 |
| 42) Toluene | 25.63 | 620 | 8017M | 15.69 | NG | |
| 45) p-Bromofluorobenzene (SURR) | 31.69 | 776 | 204308 | 302.72 | NG | 80 |

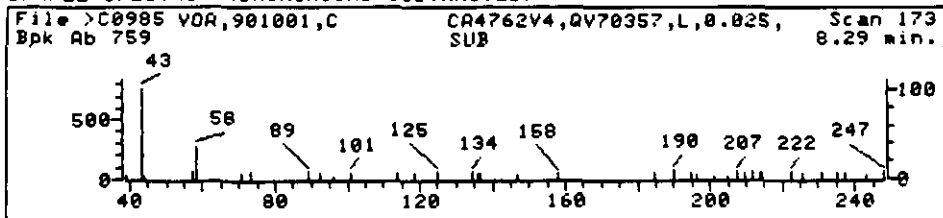
* Compound is ISTD

CA
 10-16-90.

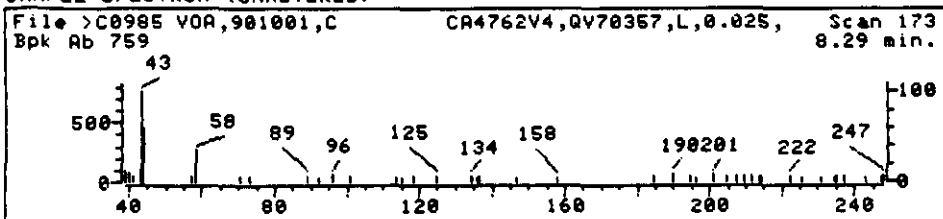
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



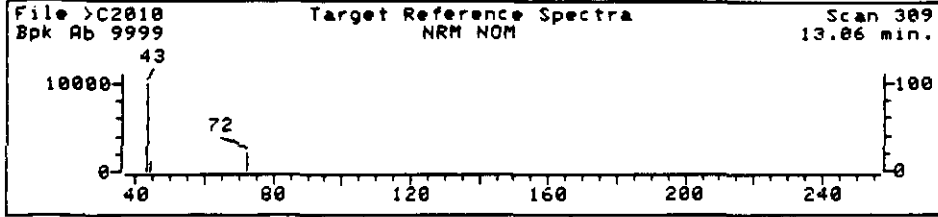
SAMPLE SPECTRUM (UNALTERED)



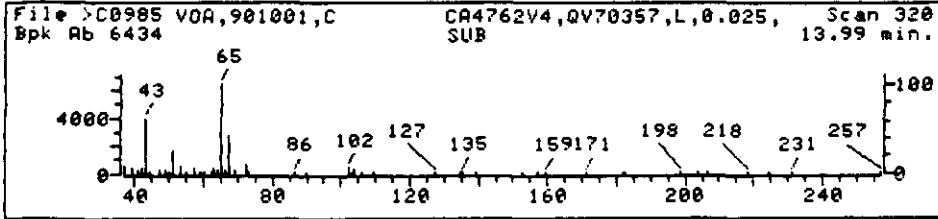
Data File: >C0985::U4 Quant Output File: ^C0985::AQ
 Name: UOA,901001,C
 Misc: CA4762V4,QV70357,L,0.025,,
 Quant Time: 901003 22:11 Quant ID File: IC1172::US
 Injected at: 901003 21:33 Last Calibration: 901003 17:09

Compound No: 9
 Compound Name: Acetone
 Scan Number: 173
 Retention Time: 8.29 min.
 Quant Ion: 43.0
 Area: 11075
 Concentration: 140.12 NG
 q-value: 80

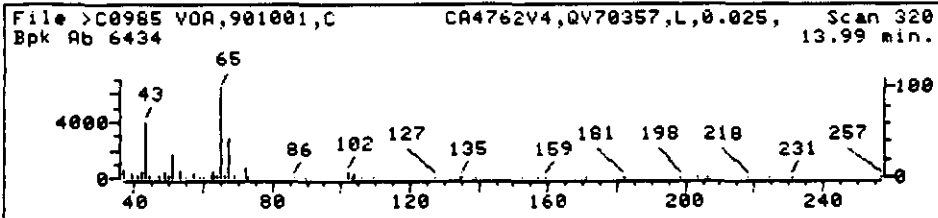
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



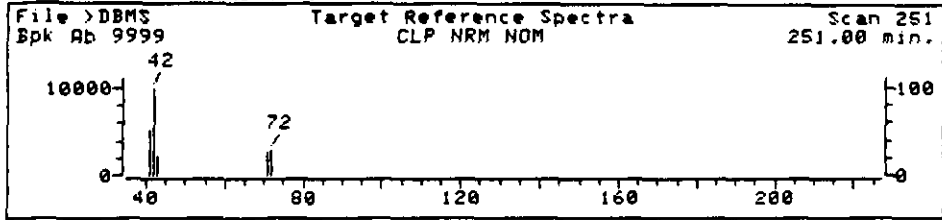
SAMPLE SPECTRUM (UNALTERED)



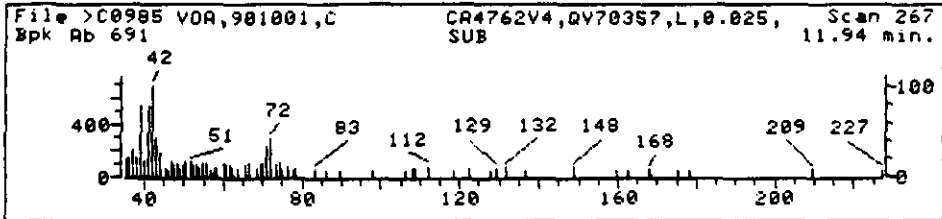
Data File: >C0985::U4 Quant Output File: ^C0985::AQ
 Name: VOA,901001,C
 Misc: CA4762U4,QV70357,L,0.025,,
 Quant Time: 901003 22:11 Quant ID File: IC1172::US
 Injected at: 901003 21:33 Last Calibration: 901003 17:09

Compound No: 21
 Compound Name: Methyl ethyl ketone
 Scan Number: 320
 Retention Time: 13.99 min.
 Quant Ion: 72.0
 Area: 9552
 Concentration: 403.76 NG
 q-value: 96

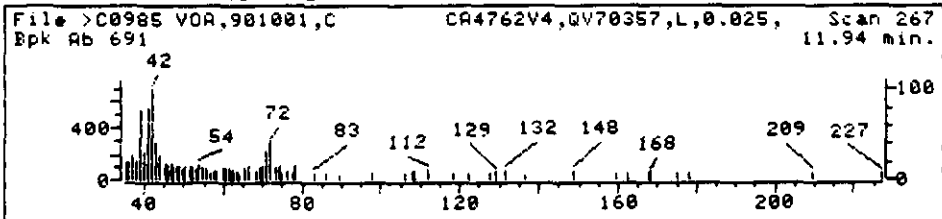
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >C0985::U4 Quant Output File: ^C0985::AQ
Name: UOA,901001,C
Misc: CA4762V4,QV70357,L,0.025,,
Quant Time: 901003 22:11 Quant ID File: IC1172::US
Injected at: 901003 21:33 Last Calibration: 901003 17:09

Compound No: 15
Compound Name: Tetrahydrofuran
Scan Number: 267
Retention Time: 11.94 min.
Quant Ion: 42.0
Area: 7213
Concentration: 249.02 NG
q-value: 100

ETC

STANDARDS DATA

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:ETCNI

Contract:

Lab Code:F2

Case No.:

SAS No.:

SDG No.:

Instrument ID:GC/MS C

Calibration Date(s) 09/26/90

09/26/90

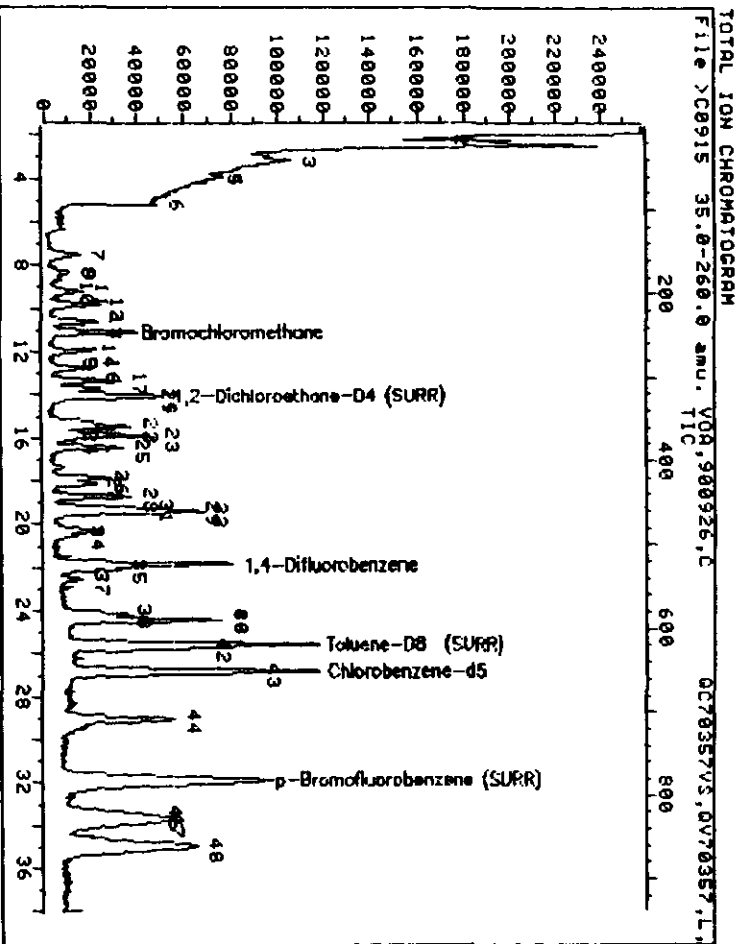
Matrix:(soil/water) WATER

Level:(low/med) LOW

Column:(pack/cap) PACK

Min RRF for SPCC(*) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

| LAB FILE ID: | RRF20 =>C0915 | RRF50 =>C0914 | | | | | |
|----------------------------|---------------|---------------|--------|--------|--------|-------|-------|
| RRF100=>C0913 | RRF150=>C0912 | RRF200=>C0911 | | | | | |
| COMPOUND | RRF20 | RRF50 | RRF100 | RRF150 | RRF200 | RRF | % RSD |
| Chloromethane | .746 | .648 | .619 | .587 | .587 | .637 | 10.3# |
| Bromomethane | .474 | .298 | .216 | .308 | .148 | .289 | 42.4 |
| Vinyl Chloride | 1.598 | 1.187 | 1.061 | 1.282 | 1.057 | 1.237 | 18.0* |
| Chloroethane | 1.014 | .827 | .797 | .869 | .967 | .895 | 10.3 |
| Methylene Chloride | 1.913 | 1.631 | 1.621 | 1.579 | 1.529 | 1.655 | 9.1 |
| Acetone | .606 | .631 | .561 | .612 | .619 | .606 | 4.5 |
| Carbon Disulfide | 5.121 | 4.577 | 4.609 | 4.680 | 4.514 | 4.700 | 5.2 |
| 1,1-Dichloroethene | 1.669 | 1.462 | 1.440 | 1.456 | 1.375 | 1.480 | 7.5* |
| 1,1-Dichloroethane | 3.113 | 2.697 | 2.675 | 2.710 | 2.570 | 2.753 | 7.6# |
| 1,2-Dichloroethene (total) | 1.677 | 1.434 | 1.409 | 1.491 | 1.356 | 1.473 | 8.4 |
| Chloroform | 5.511 | 4.748 | 4.641 | 4.702 | 4.508 | 4.822 | 8.2* |
| 1,2-Dichloroethane | 4.718 | 4.074 | 3.899 | 3.949 | 3.856 | 4.099 | 8.7 |
| 2-Butanone | .027 | .024 | .023 | .023 | .021 | .024 | 8.0 |
| 1,1,1-Trichloroethane | 1.082 | .953 | .915 | .932 | .979 | .972 | 6.8 |
| Carbon Tetrachloride | 1.070 | .936 | .908 | .933 | .993 | .968 | 6.7 |
| Vinyl Acetate | .505 | .461 | .405 | .461 | .463 | .459 | 7.8 |
| Bromodichloromethane | .901 | .801 | .773 | .795 | .806 | .815 | 6.1 |
| 1,2-Dichloropropane | .283 | .253 | .248 | .249 | .238 | .254 | 6.7* |
| cis-1,3-Dichloropropene | .526 | .479 | .462 | .464 | .462 | .478 | 5.7 |
| Trichloroethene | .482 | .420 | .454 | .413 | .414 | .437 | 7.0 |
| Dibromochloromethane | .666 | .619 | .598 | .614 | .628 | .625 | 4.1 |
| 1,1,2-Trichloroethane | .345 | .312 | .308 | .318 | .321 | .321 | 4.5 |
| Benzene | 1.003 | .898 | .939 | .978 | .952 | .954 | 4.2 |
| trans-1,3-Dichloropropene | .703 | .645 | .634 | .653 | .668 | .661 | 4.1 |
| Bromoform | .603 | .574 | .551 | .568 | .565 | .572 | 3.3# |
| 4-Methyl-2-Pentanone | .347 | .322 | .277 | .281 | .249 | .295 | 13.2 |
| 2-Hexanone | .280 | .278 | .233 | .246 | .227 | .253 | 9.9 |
| Tetrachloroethene | .537 | .469 | .447 | .448 | .446 | .470 | 8.3 |
| 1,1,2,2-Tetrachloroethane | .553 | .517 | .404 | .528 | .491 | .498 | 11.5# |
| Toluene | .878 | .789 | .784 | .796 | .780 | .805 | 5.1* |
| Chlorobenzene | 1.302 | 1.157 | 1.121 | 1.134 | 1.111 | 1.165 | 6.7# |
| Ethylbenzene | .715 | .644 | .630 | .647 | .641 | .655 | 5.2* |
| Styrene | 1.776 | 1.594 | 1.560 | 1.622 | 1.641 | 1.639 | 5.0 |
| Xylene (total) | .983 | .892 | .878 | .916 | .936 | .921 | 4.4 |
| Toluene-d8 | 1.475 | 1.457 | 1.461 | 1.470 | 1.498 | 1.472 | 1.1 |
| Bromofluorobenzene | 1.077 | 1.078 | 1.070 | 1.085 | 1.130 | 1.088 | 2.2 |
| 1,2-Dichloroethane-d4 | 3.407 | 3.393 | 3.454 | 3.586 | 3.610 | 3.490 | 2.9 |



Data File: >C0915::U0

Quant Output File: >C0915::AQ

Name: VOA,900926,C

Misc: QC70357V5, QV70357, L, S, ,

2UL CAL 1, 11, XVOA

ID File: IC1171::US

Title: IF8, PP/VOA, TCL, XVOA13

Last Calibration: 900926 15:42

Operator ID: MGRMS

Quant Time: 900926 15:57

Injected at: 900926 14:41

QUANT REPORT

Page 2

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900926 15:57
 Output File: ^C0915::AQ Injected at: 900926 14:41
 Data File: >C0915::U0 Dilution Factor: 1.00000
 Name: UOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 2UL CAL I,II,XVOA

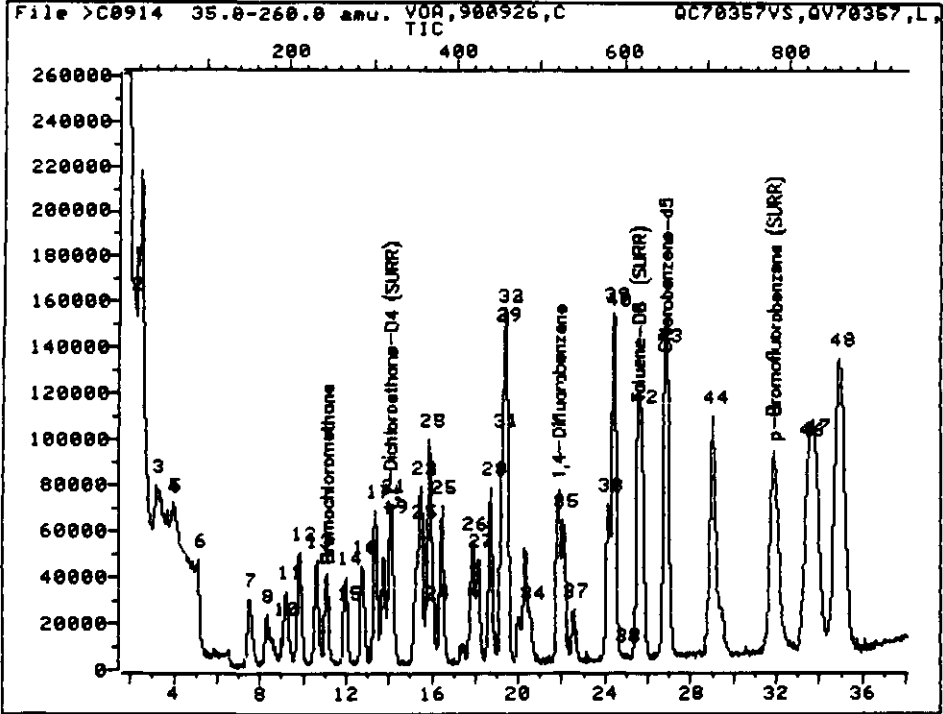
ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 43) | Chlorobenzene | 26.90 | 652 | 170050 | 92.20 | NG | 97 |
| 44) | Ethylbenzene | 29.04 | 707 | 93449 | 27.15 | NG | 79 |
| 45) | p-Bromofluorobenzene (SURR) | 31.84 | 779 | 351619 | 263.74 | NG | 85 |
| 46) | Styrene | 33.51 | 822 | 231937 | 85.74 | NG | 89 |
| 47) | m-Xylene | 33.82 | 830 | 125328 | 37.63 | NG | 90 |
| 48) | o+p-Xylenes | 34.98 | 860 | 256790 | 40.41 | NG | 88 |

* Compound is ISTD

EAK 10/15/90

TOTAL ION CHROMATOGRAM



Data File: >C0914::UO Quant Output File: ^C0914::AQ
 Name: UOA,900926,C
 Misc: QC70357US,QU70357,L,5,, SUL CAL I,II,XUOA

Id File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900926 16:16

Operator ID: MGRMS
 Quant Time: 900926 16:20
 Injected at: 900926 13:52

QUANT REPORT

Operator ID: MGRMS
 Output File: ^C0914::AQ
 Data File: >C0914::U0
 Name: UOA,900926,C

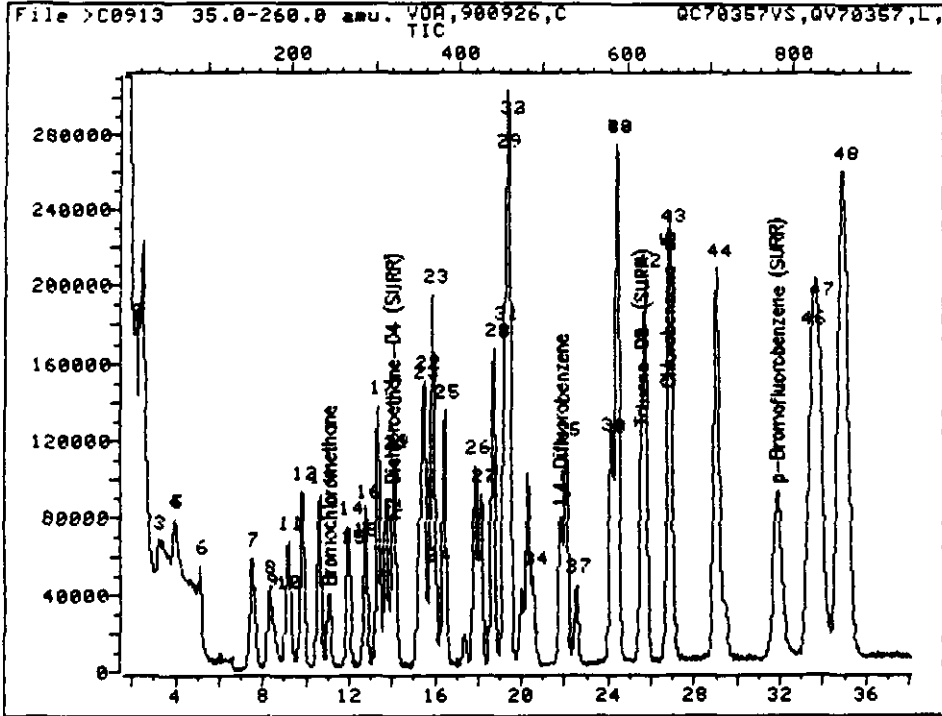
Quant Rev: 7 Quant Time: 900926 16:20
 Injected at: 900926 13:52
 Dilution Factor: 1.00000

Misc: QC70357US,QU70357,L,5,, 5UL CAL I,II,XVOA

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 16:16

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|---------------------------------------|------------------|----------------|------------------|-------------------|---------------|----------------|
| 1) *Bromochloromethane | 11.09 | 245 | 70232 | 250.00 | NG | 97 |
| 2) Methyl chloride | 2.16 | 15 | 45532 | 254.27 | NG | 98 |
| 3) Methyl bromide | 3.17 | 41 | 20954 | 258.19 | NG | 88 |
| 4) Dichlorodifluoromethane | 3.91 | 60 | 52592 | 233.25 | NG | 96 |
| 5) Vinyl chloride | 3.99 | 62 | 83361 | 239.87 | NG | 95 |
| 6) Chloroethane | 5.15 | 92 | 58101 | 231.15 | NG | 99 |
| 7) Methylene chloride | 7.48 | 152 | 114536 | 246.38 | NG | 95 |
| 8) Acrolein | 8.30 | 173 | 139391 | 3727.47 | NG | 86 |
| 9) Acetone | 8.33 | 174 | 44351 | 260.62 | NG | 82 |
| 10) Acrylonitrile | 9.07 | 193 | 29308 | 401.95 | NG | 83 |
| 11) Carbon disulfide | 9.23 | 197 | 321471 | 243.46 | NG | 99 |
| 12) Trichlorofluoromethane | 9.85 | 213 | 290849 | 241.86 | NG | 93 |
| 13) 1,1-Dichloroethylene | 10.62 | 233 | 102651 | 246.84 | NG | 91 |
| 14) 1,1-Dichloroethane | 11.98 | 268 | 189423 | 244.94 | NG | 94 |
| 14) 1,1-Dichloroethane | 12.76 | 288 | 54500 | 78.48 | NG | 58 |
| 15) Tetrahydrofuran | 12.06 | 270 | 12343 | 248.30 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.76 | 288 | 100678 | 243.26 | NG | 96 |
| 17) Chloroform | 13.34 | 303 | 333431 | 246.15 | NG | 98 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.04 | 321 | 238291 | 243.05 | NG | 97 |
| 19) 1,2-Dichloroethane | 14.16 | 324 | 286158 | 248.48 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.84 | 522 | 342312 | 250.00 | NG | 98 |
| 21) Methyl ethyl ketone | 14.12 | 323 | 8313 | 256.28 | NG | 87 |
| 22) 1,1,1-Trichloroethane | 15.48 | 358 | 326236 | 245.09 | NG | 95 |
| 23) Carbon tetrachloride | 15.48 | 358 | 34127 | 25.75 | NG | 97 |
| 23) Carbon tetrachloride | 15.90 | 369 | 320401 | 241.74 | NG | 98 |
| 24) Vinyl acetate | 16.10 | 374 | 157868 | 251.11 | NG | 95 |
| 25) Dichlorobromomethane | 15.55 | 368 | 2699 | 2.42 | NG | 86 |
| 25) Dichlorobromomethane | 15.98 | 369 | 4779 | 4.28 | NG | 92 |
| 25) Dichlorobromomethane | 16.45 | 383 | 274289 | 245.70 | NG | 99 |
| 26) 1,2-Dichloropropane | 17.88 | 420 | 86444 | 248.47 | NG | 98 |
| 27) cis-1,3-Dichloropropylene | 18.11 | 426 | 163824 | 250.08 | NG | 96 |
| 28) Trichloroethylene | 18.70 | 441 | 143838 | 240.61 | NG | 91 |
| 29) Chlorodibromomethane | 19.36 | 458 | 212019 | 247.68 | NG | 98 |
| 30) bis(Chloromethyl)ether | 19.36 | 458 | 83203 | 245.15 | NG | 100 |
| 31) Benzene | 19.24 | 455 | 307266 | 235.22 | NG | 90 |
| 32) 1,1,2-Trichloroethane | 19.43 | 460 | 106714 | 242.96 | NG | 91 |
| 33) trans-1,3-Dichloropropylene | 19.43 | 460 | 220924 | 244.17 | NG | 90 |
| 34) 2-Chloroethylvinyl ether | 20.52 | 488 | 62156 | 256.96 | NG | 100 |
| 35) Bromoform | 22.07 | 528 | 196579 | 250.82 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.78 | 649 | 318334 | 250.00 | NG | 77 |
| 37) Methyl-iso-butyl ketone | 22.54 | 540 | 102645 | 272.99 | NG | 95 |
| 38) 2-Hexanone | 24.13 | 581 | 88566 | 275.20 | NG | 87 |
| 38) 2-Hexanone | 24.95 | 682 | 2369 | 7.36 | NG | 93 |

TOTAL ION CHROMATOGRAM



Data File: >C0913::U0 Quant Output File: ^C0913::AQ
 Name: VOA,900926,C
 Misc: QC70357VS,QV70357,L,5,, 10UL CAL I,II,XVOA

Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 15:42

Operator ID: MGRMS
 Quant Time: 900926 15:53
 Injected at: 900926 13:03

QUANT REPORT

Operator ID: MGRMS
 Output File: ^C0913::AQ
 Data File: >C0913::U0
 Name: UOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 10UL CAL I,II,XVOA

Quant Rev: 7
 Quant Time: 900926 15:53
 Injected at: 900926 13:03
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|---------------------------------------|------------------|----------------|-------------------|--------------------|---------------|----------------|
| 1) *Bromochloromethane | 11.05 | 244 | 68225 | 250.00 | NG | 99 |
| 2) Methyl chloride | 2.16 | 15 | 84438 | 477.26 | NG | 97 |
| 3) Methyl bromide | 3.21 | 42 | 29462 | 361.85 | NG | 90 |
| 4) Dichlorodifluoromethane | 3.95 | 61 | 95209 | 465.90 | NG | 97 |
| 5) Vinyl chloride | 3.99 | 62 | 144780 | 446.97 | NG | 97 |
| 6) Chloroethane | 5.15 | 92 | 108717 | 481.55 | NG | 97 |
| 7) Methylene chloride | 7.48 | 152 | 221249 | 497.13 | NG | 98 |
| 8) Acrolein | 8.30 | 173 | 265559 | 7844.73 | NG | 91 |
| 9) Acetone | 8.37 | 175 | 76485 | 443.82 | NG | 80 |
| 10) Acrylonitrile | 9.07 | 193 | 57919 | 813.74 | NG | 78 |
| 11) Carbon disulfide | 9.19 | 196 | 628910 | 503.48 | NG | 98 |
| 12) Trichlorofluoromethane | 9.81 | 212 | 550069 | 486.72 | NG | 95 |
| 13) 1,1-Dichloroethylene | 10.62 | 233 | 196535 | 492.73 | NG | 90 |
| 14) 1,1-Dichloroethane | 11.94 | 267 | 364980 | 470.16 | NG | 95 |
| 15) Tetrahydrofuran | 12.02 | 269 | 22793 | 12645.10 | NG | 100 |
| 15) Tetrahydrofuran | 12.68 | 286 | 3651 | 2825.58 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.72 | 287 | 192283 | 352.18 | NG | 96 |
| 17) Chloroform | 13.34 | 303 | 633220 | 639.36 | NG | 96 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.00 | 320 | 235624 | 338.36 | NG | 99 |
| 19) 1,2-Dichloroethane | 14.12 | 323 | 532039 | 595.89 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.80 | 521 | 340907 | 250.00 | NG | 99 |
| 21) Methyl ethyl ketone | 14.08 | 322 | 15681 | 60.98 | NG | 86 |
| 22) 1,1,1-Trichloroethane | 15.48 | 358 | 623521 | 586.57 | NG | 95 |
| 23) Carbon tetrachloride | 15.44 | 357 | 66844 | 78.86 | NG | 98 |
| 23) Carbon tetrachloride | 15.86 | 368 | 619317 | 656.56 | NG | 97 |
| 24) Vinyl acetate | 16.06 | 373 | 276192 | 353.33 | NG | 95 |
| 25) Dichlorobromomethane | 16.45 | 383 | 526976 | 581.86 | NG | 98 |
| 26) 1,2-Dichloropropane | 17.84 | 419 | 168757 | 310.51 | NG | 98 |
| 27) cis-1,3-Dichloropropylene | 18.12 | 426 | 315197 | 405.77 | NG | 94 |
| 28) Trichloroethylene | 18.70 | 441 | 309255 | 494.10 | NG | 88 |
| 29) Chlorodibromomethane | 19.32 | 457 | 408046 | 604.43 | NG | 92 |
| 30) bis(Chloromethyl)ether | 19.32 | 457 | 159828 | 1985.88 | NG | 100 |
| 31) Benzene | 19.20 | 454 | 640063 | 378.77 | NG | 91 |
| 32) 1,1,2-Trichloroethane | 19.44 | 460 | 209992 | 445.98 | NG | 88 |
| 33) trans-1,3-Dichloropropylene | 19.44 | 460 | 432174 | 531.71 | NG | 91 |
| 34) 2-Chloroethylvinyl ether | 20.52 | 488 | 118333 | 402.29 | NG | 100 |
| 35) Bromoform | 22.08 | 528 | 375897 | 972.35 | NG | 97 |
| 36) *Chlorobenzene-d5 | 26.74 | 648 | 322654 | 250.00 | NG | 78 |
| 37) Methyl-iso-butyl ketone | 22.54 | 540 | 178675 | 324.43 | NG | 93 |
| 38) 2-Hexanone | 24.09 | 580 | 150372 | 302.24 | NG | 88 |
| 39) 1,1,2,2-Tetrachloroethane | 24.40 | 588 | 260722 | 298.50 | NG | 97 |
| 40) Tetrachloroethylene | 24.40 | 588 | 288656 | 268.45 | NG | 99 |
| 41) Toluene-D8 (SURR) | 25.53 | 617 | 471320 | 229.73 | NG | 90 |

QUANT REPORT

Operator ID: MGRMS
Output File: ^C0913::AQ
Data File: >C0913::U0
Name: VDA,900926,C
Misc: QC70357US,QU70357,L,5,,

Quant Rev: 7 Quant Time: 900926 15:53
 Injected at: 900926 13:03
 Dilution Factor: 1.00000

10UL CAL I,II,XVDA

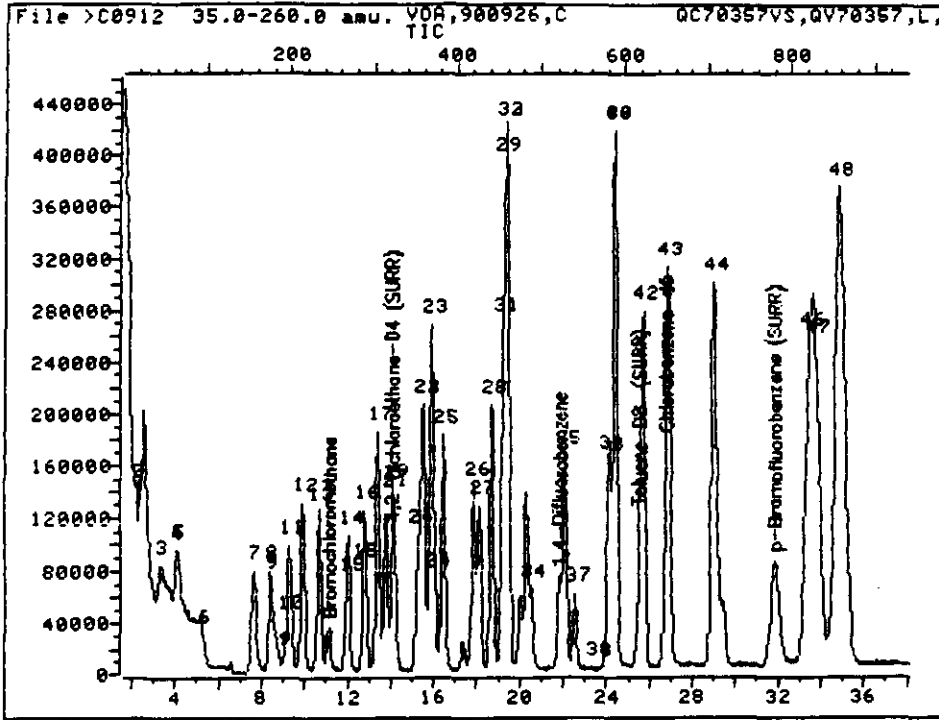
ID File: IC1171::US
Title: IFB, PP/VDA, TCL, XVDA13
Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|--------|-------|----|
| 42) | Toluene | 25.72 | 622 | 505696 | 204.72 | NG | 95 |
| 43) | Chlorobenzene | 26.90 | 652 | 723198 | 396.83 | NG | 98 |
| 44) | Ethylbenzene | 29.03 | 707 | 406333 | 119.47 | NG | 78 |
| 45) | p-Bromofluorobenzene (SURR) | 31.83 | 779 | 345303 | 262.12 | NG | 86 |
| 46) | Styrene | 33.46 | 821 | 1006763 | 376.65 | NG | 90 |
| 47) | m-Xylene | 33.77 | 829 | 541251 | 164.48 | NG | 91 |
| 48) | o+p-Xylenes | 34.94 | 859 | 1133798 | 180.56 | NG | 89 |

* Compound is ISTD

EMC N/15/90

TOTAL ION CHROMATOGRAM



Data File: >C0912::U0 Quant Output File: ^C0912::AQ
 Name: VOA,900926,C
 Misc: QC70357VS,QV70357,L,5,, 15UL CAL 1,11,XVOA

 Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 15:42

 Operator ID: MGRMS
 Quant Time: 900926 15:49
 Injected at: 900926 12:13

QUANT REPORT

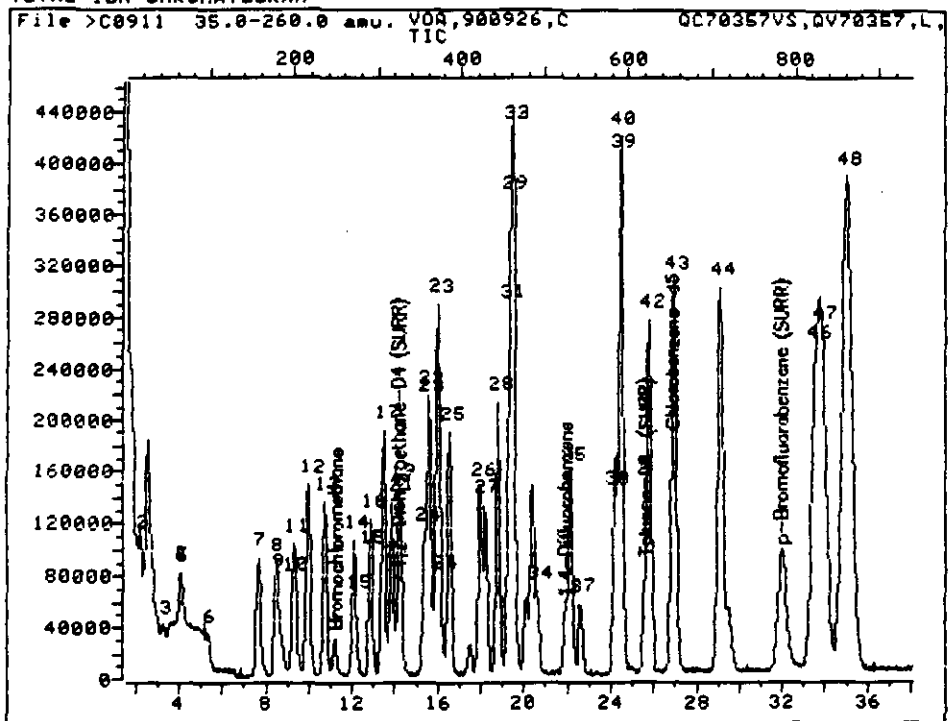
Operator ID: MGRMS
 Output File: ^C0912::AQ
 Data File: >C0912::U0
 Name: VOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 15UL CAL I,II,XVOA

Quant Rev: 7
 Quant Time: 900926 15:49
 Injected at: 900926 12:13
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|---------------------------------------|------------------|----------------|-------------------|--------------------|---------------|----------------|
| 1) *Bromochloromethane | 11.12 | 245 | 61517 | 250.00 | NG | 98 |
| 2) Methyl chloride | 2.23 | 16 | 108400 | 679.50 | NG | 97 |
| 3) Methyl bromide | 3.28 | 43 | 56817 | 773.91 | NG | 93 |
| 4) Dichlorodifluoromethane | 4.05 | 63 | 151377 | 821.52 | NG | 96 |
| 5) Vinyl chloride | 4.09 | 64 | 236659 | 810.29 | NG | 99 |
| 6) Chloroethane | 5.30 | 95 | 160295 | 787.44 | NG | 98 |
| 7) Methylene chloride | 7.63 | 155 | 291356 | 726.04 | NG | 97 |
| 8) Acrolein | 8.40 | 175 | 441660 | 14469.49 | NG | 89 |
| 9) Acetone | 8.44 | 176 | 112895 | 726.53 | NG | 86 |
| 9) Acetone | 9.02 | 191 | 2503 | 16.11 | NG | 22 |
| 10) Acrylonitrile | 9.14 | 194 | 77255 | 1203.76 | NG | 92 |
| 11) Carbon disulfide | 9.29 | 198 | 863774 | 766.90 | NG | 98 |
| 12) Trichlorofluoromethane | 9.92 | 214 | 753293 | 739.22 | NG | 95 |
| 13) 1,1-Dichloroethylene | 10.65 | 233 | 268737 | 747.21 | NG | 91 |
| 14) 1,1-Dichloroethane | 12.01 | 268 | 500082 | 714.45 | NG | 95 |
| 15) Tetrahydrofuran | 12.09 | 270 | 31236 | 19218.72 | NG | 100 |
| 15) Tetrahydrofuran | 12.71 | 286 | 6132 | 3772.07 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.79 | 288 | 275091 | 558.79 | NG | 97 |
| 17) Chloroform | 13.37 | 303 | 867721 | 971.67 | NG | 98 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.07 | 321 | 220621 | 351.37 | NG | 99 |
| 19) 1,2-Dichloroethane | 14.15 | 323 | 728878 | 905.38 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.83 | 521 | 307690 | 250.00 | NG | 99 |
| 21) Methyl ethyl ketone | 14.11 | 322 | 21496 | 92.62 | NG | 85 |
| 22) 1,1,1-Trichloroethane | 15.51 | 358 | 860583 | 896.98 | NG | 95 |
| 23) Carbon tetrachloride | 15.51 | 358 | 98300 | 106.07 | NG | 93 |
| 23) Carbon tetrachloride | 15.89 | 368 | 861224 | 1011.59 | NG | 96 |
| 24) Vinyl acetate | 15.27 | 352 | 198487 | 201.34 | NG | 75 |
| 24) Vinyl acetate | 16.09 | 373 | 425268 | 602.78 | NG | 96 |
| 25) Dichlorobromomethane | 16.44 | 382 | 733919 | 897.83 | NG | 97 |
| 26) 1,2-Dichloropropane | 17.83 | 418 | 229684 | 468.24 | NG | 99 |
| 27) cis-1,3-Dichloropropylene | 18.11 | 425 | 428246 | 610.82 | NG | 92 |
| 28) Trichloroethylene | 18.69 | 440 | 381133 | 674.68 | NG | 89 |
| 29) Chlorodibromomethane | 19.35 | 457 | 566804 | 930.24 | NG | 96 |
| 30) bis(Chloromethyl)ether | 19.35 | 457 | 222702 | 3065.02 | NG | 100 |
| 31) Benzene | 19.23 | 454 | 903092 | 592.12 | NG | 92 |
| 32) 1,1,2-Trichloroethane | 19.43 | 459 | 293592 | 690.84 | NG | 89 |
| 33) trans-1,3-Dichloropropylene | 19.43 | 459 | 603037 | 822.02 | NG | 93 |
| 34) 2-Chloroethylvinyl ether | 20.51 | 487 | 161877 | 609.74 | NG | 100 |
| 35) Bromoform | 22.07 | 527 | 524398 | 1502.93 | NG | 98 |
| 36) *Chlorobenzene-d5 | 26.74 | 647 | 298374 | 250.00 | NG | 77 |
| 37) Methyl-iso-butyl ketone | 22.53 | 539 | 251487 | 493.79 | NG | 94 |
| 38) 2-Hexanone | 23.50 | 564 | 2970 | 6.47 | NG | 87 |
| 38) 2-Hexanone | 24.09 | 579 | 220079 | 478.34 | NG | 91 |

TOTAL ION CHROMATOGRAM



Data File: >C0911::U0 Quant Output File: ^C0911::AQ
Name: VOA,900926,C
Misc: QC70357VS,QV70357,L,5,, 20UL CAL I,II,XVOA

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900926 15:42

Operator ID: MGRMS
Quant Time: 900926 15:43
Injected at: 900926 11:24

QUANT REPORT

Operator ID: MGRMS
 Output File: ^C0911::AQ
 Data File: >C0911::U0
 Name: UOA,900926,C

Quant Rev: 7 Quant Time: 900926 15:43
 Injected at: 900926 11:24
 Dilution Factor: 1.00000

Misc: QC70357VS,QU70357,L,5,, 20UL CAL I,II,XVOA

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|---------------------------------------|------------------|----------------|-------------------|--------------------|---------------|----------------|
| 1) *Bromochloromethane | 11.23 | 249 | 49370 | 250.00 | NG | 97 |
| 2) Methyl chloride | 2.25 | 18 | 115927 | 905.48 | NG | 98 |
| 3) Methyl bromide | 3.30 | 45 | 29185 | 495.34 | NG | 99 |
| 4) Dichlorodifluoromethane | 4.04 | 64 | 139319 | 942.11 | NG | 95 |
| 5) Vinyl chloride | 4.08 | 65 | 208786 | 890.74 | NG | 96 |
| 6) Chloroethane | 5.32 | 97 | 191040 | 1169.37 | NG | 98 |
| 7) Methylene chloride | 7.65 | 157 | 302039 | 937.85 | NG | 97 |
| 8) Acrolein | 8.47 | 178 | 453160 | 18499.02 | NG | 88 |
| 9) Acetone | 8.55 | 180 | 122291 | 980.63 | NG | 81 |
| 10) Acrylonitrile | 9.25 | 198 | 70170 | 1362.38 | NG | 84 |
| 11) Carbon disulfide | 9.32 | 200 | 891360 | 986.11 | NG | 99 |
| 12) Trichlorofluoromethane | 9.98 | 217 | 862656 | 1054.83 | NG | 93 |
| 13) 1,1-Dichloroethylene | 10.76 | 237 | 271510 | 940.66 | NG | 89 |
| 14) 1,1-Dichloroethane | 12.12 | 272 | 507563 | 903.54 | NG | 94 |
| 15) Tetrahydrofuran | 12.20 | 274 | 32334 | 24789.09 | NG | 100 |
| 15) Tetrahydrofuran | 12.82 | 298 | 9922 | 7686.77 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.90 | 292 | 267830 | 677.90 | NG | 99 |
| 17) Chloroform | 13.48 | 307 | 890236 | 1242.16 | NG | 98 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.14 | 324 | 178226 | 353.68 | NG | 96 |
| 19) 1,2-Dichloroethane | 14.26 | 327 | 761552 | 1178.71 | NG | 96 |
| 20) *1,4-Difluorobenzene | 21.90 | 524 | 232529 | 250.00 | NG | 99 |
| 21) Methyl ethyl ketone | 14.22 | 326 | 19882 | 113.35 | NG | 87 |
| 22) 1,1,1-Trichloroethane | 15.61 | 362 | 910803 | 1256.17 | NG | 94 |
| 23) Carbon tetrachloride | 15.58 | 361 | 88864 | 136.87 | NG | 97 |
| 23) Carbon tetrachloride | 16.00 | 372 | 923576 | 1435.47 | NG | 95 |
| 24) Vinyl acetate | 15.38 | 356 | 286336 | 387.08 | NG | 76 |
| 24) Vinyl acetate | 16.20 | 377 | 430938 | 808.26 | NG | 97 |
| 25) Dichlorobromomethane | 16.55 | 386 | 749680 | 1213.56 | NG | 96 |
| 26) 1,2-Dichloropropane | 17.98 | 423 | 221792 | 598.30 | NG | 95 |
| 27) cis-1,3-Dichloropropylene | 18.22 | 429 | 429536 | 810.69 | NG | 95 |
| 28) Trichloroethylene | 18.80 | 444 | 385137 | 902.14 | NG | 89 |
| 29) Chlorodibromomethane | 19.42 | 460 | 583657 | 1267.52 | NG | 98 |
| 30) bis(Chloromethyl)ether | 19.42 | 460 | 228974 | 4171.04 | NG | 100 |
| 31) Benzene | 19.34 | 458 | 885817 | 768.53 | NG | 91 |
| 32) 1,1,2-Trichloroethane | 19.54 | 463 | 298426 | 929.20 | NG | 89 |
| 33) trans-1,3-Dichloropropylene | 19.54 | 463 | 621566 | 1121.14 | NG | 90 |
| 34) 2-Chloroethylvinyl ether | 20.62 | 491 | 160736 | 801.14 | NG | 100 |
| 35) Bromoform | 22.14 | 530 | 525932 | 1994.55 | NG | 98 |
| 36) *Chlorobenzene-d5 | 26.85 | 650 | 228594 | 250.00 | NG | 81 |
| 37) Methyl-iso-butyl ketone | 22.60 | 542 | 227996 | 584.32 | NG | 91 |
| 38) 2-Hexanone | 24.16 | 582 | 207178 | 587.76 | NG | 86 |
| 39) 1,1,2,2-Tetrachloroethane | 24.47 | 590 | 449164 | 725.84 | NG | 97 |
| 40) Tetrachloroethylene | 24.51 | 591 | 408254 | 535.89 | NG | 97 |

QUANT REPORT

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900926 15:43
 Output File: ^C0911::AQ Injected at: 900926 11:24
 Data File: >C0911::U0 Dilution Factor: 1.00000
 Name: UOA,900926,C
 Misc: QC70357VS,QU70357,L,5,, 20UL CAL 1,11,XVOA

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|--------|-------|----|
| 41) | Toluene-D8 (SURR) | 25.59 | 619 | 342513 | 235.64 | NG | 92 |
| 42) | Toluene | 25.79 | 624 | 712783 | 407.28 | NG | 94 |
| 43) | Chlorobenzene | 26.97 | 653 | 1015905 | 786.82 | NG | 96 |
| 44) | Ethylbenzene | 29.10 | 708 | 586369 | 243.34 | NG | 77 |
| 45) | p-Bromofluorobenzene (SURR) | 31.94 | 781 | 258347 | 276.81 | NG | 82 |
| 46) | Styrene | 33.57 | 823 | 1500542 | 792.37 | NG | 92 |
| 47) | m-Xylene | 33.88 | 831 | 817849 | 350.81 | NG | 91 |
| 48) | o+p-Xylenes | 35.05 | 861 | 1711432 | 384.69 | NG | 88 |

* Compound is ISTD

EAL 10/15/90

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name:ETCNJ

Contract:

Lab Code:F2

Case No.:

SAS No.:

SDG No.:

Instrument ID:GC/MS C

Calibration Date: 09/27/90

Time: 1656

Lab File ID: >C0927

Init Calib. Dates(s): 09/26/90

09/26/90

Matrix:(soil/water) WATER

Level:(low/med) LOW

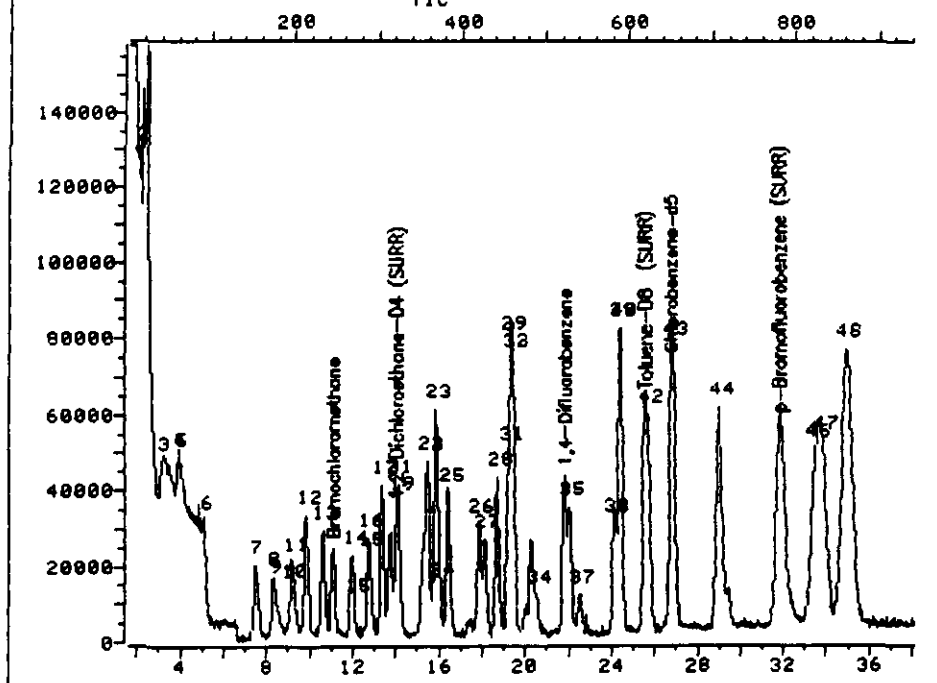
Column:(pack/cap) PACK

Min RRF50 for SPCC(#) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 25.0%

| COMPOUND | RRF | RRF50 | %D |
|---------------------------------|-------|-------|-------|
| Chloromethane_____# | .637 | 1.077 | 69.0# |
| Bromomethane_____ | .289 | .346 | 19.8 |
| Vinyl Chloride_____* | 1.237 | 1.461 | 18.1* |
| Chloroethane_____ | .895 | 1.032 | 15.4 |
| Methylene Chloride_____ | 1.655 | 1.661 | .4 |
| Acetone_____ | .606 | .713 | 17.7 |
| Carbon Disulfide_____ | 4.700 | 4.402 | 6.3 |
| 1,1-Dichloroethene_____* | 1.480 | 1.375 | 7.1* |
| 1,1-Dichloroethane_____* | 2.753 | 2.529 | 8.1# |
| 1,2-Dichloroethene (total)_____ | 1.473 | 1.335 | 9.4 |
| Chloroform_____* | 4.822 | 4.530 | 6.1* |
| 1,2-Dichloroethane_____ | 4.099 | 3.918 | 4.4 |
| 2-Butanone_____ | .024 | .024 | 2.2 |
| 1,1,1-Trichloroethene_____ | .972 | 1.081 | 11.2 |
| Carbon Tetrachloride_____ | .968 | 1.119 | 15.6 |
| Vinyl Acetate_____ | .459 | .421 | 8.3 |
| Bromodichloromethane_____ | .815 | .836 | 2.5 |
| 1,2-Dichloropropane_____* | .254 | .238 | 6.2* |
| cis-1,3-Dichloropropene_____ | .478 | .465 | 2.7 |
| Trichloroethene_____ | .437 | .442 | 1.3 |
| Dibromochloromethane_____ | .625 | .654 | 4.7 |
| 1,1,2-Trichloroethane_____ | .321 | .301 | 6.1 |
| Benzene_____ | .954 | .840 | 11.9 |
| trans-1,3-Dichloropropene_____ | .661 | .650 | 1.7 |
| Bromoform_____# | .572 | .584 | 2.1# |
| 4-Methyl-2-Pentanone_____ | .295 | .283 | 4.0 |
| 2-Hexanone_____ | .253 | .264 | 4.4 |
| Tetrachloroethene_____ | .470 | .491 | 4.6 |
| 1,1,2,2-Tetrachloroethane_____* | .498 | .437 | 12.3# |
| Toluene_____* | .805 | .749 | 7.0* |
| Chlorobenzene_____# | 1.165 | 1.161 | .3# |
| Ethylbenzene_____* | .655 | .671 | 2.4* |
| Styrene_____ | 1.639 | 1.583 | 3.4 |
| Xylene (total)_____ | .921 | .970 | 5.3 |
| Toluene-d8_____ | 1.472 | 1.460 | .8 |
| Bromofluorobenzene_____ | 1.088 | 1.129 | 3.8 |
| 1,2-Dichloroethane-d4_____ | 3.490 | 3.231 | 7.4 |

TOTAL ION CHROMATOGRAM

File >C0927 35.0-260.0 amu. VOA,900927,C DC70357VS,0V70357,L,
TIC



Data File: >C0927::U4

Quant Output File: ^C0927::AQ

Name: VOA,900927,C

Misc: QC70357VS,QU70357,L,5,,

5UL CAL I,II,XVOA

Id File: IC1171::US

Title: IFB, PP/VOA, TCL, XVOA13

Last Calibration: 900926 16:16

Operator ID: MGRMS

Quant Time: 900927 17:35

Injected at: 900927 16:56

QUANT REPORT

Page 2

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900927 17:35
 Output File: ^C0927::AQ Injected at: 900927 16:56
 Data File: >C0927::U4 Dilution Factor: 1.00000
 Name: UOA,900927,C
 Misc: QC70357US,QV70357,L,5,, SUL CAL I,II,XUOA

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900926 16:16

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 42) | Toluene | 25.71 | 622 | 128494 | 232.51 | NG | 97 |
| 43) | Chlorobenzene | 26.89 | 652 | 199297 | 249.22 | NG | 98 |
| 44) | Ethylbenzene | 29.02 | 707 | 115137 | 255.93 | NG | 77 |
| 45) | p-Bromofluorobenzene (SURR) | 31.82 | 779 | 193769 | 259.47 | NG | 92 |
| 46) | Styrene | 33.53 | 823 | 271632 | 241.51 | NG | 85 |
| 47) | m-Xylene | 33.84 | 831 | 153657 | 253.53 | NG | 94 |
| 48) | o+p-Xylenes | 34.96 | 860 | 333036 | 526.71 | NG | 87 |

* Compound is ISTD

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:ETCNJ

Contract:

Lab Code:F2

Case No.:

SAS No.:

SDG No.:

Instrument ID:GC/MS C

Calibration Date(s) 10/02/90

10/02/90

Matrix:(soil/water) WATER

Level:(low/med) LOW

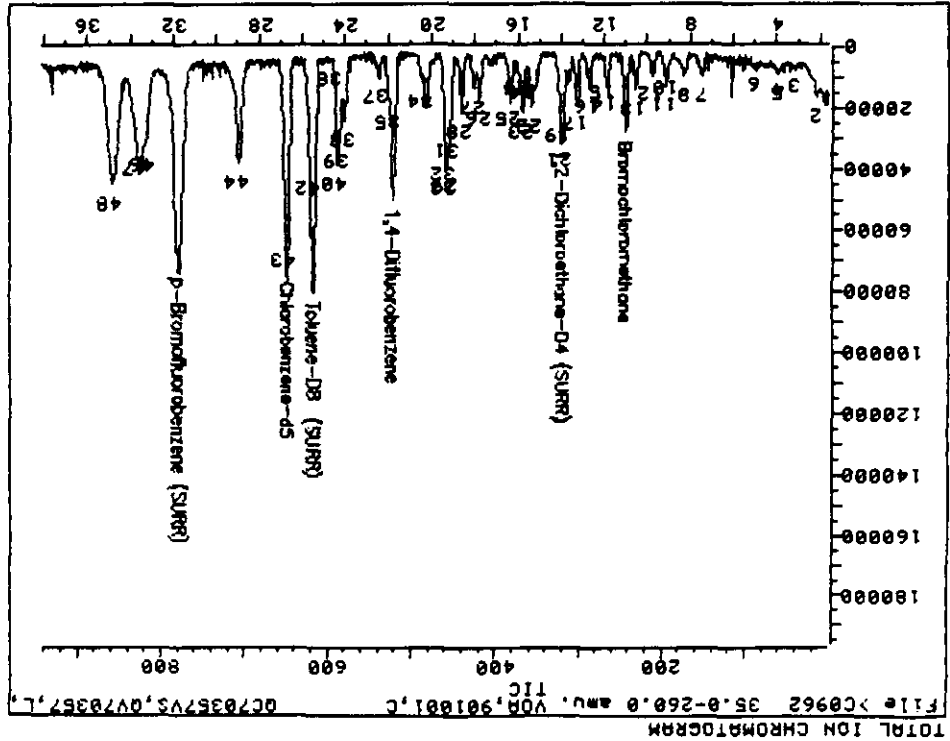
Column:(pack/cap) PACK

Min RRF for SPCC(#) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

| | | |
|----------------|---------------|---------------|
| LAB FILE ID: | RRF20 =>C0962 | RRF50 =>C0961 |
| IRRF100=>C0960 | RRF150=>C0959 | RRF200=>C0958 |

| COMPOUND | RRF20 | RRF50 | RRF100 | RRF150 | RRF200 | RRF | % RSD |
|----------------------------|-------|-------|--------|--------|--------|-------|-------|
| Chloromethane | 1.098 | .925 | .925 | .907 | 1.315 | 1.034 | 17.0 |
| Bromomethane | .270 | .171 | .127 | .119 | .136 | .165 | 37.8 |
| Vinyl Chloride | .860 | .684 | .567 | .529 | .702 | .669 | 19.5 |
| Chloroethane | .495 | .411 | .403 | .403 | .474 | .437 | 10.0 |
| Methylene Chloride | 1.490 | 1.165 | .884 | .941 | 1.059 | 1.108 | 21.6 |
| Acetone | .803 | .607 | .487 | .497 | .483 | .575 | 23.8 |
| Carbon Disulfide | 6.833 | 6.678 | 3.834 | 4.540 | 7.107 | 5.799 | 25.9 |
| 1,1-Dichloroethene | 1.649 | 1.622 | 1.672 | 1.694 | 1.825 | 1.692 | 4.7 |
| 1,1-Dichloroethane | 3.626 | 3.263 | 3.237 | 3.308 | 3.598 | 3.407 | 5.6 |
| 1,2-Dichloroethene (total) | 1.624 | 1.576 | 1.596 | 1.661 | 1.732 | 1.638 | 3.8 |
| Chloroform | 5.077 | 4.498 | 4.485 | 4.561 | 4.877 | 4.700 | 5.6 |
| 1,2-Dichloroethane | 4.047 | 3.642 | 3.430 | 3.418 | 3.660 | 3.639 | 7.0 |
| 2-Butanone | .038 | .038 | .034 | .034 | .029 | .035 | 10.8 |
| 1,1,1-Trichloroethane | .868 | .756 | .695 | .711 | .738 | .754 | 9.0 |
| Carbon Tetrachloride | .769 | .677 | .643 | .665 | .705 | .692 | 7.0 |
| Vinyl Acetate | .493 | .521 | .456 | .538 | .544 | .510 | 7.1 |
| Bromodichloromethane | .801 | .710 | .652 | .671 | .686 | .704 | 8.2 |
| 1,2-Dichloropropane | .386 | .343 | .330 | .342 | .346 | .349 | 6.1 |
| cis-1,3-Dichloropropene | .532 | .493 | .467 | .488 | .494 | .495 | 4.7 |
| Trichloroethene | .400 | .365 | .368 | .359 | .355 | .369 | 4.9 |
| Dibromochloromethane | .552 | .514 | .492 | .507 | .511 | .515 | 4.3 |
| 1,1,2-Trichloroethane | .346 | .323 | .322 | .326 | .330 | .329 | 3.1 |
| Benzene | 1.179 | 1.114 | 1.118 | 1.187 | 1.223 | 1.164 | 4.0 |
| trans-1,3-Dichloropropene | .713 | .650 | .603 | .627 | .633 | .645 | 6.4 |
| Bromoform | .477 | .439 | .417 | .416 | .383 | .426 | 8.1 |
| 4-Methyl-2-Pentanone | .380 | .373 | .344 | .320 | .292 | .342 | 10.7 |
| 2-Hexanone | .387 | .345 | .336 | .307 | .277 | .331 | 12.5 |
| Tetrachloroethene | .367 | .349 | .344 | .334 | .330 | .345 | 4.2 |
| 1,1,2,2-Tetrachloroethane | .534 | .544 | .488 | .525 | .505 | .519 | 4.3 |
| Toluene | .858 | .809 | .814 | .817 | .852 | .830 | 2.8 |
| Chlorobenzene | 1.226 | 1.170 | 1.128 | 1.121 | 1.162 | 1.161 | 3.6 |
| Ethylbenzene | .758 | .692 | .672 | .677 | .695 | .699 | 5.0 |
| Styrene | 1.823 | 1.663 | 1.578 | 1.542 | 1.625 | 1.646 | 6.6 |
| Xylene (total) | 1.050 | .948 | .901 | .886 | .940 | .945 | 6.8 |
| Toluene-d8 | 1.478 | 1.471 | 1.455 | 1.455 | 1.462 | 1.464 | .7 |
| Bromofluorobenzene | 1.082 | 1.038 | 1.002 | 1.003 | 1.045 | 1.034 | 3.2 |
| 1,2-Dichloroethane-d4 | 3.212 | 3.041 | 3.093 | 3.263 | 3.539 | 3.229 | 6.0 |

Data File: >C0962:U0
 Name: VDA,901001,C
 Misc: DC70357VS,QV70357,L,5,, 2UL
 Id File: IC1171:US
 Title: IFB, PP/UDA, TCL, XUD13
 Last Calibration: 901002 16:53
 Operator ID: JAB781
 Quant Time: 901002 17:06
 Injected at: 901002 15:32



QUANT REPORT

Page 1

Operator ID: JA8781
 Output File: ^C0962::AQ
 Data File: >C0962::U0
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5,, 2UL

Quant Rev: 7 Quant Time: 901002 17:06
 Injected at: 901002 15:32
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|---------|-------|-----|
| 1) *Bromochloromethane | 11.01 | 243 | 40870 | 250.00 | NG | 95 |
| 2) Methyl chloride | 2.16 | 15 | 17949 | 106.18 | NG | 98 |
| 3) Methyl bromide | 3.17 | 41 | 4412 | 164.03 | NG | 99 |
| 4) Dichlorodifluoromethane | 3.95 | 61 | 9787 | 128.58 | NG | 98 |
| 5) Vinyl chloride | 3.99 | 62 | 14067 | 128.70 | NG | 92 |
| 6) Chloroethane | 5.11 | 91 | 8097 | 113.28 | NG | 87 |
| 7) Methylene chloride | 7.48 | 152 | 24363 | 134.53 | NG | 99 |
| 8) Acrolein | 8.30 | 173 | 42501 | 2269.06 | NG | 97 |
| 9) Acetone | 8.30 | 173 | 13122 | 139.52 | NG | 99 |
| 10) Acrylonitrile | 9.03 | 192 | 16223 | 197.03 | NG | 96 |
| 11) Carbon disulfide | 9.15 | 195 | 111713 | 117.85 | NG | 99 |
| 12) Trichlorofluoromethane | 9.77 | 211 | 57950 | 98.14 | NG | 89 |
| 13) 1,1-Dichloroethylene | 10.54 | 231 | 26951 | 97.43 | NG | 92 |
| 14) 1,1-Dichloroethane | 11.90 | 266 | 59283 | 106.45 | NG | 96 |
| 15) Tetrahydrofuran | 11.98 | 268 | 3236 | 75.33 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.72 | 287 | 26552 | 99.17 | NG | 92 |
| 17) Chloroform | 13.26 | 301 | 82994 | 108.03 | NG | 98 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.96 | 319 | 131291 | 248.68 | NG | 95 |
| 19) 1,2-Dichloroethane | 14.07 | 322 | 66160 | 111.20 | NG | 96 |
| 20) *1,4-Difluorobenzene | 21.79 | 521 | 203524 | 250.00 | NG | 98 |
| 21) Methyl ethyl ketone | 14.03 | 321 | 3125 | 110.25 | NG | 91 |
| 22) 1,1,1-Trichloroethane | 15.39 | 356 | 70670 | 115.18 | NG | 91 |
| 23) Carbon tetrachloride | 15.39 | 356 | 7889 | 14.01 | NG | 83 |
| 23) Carbon tetrachloride | 15.82 | 367 | 62573 | 111.13 | NG | 99 |
| 24) Vinyl acetate | 15.66 | 363 | 24580 | 59.15 | NG | 80 |
| 24) Vinyl acetate | 16.01 | 372 | 40164 | 96.66 | NG | 97 |
| 25) Dichlorobromomethane | 16.40 | 382 | 65190 | 113.74 | NG | 98 |
| 26) 1,2-Dichloropropane | 17.80 | 418 | 31426 | 110.50 | NG | 94 |
| 27) cis-1,3-Dichloropropylene | 18.03 | 424 | 43309 | 107.51 | NG | 91 |
| 28) Trichloroethylene | 18.65 | 440 | 32596 | 108.40 | NG | 82 |
| 29) Chlorodibromomethane | 19.27 | 456 | 44953 | 107.21 | NG | 98 |
| 30) bis(Chloromethyl)ether | 19.23 | 455 | 19135 | 112.10 | NG | 100 |
| 31) Benzene | 19.15 | 453 | 95987 | 101.27 | NG | 96 |
| 32) 1,1,2-Trichloroethane | 19.39 | 459 | 28182 | 105.13 | NG | 82 |
| 33) trans-1,3-Dichloropropylene | 19.35 | 458 | 58007 | 110.46 | NG | 86 |
| 34) 2-Chloroethylvinyl ether | 20.47 | 487 | 21381 | 105.47 | NG | 100 |
| 35) Bromoform | 21.98 | 526 | 38856 | 111.92 | NG | 96 |
| 36) *Chlorobenzene-d5 | 26.69 | 647 | 242723 | 250.00 | NG | 86 |
| 37) Methyl-iso-butyl ketone | 22.49 | 539 | 36908 | 111.11 | NG | 95 |
| 38) 2-Hexanone | 24.04 | 579 | 37569 | 117.06 | NG | 91 |
| 38) 2-Hexanone | 24.70 | 596 | 2378 | 7.41 | NG | 89 |
| 39) 1,1,2,2-Tetrachloroethane | 24.27 | 585 | 51853 | 102.83 | NG | 93 |
| 40) Tetrachloroethylene | 24.35 | 587 | 35670 | 106.50 | NG | 99 |

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QUANT REPORT

Page 2

Operator ID: JAB781
 Output File: ^C0962::AQ
 Data File: >C0962::U0
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5,, 2UL

Quant Rev: 7 Quant Time: 901002 17:06
 Injected at: 901002 15:32
 Dilution Factor: 1.00000

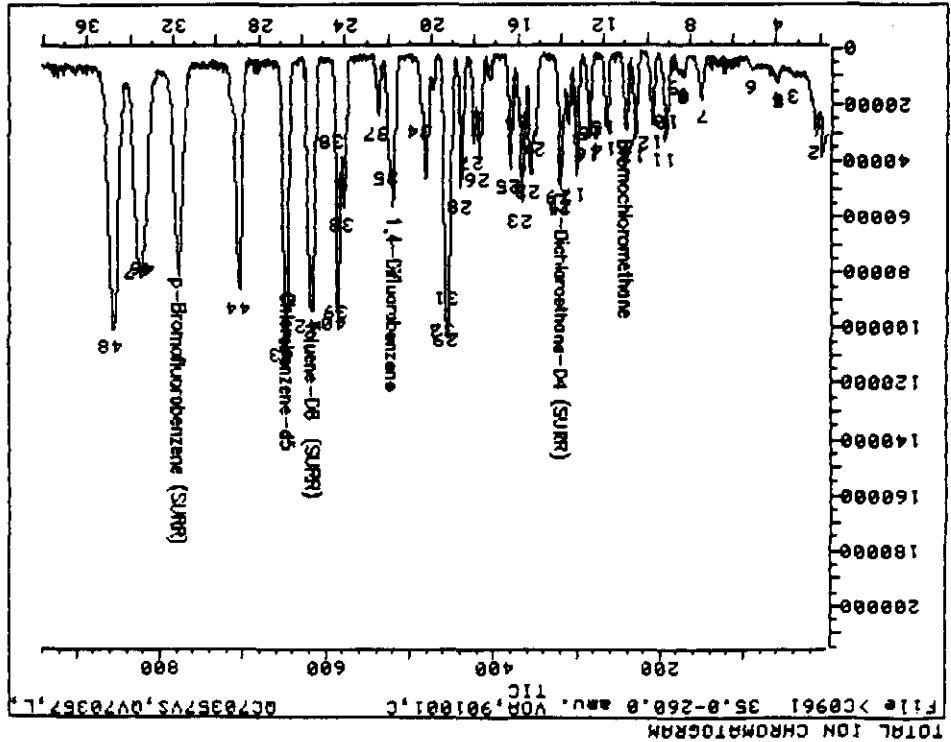
ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 41) | Toluene-D8 (SURR) | 25.48 | 616 | 358633 | 252.30 | NG | 93 |
| 42) | Toluene | 25.67 | 621 | 83277 | 103.36 | NG | 91 |
| 43) | Chlorobenzene | 26.80 | 650 | 119003 | 105.55 | NG | 97 |
| 44) | Ethylbenzene | 28.94 | 705 | 73613 | 108.50 | NG | 81 |
| 45) | p-Bromofluorobenzene (SURR) | 31.73 | 777 | 262700 | 261.64 | NG | 92 |
| 46) | Styrene | 33.36 | 819 | 176969 | 110.72 | NG | 93 |
| 47) | m-Xylene | 33.71 | 828 | 97183 | 109.35 | NG | 99 |
| 48) | o+p-Xylenes | 34.87 | 858 | 203958 | 222.26 | NG | 88 |

* Compound is ISTD

Data File: >C0961:U0
 Name: VOA,901001,C
 Misc: DC70357VS,QV70357,L,5,, 5UL
 ID File: IC1121:US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53
 Operator ID: JAB781
 Quant Time: 901002 16:55
 Injected at: 901002 14:43

Quant Output File: >C0961:AD



QUANT REPORT

Operator ID: JA8781
 Output File: ^C0961::AQ
 Data File: >C0961::U0
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5,, 5UL

Quant Rev: 7 Quant Time: 901002 16:55
 Injected at: 901002 14:43
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|---------|-------|-----|
| 1) | *Bromochloromethane | 10.98 | 242 | 45268 | 250.00 | NG | 94 |
| 2) | Methyl chloride | 2.17 | 15 | 41886 | 223.72 | NG | 99 |
| 3) | Methyl bromide | 3.18 | 41 | 7729 | 259.43 | NG | 99 |
| 4) | Dichlorodifluoromethane | 3.92 | 60 | 21384 | 253.64 | NG | 97 |
| 5) | Vinyl chloride | 3.95 | 61 | 30956 | 255.71 | NG | 94 |
| 6) | Chloroethene | 5.16 | 92 | 18624 | 235.25 | NG | 89 |
| 7) | Methylene chloride | 7.45 | 151 | 52721 | 262.83 | NG | 96 |
| 8) | Acrolein | 8.30 | 173 | 49447 | 2383.42 | NG | 91 |
| 9) | Acetone | 8.34 | 174 | 27481 | 263.81 | NG | 93 |
| 9) | Acetone | 8.77 | 185 | 2024 | 19.43 | NG | 41 |
| 10) | Acrylonitrile | 9.00 | 191 | 45608 | 500.09 | NG | 91 |
| 11) | Carbon disulfide | 9.12 | 194 | 302293 | 287.91 | NG | 97 |
| 12) | Trichlorofluoromethane | 9.74 | 210 | 158968 | 243.07 | NG | 91 |
| 13) | 1,1-Dichloroethylene | 10.51 | 230 | 73422 | 239.63 | NG | 93 |
| 14) | 1,1-Dichloroethane | 11.87 | 265 | 147715 | 239.48 | NG | 97 |
| 15) | Tetrahydrofuran | 11.95 | 267 | 7108 | 149.39 | NG | 100 |
| 15) | Tetrahydrofuran | 12.60 | 284 | 2152 | 45.23 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.64 | 285 | 71329 | 240.52 | NG | 98 |
| 17) | Chloroform | 13.26 | 301 | 203625 | 239.29 | NG | 96 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.96 | 319 | 137642 | 235.38 | NG | 96 |
| 19) | 1,2-Dichloroethane | 14.04 | 321 | 164881 | 250.20 | NG | 97 |
| 20) | *1,4-Difluorobenzene | 21.76 | 520 | 234953 | 250.00 | NG | 96 |
| 21) | Methyl ethyl ketone | 14.00 | 320 | 9007 | 275.25 | NG | 84 |
| 22) | 1,1,1-Trichloroethane | 15.40 | 356 | 177556 | 250.68 | NG | 96 |
| 23) | Carbon tetrachloride | 15.40 | 356 | 19550 | 30.08 | NG | 98 |
| 23) | Carbon tetrachloride | 15.79 | 366 | 159011 | 244.63 | NG | 96 |
| 24) | Vinyl acetate | 15.16 | 350 | 47100 | 98.18 | NG | 81 |
| 24) | Vinyl acetate | 15.98 | 371 | 122478 | 255.32 | NG | 99 |
| 25) | Dichlorobromomethane | 16.37 | 381 | 166787 | 252.07 | NG | 94 |
| 26) | 1,2-Dichloropropane | 17.76 | 417 | 80498 | 245.18 | NG | 91 |
| 27) | cis-1,3-Dichloropropylene | 18.03 | 424 | 115898 | 249.21 | NG | 92 |
| 28) | Trichloroethylene | 18.62 | 439 | 85764 | 247.05 | NG | 87 |
| 29) | Chlorodibromomethane | 19.24 | 455 | 120679 | 249.32 | NG | 99 |
| 30) | bis(Chloromethyl)ether | 19.24 | 455 | 50123 | 254.35 | NG | 100 |
| 31) | Benzene | 19.16 | 453 | 261723 | 239.18 | NG | 93 |
| 32) | 1,1,2-Trichloroethane | 19.35 | 458 | 75776 | 244.87 | NG | 84 |
| 33) | trans-1,3-Dichloropropylene | 19.35 | 458 | 152620 | 251.75 | NG | 89 |
| 34) | 2-Chloroethylvinyl ether | 20.44 | 486 | 58207 | 248.73 | NG | 100 |
| 35) | Bromoform | 21.95 | 525 | 103057 | 257.13 | NG | 94 |
| 36) | *Chlorobenzene-d5 | 26.65 | 646 | 265376 | 250.00 | NG | 86 |
| 37) | Methyl-iso-butyl ketone | 22.46 | 538 | 99055 | 272.76 | NG | 99 |
| 38) | 2-Hexanone | 24.01 | 578 | 91684 | 261.29 | NG | 88 |
| 38) | 2-Hexanone | 24.47 | 590 | 3200 | 9.12 | NG | 86 |

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QUANT REPORT

Page 2

Operator ID: JAB781
 Output File: ^C0961::AQ
 Data File: >C0961::U0
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5,, 5UL

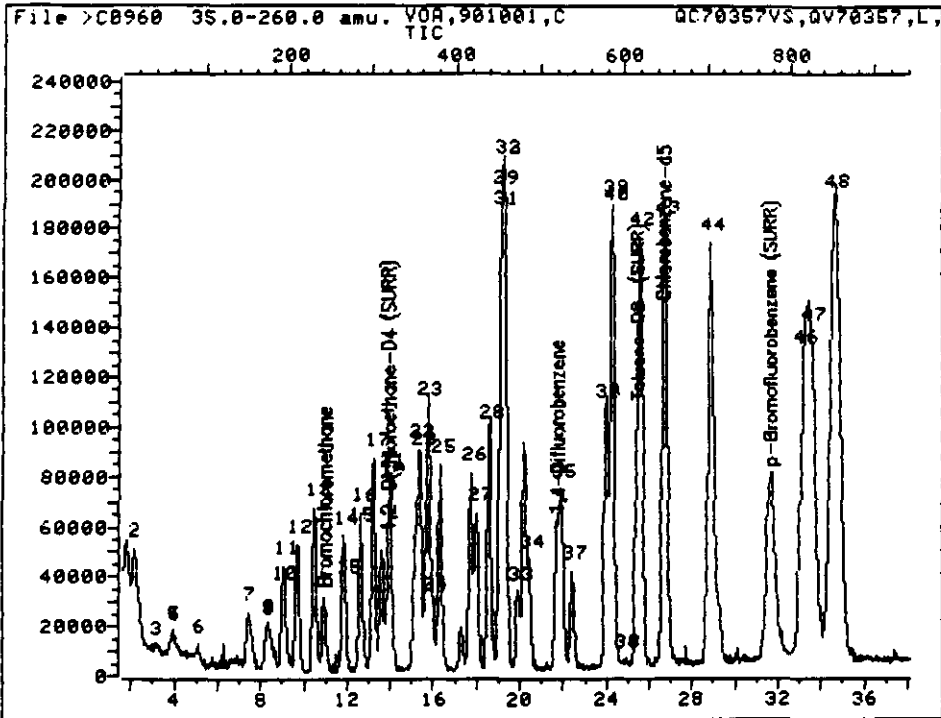
Quant Rev: 7 Quant Time: 901002 16:55
 Injected at: 901002 14:43
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 39) | 1,1,2,2-Tetrachloroethane | 24.28 | 585 | 144422 | 261.96 | NG | 99 |
| 40) | Tetrachloroethylene | 24.36 | 587 | 92698 | 253.15 | NG | 96 |
| 41) | Toluene-D8 (SURR) | 25.48 | 616 | 390359 | 251.18 | NG | 89 |
| 42) | Toluene | 25.64 | 620 | 214769 | 243.82 | NG | 91 |
| 43) | Chlorobenzene | 26.81 | 650 | 310398 | 251.80 | NG | 94 |
| 44) | Ethylbenzene | 28.90 | 704 | 183632 | 247.57 | NG | 79 |
| 45) | p-Bromofluorobenzene (SURR) | 31.74 | 777 | 275490 | 250.96 | NG | 83 |
| 46) | Styrene | 33.37 | 819 | 441269 | 252.51 | NG | 93 |
| 47) | m-Xylene | 33.68 | 827 | 243938 | 251.05 | NG | 97 |
| 48) | o+p-Xylenes | 34.76 | 855 | 503073 | 501.43 | NG | 89 |

* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >C0960::U0 Quant Output File: ^C0960::AQ
 Name: VOA,901001,C
 Misc: QC70357VS,QU70357,L,5,, 10UL

Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

Operator ID: JAB781
 Quant Time: 901002 17:03
 Injected at: 901002 13:53

QUANT REPORT

Page 1

Operator ID: JAB781 Quant Rev: 7 Quant Time: 901002 17:03
 Output File: ^C0960::AQ Injected at: 901002 13:53
 Data File: >C0960::U0 Dilution Factor: 1.00000
 Name: VOA,901001,C
 Misc: QC70357US,QV70357,L,5,, 10UL

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|---------|-------|-----|
| 1) *Bromochloromethane | 10.93 | 241 | 46511 | 250.00 | NG | 92 |
| 2) Methyl chloride | 2.12 | 14 | 86009 | 447.10 | NG | 98 |
| 3) Methyl bromide | 3.13 | 40 | 11787 | 385.06 | NG | 93 |
| 4) Dichlorodifluoromethane | 3.87 | 59 | 36001 | 415.60 | NG | 95 |
| 5) Vinyl chloride | 3.95 | 61 | 52754 | 424.12 | NG | 97 |
| 6) Chloroethane | 5.11 | 91 | 37464 | 460.58 | NG | 94 |
| 7) Methylene chloride | 7.44 | 151 | 82230 | 398.98 | NG | 99 |
| 8) Acrolein | 8.26 | 172 | 130129 | 6104.78 | NG | 87 |
| 9) Acetone | 8.33 | 174 | 45276 | 423.02 | NG | 94 |
| 10) Acrylonitrile | 8.99 | 191 | 57272 | 611.21 | NG | 93 |
| 11) Carbon disulfide | 9.07 | 193 | 356632 | 330.59 | NG | 99 |
| 12) Trichlorofluoromethane | 9.69 | 209 | 314827 | 468.52 | NG | 96 |
| 13) 1,1-Dichloroethylene | 10.51 | 230 | 155521 | 494.02 | NG | 96 |
| 14) 1,1-Dichloroethane | 11.82 | 264 | 301095 | 475.09 | NG | 95 |
| 15) Tetrahydrofuran | 11.94 | 267 | 30274 | 619.26 | NG | 100 |
| 15) Tetrahydrofuran | 12.60 | 284 | 4645 | 95.01 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.64 | 285 | 148502 | 487.36 | NG | 98 |
| 17) Chloroform | 13.26 | 301 | 417209 | 477.18 | NG | 94 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.92 | 318 | 143846 | 239.42 | NG | 98 |
| 19) 1,2-Dichloroethane | 14.03 | 321 | 319071 | 471.23 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.72 | 519 | 256903 | 250.00 | NG | 99 |
| 21) Methyl ethyl ketone | 14.00 | 320 | 17574 | 491.17 | NG | 92 |
| 22) 1,1,1-Trichloroethane | 15.35 | 355 | 357272 | 461.31 | NG | 93 |
| 23) Carbon tetrachloride | 15.39 | 356 | 40615 | 57.15 | NG | 97 |
| 23) Carbon tetrachloride | 15.74 | 365 | 330550 | 465.09 | NG | 97 |
| 24) Vinyl acetate | 16.01 | 372 | 234265 | 446.62 | NG | 96 |
| 25) Dichlorobromomethane | 16.32 | 380 | 335186 | 463.29 | NG | 98 |
| 26) 1,2-Dichloropropane | 17.76 | 417 | 169497 | 472.15 | NG | 97 |
| 27) cis-1,3-Dichloropropylene | 17.99 | 423 | 239863 | 471.71 | NG | 91 |
| 28) Trichloroethylene | 18.57 | 438 | 188967 | 497.83 | NG | 92 |
| 29) Chlorodibromomethane | 19.19 | 454 | 252858 | 477.77 | NG | 96 |
| 30) bis(Chloromethyl)ether | 19.19 | 454 | 100814 | 467.88 | NG | 100 |
| 31) Benzene | 19.15 | 453 | 574627 | 480.27 | NG | 94 |
| 32) 1,1,2-Trichloroethane | 19.31 | 457 | 165271 | 488.44 | NG | 88 |
| 33) trans-1,3-Dichloropropylene | 19.31 | 457 | 309782 | 467.34 | NG | 92 |
| 33) trans-1,3-Dichloropropylene | 19.93 | 473 | 38584 | 58.21 | NG | 58 |
| 34) 2-Chloroethylvinyl ether | 20.44 | 486 | 126916 | 495.99 | NG | 100 |
| 35) Bromoform | 21.91 | 524 | 214387 | 489.21 | NG | 95 |
| 36) *Chlorobenzene-d5 | 26.66 | 646 | 285650 | 250.00 | NG | 87 |
| 37) Methyl-iso-butyl ketone | 22.41 | 537 | 196718 | 503.24 | NG | 94 |
| 38) 2-Hexanone | 23.97 | 577 | 191773 | 507.75 | NG | 90 |
| 38) 2-Hexanone | 24.86 | 600 | 4253 | 11.26 | NG | 95 |
| 39) 1,1,2,2-Tetrachloroethane | 24.28 | 585 | 278991 | 470.12 | NG | 93 |

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QUANT REPORT

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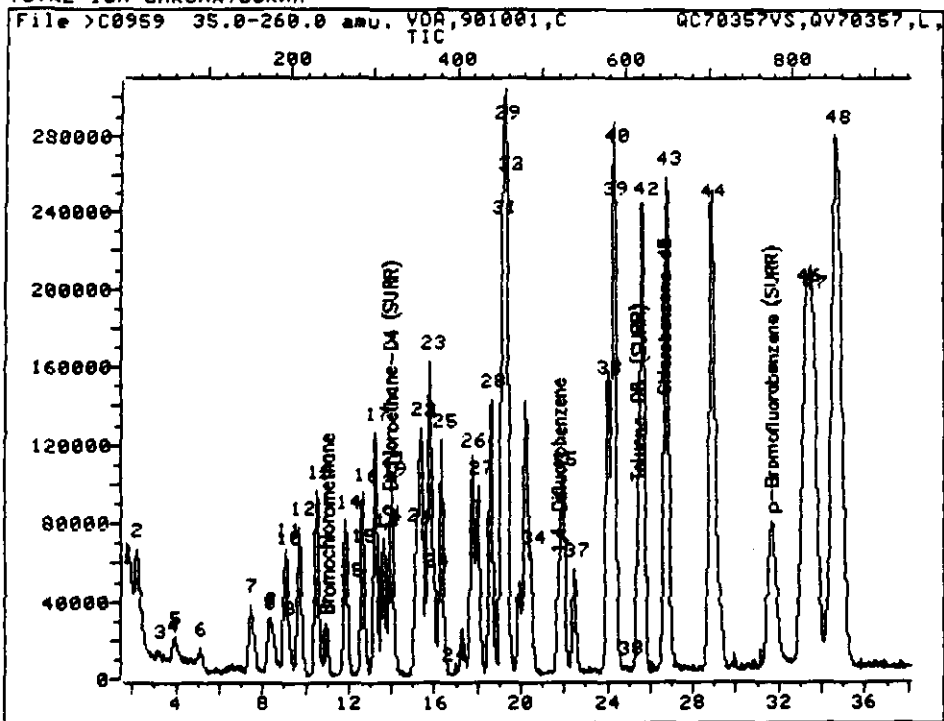
Operator ID: JAB781 Quant Rev: 7 Quant Time: 901002 17:03
 Output File: ^C0960::AQ Injected at: 901002 13:53
 Data File: >C0960::U0 Dilution Factor: 1.00000
 Name: VOA,901001,C
 Misc: QC70357VS,QU70357,L,5,, 10UL

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|--------|-------|----|
| 40) | Tetrachloroethylene | 24.31 | 586 | 196320 | 498.08 | NG | 99 |
| 41) | Toluene-DB (SURR) | 25.44 | 615 | 415602 | 248.44 | NG | 91 |
| 42) | Toluene | 25.60 | 619 | 464802 | 490.22 | NG | 93 |
| 43) | Chlorobenzene | 26.77 | 649 | 644676 | 485.85 | NG | 93 |
| 44) | Ethylbenzene | 28.87 | 703 | 383763 | 480.65 | NG | 79 |
| 45) | p-Bromofluorobenzene (SURR) | 31.62 | 774 | 286218 | 242.23 | NG | 91 |
| 46) | Styrene | 33.21 | 815 | 901649 | 479.34 | NG | 92 |
| 47) | m-Xylene | 33.56 | 824 | 497605 | 475.77 | NG | 96 |
| 48) | o+p-Xylenes | 34.65 | 852 | 1029765 | 953.55 | NG | 87 |

* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >C0959::U0 Quant Output File: ^C0959::AQ
 Name: VOA,901001,C
 Misc: QC70357VS,QV70357,L,5,, 15UL

Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

Operator ID: JA8781
 Quant Time: 901002 17:01
 Injected at: 901002 13:04

QUANT REPORT

Page 1

Operator ID: JAB781
 Output File: ^C0959::AQ
 Data File: >C0959::U0
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5,, 15UL

Quant Rev: 7 Quant Time: 901002 17:01
 Injected at: 901002 13:04
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVQA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|----------|-------|-----|
| 1) | *Bromochloromethane | 10.97 | 242 | 43946 | 250.00 | NG | 90 |
| 2) | Methyl chloride | 2.12 | 14 | 119589 | 657.95 | NG | 97 |
| 3) | Methyl bromide | 3.17 | 41 | 15696 | 542.69 | NG | 99 |
| 4) | Dichlorodifluoromethane | 3.83 | 58 | 47936 | 585.67 | NG | 96 |
| 5) | Vinyl chloride | 3.91 | 60 | 69767 | 593.64 | NG | 93 |
| 6) | Chloroethane | 5.11 | 91 | 53146 | 691.50 | NG | 99 |
| 7) | Methylene chloride | 7.44 | 151 | 124051 | 637.03 | NG | 99 |
| 8) | Acrolein | 8.26 | 172 | 209088 | 10381.53 | NG | 92 |
| 8) | Acrolein | 9.23 | 197 | 4123 | 204.71 | NG | 94 |
| 9) | Acetone | 8.33 | 174 | 65468 | 647.37 | NG | 95 |
| 10) | Acrylonitrile | 9.03 | 192 | 82560 | 932.51 | NG | 96 |
| 11) | Carbon disulfide | 9.07 | 193 | 598559 | 587.23 | NG | 99 |
| 12) | Trichlorofluoromethane | 9.69 | 209 | 474825 | 747.88 | NG | 91 |
| 13) | 1,1-Dichloroethylene | 10.51 | 230 | 223289 | 750.68 | NG | 93 |
| 14) | 1,1-Dichloroethane | 11.86 | 265 | 436142 | 728.35 | NG | 95 |
| 15) | Tetrahydrofuran | 11.94 | 267 | 39683 | 859.10 | NG | 100 |
| 15) | Tetrahydrofuran | 12.56 | 283 | 4728 | 102.36 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.64 | 285 | 218918 | 760.39 | NG | 98 |
| 17) | Chloroform | 13.22 | 300 | 601279 | 727.85 | NG | 98 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.92 | 318 | 143382 | 252.58 | NG | 95 |
| 19) | 1,2-Dichloroethane | 13.61 | 310 | 3987 | 6.23 | NG | 66 |
| 19) | 1,2-Dichloroethane | 14.03 | 321 | 450613 | 704.35 | NG | 98 |
| 20) | *1,4-Difluorobenzene | 21.72 | 519 | 242651 | 250.00 | NG | 99 |
| 21) | Methyl ethyl ketone | 14.00 | 320 | 24669 | 729.96 | NG | 89 |
| 22) | 1,1,1-Trichloroethane | 15.35 | 355 | 517516 | 707.46 | NG | 98 |
| 23) | Carbon tetrachloride | 15.35 | 355 | 56892 | 84.75 | NG | 96 |
| 23) | Carbon tetrachloride | 15.78 | 366 | 484005 | 721.00 | NG | 97 |
| 24) | Vinyl acetate | 15.12 | 349 | 127578 | 257.51 | NG | 76 |
| 24) | Vinyl acetate | 15.97 | 371 | 391627 | 790.49 | NG | 98 |
| 24) | Vinyl acetate | 16.79 | 392 | 2236 | 4.51 | NG | 50 |
| 25) | Dichlorobromomethane | 16.32 | 380 | 488465 | 714.81 | NG | 96 |
| 26) | 1,2-Dichloropropane | 17.72 | 416 | 249216 | 734.99 | NG | 95 |
| 27) | cis-1,3-Dichloropropylene | 17.99 | 423 | 355519 | 740.22 | NG | 92 |
| 28) | Trichloroethylene | 18.61 | 439 | 261235 | 728.64 | NG | 87 |
| 29) | Chlorodibromomethane | 19.19 | 454 | 368760 | 737.69 | NG | 99 |
| 30) | bis(Chloromethyl)ether | 19.19 | 454 | 145087 | 712.90 | NG | 100 |
| 31) | Benzene | 19.12 | 452 | 864067 | 764.60 | NG | 96 |
| 32) | 1,1,2-Trichloroethane | 19.35 | 458 | 236975 | 741.48 | NG | 84 |
| 33) | trans-1,3-Dichloropropylene | 19.35 | 458 | 456365 | 728.91 | NG | 90 |
| 34) | 2-Chloroethylvinyl ether | 20.44 | 486 | 177590 | 734.79 | NG | 100 |
| 35) | Bromoform | 21.91 | 524 | 302707 | 731.31 | NG | 98 |
| 36) | *Chlorobenzene-d5 | 26.62 | 645 | 281445 | 250.00 | NG | 86 |
| 37) | Methyl-iso-butyl ketone | 22.45 | 538 | 270605 | 702.59 | NG | 94 |

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QUANT REPORT

Page 2

Operator ID: JA8781
 Output File: ^C0959::AQ
 Data File: >C0959::U0
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5,, 15UL

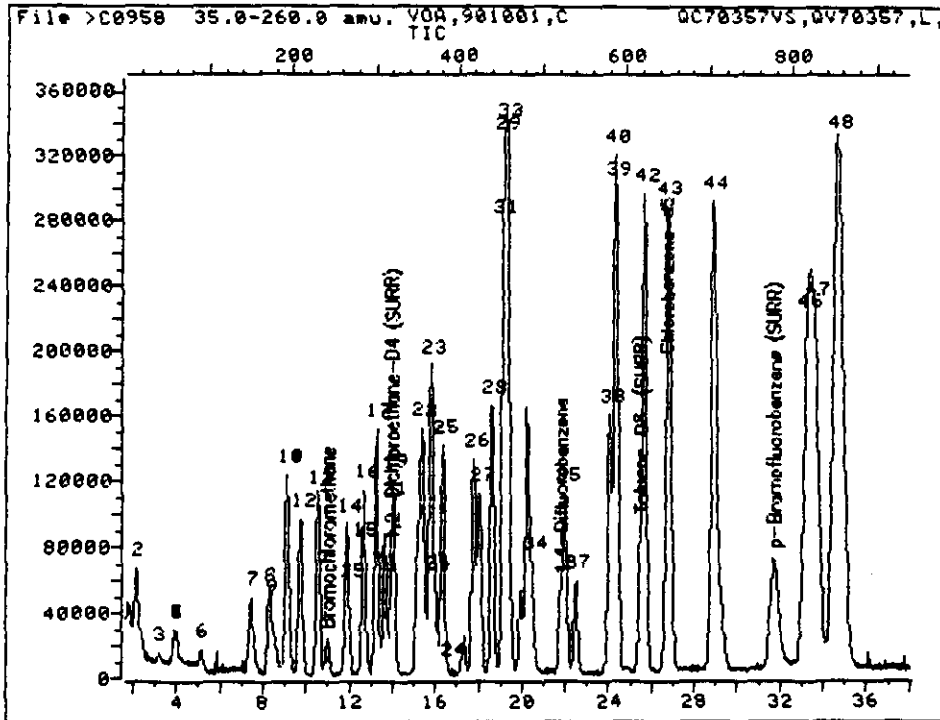
Quant Rev: 7 Quant Time: 901002 17:01
 Injected at: 901002 13:04
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|---------|-------|----|
| 38) | 2-Hexanone | 23.97 | 577 | 259439 | 697.16 | NG | 99 |
| 38) | 2-Hexanone | 24.94 | 602 | 3828 | 10.29 | NG | 99 |
| 39) | 1,1,2,2-Tetrachloroethane | 24.24 | 584 | 443558 | 758.60 | NG | 98 |
| 40) | Tetrachloroethylene | 24.32 | 586 | 282256 | 726.81 | NG | 98 |
| 41) | Toluene-D8 (SURR) | 25.40 | 614 | 409590 | 248.51 | NG | 91 |
| 42) | Toluene | 25.64 | 620 | 689511 | 738.08 | NG | 91 |
| 43) | Chlorobenzene | 26.74 | 648 | 946510 | 723.99 | NG | 92 |
| 44) | Ethylbenzene | 28.83 | 702 | 571616 | 726.63 | NG | 81 |
| 45) | p-Bromofluorobenzene (SURR) | 31.63 | 774 | 282323 | 242.50 | NG | 86 |
| 46) | Styrene | 33.26 | 816 | 1302237 | 702.65 | NG | 90 |
| 47) | m-Xylene | 33.57 | 824 | 734713 | 712.98 | NG | 98 |
| 48) | o+p-Xylenes | 34.65 | 852 | 1496032 | 1406.00 | NG | 88 |

* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >C0958::U0 Quant Output File: ^C0958::AQ
Name: VOA,901001,C
Misc: QC70357VS,QU70357,L,5,, 20UL

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901002 16:53

Operator ID: JA8781
Quant Time: 901002 16:58
Injected at: 901002 12:15

QUANT REPORT

Page 1

Operator ID: JAB781
 Output File: ^C0958::AQ
 Data File: >C0958::U0
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5,, 20UL

Quant Rev: 7 Quant Time: 901002 16:58
 Injected at: 901002 12:15
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|---------|----------|-------|-----|
| 1) *Bromochloromethane | 10.96 | 241 | 35888 | 250.00 | NG | 98 |
| 2) Methyl chloride | 2.12 | 13 | 188780 | 1271.83 | NG | 98 |
| 3) Methyl bromide | 3.16 | 40 | 19565 | 828.34 | NG | 99 |
| 4) Dichlorodifluoromethane | 3.90 | 59 | 72696 | 1087.61 | NG | 97 |
| 5) Vinyl chloride | 3.90 | 59 | 100809 | 1050.37 | NG | 98 |
| 6) Chloroethane | 5.14 | 91 | 67975 | 1083.03 | NG | 99 |
| 7) Methylene chloride | 7.43 | 150 | 152042 | 956.08 | NG | 97 |
| 8) Acrolein | 8.25 | 171 | 357306 | 21724.14 | NG | 93 |
| 9) Acetone | 8.37 | 174 | 69403 | 840.37 | NG | 95 |
| 10) Acrylonitrile | 9.06 | 192 | 113050 | 1563.59 | NG | 99 |
| 11) Carbon disulfide | 9.10 | 193 | 1020286 | 1225.73 | NG | 99 |
| 12) Trichlorofluoromethane | 9.72 | 209 | 576578 | 1112.05 | NG | 93 |
| 13) 1,1-Dichloroethylene | 10.50 | 229 | 261922 | 1078.28 | NG | 95 |
| 14) 1,1-Dichloroethane | 11.86 | 264 | 516512 | 1056.24 | NG | 96 |
| 15) Tetrahydrofuran | 11.97 | 267 | 36980 | 980.34 | NG | 100 |
| 15) Tetrahydrofuran | 12.59 | 283 | 11239 | 297.95 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.67 | 285 | 248677 | 1057.69 | NG | 97 |
| 17) Chloroform | 13.25 | 300 | 700113 | 1037.77 | NG | 98 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.95 | 318 | 126995 | 273.94 | NG | 97 |
| 19) 1,2-Dichloroethane | 14.03 | 320 | 525411 | 1005.66 | NG | 94 |
| 20) *1,4-Difluorobenzene | 21.79 | 520 | 205564 | 250.00 | NG | 98 |
| 21) Methyl ethyl ketone | 13.99 | 319 | 24075 | 840.91 | NG | 89 |
| 22) 1,1,1-Trichloroethane | 15.39 | 355 | 607067 | 979.60 | NG | 97 |
| 23) Carbon tetrachloride | 15.39 | 355 | 65406 | 115.01 | NG | 93 |
| 23) Carbon tetrachloride | 15.81 | 366 | 579302 | 1018.65 | NG | 99 |
| 24) Vinyl acetate | 16.01 | 371 | 446961 | 1064.94 | NG | 96 |
| 24) Vinyl acetate | 16.74 | 390 | 2108 | 5.02 | NG | 59 |
| 25) Dichlorobromomethane | 16.36 | 380 | 564259 | 974.70 | NG | 97 |
| 26) 1,2-Dichloropropane | 17.75 | 416 | 284368 | 989.97 | NG | 88 |
| 27) cis-1,3-Dichloropropylene | 18.02 | 423 | 405944 | 997.69 | NG | 94 |
| 28) Trichloroethylene | 18.61 | 438 | 291786 | 960.68 | NG | 86 |
| 29) Chlorodibromomethane | 19.23 | 454 | 419860 | 991.44 | NG | 96 |
| 30) bis(Chloromethyl)ether | 19.23 | 454 | 168159 | 975.33 | NG | 100 |
| 31) Benzene | 19.15 | 452 | 1005812 | 1050.60 | NG | 98 |
| 32) 1,1,2-Trichloroethane | 19.34 | 457 | 271744 | 1003.68 | NG | 82 |
| 33) trans-1,3-Dichloropropylene | 19.34 | 457 | 520749 | 981.81 | NG | 94 |
| 34) 2-Chloroethylvinyl ether | 20.47 | 486 | 200384 | 978.68 | NG | 100 |
| 35) Bromoform | 21.98 | 525 | 315163 | 898.77 | NG | 94 |
| 36) *Chlorobenzene-d5 | 26.70 | 646 | 236905 | 250.00 | NG | 85 |
| 37) Methyl-iso-butyl ketone | 22.49 | 538 | 277047 | 854.56 | NG | 98 |
| 38) 2-Hexanone | 24.04 | 578 | 262866 | 839.18 | NG | 95 |
| 39) 1,1,2,2-Tetrachloroethane | 24.31 | 585 | 478478 | 972.17 | NG | 97 |
| 40) Tetrachloroethylene | 24.35 | 586 | 312878 | 957.13 | NG | 97 |

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QUANT REPORT

Page 2

Operator ID: JA8781 Quant Rev: 7 Quant Time: 901002 16:58
 Output File: ^C0958::AQ Injected at: 901002 12:15
 Data File: >C0958::U0 Dilution Factor: 1.00000
 Name: VOA,901001,C
 Misc: QC70357US,QV70357,L,5,, 20UL

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|---------|-------|----|
| 41) | Toluene-D8 (SURR) | 25.47 | 615 | 346251 | 249.57 | NG | 95 |
| 42) | Toluene | 25.67 | 620 | 807213 | 1026.53 | NG | 90 |
| 43) | Chlorobenzene | 26.78 | 648 | 1100799 | 1000.31 | NG | 89 |
| 44) | Ethylbenzene | 28.91 | 703 | 658561 | 994.55 | NG | 80 |
| 45) | p-Bromofluorobenzene (SURR) | 31.71 | 775 | 247619 | 252.68 | NG | 80 |
| 46) | Styrene | 33.26 | 815 | 1540003 | 987.17 | NG | 91 |
| 47) | m-Xylene | 33.61 | 824 | 867468 | 1000.07 | NG | 95 |
| 48) | o+p-Xylenes | 34.73 | 853 | 1782225 | 1989.88 | NG | 87 |

* Compound is ISTD

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:ETCNJ

Contract:

Lab Code:F2

Case No.:

SAS No.:

SDG No.:

Instrument ID:GC/MS C

Calibration Date(s) 10/03/90

10/03/90

Matrix:(soil/water) WATER

Level:(low/med) LOW

Column:(pack/cap) PACK

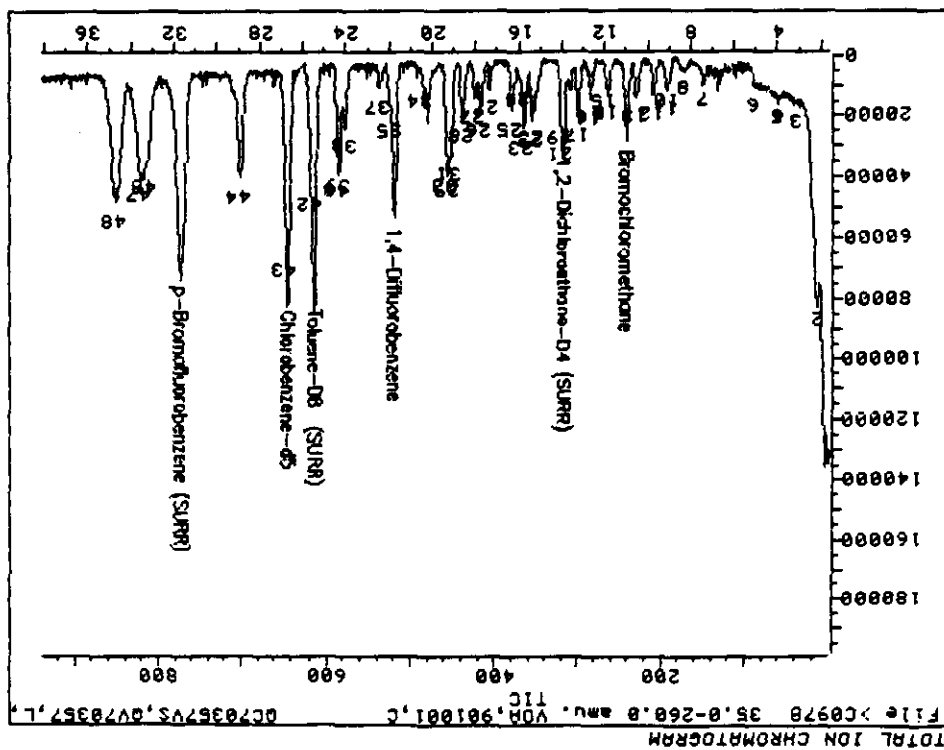
Min RRF for SPCC(*) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

| LAB FILE ID: | RRF20 =>C0978 | RRF50 =>C0972 | | | RRF100 =>C0977 | RRF150 =>C0976 | RRF200 =>C0975 | | |
|----------------------------|---------------|---------------|--------|--------|----------------|----------------|----------------|-----|--|
| COMPOUND | RRF20 | RRF50 | RRF100 | RRF150 | RRF200 | RRF | % | RSD | |
| Chloromethane | .573 | .735 | 1.123 | .521 | 1.488 | .888 | 46.2 | * | |
| Bromomethane | .240 | .302 | .140 | .103 | .165 | .190 | 42.3 | | |
| Vinyl Chloride | .705 | .854 | .683 | .478 | .857 | .715 | 21.8 | * | |
| Chloroethane | .426 | .567 | .460 | .366 | .622 | .488 | 21.4 | | |
| Methylene Chloride | 1.204 | 1.588 | 1.176 | 1.067 | 1.516 | 1.310 | 17.4 | | |
| Acetone | .957 | .896 | .582 | .544 | .668 | .729 | 25.6 | | |
| Carbon Disulfide | 5.944 | 7.468 | 4.613 | 7.465 | 6.830 | 6.464 | 18.7 | | |
| 1,1-Dichloroethene | 1.678 | 1.666 | 1.635 | 1.717 | 1.794 | 1.698 | 3.6 | * | |
| 1,1-Dichloroethane | 3.608 | 3.578 | 3.360 | 3.473 | 3.611 | 3.526 | 3.1 | * | |
| 1,2-Dichloroethene (total) | 1.660 | 1.572 | 1.556 | 1.657 | 1.641 | 1.617 | 3.0 | | |
| Chloroform | 5.094 | 5.263 | 4.600 | 4.849 | 5.013 | 4.964 | 5.1 | * | |
| 1,2-Dichloroethane | 4.185 | 4.302 | 3.690 | 3.764 | 3.965 | 3.981 | 6.6 | | |
| 2-Butanone | .058 | .072 | .032 | .032 | .029 | .045 | 43.2 | | |
| 1,1,1-Trichloroethane | .911 | 1.039 | .750 | .799 | .812 | .862 | 13.3 | | |
| Carbon Tetrachloride | .826 | .940 | .699 | .736 | .765 | .793 | 11.9 | | |
| Vinyl Acetate | .568 | .570 | .581 | .623 | .528 | .574 | 5.9 | | |
| Bromodichloromethane | .825 | .899 | .695 | .738 | .736 | .779 | 10.5 | | |
| 1,2-Dichloropropane | .356 | .331 | .323 | .347 | .323 | .336 | 4.5 | * | |
| cis-1,3-Dichloropropene | .509 | .553 | .485 | .513 | .491 | .510 | 5.2 | | |
| Trichloroethene | .408 | .373 | .353 | .366 | .362 | .372 | 5.7 | | |
| Dibromochloromethane | .545 | .538 | .511 | .528 | .520 | .528 | 2.6 | | |
| 1,1,2-Trichloroethane | .321 | .304 | .322 | .335 | .330 | .322 | 3.7 | | |
| Benzene | 1.159 | 1.120 | 1.118 | 1.243 | 1.203 | 1.169 | 4.6 | | |
| trans-1,3-Dichloropropene | .686 | .732 | .650 | .663 | .673 | .681 | 4.6 | | |
| Bromoform | .477 | .403 | .445 | .429 | .409 | .433 | 6.9 | * | |
| 4-Methyl-2-Pentanone | .426 | .298 | .342 | .340 | .289 | .339 | 16.1 | | |
| 2-Hexanone | .389 | .267 | .323 | .329 | .274 | .316 | 15.6 | | |
| Tetrachloroethene | .403 | .345 | .349 | .330 | .329 | .351 | 8.6 | | |
| 1,1,2,2-Tetrachloroethane | .548 | .412 | .537 | .535 | .458 | .498 | 12.1 | * | |
| Toluene | .849 | .845 | .817 | .867 | .888 | .853 | 3.1 | * | |
| Chlorobenzene | 1.214 | 1.141 | 1.135 | 1.119 | 1.161 | 1.154 | 3.2 | * | |
| Ethylbenzene | .743 | .730 | .699 | .678 | .738 | .718 | 3.9 | * | |
| Styrene | 1.859 | 1.866 | 1.696 | 1.586 | 1.760 | 1.753 | 6.7 | | |
| Xylene (total) | 1.092 | 1.143 | .987 | .921 | 1.026 | 1.034 | 8.4 | | |
| Toluene-d8 | 1.456 | 1.544 | 1.498 | 1.432 | 1.463 | 1.478 | 3.0 | | |
| Bromofluorobenzene | 1.116 | 1.275 | 1.085 | 1.042 | 1.116 | 1.127 | 7.8 | | |
| 1,2-Dichloroethane-d4 | 3.274 | 3.769 | 3.321 | 3.511 | 3.837 | 3.543 | 7.2 | | |

FORM VI UOA

1/87 Rev.

Data File: >C0978:U4
Name: V0A,901001,C
Misc: DC70357VS,QV70357,L,5, 2UL
ID File: IC1171:US
Title: IFB, RP/V0A, TCL, XU0A13
Last Calibration: 901002 16:53
Operator ID: KB6656
Quant Time: 901003 16:40
Injected at: 901003 15:35



QUANT REPORT

Operator ID: KB6656
 Output File: ^C0978::AQ
 Data File: >C0978::U4
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5, 2UL

Quant Rev: 7 Quant Time: 901003 16:40
 Injected at: 901003 15:35
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|---------|-------|-----|
| 1) *Bromochloromethane | 10.94 | 241 | 42566 | 250.00 | NG | 97 |
| 2) Methyl chloride | 2.14 | 14 | 9755 | 55.41 | NG | 92 |
| 3) Methyl bromide | 3.15 | 40 | 4089 | 145.96 | NG | 84 |
| 4) Dichlorodifluoromethane | 3.88 | 59 | 8620 | 108.73 | NG | 92 |
| 5) Vinyl chloride | 3.96 | 61 | 11997 | 105.39 | NG | 94 |
| 6) Chloroethane | 5.12 | 91 | 7254 | 97.44 | NG | 84 |
| 7) Methylene chloride | 7.42 | 150 | 20507 | 108.72 | NG | 99 |
| 8) Acrolein | 8.31 | 173 | 23677 | 1213.71 | NG | 93 |
| 9) Acetone | 8.31 | 173 | 16293 | 166.33 | NG | 89 |
| 10) Acrylonitrile | 9.01 | 191 | 14008 | 163.35 | NG | 90 |
| 11) Carbon disulfide | 9.08 | 193 | 101202 | 102.51 | NG | 97 |
| 12) Trichlorofluoromethane | 9.70 | 209 | 69603 | 113.18 | NG | 91 |
| 13) 1,1-Dichloroethylene | 10.48 | 229 | 28574 | 99.18 | NG | 93 |
| 14) 1,1-Dichloroethane | 11.84 | 264 | 61430 | 105.91 | NG | 97 |
| 14) 1,1-Dichloroethane | 12.65 | 285 | 14699 | 25.34 | NG | 65 |
| 15) Tetrahydrofuran | 11.87 | 265 | 2073 | 46.33 | NG | 100 |
| 15) Tetrahydrofuran | 11.95 | 267 | 2920 | 65.26 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.65 | 285 | 28272 | 101.38 | NG | 94 |
| 17) Chloroform | 13.23 | 300 | 86729 | 108.39 | NG | 97 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.89 | 317 | 139355 | 253.44 | NG | 95 |
| 19) 1,2-Dichloroethane | 14.05 | 321 | 71261 | 115.00 | NG | 96 |
| 20) *1,4-Difluorobenzene | 21.73 | 519 | 210589 | 250.00 | NG | 98 |
| 21) Methyl ethyl ketone | 14.01 | 320 | 4914 | 167.54 | NG | 99 |
| 22) 1,1,1-Trichloroethane | 15.37 | 355 | 76758 | 120.91 | NG | 95 |
| 23) Carbon tetrachloride | 15.33 | 354 | 8892 | 15.26 | NG | 81 |
| 23) Carbon tetrachloride | 15.75 | 365 | 69546 | 119.37 | NG | 96 |
| 24) Vinyl acetate | 15.99 | 371 | 47815 | 111.21 | NG | 97 |
| 25) Dichlorobromomethane | 16.34 | 380 | 69506 | 117.20 | NG | 98 |
| 26) 1,2-Dichloropropane | 17.73 | 416 | 30006 | 101.97 | NG | 99 |
| 27) cis-1,3-Dichloropropylene | 17.38 | 407 | 2119 | 5.08 | NG | 71 |
| 27) cis-1,3-Dichloropropylene | 18.00 | 423 | 42910 | 102.94 | NG | 92 |
| 28) Trichloroethylene | 18.59 | 438 | 34378 | 110.49 | NG | 86 |
| 29) Chlorodibromomethane | 19.21 | 454 | 45927 | 105.86 | NG | 99 |
| 30) bis(Chloromethyl)ether | 19.17 | 453 | 18986 | 107.49 | NG | 100 |
| 31) Benzene | 19.13 | 452 | 97590 | 99.50 | NG | 90 |
| 32) 1,1,2-Trichloroethane | 19.32 | 457 | 27035 | 97.47 | NG | 84 |
| 33) trans-1,3-Dichloropropylene | 19.32 | 457 | 57788 | 106.35 | NG | 88 |
| 34) 2-Chloroethylvinyl ether | 20.45 | 486 | 21125 | 100.71 | NG | 100 |
| 35) Bromoform | 21.92 | 524 | 40197 | 111.90 | NG | 98 |
| 36) *Chlorobenzene-d5 | 26.65 | 643 | 243193 | 250.00 | NG | 80 |
| 37) Methyl-iso-butyl ketone | 22.43 | 537 | 41488 | 124.66 | NG | 95 |
| 38) 2-Hexanone | 23.98 | 577 | 37852 | 117.71 | NG | 83 |
| 39) 1,1,2,2-Tetrachloroethane | 24.25 | 584 | 53349 | 105.59 | NG | 97 |

QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0978::AQ
 Data File: >C0978::U4
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5, 2UL

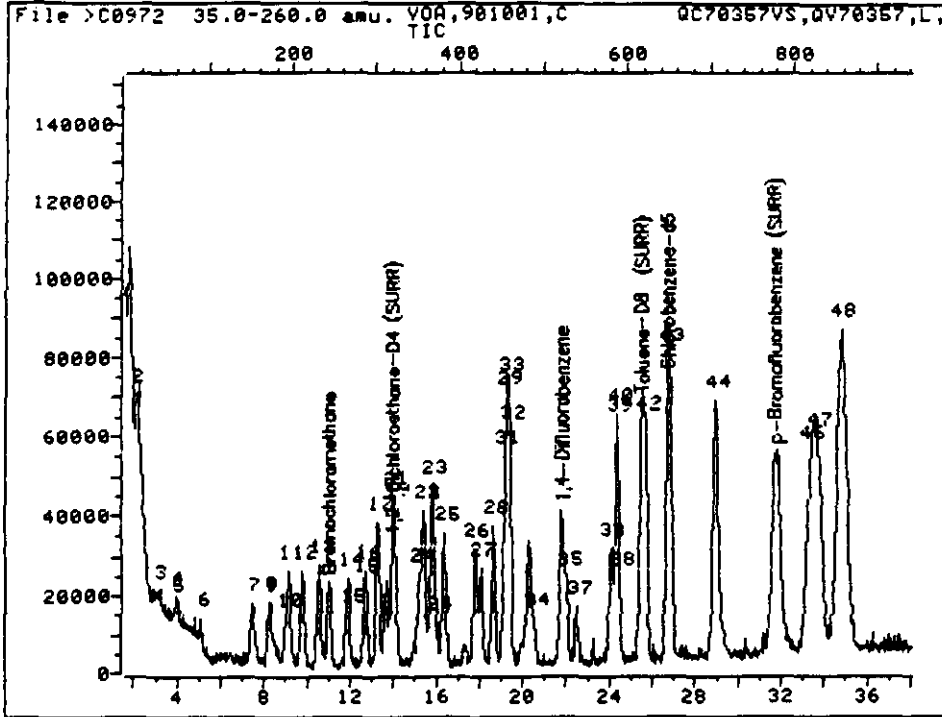
Quant Rev: 7 Quant Time: 901003 16:40
 . Injected at: 901003 15:35
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 40) | Tetrachloroethylene | 24.33 | 586 | 39161 | 116.70 | NG | 93 |
| 41) | Toluene-D8 (SURR) | 25.42 | 614 | 354032 | 248.58 | NG | 92 |
| 42) | Toluene | 25.61 | 619 | 82575 | 102.30 | NG | 93 |
| 43) | Chlorobenzene | 26.76 | 646 | 118136 | 104.58 | NG | 96 |
| 44) | Ethylbenzene | 28.86 | 700 | 72302 | 106.37 | NG | 78 |
| 45) | p-Bromofluorobenzene (SURR) | 31.65 | 772 | 271522 | 269.90 | NG | 83 |
| 46) | Styrene | 33.28 | 814 | 180878 | 112.95 | NG | 98 |
| 47) | m-Xylene | 33.56 | 821 | 102366 | 114.96 | NG | 96 |
| 48) | o+p-Xylenes | 34.72 | 851 | 212529 | 231.16 | NG | 90 |

* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >C0972::U4
Name: VOA,901001,C
Misc: QC70357VS,QV70357,L,5,

Quant Output File: ^C0972::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901002 16:53

Operator ID: KB6656
Quant Time: 901003 10:39
Injected at: 901003 10:01

QUANT REPORT

Page 1

Operator ID: KB6656
 Output File: ^C0972::AQ
 Data File: >C0972::U4
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5,

Quant Rev: 7 Quant Time: 901003 10:39
 Injected at: 901003 10:01
 Dilution Factor: 1.00000

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ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|---------|-------|-----|
| 1) *Bromochloromethane | 11.00 | 243 | 33439 | 250.00 | NG | 92 |
| 2) Methyl chloride | 2.12 | 14 | 24563 | 177.60 | NG | 92 |
| 3) Methyl bromide | 3.17 | 41 | 10088 | 458.39 | NG | 89 |
| 4) Dichlorodifluoromethane | 3.94 | 61 | 20601 | 330.79 | NG | 95 |
| 5) Vinyl chloride | 3.98 | 62 | 28559 | 319.36 | NG | 99 |
| 6) Chloroethane | 5.14 | 92 | 18945 | 323.95 | NG | 97 |
| 7) Methylene chloride | 7.47 | 152 | 53114 | 358.46 | NG | 92 |
| 8) Acrolein | 8.25 | 172 | 105297 | 6870.91 | NG | 88 |
| 9) Acetone | 8.29 | 173 | 29974 | 389.52 | NG | 97 |
| 10) Acrylonitrile | 9.02 | 192 | 30914 | 458.88 | NG | 93 |
| 11) Carbon disulfide | 9.14 | 195 | 249726 | 321.98 | NG | 97 |
| 12) Trichlorofluoromethane | 9.80 | 212 | 145098 | 300.35 | NG | 93 |
| 13) 1,1-Dichloroethylene | 10.57 | 232 | 55695 | 246.08 | NG | 87 |
| 14) 1,1-Dichloroethane | 11.93 | 267 | 119653 | 262.60 | NG | 98 |
| 15) Tetrahydrofuran | 12.01 | 269 | 10461 | 297.63 | NG | 100 |
| 15) Tetrahydrofuran | 12.67 | 286 | 2117 | 60.23 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.71 | 287 | 52582 | 240.03 | NG | 89 |
| 17) Chloroform | 13.29 | 302 | 175991 | 279.98 | NG | 97 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.95 | 319 | 126037 | 291.78 | NG | 99 |
| 19) 1,2-Dichloroethane | 14.06 | 322 | 143848 | 295.50 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.78 | 521 | 154748 | 250.00 | NG | 95 |
| 21) Methyl ethyl ketone | 14.03 | 321 | 11158 | 517.72 | NG | 91 |
| 22) 1,1,1-Trichloroethane | 15.42 | 357 | 160852 | 344.79 | NG | 98 |
| 23) Carbon tetrachloride | 15.42 | 357 | 16473 | 38.48 | NG | 93 |
| 23) Carbon tetrachloride | 15.81 | 367 | 145452 | 339.75 | NG | 97 |
| 24) Vinyl acetate | 15.23 | 352 | 49264 | 155.92 | NG | 78 |
| 24) Vinyl acetate | 16.04 | 373 | 88161 | 279.03 | NG | 89 |
| 25) Dichlorobromomethane | 16.39 | 382 | 139050 | 319.07 | NG | 95 |
| 26) 1,2-Dichloropropane | 17.79 | 418 | 51182 | 236.69 | NG | 98 |
| 27) cis-1,3-Dichloropropylene | 18.06 | 425 | 85652 | 279.63 | NG | 91 |
| 28) Trichloroethylene | 18.64 | 440 | 57747 | 252.56 | NG | 78 |
| 29) Chlorodibromomethane | 19.26 | 456 | 83180 | 260.92 | NG | 99 |
| 30) bis(Chloromethyl)ether | 19.26 | 456 | 38771 | 298.72 | NG | 100 |
| 31) Benzene | 19.18 | 454 | 173359 | 240.54 | NG | 89 |
| 32) 1,1,2-Trichloroethane | 19.42 | 460 | 47019 | 230.69 | NG | 88 |
| 33) trans-1,3-Dichloropropylene | 19.38 | 459 | 113305 | 283.77 | NG | 85 |
| 34) 2-Chloroethylvinyl ether | 20.50 | 488 | 32319 | 209.68 | NG | 100 |
| 35) Bromoform | 22.02 | 527 | 62340 | 236.16 | NG | 97 |
| 36) *Chlorobenzene-d5 | 26.72 | 648 | 176233 | 250.00 | NG | 65 |
| 37) Methyl-iso-butyl ketone | 22.48 | 539 | 52446 | 217.46 | NG | 91 |
| 38) 2-Hexanone | 24.07 | 580 | 47055 | 201.94 | NG | 80 |
| 38) 2-Hexanone | 24.50 | 591 | 2192 | 9.41 | NG | 96 |
| 39) 1,1,2,2-Tetrachloroethane | 24.34 | 587 | 72692 | 198.54 | NG | 99 |

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QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0972::AQ
 Data File: >C0972::U4
 Name: UQA,901001,C
 Misc: QC70357US,QU70357,L,5,

Quant Rev: 7 Quant Time: 901003 10:39
 Injected at: 901003 10:01
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UQA, TCL, XUQA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 40) | Tetrachloroethylene | 24.38 | 588 | 60842 | 250.20 | NG | 98 |
| 41) | Toluene-D8 (SURR) | 25.51 | 617 | 272171 | 263.72 | NG | 93 |
| 42) | Toluene | 25.70 | 622 | 148919 | 254.58 | NG | 99 |
| 43) | Chlorobenzene | 26.83 | 651 | 201006 | 245.54 | NG | 96 |
| 44) | Ethylbenzene | 28.97 | 706 | 128615 | 261.10 | NG | 78 |
| 45) | p-Bromofluorobenzene (SURR) | 31.72 | 777 | 224676 | 308.20 | NG | 73 |
| 46) | Styrene | 33.39 | 820 | 328857 | 283.38 | NG | 98 |
| 47) | m-Xylene | 33.70 | 828 | 189326 | 293.41 | NG | 96 |
| 48) | o+p-Xylenes | 34.83 | 857 | 402974 | 604.82 | NG | 90 |

* Compound is ISTD

QUANT REPORT

Operator ID: KB6656
 Output File: ^C0977::AQ
 Data File: >C0977::U4
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5, 10UL

Quant Rev: 7 Quant Time: 901003 16:43
 Injected at: 901003 14:45
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|----------|-------|-----|
| 1) *Bromochloromethane | 10.94 | 240 | 40449 | 250.00 | NG | 95 |
| 2) Methyl chloride | 2.17 | 14 | 90814 | 542.83 | NG | 97 |
| 3) Methyl bromide | 3.22 | 41 | 11319 | 425.19 | NG | 98 |
| 4) Dichlorodifluoromethane | 3.92 | 59 | 37369 | 496.04 | NG | 93 |
| 5) Vinyl chloride | 3.96 | 60 | 55220 | 510.48 | NG | 97 |
| 6) Chloroethane | 5.16 | 91 | 37245 | 526.51 | NG | 97 |
| 7) Methylene chloride | 7.45 | 150 | 95143 | 530.82 | NG | 98 |
| 8) Acrolein | 8.31 | 172 | 244743 | 13202.43 | NG | 95 |
| 9) Acetone | 8.38 | 174 | 47100 | 506.01 | NG | 94 |
| 10) Acrylonitrile | 9.04 | 191 | 58471 | 717.52 | NG | 97 |
| 11) Carbon disulfide | 9.12 | 193 | 373170 | 397.76 | NG | 98 |
| 12) Trichlorofluoromethane | 9.74 | 209 | 293996 | 503.09 | NG | 95 |
| 13) 1,1-Dichloroethylene | 10.52 | 229 | 132241 | 483.02 | NG | 93 |
| 14) 1,1-Dichloroethane | 11.87 | 264 | 271805 | 493.15 | NG | 95 |
| 15) Tetrahydrofuran | 11.95 | 266 | 21639 | 508.97 | NG | 100 |
| 15) Tetrahydrofuran | 12.65 | 284 | 2620 | 61.62 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.69 | 285 | 125902 | 475.11 | NG | 96 |
| 17) Chloroform | 13.27 | 300 | 372092 | 489.36 | NG | 95 |
| 18) 1,2-Dichloroethane-D4 (SUWR) | 13.97 | 318 | 134320 | 257.07 | NG | 98 |
| 19) 1,2-Dichloroethane | 14.05 | 320 | 298525 | 506.96 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.77 | 519 | 216201 | 250.00 | NG | 98 |
| 21) Methyl ethyl ketone | 14.01 | 319 | 14033 | 466.04 | NG | 95 |
| 22) 1,1,1-Trichloroethane | 15.40 | 355 | 324364 | 497.66 | NG | 96 |
| 23) Carbon tetrachloride | 15.36 | 354 | 35214 | 58.87 | NG | 95 |
| 23) Carbon tetrachloride | 15.79 | 365 | 302117 | 505.11 | NG | 94 |
| 24) Vinyl acetate | 15.21 | 350 | 80279 | 181.86 | NG | 78 |
| 24) Vinyl acetate | 16.02 | 371 | 251350 | 569.41 | NG | 93 |
| 25) Dichlorobromomethane | 16.37 | 380 | 300604 | 493.71 | NG | 97 |
| 26) 1,2-Dichloropropane | 17.77 | 416 | 139502 | 461.75 | NG | 91 |
| 27) cis-1,3-Dichloropropylene | 18.04 | 423 | 209777 | 490.21 | NG | 93 |
| 28) Trichloroethylene | 18.66 | 439 | 152596 | 477.69 | NG | 87 |
| 29) Chlorodibromomethane | 19.24 | 454 | 220994 | 496.17 | NG | 97 |
| 30) bis(Chloromethyl)ether | 19.20 | 453 | 89735 | 494.86 | NG | 100 |
| 31) Benzene | 19.17 | 452 | 483500 | 480.18 | NG | 94 |
| 32) 1,1,2-Trichloroethane | 19.36 | 457 | 139077 | 488.40 | NG | 82 |
| 33) trans-1,3-Dichloropropylene | 19.36 | 457 | 281099 | 503.90 | NG | 89 |
| 34) 2-Chloroethylvinyl ether | 20.49 | 486 | 105157 | 488.32 | NG | 100 |
| 35) Bromoform | 21.96 | 524 | 192370 | 521.61 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.67 | 645 | 245701 | 250.00 | NG | 82 |
| 37) Methyl-iso-butyl ketone | 22.46 | 537 | 167817 | 499.10 | NG | 94 |
| 38) 2-Hexanone | 24.02 | 577 | 158671 | 488.41 | NG | 94 |
| 39) 1,1,2,2-Tetrachloroethane | 24.29 | 584 | 264065 | 517.32 | NG | 97 |
| 40) Tetrachloroethylene | 24.33 | 585 | 171398 | 505.56 | NG | 97 |

QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0977::AQ
 Data File: >C0977::U4
 Name: VDA,901001,C
 Misc: QC70357US,QU70357,L,5, 10UL

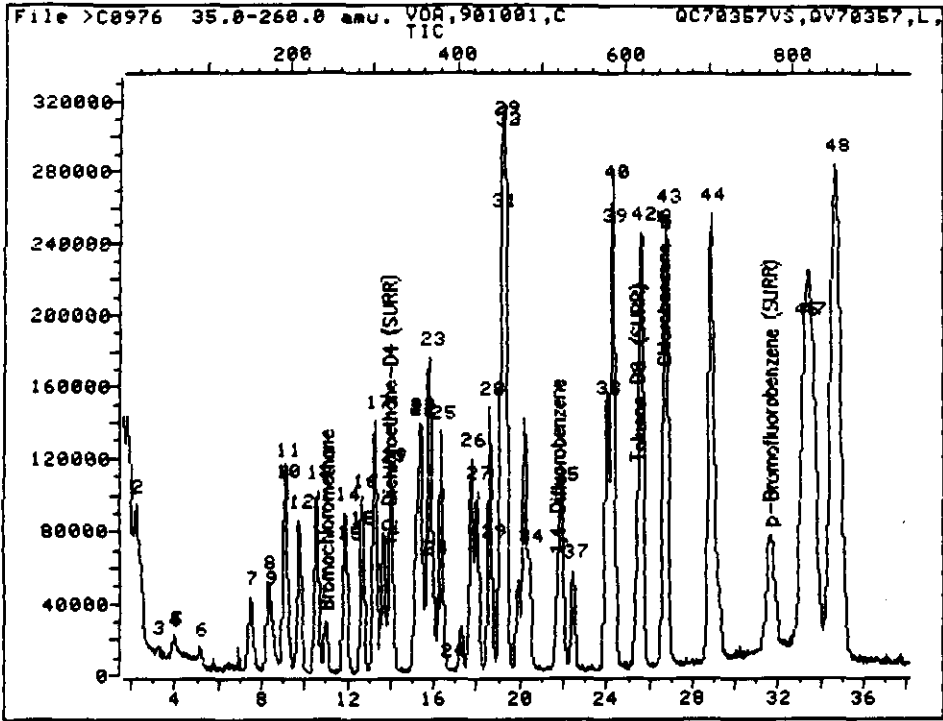
Quant Rev: 7 Quant Time: 901003 16:43
 Injected at: 901003 14:45
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VDA, TCL, XVDA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|---------|-------|----|
| 41) | Toluene-DB (SURR) | 25.49 | 615 | 367994 | 255.75 | NG | 91 |
| 42) | Toluene | 25.65 | 619 | 401656 | 492.50 | NG | 95 |
| 43) | Chlorobenzene | 26.79 | 648 | 557897 | 488.82 | NG | 94 |
| 44) | Ethylbenzene | 28.88 | 702 | 343471 | 500.13 | NG | 79 |
| 45) | p-Bromofluorobenzene (SURR) | 31.67 | 774 | 266698 | 262.40 | NG | 84 |
| 46) | Styrene | 33.31 | 816 | 833355 | 515.07 | NG | 94 |
| 47) | m-Xylene | 33.62 | 824 | 460640 | 512.04 | NG | 97 |
| 48) | o+p-Xylenes | 34.70 | 852 | 969564 | 1043.78 | NG | 87 |

* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >C0976::U4 Quant Output File: ^C0976::AQ
Name: VOA,901001,C
Misc: QC70357US,QU70357,L,5, 15UL

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901002 16:53

Operator ID: KB6656
Quant Time: 901003 16:46
Injected at: 901003 13:55

QUANT REPORT

Operator ID: KB6656
 Output File: ^C0976::AQ
 Data File: >C0976::U4
 Name: VOA,901001,C
 Misc: QC70357VS,QU70357,L,5, 15UL

Quant Rev: 7 Quant Time: 901003 16:46
 Injected at: 901003 13:55
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|---------|----------|-------|-----|
| 1) *Bromochloromethane | 10.98 | 242 | 44783 | 250.00 | NG | 99 |
| 2) Methyl chloride | 2.21 | 16 | 69987 | 377.85 | NG | 99 |
| 3) Methyl bromide | 3.22 | 42 | 13798 | 468.15 | NG | 97 |
| 4) Dichlorodifluoromethane | 3.92 | 60 | 44312 | 531.28 | NG | 93 |
| 5) Vinyl chloride | 4.00 | 62 | 64172 | 535.83 | NG | 98 |
| 6) Chloroethane | 5.16 | 92 | 49183 | 627.98 | NG | 98 |
| 7) Methylene chloride | 7.49 | 152 | 143319 | 722.22 | NG | 98 |
| 8) Acrolein | 8.31 | 173 | 370346 | 18044.55 | NG | 90 |
| 9) Acetone | 8.38 | 175 | 73037 | 708.72 | NG | 95 |
| 10) Acrylonitrile | 9.04 | 192 | 138313 | 1533.03 | NG | 90 |
| 11) Carbon disulfide | 9.12 | 194 | 1002979 | 965.61 | NG | 99 |
| 12) Trichlorofluoromethane | 9.74 | 210 | 504835 | 780.28 | NG | 96 |
| 13) 1,1-Dichloroethylene | 10.56 | 231 | 230701 | 761.10 | NG | 92 |
| 14) 1,1-Dichloroethane | 11.87 | 265 | 466532 | 764.54 | NG | 96 |
| 15) Tetrahydrofuran | 11.95 | 267 | 42526 | 903.44 | NG | 100 |
| 15) Tetrahydrofuran | 12.61 | 284 | 8088 | 171.83 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.65 | 285 | 222632 | 758.83 | NG | 98 |
| 17) Chloroform | 13.27 | 301 | 651450 | 773.84 | NG | 98 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.93 | 318 | 157254 | 271.83 | NG | 97 |
| 19) 1,2-Dichloroethane | 14.05 | 321 | 505720 | 775.71 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.73 | 519 | 235893 | 250.00 | NG | 99 |
| 21) Methyl ethyl ketone | 14.01 | 320 | 22639 | 689.08 | NG | 94 |
| 22) 1,1,1-Trichloroethane | 15.40 | 356 | 565553 | 795.28 | NG | 97 |
| 23) Carbon tetrachloride | 15.37 | 355 | 57679 | 88.38 | NG | 92 |
| 23) Carbon tetrachloride | 15.79 | 366 | 520609 | 797.74 | NG | 91 |
| 24) Vinyl acetate | 15.99 | 371 | 440600 | 914.82 | NG | 97 |
| 24) Vinyl acetate | 16.84 | 393 | 2264 | 4.70 | NG | 73 |
| 25) Dichlorobromomethane | 16.34 | 380 | 522200 | 786.07 | NG | 95 |
| 26) 1,2-Dichloropropane | 17.73 | 416 | 245780 | 745.62 | NG | 99 |
| 27) cis-1,3-Dichloropropylene | 18.00 | 423 | 362807 | 777.03 | NG | 93 |
| 28) Trichloroethylene | 18.59 | 438 | 258713 | 742.27 | NG | 87 |
| 29) Chlorodibromomethane | 18.70 | 441 | 2087 | 4.29 | NG | 57 |
| 29) Chlorodibromomethane | 19.21 | 454 | 373799 | 769.19 | NG | 96 |
| 30) bis(Chloromethyl)ether | 19.21 | 454 | 152055 | 768.54 | NG | 100 |
| 31) Benzene | 19.13 | 452 | 879768 | 800.80 | NG | 94 |
| 32) 1,1,2-Trichloroethane | 19.32 | 457 | 237165 | 763.34 | NG | 87 |
| 33) trans-1,3-Dichloropropylene | 19.32 | 457 | 469529 | 771.42 | NG | 92 |
| 34) 2-Chloroethylvinyl ether | 20.41 | 485 | 176943 | 753.08 | NG | 100 |
| 35) Bromoform | 21.96 | 525 | 303909 | 755.25 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.65 | 644 | 269685 | 250.00 | NG | 85 |
| 37) Methyl-iso-butyl ketone | 22.43 | 537 | 275155 | 745.56 | NG | 94 |
| 38) 2-Hexanone | 23.98 | 577 | 266545 | 747.49 | NG | 95 |
| 39) 1,1,2,2-Tetrachloroethane | 24.25 | 584 | 432936 | 772.72 | NG | 98 |

QUANT REPORT

Page 2

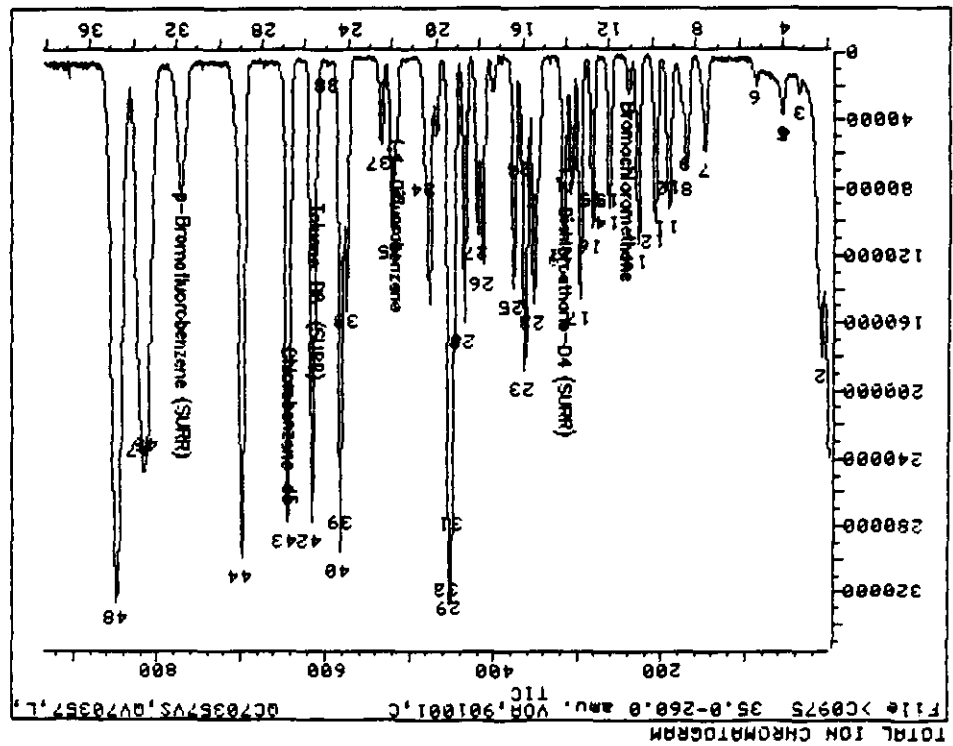
Operator ID: KB6656 Quant Rev: 7 Quant Time: 901003 16:46
 Output File: ^C0976::AQ Injected at: 901003 13:55
 Data File: >C0976::U4 Dilution Factor: 1.00000
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5, 15UL

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|---------|-------|----|
| 40) | Tetrachloroethylene | 24.33 | 586 | 267038 | 717.61 | NG | 98 |
| 41) | Toluene-DB (SURR) | 25.45 | 615 | 386124 | 244.48 | NG | 93 |
| 42) | Toluene | 25.61 | 619 | 701550 | 783.72 | NG | 97 |
| 43) | Chlorobenzene | 26.77 | 647 | 905631 | 722.93 | NG | 94 |
| 44) | Ethylbenzene | 28.86 | 701 | 548559 | 727.73 | NG | 80 |
| 45) | p-Bromofluorobenzene (SURR) | 31.62 | 772 | 280987 | 251.88 | NG | 83 |
| 46) | Styrene | 33.21 | 813 | 1283190 | 722.57 | NG | 91 |
| 47) | m-Xylene | 33.60 | 823 | 716403 | 725.52 | NG | 96 |
| 48) | o+p-Xylenes | 34.65 | 850 | 1489830 | 1461.23 | NG | 89 |

* Compound is ISTD

Data File: <C0975>:U4
 Name: VOA,901001,C
 Misc: Q070357US,Q070357,L,5, 20UL
 ID File: IC1171:US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53
 Operator ID: KB6656
 Quant Time: 901003 14:45
 Injected at: 901003 13:05



QUANT REPORT

Operator ID: KB6656
 Output File: ^C0975::AQ
 Data File: >C0975::U4
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5, 20UL

Quant Rev: 7 Quant Time: 901003 14:45
 Injected at: 901003 13:05
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|----------|-------|-----|
| 1) *Bromochloromethane | 11.07 | 240 | 33082 | 250.00 | NG | 93 |
| 2) Methyl chloride | 2.19 | 11 | 196957 | 1439.46 | NG | 97 |
| 3) Methyl bromide | 3.20 | 37 | 21780 | 1000.34 | NG | 96 |
| 4) Dichlorodifluoromethane | 3.93 | 56 | 79248 | 1286.20 | NG | 93 |
| 5) Vinyl chloride | 3.97 | 57 | 113382 | 1281.58 | NG | 94 |
| 6) Chloroethane | 5.21 | 89 | 82361 | 1423.55 | NG | 96 |
| 7) Methylene chloride | 7.50 | 148 | 200574 | 1368.24 | NG | 99 |
| 8) Acrolein | 8.36 | 170 | 442510 | 29186.55 | NG | 91 |
| 9) Acetone | 8.44 | 172 | 88336 | 1160.35 | NG | 92 |
| 10) Acrylonitrile | 9.10 | 189 | 107700 | 1615.94 | NG | 99 |
| 11) Carbon disulfide | 9.17 | 191 | 903789 | 1177.87 | NG | 99 |
| 12) Trichlorofluoromethane | 9.83 | 208 | 590028 | 1234.52 | NG | 92 |
| 13) 1,1-Dichloroethylene | 10.61 | 228 | 237397 | 1060.21 | NG | 87 |
| 14) 1,1-Dichloroethane | 11.93 | 262 | 477859 | 1060.08 | NG | 95 |
| 15) Tetrahydrofuran | 12.00 | 264 | 35476 | 1020.24 | NG | 100 |
| 15) Tetrahydrofuran | 12.66 | 281 | 8973 | 258.05 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.74 | 283 | 217086 | 1001.64 | NG | 92 |
| 17) Chloroform | 13.32 | 298 | 663299 | 1066.60 | NG | 96 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.98 | 315 | 126950 | 297.07 | NG | 98 |
| 19) 1,2-Dichloroethane | 14.10 | 318 | 524656 | 1089.39 | NG | 96 |
| 20) *1,4-Difluorobenzene | 21.74 | 515 | 183065 | 250.00 | NG | 98 |
| 21) Methyl ethyl ketone | 14.06 | 317 | 21244 | 833.22 | NG | 93 |
| 22) 1,1,1-Trichloroethane | 15.42 | 352 | 594519 | 1077.26 | NG | 97 |
| 23) Carbon tetrachloride | 15.42 | 352 | 58362 | 115.24 | NG | 99 |
| 23) Carbon tetrachloride | 15.84 | 363 | 560082 | 1105.89 | NG | 96 |
| 24) Vinyl acetate | 16.04 | 368 | 386372 | 1033.72 | NG | 97 |
| 25) Dichlorobromomethane | 16.39 | 377 | 539072 | 1045.63 | NG | 96 |
| 26) 1,2-Dichloropropane | 17.78 | 413 | 236212 | 923.39 | NG | 96 |
| 27) cis-1,3-Dichloropropylene | 18.06 | 420 | 359779 | 992.91 | NG | 93 |
| 28) Trichloroethylene | 18.64 | 435 | 265049 | 979.90 | NG | 84 |
| 29) Chlorodibromomethane | 19.26 | 451 | 380714 | 1009.49 | NG | 99 |
| 30) bis(Chloromethyl)ether | 19.26 | 451 | 161033 | 1048.79 | NG | 100 |
| 31) Benzene | 19.18 | 449 | 880604 | 1032.87 | NG | 94 |
| 32) 1,1,2-Trichloroethane | 19.37 | 454 | 242009 | 1003.71 | NG | 81 |
| 33) trans-1,3-Dichloropropylene | 19.37 | 454 | 492760 | 1043.22 | NG | 89 |
| 34) 2-Chloroethylvinyl ether | 20.46 | 482 | 172878 | 948.11 | NG | 100 |
| 35) Bromoform | 21.97 | 521 | 299168 | 958.02 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.69 | 641 | 214016 | 250.00 | NG | 80 |
| 37) Methyl-iso-butyl ketone | 22.48 | 534 | 247124 | 843.78 | NG | 96 |
| 38) 2-Hexanone | 24.03 | 574 | 234177 | 827.55 | NG | 95 |
| 38) 2-Hexanone | 24.92 | 597 | 6261 | 22.13 | NG | 44 |
| 39) 1,1,2,2-Tetrachloroethane | 24.30 | 581 | 391687 | 880.94 | NG | 96 |
| 40) Tetrachloroethylene | 24.34 | 582 | 281542 | 953.38 | NG | 97 |

QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0975::AQ
 Data File: >C0975::U4
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5, 20UL

Quant Rev: 7 Quant Time: 901003 14:45
 Injected at: 901003 13:05
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|---------|-------|----|
| 41) | Toluene-D8 (SURR) | 25.47 | 611 | 313012 | 249.74 | NG | 91 |
| 42) | Toluene | 25.66 | 616 | 760493 | 1070.55 | NG | 91 |
| 43) | Chlorobenzene | 26.81 | 644 | 993619 | 999.48 | NG | 94 |
| 44) | Ethylbenzene | 28.90 | 698 | 631612 | 1055.86 | NG | 77 |
| 45) | p-Bromofluorobenzene (SURR) | 31.66 | 769 | 238765 | 269.70 | NG | 80 |
| 46) | Styrene | 33.29 | 811 | 1506277 | 1068.82 | NG | 93 |
| 47) | m-Xylene | 33.64 | 820 | 842266 | 1074.86 | NG | 95 |
| 48) | o+p-Xylenes | 34.73 | 848 | 1756527 | 2170.94 | NG | 88 |

* Compound is ISTD

VOLATILE INTERNAL ^{8A} STANDARD AREA SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID (Standard): >C0914 Data Analyzed: 09/26/90
 Instrument ID: GC/MS C Time Analyzed: 1352
 Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) *PACK*

| | IS1(BCM) | RT | IS2(DFB) | RT | IS3(CBZ) | RT |
|----------------|----------|-------|----------|-------|----------|-------|
| | AREA # | | AREA # | | AREA # | |
| 12 HOUR STD | 70232 | 11.09 | 342312 | 21.84 | 318334 | 26.78 |
| UPPER LIMIT | 140464 | | 684624 | | 636668 | |
| LOWER LIMIT | 35116 | | 171156 | | 159167 | |
| EPA SAMPLE NO. | | | | | | |
| 01 UBLK01 | 47287 | 11.08 | 219376 | 21.83 | 199092 | 26.79 |
| 02 A4753 | 46998 | 11.19 | 224774 | 21.94 | 207367 | 26.92 |
| 03 A4755 | 51014 | 11.20 | 234507 | 21.98 | 216956 | 26.92 |
| 04 A4763 | 45139 | 11.24 | 219122 | 21.95 | 207394 | 26.88 |
| 05 A4756 | 44376 | 11.26 | 215254 | 21.96 | 204163 | 26.90 |
| 06 A4758 | 41198 | 11.30 | 192371 | 21.96 | 182527 | 26.90 |
| 07 | | | | | | |
| 08 | | | | | | |
| 09 | | | | | | |
| 10 | | | | | | |
| 11 | | | | | | |
| 12 | | | | | | |
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| 16 | | | | | | |
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| 18 | | | | | | |
| 19 | | | | | | |
| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene-d5

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

* Column used to flag internal standard area values with an asterisk.

page 1 of 1

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Lab File ID (Standard): >C0927

Data Analyzed: 09/27/90

Instrument ID: GC/MS C

Time Analyzed: 1656

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) *PALC*

| | IS1(BCM) | RT | IS2(DFB) | RT | IS3(CBZ) | RT |
|---------------------------------|----------|-------|----------|-------|----------|-------|
| | AREA # | | AREA # | | AREA # | |
| 12 HOUR STD | 40844 | 11.08 | 179047 | 21.83 | 171595 | 26.77 |
| UPPER LIMIT | 81688 | | 358094 | | 343190 | |
| LOWER LIMIT | 20422 | | 89523 | | 85798 | |
| EPA SAMPLE NO. | | | | | | |
| 01 UBLK2 | 36606 | 11.11 | 178010 | 21.86 | 167977 | 26.79 |
| 02 A4754 | 42061 | 11.10 | 196522 | 21.85 | 185259 | 26.82 |
| <i>(CA)</i> 03 <i>UBA A4757</i> | 39142 | 11.10 | 183849 | 21.84 | 173277 | 26.78 |
| <i>10-16-90</i> 04 A4759 | 42889 | 11.11 | 192985 | 21.85 | 186569 | 26.79 |
| 05 | | | | | | |
| 06 | | | | | | |
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| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene-d5

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk.

page $\frac{+}{+}$ of $\frac{+}{+}$

FORM VIII UOA

1/87 Rev.

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID (Standard): >C0961 Data Analyzed: 10/02/90
 Instrument ID: GC/MS C Time Analyzed: 1443
 Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) PACK

| | IS1(BCM) | RT | IS2(DFB) | RT | IS3(CBZ) | RT |
|----------------|----------|-------|----------|-------|----------|-------|
| | AREA # | | AREA # | | AREA # | |
| 12 HOUR STD | 45268 | 10.98 | 234953 | 21.76 | 265376 | 26.65 |
| UPPER LIMIT | 90536 | | 469906 | | 530752 | |
| LOWER LIMIT | 22634 | | 117476 | | 132688 | |
| EPA SAMPLE NO. | | | | | | |
| 01 UBLK4 | 36061 | 11.01 | 187095 | 21.79 | 221523 | 26.72 |
| 02 A4762 | 29250 | 10.96 | 152364 | 21.78 | 179637 | 26.75 |
| 03 A4761 | 28016 | 11.01 | 145859 | 21.75 | 173819 | 26.73 |
| 04 | | | | | | |
| 05 | | | | | | |
| 06 | | | | | | |
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| 18 | | | | | | |
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| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene-d5

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

* Column used to flag internal standard area values with an asterisk.

page 1 of 1

BA
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID (Standard): >C0972 Data Analyzed: 10/03/90
 Instrument ID: GC/MS C Time Analyzed: 1001
 Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) PAC/C

| | IS1(BCM) | IS2(DFB) | IS3(CBZ) |
|---------------------------------------|----------|----------|----------|
| | AREA # | AREA # | AREA # |
| | RT | RT | RT |
| 12 HOUR STD | 33439 | 154748 | 176233 |
| | 11.00 | 21.78 | 26.72 |
| UPPER LIMIT | 66878 | 309496 | 352466 |
| LOWER LIMIT | 16719 | 77374 | 88117 |
| EPA SAMPLE NO. | | | |
| 01 UBLK5 | 32787 | 159758 | 189232 |
| | 10.99 | 21.73 | 26.68 |
| 02 A4755MS | 31333 | 158395 | 191425 |
| | 11.02 | 21.80 | 26.75 |
| 03 A4755MSD | 33706 | 166042 | 193787 |
| | 10.96 | 21.74 | 26.68 |
| 04 A4760 PL ^(C) | 26590 | 126921 | 146632 |
| | 11.02 | 21.76 | 26.69 |
| 05 <u>6-16-90</u> | | | |
| 06 | | | |
| 07 | | | |
| 08 | | | |
| 09 | | | |
| 10 | | | |
| 11 | | | |
| 12 | | | |
| 13 | | | |
| 14 | | | |
| 15 | | | |
| 16 | | | |
| 17 | | | |
| 18 | | | |
| 19 | | | |
| 20 | | | |
| 21 | | | |
| 22 | | | |

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene-d5

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

* Column used to flag internal standard area values with an asterisk.

page 1 of 1

ETC

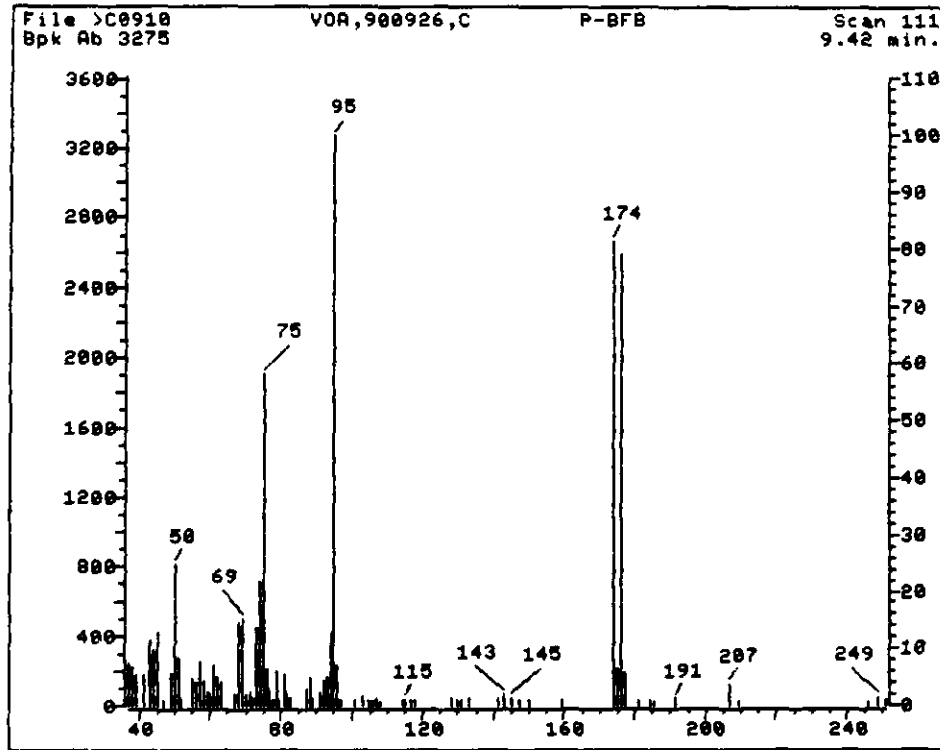
RAW QC DATA

MS Data File: >C0910::U4

Name: VOA,900926,C
Misc: P-BFB

Operator: JAB781

Date/Time: 9/26/90 8:54



MS Data File: >C0910::U4

Name: VOA,900926,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/26/90 8:54

>C0910 VOA,900926,C P-BFB
111 NRM

File: >C0910 Scan #: 111 Retn. time: 9.42

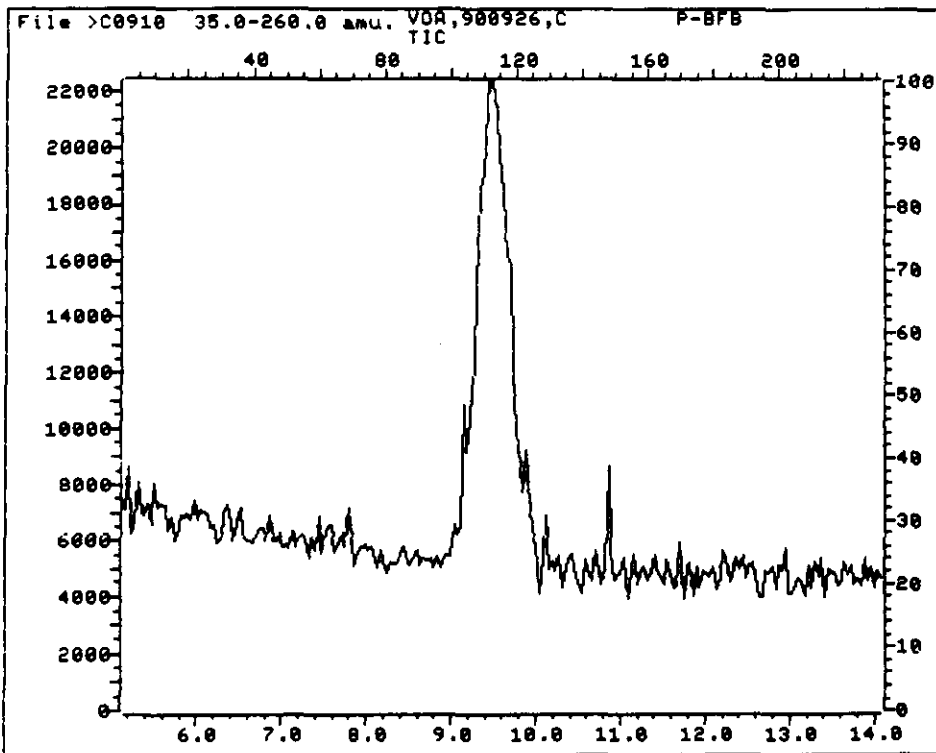
| m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. |
|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 37.00 | 7.481 | 61.00 | 6.931 | 79.70 | 1.008 | 104.70 | .977 | 145.15 | 1.527 |
| 38.00 | 6.595 | 62.00 | 5.191 | 80.90 | 5.435 | 105.30 | .794 | 147.05 | 1.130 |
| 39.00 | 5.374 | 63.00 | 4.031 | 81.90 | 2.137 | 105.60 | .977 | 149.95 | 1.008 |
| 41.10 | 5.282 | 67.10 | 2.107 | 82.50 | 1.038 | 105.80 | .916 | 158.95 | .947 |
| 43.10 | 11.328 | 68.00 | 14.443 | 83.00 | 1.435 | 106.90 | 1.374 | 173.95 | 81.405 |
| 44.00 | 9.802 | 69.00 | 15.298 | 87.00 | 3.084 | 108.00 | .733 | 175.05 | 6.504 |
| 45.00 | 12.855 | 70.00 | 2.076 | 88.00 | 5.069 | 114.60 | .885 | 175.95 | 78.870 |
| 46.90 | 1.069 | 70.30 | 1.893 | 89.00 | 1.221 | 114.90 | 1.130 | 176.95 | 5.863 |
| 49.00 | 5.802 | 71.10 | 2.351 | 91.00 | 2.382 | 116.80 | 1.130 | 180.95 | .702 |
| 50.00 | 24.947 | 72.00 | 1.221 | 92.00 | 4.275 | 117.80 | 1.008 | 184.05 | 1.069 |
| 51.00 | 8.519 | 73.00 | 13.802 | 93.00 | 5.130 | 128.05 | 1.435 | 185.55 | .824 |
| 51.90 | 1.405 | 74.00 | 21.893 | 94.00 | 12.947 | 129.85 | 1.069 | 191.15 | 1.221 |
| 55.10 | 4.733 | 75.00 | 58.290 | 95.00 | 100.000 | 131.05 | .947 | 206.95 | 3.878 |
| 56.00 | 4.061 | 76.00 | 6.321 | 96.00 | 7.023 | 132.95 | 1.496 | 209.15 | .641 |
| 57.00 | 7.847 | 77.00 | 2.321 | 97.00 | 1.160 | 140.95 | 1.221 | 245.75 | .794 |
| 58.00 | 4.305 | 78.10 | 1.191 | 101.00 | 1.160 | 141.15 | 1.252 | 248.95 | 1.405 |
| 59.10 | 2.351 | 78.90 | 5.924 | 103.00 | 1.740 | 142.85 | 1.985 | 250.95 | .977 |
| 60.00 | 2.168 | | | | | | | | |

MS Data File: >C0910::U4

Name: UOA,900926,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/26/90 8:54

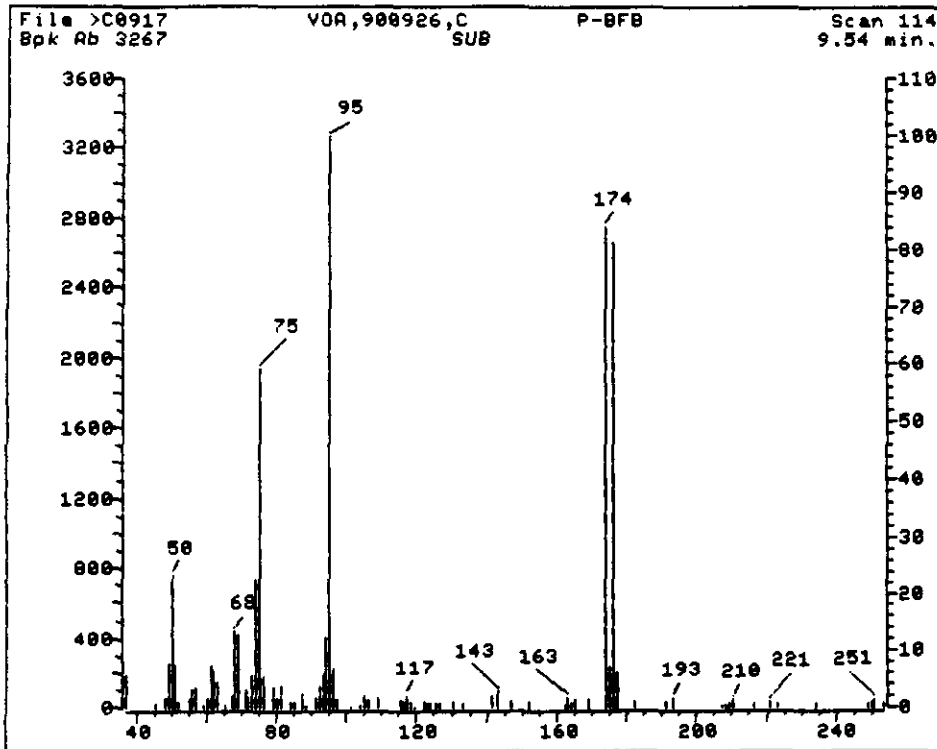


MS Data File: >C0917::U4

Name: VOA,900926,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/26/90 16:25



MS Data File: >C0917::U4

Name: VOA,900926,C Operator: JA8781 Date/Time: 9/26/90 16:25
Misc: P-BFB

/C0917 VOA,900926,C P-BFB
114 SUB NRM

File: >C0917 Scan #: 114 Retn. time: 9.54

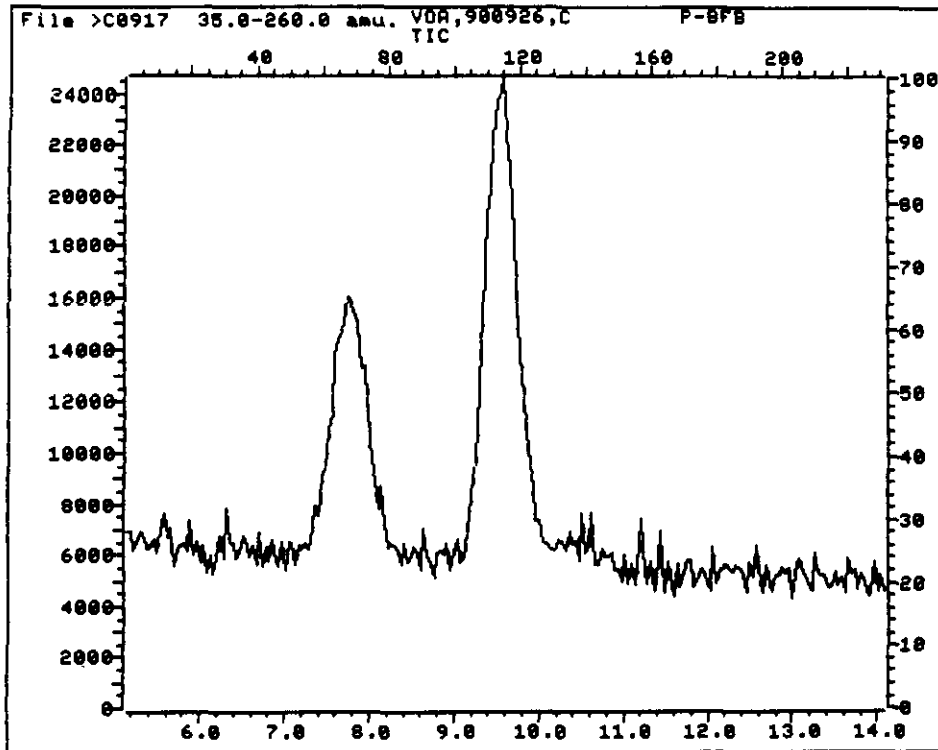
| m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. |
|-------|--------|-------|--------|--------|---------|--------|--------|--------|-------|
| 37.00 | 5.908 | 69.00 | 12.856 | 95.00 | 100.000 | 123.90 | .765 | 177.05 | 6.030 |
| 45.10 | .857 | 71.10 | 3.122 | 96.00 | 6.856 | 125.20 | .673 | 182.05 | .979 |
| 48.00 | 1.592 | 72.00 | 1.622 | 97.30 | 1.408 | 126.60 | .857 | 190.95 | .949 |
| 49.00 | 7.836 | 73.00 | 5.510 | 101.10 | .184 | 130.05 | 1.194 | 192.95 | 1.561 |
| 50.00 | 23.263 | 74.00 | 22.467 | 103.90 | .275 | 132.95 | .796 | 207.05 | .429 |
| 51.00 | 7.530 | 75.00 | 59.443 | 105.10 | 1.928 | 141.05 | 2.173 | 208.05 | .490 |
| 51.80 | 1.071 | 76.00 | 5.448 | 106.00 | 1.439 | 142.95 | 2.694 | 209.05 | .582 |
| 55.10 | 1.377 | 79.00 | 3.520 | 109.00 | 1.714 | 147.05 | 1.133 | 210.05 | 1.377 |
| 56.00 | 3.183 | 79.90 | 1.653 | 115.10 | 1.500 | 152.05 | .888 | 216.15 | .735 |
| 57.00 | 3.275 | 80.90 | 3.826 | 115.80 | 1.163 | 162.15 | .398 | 221.05 | .888 |
| 59.00 | .367 | 84.10 | .918 | 116.20 | .796 | 162.95 | 1.622 | 222.85 | .735 |
| 59.90 | 1.745 | 85.10 | .888 | 117.00 | 2.173 | 163.85 | .765 | 233.85 | .765 |
| 61.00 | 7.254 | 87.00 | 2.418 | 118.00 | 1.163 | 164.95 | 1.408 | 234.05 | .826 |
| 62.00 | 6.367 | 88.10 | .429 | 119.00 | .122 | 169.15 | 1.224 | 249.05 | .551 |
| 63.00 | 4.561 | 91.10 | 1.622 | 121.90 | .888 | 173.95 | 84.022 | 249.85 | 1.071 |
| 65.00 | .398 | 92.00 | 3.887 | 122.40 | .643 | 175.05 | 7.040 | 250.85 | 1.224 |
| 67.10 | 2.204 | 93.00 | 5.877 | 123.00 | .367 | 175.95 | 81.267 | 253.45 | .857 |
| 68.00 | 13.315 | 94.00 | 12.489 | | | | | | |

MS Data File: >C0917::U4

Name: UOA,900926,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/26/90 16:25

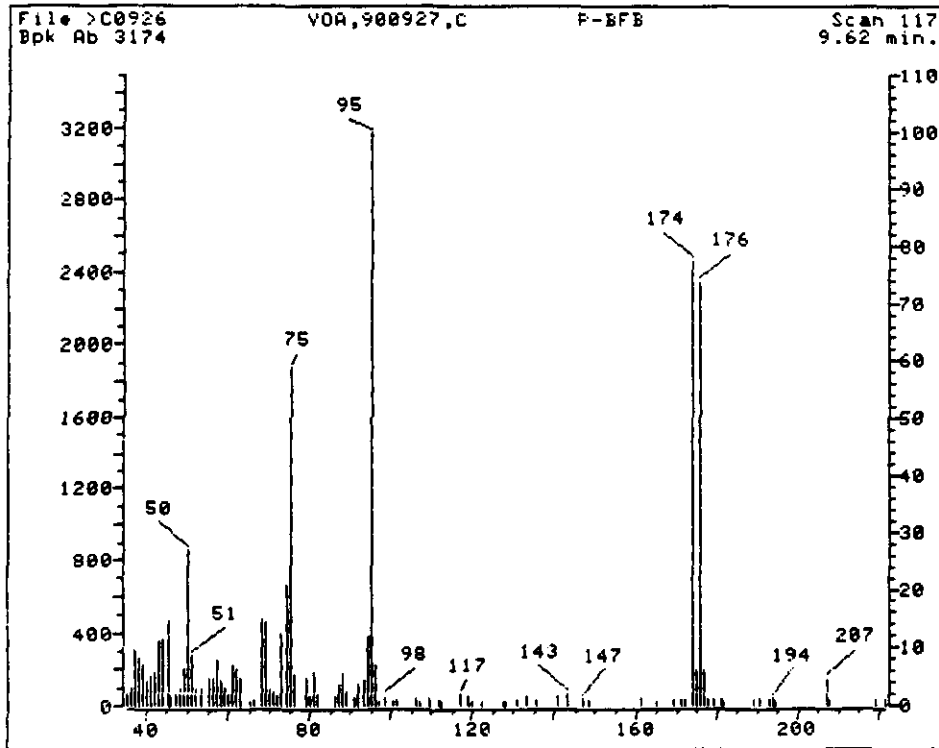


MS Data File: >C0926::U8

Name: VOA,900927,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/27/90 8:58



MS Data File: >C0926::U8

Name: VOA,900927,C Operator: JA8781 Date/Time: 9/27/90 8:58
Misc: P-BFB

>C0926 VOA,900927,C P-BFB
117 NRM

File: >C0926 Scan #: 117 Retn. time: 9.62

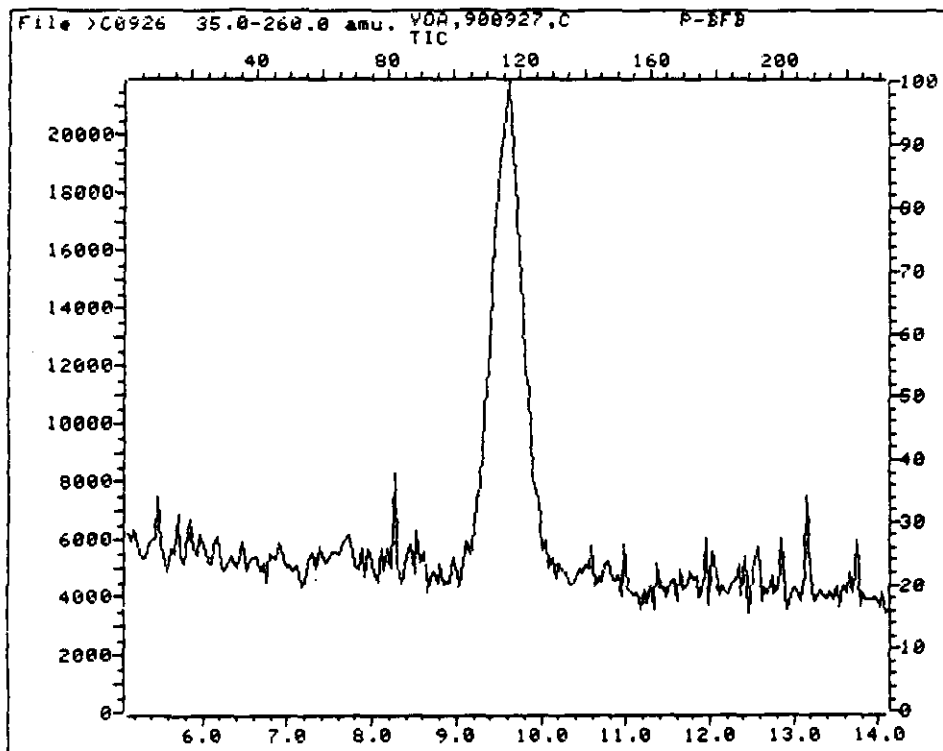
| m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. |
|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 35.30 | 2.048 | 56.00 | 4.820 | 78.90 | 4.789 | 109.20 | 1.229 | 171.95 | .977 |
| 36.00 | 3.214 | 57.00 | 8.255 | 80.10 | 1.701 | 111.80 | 1.040 | 172.25 | .914 |
| 37.00 | 9.704 | 58.00 | 4.505 | 80.90 | 5.577 | 112.40 | .630 | 173.95 | 77.347 |
| 38.00 | 8.349 | 59.00 | 2.962 | 81.90 | 2.016 | 117.00 | 1.953 | 175.05 | 5.986 |
| 39.10 | 7.089 | 60.00 | 2.016 | 85.90 | 1.638 | 119.00 | 1.796 | 175.95 | 73.661 |
| 39.90 | 4.348 | 61.00 | 7.057 | 86.90 | 3.749 | 119.80 | .756 | 176.95 | 5.923 |
| 41.10 | 5.230 | 62.00 | 6.364 | 88.00 | 5.640 | 122.30 | .819 | 177.95 | 1.197 |
| 42.00 | 5.577 | 63.00 | 4.789 | 89.10 | 2.331 | 127.25 | .819 | 179.45 | .945 |
| 43.00 | 11.374 | 65.10 | .851 | 90.90 | 1.481 | 128.05 | .756 | 180.95 | 1.134 |
| 44.00 | 11.815 | 66.00 | .882 | 91.90 | 3.718 | 130.95 | .977 | 181.75 | .788 |
| 45.10 | 14.808 | 66.30 | .851 | 93.00 | 4.505 | 132.95 | 1.670 | 189.05 | .882 |
| 45.90 | 1.701 | 68.00 | 15.028 | 94.00 | 12.130 | 135.45 | 1.040 | 190.55 | 1.071 |
| 47.10 | 2.048 | 69.00 | 14.556 | 95.00 | 100.000 | 140.95 | 1.638 | 190.85 | 1.260 |
| 47.90 | 2.615 | 70.00 | 2.804 | 96.00 | 7.120 | 142.95 | 1.953 | 193.05 | 1.008 |
| 49.00 | 6.301 | 71.10 | 2.268 | 97.10 | .851 | 146.85 | 1.292 | 193.85 | 1.323 |
| 50.00 | 27.221 | 72.00 | 1.575 | 98.30 | 1.512 | 148.25 | .914 | 194.55 | .756 |
| 51.00 | 8.727 | 73.00 | 12.508 | 100.20 | .662 | 161.25 | 1.292 | 207.05 | 4.474 |
| 52.00 | 2.615 | 74.00 | 20.983 | 101.20 | .977 | 164.95 | .693 | 207.85 | 1.103 |
| 53.10 | 3.088 | 75.00 | 59.074 | 106.00 | 1.260 | 169.45 | .945 | 219.05 | .977 |
| 55.00 | 4.789 | 76.00 | 5.325 | 106.90 | .788 | 171.15 | .977 | 221.55 | 1.134 |

MS Data File: >C0926::UB

Name: VOA,900927,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/27/90 8:58

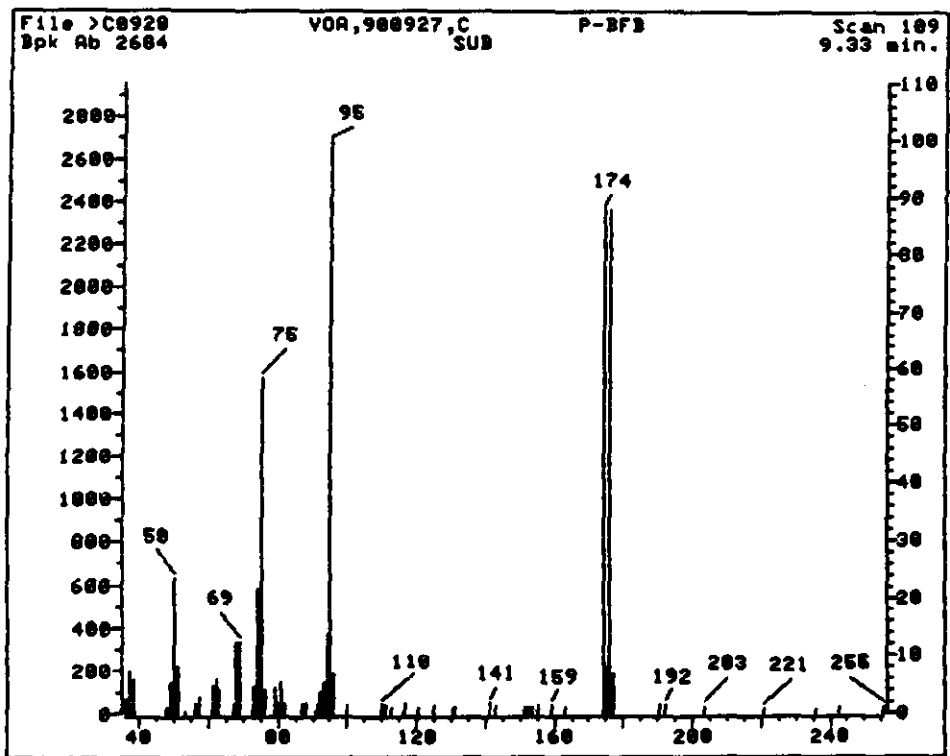


MS Data File: >C0928::U4

Name: UDA,900927,C
Misc: P-BFB

Operator: MGRMS

Date/Time: 9/27/90 17:41



MS Data File: >C0928::U4

Name: VOA,900927,C
Misc: P-BFB

Operator: MGRMS

Date/Time: 9/27/90 17:41

>C0928 VOA,900927,C P-BFB
109 SUB NRM

File: >C0928 Scan #: 109 Retn. time: 9.33

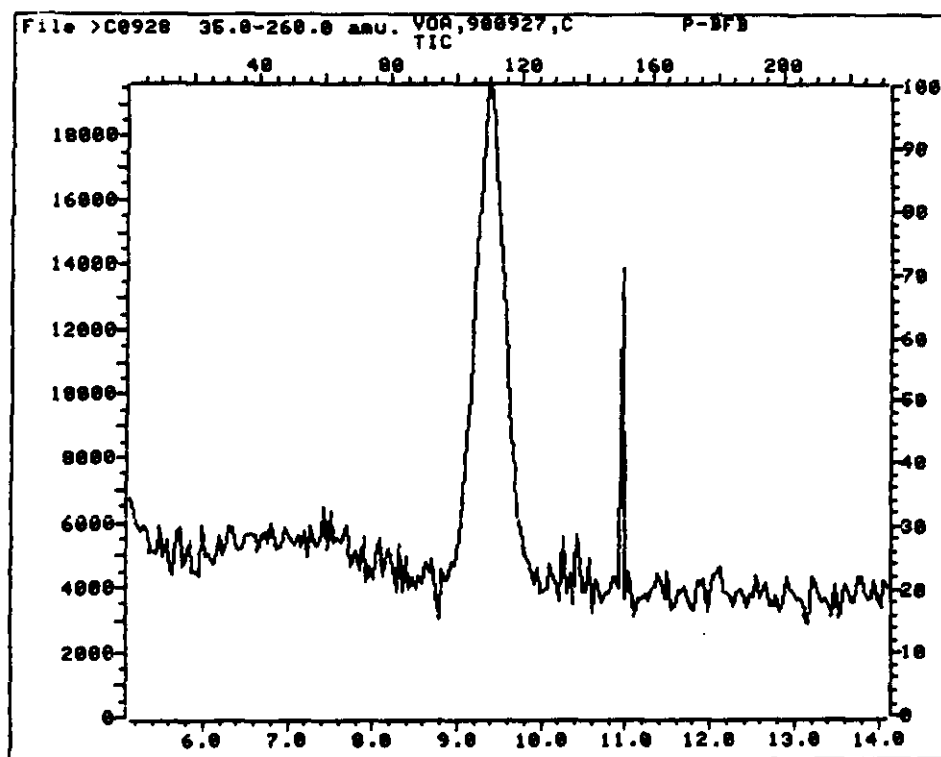
| m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. |
|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 36.00 | 2.496 | 63.00 | 3.949 | 88.00 | 2.086 | 116.80 | 1.639 | 159.05 | 1.528 |
| 37.00 | 7.377 | 67.20 | 1.677 | 90.80 | .447 | 120.60 | .894 | 162.85 | 1.006 |
| 38.00 | 6.222 | 68.00 | 12.444 | 91.20 | 1.677 | 125.20 | 1.379 | 173.95 | 87.928 |
| 39.00 | .522 | 69.00 | 12.556 | 92.00 | 3.875 | 130.55 | .782 | 174.95 | 7.787 |
| 47.50 | 1.155 | 73.00 | 4.657 | 93.00 | 5.477 | 130.75 | .894 | 175.95 | 87.779 |
| 49.00 | 5.551 | 74.00 | 21.610 | 94.00 | 13.748 | 131.05 | 1.006 | 177.05 | 6.781 |
| 50.00 | 23.547 | 75.00 | 58.569 | 95.00 | 100.000 | 141.05 | 2.012 | 190.35 | 1.230 |
| 51.00 | 7.973 | 76.00 | 3.912 | 96.00 | 6.706 | 143.05 | 1.528 | 192.05 | 1.490 |
| 53.10 | .447 | 78.90 | 4.247 | 99.50 | 1.490 | 151.45 | 1.118 | 203.35 | 1.155 |
| 56.00 | 1.751 | 79.90 | 1.937 | 110.10 | 1.677 | 152.15 | .969 | 220.95 | .782 |
| 57.10 | 2.720 | 80.90 | 5.477 | 110.90 | 1.267 | 153.15 | 1.155 | 235.95 | .782 |
| 60.90 | 4.844 | 81.80 | 1.639 | 112.50 | .931 | 153.35 | 1.155 | 242.35 | 1.006 |
| 62.00 | 5.961 | 87.00 | 1.826 | 116.00 | .112 | 154.95 | 1.267 | 255.45 | 1.080 |

MS Data File: >C0928::U4

Name: UOA,900927,C
Misc: P-BFB

Operator: MGRMS

Date/Time: 9/27/90 17:41

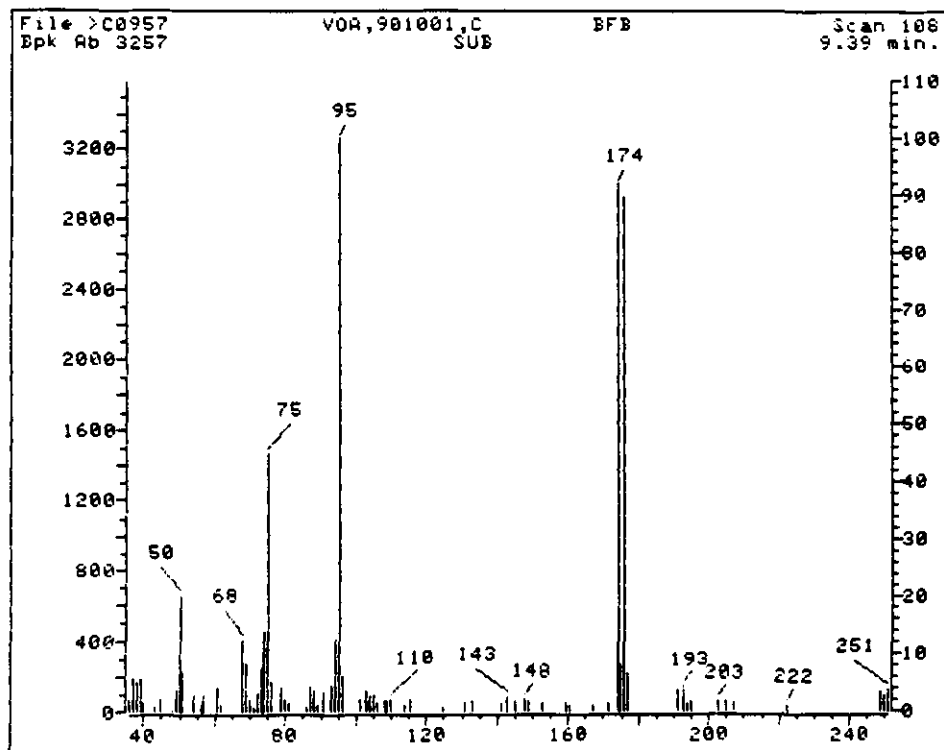


MS Data File: >C0957::U8

Name: VDA,901001,C
Misc: BFB

Operator: JA8781

Date/Time: 10/02/90 11:48



MS Data File: >C0957::U8

Name: UOA,901001,C
Misc: BFB

Operator: JA8781

Date/Time: 10/02/90 11:48

>C0957 UOA,901001,C BFB
 108 SUB NRM

File: >C0957 Scan #: 108 Retn. time: 9.39

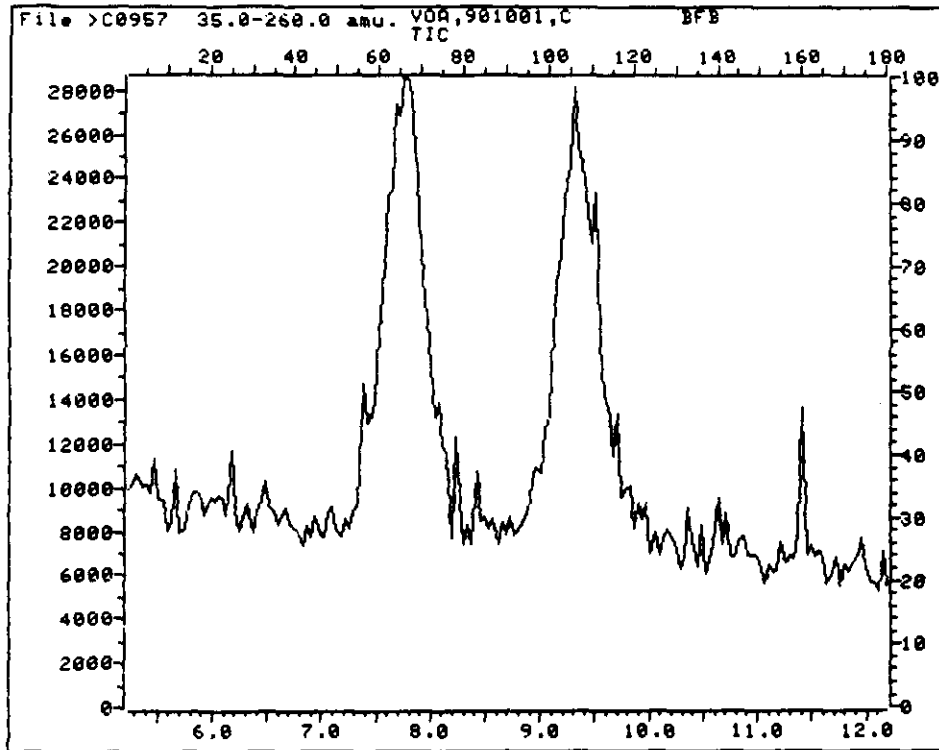
| m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. |
|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 36.00 | 2.119 | 68.00 | 12.558 | 91.00 | 3.377 | 115.20 | 1.904 | 173.95 | 92.017 |
| 37.00 | 5.711 | 69.00 | 8.443 | 93.00 | 4.483 | 124.80 | .645 | 174.95 | 8.014 |
| 38.00 | 5.097 | 69.80 | 2.180 | 94.00 | 12.312 | 130.80 | 1.320 | 175.95 | 89.561 |
| 39.00 | 5.834 | 71.00 | .706 | 95.00 | 100.000 | 132.90 | 1.597 | 176.95 | 6.478 |
| 39.90 | 1.412 | 72.00 | 3.132 | 96.00 | 6.049 | 141.45 | 1.228 | 191.05 | 3.562 |
| 43.00 | .553 | 73.00 | 7.277 | 101.00 | 2.149 | 142.85 | 2.487 | 192.95 | 4.237 |
| 45.00 | 1.965 | 74.00 | 13.755 | 102.90 | 3.531 | 145.05 | 1.873 | 193.95 | 1.351 |
| 48.10 | .154 | 75.00 | 45.164 | 104.00 | 2.272 | 147.95 | 2.088 | 194.75 | 1.873 |
| 48.90 | 3.776 | 76.00 | 4.943 | 105.10 | 2.733 | 148.95 | 1.842 | 202.85 | 1.811 |
| 50.00 | 20.233 | 78.90 | 4.022 | 106.20 | 1.412 | 152.95 | 1.228 | 205.05 | 1.597 |
| 51.00 | 6.785 | 79.80 | 2.149 | 108.50 | 1.689 | 159.15 | .123 | 207.05 | 1.320 |
| 54.00 | 2.671 | 80.90 | 1.259 | 108.80 | 1.719 | 159.35 | 1.535 | 222.15 | 1.075 |
| 56.00 | 1.197 | 85.90 | .860 | 109.00 | 1.627 | 160.45 | 1.167 | 248.95 | 3.408 |
| 57.00 | 2.733 | 87.00 | 4.513 | 110.00 | 2.057 | 167.15 | 1.075 | 250.05 | 2.671 |
| 60.90 | 3.991 | 88.00 | 3.592 | 113.70 | 1.075 | 171.45 | 1.290 | 250.95 | 3.869 |
| 62.00 | .952 | 89.00 | 1.167 | | | | | | |

MS Data File: >C0957::U8

Name: VDA,901001,C
Misc: BFB

Operator: JAB781

Date/Time: 10/02/90 11:48

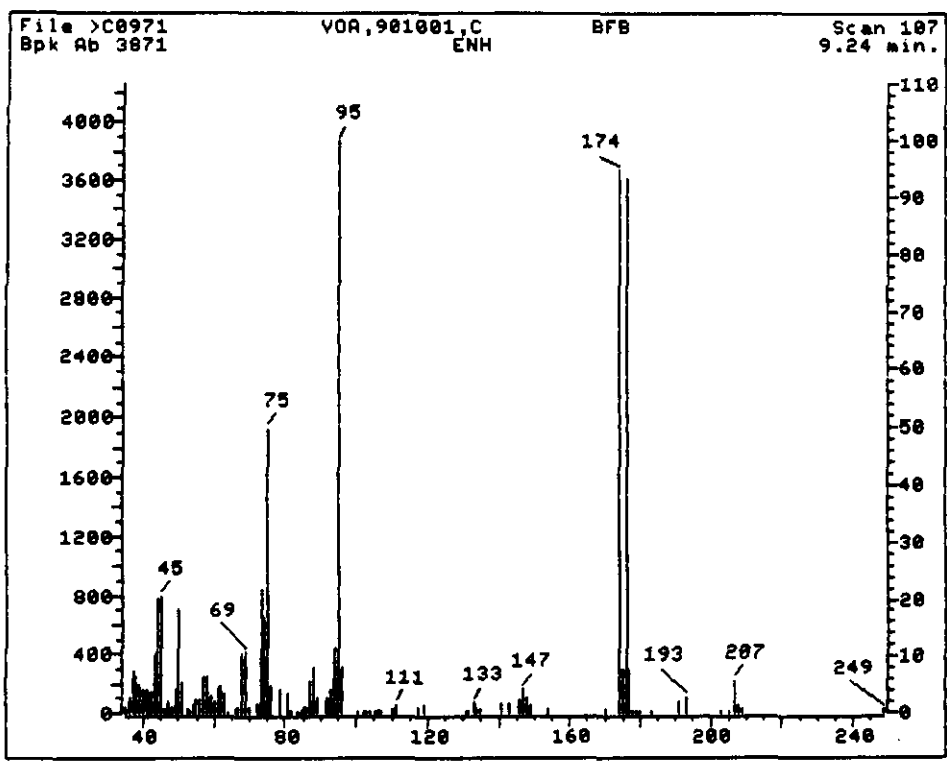


MS Data File: >C0971::U5

Name: VOA,901001,C
Misc: BFB

Operator: KB6656

Date/Time: 10/03/90 9:30



MS Data File: >C0971::U5

Name: UDA,901001,C Operator: KB6656 Date/Time: 10/03/90 9:30
Misc: BFB

>C0971 UDA,901001,C BFB
107 NRM ENH

File: >C0971 Scan #: 107 Retn. time: 9.24

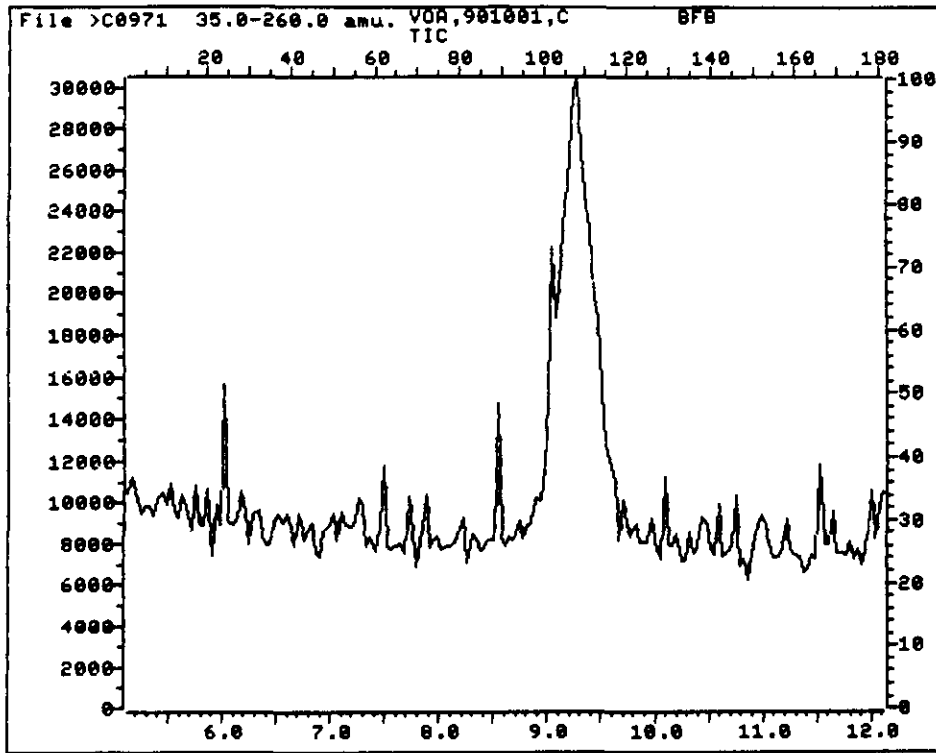
| m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. | m/z | Int. |
|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 35.10 | 1.121 | 54.90 | 2.423 | 78.80 | 4.128 | 105.60 | .377 | 148.05 | 2.640 |
| 35.90 | 2.743 | 55.90 | 2.392 | 80.90 | 3.250 | 105.80 | .356 | 149.05 | 1.271 |
| 37.00 | 7.316 | 57.00 | 6.494 | 81.50 | .232 | 106.50 | .269 | 153.85 | .661 |
| 38.00 | 6.396 | 57.90 | 6.256 | 82.00 | .393 | 107.10 | .382 | 170.15 | .568 |
| 39.00 | 5.228 | 59.10 | 3.079 | 83.50 | .258 | 110.10 | .599 | 173.95 | 94.932 |
| 39.90 | 4.133 | 59.90 | 2.191 | 84.50 | .284 | 110.40 | .372 | 174.95 | 7.889 |
| 41.00 | 4.112 | 61.00 | 4.422 | 85.00 | .951 | 111.10 | 1.503 | 175.95 | 93.521 |
| 42.00 | 3.653 | 61.90 | 4.707 | 85.90 | 1.142 | 111.40 | .537 | 176.95 | 7.579 |
| 43.00 | 10.297 | 63.00 | 3.270 | 87.00 | 5.735 | 117.10 | .914 | 177.95 | .517 |
| 44.00 | 20.149 | 64.10 | .196 | 88.00 | 8.209 | 118.80 | 1.209 | 178.85 | .372 |
| 45.00 | 20.660 | 65.90 | .837 | 88.90 | 2.692 | 130.90 | .336 | 180.25 | .248 |
| 45.90 | .770 | 66.90 | .997 | 92.00 | 2.749 | 131.20 | .258 | 183.45 | .511 |
| 47.00 | 2.067 | 68.00 | 10.519 | 92.90 | 4.180 | 131.50 | .232 | 190.85 | 2.206 |
| 48.20 | 1.199 | 69.00 | 10.710 | 94.00 | 11.314 | 133.00 | 2.025 | 192.95 | 2.795 |
| 49.00 | 4.221 | 70.20 | .883 | 95.00 | 100.000 | 133.80 | .610 | 202.85 | .253 |
| 49.90 | 18.056 | 71.90 | 1.679 | 96.00 | 7.992 | 134.80 | .672 | 205.15 | .449 |
| 51.00 | 5.409 | 73.00 | 21.766 | 100.70 | .315 | 141.05 | 1.875 | 206.95 | 5.884 |
| 52.60 | .599 | 74.00 | 16.956 | 101.90 | .253 | 142.95 | 1.571 | 207.95 | 1.297 |
| 53.00 | .387 | 75.00 | 49.752 | 102.90 | .537 | 145.95 | 2.304 | 209.15 | .873 |
| 53.90 | 1.374 | 75.90 | 4.686 | 103.50 | .269 | 146.95 | 4.402 | 248.75 | .594 |

MS Data File: >C0971::U5

Name: VOA,901001,C
Misc: BFB

Operator: KB6656

Date/Time: 10/03/90 9:30



1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

UBLK01

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: QC70357U

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0918

Level: (low/med) LOW

Date Received: ~~10/17/90~~ 09/26/90

% Moisture: not dec.

Date Analyzed: 09/26/90 ^{CA} 10-16-90

Column: (pack/cap) PACK

Dilution Factor: 1

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

| | | | |
|------------|----------------------------|----|---|
| 74-87-3 | Chloromethane | 10 | U |
| 74-83-9 | Bromomethane | 10 | U |
| 75-01-4 | Vinyl Chloride | 10 | U |
| 75-00-3 | Chloroethane | 10 | U |
| 75-09-2 | Methylene Chloride | 5 | U |
| 67-64-1 | Acetone | 10 | U |
| 75-15-0 | Carbon Disulfide | 5 | U |
| 75-35-4 | 1,1-Dichloroethene | 5 | U |
| 75-34-3 | 1,1-Dichloroethane | 5 | U |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | U |
| 67-66-3 | Chloroform | 5 | U |
| 107-06-2 | 1,2-Dichloroethane | 5 | U |
| 78-93-3 | 2-Butanone | 10 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | U |
| 56-23-5 | Carbon Tetrachloride | 5 | U |
| 108-05-4 | Vinyl Acetate | 10 | U |
| 75-27-4 | Bromodichloromethane | 5 | U |
| 78-87-5 | 1,2-Dichloropropane | 5 | U |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | U |
| 79-01-6 | Trichloroethene | 5 | U |
| 124-48-1 | Dibromochloromethane | 5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | U |
| 71-43-2 | Benzene | 5 | U |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | U |
| 75-25-2 | Bromoform | 5 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | U |
| 591-78-6 | 2-Hexanone | 10 | U |
| 127-18-4 | Tetrachloroethene | 5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | U |
| 108-88-3 | Toluene | 5 | U |
| 108-90-7 | Chlorobenzene | 5 | U |
| 100-41-4 | Ethylbenzene | 5 | U |
| 100-42-5 | Styrene | 5 | U |
| 1330-20-7 | Xylene (total) | 5 | U |

12
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETCNS Contract: _____

VBLK01

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: GC70357

Sample wt/vol: 5 (g/mL) HL Lab File ID: 7C0918

Level: (low/med) LOW Date Received: 09/26/90

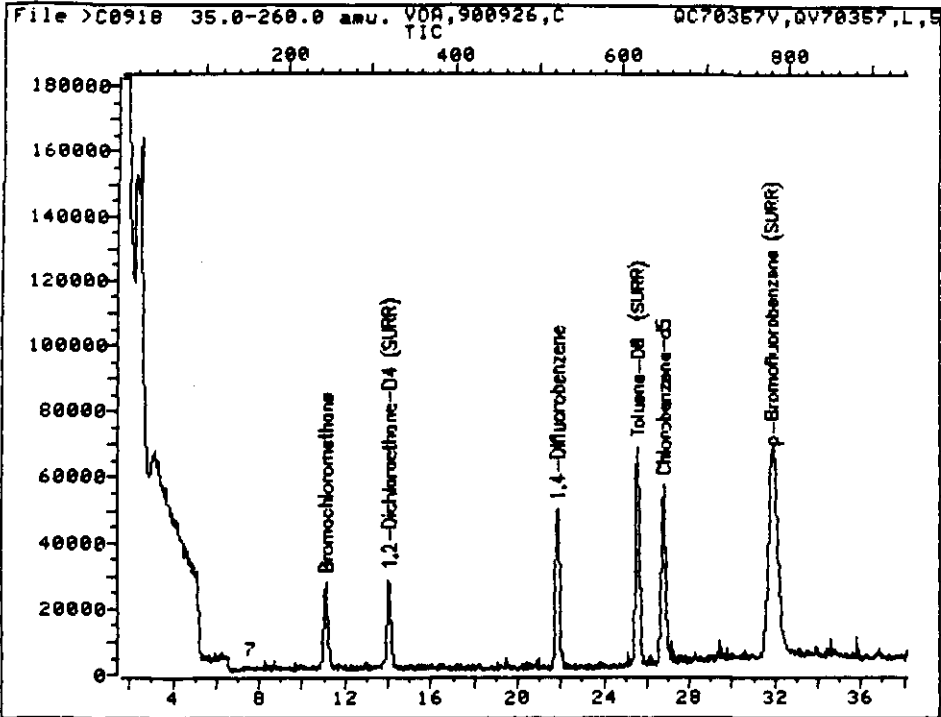
% Moisture: not dec. _____ Date Analyzed: 09/26/90

Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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| 30. | | | | |

TOTAL ION CHROMATOGRAM



Data File: >C0918::U0 Quant Output File: ^C0918::AQ
 Name: UOA,900926,C
 Misc: QC70357V,QV70357,L,5,, BLANK

Id File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 16:16

Operator ID: JA8781
 Quant Time: 900926 18:05
 Injected at: 900926 17:25

QUANT REPORT

Page 1

Operator ID: JA8781
 Output File: ^C0918::AQ
 Data File: >C0918::U0
 Name: VOA,900926,C

Quant Rev: 7 Quant Time: 900926 18:05
 Injected at: 900926 17:25
 Dilution Factor: 1.00000

Misc: QC70357U,QU70357,L,5,, BLANK

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 16:16

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.08 | 245 | 47287 | 250.00 | NG | 97 |
| 7) Methylene chloride | 7.48 | 152 | 3530 | 11.20 | NG | 96 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.03 | 321 | 157251 | 238.22 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.83 | 522 | 219376 | 250.00 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.79 | 649 | 199092 | 250.00 | NG | 80 |
| 41) Toluene-D8 (SURR) | 25.55 | 618 | 295439 | 251.96 | NG | 93 |
| 45) p-Bromofluorobenzene (SURR) | 31.83 | 779 | 228780 | 264.05 | NG | 85 |

* Compound is ISTD

CD
 10-17-90

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

UBLK2

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: QC70357U2

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0929

Level: (low/med) LOW

Date Received: ~~10/1/90~~ ^{09/27/90} *CP*

% Moisture: not dec.

Date Analyzed: 09/27/90 ¹⁰⁻¹⁶⁻⁹⁰

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | | Q |
|------------|----------------------------|----------------------|------|----|
| | | (ug/L or ug/Kg) | UG/L | |
| 74-87-3 | Chloromethane | | 10 | 10 |
| 74-83-9 | Bromomethane | | 10 | 10 |
| 75-01-4 | Vinyl Chloride | | 10 | 10 |
| 75-00-3 | Chloroethane | | 10 | 10 |
| 75-09-2 | Methylene Chloride | | 5 | 10 |
| 67-64-1 | Acetone | | 10 | 10 |
| 75-15-0 | Carbon Disulfide | | 5 | 10 |
| 75-35-4 | 1,1-Dichloroethane | | 5 | 10 |
| 75-34-3 | 1,1-Dichloroethane | | 5 | 10 |
| 540-59-0 | 1,2-Dichloroethene (total) | | 5 | 10 |
| 67-66-3 | Chloroform | | 5 | 10 |
| 107-06-2 | 1,2-Dichloroethane | | 5 | 10 |
| 78-93-3 | 2-Butanone | | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | | 5 | 10 |
| 56-23-5 | Carbon Tetrachloride | | 5 | 10 |
| 108-05-4 | Vinyl Acetate | | 10 | 10 |
| 75-27-4 | Bromodichloromethane | | 5 | 10 |
| 78-87-5 | 1,2-Dichloropropane | | 5 | 10 |
| 10061-01-5 | cis-1,3-Dichloropropene | | 5 | 10 |
| 79-01-6 | Trichloroethene | | 5 | 10 |
| 124-48-1 | Dibromochloromethane | | 5 | 10 |
| 79-00-5 | 1,1,2-Trichloroethane | | 5 | 10 |
| 71-43-2 | Benzene | | 5 | 10 |
| 10061-02-6 | trans-1,3-Dichloropropene | | 5 | 10 |
| 75-25-2 | Bromoform | | 5 | 10 |
| 108-10-1 | 4-Methyl-2-Pentanone | | 10 | 10 |
| 591-78-6 | 2-Hexanone | | 10 | 10 |
| 127-18-4 | Tetrachloroethene | | 5 | 10 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5 | 10 |
| 108-88-3 | Toluene | | 5 | 10 |
| 108-90-7 | Chlorobenzene | | 5 | 10 |
| 100-41-4 | Ethylbenzene | | 5 | 10 |
| 100-42-5 | Styrene | | 5 | 10 |
| 1330-20-7 | Xylene (total) | | 5 | 10 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETC NJ Contract: _____

VBLK 2

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: QC70357

Sample wt/vol: 5 (g/mL) ML Lab File ID: 7C0929

Level: (low/med) LOW Date Received: 09/27/90

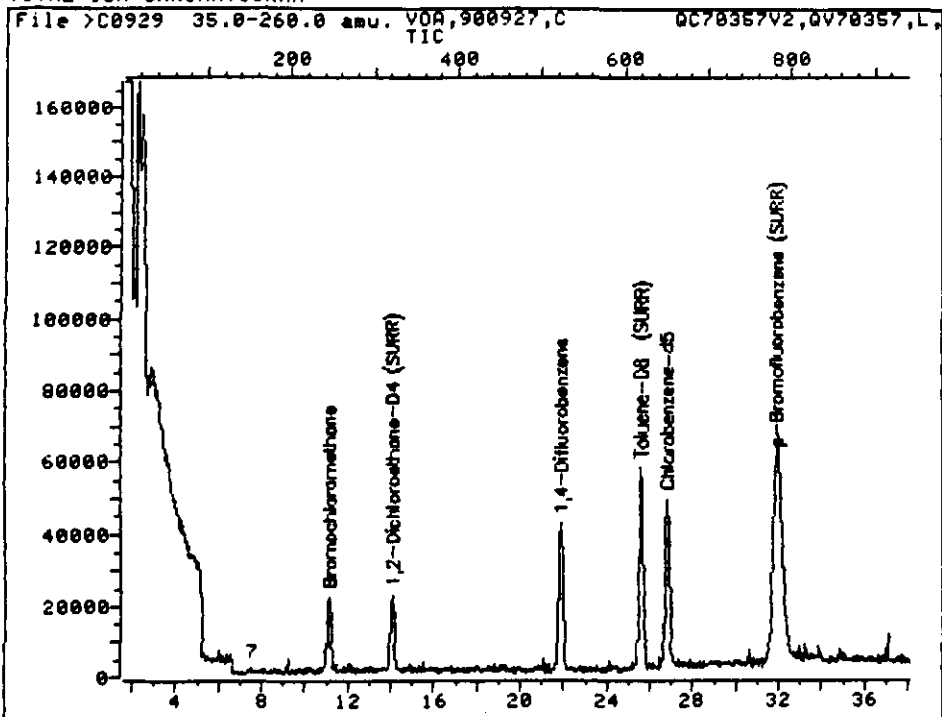
% Moisture: not dec. _____ Date Analyzed: 09/27/90

Column: (pack/cap) PACIS Dilution Factor: 1

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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TOTAL ION CHROMATOGRAM



Data File: >C0929::U4

Quant Output File: ^C0929::AQ

Name: VOA,900927,C

Misc: QC70357V2,QU70357,L,5,, BLANK

Id File: IC1171::US

Title: IFB, PP/VOA, TCL, XVOA13

Last Calibration: 900927 17:41

Operator ID: MGRMS

Quant Time: 900927 18:53

Injected at: 900927 18:14

QUANT REPORT

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900927 18:53
 Output File: ^C0929::AQ Injected at: 900927 18:14
 Data File: >C0929::U4 Dilution Factor: 1.00000
 Name: VOA,900927,C
 Misc: QC70357U2,QU70357,L,5,, BLANK

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900927 17:41

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.11 | 245 | 36606 | 250.00 | NG | 94 |
| 7) Methylene chloride | 7.47 | 151 | 3941 | 16.21 | NG | 82 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.06 | 321 | 128055 | 270.65 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.86 | 522 | 178010 | 250.00 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.79 | 649 | 167977 | 250.00 | NG | 82 |
| 41) Toluene-D8 (SURR) | 25.58 | 618 | 243972 | 248.68 | NG | 92 |
| 45) p-Bromofluorobenzene (SURR) | 31.88 | 780 | 195124 | 257.17 | NG | 85 |

* Compound is ISTD

CA
10-16-90.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLK4

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: QC70357U4

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0963

Level: (low/med) LOW

Date Received: ~~09/11/90~~ 10/02/90

% Moisture: not dec.

Date Analyzed: 10/02/90

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|----|
| 74-87-3 | Chloromethane | 10 | 1U |
| 74-83-9 | Bromomethane | 10 | 1U |
| 75-01-4 | Vinyl Chloride | 10 | 1U |
| 75-00-3 | Chloroethane | 10 | 1U |
| 75-09-2 | Methylene Chloride | 5 | 1U |
| 67-64-1 | Acetone | 10 | 1U |
| 75-15-0 | Carbon Disulfide | 5 | 1U |
| 75-35-4 | 1,1-Dichloroethene | 5 | 1U |
| 75-34-3 | 1,1-Dichloroethane | 5 | 1U |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | 1U |
| 67-66-3 | Chloroform | 5 | 1U |
| 107-06-2 | 1,2-Dichloroethane | 5 | 1U |
| 78-93-3 | 2-Butanone | 10 | 1U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | 1U |
| 56-23-5 | Carbon Tetrachloride | 5 | 1U |
| 108-05-4 | Vinyl Acetate | 10 | 1U |
| 75-27-4 | Bromodichloromethane | 5 | 1U |
| 78-87-5 | 1,2-Dichloropropane | 5 | 1U |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | 1U |
| 79-01-6 | Trichloroethene | 5 | 1U |
| 124-48-1 | Dibromochloromethane | 5 | 1U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | 1U |
| 71-43-2 | Benzene | 5 | 1U |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | 1U |
| 75-25-2 | Bromoform | 5 | 1U |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 1U |
| 591-78-6 | 2-Hexanone | 10 | 1U |
| 127-18-4 | Tetrachloroethene | 5 | 1U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | 1U |
| 108-88-3 | Toluene | 5 | 1U |
| 108-90-7 | Chlorobenzene | 5 | 1U |
| 100-41-4 | Ethylbenzene | 5 | 1U |
| 100-42-5 | Styrene | 5 | 1U |
| 1330-20-7 | Xylene (total) | 5 | 1U |

18
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLK04

Lab Name: ETC NJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDS No.: _____

Matrix: (soil/water) WATER Lab Sample ID: QC70257

Sample wt/vol: 5 (g/mL) ML Lab File ID: >C0963

Level: (low/med) LOW Date Received: 10/02/90

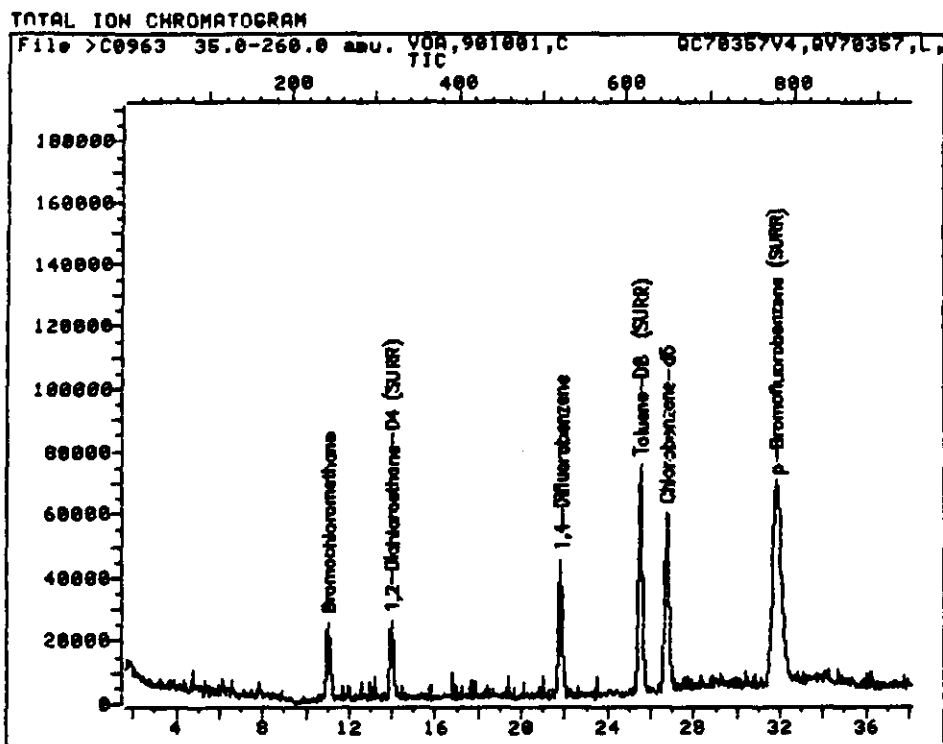
Moisture: not dec. _____ Date Analyzed: 10/02/90

Column: (pack/cap) PACK Dilution Factor: 1

CONCENTRATION UNITS:
 (ug/L or ug/kg) ug/L

Number TICs found: 0

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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Data File: >C0963::U0 Quant Output File: ^C0963::AQ
 Name: UBLK4
 Misc: QC70357V4,QU70357,L:M4,5,, BLANK

Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

Operator ID: JA8781
 Quant Time: 901002 17:10
 Injected at: 901002 16:21

QUANT REPORT

Operator ID: JA8781
 Output File: ^C0963::AQ
 Data File: >C0963::U0
 Name: UBLK4
 Misc: QC70357U4,QU70357,L:M4,5,, BLANK

Quant Rev: 7 Quant Time: 901002 17:10
 Injected at: 901002 16:21
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XV0A13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|---------|--------|-------|----|
| 1) | *Bromochloromethane | 11.01 | 243 | 36061 | 250.00 | NG | 99 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.95 | 319 | 122247 | 262.43 | NG | 91 |
| 20) | *1,4-Difluorobenzene | 21.79 | 521 | 187095 | 250.00 | NG | 99 |
| 36) | *Chlorobenzene-d5 | 26.72 | 646 | 221523 | 250.00 | NG | 76 |
| 41) | Toluene-D8 (SURR) | 25.51 | 617 | 327258M | 252.26 | NG | 88 |
| 45) | p-Bromofluorobenzene (SURR) | 31.80 | 777 | 255405 | 278.72 | NG | 80 |

* Compound is ISTD

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

UBLK5

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: QC7035705

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0979

Level: (low/med) LOW

Date Received: 10/03/90

% Moisture: not dec.

Date Analyzed: 10/03/90

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | |
|------------|----------------------------|----------------------|------|
| | | (ug/L or ug/Kg) | UG/L |
| 74-87-3 | Chloromethane | 10 | 10 |
| 74-83-9 | Bromomethane | 10 | 10 |
| 75-01-4 | Vinyl Chloride | 10 | 10 |
| 75-00-3 | Chloroethane | 10 | 10 |
| 75-09-2 | Methylene Chloride | 5 | 10 |
| 67-64-1 | Acetone | 10 | 10 |
| 75-15-0 | Carbon Disulfide | 5 | 10 |
| 75-35-4 | 1,1-Dichloroethene | 5 | 10 |
| 75-34-3 | 1,1-Dichloroethane | 5 | 10 |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | 10 |
| 67-66-3 | Chloroform | 5 | 10 |
| 107-06-2 | 1,2-Dichloroethane | 5 | 10 |
| 78-93-3 | 2-Butanone | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | 10 |
| 56-23-5 | Carbon Tetrachloride | 5 | 10 |
| 108-05-4 | Vinyl Acetate | 10 | 10 |
| 75-27-4 | Bromodichloromethane | 5 | 10 |
| 78-87-5 | 1,2-Dichloropropane | 5 | 10 |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | 10 |
| 79-01-6 | Trichloroethene | 5 | 10 |
| 124-48-1 | Dibromochloromethane | 5 | 10 |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | 10 |
| 71-43-2 | Benzene | 5 | 10 |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | 10 |
| 75-25-2 | Bromoform | 5 | 10 |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 10 |
| 591-78-6 | 2-Hexanone | 10 | 10 |
| 127-18-4 | Tetrachloroethene | 5 | 10 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | 10 |
| 108-88-3 | Toluene | 5 | 10 |
| 108-90-7 | Chlorobenzene | 5 | 10 |
| 100-41-4 | Ethylbenzene | 5 | 10 |
| 100-42-5 | Styrene | 5 | 10 |
| 1330-20-7 | Xylene (total) | 5 | 10 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBK5

Lab Name: ETC NJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: GC70357

Sample wt/vol: 5 (g/mL) ML Lab File ID: 700979

Level: (low/med) LOW Date Received: 10/03/90

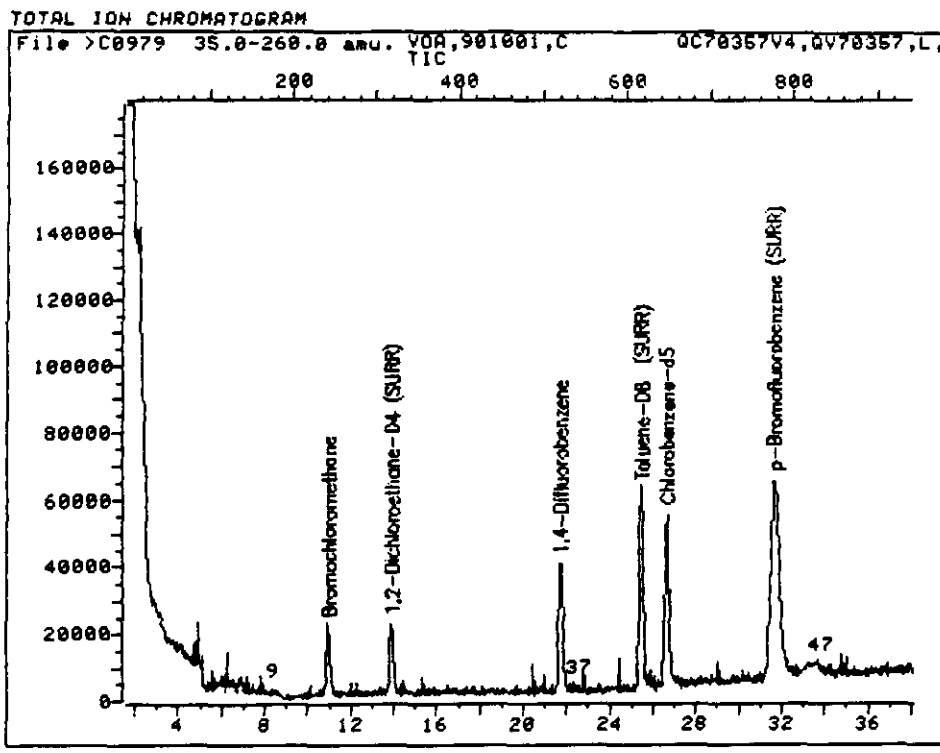
% Moisture: not dec. _____ Date Analyzed: 10/03/90

Column: (pack/cap) PALIS Dilution Factor: 1

CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

Number TICs found: 0

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
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| 30. | | | | |



Data File: >C0979::U4 Quant Output File: ^C0979::AQ
Name: VOA,901001,C
Misc: QC70357V4,QV70357,L,5,,

Id File: IC1172::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901003 17:09

Operator ID: KB6656
Quant Time: 901003 17:10
Injected at: 901003 16:25

QUANT REPORT

Page 1

Operator ID: KB6656
 Jutput File: ^C0979::AQ
 Data File: >C0979::U4
 Name: VOA,901001,C
 Misc: QC70357U4,QU70357,L,5,,

Quant Rev: 7 Quant Time: 901003 17:10
 Injected at: 901003 16:25
 Dilution Factor: 1.00000

ID File: IC1172::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901003 17:09

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) | *Bromochloromethane | 10.99 | 242 | 32787 | 250.00 | NG | 98 |
| 9) | Acetone | 8.31 | 173 | 2343 | 24.50 | NG | 94 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.93 | 318 | 112633 | 242.43 | NG | 90 |
| 20) | *1,4-Difluorobenzene | 21.73 | 519 | 159758 | 250.00 | NG | 95 |
| 36) | *Chlorobenzene-d5 | 26.68 | 646 | 189232 | 250.00 | NG | 75 |
| 37) | Methyl-iso-butyl ketone | 22.47 | 538 | 3765 | 14.68 | NG | 82 |
| 41) | Toluene-D8 (SURR) | 25.49 | 616 | 272990 | 243.94 | NG | 86 |
| 45) | p-Bromofluorobenzene (SURR) | 31.69 | 775 | 239059 | 280.27 | NG | 83 |
| 47) | m-Xylene | 33.71 | 827 | 2049 | 2.74 | NG | 91 |

* Compound is ISTD

Q
 10-16-90.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4755MS

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4755US

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0981

Level: (low/med) LOW

Date Received: ^{24/23/90} ~~10/1990~~ ^{CR} 10/11/90

% Moisture: not dec.

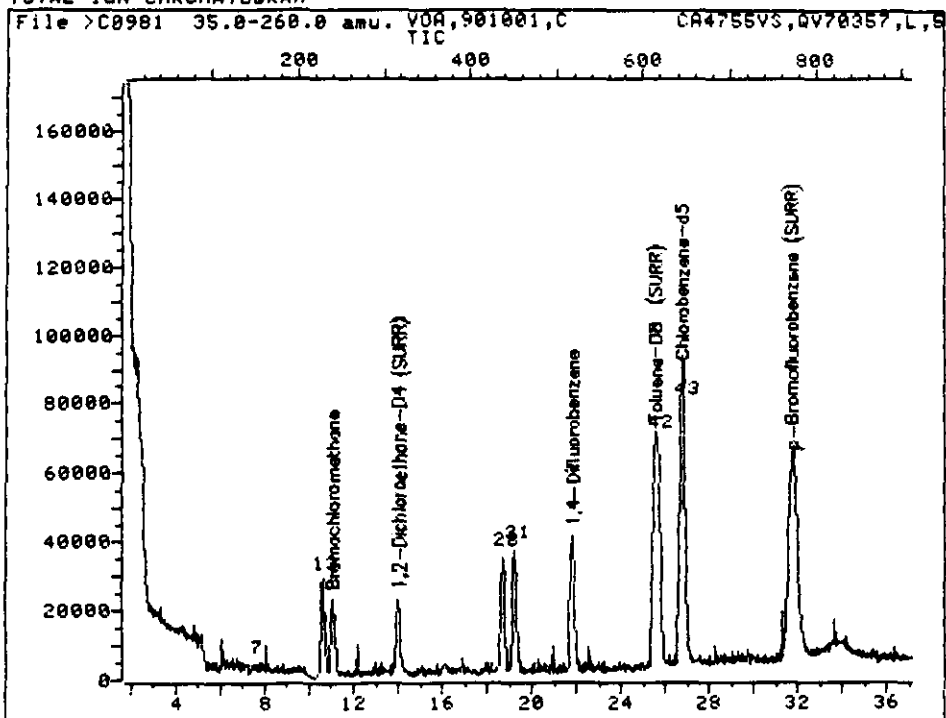
Date Analyzed: 10/03/90

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | | Q |
|------------|----------------------------|----------------------|------|----|
| | | (ug/L or ug/Kg) | UG/L | |
| 74-87-3 | Chloromethane | 10 | 10 | 10 |
| 74-83-9 | Bromomethane | 10 | 10 | 10 |
| 75-01-4 | Vinyl Chloride | 10 | 10 | 10 |
| 75-00-3 | Chloroethane | 10 | 10 | 10 |
| 75-09-2 | Methylene Chloride | 5 | 5 | 5 |
| 67-64-1 | Acetone | 10 | 10 | 10 |
| 75-15-0 | Carbon Disulfide | 5 | 5 | 5 |
| 75-35-4 | 1,1-Dichloroethene | 54 | 54 | 54 |
| 75-34-3 | 1,1-Dichloroethane | 5 | 5 | 5 |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | 5 | 5 |
| 67-66-3 | Chloroform | 5 | 5 | 5 |
| 107-06-2 | 1,2-Dichloroethane | 5 | 5 | 5 |
| 78-93-3 | 2-Butanone | 10 | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | 5 | 5 |
| 56-23-5 | Carbon Tetrachloride | 5 | 5 | 5 |
| 108-05-4 | Vinyl Acetate | 10 | 10 | 10 |
| 75-27-4 | Bromodichloromethane | 5 | 5 | 5 |
| 78-87-5 | 1,2-Dichloropropane | 5 | 5 | 5 |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | 5 | 5 |
| 79-01-6 | Trichloroethene | 48 | 48 | 48 |
| 124-48-1 | Dibromochloromethane | 5 | 5 | 5 |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | 5 | 5 |
| 71-43-2 | Benzene | 42 | 42 | 42 |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | 5 | 5 |
| 75-25-2 | Bromoform | 5 | 5 | 5 |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 10 | 10 |
| 591-78-6 | 2-Hexanone | 10 | 10 | 10 |
| 127-18-4 | Tetrachloroethene | 5 | 5 | 5 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | 5 | 5 |
| 108-88-3 | Toluene | 49 | 49 | 49 |
| 108-90-7 | Chlorobenzene | 48 | 48 | 48 |
| 100-41-4 | Ethylbenzene | 5 | 5 | 5 |
| 100-42-5 | Styrene | 5 | 5 | 5 |
| 1330-20-7 | Xylene (total) | 5 | 5 | 5 |

TOTAL ION CHROMATOGRAM



Data File: >C0981::U4
Name: VOA,901001,C
Misc: CA4755VS,QV70357,L,5,,

Quant Output File: ^C0981::AQ

Id File: IC1172::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901003 17:09

Operator ID: PT1575
Quant Time: 901003 20:28
Injected at: 901003 18:19

QUANT REPORT

Operator ID: PT1575
 Output File: ^C0981::AQ
 Data File: >C0981::U4
 Name: VOA,901001,C
 Misc: CA4755U, QU70357,L,5,,
 S

Quant Rev: 7 Quant Time: 901003 20:28
 Injected at: 901003 18:19
 Dilution Factor: 1.00000

ID File: IC1172::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|-----|
| 1) *Bromochloromethane | 11.02 | 240 | 31333 | 250.00 | NG | 98 |
| 7) Methylene chloride | 7.53 | 150 | 4026 | 24.52 | NG | 76 |
| 13) 1,1-Dichloroethylene | 10.56 | 228 | 57573 | 270.54 | NG | 89 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.01 | 317 | 120617 | 271.66 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.80 | 518 | 158395 | 250.00 | NG | 98 |
| 28) Trichloroethylene | 18.70 | 438 | 56640 | 240.09 | NG | 79 |
| 30) bis(Chloromethyl)ether | 19.21 | 451 | 10705 | 75.34 | NG | 100 |
| 31) Benzene | 19.21 | 451 | 153911 | 207.88 | NG | 87 |
| 36) *Chlorobenzene-d5 | 26.75 | 643 | 191425 | 250.00 | NG | 78 |
| 41) Toluene-D8 (SURR) | 25.53 | 614 | 297943 | 263.19 | NG | 90 |
| 42) Toluene | 25.68 | 618 | 161625 | 247.36 | NG | 93 |
| 43) Chlorobenzene | 26.86 | 646 | 211336 | 239.16 | NG | 95 |
| 45) p-Bromofluorobenzene (SURR) | 31.71 | 771 | 241906 | 280.36 | NG | 81 |

* Compound is ISTD

CA
 10-16-90

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4755MSD

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4755VR

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0982

Level: (low/med) LOW

Date Received: ~~10/1/90~~ ^{10/20/90} ③

% Moisture: not dec.

Date Analyzed: 10/03/90 ^{10/6/90}

Column: (pack/cap) PACK

Dilution Factor: 1

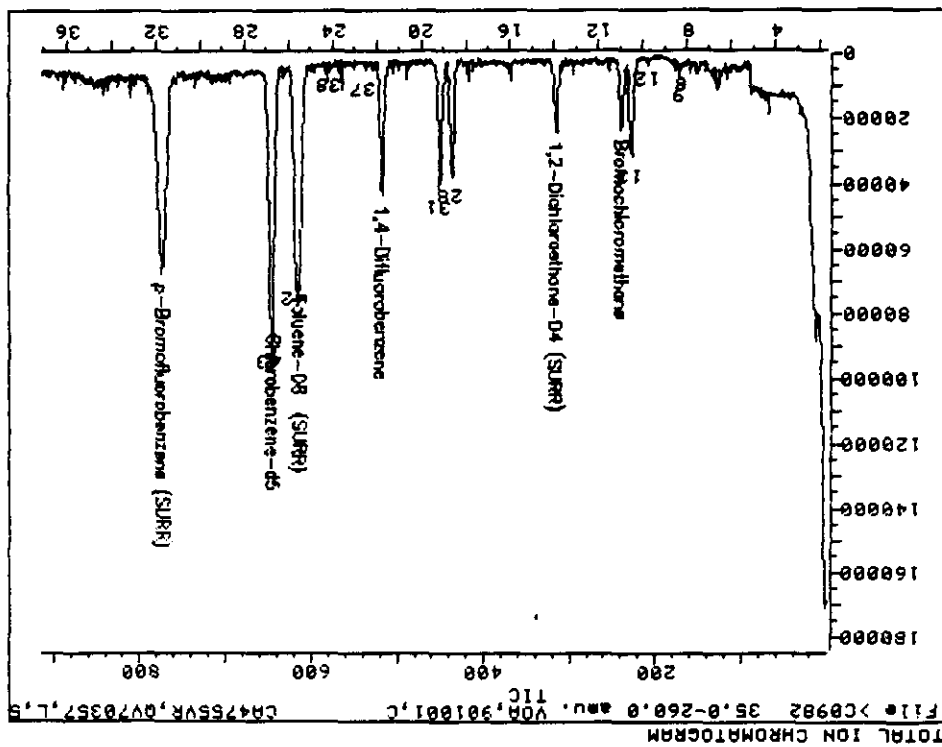
CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

| | | | |
|------------|----------------------------|-----|----|
| 74-87-3 | Chloromethane | 110 | 1U |
| 74-83-9 | Bromomethane | 110 | 1U |
| 75-01-4 | Vinyl Chloride | 110 | 1U |
| 75-00-3 | Chloroethane | 110 | 1U |
| 75-09-2 | Methylene Chloride | 15 | 1U |
| 67-64-1 | Acetone | 110 | 1U |
| 75-15-0 | Carbon Disulfide | 15 | 1U |
| 75-35-4 | 1,1-Dichloroethane | 151 | 1U |
| 75-34-3 | 1,1-Dichloroethane | 15 | 1U |
| 540-59-0 | 1,2-Dichloroethene (total) | 15 | 1U |
| 67-66-3 | Chloroform | 15 | 1U |
| 107-06-2 | 1,2-Dichloroethane | 15 | 1U |
| 78-93-3 | 2-Butanone | 110 | 1U |
| 71-55-6 | 1,1,1-Trichloroethane | 15 | 1U |
| 56-23-5 | Carbon Tetrachloride | 15 | 1U |
| 108-05-4 | Vinyl Acetate | 110 | 1U |
| 75-27-4 | Bromodichloromethane | 15 | 1U |
| 78-87-5 | 1,2-Dichloropropane | 15 | 1U |
| 10061-01-5 | cis-1,3-Dichloropropene | 15 | 1U |
| 79-01-6 | Trichloroethene | 150 | 1U |
| 124-48-1 | Dibromochloromethane | 15 | 1U |
| 79-00-5 | 1,1,2-Trichloroethane | 15 | 1U |
| 71-43-2 | Benzene | 143 | 1U |
| 10061-02-6 | trans-1,3-Dichloropropene | 15 | 1U |
| 75-25-2 | Bromoform | 15 | 1U |
| 108-10-1 | 4-Methyl-2-Pentanone | 110 | 1U |
| 591-78-6 | 2-Hexanone | 110 | 1U |
| 127-18-4 | Tetrachloroethene | 15 | 1U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 15 | 1U |
| 108-88-3 | Toluene | 146 | 1U |
| 108-90-7 | Chlorobenzene | 149 | 1U |
| 100-41-4 | Ethylbenzene | 15 | 1U |
| 100-42-5 | Styrene | 15 | 1U |
| 1330-20-7 | Xylene (total) | 15 | 1U |

Data File: >C0982:U4
 Name: U0A,901001,C
 Misc: C44755UR,0V70357,L,5,,
 ID File: IC1172:US
 Title: IFB, PP/UDA, TCL, XUDAI3
 Last Calibration: 901003 17:09
 Operator ID: PT1575
 Quant Time: 901003 20:38
 Injected at: 901003 19:07

Quant Output File: >C0982:AD



QUANT REPORT

Operator ID: PT1575
 Output File: ^C0982::AQ
 Data File: >C0982::U4
 Name: UOA,901001,C
 Misc: CA4755UR,QU70357,L,5,,

Quant Rev: 7 Quant Time: 901003 20:38
 Injected at: 901003 19:07
 Dilution Factor: 1.00000

ID File: IC1172::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|-------------------|---------------|-----|
| 1) *Bromochloromethane | 10.96 | 240 | 33706 | 250.00 | NG | 93 |
| 8) Acrolein | 8.28 | 171 | 3252 | 141.20 | NG | 77 |
| 9) Acetone | 8.36 | 173 | 3871 | 39.37 | NG | 77 |
| 12) Trichlorofluoromethane | 9.72 | 208 | 3383 | 6.19 | NG | 72 |
| 13) 1,1-Dichloroethylene | 10.49 | 228 | 58885 | 257.23 | NG | 84 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.94 | 317 | 126631 | 265.13 | NG | 98 |
| 20) *1,4-Difluorobenzene | 21.74 | 518 | 166042 | 250.00 | NG | 96 |
| 28) Trichloroethylene | 18.60 | 437 | 61635 | 249.23 | NG | 81 |
| 30) bis(Chloromethyl)ether | 19.14 | 451 | 10810 | 72.57 | NG | 100 |
| 31) Benzene | 19.14 | 451 | 166534 | 214.57 | NG | 86 |
| 36) *Chlorobenzene-d5 | 26.68 | 644 | 193787 | 250.00 | NG | 73 |
| 37) Methyl-iso-butyl ketone | 22.47 | 537 | 3213 | 12.23 | NG | 97 |
| 38) 2-Hexanone | 24.07 | 578 | 4067 | 16.58 | NG | 95 |
| 41) Toluene-D8 (SURR) | 25.46 | 614 | 281173 | 245.35 | NG | 86 |
| 42) Toluene | 25.66 | 619 | 151333 | 228.78 | NG | 93 |
| 43) Chlorobenzene | 26.79 | 647 | 219237 | 245.07 | NG | 91 |
| 45) p-Bromofluorobenzene (SURR) | 31.68 | 773 | 243679 | 278.97 | NG | 81 |

* Compound is ISTD

CA
 10-16-90.

ETC

CHAIN OF CUSTODY

Company: WASTE MANAGEMENT, INC. Attn.: DEE BENCICH

Facility/Site: _____ Phone: (708) 409 - 0700

Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 405 Facility Site Code (Optional Sample Point Descriptions)

Sample Point: WI-26WPZ110 900924 1150
Source Code Your Sample Point ID Start Date Start Time Elapsed Hours
(left justify) (YY/MM/DD) (2400 hr. clock) (composite)

Source Codes
 Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|-------------|----------------|--------------------|
| No | Type | Size | Preserv. | | Filt. (Y/N) | Observations | Observations |
| 1 | TE | 40 | GC/MS | TRIP BLANK | | | ✓ |
| 2 | VOA | 40 | HCL | VOLATILES | N | Sampled for | ✓ |
| 1 | VS | 40 | HCL | VOA SCREEN | N | VOLATILES ONLY | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | | | |
| 1 | CONS | 125 | H2SO4 | TOC2 | | | Bottles Recd Empty |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | | | |
| 1 | MET | 1000 | HNO3 | METALS | | | |
| 1 | CYAN | 500 | NaOH | CY/T | | | |
| 1 | FL | 125 | NONE | FLUORIDE | | | |

CHAIN OF CUSTODY CHRONICLE

1 Shuttle Opened By: (print) DAN GREEN Date: 9-3-90 Time: 14:10
 Signature: [Signature] Seal #: 0175397 Intact: ✓

2 I have received these materials in good condition from the above person.
 Name: JOHN Y. RUDD Signature: [Signature]
 Date: 9/18/90 Time: 10:00am Remarks: OK

3 I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4 Shuttle Sealed By: (print) JOHN Y. RUDD Date: 9/24/90 Time: 12:30
 Signature: _____ Seal #: 0175386 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/25/90 Time: 9:15
 SHUTTLE # 590 TEMP. °C 3 SEAL # 175396 COND. Intact 206

CHAIN OF CUSTODY FORM (CC1) ORIGINAL Date Sealed 90/08/30 By: MM

Company: WASTE MANAGEMENT, INC.

Attn.: DEE BENCICH

Facility/Site: _____

Phone: (708) 409 - 0700

Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 405

Facility/Site Code

(Optional Sample Point Descriptions)

Sample Point: W-26WP75

Source Code (from below)

Your Sample Point ID (left justify)

900924

Start Date (YY/MM/DD)

1340

Start Time (2400 hr clock)

Elapsed Hours (composite)

Source Codes

Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|------------|----------------|---------------------|
| No | Type | Size | Preserv. | | Fill (Y/N) | Observations | Observations |
| 1 | TE | 40 | GC/MS | TRIP BLANK | | | ✓ |
| 2 | VOA | 40 | HCL | VOLATILES | N | VOLATILES ONLY | ✓ |
| 1 | VS | 40 | HCL | VOA SCREEN | N | SAMPLED | ✓ |
| 1 | COND | 125 | NONE | CHLORIDE | | | Bottles Rec'd Empty |
| 1 | CONS | 125 | H2SO4 | TOC2 | | | |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | | | |
| 1 | MET | 1000 | HNO3 | METALS | | | |
| 1 | CYAN | 500 | NaOH | CY/T | | | |
| 1 | FL | 125 | NONE | FLUORIDE | | | |
| | | | | | | | |
| | | | | | | | |

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) DAN GREEN Date: 9-3-90 Time: 14:20
 Signature: [Signature] Seal #: 0175391 Intact: No

2. I have received these materials in good condition from the above person.
 Name: John V. Rudd Signature: [Signature]
 Date: 9/24/90 Time: 09:00 Remarks: OK

3. I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John V. Rudd Date: 9/24/90 Time: 1600
 Signature: [Signature] Seal #: 0175350 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/25/90 Time: 910
 SHUTTLE # 228 TEMP. °C 40 SEAL # 175350 COND. Intact



FIELD PARAMETER FORM (CC2)

Form 0002
Sample Management
12/89

ETC JOB # CA4758

Sample Point

Source Code

Sample Point I.D.

FIELD PROCEDURES

9.0.0.9.2.4

PURGE DATE
(YY MM DD)

1.3.4.0

START PURGE
(2400 Hr Clock)

1.0.0

ELAPSED HRS

1.8.1

WATER VOL. IN CASING
(Gallons)

9.0

VOLUME PURGED
(Gallons)

SAMPLING METHOD: BALLER

Sampler Type

E

A-Submersible Pump
B-ISCO
C-Bladder Pump

D-Dipper/Bottle
E-Bailer
F-Scoop/Shovel

X-Other

(SPECIFY OTHER)

Sampler Material

A

A-Teflon
B-Metal

C-PVC
D-Plastic

X-Other

(SPECIFY OTHER)

Tubing Material

A-Teflon
B-Tygon

C-Polyethylene
D-Silicon

X-Other

(SPECIFY OTHER)

Sample Composited

(Y/N)

N

Procedure/Proportions

FIELD MEASUREMENTS

Well Elevation (ft/msl)

196.6 196 *

Well Depth (ft)

154 46

Depth to Ground water (ft)

4 3 3 9

Sample Depth (non-well) (ft)

Groundwater Elevation (ft msl)

192 3 5 7

1st

7.22

(STD)

ph

1st

610

spec. cond.

um/cm
at 25 °C

(other parameter)

value

units

2nd

7.22

(STD)

ph

2nd

610

spec. cond.

um/cm
at 25 °C

(other parameter)

value

units

3rd

7.23

(STD)

ph

3rd

610

spec. cond.

um/cm
at 25 °C

(other parameter)

value

units

4th

7.22

(STD)

ph

4th

610

spec. cond.

um/cm
at 25 °C

(other parameter)

value

units

12.1

Sample Temp

Turbidity

NTU

FIELD COMMENTS

Sample Appearance: CLEAR / NO ODOOR

Weather Conditions: WINDY - CLEAR - 60s °F - WINDS SW 10-12 MPH

Other: SAMPLED ONLY FOR VOLATILES

FILTERING: Use Chain of Custody (CC1) to indicate which bottles were filtered

Sampler:

John V. Rudd

Employer:

PELA

I certify that sampling procedures were in accordance with applicable EPA state and corporate protocols.

5/28/90

(Date)

John V. Rudd

(Signature)

ORIGINAL

209

CHAIN OF CUSTODY FORM (CC1)

ORIGINAL Date Sealed 90/08/30 By: MM

Company: WASTE MANAGEMENT, INC. Attn.: DEE BENCICH
 Facility/Site: _____ Phone: (708) 409 - 0700
 Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 4105 _____ (Optional Sample Point Descriptions)
 Sample Point: W-26WP29 900925 0845 _____
Source Code (from below) Your Sample Point ID (left justify) Start Date (YY/MM/DD) Start Time (2400 hr. clock) Elapsed Hours (composite)
 Source Codes
 Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|------------|------------------|---------------------|
| No | Type | Size | Preserv. | | FILL (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | | | ✓ small bubble |
| 2 | VOA | 40 | HCL | VOLATILES | N | SAMPLED EOD ✓ | |
| 1 | VS | 40 | HCL | VOA SCREEN | N | VOLATILES ONLY ✓ | |
| 1 | CONU | 125 | NONE | CHLORIDE | | | Bottles Rec'd empty |
| 1 | CONS | 125 | H2SO4 | TOC2 | | | |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | | | |
| 1 | MET | 1000 | HNO3 | METALS | | | |
| 1 | CYAN | 500 | NaOH | CY/T | | | |
| 1 | FL | 125 | NONE | FLUORIDE | | | |

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) JAN GERON Date: 9-3-90 Time: 14:30
 Signature: [Signature] Seal #: 0175369 Intact: ✓

I have received these materials in good condition from the above person.
 2. Name: John V. RUDD Signature: [Signature]
 Date: 9/25/90 Time: 07:00 Remarks: OK

I have received these materials in good condition from the above person.
 3. Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John V. RUDD Date: 9/25/90 Time: 10:00
 Signature: [Signature] Seal #: 0175368 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/26/90 Time: 11:30
 BOTTLE # 993 TEMP. °C 40 SEAL # 175368 COND. [Signature]



FIELD PARAMETER FORM (CC2)

Form 0002
Sample Management
12/89

ETC JOB # CA2759
Sample Point W 2 GWP 29
Source Code Sample Point I.D.

FIELD PROCEDURES

PURGE DATE (YY MM DD) 9 10 19 2 5 START PURGE (2400 Hr Clock) 0 18 4 1 5 ELAPSED HRS 1 5 WATER VOL IN CASING (Gallons) 1 9 9 VOLUME PURGED (Gallons) 9 0

SAMPLING METHOD: BAILER

Sampler Type E A-Submersible Pump D-Dipper/Bottle
B-ISCO E-Bailer X-Other _____ (SPECIFY OTHER)
C-Bladder Pump F-Scoop/Shovel

Sampler Material A A-Teflon C-PVC X-Other _____ (SPECIFY OTHER)
B-Metal D-Plastic

Tubing Material A-Teflon C-Polyethylene X-Other _____ (SPECIFY OTHER)
B-Tygon D-Silicon

Sample Compositing Y/N N Procedure/Proportions

FIELD MEASUREMENTS

Well Elevation (ft/msl) 9 4 4 0 1 * Well Depth (ft) 3 2 5 1
Depth to Ground water (ft) 2 0 3 1 Sample Depth (non-well) (ft)
Groundwater Elevation (ft msl) 9 2 3 7 0

| | | | | | |
|-------------------------------|------------------------------|----------------|-----------|-----------|-----------|
| 1st <u>7 0 2</u> (STD) pH | 1st <u>7 4 8</u> spec. cond. | um/cm at 25 °C | <u> </u> | <u> </u> | <u> </u> |
| 2nd <u>7 0 2</u> (STD) pH | 2nd <u>7 4 8</u> spec. cond. | um/cm at 25 °C | <u> </u> | <u> </u> | <u> </u> |
| 3rd <u>7 0 2</u> (STD) pH | 3rd <u>7 4 6</u> spec. cond. | um/cm at 25 °C | <u> </u> | <u> </u> | <u> </u> |
| 4th <u>7 0 0</u> (STD) pH | 4th <u>7 4 8</u> spec. cond. | um/cm at 25 °C | <u> </u> | <u> </u> | <u> </u> |
| <u>1 2 7</u> (°C) Sample Temp | <u> </u> NTU Turbidity | | | | |

FIELD COMMENTS

Sample Appearance: COLOR - PALE YEL. ORANGE NO ODOOR
Weather Conditions: CLEAR - 60s°F - WIND NW 4-6 mph
Other: * WELL RECORDS
WELL SAMPLED FOR VOLATILES ONLY

FILTERING: Use Chain of Custody (CC1) to indicate which bottles were filtered

Sampler: John Y. Rudd Employer: PELA

I certify that sampling procedures were in accordance with applicable EPA state and corporate protocols.
9/25/90 John Y. Rudd
(Date) (Signature)

ORIGINAL

CHAIN OF CUSTODY FORM (CC1) ORIGINAL Date Sealed 90/08/30 By: MM

Company: WASTE MANAGEMENT, INC.

Attn.: DEE BENCICH

Facility/Site: _____

Phone: (708) 409 - 0700

Address: DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 405

Optional Sample Point Descriptions: _____

Sample Point: W-26WPZ1115

900925

11020

Source Codes

Well (W) Duffall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|-----------|----------------|----------------|
| No | Type | Size | Preserv. | | FIL (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | | | ✓ small bubble |
| 2 | VOA | 40 | HCL | VOLATILES | N | SAMPLED FOR | ✓ |
| 1 | VS | 40 | HCL | VOA SCREEN | N | VOLATILES ONLY | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | | | |
| 1 | CONS | 125 | H2SO4 | TOC2 | | | |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | | | |
| 1 | MET | 1000 | HNO3 | METALS | | | |
| 1 | CYAN | 500 | NaOH | CY/T | | | |
| 1 | FL | 125 | NONE | FLUORIDE | | | |

Bottles Rec'd Empty

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) DAN GREEN Date: 9-3-90 Time: 14:35
Signature: [Signature] Seal #: 0175363 Intact: No

2. I have received these materials in good condition from the above person.
Name: John Y Rudd Signature: [Signature]
Date: 9/18/90 Time: 10:00 am Remarks: OK

3. I have received these materials in good condition from the above person.
Name: _____ Signature: _____
Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John Y. Rudd Date: 9/18/90 Time: 1200
Signature: [Signature] Seal #: 0175362 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/26/90 Time: 1145
BOTTLE # 693 TEMP. °C 30 SEAL # Intact RECOND # 175362



FIELD PARAMETER FORM (CC2)

Form 0002
Sample Management
12/89

ETC JOB # CA4760

Sample Point W 2 GWP 21115
Source Code Sample Point I.D.

FIELD PROCEDURES

91010925
PURGE DATE
(YY MM DD)

110120
START PURGE
(2400 H Clock)

7.5
ELAPSED HRS

11116
WATER VOL. IN CASING
(Gallons)

16010
VOLUME PURGED
(Gallons)

SAMPLING METHOD: BAILER

Sampler Type E A-Submersible Pump D-Dipper/Bottle
B-ISCO E-Bailer X-Other _____
C-Bladder Pump F-Scoop/Shovel (SPECIFY OTHER)

Sampler Material A A-Teflon C-PVC X-Other _____
B-Metal D-Plastic (SPECIFY OTHER)

Tubing Material A-Teflon C-Polyethylene X-Other _____
B-Tygon D-Silicon (SPECIFY OTHER)

Sample Composited Y/N

Procedure/Proportions

FIELD MEASUREMENTS

Well Elevation (ft/msl) 1940.34 * Well Depth (ft) 25.95

Depth to Ground water (ft) 118.85 Sample Depth (non-well) (ft)

Groundwater Elevation (ft msl) 1921.49

1st 6.77 (STD) 1st 1125 um/cm at 25°C
ph spec. cond. (other parameter) value units

2nd 6.77 (STD) 2nd 1125 um/cm at 25°C
ph spec. cond. (other parameter) value units

3rd 6.76 (STD) 3rd 1125 um/cm at 25°C
ph spec. cond. (other parameter) value units

4th 6.77 (STD) 4th 1125 um/cm at 25°C
ph spec. cond. (other parameter) value units

12.5 (°C)
Sample Temp

 NTU
Turbidity

FIELD COMMENTS

Sample Appearance: COLOR - PINK VEL. ORANGE - GARBAGE ODOR

Weather Conditions: CLEAR - 60° F - WIND N.W. 4-6 MPH

Other: * WELL RECORDS

SAMPLED FOR VOLATILES ONLY

FILTERING: Use Chain of Custody (CC1) to indicate which bottles were filtered

Sampler: John V. Rudd
(Print)

Employer: DELA

I certify that sampling procedures were in accordance with applicable EPA state and corporate protocols.

9/25/90 John V. Rudd
(Date) (Signature)

ORIGINAL

Company: WASTE MANAGEMENT, INC.

Attn.: DEE BENCICH

Facility/Site: _____

Phone: (708) 409 - 0700

Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 4 0 5 Facility/Site Code (Optional Sample Point Descriptions)

Sample Point: W-26WP2111I 9010925 113110
Source Code (from below) Your Sample Point ID (left justify) Start Date (YY/MM/DD) Start Time (2400 hr clock) Elapsed Hours (composite)

Source Codes:
 Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|------------|------------------|--------------------|
| No | Type | Size | Preserv. | | Fill (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | | | ✓ small bubble |
| 2 | VOA | 40 | HCL | VOLATILES | N | SAMPLED FOR ✓ | |
| 1 | US | 40 | HCL | VOA SCREEN | N | VOLATILES ONLY ✓ | |
| 1 | CONU | 125 | NONE | CHLORIDE | | | Bottles Recd empty |
| 1 | CONS | 125 | H2SO4 | TOC2 | | | |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | | | |
| 1 | MET | 1000 | HNO3 | METALS | | | |
| 1 | CYAN | 500 | NaOH | CY/T | | | |
| 1 | FL | 125 | NONE | FLUORIDE | | | |

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) Dan Green Date: 9-3-90 Time: 14:40
 Signature: [Signature] Seal #: 0175379 Intact: ✓

2. I have received these materials in good condition from the above person.
 Name: John Y. Rudd Signature: [Signature]
 Date: 9/18/90 Time: 1000 Remarks: OK

3. I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John Y. Rudd Date: 9/25/90 Time: 1500
 Signature: [Signature] Seal #: 0175379 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/26/90 Time: 1140
 BOTTLES 857 SEAL 0175379 COMP. [Signature]

CHAIN OF CUSTODY FORM (CC1) ORIGINAL Date Sealed 9/08/90 By: MM

Company: WASTE MANAGEMENT, INC.

Attn.: DEE BENCICH

Facility/Site: _____

Phone: (708) 409 - 0700

Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 405 Facility/Site Code (Optional Sample Point Descriptions)

Sample Point: W-26WP2111D1 900925 115315
Source Code Your Sample Point ID Start Date Start Time Elapsed Hours
(from below) (left justify) (YY/MM/DD) (2400 hr. clock) (composite)

Source Codes:

Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|------------|----------------|----------------|
| No | Type | Size | Preserv. | | FIL. (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | | | ✓ small bubble |
| 2 | VDA | 40 | HCL | VOLATILES | N | SAMPLED FOR | ✓ |
| 1 | VS | 40 | HCL | VDA SCREEN | N | VOLATILES ONLY | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | | | |
| 1 | CONS | 125 | H2SO4 | TOC2 | | | |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | | | |
| 1 | MET | 1000 | HNO3 | METALS | | | |
| 1 | CYAN | 500 | NaOH | CY/T | | | |
| 1 | FL | 125 | NONE | FLUORIDE | | | |

Bottles Rec'd empty

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) DAN GREEN Date: 9-3-90 Time: 14:50
 Signature: [Signature] Seal #: 0175395 Intact: ✓

2. I have received these materials in good condition from the above person.
 Name: John V. Rudd Signature: [Signature]
 Date: 9/18/90 Time: 10:00am Remarks: OK

3. I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John V. Rudd Date: 9/25/90 Time: 1700
 Signature: [Signature] Seal #: 0175394 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/26/90 Time: 11:35
 TEM: 18.7 TEM °C: 30 SEAL: 175394 COMP: [Signature]

Date 9/26/90 Instr. C

A-Type PP/NOA, TCL / IFB XVOA13
 Tune file APEI01
 Seq. file MI
 Method file VORC
 ID file IC1171
 CB file CC1171
 Analyst JR
 Reviewed by/Date [Signature]

Batch #'s QV 70357

| Standard | Conc ppm | Lot No. | Lot Vol. (ul) |
|------------------|----------|---------|------------------|
| P-BFB | 50 | 4871 | 10 |
| INT STD/SUPP | 25 | 4972 | 10.0 |
| CAL I | 50 | 9916 | 20, 15, 10, 5, 2 |
| CAL II | ↓ | 9997 | ↓ |
| XVOA | ↓ | 9505 | ↓ |
| IFB MATRIX SPILE | 25 | 9673 | 10.0 |

Standards Updated CAL I, CAL II, XVOA

Date 9/26 By [Signature]

Tap # Inj. ul

| RI | NAME | DATA File | 10it. (ml) g | ALS # | Dil | Inj. Time | Comments | P |
|------|------------|-----------|--------------|-------|------------|-----------|-------------------------|---|
| 9/26 | P-BFB | >C0910 | 1ul | - | | 854 | ok 111 | Q |
| | QC 70357VS | >C0911 | 5ml | 6 | | 1124 | ok 20ul CAL I, II, XVOA | Q |
| | QC 70357VS | >C0912 | | 7 | CBUPD | 1213 | ok 15ul IC1171 | Q |
| | QC 70357VS | >C0913 | | 8 | CBACN | 1303 | ok 10ul CC1172 | Q |
| | QC 70357VS | >C0914 | | 9 | CBMUS | 1352 | ok 5ul | Q |
| | QC 70357VS | >C0915 | | 10 | | 1441 | ok 2ul | Q |
| | QC 70357V | >C0916 | ↓ | 1 | - NOT USED | | BLANK | |
| | P-BFB | >C0917 | 1ul | - | | 1625 | (114-104) | Q |
| | QC 70357V | >C0918 | 5ml | 6 | | 1725 | ok | Q |
| | QC 70357V | >C0919 | | 7 | | | NOT USED | |
| | CA 4753V | >C0920 | | 8 | | 1901 | ok | Q |
| | CA 4754V | >C0921 | | 9 | | 1954 | ok re-run, pos leak | Q |
| | CA 4755V | >C0922 | | 10 | | 2043 | ok | Q |
| | CA 4763V | >C0923 | | 1 | | 2152 | ok | Q |
| | CA 4756V | >C0924 | | 2 | | 2221 | ok | Q |
| | CA 4758V | >C0925 | | 3 | | 2310 | ok | Q |
| 9/27 | P-BFB | >C0926 | 1ul | - | | 858 | ok | Q |
| | QC 70357VS | >C0927 | 5ml | 1 | QUR | 1656 | ok 5ul CAL I, II, XVOA | Q |

R : redo , P : Anoclor search , * : Plus search



OHM Corporation

ETC Environmental Testing
and Certification Corp.

Technical Report
for

CITY DISPOSAL CORPORATION LANDFILL

| <i>Chain of Custody Data Required for ETC Data Management Summary Reports</i> | | | | | | |
|---|------------------------|-----------------|---------------------|-------------|-------------|----------------------|
| <i>ETC Sample No.</i> | <i>Company</i> | <i>Facility</i> | <i>Sample Point</i> | <i>Date</i> | <i>Time</i> | <i>Elapsed Hours</i> |
| CA4753-CA4756, CA4763 | WASTE MANAGEMENT, INC. | | | | | |
| | | 405 | | | | |

Richard P. Albert
Vice President, General Manager

This Technical Report is an INSITE™ service generated by LODESTAR™ Data Management Software.

ETC

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| Standard Data | 42 |
| Raw QC Data | 85 |
| Chain of Custody | 121 |

ETC

CASE NARRATIVE

SDG NARRATIVE

This technical report submitted by ETC Corporation contains the analytical results and required deliverables for WASTE MANAGEMENT, INC. Site 405 samples as identified below:

| <u>ETC ID</u> | <u>CLIENT ID</u> | <u>ETC ID</u> | <u>CLIENT ID</u> |
|---------------|------------------|---------------|------------------|
| CA4753 | 2GWBL | CA4756 | 2GWIH |
| CA4754 | 2GWLA | CA4763 | 2FB |
| CA4755 | 2GWSW | | |

During the preparation and analysis of these samples, the following was observed:

VOLATILES:

(OV70357): Problems were not observed during the analysis of these samples.

Release of the data contained in this hardcopy data package has been authorized by the following signature.

Gregory G. Morrison
Gregory G. Morrison
Laboratory Manager

10-17-90
Date

ETC

METHODOLOGY

INTRODUCTION

Environmental Testing and Certification Corp.
284 Raritan Center Parkway, CN 7808
Edison, New Jersey 08818-7808
(201) 225-6700

This report contains analytical results in tabular form for your sample. It includes comprehensive data for each analytical process. Associated quality control data is also presented, including QC batch results for the laboratory blank, spiked blank, matrix spike and a replicated sample spike, as well as results from surrogate compound analyses. Quality control data for instrument performance is also included. Other appendices may include data system printouts and chain of custody records.

Analytical Methodology

ETC utilizes a wide variety of methodologies and EPA approved procedures which are listed below. Analytical results and Quality Assurance protocols are based upon the appropriate guideline dependent upon the level of deliverable requested.

- "Methods of Organic Chemical Analysis of Municipal and Industrial Wastewater", Federal Register Vol. 49, No. 209, October 26, 1984;
- "Test Methods for Evaluating Solid Waste", SW-846 Third Edition, September 1986, USEPA;
- "Standard Methods for the Examination of Water and Wastewater" 1985, 14th, 15th and 16th Edition;
- "Methods for Chemical Analysis of Water and Wastes" March 1983, EMSL, EPA 600 4-79-020;
- Organic Analysis: Multi-media, Multi-Concentration-IFB-CLP, SOW February 1988;
- Inorganic Analysis: Multi-media, Multi-Concentration IFB-CLP, SOW July 1987;
- Dioxin Analysis: Soil/Sediment/Water Matrix; Multi-Concentration, Selected Ion Monitoring with Jar Extraction Procedure; IFB WA86-K357.
- "Methods for the Determination of Organic Compounds in Drinking Water" December 1988, EPA-600/4-88/039;
- "Handbook for Analytical Quality Control in Water and Wastewater Laboratories", EPA-600/4-79-019, March 1979;
- "National Enforcement Investigation Center Policies and Procedures Manual, EPA-330/9/78/001-R, Revised May 1986.

ETC

ETC Network Laboratories are certified to perform analytical analyses upon samples collected throughout the United States. It may be required that this environmental data be reported with reference to a certified laboratory. The Laboratory Identification Numbers for ETC Corp., Edison NJ Laboratory are summarized below for your information and reference. Please contact your Program Manager should you require certification verification for an ETC Network Laboratory.

Certification Summary, January 1990

| <u>State Agency</u> | <u>Certification Number</u> |
|---|------------------------------------|
| Alabama Department of Environmental Management | 40280 |
| Arizona Department of Environmental Quality | 322 |
| California Department of Health Services | 162 |
| Connecticut Department of Health Services | 0511 |
| Florida Department of Health and Rehabilitative Services | E87074 & 87262 |
| Illinois Environmental Protection Agency | 100224 |
| Kansas Department of Health and Environment | E-148 & E1122 |
| New Hampshire Department of Environmental Services | 198948-B |
| New Jersey Department of Environmental Protection | 12257 |
| New York Department of Public Health | 10586 |
| Oklahoma Water Resources Board | 8703 |
| Pennsylvania Department of Environmental Regulation | 68-323 |
| South Carolina Department of Health and Environmental Control | 94002 |
| Tennessee Department of Health and Environmental Laboratory Services | 02915 |
| Utah Department of Health | E-91 |
| Virginia Department of General Services | 00113 |
| Wisconsin Department of Natural Resources | 999464070 |

ETC

VOLATILE DATA

ETC

QC SUMMARY

ETC

END OF PAGE INSERTION

2A
WATER VOLATILE SURROGATE RECOVERY

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

| | EPA SAMPLE NO. | S1 (TOL)# | S2 (BFB)# | S3 (DCE)# | OTHER | TOT (OUT) |
|----|-------------------|--------------|--------------|--------------|-------|--------------|
| 01 | UBLK01 | 101 | 106 | 95 | | 0 |
| 02 | A4753 | 102 | 107 | 99 | | 0 |
| 03 | A4755 | 101 | 107 | 101 | | 0 |
| 04 | A4763 | 100 | 107 | 103 | | 0 |
| 05 | A4756 | 100 | 107 | 102 | | 0 |
| 06 | A4758 | 100 | 110 | 103 | | 0 |
| 07 | UBLK2 | 99 | 103 | 108 | | 0 |
| 08 | A4754 | 103 | 105 | 112 | | 0 |
| 09 | UBLK5 | 98 | 112 | 97 | | 0 |
| 10 | A4755MS | 105 | 112 | 109 | | 0 |
| 11 | A4755MSD | 98 | 112 | 106 | | 0 |
| 12 | | | | | | |
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QC LIMITS

S1 (TOL) = Toluene-d8 (88-110)
 S2 (BFB) = Bromofluorobenzene (86-115)
 S3 (DCE) = 1,2-Dichloroethane-d4 (76-114)

Column to be used to flag recovery values
 * Values outside QC limits
 D- Surrogates diluted out

3A
WATER VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix Spike - EPA Sample No.: A4755

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC # | QC LIMITS REC. |
|--------------------|--------------------------|-----------------------------------|-------------------------------|------------------|----------------------|
| 1,1-Dichloroethene | 50.000 | 0.000 | 54.109 | 108 | 161-145 |
| Trichloroethene | 50.000 | 0.000 | 48.018 | 96 | 171-120 |
| Benzene | 50.000 | 0.000 | 41.577 | 83 | 176-127 |
| Toluene | 50.000 | 0.000 | 49.472 | 99 | 176-125 |
| Chlorobenzene | 50.000 | 0.000 | 47.831 | 96 | 175-130 |

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC # | % RPD # | QC LIMITS RPD REC. |
|--------------------|--------------------------|--------------------------------|-------------------|------------|-------------------------|
| 1,1-Dichloroethene | 50.000 | 51.446 | 103 | 5 | 14 161-145 |
| Trichloroethene | 50.000 | 49.846 | 100 | 4 | 14 171-120 |
| Benzene | 50.000 | 42.915 | 86 | 3 | 11 176-127 |
| Toluene | 50.000 | 45.757 | 92 | 8 | 13 176-125 |
| Chlorobenzene | 50.000 | 49.015 | 98 | 2 | 13 175-130 |

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of limits

RPD: 0 out of 5 outside limits

Spike Recovery: 0 out of 10 outside limits

Comments:

VOLATILE METHOD ^{4A} BLANK SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0918 Lab Sample ID: VBLK01
 Date Analyzed 09/26/90 Time Analyzed: 1725
 Matrix: (soil/water) WATER Level:(low/med) LOW
 Instrument ID: GC/MS C

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | TIME ANALYZED |
|----|-------------------|------------------|----------------|------------------|
| 01 | A4753 | CA4753U | >C0920 | 1904 |
| 02 | A4755 | CA4755U | >C0922 | 2043 |
| 03 | A4763 | CA4763U | >C0923 | 2132 |
| 04 | A4756 | CA4756U | >C0924 | 2221 |
| 05 | A4758 | CA4758U | >C0925 | 2310 |
| 06 | | | | |
| 07 | | | | |
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| 09 | | | | |
| 10 | | | | |
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| 30 | | | | |

Comments: _____

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0979 Lab Sample ID: UBLK5
 Date Analyzed 10/03/90 Time Analyzed: 1625
 Matrix: (soil/water) WATER Level:(low/med) LOW
 Instrument ID: GC/MS C

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | TIME ANALYZED |
|----|-------------------|------------------|----------------|------------------|
| 01 | A4755MS | CA4755US | >C0981 | 1819 |
| 02 | A4755MSD | CA4755UR | >C0982 | 1907 |
| 03 | | | | |
| 04 | | | | |
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| 30 | | | | |

Comments: _____

5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID: >C0917 BFB Injection Date: 09/26/90
 Instrument ID: GC/MS C BFB Injection Time: 1625
 Matrix: (soil/water) WATER Level: (low/med) LOW Column: (pack/cap) PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 23.3 |
| 75 | 30.0 - 60.0% of mass 95 | 59.4 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 6.9 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 84.0 |
| 175 | 5.0 - 9.0% of mass 174 | 7.0 (8.4)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 81.3 (96.7)1 |
| 177 | 5.0 - 9.0% of mass 176 | 6.0 (7.4)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------|---------------|-------------|---------------|---------------|
| 01 | UBLK01 | QC70357U | >C0918 | 09/26/90 | 1725 |
| 02 | A4753 | CA4753U | >C0920 | 09/26/90 | 1904 |
| 03 | A4755 | CA4755U | >C0922 | 09/26/90 | 2043 |
| 04 | A4763 | CA4763U | >C0923 | 09/26/90 | 2132 |
| 05 | A4756 | CA4756U | >C0924 | 09/26/90 | 2221 |
| 06 | A4758 | CA4758U | >C0925 | 09/26/90 | 2310 |
| 07 | | | | | |
| 08 | | | | | |
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5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Lab File ID: >C0926

BFB Injection Date: 09/27/90

Instrument ID: GC/MS C

BFB Injection Time: 0858

Matrix: (soil/water) WATER Level: (low/med) LOW Column: (pack/cap) PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 27.2 |
| 75 | 30.0 - 60.0% of mass 95 | 59.1 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 7.1 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 77.4 |
| 175 | 5.0 - 9.0% of mass 174 | 6.0 (7.7)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 73.7 (95.2)1 |
| 177 | 5.0 - 9.0% of mass 176 | 5.9 (8.0)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------|---------------|-------------|---------------|---------------|
| 01 | USTD50 | QC70357US | >C0927 | 09/27/90 | 1656 |
| 02 | UBLK2 | QC70357U2 | >C0929 | 09/27/90 | 1814 |
| 03 | A4754 | CA4754U2 | >C0931 | 09/27/90 | 1952 |
| 04 | | | | | |
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14

5A
VOLATILE ORGANIC GC/MS TUNING AND MASS
CALIBRATION - BROMOFLUOROBENZENE (BFB)

Lab Name:ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Lab File ID: >C0971

BFB Injection Date:10/03/90

Instrument ID: GC/MS C

BFB Injection Time:0930

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap)PACK

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|--|----------------------|
| 50 | 15.0 - 40.0% of mass 95 | 18.1 |
| 75 | 30.0 - 60.0% of mass 95 | 49.8 |
| 95 | Base peak, 100% relative abundance | 100.0 |
| 96 | 5.0 - 9.0% of mass 95 | 8.0 |
| 173 | Less than 2.0% of mass 174 | 0.0 (0.0)1 |
| 174 | Greater than 50.0% of mass 95 | 94.9 |
| 175 | 5.0 - 9.0% of mass 174 | 7.9 (8.3)1 |
| 176 | Greater than 95.0%, but less than 101.0% of mass 174 | 93.5 (98.5)1 |
| 177 | 5.0 - 9.0% of mass 176 | 7.6 (8.1)2 |

1-Value is % mass 174

2-Value is % mass 176

THIS TUNE APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| | EPA SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------|---------------|-------------|---------------|---------------|
| 01 | USTD50 | QC70357US | >C0972 | 10/03/90 | 1001 |
| 02 | USTD200 | QC70357US | >C0975 | 10/03/90 | 1305 |
| 03 | USTD150 | QC70357US | >C0976 | 10/03/90 | 1355 |
| 04 | USTD100 | QC70357US | >C0977 | 10/03/90 | 1445 |
| 05 | USTD20 | QC70357US | >C0978 | 10/03/90 | 1535 |
| 06 | UCLK5 | QC70357U5 | >C0979 | 10/03/90 | 1625 |
| 07 | A4755MS | CA4755US | >C0981 | 10/03/90 | 1819 |
| 08 | A4755MSD | CA4755UR | >C0982 | 10/03/90 | 1907 |
| 09 | | | | | |
| 10 | | | | | |
| 11 | | | | | |
| 12 | | | | | |
| 13 | | | | | |
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ETC

SAMPLE DATA

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4753

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4753U

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0920

Level: (low/med) LDW

Date Received: ~~10/1/90~~ ^{09/20/90} *CA*

% Moisture: not dec.

Date Analyzed: 09/26/90 ¹⁰⁻¹⁶⁻⁹⁰

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|----|
| 74-87-3 | Chloromethane | 10 | 10 |
| 74-83-9 | Bromomethane | 10 | 10 |
| 75-01-4 | Vinyl Chloride | 10 | 10 |
| 75-00-3 | Chloroethane | 10 | 10 |
| 75-09-2 | Methylene Chloride | 15 | 10 |
| 67-64-1 | Acetone | 10 | 10 |
| 75-15-0 | Carbon Disulfide | 15 | 10 |
| 75-35-4 | 1,1-Dichloroethene | 15 | 10 |
| 75-34-3 | 1,1-Dichloroethane | 15 | 10 |
| 540-59-0 | 1,2-Dichloroethene (total) | 15 | 10 |
| 67-66-3 | Chloroform | 15 | 10 |
| 107-06-2 | 1,2-Dichloroethane | 15 | 10 |
| 78-93-3 | 2-Butanone | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | 15 | 10 |
| 56-23-5 | Carbon Tetrachloride | 15 | 10 |
| 108-05-4 | Vinyl Acetate | 10 | 10 |
| 75-27-4 | Bromodichloromethane | 15 | 10 |
| 78-87-5 | 1,2-Dichloropropane | 15 | 10 |
| 10061-01-5 | cis-1,3-Dichloropropene | 15 | 10 |
| 79-01-6 | Trichloroethene | 15 | 10 |
| 124-48-1 | Dibromochloromethane | 15 | 10 |
| 79-00-5 | 1,1,2-Trichloroethane | 15 | 10 |
| 71-43-2 | Benzene | 15 | 10 |
| 10061-02-6 | trans-1,3-Dichloropropene | 15 | 10 |
| 75-25-2 | Bromoform | 15 | 10 |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 10 |
| 591-78-6 | 2-Hexanone | 10 | 10 |
| 127-18-4 | Tetrachloroethene | 15 | 10 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 15 | 10 |
| 108-88-3 | Toluene | 15 | 10 |
| 108-90-7 | Chlorobenzene | 15 | 10 |
| 100-41-4 | Ethylbenzene | 15 | 10 |
| 100-42-5 | Styrene | 15 | 10 |
| 1330-20-7 | Xylene (total) | 15 | 10 |

17

12
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETCNS Contract: _____

A4753

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: CA4753

Sample wt/vol: 5 (g/mL) HL Lab File ID: 7C0920

Level: (low/med) LOW Date Received: 09/26/90

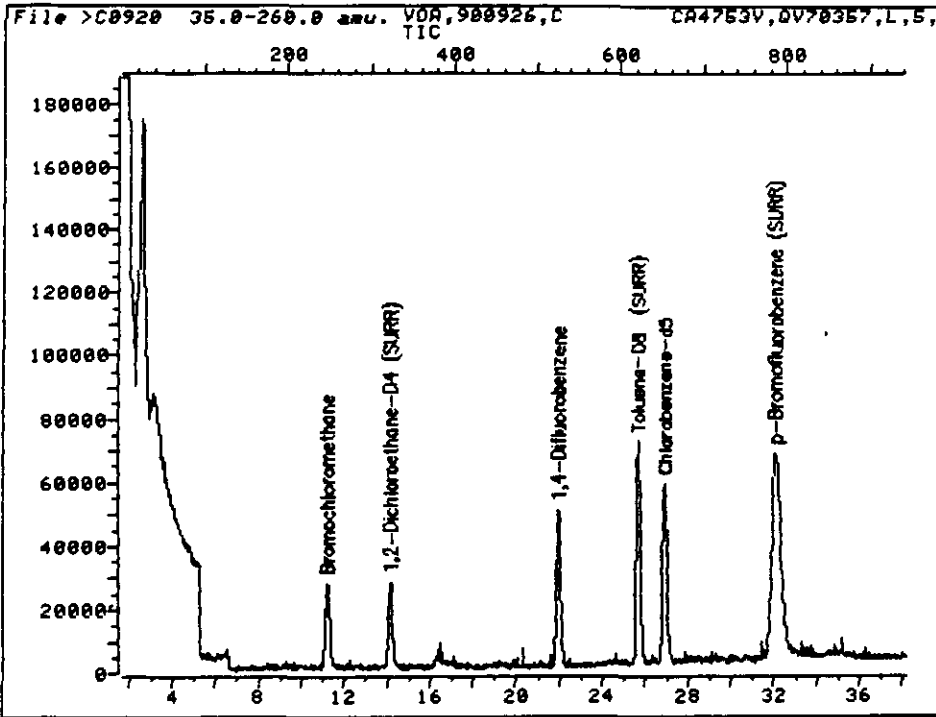
% Moisture: not dec. _____ Date Analyzed: 09/26/90

Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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| 30. | | | | |

TOTAL ION CHROMATOGRAM



Data File: >C0920::U0
Name: VOA,900926;C
Misc: CA4753V,QU70357,L,5,,

Quant Output File: ^C0920::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900926 16:16

Operator ID: JA8781
Quant Time: 900926 19:44
Injected at: 900926 19:04

QUANT REPORT

Page 1

Operator ID: JA8781
 Output File: ^C0920::AQ
 Data File: >C0920::U0
 Name: UOA,900926,C
 Misc: CA4753U,QV70357,L,5,,

Quant Rev: 7 Quant Time: 900926 19:44
 Injected at: 900926 19:04
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUQA13
 Last Calibration: 900926 16:16

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|--------|-------|----|
| 1) *Bromochloromethane | 11.19 | 246 | 46998 | 250.00 | NG | 95 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.14 | 322 | 163095 | 248.59 | NG | 96 |
| 20) *1,4-Difluorobenzene | 21.94 | 523 | 224774 | 250.00 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.92 | 651 | 207367 | 250.00 | NG | 75 |
| 41) Toluene-D8 (SURR) | 25.66 | 619 | 312662 | 256.01 | NG | 89 |
| 45) p-Bromofluorobenzene (SURR) | 32.08 | 784 | 242017 | 268.18 | NG | 81 |

* Compound is ISTD

EP
10-17-10.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4754

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4754U2

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0931

Level: (low/med) LOW

Date Received: ~~10/1/90~~ ^{09/20/90} CP

% Moisture: not dec.

Date Analyzed: 09/27/90 ¹⁰⁻¹⁶⁻⁹⁰

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) | UG/L | Q |
|------------|----------------------------|---|------|---|
| 74-87-3 | Chloromethane | | 10 | U |
| 74-83-9 | Bromomethane | | 10 | U |
| 75-01-4 | Vinyl Chloride | | 10 | U |
| 75-00-3 | Chloroethane | | 10 | U |
| 75-09-2 | Methylene Chloride | | 5 | U |
| 67-64-1 | Acetone | | 10 | U |
| 75-15-0 | Carbon Disulfide | | 5 | U |
| 75-35-4 | 1,1-Dichloroethane | | 5 | U |
| 75-34-3 | 1,1-Dichloroethane | | 5 | U |
| 540-59-0 | 1,2-Dichloroethane (total) | | 5 | U |
| 67-66-3 | Chloroform | | 5 | U |
| 107-06-2 | 1,2-Dichloroethane | | 5 | U |
| 78-93-3 | 2-Butanone | | 10 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 5 | U |
| 56-23-5 | Carbon Tetrachloride | | 5 | U |
| 108-05-4 | Vinyl Acetate | | 10 | U |
| 75-27-4 | Bromodichloromethane | | 5 | U |
| 78-87-5 | 1,2-Dichloropropane | | 5 | U |
| 10061-01-5 | cis-1,3-Dichloropropene | | 5 | U |
| 79-01-6 | Trichloroethene | | 5 | U |
| 124-48-1 | Dibromochloromethane | | 5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 5 | U |
| 71-43-2 | Benzene | | 5 | U |
| 10061-02-6 | trans-1,3-Dichloropropene | | 5 | U |
| 75-25-2 | Bromoform | | 5 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | | 10 | U |
| 591-78-6 | 2-Hexanone | | 10 | U |
| 127-18-4 | Tetrachloroethene | | 5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5 | U |
| 108-88-3 | Toluene | | 5 | U |
| 108-90-7 | Chlorobenzene | | 5 | U |
| 100-41-4 | Ethylbenzene | | 5 | U |
| 100-42-5 | Styrene | | 5 | U |
| 1330-20-7 | Xylene (total) | | 5 | U |

22

12
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

A4754

Lab Name: ETC NJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: C44754

Sample wt/vol: 5 (g/mL) HL Lab File ID: 200931

Level: (low/med) LOW Date Received: 09/01/90

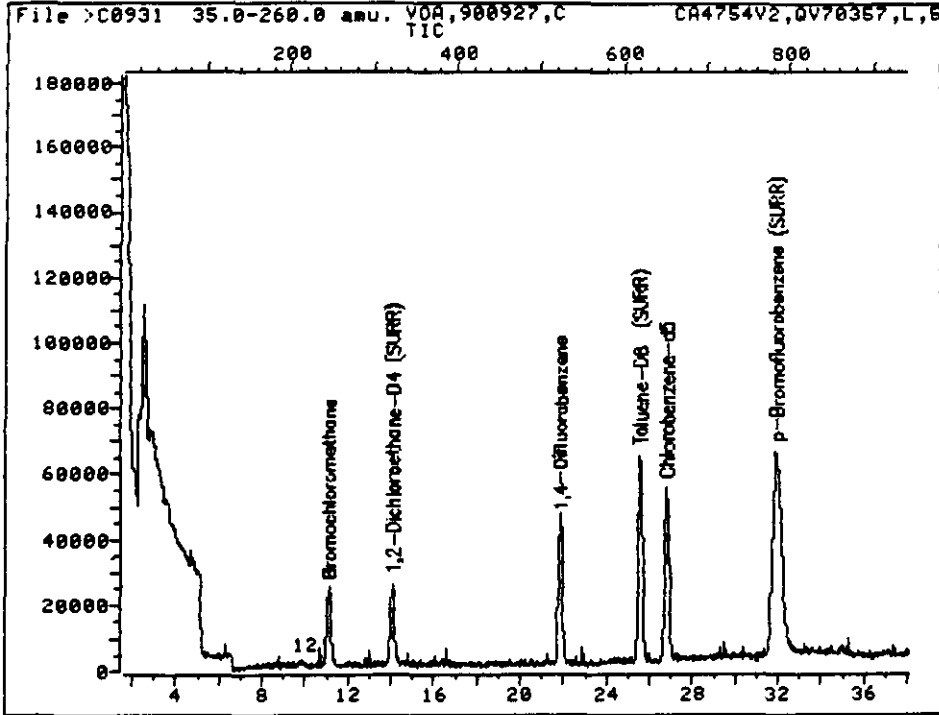
% Moisture: not dec. _____ Date Analyzed: 09/17/90

Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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| 30. | | | | |

TOTAL ION CHROMATOGRAM



Data File: >C0931::U4
Name: VOA,900927,C
Misc: CA4754V2,QV70357,L,5,,

Quant Output File: ^C0931::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900927 17:41

Operator ID: MGRMS
Quant Time: 900927 20:31
Injected at: 900927 19:52

QUANT REPORT

Page 1

Operator ID: MGRMS
 Output File: ^C0931::AQ
 Data File: >C0931::U4
 Name: UOA,900927,C
 Misc: CA4754U2,QU70357,L,5,,

Quant Rev: 7 Quant Time: 900927 20:31
 Injected at: 900927 19:52
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900927 17:41

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.10 | 245 | 42061 | 250.00 | NG | 98 |
| 12) Trichlorofluoromethane | 9.90 | 214 | 9391 | 12.47 | NG | 85 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.09 | 322 | 152876 | 281.20 | NG | 95 |
| 20) *1,4-Difluorobenzene | 21.85 | 522 | 196522 | 250.00 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.82 | 650 | 185259 | 250.00 | NG | 76 |
| 41) Toluene-D8 (SURR) | 25.57 | 618 | 278543 | 257.44 | NG | 90 |
| 45) p-Bromofluorobenzene (SURR) | 31.94 | 782 | 220655 | 263.69 | NG | 87 |

* Compound is ISTD

(a)
 10-17-90.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4755

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4755U

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0922

Level: (low/med) LOW

Date Received: ~~10/19/90~~ ^{04/10/90} *EA*

% Moisture: not dec.

Date Analyzed: 09/26/90 ¹⁰⁻¹⁶⁻⁹⁰

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | | Q |
|------------|----------------------------|----------------------|------|----|
| | | (ug/L or ug/Kg) | UG/L | |
| 74-87-3 | Chloromethane | 10 | 10 | 10 |
| 74-83-9 | Bromomethane | 10 | 10 | 10 |
| 75-01-4 | Vinyl Chloride | 10 | 10 | 10 |
| 75-00-3 | Chloroethane | 10 | 10 | 10 |
| 75-09-2 | Methylene Chloride | 5 | 5 | 10 |
| 67-64-1 | Acetone | 10 | 10 | 10 |
| 75-15-0 | Carbon Disulfide | 5 | 5 | 10 |
| 75-35-4 | 1,1-Dichloroethene | 5 | 5 | 10 |
| 75-34-3 | 1,1-Dichloroethane | 5 | 5 | 10 |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | 5 | 10 |
| 67-66-3 | Chloroform | 5 | 5 | 10 |
| 107-06-2 | 1,2-Dichloroethane | 5 | 5 | 10 |
| 78-93-3 | 2-Butanone | 10 | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | 5 | 10 |
| 56-23-5 | Carbon Tetrachloride | 5 | 5 | 10 |
| 108-05-4 | Vinyl Acetate | 10 | 10 | 10 |
| 75-27-4 | Bromodichloromethane | 5 | 5 | 10 |
| 78-87-5 | 1,2-Dichloropropane | 5 | 5 | 10 |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | 5 | 10 |
| 79-01-6 | Trichloroethene | 5 | 5 | 10 |
| 124-48-1 | Dibromochloromethane | 5 | 5 | 10 |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | 5 | 10 |
| 71-43-2 | Benzene | 5 | 5 | 10 |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | 5 | 10 |
| 75-25-2 | Bromoform | 5 | 5 | 10 |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 10 | 10 |
| 591-78-6 | 2-Hexanone | 10 | 10 | 10 |
| 127-18-4 | Tetrachloroethene | 5 | 5 | 10 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | 5 | 10 |
| 108-88-3 | Toluene | 5 | 5 | 10 |
| 108-90-7 | Chlorobenzene | 5 | 5 | 10 |
| 100-41-4 | Ethylbenzene | 5 | 5 | 10 |
| 100-42-5 | Styrene | 5 | 5 | 10 |
| 1330-20-7 | Xylene (total) | 5 | 5 | 10 |

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12
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

A4765

Lab Name: ETC NJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: CA4765

Sample wt/vol: 5 (g/mL) ML Lab File ID: 7CC942

Level: (low/med) LOW Date Received: 09/20/90

% Moisture: not dec. _____ Date Analyzed: 12/12/90

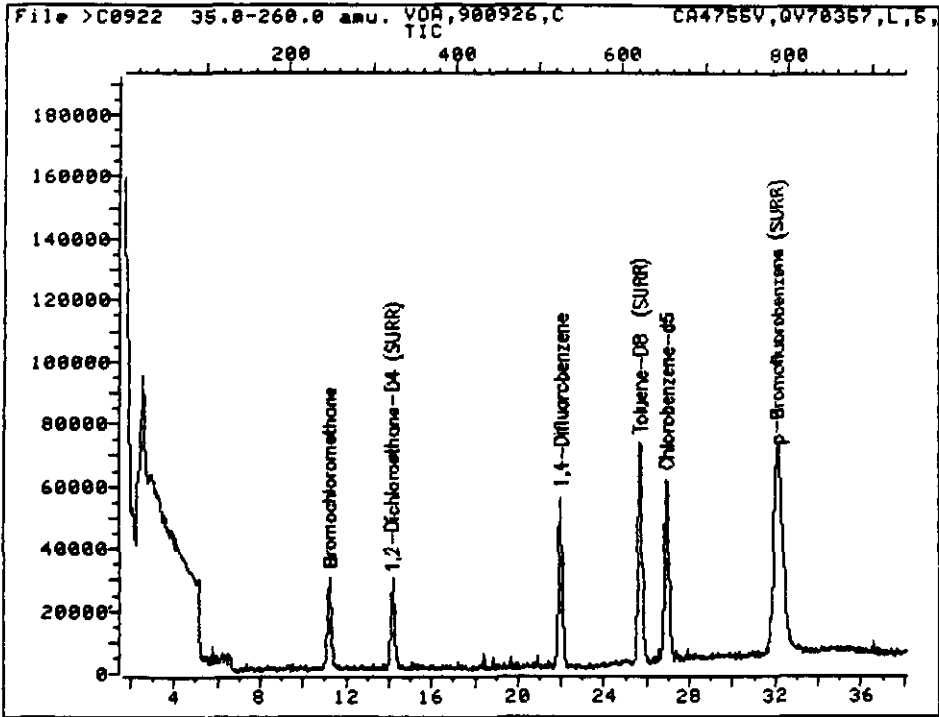
Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 0

CONCENTRATION UNITS:
 (ug/L or ug/kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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| 30. | | | | |

TOTAL ION CHROMATOGRAM



Data File: >C0922::U0
Name: UOA,900926,C
Misc: CA4755V,QV70357,L,5,,

Quant Output File: ^C0922::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900926 16:16

Operator ID: JA8781
Quant Time: 900926 21:22
Injected at: 900926 20:43

QUANT REPORT

Page 1

Operator ID: JAB781
 Output File: ^C0922::AQ
 Data File: >C0922::U0
 Name: VOA,900926,C
 Misc: CA4755U,QU70357,L,5,,

Quant Rev: 7 Quant Time: 900926 21:22
 Injected at: 900926 20:43
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 16:16

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|--------|-------|----|
| 1) | *Bromochloromethane | 11.20 | 247 | 51014 | 250.00 | NG | 95 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 14.15 | 323 | 180352 | 253.25 | NG | 99 |
| 20) | *1,4-Difluorobenzene | 21.98 | 525 | 234507 | 250.00 | NG | 96 |
| 36) | *Chlorobenzene-d5 | 26.92 | 652 | 216956 | 250.00 | NG | 75 |
| 41) | Toluene-D8 (SURR) | 25.67 | 620 | 321910 | 251.93 | NG | 88 |
| 45) | p-Bromofluorobenzene (SURR) | 32.08 | 785 | 251598 | 266.47 | NG | 81 |

* Compound is ISTD

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CA4756

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4756U

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0924

Level: (low/med) LOW

Date Received: ~~10/1/90~~ ^{09/26/90} *CP*

% Moisture: not dec.

Date Analyzed: 09/26/90 ¹⁰⁻¹⁶⁻⁹⁰

Column: (pack/cap) PACK

Dilution Factor: 1

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

| | | | |
|------------|----------------------------|----|----|
| 74-87-3 | Chloromethane | 10 | IU |
| 74-83-9 | Bromomethane | 10 | IU |
| 75-01-4 | Vinyl Chloride | 10 | IU |
| 75-00-3 | Chloroethane | 10 | IU |
| 75-09-2 | Methylene Chloride | 15 | IU |
| 67-64-1 | Acetone | 10 | IU |
| 75-15-0 | Carbon Disulfide | 15 | IU |
| 75-35-4 | 1,1-Dichloroethane | 15 | IU |
| 75-34-3 | 1,1-Dichloroethane | 15 | IU |
| 540-59-0 | 1,2-Dichloroethane (total) | 15 | IU |
| 67-66-3 | Chloroform | 15 | IU |
| 107-06-2 | 1,2-Dichloroethane | 15 | IU |
| 78-93-3 | 2-Butanone | 10 | IU |
| 71-55-6 | 1,1,1-Trichloroethane | 15 | IU |
| 56-23-5 | Carbon Tetrachloride | 15 | IU |
| 108-05-4 | Vinyl Acetate | 10 | IU |
| 75-27-4 | Bromodichloromethane | 15 | IU |
| 78-87-5 | 1,2-Dichloropropane | 15 | IU |
| 10061-01-5 | cis-1,3-Dichloropropene | 15 | IU |
| 79-01-6 | Trichloroethene | 15 | IU |
| 124-48-1 | Dibromochloromethane | 15 | IU |
| 79-00-5 | 1,1,2-Trichloroethane | 15 | IU |
| 71-43-2 | Benzene | 15 | IU |
| 10061-02-6 | trans-1,3-Dichloropropene | 15 | IU |
| 75-25-2 | Bromoform | 15 | IU |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | IU |
| 591-78-6 | 2-Hexanone | 10 | IU |
| 127-18-4 | Tetrachloroethene | 15 | IU |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 15 | IU |
| 108-88-3 | Toluene | 15 | IU |
| 108-90-7 | Chlorobenzene | 15 | IU |
| 100-41-4 | Ethylbenzene | 15 | IU |
| 100-42-5 | Styrene | 15 | IU |
| 1330-20-7 | Xylene (total) | 15 | IU |

32

18
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

A4756

Lab Name: ETCNS Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: CA4756

Sample wt/vol: 5 (g/mL) HL Lab File ID: >C0924

Level: (low/med) LOW Date Received: 09/21/90

% Moisture: not dec. _____ Date Analyzed: 09/26/90

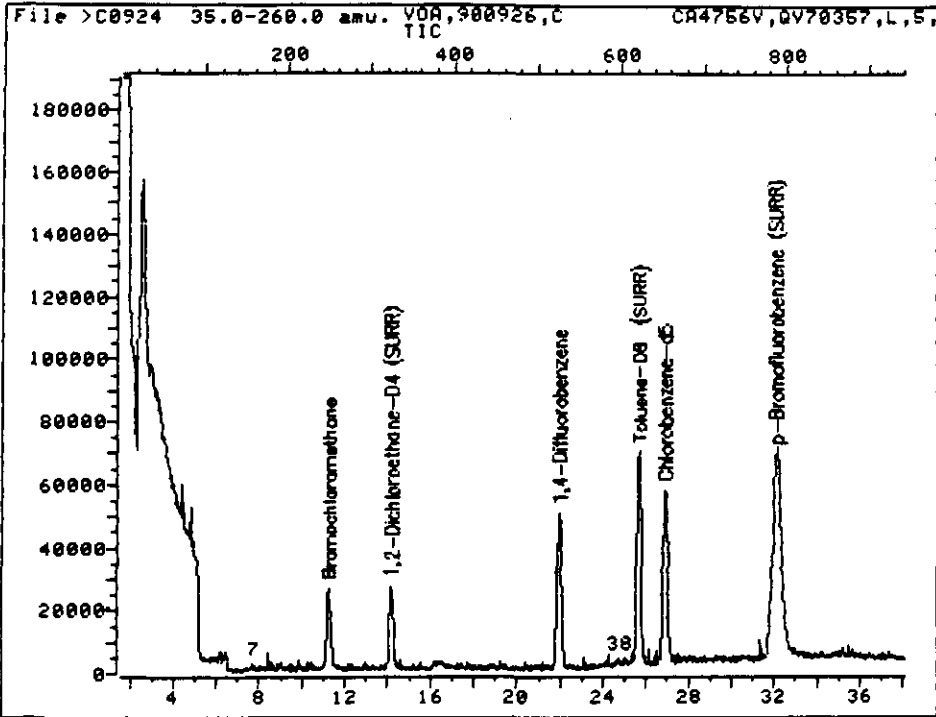
Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 2

CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
| 3. | | | | |
| 4. | | | | |
| 5. | | | | |
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| 26. | | | | |
| 27. | | | | |
| 28. | | | | |
| 29. | | | | |
| 30. | | | | |

TOTAL ION CHROMATOGRAM



Data File: >C0924::U0
Name: VOA,900926,C
Misc: CA4756V,QU70357,L,5,,

Quant Output File: ^C0924::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900926 16:16

Operator ID: JA8781
Quant Time: 900926 23:00
Injected at: 900926 22:21

QUANT REPORT

Page 1

Operator ID: JA8781
 Output File: ^C0924::AQ
 Data File: >C0924::U0
 Name: UOA,900926,C
 Misc: CA4756U,QU70357,L,5,,

Quant Rev: 7 Quant Time: 900926 23:00
 Injected at: 900926 22:21
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900926 16:16

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.26 | 250 | 44376 | 250.00 | NG | 95 |
| 7) Methylene chloride | 7.65 | 157 | 3431 | 11.60 | NG | 90 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.17 | 325 | 157436 | 254.14 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.96 | 526 | 215254 | 250.00 | NG | 96 |
| 36) *Chlorobenzene-d5 | 26.90 | 653 | 204163 | 250.00 | NG | 76 |
| 38) 2-Hexanone | 24.68 | 596 | 2711 | 13.13 | NG | 53 |
| 41) Toluene-D8 (SURR) | 25.69 | 622 | 300152 | 249.62 | NG | 89 |
| 45) p-Bromofluorobenzene (SURR) | 32.10 | 787 | 237983 | 267.85 | NG | 83 |

* Compound is ISTD

CA
 1017-90.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

1A4763

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4763U

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0923

Level: (low/med) LOW

Date Received: ~~10/19/90~~ ^{04/28/90} (CP) 10/19/90

% Moisture: not dec.

Date Analyzed: 09/26/90

Column: (pack/cap) PACK

Dilution Factor: 1

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|----------------------------|-----------------|------|----|
| 74-87-3 | Chloromethane | | 10 | 1U |
| 74-83-9 | Bromomethane | | 10 | 1U |
| 75-01-4 | Vinyl Chloride | | 10 | 1U |
| 75-00-3 | Chloroethane | | 10 | 1U |
| 75-09-2 | Methylene Chloride | | 15 | 1U |
| 67-64-1 | Acetone | | 10 | 1U |
| 75-15-0 | Carbon Disulfide | | 15 | 1U |
| 75-35-4 | 1,1-Dichloroethene | | 15 | 1U |
| 75-34-3 | 1,1-Dichloroethane | | 15 | 1U |
| 540-59-0 | 1,2-Dichloroethene (total) | | 15 | 1U |
| 67-66-3 | Chloroform | | 15 | 1U |
| 107-06-2 | 1,2-Dichloroethane | | 15 | 1U |
| 78-93-3 | 2-Butanone | | 10 | 1U |
| 71-55-6 | 1,1,1-Trichloroethane | | 15 | 1U |
| 56-23-5 | Carbon Tetrachloride | | 15 | 1U |
| 108-05-4 | Vinyl Acetate | | 10 | 1U |
| 75-27-4 | Bromodichloromethane | | 15 | 1U |
| 78-87-5 | 1,2-Dichloropropane | | 15 | 1U |
| 10061-01-5 | cis-1,3-Dichloropropene | | 15 | 1U |
| 79-01-6 | Trichloroethene | | 15 | 1U |
| 124-48-1 | Dibromochloromethane | | 15 | 1U |
| 79-00-5 | 1,1,2-Trichloroethane | | 15 | 1U |
| 71-43-2 | Benzene | | 15 | 1U |
| 10061-02-6 | trans-1,3-Dichloropropene | | 15 | 1U |
| 75-25-2 | Bromoform | | 15 | 1U |
| 108-10-1 | 4-Methyl-2-Pentanone | | 10 | 1U |
| 591-78-6 | 2-Hexanone | | 10 | 1U |
| 127-18-4 | Tetrachloroethene | | 15 | 1U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 15 | 1U |
| 108-88-3 | Toluene | | 15 | 1U |
| 108-90-7 | Chlorobenzene | | 15 | 1U |
| 100-41-4 | Ethylbenzene | | 15 | 1U |
| 100-42-5 | Styrene | | 15 | 1U |
| 1330-20-7 | Xylene (total) | | 15 | 1U |

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1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETC NJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: CA4763

Sample wt/vol: 5 (g/mL) HL Lab File ID: >C0903

Level: (low/med) LOW Date Received: 09/20/90

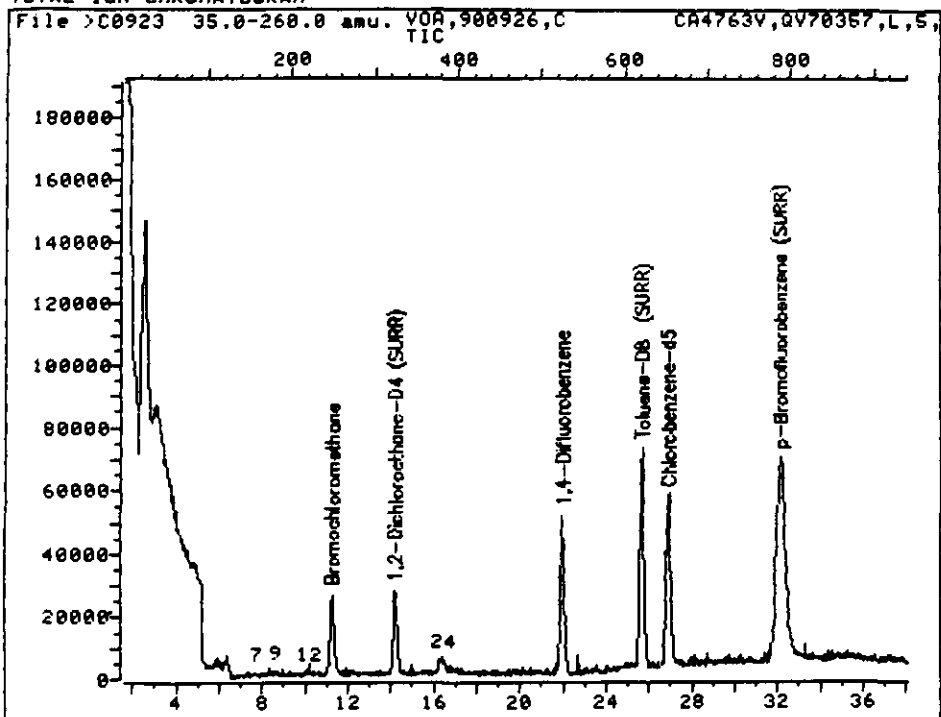
% Moisture: not dec. _____ Date Analyzed: 09/26/90

Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
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TOTAL ION CHROMATOGRAM



Data File: >C0923::U0
Name: VOA,900926,C
Misc: CA4763V,QV70357,L,5,,

Quant Output File: ^C0923::AQ

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900926 16:16

Operator ID: JA8781
Quant Time: 900926 22:11
Injected at: 900926 21:32

QUANT REPORT

Operator ID: JA8781
 Output File: ^C0923::AQ
 Data File: >C0923::U0
 Name: UOA,900926,C
 Misc: CA4763U,QU70357,L,5,,

Quant Rev: 7 Quant Time: 900926 22:11
 Injected at: 900926 21:32
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 16:16

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.24 | 249 | 45139 | 250.00 | NG | 96 |
| 7) Methylene chloride | 7.64 | 156 | 2379 | 7.96 | NG | 84 |
| 9) Acetone | 8.53 | 179 | 5165 | 47.22 | NG | 74 |
| 12) Trichlorofluoromethane | 10.00 | 217 | 4109 | 5.32 | NG | 94 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.19 | 325 | 162300 | 257.57 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.95 | 525 | 219122 | 250.00 | NG | 99 |
| 24) Vinyl acetate | 16.29 | 379 | 3767 | 9.98 | NG | 73 |
| 36) *Chlorobenzene-d5 | 26.88 | 652 | 207394 | 250.00 | NG | 77 |
| 41) Toluene-DB (SURR) | 25.67 | 621 | 306516 | 250.94 | NG | 86 |
| 45) p-Bromofluorobenzene (SURR) | 32.12 | 787 | 241814 | 267.92 | NG | 84 |

* Compound is ISTD

EA
 10-17-90.

ETC

STANDARDS DATA

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:ETCNJ

Contract:

Lab Code:F2

Case No.:

SAS No.:

SDG No.:

Instrument ID:GC/MS C

Calibration Date(s) 09/26/90

09/26/90

Matrix:(soil/water) WATER

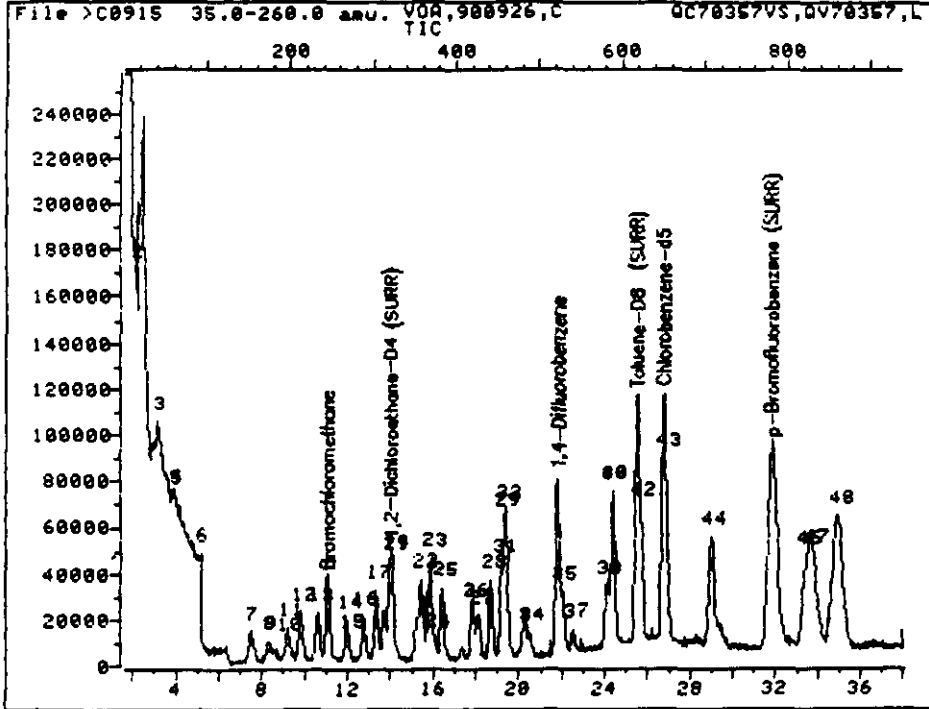
Level:(low/med) LOW

Column:(pack/cap) PACK

Min RRF for SPCC(*) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

| LAB FILE ID: | RRF20 =>C0915 | RRF50 =>C0914 | | | | | |
|------------------------------|---------------|---------------|--------|--------|--------|-------|-------|
| RRF100=>C0913 | RRF150=>C0912 | RRF200=>C0911 | | | | | |
| COMPOUND | RRF20 | RRF50 | RRF100 | RRF150 | RRF200 | RRF | % RSD |
| Chloromethane_____# | .746 | .648 | .619 | .587 | .587 | .637 | 10.3# |
| Bromomethane_____ | .474 | .298 | .216 | .308 | .148 | .289 | 42.4 |
| Vinyl Chloride_____* | 1.598 | 1.187 | 1.061 | 1.282 | 1.057 | 1.237 | 18.0* |
| Chloroethane_____ | 1.014 | .827 | .797 | .869 | .967 | .895 | 10.3 |
| Methylene Chloride____ | 1.913 | 1.631 | 1.621 | 1.579 | 1.529 | 1.655 | 9.1 |
| Acetone_____ | .606 | .631 | .561 | .612 | .619 | .606 | 4.5 |
| Carbon Disulfide_____ | 5.121 | 4.577 | 4.609 | 4.680 | 4.514 | 4.700 | 5.2 |
| 1,1-Dichloroethane____* | 1.669 | 1.462 | 1.440 | 1.456 | 1.375 | 1.480 | 7.5# |
| 1,1-Dichloroethane____# | 3.113 | 2.697 | 2.675 | 2.710 | 2.570 | 2.753 | 7.6# |
| 1,2-Dichloroethane (total)_ | 1.677 | 1.434 | 1.409 | 1.491 | 1.356 | 1.473 | 8.4 |
| Chloroform_____* | 5.511 | 4.748 | 4.641 | 4.702 | 4.508 | 4.822 | 8.2* |
| 1,2-Dichloroethane____ | 4.718 | 4.074 | 3.899 | 3.949 | 3.856 | 4.099 | 8.7 |
| 2-Butanone_____ | .027 | .024 | .023 | .023 | .021 | .024 | 8.0 |
| 1,1,1-Trichloroethane____ | 1.082 | .953 | .915 | .932 | .979 | .972 | 6.8 |
| Carbon Tetrachloride____ | 1.070 | .936 | .908 | .933 | .993 | .968 | 6.7 |
| Vinyl Acetate_____ | .505 | .461 | .405 | .461 | .463 | .459 | 7.8 |
| Bromodichloromethane____ | .901 | .801 | .773 | .795 | .806 | .815 | 6.1 |
| 1,2-Dichloropropene____* | .283 | .253 | .248 | .249 | .238 | .254 | 6.7* |
| cis-1,3-Dichloropropene___ | .526 | .479 | .462 | .464 | .462 | .478 | 5.7 |
| Trichloroethene_____ | .482 | .420 | .454 | .413 | .414 | .437 | 7.0 |
| Dibromochloromethane____ | .666 | .619 | .598 | .614 | .628 | .625 | 4.1 |
| 1,1,2-Trichloroethane____ | .345 | .312 | .308 | .318 | .321 | .321 | 4.5 |
| Benzene_____ | 1.003 | .898 | .939 | .978 | .952 | .954 | 4.2 |
| trans-1,3-Dichloropropene__ | .703 | .645 | .634 | .653 | .668 | .661 | 4.1 |
| Bromoform_____# | .603 | .574 | .551 | .568 | .565 | .572 | 3.3# |
| 4-Methyl-2-Pentanone____ | .347 | .322 | .277 | .281 | .249 | .295 | 13.2 |
| 2-Hexanone_____ | .280 | .278 | .233 | .246 | .227 | .253 | 9.9 |
| Tetrachloroethene_____ | .537 | .469 | .447 | .448 | .446 | .470 | 8.3 |
| 1,1,2,2-Tetrachloroethane__# | .553 | .517 | .404 | .528 | .491 | .498 | 11.5# |
| Toluene_____* | .878 | .789 | .784 | .796 | .780 | .805 | 5.1* |
| Chlorobenzene_____# | 1.302 | 1.157 | 1.121 | 1.134 | 1.111 | 1.165 | 6.7# |
| Ethylbenzene_____* | .715 | .644 | .630 | .647 | .641 | .655 | 5.2* |
| Styrene_____ | 1.776 | 1.594 | 1.560 | 1.622 | 1.641 | 1.639 | 5.0 |
| Xylene (total)_____ | .983 | .892 | .878 | .916 | .936 | .921 | 4.4 |
| Toluene-d8_____ | 1.475 | 1.457 | 1.461 | 1.470 | 1.498 | 1.472 | 1.1 |
| Bromofluorobenzene____ | 1.077 | 1.078 | 1.070 | 1.085 | 1.130 | 1.088 | 2.2 |
| 1,2-Dichloroethane-d4____ | 3.407 | 3.393 | 3.454 | 3.586 | 3.610 | 3.490 | 2.9 |

TOTAL ION CHROMATOGRAM



Data File: >C0915::U0

Quant Output File: ^C0915::AQ

Name: UDA,900926,C

Misc: QC70357US,QU70357,L,5,,

2UL CAL I,II,XUOA

Id File: IC1171::US

Title: IFB, PP/UDA, TCL, XUOA13

Last Calibration: 900926 15:42

Operator ID: MGRMS

Quant Time: 900926 15:57

Injected at: 900926 14:41

QUANT REPORT

Operator ID: MGRMS
 Quant Rev: 7 Quant Time: 900926 15:57
 Output File: ^C0915::AQ Injected at: 900926 14:41
 Data File: >C0915::U0 Dilution Factor: 1.00000
 Name: UOA,900926,E VS'DJ
 Misc: QC70357US,QU70357,L,5,, 2UL CAL I,II,XUOA

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------|-----------------------------------|------------------|----------------|------------------|-------------------|---------------|----------------|
| 1) | *Bromochloromethane | 11.05 | 244 | 67956 | 250.00 | NG | 99 |
| 2) | Methyl chloride | 2.16 | 15 | 20268 | 115.01 | NG | 98 |
| 3) | Methyl bromide | 3.17 | 41 | 12898 | 159.04 | NG | 95 |
| 4) | Dichlorodifluoromethane | 3.95 | 61 | 28289 | 138.98 | NG | 93 |
| 5) | Vinyl chloride | 3.99 | 62 | 43431 | 134.61 | NG | 94 |
| 6) | Chloroethane | 5.15 | 92 | 27553 | 122.53 | NG | 93 |
| 7) | Methylene chloride | 7.52 | 153 | 52013 | 117.33 | NG | 95 |
| 8) | Acrolein | 8.30 | 173 | 55190 | 1636.79 | NG | 86 |
| 9) | Acetone | 8.34 | 174 | 16459M | 95.88 | NG | |
| 10) | Acrylonitrile | 9.11 | 194 | 12522 | 176.63 | NG | 92 |
| 11) | Carbon disulfide | 9.23 | 197 | 139206 | 111.88 | NG | 97 |
| 12) | Trichlorofluoromethane | 9.85 | 213 | 129947 | 115.44 | NG | 93 |
| 13) | 1,1-Dichloroethylene | 10.63 | 233 | 45358 | 114.17 | NG | 91 |
| 14) | 1,1-Dichloroethane | 11.98 | 268 | 84606 | 109.42 | NG | 94 |
| 15) | Tetrahydrofuran | 12.06 | 270 | 5680 | 3163.62 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.76 | 288 | 45577 | 83.81 | NG | 99 |
| 17) | Chloroform | 13.38 | 304 | 149809 | 151.86 | NG | 97 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 14.04 | 321 | 231517 | 333.78 | NG | 95 |
| 19) | 1,2-Dichloroethane | 14.12 | 323 | 128237 | 144.20 | NG | 97 |
| 20) | *1,4-Difluorobenzene | 21.84 | 522 | 339584 | 250.00 | NG | 99 |
| 21) | Methyl ethyl ketone | 14.12 | 323 | 3600 | 14.05 | NG | 67 |
| 22) | 1,1,1-Trichloroethane | 15.48 | 358 | 146924 | 138.75 | NG | 96 |
| 22) | Carbon tetrachloride | 15.48 | 358 | 15976 | 17.88 | NG | 93 |
| 23) | Carbon tetrachloride | 15.90 | 369 | 145279 | 154.62 | NG | 98 |
| 24) | Vinyl acetate | 16.10 | 374 | 68655 | 88.17 | NG | 99 |
| 25) | Dichlorobromomethane | 16.45 | 383 | 122422 | 135.70 | NG | 98 |
| 26) | 1,2-Dichloropropane | 17.88 | 420 | 38455 | 71.03 | NG | 92 |
| 27) | cis-1,3-Dichloropropylene | 18.16 | 427 | 71381 | 92.25 | NG | 92 |
| 28) | Trichloroethylene | 18.70 | 441 | 65499 | 105.06 | NG | 92 |
| 29) | Chlorodibromomethane | 19.36 | 458 | 90526 | 134.62 | NG | 99 |
| 30) | bis(Chloromethyl)ether | 19.36 | 458 | 37275 | 464.95 | NG | 100 |
| 31) | Benzene | 19.24 | 455 | 136244 | 80.94 | NG | 89 |
| 32) | 1,1,2-Trichloroethane | 19.44 | 460 | 46896 | 99.99 | NG | 93 |
| 33) | trans-1,3-Dichloropropylene | 19.44 | 460 | 95520 | 117.98 | NG | 89 |
| 34) | 2-Chloroethylvinyl ether | 20.52 | 488 | 24448 | 83.44 | NG | 100 |
| 35) | Bromoform | 22.08 | 528 | 81887 | 212.65 | NG | 95 |
| 36) | *Chlorobenzene-d5 | 26.79 | 649 | 326537 | 250.00 | NG | 78 |
| 37) | Methyl-iso-butyl ketone | 22.54 | 540 | 45299 | 81.27 | NG | 91 |
| 38) | 2-Hexanone | 24.10 | 580 | 36575 | 72.64 | NG | 87 |
| 39) | 1,1,2,2-Tetrachloroethane | 24.41 | 588 | 72212 | 81.69 | NG | 99 |
| 40) | Tetrachloroethylene | 24.41 | 588 | 70134 | 64.45 | NG | 95 |
| 41) | Toluene-DB (SURR) | 25.53 | 617 | 481657 | 231.97 | NG | 90 |
| 42) | Toluene | 25.73 | 622 | 114626 | 45.85 | NG | 97 |

QUANT REPORT

Page 2

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900926 15:57
 Output File: ^C0915::AQ Injected at: 900926 14:41
 Data File: >C0915::U0 Dilution Factor: 1.00000
 Name: VOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 2UL CAL 1,11,XVOA

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 43) | Chlorobenzene | 26.90 | 652 | 170050 | 92.20 | NG | 97 |
| 44) | Ethylbenzene | 29.04 | 707 | 93449 | 27.15 | NG | 79 |
| 45) | p-Bromofluorobenzene (SURR) | 31.84 | 779 | 351619 | 263.74 | NG | 85 |
| 46) | Styrene | 33.51 | 822 | 231937 | 85.74 | NG | 89 |
| 47) | m-Xylene | 33.82 | 830 | 125328 | 37.63 | NG | 90 |
| 48) | o+p-Xylenes | 34.98 | 860 | 256790 | 40.41 | NG | 88 |

* Compound is ISTD

Eak 12/15/90

QUANT REPORT

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900926 16:20
 Output File: ^C0914::AQ Injected at: 900926 13:52
 Data File: >C0914::U0 Dilution Factor: 1.00000
 Name: UOA,900926,C
 Misc: QC70357VS,QU70357,L,5,, 5UL CAL I,II,XVOA

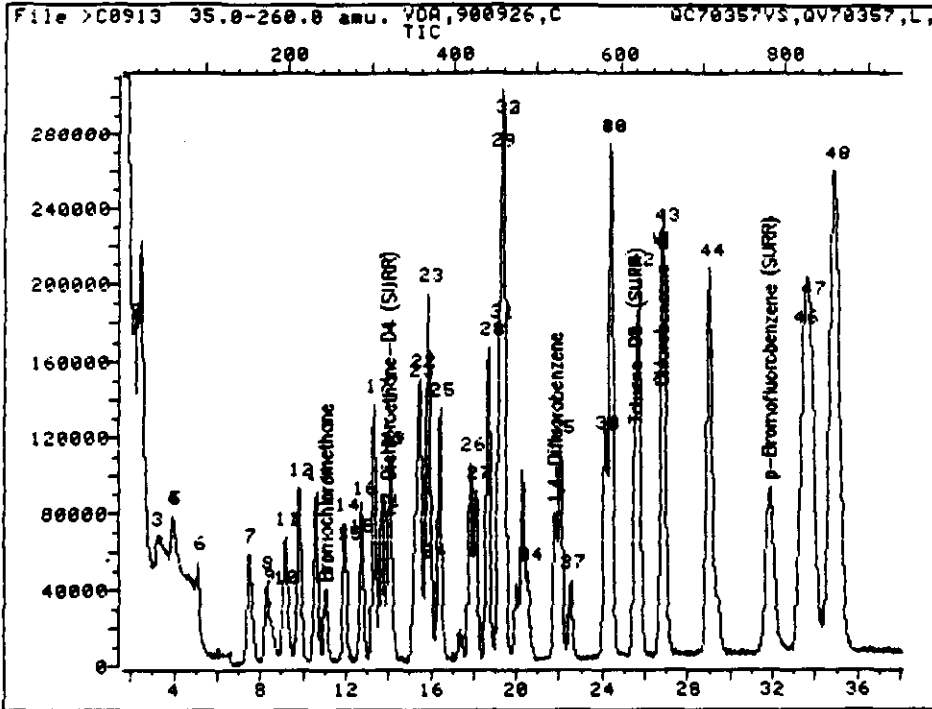
ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 16:16

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 39) | 1,1,2,2-Tetrachloroethane | 24.40 | 588 | 164492 | 259.15 | NG | 97 |
| 40) | Tetrachloroethylene | 24.44 | 589 | 149308 | 249.68 | NG | 98 |
| 41) | Toluene-D8 (SURR) | 25.53 | 617 | 463917 | 247.45 | NG | 90 |
| 42) | Toluene | 25.72 | 622 | 251169 | 244.99 | NG | 95 |
| 43) | Chlorobenzene | 26.90 | 652 | 368430 | 248.34 | NG | 98 |
| 44) | Ethylbenzene | 29.03 | 707 | 205107 | 245.75 | NG | 78 |
| 45) | p-Bromofluorobenzene (SURR) | 31.87 | 780 | 343104 | 247.66 | NG | 85 |
| 46) | Styrene | 33.50 | 822 | 507414 | 243.18 | NG | 90 |
| 47) | m-Xylene | 33.81 | 830 | 272301 | 242.18 | NG | 92 |
| 48) | o+p-Xylenes | 34.94 | 859 | 568101 | 484.31 | NG | 87 |

* Compound is ISTD

En 10/15/02

TOTAL ION CHROMATOGRAM



Data File: >C0913::U0

Quant Output File: ^C0913::AQ

Name: VOA,900926,C

Misc: QC70357VS,QU70357,L,5,,

10UL CAL I,II,XVOA

Id File: IC1171::US

Title: IFB, PP/VOA, TCL, XVOA13

Last Calibration: 900926 15:42

Operator ID: MGRMS

Quant Time: 900926 15:53

Injected at: 900926 13:03

QUANT REPORT

Operator ID: MGRMS
 Output File: ^C0913::AQ
 Data File: >C0913::U0
 Name: VOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 10UL CAL 1,11,XVOA

Quant Rev: 7
 Quant Time: 900926 15:53
 Injected at: 900926 13:03
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|---------------------------------------|------------------|----------------|-------------------|--------------------|---------------|----------------|
| 1) *Bromochloromethane | 11.05 | 244 | 68225 | 250.00 | NG | 99 |
| 2) Methyl chloride | 2.16 | 15 | 84438 | 477.26 | NG | 97 |
| 3) Methyl bromide | 3.21 | 42 | 29462 | 361.85 | NG | 90 |
| 4) Dichlorodifluoromethane | 3.95 | 61 | 95209 | 465.90 | NG | 97 |
| 5) Vinyl chloride | 3.99 | 62 | 144780 | 446.97 | NG | 97 |
| 6) Chloroethane | 5.15 | 92 | 108717 | 481.55 | NG | 97 |
| 7) Methylene chloride | 7.48 | 152 | 221249 | 497.13 | NG | 98 |
| 8) Acrolein | 8.30 | 173 | 265559 | 7844.73 | NG | 91 |
| 9) Acetone | 8.37 | 175 | 76485 | 443.82 | NG | 80 |
| 10) Acrylonitrile | 9.07 | 193 | 57919 | 813.74 | NG | 78 |
| 11) Carbon disulfide | 9.19 | 196 | 628910 | 503.48 | NG | 98 |
| 12) Trichlorofluoromethane | 9.81 | 212 | 550069 | 486.72 | NG | 95 |
| 13) 1,1-Dichloroethylene | 10.62 | 233 | 196535 | 492.73 | NG | 90 |
| 14) 1,1-Dichloroethane | 11.94 | 267 | 364980 | 470.16 | NG | 95 |
| 15) Tetrahydrofuran | 12.02 | 269 | 22793 | 12645.10 | NG | 100 |
| 15) Tetrahydrofuran | 12.68 | 286 | 3651 | 2025.58 | NG | 100 |
| 16) 1,2-Trans-dichloroethylene | 12.72 | 287 | 192283 | 352.18 | NG | 96 |
| 17) Chloroform | 13.34 | 303 | 633220 | 639.36 | NG | 96 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.00 | 320 | 235624 | 338.36 | NG | 99 |
| 19) 1,2-Dichloroethane | 14.12 | 323 | 532039 | 595.89 | NG | 97 |
| 20) *1,4-Difluorobenzene | 21.80 | 521 | 340907 | 250.00 | NG | 99 |
| 21) Methyl ethyl ketone | 14.08 | 322 | 15681 | 60.98 | NG | 86 |
| 22) 1,1,1-Trichloroethane | 15.48 | 358 | 623521 | 586.57 | NG | 95 |
| 23) Carbon tetrachloride | 15.44 | 357 | 66844 | 78.86 | NG | 98 |
| 23) Carbon tetrachloride | 15.86 | 368 | 619317 | 656.56 | NG | 97 |
| 24) Vinyl acetate | 16.06 | 373 | 276192 | 353.33 | NG | 95 |
| 25) Dichlorobromomethane | 16.45 | 383 | 526976 | 581.86 | NG | 98 |
| 26) 1,2-Dichloropropane | 17.84 | 419 | 168757 | 310.51 | NG | 98 |
| 27) cis-1,3-Dichloropropylene | 18.12 | 426 | 315197 | 405.77 | NG | 94 |
| 28) Trichloroethylene | 18.70 | 441 | 309255 | 494.10 | NG | 88 |
| 29) Chlorodibromomethane | 19.32 | 457 | 408046 | 604.43 | NG | 92 |
| 30) bis(Chloromethyl)ether | 19.32 | 457 | 159028 | 1985.88 | NG | 100 |
| 31) Benzene | 19.20 | 454 | 640063 | 378.77 | NG | 91 |
| 32) 1,1,2-Trichloroethane | 19.44 | 460 | 209992 | 445.98 | NG | 88 |
| 33) trans-1,3-Dichloropropylene | 19.44 | 460 | 432174 | 531.71 | NG | 91 |
| 34) 2-Chloroethylvinyl ether | 20.52 | 488 | 118333 | 402.29 | NG | 100 |
| 35) Bromoform | 22.08 | 528 | 375897 | 972.35 | NG | 97 |
| 36) *Chlorobenzene-d5 | 26.74 | 648 | 322654 | 250.00 | NG | 78 |
| 37) Methyl-iso-butyl ketone | 22.54 | 540 | 178675 | 324.43 | NG | 93 |
| 38) 2-Hexanone | 24.09 | 580 | 150372 | 302.24 | NG | 88 |
| 39) 1,1,2,2-Tetrachloroethane | 24.40 | 588 | 260722 | 298.50 | NG | 97 |
| 40) Tetrachloroethylene | 24.40 | 588 | 288656 | 268.45 | NG | 99 |
| 41) Toluene-D8 (SURR) | 25.53 | 617 | 471320 | 229.73 | NG | 90 |

QUANT REPORT

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900926 15:53
 Jutput File: ^C0913::AQ Injected at: 900926 13:03
 Data File: >C0913::U0 Dilution Factor: 1.00000
 Name: VOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 10UL CAL I,II,XVOA

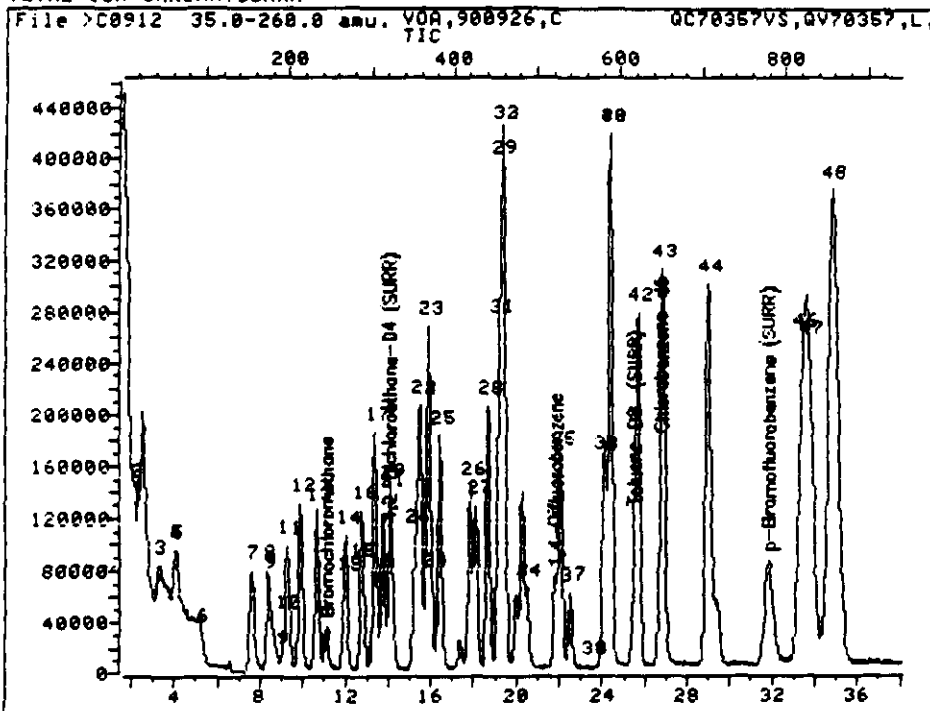
ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|--------|-------|----|
| 42) | Toluene | 25.72 | 622 | 505696 | 204.72 | NG | 95 |
| 43) | Chlorobenzene | 26.90 | 652 | 723198 | 396.83 | NG | 98 |
| 44) | Ethylbenzene | 29.03 | 707 | 406333 | 119.47 | NG | 78 |
| 45) | p-Bromofluorobenzene (SURR) | 31.83 | 779 | 345303 | 262.12 | NG | 86 |
| 46) | Styrene | 33.46 | 821 | 1006763 | 376.65 | NG | 90 |
| 47) | m-Xylene | 33.77 | 829 | 541251 | 164.48 | NG | 91 |
| 48) | o+p-Xylenes | 34.94 | 859 | 1133798 | 180.56 | NG | 89 |

* Compound is ISTD

EMC 4/15/90

TOTAL ION CHROMATOGRAM



Data File: >C0912::U0 Quant Output File: ^C0912::AQ
 Name: VOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 15UL CAL I,II,XVOA

Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 15:42

Operator ID: MGRMS
 Quant Time: 900926 15:49
 Injected at: 900926 12:13

QUANT REPORT

Page 2

Operator ID: MGRMS
 Output File: ^C0912::AQ
 Data File: >C0912::U0
 Name: UDA,900926,C
 Misc: QC70357US,QU70357,L,5,, 15UL CAL I,II,XUOA

Quant Rev: 7 Quant Time: 900926 15:49
 Injected at: 900926 12:13
 Dilution Factor: 1.00000

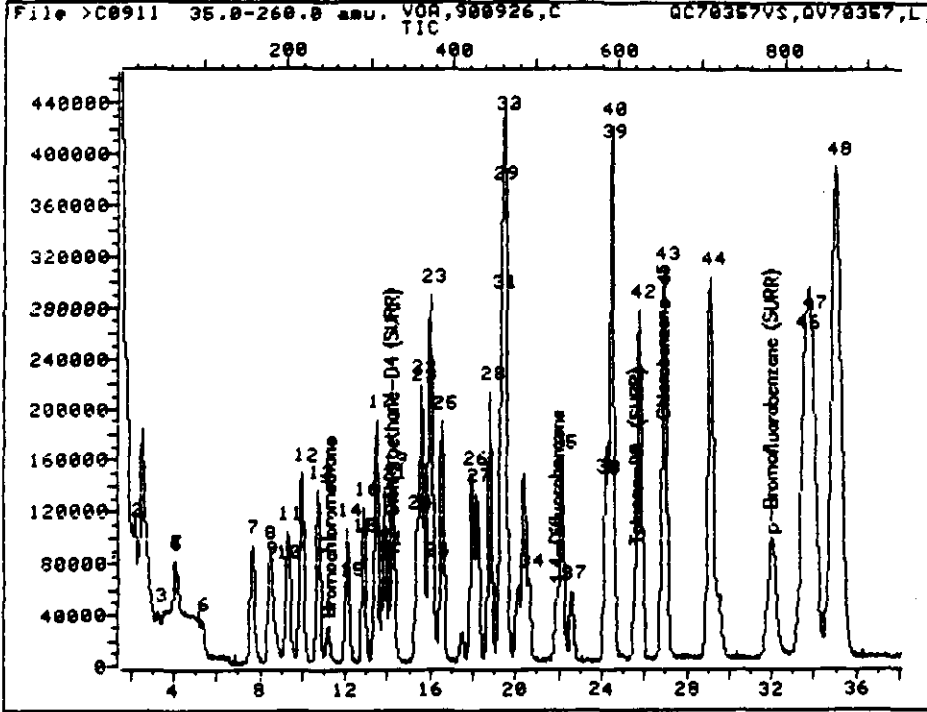
ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|--------|-------|----|
| 39) | 1,1,2,2-Tetrachloroethane | 24.40 | 587 | 472274 | 584.70 | NG | 97 |
| 40) | Tetrachloroethylene | 24.40 | 587 | 401360 | 403.63 | NG | 97 |
| 41) | Toluene-D8 (SURR) | 25.52 | 616 | 438728 | 231.24 | NG | 95 |
| 42) | Toluene | 25.72 | 621 | 712541 | 311.93 | NG | 95 |
| 43) | Chlorobenzene | 26.86 | 650 | 1015436 | 602.53 | NG | 98 |
| 44) | Ethylbenzene | 28.99 | 705 | 578703 | 183.99 | NG | 78 |
| 45) | p-Bromofluorobenzene (SURR) | 31.83 | 778 | 323728 | 265.74 | NG | 84 |
| 46) | Styrene | 33.46 | 820 | 1452199 | 587.50 | NG | 91 |
| 47) | m-Xylene | 33.77 | 828 | 776028 | 255.02 | NG | 94 |
| 48) | o+p-Xylenes | 34.86 | 856 | 1640544 | 282.52 | NG | 89 |

* Compound is ISTD

EAK 10/15/90

TOTAL ION CHROMATOGRAM



Data File: >C0911::U0

Quant Output File: ^C0911::AQ

Name: UOA,900926,C

Misc: QC70357VS,QU70357,L,5,,

20UL CAL 1,11,XVOA

Id File: IC1171::US

Title: IFB, PP/VOA, TCL, XVOA13

Last Calibration: 900926 15:42

Operator ID: MGRMS

Quant Time: 900926 15:43

Injected at: 900926 11:24

QUANT REPORT

Page 2

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900926 15:43
 Output File: ^C0911::AQ Injected at: 900926 11:24
 Data File: >C0911::U0 Dilution Factor: 1.00000
 Name: UOA,900926,C
 Misc: QC70357US,QU70357,L,5,, 20UL CAL I,II,XVOA

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900926 15:42

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|--------|-------|----|
| 41) | Toluene-D8 (SURR) | 25.59 | 619 | 342513 | 235.64 | NG | 92 |
| 42) | Toluene | 25.79 | 624 | 712783 | 407.28 | NG | 94 |
| 43) | Chlorobenzene | 26.97 | 653 | 1015905 | 786.82 | NG | 96 |
| 44) | Ethylbenzene | 29.10 | 708 | 586369 | 243.34 | NG | 77 |
| 45) | p-Bromofluorobenzene (SURR) | 31.94 | 781 | 258347 | 276.81 | NG | 82 |
| 46) | Styrene | 33.57 | 823 | 1500542 | 792.37 | NG | 92 |
| 47) | m-Xylene | 33.88 | 831 | 817849 | 350.81 | NG | 91 |
| 48) | o+p-Xylenes | 35.05 | 861 | 1711432 | 384.69 | NG | 88 |

* Compound is ISTD

EAW 10/15/20

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name:ETCNJ

Contract:

Lab Code:F2

Case No.:

SAS No.:

SDG No.:

Instrument ID:GC/MS C

Calibration Date: 09/27/90

Time: 1656

Lab File ID: >C0927

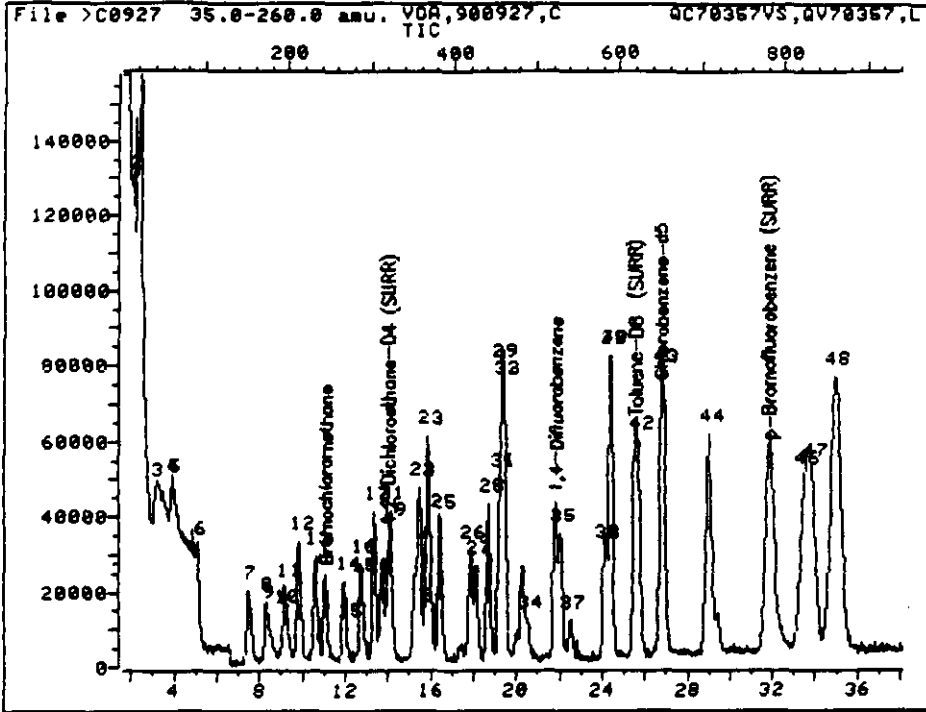
Init Calib. Dates(s): 09/26/90 09/26/90

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) PACK

Min RRF50 for SPCC(*) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 25.0%

| COMPOUND | RRF | RRF50 | %D |
|----------------------------|-------|-------|-------|
| Chloromethane | .637 | 1.077 | 69.0* |
| Bromomethane | .289 | .346 | 19.8 |
| Vinyl Chloride | 1.237 | 1.461 | 18.1* |
| Chloroethane | .895 | 1.032 | 15.4 |
| Methylene Chloride | 1.655 | 1.661 | .4 |
| Acetone | .606 | .713 | 17.7 |
| Carbon Disulfide | 4.700 | 4.402 | 6.3 |
| 1,1-Dichloroethene | 1.480 | 1.375 | 7.1* |
| 1,1-Dichloroethane | 2.753 | 2.529 | 8.1* |
| 1,2-Dichloroethene (total) | 1.473 | 1.335 | 9.4 |
| Chloroform | 4.822 | 4.530 | 6.1* |
| 1,2-Dichloroethane | 4.099 | 3.918 | 4.4 |
| 2-Butanone | .024 | .024 | 2.2 |
| 1,1,1-Trichloroethane | .972 | 1.081 | 11.2 |
| Carbon Tetrachloride | .968 | 1.119 | 15.6 |
| Vinyl Acetate | .459 | .421 | 8.3 |
| Bromodichloromethane | .815 | .836 | 2.5 |
| 1,2-Dichloropropane | .254 | .238 | 6.2* |
| cis-1,3-Dichloropropene | .478 | .465 | 2.7 |
| Trichloroethene | .437 | .442 | 1.3 |
| Dibromochloromethane | .625 | .654 | 4.7 |
| 1,1,2-Trichloroethane | .321 | .301 | 6.1 |
| Benzene | .954 | .840 | 11.9 |
| trans-1,3-Dichloropropene | .661 | .650 | 1.7 |
| Bromoform | .572 | .584 | 2.1* |
| 4-Methyl-2-Pentanone | .295 | .283 | 4.0 |
| 2-Hexanone | .253 | .264 | 4.4 |
| Tetrachloroethene | .470 | .491 | 4.6 |
| 1,1,2,2-Tetrachloroethane | .498 | .437 | 12.3* |
| Toluene | .805 | .749 | 7.0* |
| Chlorobenzene | 1.165 | 1.161 | .3* |
| Ethylbenzene | .655 | .671 | 2.4* |
| Styrene | 1.639 | 1.583 | 3.4 |
| Xylene (total) | .921 | .970 | 5.3 |
| Toluene-d8 | 1.472 | 1.460 | .8 |
| Bromofluorobenzene | 1.088 | 1.129 | 3.8 |
| 1,2-Dichloroethane-d4 | 3.490 | 3.231 | 7.4 |

TOTAL ION CHROMATOGRAM



Data File: >C0927::U4 Quant Output File: ^C0927::AQ
 Name: UOA,900927,C
 Misc: QC70357US,QU70357,L,5,, 5UL CAL I,II,XUOA

Id File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900926 16:16

Operator ID: MGRMS
 Quant Time: 900927 17:35
 Injected at: 900927 16:56

QUANT REPORT

Page 2

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900927 17:35
 Output File: ^C0927::AQ Injected at: 900927 16:56
 Data File: >C0927::U4 Dilution Factor: 1.00000
 Name: VOA,900927,C
 Misc: QC70357US,QU70357,L,5,, 5UL CAL I,II,XVOA

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 900926 16:16

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 42) | Toluene | 25.71 | 622 | 128494 | 232.51 | NG | 97 |
| 43) | Chlorobenzene | 26.89 | 652 | 199297 | 249.22 | NG | 98 |
| 44) | Ethylbenzene | 29.02 | 707 | 115137 | 255.93 | NG | 77 |
| 45) | p-Bromofluorobenzene (SURR) | 31.82 | 779 | 193769 | 259.47 | NG | 92 |
| 46) | Styrene | 33.53 | 823 | 271632 | 241.51 | NG | 85 |
| 47) | m-Xylene | 33.84 | 831 | 153657 | 253.53 | NG | 94 |
| 48) | o+p-Xylenes | 34.96 | 860 | 333036 | 526.71 | NG | 87 |

* Compound is ISTD

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:ETCNJ

Contract:

Lab Code:F2

Case No.:

SAS No.:

SDG No.:

Instrument ID:GC/MS C

Calibration Date(s) 10/03/90

10/03/90

Matrix:(soil/water) WATER

Level:(low/med) LOW

Column:(pack/cap) PACK

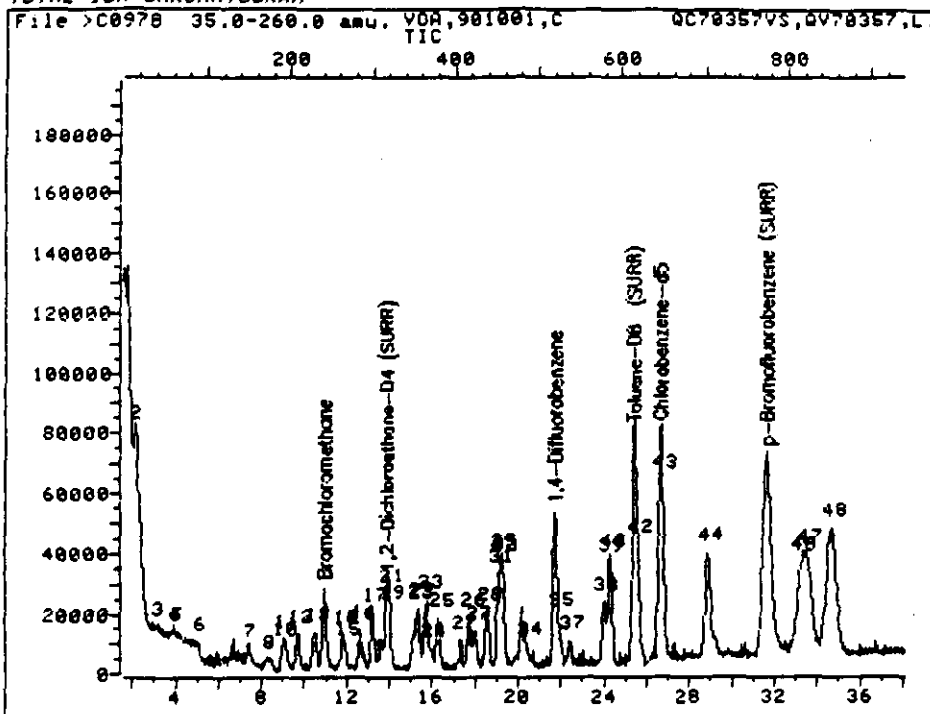
Min RRF for SPCC(*) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

| LAB FILE ID: | RRF20 =>C0978 | RRF50 =>C0972 | | | | | |
|-----------------------------|---------------|---------------|--------|--------|--------|-------|-------|
| IRRF100=>C0977 | RRF150=>C0976 | RRF200=>C0975 | | | | | |
| COMPOUND | RRF20 | RRF50 | RRF100 | RRF150 | RRF200 | RRF | % RSD |
| Chloromethane_____* | .573 | .735 | 1.123 | .521 | 1.488 | .888 | 46.2# |
| Bromomethane_____ | .240 | .302 | .140 | .103 | .165 | .190 | 42.3 |
| Vinyl Chloride_____* | .705 | .854 | .683 | .478 | .857 | .715 | 21.8* |
| Chloroethane_____ | .426 | .567 | .460 | .366 | .622 | .488 | 21.4 |
| Methylene Chloride____ | 1.204 | 1.588 | 1.176 | 1.067 | 1.516 | 1.310 | 17.4 |
| Acetone_____ | .957 | .896 | .582 | .544 | .668 | .729 | 25.6 |
| Carbon Disulfide_____ | 5.944 | 7.468 | 4.613 | 7.465 | 6.830 | 6.464 | 18.7 |
| 1,1-Dichloroethene____* | 1.678 | 1.666 | 1.635 | 1.717 | 1.794 | 1.698 | 3.6* |
| 1,1-Dichloroethane____* | 3.608 | 3.578 | 3.360 | 3.473 | 3.611 | 3.526 | 3.1# |
| 1,2-Dichloroethene (total)_ | 1.660 | 1.572 | 1.556 | 1.657 | 1.641 | 1.617 | 3.0 |
| Chloroform_____* | 5.094 | 5.263 | 4.600 | 4.849 | 5.013 | 4.964 | 5.1* |
| 1,2-Dichloroethane____ | 4.185 | 4.302 | 3.690 | 3.764 | 3.965 | 3.981 | 6.6 |
| 2-Butanone_____ | .058 | .072 | .032 | .032 | .029 | .045 | 43.2 |
| 1,1,1-Trichloroethane____ | .911 | 1.039 | .750 | .799 | .812 | .862 | 13.3 |
| Carbon Tetrachloride____ | .826 | .940 | .699 | .736 | .765 | .793 | 11.9 |
| Vinyl Acetate_____ | .568 | .570 | .581 | .623 | .528 | .574 | 5.9 |
| Bromodichloromethane____ | .825 | .899 | .695 | .738 | .736 | .779 | 10.5 |
| 1,2-Dichloropropane____* | .356 | .331 | .323 | .347 | .323 | .336 | 4.5* |
| cis-1,3-Dichloropropene___ | .509 | .553 | .485 | .513 | .491 | .510 | 5.2 |
| Trichloroethene_____ | .408 | .373 | .353 | .366 | .362 | .372 | 5.7 |
| Dibromochloromethane____ | .545 | .538 | .511 | .528 | .520 | .528 | 2.6 |
| 1,1,2-Trichloroethane____ | .321 | .304 | .322 | .335 | .330 | .322 | 3.7 |
| Benzene_____ | 1.159 | 1.120 | 1.118 | 1.243 | 1.203 | 1.169 | 4.6 |
| trans-1,3-Dichloropropene_ | .686 | .732 | .650 | .663 | .673 | .681 | 4.6 |
| Bromoform_____* | .477 | .403 | .445 | .429 | .409 | .433 | 6.9# |
| 4-Methyl-2-Pentanone____ | .426 | .298 | .342 | .340 | .289 | .339 | 16.1 |
| 2-Hexanone_____ | .389 | .267 | .323 | .329 | .274 | .316 | 15.6 |
| Tetrachloroethene____ | .403 | .345 | .349 | .330 | .329 | .351 | 8.6 |
| 1,1,2,2-Tetrachloroethane_* | .548 | .412 | .537 | .535 | .458 | .498 | 12.1# |
| Toluene_____* | .849 | .845 | .817 | .867 | .888 | .853 | 3.1* |
| Chlorobenzene_____* | 1.214 | 1.141 | 1.135 | 1.119 | 1.161 | 1.154 | 3.2# |
| Ethylbenzene_____* | .743 | .730 | .699 | .678 | .738 | .718 | 3.9* |
| Styrene_____ | 1.859 | 1.866 | 1.696 | 1.586 | 1.760 | 1.753 | 6.7 |
| Xylene (total)_____ | 1.092 | 1.143 | .987 | .921 | 1.026 | 1.034 | 8.4 |
| Toluene-d8_____ | 1.456 | 1.544 | 1.498 | 1.432 | 1.463 | 1.478 | 3.0 |
| Bromofluorobenzene____ | 1.116 | 1.275 | 1.085 | 1.042 | 1.116 | 1.127 | 7.8 |
| 1,2-Dichloroethane-d4____ | 3.274 | 3.769 | 3.321 | 3.511 | 3.837 | 3.543 | 7.2 |

FORM UI UOA

1/87 Rev.

TOTAL ION CHROMATOGRAM



Data File: >C0978::U4

Quant Output File: ^C0978::AQ

Name: VOA,901001,C

Misc: QC70357VS,QV70357,L,5, 2UL

Id File: IC1171::US

Title: IFB, PP/VOA, TCL, XVOA13

Last Calibration: 901002 16:53

Operator ID: KB6656

Quant Time: 901003 16:40

Injected at: 901003 15:35

QUANT REPORT

Operator ID: KB6656
 Output File: ^C0978::AQ
 Data File: >C0978::U4
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5, 2UL

Quant Rev: 7 Quant Time: 901003 16:40
 Injected at: 901003 15:35
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|---------|-------|-----|
| 1) | *Bromochloromethane | 10.94 | 241 | 42566 | 250.00 | NG | 97 |
| 2) | Methyl chloride | 2.14 | 14 | 9755 | 55.41 | NG | 92 |
| 3) | Methyl bromide | 3.15 | 40 | 4089 | 145.96 | NG | 84 |
| 4) | Dichlorodifluoromethane | 3.88 | 59 | 8620 | 108.73 | NG | 92 |
| 5) | Vinyl chloride | 3.96 | 61 | 11997 | 105.39 | NG | 94 |
| 6) | Chloroethane | 5.12 | 91 | 7254 | 97.44 | NG | 84 |
| 7) | Methylene chloride | 7.42 | 150 | 20507 | 108.72 | NG | 99 |
| 8) | Acrolein | 8.31 | 173 | 23677 | 1213.71 | NG | 93 |
| 9) | Acetone | 8.31 | 173 | 16293 | 166.33 | NG | 89 |
| 10) | Acrylonitrile | 9.01 | 191 | 14008 | 163.35 | NG | 90 |
| 11) | Carbon disulfide | 9.08 | 193 | 101202 | 102.51 | NG | 97 |
| 12) | Trichlorofluoromethane | 9.70 | 209 | 69603 | 113.18 | NG | 91 |
| 13) | 1,1-Dichloroethylene | 10.48 | 229 | 28574 | 99.18 | NG | 93 |
| 14) | 1,1-Dichloroethane | 11.84 | 264 | 61430 | 105.91 | NG | 97 |
| 14) | 1,1-Dichloroethane | 12.65 | 285 | 14699 | 25.34 | NG | 65 |
| 15) | Tetrahydrofuran | 11.87 | 265 | 2073 | 46.33 | NG | 100 |
| 15) | Tetrahydrofuran | 11.95 | 267 | 2920 | 65.26 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.65 | 285 | 28272 | 101.38 | NG | 94 |
| 17) | Chloroform | 13.23 | 300 | 86729 | 108.39 | NG | 97 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.89 | 317 | 139355 | 253.44 | NG | 95 |
| 19) | 1,2-Dichloroethane | 14.05 | 321 | 71261 | 115.00 | NG | 96 |
| 20) | *1,4-Difluorobenzene | 21.73 | 519 | 210589 | 250.00 | NG | 98 |
| 21) | Methyl ethyl ketone | 14.01 | 320 | 4914 | 167.54 | NG | 99 |
| 22) | 1,1,1-Trichloroethane | 15.37 | 355 | 76758 | 120.91 | NG | 95 |
| 23) | Carbon tetrachloride | 15.33 | 354 | 8892 | 15.26 | NG | 81 |
| 23) | Carbon tetrachloride | 15.75 | 365 | 69546 | 119.37 | NG | 96 |
| 24) | Vinyl acetate | 15.99 | 371 | 47815 | 111.21 | NG | 97 |
| 25) | Dichlorobromomethane | 16.34 | 380 | 69506 | 117.20 | NG | 98 |
| 26) | 1,2-Dichloropropane | 17.73 | 416 | 30006 | 101.97 | NG | 99 |
| 27) | cis-1,3-Dichloropropylene | 17.38 | 407 | 2119 | 5.08 | NG | 71 |
| 27) | cis-1,3-Dichloropropylene | 18.00 | 423 | 42910 | 102.94 | NG | 92 |
| 28) | Trichloroethylene | 18.59 | 438 | 34378 | 110.49 | NG | 86 |
| 29) | Chlorodibromomethane | 19.21 | 454 | 45927 | 105.86 | NG | 99 |
| 30) | bis(Chloromethyl)ether | 19.17 | 453 | 18986 | 107.49 | NG | 100 |
| 31) | Benzene | 19.13 | 452 | 97590 | 99.50 | NG | 90 |
| 32) | 1,1,2-Trichloroethane | 19.32 | 457 | 27035 | 97.47 | NG | 84 |
| 33) | trans-1,3-Dichloropropylene | 19.32 | 457 | 57788 | 106.35 | NG | 88 |
| 34) | 2-Chloroethylvinyl ether | 20.45 | 486 | 21125 | 100.71 | NG | 100 |
| 35) | Bromoform | 21.92 | 524 | 40197 | 111.90 | NG | 98 |
| 36) | *Chlorobenzene-d5 | 26.65 | 643 | 243193 | 250.00 | NG | 80 |
| 37) | Methyl-iso-butyl ketone | 22.43 | 537 | 41488 | 124.66 | NG | 95 |
| 38) | 2-Hexanone | 23.98 | 577 | 37852 | 117.71 | NG | 83 |
| 39) | 1,1,2,2-Tetrachloroethane | 24.25 | 584 | 53349 | 105.59 | NG | 97 |

QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0978::AQ
 Data File: >C0978::U4
 Name: UOA,901001,C
 Misc: QC70357US,QV70357,L,5, 2UL

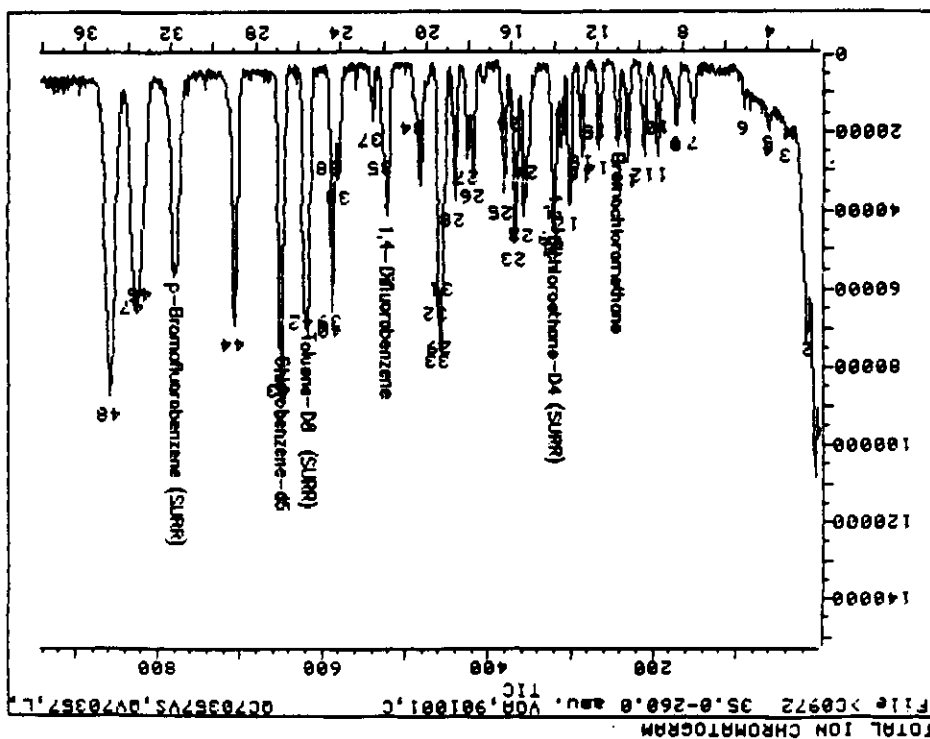
Quant Rev: 7 Quant Time: 901003 16:40
 Injected at: 901003 15:35
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|--------|-------|----|
| 40) | Tetrachloroethylene | 24.33 | 586 | 39161 | 116.70 | NG | 93 |
| 41) | Toluene-D8 (SURR) | 25.42 | 614 | 354032 | 248.58 | NG | 92 |
| 42) | Toluene | 25.61 | 619 | 82575 | 102.30 | NG | 93 |
| 43) | Chlorobenzene | 26.76 | 646 | 118136 | 104.58 | NG | 96 |
| 44) | Ethylbenzene | 28.86 | 700 | 72302 | 106.37 | NG | 78 |
| 45) | p-Bromofluorobenzene (SURR) | 31.65 | 772 | 271522 | 269.90 | NG | 83 |
| 46) | Styrene | 33.28 | 814 | 180878 | 112.95 | NG | 98 |
| 47) | m-Xylene | 33.56 | 821 | 102366 | 114.96 | NG | 96 |
| 48) | o+p-Xylenes | 34.72 | 851 | 212529 | 231.16 | NG | 90 |

* Compound is ISTD

Data File: >C0972:U4
 Name: V04,901001,C
 Misc: DC70357VS,QV70357,L,5,
 ID File: IC1171:US
 Title: IFB, RP/V04, TCL, XVD013
 Last Calibration: 901002 16:53
 Operator ID: KB6656
 Quant Time: 901003 10:39
 Injected at: 901003 10:01



QUANT REPORT

Operator ID: KB6656
 Output File: ^C0972::AQ
 Data File: >C0972::U4
 Name: UDA,901001,C
 Misc: QC70357US,QU70357,L,5,

Quant Rev: 7 Quant Time: 901003 10:39
 Injected at: 901003 10:01
 Dilution Factor: 1.00000

501

ID File: IC1171::US
 Title: IFB, PP/UDA, TCL, XUOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|---------|-------|-----|
| 1) | *Bromochloromethane | 11.00 | 243 | 33439 | 250.00 | NG | 92 |
| 2) | Methyl chloride | 2.12 | 14 | 24563 | 177.60 | NG | 92 |
| 3) | Methyl bromide | 3.17 | 41 | 10088 | 458.39 | NG | 89 |
| 4) | Dichlorodifluoromethane | 3.94 | 61 | 20601 | 330.79 | NG | 95 |
| 5) | Vinyl chloride | 3.98 | 62 | 28559 | 319.36 | NG | 99 |
| 6) | Chloroethane | 5.14 | 92 | 18945 | 323.95 | NG | 97 |
| 7) | Methylene chloride | 7.47 | 152 | 53114 | 358.46 | NG | 92 |
| 8) | Acrolein | 8.25 | 172 | 105297 | 6870.91 | NG | 88 |
| 9) | Acetone | 8.29 | 173 | 29974 | 389.52 | NG | 97 |
| 10) | Acrylonitrile | 9.02 | 192 | 30914 | 458.88 | NG | 93 |
| 11) | Carbon disulfide | 9.14 | 195 | 249726 | 321.98 | NG | 97 |
| 12) | Trichlorofluoromethane | 9.80 | 212 | 145098 | 300.35 | NG | 93 |
| 13) | 1,1-Dichloroethylene | 10.57 | 232 | 55695 | 246.08 | NG | 87 |
| 14) | 1,1-Dichloroethane | 11.93 | 267 | 119653 | 262.60 | NG | 98 |
| 15) | Tetrahydrofuran | 12.01 | 269 | 10461 | 297.63 | NG | 100 |
| 15) | Tetrahydrofuran | 12.67 | 286 | 2117 | 60.23 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.71 | 287 | 52582 | 240.03 | NG | 89 |
| 17) | Chloroform | 13.29 | 302 | 175991 | 279.98 | NG | 97 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.95 | 319 | 126037 | 291.78 | NG | 99 |
| 19) | 1,2-Dichloroethane | 14.06 | 322 | 143848 | 295.50 | NG | 97 |
| 20) | *1,4-Difluorobenzene | 21.78 | 521 | 154748 | 250.00 | NG | 95 |
| 21) | Methyl ethyl ketone | 14.03 | 321 | 11158 | 517.72 | NG | 91 |
| 22) | 1,1,1-Trichloroethane | 15.42 | 357 | 160852 | 344.79 | NG | 98 |
| 23) | Carbon tetrachloride | 15.42 | 357 | 16473 | 38.48 | NG | 93 |
| 23) | Carbon tetrachloride | 15.81 | 367 | 145452 | 339.75 | NG | 97 |
| 24) | Vinyl acetate | 15.23 | 352 | 49264 | 155.92 | NG | 78 |
| 24) | Vinyl acetate | 16.04 | 373 | 88161 | 279.03 | NG | 89 |
| 25) | Dichlorobromomethane | 16.39 | 382 | 139050 | 319.07 | NG | 95 |
| 26) | 1,2-Dichloropropane | 17.79 | 418 | 51182 | 236.69 | NG | 98 |
| 27) | cis-1,3-Dichloropropylene | 18.06 | 425 | 85652 | 279.63 | NG | 91 |
| 28) | Trichloroethylene | 18.64 | 440 | 57747 | 252.56 | NG | 78 |
| 29) | Chlorodibromomethane | 19.26 | 456 | 83180 | 260.92 | NG | 99 |
| 30) | bis(Chloromethyl)ether | 19.26 | 456 | 38771 | 298.72 | NG | 100 |
| 31) | Benzene | 19.18 | 454 | 173359 | 240.54 | NG | 89 |
| 32) | 1,1,2-Trichloroethane | 19.42 | 460 | 47019 | 230.69 | NG | 88 |
| 33) | trans-1,3-Dichloropropylene | 19.38 | 459 | 113305 | 283.77 | NG | 85 |
| 34) | 2-Chloroethylvinyl ether | 20.50 | 488 | 32319 | 209.68 | NG | 100 |
| 35) | Bromoform | 22.02 | 527 | 62340 | 236.16 | NG | 97 |
| 36) | *Chlorobenzene-d5 | 26.72 | 648 | 176233 | 250.00 | NG | 65 |
| 37) | Methyl-iso-butyl ketone | 22.48 | 539 | 52446 | 217.46 | NG | 91 |
| 38) | 2-Hexanone | 24.07 | 580 | 47055 | 201.94 | NG | 80 |
| 38) | 2-Hexanone | 24.50 | 591 | 2192 | 9.41 | NG | 96 |
| 39) | 1,1,2,2-Tetrachloroethane | 24.34 | 587 | 72692 | 198.54 | NG | 99 |

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QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0972::AQ
 Data File: >C0972::U4
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5,

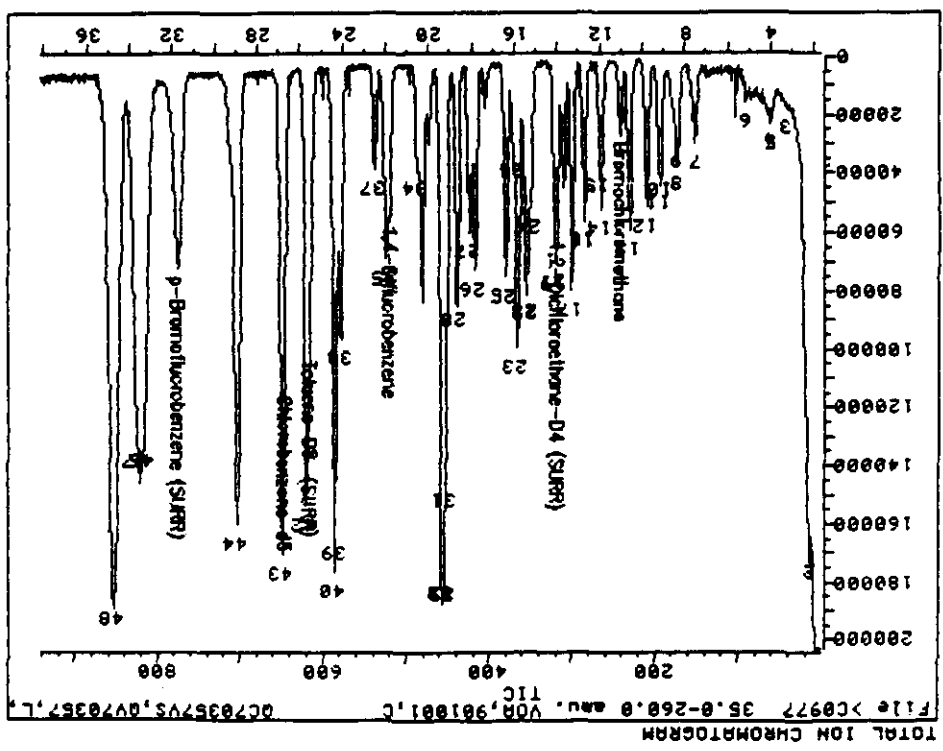
Quant Rev: 7 Quant Time: 901003 10:39
 Injected at: 901003 10:01
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|---------------------------------|-------|-------|--------|--------|-------|----|
| 40) Tetrachloroethylene | 24.38 | 588 | 60842 | 250.20 | NG | 98 |
| 41) Toluene-D8 (SURR) | 25.51 | 617 | 272171 | 263.72 | NG | 93 |
| 42) Toluene | 25.70 | 622 | 148919 | 254.58 | NG | 99 |
| 43) Chlorobenzene | 26.83 | 651 | 201006 | 245.54 | NG | 96 |
| 44) Ethylbenzene | 28.97 | 706 | 128615 | 261.10 | NG | 78 |
| 45) p-Bromofluorobenzene (SURR) | 31.72 | 777 | 224676 | 308.20 | NG | 73 |
| 46) Styrene | 33.39 | 820 | 328857 | 283.38 | NG | 98 |
| 47) m-Xylene | 33.70 | 828 | 189326 | 293.41 | NG | 96 |
| 48) o+p-Xylenes | 34.83 | 857 | 402974 | 604.82 | NG | 90 |

* Compound is ISTD

Data File: <C0977>:U4
 Name: VOA,901001,C
 Misc: DC70357VS,QV70357,L,5, 10UL
 ID File: IC1171:US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53
 Operator ID: KB6656
 Quant Time: 901003 16:43
 Injected at: 901003 14:45



QUANT REPORT

Operator ID: KB6656
 Output File: ^C0977::AQ
 Data File: >C0977::U4
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5, 10UL

Quant Rev: 7 Quant Time: 901003 16:43
 Injected at: 901003 14:45
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|----------|-------|-----|
| 1) | *Bromochloromethane | 10.94 | 240 | 40449 | 250.00 | NG | 95 |
| 2) | Methyl chloride | 2.17 | 14 | 90814 | 542.83 | NG | 97 |
| 3) | Methyl bromide | 3.22 | 41 | 11319 | 425.19 | NG | 98 |
| 4) | Dichlorodifluoromethane | 3.92 | 59 | 37369 | 496.04 | NG | 93 |
| 5) | Vinyl chloride | 3.96 | 60 | 55220 | 510.48 | NG | 97 |
| 6) | Chloroethane | 5.16 | 91 | 37245 | 526.51 | NG | 97 |
| 7) | Methylene chloride | 7.45 | 150 | 95143 | 530.82 | NG | 98 |
| 8) | Acrolein | 8.31 | 172 | 244743 | 13202.43 | NG | 95 |
| 9) | Acetone | 8.38 | 174 | 47100 | 506.01 | NG | 94 |
| 10) | Acrylonitrile | 9.04 | 191 | 58471 | 717.52 | NG | 97 |
| 11) | Carbon disulfide | 9.12 | 193 | 373170 | 397.76 | NG | 98 |
| 12) | Trichlorofluoromethane | 9.74 | 209 | 293996 | 503.09 | NG | 95 |
| 13) | 1,1-Dichloroethylene | 10.52 | 229 | 132241 | 483.02 | NG | 93 |
| 14) | 1,1-Dichloroethane | 11.87 | 264 | 271805 | 493.15 | NG | 95 |
| 15) | Tetrahydrofuran | 11.95 | 266 | 21639 | 508.97 | NG | 100 |
| 15) | Tetrahydrofuran | 12.65 | 284 | 2620 | 61.62 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.69 | 285 | 125902 | 475.11 | NG | 96 |
| 17) | Chloroform | 13.27 | 300 | 372092 | 489.36 | NG | 95 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.97 | 318 | 134320 | 257.07 | NG | 98 |
| 19) | 1,2-Dichloroethane | 14.05 | 320 | 298525 | 506.96 | NG | 97 |
| 20) | *1,4-Difluorobenzene | 21.77 | 519 | 216201 | 250.00 | NG | 98 |
| 21) | Methyl ethyl ketone | 14.01 | 319 | 14033 | 466.04 | NG | 95 |
| 22) | 1,1,1-Trichloroethane | 15.40 | 355 | 324364 | 497.66 | NG | 96 |
| 23) | Carbon tetrachloride | 15.36 | 354 | 35214 | 58.87 | NG | 95 |
| 23) | Carbon tetrachloride | 15.79 | 365 | 302117 | 505.11 | NG | 94 |
| 24) | Vinyl acetate | 15.21 | 350 | 80279 | 181.86 | NG | 78 |
| 24) | Vinyl acetate | 16.02 | 371 | 251350 | 569.41 | NG | 93 |
| 25) | Dichlorobromomethane | 16.37 | 380 | 300604 | 493.71 | NG | 97 |
| 26) | 1,2-Dichloropropane | 17.77 | 416 | 139502 | 461.75 | NG | 91 |
| 27) | cis-1,3-Dichloropropylene | 18.04 | 423 | 209777 | 490.21 | NG | 93 |
| 28) | Trichloroethylene | 18.66 | 439 | 152596 | 477.69 | NG | 87 |
| 29) | Chlorodibromomethane | 19.24 | 454 | 220994 | 496.17 | NG | 97 |
| 30) | bis(Chloromethyl)ether | 19.20 | 453 | 89735 | 494.86 | NG | 100 |
| 31) | Benzene | 19.17 | 452 | 483500 | 480.18 | NG | 94 |
| 32) | 1,1,2-Trichloroethane | 19.36 | 457 | 139077 | 488.40 | NG | 82 |
| 33) | trans-1,3-Dichloropropylene | 19.36 | 457 | 281099 | 503.90 | NG | 89 |
| 34) | 2-Chloroethylvinyl ether | 20.49 | 486 | 105157 | 488.32 | NG | 100 |
| 35) | Bromoform | 21.96 | 524 | 192370 | 521.61 | NG | 99 |
| 36) | *Chlorobenzene-d5 | 26.67 | 645 | 245701 | 250.00 | NG | 82 |
| 37) | Methyl-iso-butyl ketone | 22.46 | 537 | 167817 | 499.10 | NG | 94 |
| 38) | 2-Hexanone | 24.02 | 577 | 158671 | 488.41 | NG | 94 |
| 39) | 1,1,2,2-Tetrachloroethane | 24.29 | 584 | 264065 | 517.32 | NG | 97 |
| 40) | Tetrachloroethylene | 24.33 | 585 | 171398 | 505.56 | NG | 97 |

QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0977::AQ
 Data File: >C0977::U4
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5, 10UL

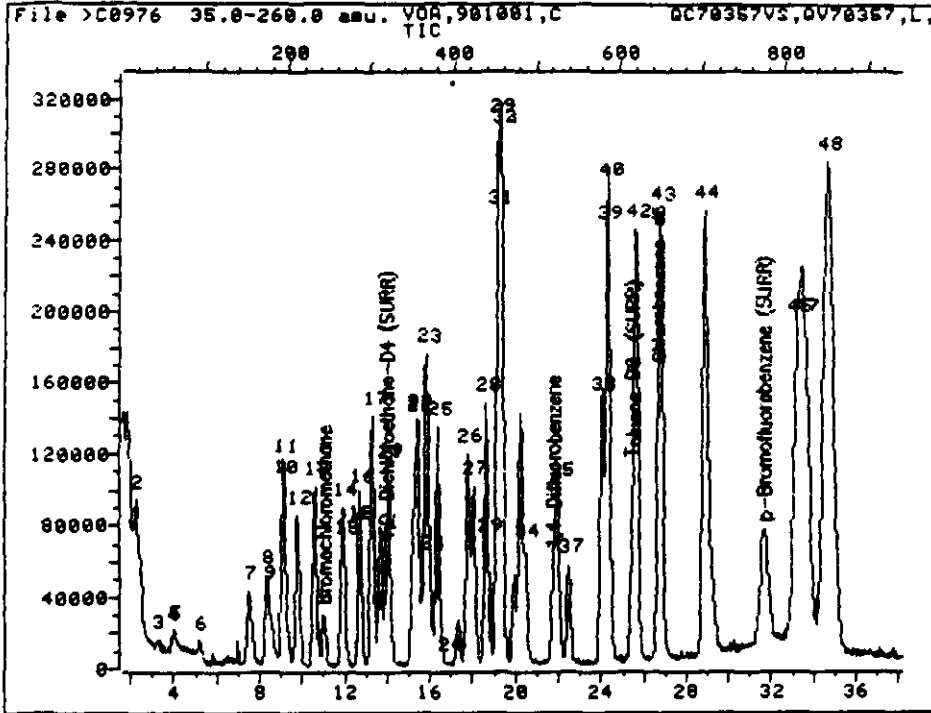
Quant Rev: 7 Quant Time: 901003 16:43
 Injected at: 901003 14:45
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|--------|---------|-------|----|
| 41) | Toluene-D8 (SURR) | 25.49 | 615 | 367994 | 255.75 | NG | 91 |
| 42) | Toluene | 25.65 | 619 | 401656 | 492.50 | NG | 95 |
| 43) | Chlorobenzene | 26.79 | 648 | 557897 | 488.82 | NG | 94 |
| 44) | Ethylbenzene | 28.88 | 702 | 343471 | 500.13 | NG | 79 |
| 45) | p-Bromofluorobenzene (SURR) | 31.67 | 774 | 266698 | 262.40 | NG | 84 |
| 46) | Styrene | 33.31 | 816 | 833355 | 515.07 | NG | 94 |
| 47) | m-Xylene | 33.62 | 824 | 460640 | 512.04 | NG | 97 |
| 48) | o+p-Xylenes | 34.70 | 852 | 969564 | 1043.78 | NG | 87 |

* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >C0976::U4

Quant Output File: ^C0976::AQ

Name: VOA,901001,C

Misc: QC70357US,QU70357,L,5, 15UL

Id File: IC1171::US

Title: IFB, PP/VOA, TCL, XVOA13

Last Calibration: 901002 16:53

Operator ID: KB6656

Quant Time: 901003 16:46

Injected at: 901003 13:55

QUANT REPORT

Operator ID: K86656
 Output File: ^C0976::AQ
 Data File: >C0976::U4
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5, 15UL

Quant Rev: 7 Quant Time: 901003 16:46
 Injected at: 901003 13:55
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|---------|----------|-------|-----|
| 1) | *Bromochloromethane | 10.98 | 242 | 44783 | 250.00 | NG | 99 |
| 2) | Methyl chloride | 2.21 | 16 | 69987 | 377.85 | NG | 99 |
| 3) | Methyl bromide | 3.22 | 42 | 13798 | 468.15 | NG | 97 |
| 4) | Dichlorodifluoromethane | 3.92 | 60 | 44312 | 531.28 | NG | 93 |
| 5) | Vinyl chloride | 4.00 | 62 | 64172 | 535.83 | NG | 98 |
| 6) | Chloroethane | 5.16 | 92 | 49183 | 627.98 | NG | 98 |
| 7) | Methylene chloride | 7.49 | 152 | 143319 | 722.22 | NG | 98 |
| 8) | Acrolein | 8.31 | 173 | 370346 | 18044.55 | NG | 90 |
| 9) | Acetone | 8.38 | 175 | 73037 | 708.72 | NG | 95 |
| 10) | Acrylonitrile | 9.04 | 192 | 138313 | 1533.03 | NG | 90 |
| 11) | Carbon disulfide | 9.12 | 194 | 1002979 | 965.61 | NG | 99 |
| 12) | Trichlorofluoromethane | 9.74 | 210 | 504835 | 780.28 | NG | 96 |
| 13) | 1,1-Dichloroethylene | 10.56 | 231 | 230701 | 761.10 | NG | 92 |
| 14) | 1,1-Dichloroethane | 11.87 | 265 | 466532 | 764.54 | NG | 96 |
| 15) | Tetrahydrofuran | 11.95 | 267 | 42526 | 903.44 | NG | 100 |
| 15) | Tetrahydrofuran | 12.61 | 284 | 8088 | 171.83 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.65 | 285 | 222632 | 758.83 | NG | 98 |
| 17) | Chloroform | 13.27 | 301 | 651450 | 773.84 | NG | 98 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.93 | 318 | 157254 | 271.83 | NG | 97 |
| 19) | 1,2-Dichloroethane | 14.05 | 321 | 505720 | 775.71 | NG | 97 |
| 20) | *1,4-Difluorobenzene | 21.73 | 519 | 235893 | 250.00 | NG | 99 |
| 21) | Methyl ethyl ketone | 14.01 | 320 | 22639 | 689.08 | NG | 94 |
| 22) | 1,1,1-Trichloroethane | 15.40 | 356 | 565553 | 795.28 | NG | 97 |
| 23) | Carbon tetrachloride | 15.37 | 355 | 57679 | 88.38 | NG | 92 |
| 23) | Carbon tetrachloride | 15.79 | 366 | 520609 | 797.74 | NG | 91 |
| 24) | Vinyl acetate | 15.99 | 371 | 440600 | 914.82 | NG | 97 |
| 24) | Vinyl acetate | 16.84 | 393 | 2264 | 4.70 | NG | 73 |
| 25) | Dichlorobromomethane | 16.34 | 380 | 522200 | 786.07 | NG | 95 |
| 26) | 1,2-Dichloropropane | 17.73 | 416 | 245780 | 745.62 | NG | 99 |
| 27) | cis-1,3-Dichloropropylene | 18.00 | 423 | 362807 | 777.03 | NG | 93 |
| 28) | Trichloroethylene | 18.59 | 438 | 258713 | 742.27 | NG | 87 |
| 29) | Chlorodibromomethane | 18.70 | 441 | 2087 | 4.29 | NG | 57 |
| 29) | Chlorodibromomethane | 19.21 | 454 | 373799 | 769.19 | NG | 96 |
| 30) | bis(Chloromethyl)ether | 19.21 | 454 | 152055 | 768.54 | NG | 100 |
| 31) | Benzene | 19.13 | 452 | 879768 | 800.80 | NG | 94 |
| 32) | 1,1,2-Trichloroethane | 19.32 | 457 | 237165 | 763.34 | NG | 87 |
| 33) | trans-1,3-Dichloropropylene | 19.32 | 457 | 469529 | 771.42 | NG | 92 |
| 34) | 2-Chloroethylvinyl ether | 20.41 | 485 | 176943 | 753.08 | NG | 100 |
| 35) | Bromoform | 21.96 | 525 | 303909 | 755.25 | NG | 99 |
| 36) | *Chlorobenzene-d5 | 26.65 | 644 | 269685 | 250.00 | NG | 85 |
| 37) | Methyl-iso-butyl ketone | 22.43 | 537 | 275155 | 745.56 | NG | 94 |
| 38) | 2-Hexanone | 23.98 | 577 | 266545 | 747.49 | NG | 95 |
| 39) | 1,1,2,2-Tetrachloroethane | 24.25 | 584 | 432936 | 772.72 | NG | 98 |

QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0976::AQ
 Data File: >C0976::U4
 Name: UOA,901001,C
 Misc: QC70357US,QU70357,L,5, 15UL

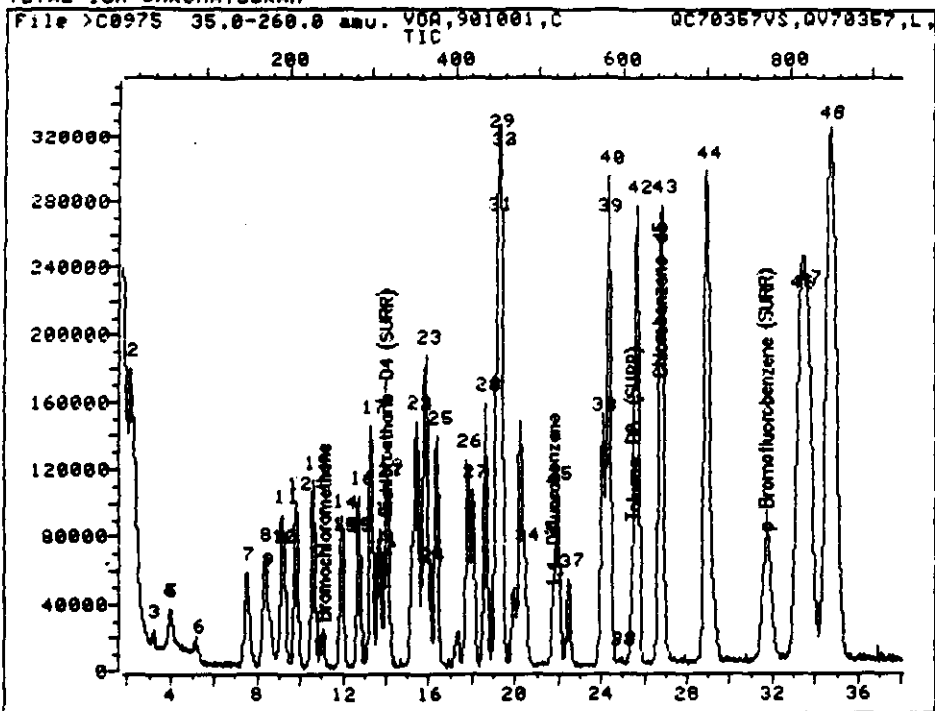
Quant Rev: 7 Quant Time: 901003 16:46
 Injected at: 901003 13:55
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|---------|-------|----|
| 40) | Tetrachloroethylene | 24.33 | 586 | 267038 | 717.61 | NG | 98 |
| 41) | Toluene-D8 (SURR) | 25.45 | 615 | 386124 | 244.48 | NG | 93 |
| 42) | Toluene | 25.61 | 619 | 701550 | 783.72 | NG | 97 |
| 43) | Chlorobenzene | 26.77 | 647 | 905631 | 722.93 | NG | 94 |
| 44) | Ethylbenzene | 28.86 | 701 | 548559 | 727.73 | NG | 80 |
| 45) | p-Bromofluorobenzene (SURR) | 31.62 | 772 | 280987 | 251.88 | NG | 83 |
| 46) | Styrene | 33.21 | 813 | 1283190 | 722.57 | NG | 91 |
| 47) | m-Xylene | 33.60 | 823 | 716403 | 725.52 | NG | 96 |
| 48) | o+p-Xylenes | 34.65 | 850 | 1489830 | 1461.23 | NG | 89 |

* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >C0975::U4 Quant Output File: ^C0975::AQ
 Name: VOA,901001,C
 Misc: QC70357US,QV70357,L,5, 20UL

Id File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

Operator ID: KB6656
 Quant Time: 901003 14:45
 Injected at: 901003 13:05

QUANT REPORT

Operator ID: KB6656 Quant Rev: 7 Quant Time: 901003 14:45
 Output File: ^C0975::AQ Injected at: 901003 13:05
 Data File: >C0975::U4 Dilution Factor: 1.00000
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5, 20UL

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|------------------------------|-------|-------|--------|----------|-------|-----|
| 1) | *Bromochloromethane | 11.07 | 240 | 33082 | 250.00 | NG | 93 |
| 2) | Methyl chloride | 2.19 | 11 | 196957 | 1439.46 | NG | 97 |
| 3) | Methyl bromide | 3.20 | 37 | 21780 | 1000.34 | NG | 96 |
| 4) | Dichlorodifluoromethane | 3.93 | 56 | 79248 | 1286.20 | NG | 93 |
| 5) | Vinyl chloride | 3.97 | 57 | 113382 | 1281.58 | NG | 94 |
| 6) | Chloroethane | 5.21 | 89 | 82361 | 1423.55 | NG | 96 |
| 7) | Methylene chloride | 7.50 | 148 | 200574 | 1368.24 | NG | 99 |
| 8) | Acrolein | 8.36 | 170 | 442510 | 29186.55 | NG | 91 |
| 9) | Acetone | 8.44 | 172 | 88336 | 1160.35 | NG | 92 |
| 10) | Acrylonitrile | 9.10 | 189 | 107700 | 1615.94 | NG | 99 |
| 11) | Carbon disulfide | 9.17 | 191 | 903789 | 1177.87 | NG | 99 |
| 12) | Trichlorofluoromethane | 9.83 | 208 | 590028 | 1234.52 | NG | 92 |
| 13) | 1,1-Dichloroethylene | 10.61 | 228 | 237397 | 1060.21 | NG | 87 |
| 14) | 1,1-Dichloroethane | 11.93 | 262 | 477859 | 1060.08 | NG | 95 |
| 15) | Tetrahydrofuran | 12.00 | 264 | 35476 | 1020.24 | NG | 100 |
| 15) | Tetrahydrofuran | 12.66 | 281 | 8973 | 258.05 | NG | 100 |
| 16) | 1,2-Trans-dichloroethylene | 12.74 | 283 | 217086 | 1001.64 | NG | 92 |
| 17) | Chloroform | 13.32 | 298 | 663299 | 1066.60 | NG | 96 |
| 18) | 1,2-Dichloroethane-D4 (SURR) | 13.98 | 315 | 126950 | 297.07 | NG | 98 |
| 19) | 1,2-Dichloroethane | 14.10 | 318 | 524656 | 1089.39 | NG | 96 |
| 20) | *1,4-Difluorobenzene | 21.74 | 515 | 183065 | 250.00 | NG | 98 |
| 21) | Methyl ethyl ketone | 14.06 | 317 | 21244 | 833.22 | NG | 93 |
| 22) | 1,1,1-Trichloroethane | 15.42 | 352 | 594519 | 1077.26 | NG | 97 |
| 23) | Carbon tetrachloride | 15.42 | 352 | 58362 | 115.24 | NG | 99 |
| 23) | Carbon tetrachloride | 15.84 | 363 | 560082 | 1105.89 | NG | 96 |
| 24) | Vinyl acetate | 16.04 | 368 | 386372 | 1033.72 | NG | 97 |
| 25) | Dichlorobromomethane | 16.39 | 377 | 539072 | 1045.63 | NG | 96 |
| 26) | 1,2-Dichloropropane | 17.78 | 413 | 236212 | 923.39 | NG | 96 |
| 27) | cis-1,3-Dichloropropylene | 18.06 | 420 | 359779 | 992.91 | NG | 93 |
| 28) | Trichloroethylene | 18.64 | 435 | 265049 | 979.90 | NG | 84 |
| 29) | Chlorodibromomethane | 19.26 | 451 | 380714 | 1009.49 | NG | 99 |
| 30) | bis(Chloromethyl)ether | 19.26 | 451 | 161033 | 1048.79 | NG | 100 |
| 31) | Benzene | 19.18 | 449 | 880604 | 1032.87 | NG | 94 |
| 32) | 1,1,2-Trichloroethane | 19.37 | 454 | 242009 | 1003.71 | NG | 81 |
| 33) | trans-1,3-Dichloropropylene | 19.37 | 454 | 492760 | 1043.22 | NG | 89 |
| 34) | 2-Chloroethylvinyl ether | 20.46 | 482 | 172878 | 948.11 | NG | 100 |
| 35) | Bromoform | 21.97 | 521 | 299168 | 958.02 | NG | 99 |
| 36) | *Chlorobenzene-d5 | 26.69 | 641 | 214016 | 250.00 | NG | 80 |
| 37) | Methyl-iso-butyl ketone | 22.48 | 534 | 247124 | 843.78 | NG | 96 |
| 38) | 2-Hexanone | 24.03 | 574 | 234177 | 827.55 | NG | 95 |
| 38) | 2-Hexanone | 24.92 | 597 | 6261 | 22.13 | NG | 44 |
| 39) | 1,1,2,2-Tetrachloroethane | 24.30 | 581 | 391687 | 880.94 | NG | 96 |
| 40) | Tetrachloroethylene | 24.34 | 582 | 281542 | 953.38 | NG | 97 |

QUANT REPORT

Page 2

Operator ID: KB6656
 Output File: ^C0975::AQ
 Data File: >C0975::U4
 Name: VOA,901001,C
 Misc: QC70357US,QU70357,L,5, 20UL

Quant Rev: 7 Quant Time: 901003 14:45
 Injected at: 901003 13:05
 Dilution Factor: 1.00000

ID File: IC1171::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901002 16:53

| | Compound | R.T. | Scan# | Area | Conc | Units | q |
|-----|-----------------------------|-------|-------|---------|---------|-------|----|
| 41) | Toluene-D8 (SURR) | 25.47 | 611 | 313012 | 249.74 | NG | 91 |
| 42) | Toluene | 25.66 | 616 | 760493 | 1070.55 | NG | 91 |
| 43) | Chlorobenzene | 26.81 | 644 | 993619 | 999.48 | NG | 94 |
| 44) | Ethylbenzene | 28.90 | 698 | 631612 | 1055.86 | NG | 77 |
| 45) | p-Bromofluorobenzene (SURR) | 31.66 | 769 | 238765 | 269.70 | NG | 80 |
| 46) | Styrene | 33.29 | 811 | 1506277 | 1068.82 | NG | 93 |
| 47) | m-Xylene | 33.64 | 820 | 842266 | 1074.86 | NG | 95 |
| 48) | o+p-Xylenes | 34.73 | 848 | 1756527 | 2170.94 | NG | 88 |

* Compound is ISTD

VOLATILE INTERNAL ^{8A} STANDARD AREA SUMMARY

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Lab File ID (Standard): >C0914

Data Analyzed: 09/26/90

Instrument ID: GC/MS C

Time Analyzed: 1352

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) JACK

| | IS1(BCM) | RT | IS2(DFB) | RT | IS3(CBZ) | RT |
|----------------|----------|-------|----------|-------|----------|-------|
| | AREA # | | AREA # | | AREA # | |
| 12 HOUR STD | 70232 | 11.09 | 342312 | 21.84 | 318334 | 26.78 |
| UPPER LIMIT | 140464 | | 684624 | | 636668 | |
| LOWER LIMIT | 35116 | | 171156 | | 159167 | |
| EPA SAMPLE NO. | | | | | | |
| 01 UBLK01 | 47287 | 11.08 | 219376 | 21.83 | 199092 | 26.79 |
| 02 A4753 | 46998 | 11.19 | 224774 | 21.94 | 207367 | 26.92 |
| 03 A4755 | 51014 | 11.20 | 234507 | 21.98 | 216956 | 26.92 |
| 04 A4763 | 45139 | 11.24 | 219122 | 21.95 | 207394 | 26.88 |
| 05 A4756 | 44376 | 11.26 | 215254 | 21.96 | 204163 | 26.90 |
| 06 A4758 | 41198 | 11.30 | 192371 | 21.96 | 182527 | 26.90 |
| 07 | | | | | | |
| 08 | | | | | | |
| 09 | | | | | | |
| 10 | | | | | | |
| 11 | | | | | | |
| 12 | | | | | | |
| 13 | | | | | | |
| 14 | | | | | | |
| 15 | | | | | | |
| 16 | | | | | | |
| 17 | | | | | | |
| 18 | | | | | | |
| 19 | | | | | | |
| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene-d5

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

* Column used to flag internal standard area values with an asterisk.

page 1 of 1

FORM VIII UOA

1/87 Rev.

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID (Standard): >C0927 Data Analyzed: 09/27/90
 Instrument ID: GC/MS C Time Analyzed: 1656
 Matrix: (soil/water) WATER Level: (low/med) LOW Column: (pack/cap) PALK

| | IS1(BCM) | | IS2(DFB) | | IS3(CBZ) | |
|-------------------|----------|-------|----------|-------|----------|-------|
| | AREA # | RT | AREA # | RT | AREA # | RT |
| 12 HOUR STD | 40844 | 11.08 | 179047 | 21.83 | 171595 | 26.77 |
| UPPER LIMIT | 81688 | | 358094 | | 343190 | |
| LOWER LIMIT | 20422 | | 89523 | | 85798 | |
| EPA SAMPLE NO. | | | | | | |
| 01 UBLK2 | 36606 | 11.11 | 178010 | 21.86 | 167977 | 26.79 |
| 02 A4754 | 42061 | 11.10 | 196522 | 21.85 | 185259 | 26.82 |
| (CA) 03 UGA A4757 | 39142 | 11.10 | 183849 | 21.84 | 173277 | 26.78 |
| 10-16-90 04 A4759 | 42889 | 11.11 | 192985 | 21.85 | 186569 | 26.79 |
| 05 | | | | | | |
| 06 | | | | | | |
| 07 | | | | | | |
| 08 | | | | | | |
| 09 | | | | | | |
| 10 | | | | | | |
| 11 | | | | | | |
| 12 | | | | | | |
| 13 | | | | | | |
| 14 | | | | | | |
| 15 | | | | | | |
| 16 | | | | | | |
| 17 | | | | | | |
| 18 | | | | | | |
| 19 | | | | | | |
| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene-d5
 UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

* Column used to flag internal standard area values with an asterisk.

page + of +

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: ETCNJ Contract:
 Lab Code: F2 Case No.: SAS No.: SDG No.:
 Lab File ID (Standard): >C0972 Data Analyzed: 10/03/90
 Instrument ID: GC/MS C Time Analyzed: 1001
 Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap) PAGE

| | IS1(BCM) | RT | IS2(DFB) | RT | IS3(CBZ) | RT |
|----------------|----------|-------|----------|-------|----------|-------|
| | AREA # | | AREA # | | AREA # | |
| 12 HOUR STD | 33439 | 11.00 | 154748 | 21.78 | 176233 | 26.72 |
| UPPER LIMIT | 66878 | | 309496 | | 352466 | |
| LOWER LIMIT | 16719 | | 77374 | | 88117 | |
| EPA SAMPLE NO. | | | | | | |
| 01 UBLK5 | 32787 | 10.99 | 159758 | 21.73 | 189232 | 26.68 |
| 02 A4755MS | 31333 | 11.02 | 158395 | 21.80 | 191425 | 26.75 |
| 03 A4755MSD | 33706 | 10.96 | 166042 | 21.74 | 193787 | 26.68 |
| 04 | | | | | | |
| 05 | | | | | | |
| 06 | | | | | | |
| 07 | | | | | | |
| 08 | | | | | | |
| 09 | | | | | | |
| 10 | | | | | | |
| 11 | | | | | | |
| 12 | | | | | | |
| 13 | | | | | | |
| 14 | | | | | | |
| 15 | | | | | | |
| 16 | | | | | | |
| 17 | | | | | | |
| 18 | | | | | | |
| 19 | | | | | | |
| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene-d5
 UPPER LIMIT = + 100% of internal standard area.
 LOWER LIMIT = - 50% of internal standard area.

* Column used to flag internal standard area values with an asterisk.

page 1 of 1

ETC

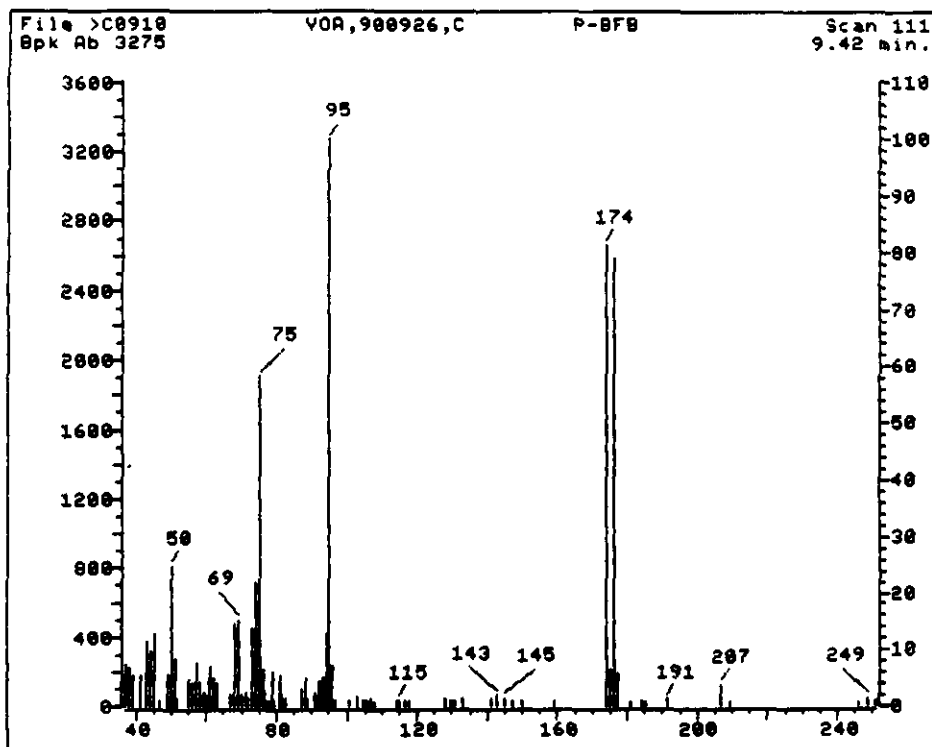
RAW QC DATA

MS Data File: >C0910::U4

Name: VOA,900926,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/26/90 8:54



MS Data File: >C0910::U4

Name: UOA,900926,C Operator: JAB781 Date/Time: 9/26/90 8:54
Misc: P-BFB

>C0910 UOA,900926,C P-BFB
111 NRM

File: >C0910 Scan #: 111 Retn. time: 9.42

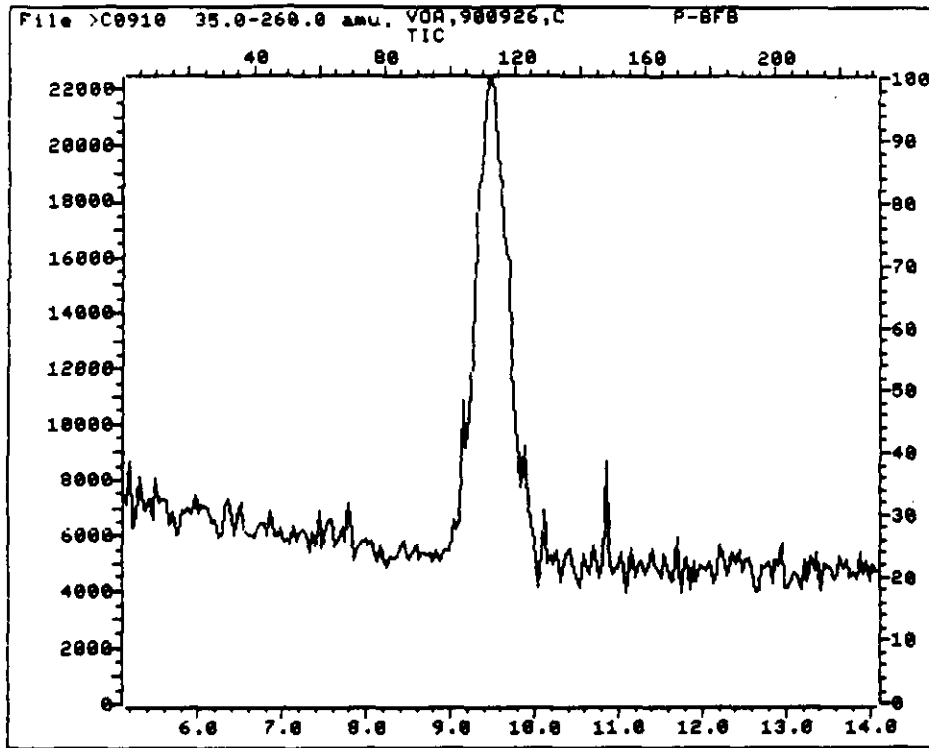
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|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 37.00 | 7.481 | 61.00 | 6.931 | 79.70 | 1.008 | 104.70 | .977 | 145.15 | 1.527 |
| 38.00 | 6.595 | 62.00 | 5.191 | 80.90 | 5.435 | 105.30 | .794 | 147.05 | 1.130 |
| 39.00 | 5.374 | 63.00 | 4.031 | 81.90 | 2.137 | 105.60 | .977 | 149.95 | 1.008 |
| 41.10 | 5.282 | 67.10 | 2.107 | 82.50 | 1.038 | 105.80 | .916 | 158.95 | .947 |
| 43.10 | 11.328 | 68.00 | 14.443 | 83.00 | 1.435 | 106.90 | 1.374 | 173.95 | 81.405 |
| 44.00 | 9.802 | 69.00 | 15.298 | 87.00 | 3.084 | 108.00 | .733 | 175.05 | 6.504 |
| 45.00 | 12.855 | 70.00 | 2.076 | 88.00 | 5.069 | 114.60 | .885 | 175.95 | 78.870 |
| 46.90 | 1.069 | 70.30 | 1.893 | 89.00 | 1.221 | 114.90 | 1.130 | 176.95 | 5.863 |
| 49.00 | 5.802 | 71.10 | 2.351 | 91.00 | 2.382 | 116.80 | 1.130 | 180.95 | .702 |
| 50.00 | 24.947 | 72.00 | 1.221 | 92.00 | 4.275 | 117.80 | 1.008 | 184.05 | 1.069 |
| 51.00 | 8.519 | 73.00 | 13.802 | 93.00 | 5.130 | 128.05 | 1.435 | 185.55 | .824 |
| 51.90 | 1.405 | 74.00 | 21.893 | 94.00 | 12.947 | 129.85 | 1.069 | 191.15 | 1.221 |
| 55.10 | 4.733 | 75.00 | 58.290 | 95.00 | 100.000 | 131.05 | .947 | 206.95 | 3.878 |
| 56.00 | 4.061 | 76.00 | 6.321 | 96.00 | 7.023 | 132.95 | 1.496 | 209.15 | .641 |
| 57.00 | 7.847 | 77.00 | 2.321 | 97.00 | 1.160 | 140.95 | 1.221 | 245.75 | .794 |
| 58.00 | 4.305 | 78.10 | 1.191 | 101.00 | 1.160 | 141.15 | 1.252 | 248.95 | 1.405 |
| 59.10 | 2.351 | 78.90 | 5.924 | 103.00 | 1.740 | 142.85 | 1.985 | 250.95 | .977 |
| 60.00 | 2.168 | | | | | | | | |

MS Data File: >C0910::U4

Name: UDA,900926,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/26/90 8:54

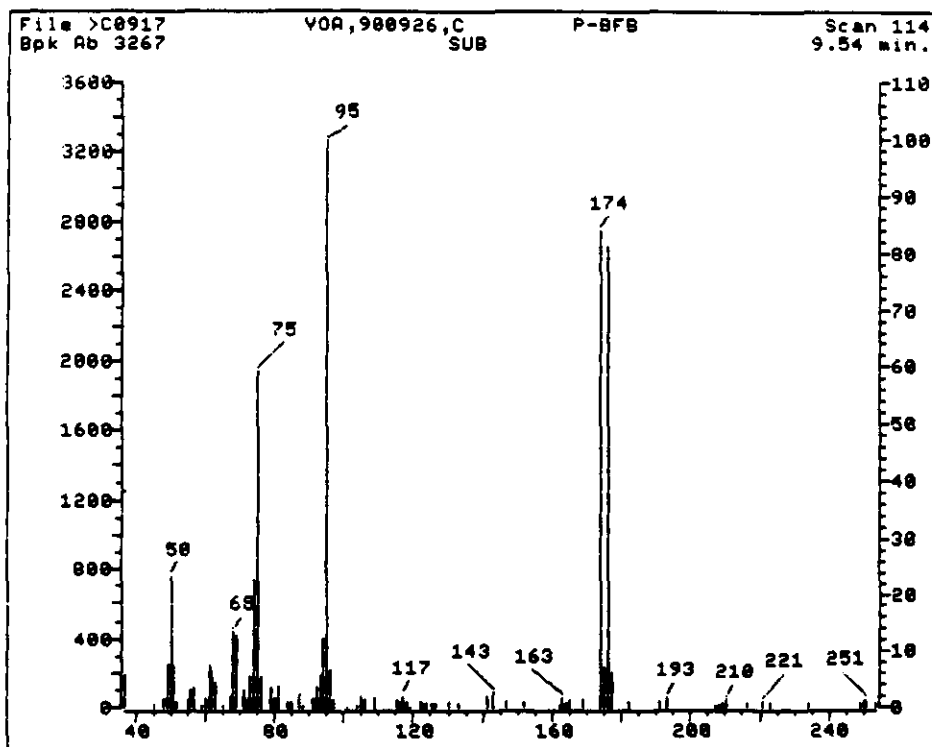


MS Data File: >C0917::U4

Name: VOA,900926,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/26/90 16:25



MS Data File: >C0917::U4

Name: UOA,900926,C Operator: JA8781 Date/Time: 9/26/90 16:25
Misc: P-BFB

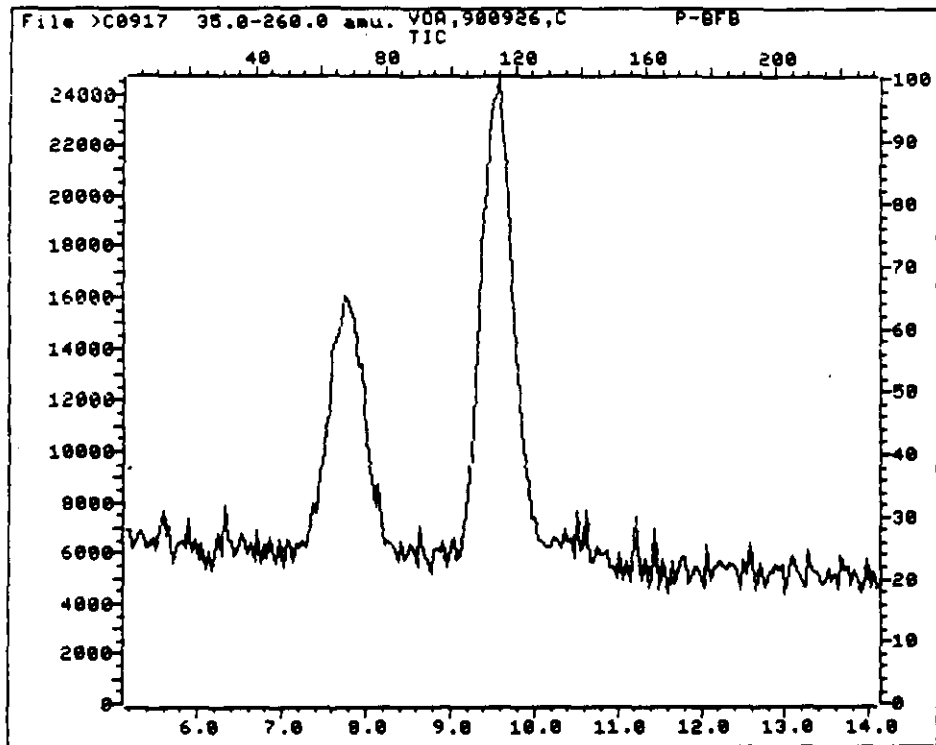
C0917 UOA,900926,C P-BFB
114 SUB NRM

File: >C0917 Scan #: 114 Retn. time: 9.54

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|-------|--------|-------|--------|--------|---------|--------|--------|--------|-------|
| 37.00 | 5.908 | 69.00 | 12.856 | 95.00 | 100.000 | 123.90 | .765 | 177.05 | 6.030 |
| 45.10 | .857 | 71.10 | 3.122 | 96.00 | 6.856 | 125.20 | .673 | 182.05 | .979 |
| 48.00 | 1.592 | 72.00 | 1.622 | 97.30 | 1.408 | 126.60 | .857 | 190.95 | .949 |
| 49.00 | 7.836 | 73.00 | 5.510 | 101.10 | .184 | 130.05 | 1.194 | 192.95 | 1.561 |
| 50.00 | 23.263 | 74.00 | 22.467 | 103.90 | .275 | 132.95 | .796 | 207.05 | .429 |
| 51.00 | 7.530 | 75.00 | 59.443 | 105.10 | 1.928 | 141.05 | 2.173 | 208.05 | .490 |
| 51.80 | 1.071 | 76.00 | 5.448 | 106.00 | 1.439 | 142.95 | 2.694 | 209.05 | .582 |
| 55.10 | 1.377 | 79.00 | 3.520 | 109.00 | 1.714 | 147.05 | 1.133 | 210.05 | 1.377 |
| 56.00 | 3.183 | 79.90 | 1.653 | 115.10 | 1.500 | 152.05 | .888 | 216.15 | .735 |
| 57.00 | 3.275 | 80.90 | 3.826 | 115.80 | 1.163 | 162.15 | .398 | 221.05 | .888 |
| 59.00 | .367 | 84.10 | .918 | 116.20 | .796 | 162.95 | 1.622 | 222.85 | .735 |
| 59.90 | 1.745 | 85.10 | .888 | 117.00 | 2.173 | 163.85 | .765 | 233.85 | .765 |
| 61.00 | 7.254 | 87.00 | 2.418 | 118.00 | 1.163 | 164.95 | 1.408 | 234.05 | .826 |
| 62.00 | 6.367 | 88.10 | .429 | 119.00 | .122 | 169.15 | 1.224 | 249.05 | .551 |
| 63.00 | 4.561 | 91.10 | 1.622 | 121.90 | .888 | 173.95 | 84.022 | 249.85 | 1.071 |
| 65.00 | .398 | 92.00 | 3.887 | 122.40 | .643 | 175.05 | 7.040 | 250.85 | 1.224 |
| 67.10 | 2.204 | 93.00 | 5.877 | 123.00 | .367 | 175.95 | 81.267 | 253.45 | .857 |
| 68.00 | 13.315 | 94.00 | 12.489 | | | | | | |

40243 cont.

MS Data File: >C0917::U4
Name: VOA,900926,C Operator: JA8781 Date/Time: 9/26/90 16:25
Misc: P-BFB

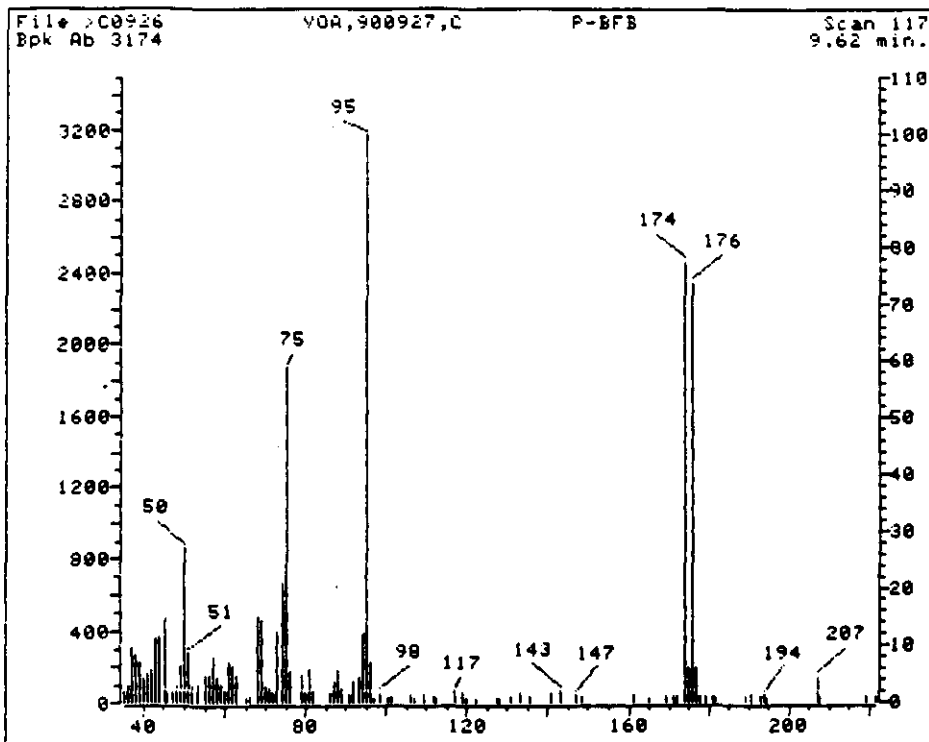


MS Data File: >C0926::U8

Name: UOA,900927,C
Misc: P-BFB

Operator: JAB781

Date/Time: 9/27/90 8:58



MS Data File: >C0926::U8

Name: UOA,900927,C Operator: JA8781 Date/Time: 9/27/90 8:58
Misc: P-BFB

>C0926 UOA,900927,C P-BFB
117 NRM

File: >C0926 Scan #: 117 Retn. time: 9.62

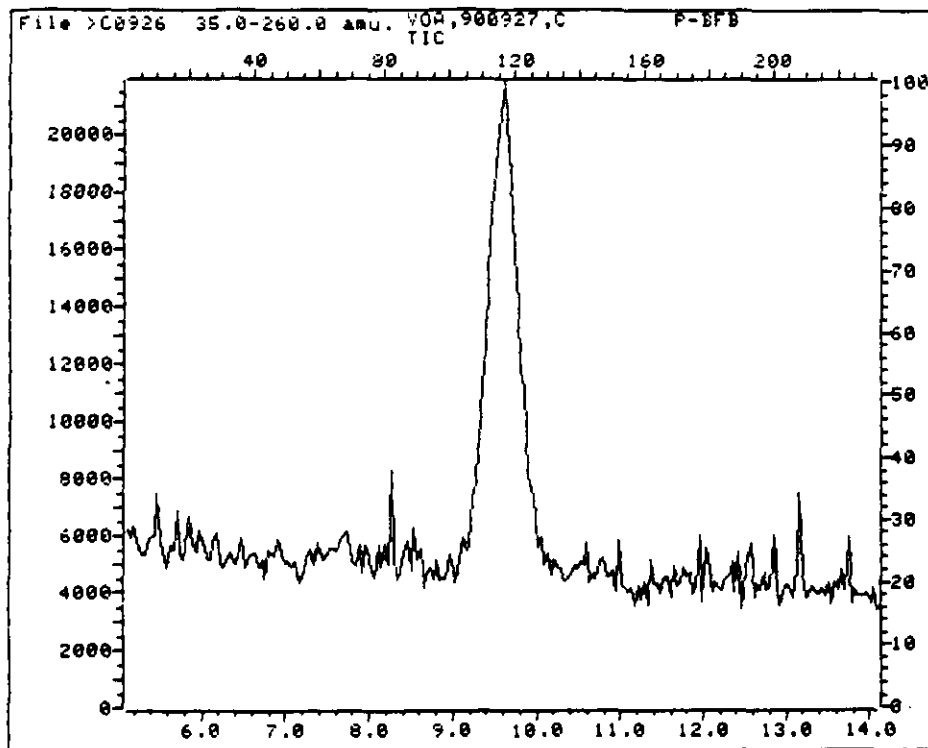
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|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 35.30 | 2.048 | 56.00 | 4.820 | 78.90 | 4.789 | 109.20 | 1.229 | 171.95 | .977 |
| 36.00 | 3.214 | 57.00 | 8.255 | 80.10 | 1.701 | 111.80 | 1.040 | 172.25 | .914 |
| 37.00 | 9.704 | 58.00 | 4.505 | 80.90 | 5.577 | 112.40 | .630 | 173.95 | 77.347 |
| 38.00 | 8.349 | 59.00 | 2.962 | 81.90 | 2.016 | 117.00 | 1.953 | 175.05 | 5.986 |
| 39.10 | 7.089 | 60.00 | 2.016 | 85.90 | 1.638 | 119.00 | 1.796 | 175.95 | 73.661 |
| 39.90 | 4.348 | 61.00 | 7.057 | 86.90 | 3.749 | 119.80 | .756 | 176.95 | 5.923 |
| 41.10 | 5.230 | 62.00 | 6.364 | 88.00 | 5.640 | 122.30 | .819 | 177.95 | 1.197 |
| 42.00 | 5.577 | 63.00 | 4.789 | 89.10 | 2.331 | 127.25 | .819 | 179.45 | .945 |
| 43.00 | 11.374 | 65.10 | .851 | 90.90 | 1.481 | 128.05 | .756 | 180.95 | 1.134 |
| 44.00 | 11.815 | 66.00 | .882 | 91.90 | 3.718 | 130.95 | .977 | 181.75 | .788 |
| 45.10 | 14.808 | 66.30 | .851 | 93.00 | 4.505 | 132.95 | 1.670 | 189.05 | .882 |
| 45.90 | 1.701 | 68.00 | 15.028 | 94.00 | 12.130 | 135.45 | 1.040 | 190.55 | 1.071 |
| 47.10 | 2.048 | 69.00 | 14.556 | 95.00 | 100.000 | 140.95 | 1.638 | 190.85 | 1.260 |
| 47.90 | 2.615 | 70.00 | 2.804 | 96.00 | 7.120 | 142.95 | 1.953 | 193.05 | 1.008 |
| 49.00 | 6.301 | 71.10 | 2.268 | 97.10 | .851 | 146.85 | 1.292 | 193.85 | 1.323 |
| 50.00 | 27.221 | 72.00 | 1.575 | 98.30 | 1.512 | 148.25 | .914 | 194.55 | .756 |
| 51.00 | 8.727 | 73.00 | 12.508 | 100.20 | .662 | 161.25 | 1.292 | 207.05 | 4.474 |
| 52.00 | 2.615 | 74.00 | 20.983 | 101.20 | .977 | 164.95 | .693 | 207.85 | 1.103 |
| 53.10 | 3.088 | 75.00 | 59.074 | 106.00 | 1.260 | 169.45 | .945 | 219.05 | .977 |
| 55.00 | 4.789 | 76.00 | 5.325 | 106.90 | .788 | 171.15 | .977 | 221.55 | 1.134 |

MS Data File: >C0926::U8

Name: UOA,900927,C
Misc: P-BFB

Operator: JA8781

Date/Time: 9/27/90 8:58

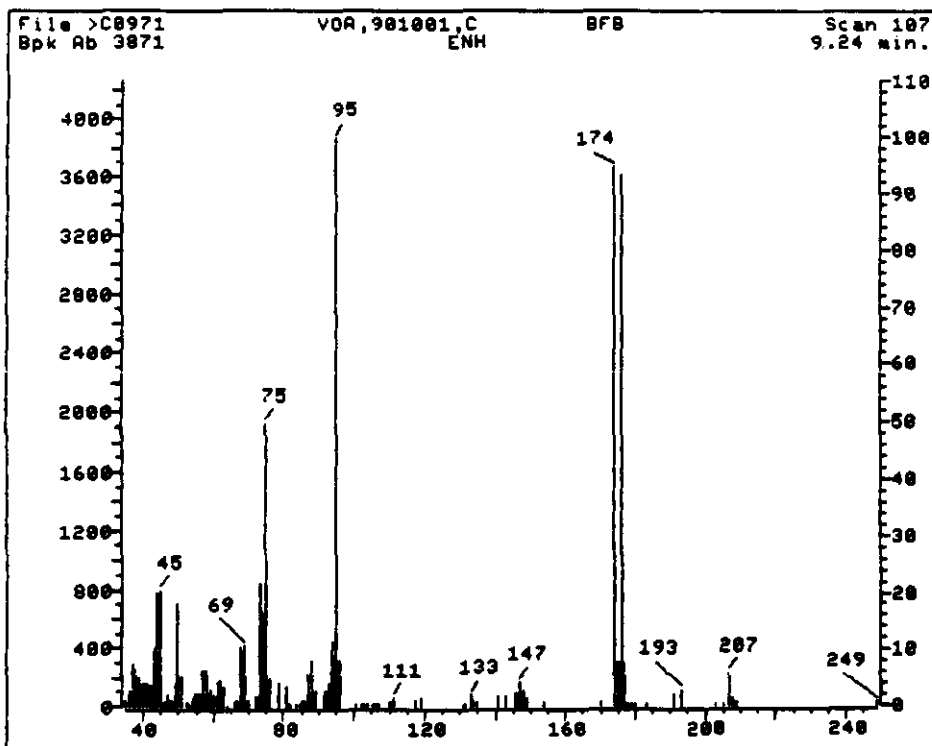


MS Data File: >C0971::U5

Name: VDA,901001,C
Misc: BFB

Operator: KB6656

Date/Time: 10/03/90 9:30



MS Data File: >C0971::U5

Name: UOA,901001,C
Misc: BFB

Operator: KB6656

Date/Time: 10/03/90 9:30

C0971 UOA,901001,C BFB
107 NRM ENH

File: >C0971 Scan #: 107 Retn. time: 9.24

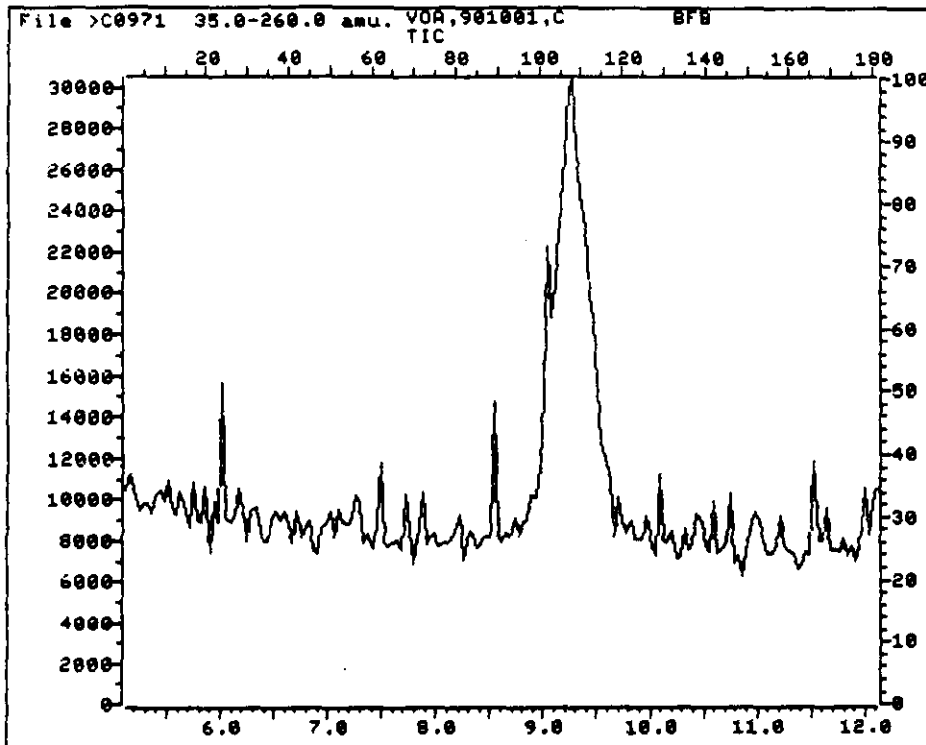
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|-------|--------|-------|--------|--------|---------|--------|-------|--------|--------|
| 35.10 | 1.121 | 54.90 | 2.423 | 78.80 | 4.128 | 105.60 | .377 | 148.05 | 2.640 |
| 35.90 | 2.743 | 55.90 | 2.392 | 80.90 | 3.250 | 105.80 | .356 | 149.05 | 1.271 |
| 37.00 | 7.316 | 57.00 | 6.494 | 81.50 | .232 | 106.50 | .269 | 153.85 | .661 |
| 38.00 | 6.396 | 57.90 | 6.256 | 82.00 | .393 | 107.10 | .382 | 170.15 | .568 |
| 39.00 | 5.228 | 59.10 | 3.079 | 83.50 | .258 | 110.10 | .599 | 173.95 | 94.932 |
| 39.90 | 4.133 | 59.90 | 2.191 | 84.50 | .284 | 110.40 | .372 | 174.95 | 7.889 |
| 41.00 | 4.112 | 61.00 | 4.422 | 85.00 | .951 | 111.10 | 1.503 | 175.95 | 93.521 |
| 42.00 | 3.653 | 61.90 | 4.707 | 85.90 | 1.142 | 111.40 | .537 | 176.95 | 7.579 |
| 43.00 | 10.297 | 63.00 | 3.270 | 87.00 | 5.735 | 117.10 | .914 | 177.95 | .517 |
| 44.00 | 20.149 | 64.10 | .196 | 88.00 | 8.209 | 118.80 | 1.209 | 178.85 | .372 |
| 45.00 | 20.660 | 65.90 | .837 | 88.90 | 2.692 | 130.90 | .336 | 180.25 | .248 |
| 45.90 | .770 | 66.90 | .997 | 92.00 | 2.749 | 131.20 | .258 | 183.45 | .511 |
| 47.00 | 2.067 | 68.00 | 10.519 | 92.90 | 4.180 | 131.50 | .232 | 190.85 | 2.206 |
| 48.20 | 1.199 | 69.00 | 10.710 | 94.00 | 11.314 | 133.00 | 2.025 | 192.95 | 2.795 |
| 49.00 | 4.221 | 70.20 | .883 | 95.00 | 100.000 | 133.80 | .610 | 202.85 | .253 |
| 49.90 | 18.056 | 71.90 | 1.679 | 96.00 | 7.992 | 134.80 | .672 | 205.15 | .449 |
| 51.00 | 5.409 | 73.00 | 21.766 | 100.70 | .315 | 141.05 | 1.875 | 206.95 | 5.884 |
| 52.60 | .599 | 74.00 | 16.956 | 101.90 | .253 | 142.95 | 1.571 | 207.95 | 1.297 |
| 53.00 | .387 | 75.00 | 49.752 | 102.90 | .537 | 145.95 | 2.304 | 209.15 | .873 |
| 53.90 | 1.374 | 75.90 | 4.686 | 103.50 | .269 | 146.95 | 4.402 | 248.75 | .594 |

MS Data File: >C0971::U5

Name: VOA,901001,C
Misc: BFB

Operator: KB6656

Date/Time: 10/03/90 9:30



1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLK01

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: QC70357U

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0918

Level: (low/med) LOW

Date Received: ~~10/1/90~~ ^{09/26/90}

% Moisture: not dec.

Date Analyzed: 09/26/90 ¹⁰⁻¹⁶⁻⁹⁰

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | |
|------------|----------------------------|----------------------|------|
| | | (ug/L or ug/Kg) | UG/L |
| 74-87-3 | Chloromethane | 10 | 10 |
| 74-83-9 | Bromomethane | 10 | 10 |
| 75-01-4 | Vinyl Chloride | 10 | 10 |
| 75-00-3 | Chloroethane | 10 | 10 |
| 75-09-2 | Methylene Chloride | 5 | 10 |
| 67-64-1 | Acetone | 10 | 10 |
| 75-15-0 | Carbon Disulfide | 5 | 10 |
| 75-35-4 | 1,1-Dichloroethene | 5 | 10 |
| 75-34-3 | 1,1-Dichloroethane | 5 | 10 |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | 10 |
| 67-66-3 | Chloroform | 5 | 10 |
| 107-06-2 | 1,2-Dichloroethane | 5 | 10 |
| 78-93-3 | 2-Butanone | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | 10 |
| 56-23-5 | Carbon Tetrachloride | 5 | 10 |
| 108-05-4 | Vinyl Acetate | 10 | 10 |
| 75-27-4 | Bromodichloromethane | 5 | 10 |
| 78-87-5 | 1,2-Dichloropropane | 5 | 10 |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | 10 |
| 79-01-6 | Trichloroethene | 5 | 10 |
| 124-48-1 | Dibromochloromethane | 5 | 10 |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | 10 |
| 71-43-2 | Benzene | 5 | 10 |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | 10 |
| 75-25-2 | Bromoform | 5 | 10 |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 10 |
| 591-78-6 | 2-Hexanone | 10 | 10 |
| 127-18-4 | Tetrachloroethene | 5 | 10 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | 10 |
| 108-88-3 | Toluene | 5 | 10 |
| 108-90-7 | Chlorobenzene | 5 | 10 |
| 100-41-4 | Ethylbenzene | 5 | 10 |
| 100-42-5 | Styrene | 5 | 10 |
| 1330-20-7 | Xylene (total) | 5 | 10 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: ETCNS Contract: _____

VBLK01

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER

Lab Sample ID: GC70357

Sample wt/vol: 5 (g/mL) HL

Lab File ID: 7C0918

Level: (low/med) LOW

Date Received: 09/26/90

Moisture: not dec. _____

Date Analyzed: 09/26/90

Column: (pack/cap) PACK

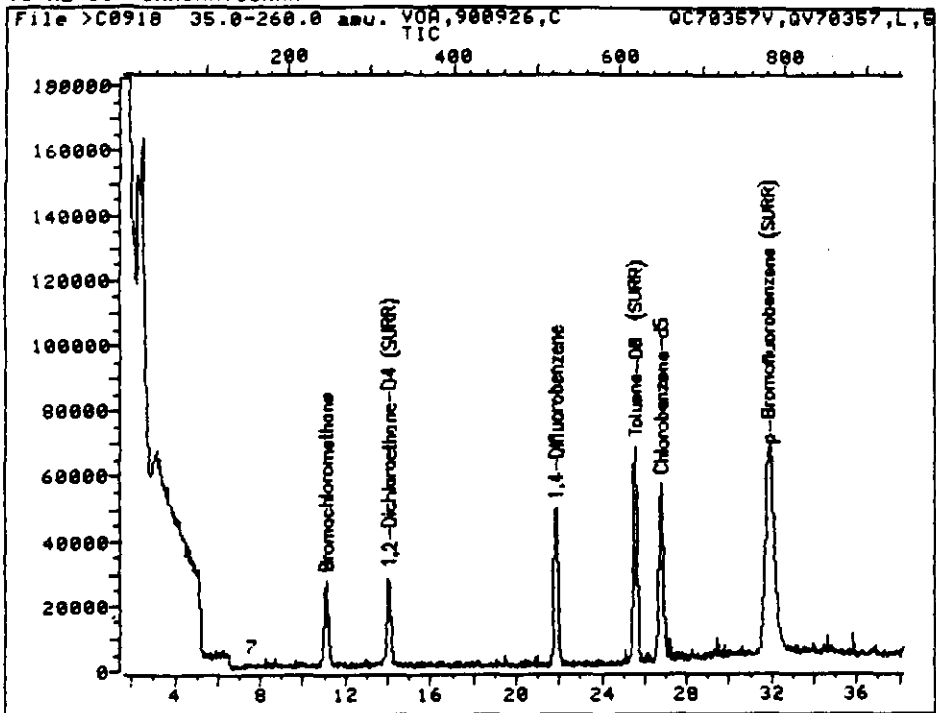
Dilution Factor: 1

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
| 2. | | | | |
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TOTAL ION CHROMATOGRAM



Data File: >C0918::U0 Quant Output File: ^C0918::AQ
Name: VOA,900926,C
Misc: QC70357U,QU70357,L,5,, BLANK

Id File: IC1171::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 900926 16:16

Operator ID: JA8781
Quant Time: 900926 18:05
Injected at: 900926 17:25

QUANT REPORT

Page 1

Operator ID: JA8781 Quant Rev: 7 Quant Time: 900926 18:05
 Output File: ^C0918::AQ Injected at: 900926 17:25
 Data File: >C0918::U0 Dilution Factor: 1.00000
 Name: UOA,900926,C
 Misc: QC70357U,QU70357,L,5,, BLANK

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 900926 16:16

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.08 | 245 | 47287 | 250.00 | NG | 97 |
| 7) Methylene chloride | 7.48 | 152 | 3530 | 11.20 | NG | 96 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.03 | 321 | 157251 | 238.22 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.83 | 522 | 219376 | 250.00 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.79 | 649 | 199092 | 250.00 | NG | 80 |
| 41) Toluene-D8 (SURR) | 25.55 | 618 | 295439 | 251.96 | NG | 93 |
| 45) p-Bromofluorobenzene (SURR) | 31.83 | 779 | 228780 | 264.05 | NG | 85 |

* Compound is ISTD

CA
 10-17-90

IA
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

UBLK2

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: QC7035702

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0929

Level: (low/med) LOW

Date Received: ~~10/1/90~~ ^{04/27/90} CR

% Moisture: not dec.

Date Analyzed: 09/27/90 ^{10/16/90}

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|---|
| 74-87-3 | Chloromethane | 10 | U |
| 74-83-9 | Bromomethane | 10 | U |
| 75-01-4 | Vinyl Chloride | 10 | U |
| 75-00-3 | Chloroethane | 10 | U |
| 75-09-2 | Methylene Chloride | 5 | U |
| 67-64-1 | Acetone | 10 | U |
| 75-15-0 | Carbon Disulfide | 5 | U |
| 75-35-4 | 1,1-Dichloroethene | 5 | U |
| 75-34-3 | 1,1-Dichloroethane | 5 | U |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | U |
| 67-66-3 | Chloroform | 5 | U |
| 107-06-2 | 1,2-Dichloroethane | 5 | U |
| 78-93-3 | Butanone | 10 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | U |
| 56-23-5 | Carbon Tetrachloride | 5 | U |
| 108-05-4 | Vinyl Acetate | 10 | U |
| 75-27-4 | Bromodichloromethane | 5 | U |
| 78-87-5 | 1,2-Dichloropropane | 5 | U |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | U |
| 79-01-6 | Trichloroethene | 5 | U |
| 124-48-1 | Dibromochloromethane | 5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | U |
| 71-43-2 | Benzene | 5 | U |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | U |
| 75-25-2 | Bromoform | 5 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | U |
| 591-78-6 | 2-Hexanone | 10 | U |
| 127-18-4 | Tetrachloroethene | 5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | U |
| 108-88-3 | Toluene | 5 | U |
| 108-90-7 | Chlorobenzene | 5 | U |
| 100-41-4 | Ethylbenzene | 5 | U |
| 100-42-5 | Styrene | 5 | U |
| 1330-20-7 | Xylene (total) | 5 | U |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLK2

Lab Name: ETCNJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: GC76227

Sample wt/vol: 5 (g/mL) HL Lab File ID: 7C0929

Level: (low/med) LOW Date Received: 04/27/90

% Moisture: not dec. _____ Date Analyzed: 04/27/90

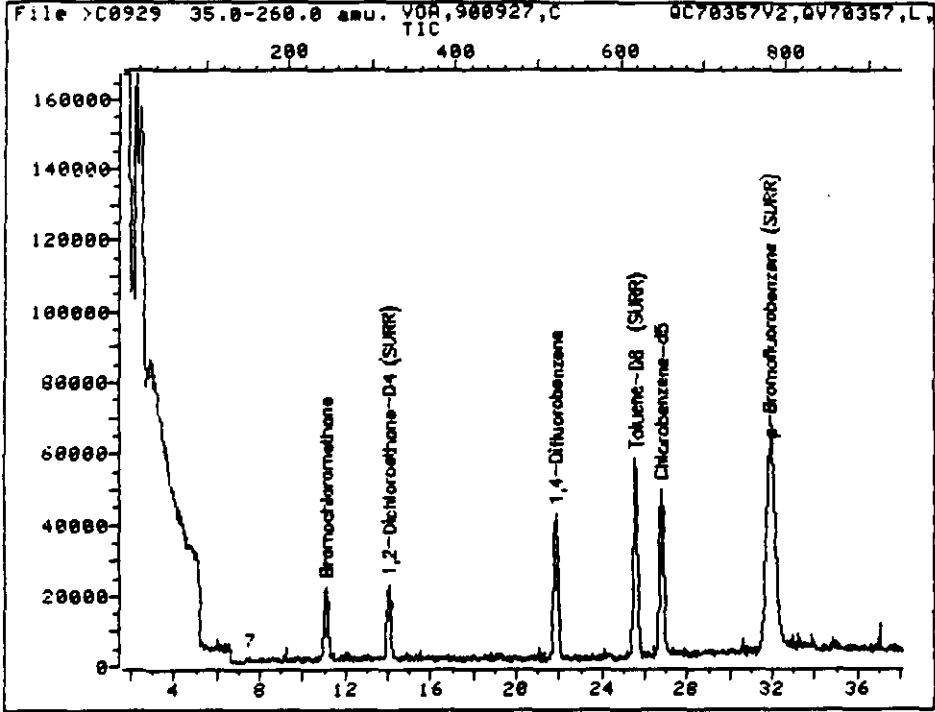
Column: (pack/cap) PACK Dilution Factor: 1

Number TICs found: 0

CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
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TOTAL ION CHROMATOGRAM



Data File: >C0929::U4

Quant Output File: ^C0929::AQ

Name: VOA,900927,C

Misc: QC70357V2,QU70357,L,5,,

BLANK

Id File: IC1171::US

Title: IFB, PP/VOA, TCL, XVOA13

Last Calibration: 900927 17:41

Operator ID: MGRMS

Quant Time: 900927 18:53

Injected at: 900927 18:14

QUANT REPORT

Page 1

Operator ID: MGRMS Quant Rev: 7 Quant Time: 900927 18:53
 Output File: ^C0929::AQ Injected at: 900927 18:14
 Data File: >C0929::U4 Dilution Factor: 1.00000
 Name: UOA,900927,C
 Misc: QC70357U2,QU70357,L,5,, BLANK

ID File: IC1171::US
 Title: IFB, PP/UOA, TCL, XVOA13
 Last Calibration: 900927 17:41

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 11.11 | 245 | 36606 | 250.00 | NG | 94 |
| 7) Methylene chloride | 7.47 | 151 | 3941 | 16.21 | NG | 82 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.06 | 321 | 128055 | 270.65 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.86 | 522 | 178010 | 250.00 | NG | 99 |
| 36) *Chlorobenzene-d5 | 26.79 | 649 | 167977 | 250.00 | NG | 82 |
| 41) Toluene-D8 (SURR) | 25.58 | 618 | 243972 | 248.68 | NG | 92 |
| 45) p-Bromofluorobenzene (SURR) | 31.88 | 780 | 195124 | 257.17 | NG | 85 |

* Compound is ISTD

CA
10-16-90.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

UBLK5

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: QC7035705

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0979

Level: (low/med) LOW

Date Received: 10/03/90

% Moisture: not dec.

Date Analyzed: 10/03/90

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | |
|------------|----------------------------|----------------------|------|
| | | (ug/L or ug/Kg) | UG/L |
| 74-87-3 | Chloromethane | 10 | 10 |
| 74-83-9 | Bromomethane | 10 | 10 |
| 75-01-4 | Vinyl Chloride | 10 | 10 |
| 75-00-3 | Chloroethane | 10 | 10 |
| 75-09-2 | Methylene Chloride | 5 | 10 |
| 67-64-1 | Acetone | 10 | 10 |
| 75-15-0 | Carbon Disulfide | 5 | 10 |
| 75-35-4 | 1,1-Dichloroethene | 5 | 10 |
| 75-34-3 | 1,1-Dichloroethane | 5 | 10 |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | 10 |
| 67-66-3 | Chloroform | 5 | 10 |
| 107-06-2 | 1,2-Dichloroethane | 5 | 10 |
| 78-93-3 | 2-Butanone | 10 | 10 |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | 10 |
| 56-23-5 | Carbon Tetrachloride | 5 | 10 |
| 108-05-4 | Vinyl Acetate | 10 | 10 |
| 75-27-4 | Bromodichloromethane | 5 | 10 |
| 78-87-5 | 1,2-Dichloropropane | 5 | 10 |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | 10 |
| 79-01-6 | Trichloroethene | 5 | 10 |
| 124-48-1 | Dibromochloromethane | 5 | 10 |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | 10 |
| 71-43-2 | Benzene | 5 | 10 |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | 10 |
| 75-25-2 | Bromoform | 5 | 10 |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | 10 |
| 591-78-6 | 2-Hexanone | 10 | 10 |
| 127-18-4 | Tetrachloroethene | 5 | 10 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | 10 |
| 108-88-3 | Toluene | 5 | 10 |
| 108-90-7 | Chlorobenzene | 5 | 10 |
| 100-41-4 | Ethylbenzene | 5 | 10 |
| 100-42-5 | Styrene | 5 | 10 |
| 1330-20-7 | Xylene (total) | 5 | 10 |

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBK5

Lab Name: ETC NJ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____

Matrix: (soil/water) WATER Lab Sample ID: 867627

Sample wt/vol: 5 (g/mL) HL Lab File ID: 700479

Level: (low/med) LOW Date Received: 10/13/90

% Moisture: not dec. _____ Date Analyzed: 10/13/90

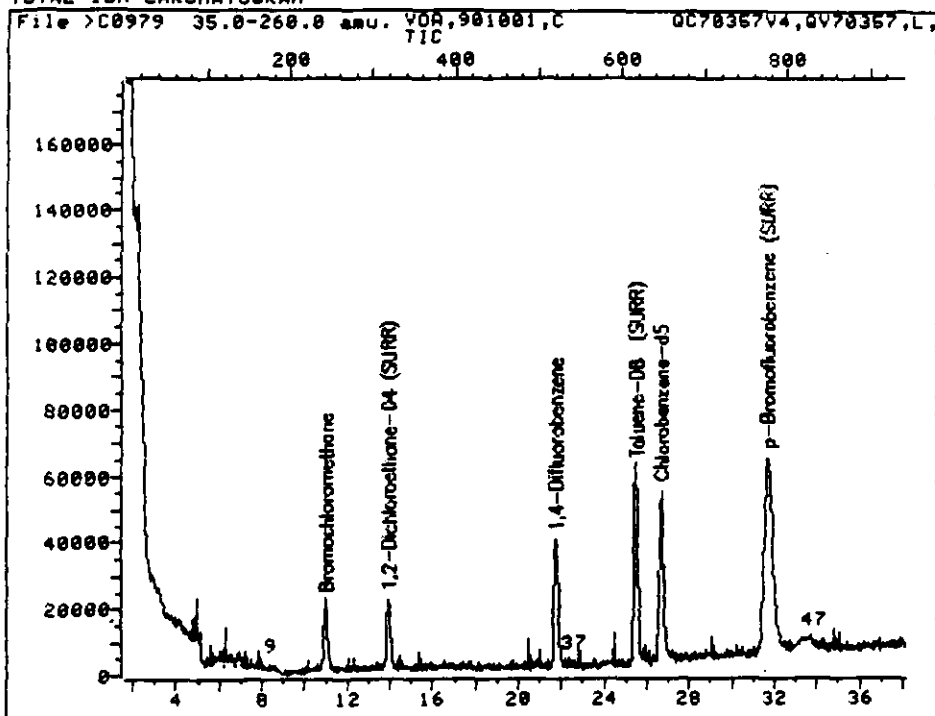
Column: (pack/cap) PACK Dilution Factor: 1

CONCENTRATION UNITS:
(ug/L or ug/kg) ug/L

Number TICs found: 0

| CAS NUMBER | COMPOUND NAME | RT | EST. CONC. | Q |
|------------|---------------|----|------------|---|
| 1. | | | | |
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TOTAL ION CHROMATOGRAM



Data File: >C0979::U4
 Name: VOA,901001,C
 Misc: QC7035704,QU70357,L,5,,

Quant Output File: ^C0979::AQ

Id File: IC1172::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901003 17:09

Operator ID: KB6656
 Quant Time: 901003 17:10
 Injected at: 901003 16:25

QUANT REPORT

Operator ID: KB6656
 Output File: ^C0979::AQ
 Data File: >C0979::U4
 Name: UOA,901001,C
 Misc: QC70357U4,QU70357,L,5,,

Quant Rev: 7 Quant Time: 901003 17:10
 Injected at: 901003 16:25
 Dilution Factor: 1.00000

ID File: IC1172::US
 Title: IFB, PP/UOA, TCL, XUOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|----|
| 1) *Bromochloromethane | 10.99 | 242 | 32787 | 250.00 | NG | 98 |
| 9) Acetone | 8.31 | 173 | 2343 | 24.50 | NG | 94 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.93 | 318 | 112633 | 242.43 | NG | 90 |
| 20) *1,4-Difluorobenzene | 21.73 | 519 | 159758 | 250.00 | NG | 95 |
| 36) *Chlorobenzene-d5 | 26.68 | 646 | 189232 | 250.00 | NG | 75 |
| 37) Methyl-iso-butyl ketone | 22.47 | 538 | 3765 | 14.68 | NG | 82 |
| 41) Toluene-D8 (SURR) | 25.49 | 616 | 272990 | 243.94 | NG | 86 |
| 45) p-Bromofluorobenzene (SURR) | 31.69 | 775 | 239059 | 280.27 | NG | 83 |
| 47) m-Xylene | 33.71 | 827 | 2049 | 2.74 | NG | 91 |

* Compound is ISTD

Q
 10-16-90.

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4755MS

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4755US

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0981

Level: (low/med) LOW

Date Received: ~~10/19/90~~ ^{04/20/90} *EP*
10-16-90

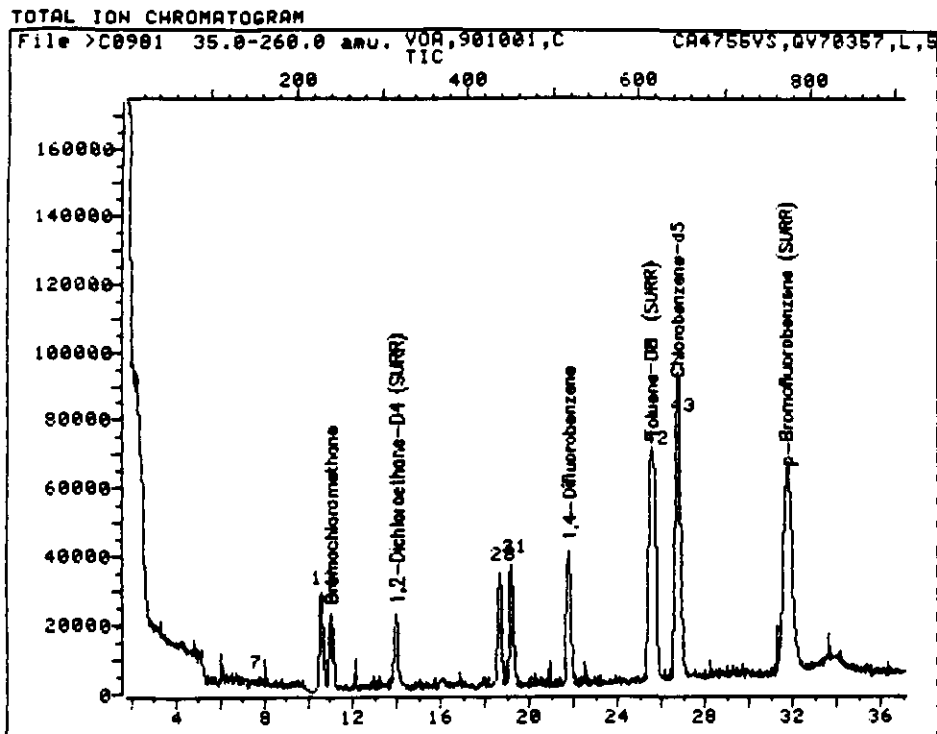
% Moisture: not dec.

Date Analyzed: 10/03/90

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|------------|----------------------------|--|---|
| 74-87-3 | Chloromethane | 10 | U |
| 74-83-9 | Bromomethane | 10 | U |
| 75-01-4 | Vinyl Chloride | 10 | U |
| 75-00-3 | Chloroethane | 10 | U |
| 75-09-2 | Methylene Chloride | 5 | U |
| 67-64-1 | Acetone | 10 | U |
| 75-15-0 | Carbon Disulfide | 5 | U |
| 75-35-4 | 1,1-Dichloroethene | 54 | U |
| 75-34-3 | 1,1-Dichloroethane | 5 | U |
| 540-59-0 | 1,2-Dichloroethene (total) | 5 | U |
| 67-66-3 | Chloroform | 5 | U |
| 107-06-2 | 1,2-Dichloroethane | 5 | U |
| 78-93-3 | 2-Butanone | 10 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | U |
| 56-23-5 | Carbon Tetrachloride | 5 | U |
| 108-05-4 | Vinyl Acetate | 10 | U |
| 75-27-4 | Bromodichloromethane | 5 | U |
| 78-87-5 | 1,2-Dichloropropane | 5 | U |
| 10061-01-5 | cis-1,3-Dichloropropene | 5 | U |
| 79-01-6 | Trichloroethene | 48 | U |
| 124-48-1 | Dibromochloromethane | 5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | U |
| 71-43-2 | Benzene | 42 | U |
| 10061-02-6 | trans-1,3-Dichloropropene | 5 | U |
| 75-25-2 | Bromoform | 5 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 10 | U |
| 591-78-6 | 2-Hexanone | 10 | U |
| 127-18-4 | Tetrachloroethene | 5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | U |
| 108-88-3 | Toluene | 49 | U |
| 108-90-7 | Chlorobenzene | 48 | U |
| 100-41-4 | Ethylbenzene | 5 | U |
| 100-42-5 | Styrene | 5 | U |
| 1330-20-7 | Xylene (total) | 5 | U |



Data File: >C0981::U4
Name: VOA,901001,C
Misc: CA4755VS,QU70357,L,5,,

Quant Output File: ^C0981::AQ

Id File: IC1172::US
Title: IFB, PP/VOA, TCL, XVDA13
Last Calibration: 901003 17:09

Operator ID: PT1575
Quant Time: 901003 20:28
Injected at: 901003 18:19

QUANT REPORT

Page 1

Operator ID: PT1575
 Output File: ^C0981::AQ
 Data File: >C0981::U4
 Name: VOA,901001,C
 Misc: CA4755U⁵QU70357,L,5,,

Quant Rev: 7 Quant Time: 901003 20:28
 Injected at: 901003 18:19
 Dilution Factor: 1.00000

ID File: IC1172::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|------------------|---------------|-----|
| 1) *Bromochloromethane | 11.02 | 240 | 31333 | 250.00 | NG | 98 |
| 7) Methylene chloride | 7.53 | 150 | 4026 | 24.52 | NG | 76 |
| 13) 1,1-Dichloroethylene | 10.56 | 228 | 57573 | 270.54 | NG | 89 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 14.01 | 317 | 120617 | 271.66 | NG | 99 |
| 20) *1,4-Difluorobenzene | 21.80 | 518 | 158395 | 250.00 | NG | 98 |
| 28) Trichloroethylene | 18.70 | 438 | 56640 | 240.09 | NG | 79 |
| 30) bis(Chloromethyl)ether | 19.21 | 451 | 10705 | 75.34 | NG | 100 |
| 31) Benzene | 19.21 | 451 | 153911 | 207.88 | NG | 87 |
| 36) *Chlorobenzene-d5 | 26.75 | 643 | 191425 | 250.00 | NG | 78 |
| 41) Toluene-D8 (SURR) | 25.53 | 614 | 297943 | 263.19 | NG | 90 |
| 42) Toluene | 25.68 | 618 | 161625 | 247.36 | NG | 93 |
| 43) Chlorobenzene | 26.86 | 646 | 211336 | 239.16 | NG | 95 |
| 45) p-Bromofluorobenzene (SURR) | 31.71 | 771 | 241906 | 280.36 | NG | 81 |

* Compound is ISTD

CA
 10-16-90

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

A4755MSD

Lab Name: ETCNJ

Contract:

Lab Code: F2

Case No.:

SAS No.:

SDG No.:

Matrix: (soil/water) WATER

Lab Sample ID: CA4755UR

Sample wt/vol: 5.0 (g/mL) ML

Lab File ID: >C0982

Level: (low/med) LOW

Date Received: ~~10/19/90~~ ^{09/20/90} (P)

% Moisture: not dec.

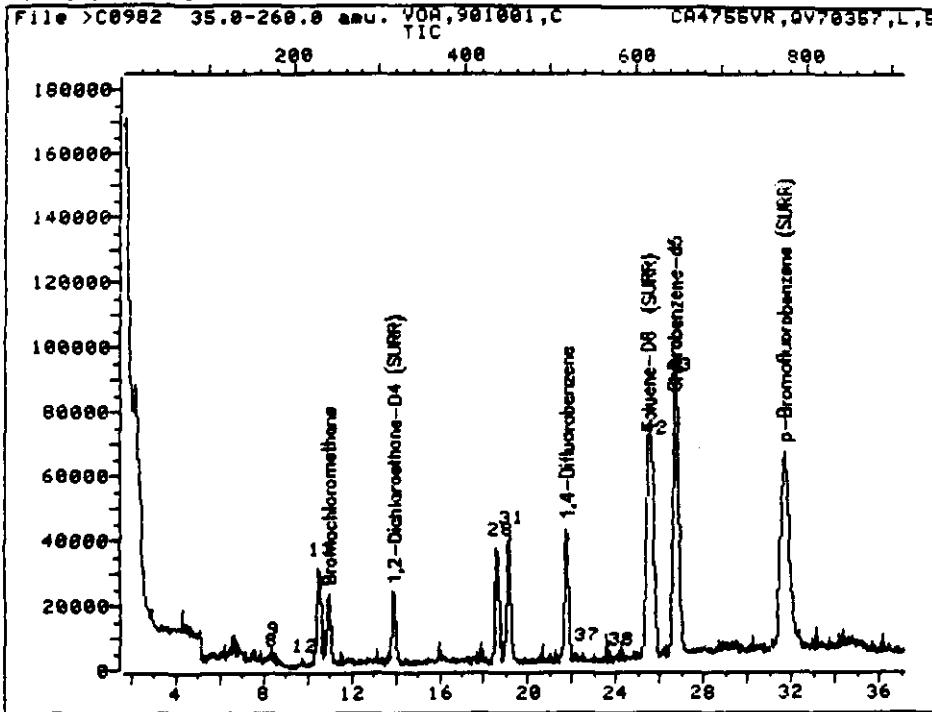
Date Analyzed: 10/03/90 ^{10/6/90}

Column: (pack/cap) PACK

Dilution Factor: 1

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) | UG/L | Q |
|------------|----------------------------|---|------|----|
| 74-87-3 | Chloromethane | | 10 | IU |
| 74-83-9 | Bromomethane | | 10 | IU |
| 75-01-4 | Vinyl Chloride | | 10 | IU |
| 75-00-3 | Chloroethane | | 10 | IU |
| 75-09-2 | Methylene Chloride | | 5 | IU |
| 67-64-1 | Acetone | | 10 | IU |
| 75-15-0 | Carbon Disulfide | | 5 | IU |
| 75-35-4 | 1,1-Dichloroethene | | 51 | I |
| 75-34-3 | 1,1-Dichloroethane | | 5 | IU |
| 540-59-0 | 1,2-Dichloroethene (total) | | 5 | IU |
| 67-66-3 | Chloroform | | 5 | IU |
| 107-06-2 | 1,2-Dichloroethane | | 5 | IU |
| 78-93-3 | 2-Butanone | | 10 | IU |
| 71-55-6 | 1,1,1-Trichloroethane | | 5 | IU |
| 56-23-5 | Carbon Tetrachloride | | 5 | IU |
| 108-05-4 | Vinyl Acetate | | 10 | IU |
| 75-27-4 | Bromodichloromethane | | 5 | IU |
| 78-87-5 | 1,2-Dichloropropane | | 5 | IU |
| 10061-01-5 | cis-1,3-Dichloropropene | | 5 | IU |
| 79-01-6 | Trichloroethene | | 50 | I |
| 124-48-1 | Dibromochloromethane | | 5 | IU |
| 79-00-5 | 1,1,2-Trichloroethane | | 5 | IU |
| 71-43-2 | Benzene | | 43 | I |
| 10061-02-6 | trans-1,3-Dichloropropene | | 5 | IU |
| 75-25-2 | Bromoform | | 5 | IU |
| 108-10-1 | 4-Methyl-2-Pentanone | | 10 | IU |
| 591-78-6 | 2-Hexanone | | 10 | IU |
| 127-18-4 | Tetrachloroethene | | 5 | IU |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5 | IU |
| 108-88-3 | Toluene | | 46 | I |
| 108-90-7 | Chlorobenzene | | 49 | I |
| 100-41-4 | Ethylbenzene | | 5 | IU |
| 100-42-5 | Styrene | | 5 | IU |
| 1330-20-7 | Xylene (total) | | 5 | IU |

TOTAL ION CHROMATOGRAM



Data File: >C0982::U4
Name: VOA,901001,C
Misc: CA4755VR,QV70357,L,5,,

Quant Output File: ^C0982::AQ

Id File: IC1172::US
Title: IFB, PP/VOA, TCL, XVOA13
Last Calibration: 901003 17:09

Operator ID: PT1575
Quant Time: 901003 20:38
Injected at: 901003 19:07

QUANT REPORT

Operator ID: PT1575
 Output File: ^C0982::AQ
 Data File: >C0982::U4
 Name: VOA,901001,C
 Misc: CA4755UR,QU70357,L,5,,

Quant Rev: 7 Quant Time: 901003 20:38
 Injected at: 901003 19:07
 Dilution Factor: 1.00000

ID File: IC1172::US
 Title: IFB, PP/VOA, TCL, XVOA13
 Last Calibration: 901003 17:09

| Compound | R.T. | Scan# | Area | Conc | Units | q |
|----------------------------------|-------|-------|--------|-------------------|---------------|-----|
| 1) *Bromochloromethane | 10.96 | 240 | 33706 | 250.00 | NG | 93 |
| 8) Acrolein | 8.28 | 171 | 3252 | 141.20 | NG | 77 |
| 9) Acetone | 8.36 | 173 | 3871 | 59.37 | NG | 77 |
| 12) Trichlorofluoromethane | 9.72 | 208 | 3383 | 6.19 | NG | 72 |
| 13) 1,1-Dichloroethylene | 10.49 | 228 | 58885 | 257.23 | NG | 84 |
| 18) 1,2-Dichloroethane-D4 (SURR) | 13.94 | 317 | 126631 | 265.13 | NG | 98 |
| 20) *1,4-Difluorobenzene | 21.74 | 518 | 166042 | 250.00 | NG | 96 |
| 28) Trichloroethylene | 18.60 | 437 | 61635 | 249.23 | NG | 81 |
| 30) bis(Chloromethyl)ether | 19.14 | 451 | 10810 | 78.57 | NG | 100 |
| 31) Benzene | 19.14 | 451 | 166534 | 214.57 | NG | 86 |
| 36) *Chlorobenzene-d5 | 26.68 | 644 | 193787 | 250.00 | NG | 73 |
| 37) Methyl-iso-butyl ketone | 22.47 | 537 | 3213 | 12.23 | NG | 97 |
| 38) 2-Hexanone | 24.07 | 578 | 4067 | 16.58 | NG | 95 |
| 41) Toluene-D8 (SURR) | 25.46 | 614 | 281173 | 245.35 | NG | 86 |
| 42) Toluene | 25.66 | 619 | 151333 | 228.78 | NG | 93 |
| 43) Chlorobenzene | 26.79 | 647 | 219237 | 245.07 | NG | 91 |
| 45) p-Bromofluorobenzene (SURR) | 31.68 | 773 | 243679 | 278.97 | NG | 81 |

* Compound is ISTD

CA
 10-16-90

ETC

CHAIN OF CUSTODY

Company: WASTE MANAGEMENT, INC.

Attn: DEE BENCICH

Facility/Site: _____

Phone: (708) 409 - 0700

Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 4015 Facility/Site Code (Optional Sample Point Descriptions)

Sample Point: W-26W8L 900919 0705
Source Code (from below) Your Sample Point ID (left justify) Start Date (YY/MM/DD) Start Time (2400 hr. clock) Elapsed Hours (composite)

Source Codes:
 Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| No | BOTTLE | | | ANALYSIS | SAMPLER | | LAB |
|----|--------|-----------------------------------|----------|-----------------|------------|--------------|--------------|
| | Type | Size | Preserv. | | FIL. (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | N | HCL | ✓ |
| 2 | VOA | 40 | HCL | VOLATILES | N | | ✓ |
| 1 | VS | 40 | HCL | VOA SCREEN | N | | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | TOC2 | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | N | | ✓ |
| 1 | MET | 1000 | HNO3 | METALS | N | | ✓ |
| 1 | CYAN | ¹⁰⁰⁰ 500 | NaOH | CY/T | N | | ✓ Rec'd IL |
| 1 | FL | 125 | NONE | FLUORIDE | N | | ✓ |
| | | | | | N | | |

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) DAN GREEN Date: 9-3-90 Time: 13:15
 Signature: [Signature] Seal #: 0175355 Intact: ✓

2. I have received these materials in good condition from the above person.
 Name: John V. Rudd Signature: [Signature]
 Date: 9/18/90 Time: 14:10 Remarks: OK

3. I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John V. Rudd Date: 9/19/90 Time: 0900
 Signature: [Signature] Seal #: 0175356 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/20/90 Time: 9:25
 SHUTTLE # 234 TEMP. °C 40 SEAL # 175356 COND. Intact 122

Company: WASTE MANAGEMENT, INC.

Attn.: DEE BENCICH

Facility/Site: _____

Phone: (708) 409 - 0700

Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 405 Facility/Site Code (Optional Sample Point Descriptions)

Sample Point: M-26WLA 900919 110110
Source Code (from below) Your Sample Point ID (left justify) Start Date (YY/MM/DD) Start Time (2400 hr clock) Elapsed Hours (composite)

Source Codes

Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|-------------|--------------|--------------|
| No | Type | Size | Preserv. | | Filt. (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | N | HCL | ✓ |
| 2 | VOA | 40 | HCL | VOLATILES | N | | ✓ |
| 1 | VS | 40 | HCL | VOA SCREEN | N | | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | TQC2 | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | N | | ✓ |
| 1 | MET | 1000 | HNO3 | METALS | N | | Rec'd empty |
| 1 | CYAN | 500 | NaOH | CY/T | N | | Not Rec'd |
| 1 | FL | 125 | NONE | FLUORIDE | N | | ✓ 9/80 |

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) DAN GREEN Date: 9-3-90 Time: 13:30
 Signature: [Signature] Seal #: 0175359 Intact: ✓

2. I have received these materials in good condition from the above person.
 Name: John Y. Rudd Signature: [Signature]
 Date: 9/18/90 Time: 1420 Remarks: OK

3. I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John Y. Rudd Date: 9/19/90 Time: 1130
 Signature: [Signature] Seal #: 1175354 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/20/90 Time: 1215
 SHUTTLE # 147 TEMP. °C 30 SEAL # 1175354 COND. Intact 26

Company: WASTE MANAGEMENT, INC.

Attn.: DEE BENCICH

Facility/Site: _____

Phone: (708) 409 - 0700

Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 405 _____ (Optional Sample Point Descriptions)

Sample Point: W-2645a _____ 900919 _____

Source Code (from below) Your Sample Point ID (left justify) Start Date (YY/MM/DD) Start Time (2400 hr. clock) Elapsed Hours (composite)

Source Codes

Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|-----------------|------------|--------------|--------------|
| No | Type | Size | Preserv. | | Fill (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | | REF | ✓ |
| 2 | VOA | 40 | HCL | VOLATILES | | | ✓✓✓✓ |
| 1 | VS | 40 | HCL | VOA SCREEN | | | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | | | ✓ |
| 1 | CONS | 125 | H2SO4 | TOC2 | | | ✓ |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | | | ✓ |
| 1 | MET | 1000 | HNO3 | METALS | | | ✓ |
| 1 | CYAN | 500 | NaOH | CY/T | | | ✓ (1) IL |
| 1 | FL | 125 | NONE | FLUORIDE | | | ✓ |

CHAIN OF CUSTODY CHRONICLE

- Shuttle Opened By: (print) DAN GROW Date: 9-3-90 Time: 13:40
 Signature: _____ Seal #: 0175361 Intact: ✓
- I have received these materials in good condition from the above person.
 Name: John Y. RUDD Signature: John Y. Rudd
 Date: 9/18/90 Time: 1430 Remarks: OK
- I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____
- Shuttle Sealed By: (print) John Y. RUDD Date: 9/19/90 Time: 1735
 Signature: John Y. Rudd Seal #: 0175360 Intact: OK

LAB USE ONLY Opened By: A. Stensler Date: 9/20/90 Time: 913
 SHUTTLE # 928 TEMP. °C 5° SEAL # 175360 COND. Intact 130



CHAIN OF CUSTODY FORM (CC1)

ORIGINAL

Date Sealed 9-5-90

By: KTC

Company: Water

Attn.: Don Green

Facility/Site: Turn of Sun

Phone: _____

Address: _____

SAMPLE IDENTIFICATION

Facility: 141015 (Optional Sample Point Descriptions)

Sample Point: W-12 FW SW 9009 9
Source Code (from below) Your Sample Point ID (left justify) Start Date (YYMM/DD) TC Start Time (2400 hr. clock) Elapsed Hours (composite)

Source Codes:

Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (L) Other (X)
Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (A) Specify _____

SHUTTLE CONTENTS

| No | BOTTLE | | | ANALYSIS | SAMPLER | | LAB |
|----|--------|------|----------|---------------------------|------------|--------------|--------------|
| | Type | Size | Preserv. | | FIL. (Y/N) | Observations | Observations |
| 1 | P | 250 | Ø | TDS, Bicarbonate, Sulfate | N | | full, clear |
| 1 | P | 250 | H2SO4 | NH3, TKN | N | | ↓ |
| | | | | | | | |
| | | | | | | | |
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CHAIN OF CUSTODY CHRONICLE

- Shuttle Opened By: (print) Don Green Date: 9-9-90 Time: 14:15
Signature: [Signature] Seal #: 01376-01377 Intact: ✓
- I have received these materials in good condition from the above person.
Name: John V. Rudd Signature: [Signature]
Date: 9/18/90 Time: 1430 Remarks: OK
- I have received these materials in good condition from the above person.
Name: _____ Signature: _____
Date: _____ Time: _____ Remarks: _____
- Shuttle Sealed By: (print) John V. Rudd Date: 9/19/90 Time: 1730
Signature: [Signature] Seal #: CA4755 Intact: OK

LAB USE ONLY Opened By: Kelly Schuck Date: 9-20-90 Time: 13:00
SHUTTLE # _____ TEMP. °C 40 SEAL # 4755 COND. Intact 132



FIELD PARAMETER FORM (CC2)

Form 0002
Sample Management
12/89

ETC JOB # CA-755
Sample Point W 216W.S1W
Source Code _____ Sample Point I.D. _____

FIELD PROCEDURES

PURGE DATE (YY MM DD) _____ START PURGE (2400 Hr Clock) _____ ELAPSED HRS _____ WATER VOL IN CASING (Gallons) _____ VOLUME PURGED (Gallons) _____

SAMPLING METHOD: PRIVATE WELL pump

Sampler Type A-Submersible Pump D-Dipper/Bottle E-Baller X-Other _____
 B-ISCO F-Scoop/Shovel (SPECIFY OTHER) _____
 C-Bladder Pump

Sampler Material A-Teflon C-PVC X-Other _____
 B-Metal D-Plastic (SPECIFY OTHER) _____

Tubing Material A-Teflon C-Polyethylene X-Other _____
 B-Tygon D-Silicon (SPECIFY OTHER) _____

Sample Compositing Y N N
Procedure/Proportions _____

FIELD MEASUREMENTS

Well Elevation (ft/msl) _____ Well Depth (ft) _____
Depth to Ground water (ft) _____ Sample Depth (non-well) (ft) _____
Groundwater Elevation (ft msl) _____

| | | | | | | | |
|-----|---------------------------------|-----|----------------------------|---------------|-------|-------|-------|
| 1st | <u>7.44</u> (STD) ph | 1st | <u>1691</u> spec. cond. | um/cm at 25°C | _____ | _____ | _____ |
| 2nd | <u>7.42</u> (STD) ph | 2nd | <u>1691</u> spec. cond. | um/cm at 25°C | _____ | _____ | _____ |
| 3rd | <u>7.41</u> (STD) ph | 3rd | <u>1690</u> spec. cond. | um/cm at 25°C | _____ | _____ | _____ |
| 4th | <u>7.41</u> (STD) ph | 4th | <u>1690</u> spec. cond. | um/cm at 25°C | _____ | _____ | _____ |
| | <u>12.2</u> (°C) Sample Temp | | _____ | Turbidity | _____ | _____ | _____ |

FIELD COMMENTS

Sample Appearance: CLEAR - NO ODDOR
Weather Conditions: PARTLY CLOUDY, 50°F
Other: _____

FILTERING: Use Chain of Custody (CC1) to indicate which bottles were filtered

Sampler: John Y. Rudd (Print) Employer: PELA

I certify that sampling procedures were in accordance with applicable EPA state and corporate protocols.

(Date) _____ (Signature) John Y. Rudd

CHAIN OF CUSTODY FORM (CC1) ORIGINAL Date Sealed 90/08/30 By: MM

Company: WASTE MANAGEMENT, INC. Attn.: DEE BENCICH
 Facility/Site: _____ Phone: (708) 409 - 0700
 Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 405 Facility/Site Code (Optional Sample Point Descriptions)
 Sample Point: W-2GWTH 900920 1738 15
Source Code (from below) Your Sample Point ID (left justify) Start Date (YYMM/DD) Start Time (2400 hr. clock) Elapsed Hours (composite)
 Source Codes:
 Well (W) Outfall (O) Bottom Sediment (B) Surface Impoundment (I) Leachate Collection Sys. (C) Other (X)
 Soil (S) River/Stream (R) Generation Point (G) Treatment Facility (T) Lake/Ocean (L) Specify _____

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|---------------------------|----------|-----------------|------------|--------------|-----------------|
| No | Type | Size | Preserv. | | Flt. (Y/N) | Observations | Observations |
| 1 | TB | 40 | GC/MS | TRIP BLANK | N | | ✓ |
| 2 | VOA | 40 | HCL | VOLATILES | N | | ✓ <u>AURALS</u> |
| 1 | VS | 40 | HCL | VOA SCREEN | N | | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | TOC2 | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | N | | ✓ |
| 1 | MET | 1000 | HNO3 | METALS | N | | ✓ |
| 1 | CYAN | <u>1000</u> <u>500</u> | NaOH | CY/T | N | | ✓ |
| 1 | FL | 125 | NONE | FLUORIDE | N | | ✓ |

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By (print) DAN GROSS Date: 9-3-90 Time: 14:00
 Signature: [Signature] Seal #: 0175399 Intact: NO

2. I have received these materials in good condition from the above person.
 Name: John V. Rudd Signature: [Signature]
 Date: 09/18/90 Time: 1400 Remarks: OK

3. I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John V. Rudd Date: 9/20/90 Time: 1900
 Signature: [Signature] Seal #: 0175399 Intact: OK

LAB USE ONLY Operated By: [Signature] Date: 9/21/90 Time: 1032
 SHUTTLE # 79 TEMP. °C 5° SEAL # 175398 COND. Intact

ETC JOB # CA4756
 Sample Point W 2 GW 1 H
Source Code Sample Point I.D.

FIELD PROCEDURES

PURGE DATE (YY MM DD) 9.0.09.20 START PURGE (2400 Hr Clock) 1738 ELAPSED HRS 1.5 WATER VOL. IN CASING (Gallons) _____ VOLUME PURGED (Gallons) _____

SAMPLING METHOD: PRIVATE well pump Plumbing

Sampler Type A-Submersible Pump D-Dipper/Bottle
 B-ISCO E-Bailer X-Other _____
 C-Bladder Pump F-Scoop/Shovel (SPECIFY OTHER)

Sampler Material A-Teflon C-PVC X-Other _____
 B-Metal D-Plastic (SPECIFY OTHER)

Tubing Material A-Teflon C-Polyethylene X-Other _____
 B-Tygon D-Silicon (SPECIFY OTHER)

Sample Compositing N Procedure/Proportions _____

FIELD MEASUREMENTS

Well Elevation (ft/msl) _____ Well Depth (ft) _____
 Depth to Ground water (ft) _____ Sample Depth (non-well) (ft) _____
 Groundwater Elevation (ft msl) _____

| | | | | | |
|-------------------------------------|-----------------------------------|----------------------|-------|-------|-------|
| 1st <u>7.017</u> (STD) <u>ph</u> | 1st <u>439</u> <u>spec. cond.</u> | <u>um/cm</u> at 25°C | _____ | _____ | _____ |
| 2nd <u>7.017</u> (STD) <u>ph</u> | 2nd <u>437</u> <u>spec. cond.</u> | <u>um/cm</u> at 25°C | _____ | _____ | _____ |
| 3rd <u>7.017</u> (STD) <u>ph</u> | 3rd <u>439</u> <u>spec. cond.</u> | <u>um/cm</u> at 25°C | _____ | _____ | _____ |
| 4th <u>7.018</u> (STD) <u>ph</u> | 4th <u>438</u> <u>spec. cond.</u> | <u>um/cm</u> at 25°C | _____ | _____ | _____ |
| <u>13.3</u> (°C) <u>Sample Temp</u> | _____ <u>Turbidity</u> | _____ <u>NTU</u> | _____ | _____ | _____ |

FIELD COMMENTS

Sample Appearance: CLEAR - NO ODR
 Weather Conditions: PC - 60° outside temp
 Other: Sample taken in Basement, approx temp 80°

FILTERING: Use Chain of Custody (CC1) to indicate which bottles were filtered

Sampler: John V. Rudd Employer: John V. Rudd PELA

I certify that sampling procedures were in accordance with applicable EPA state and corporate protocols.

9/20/90 John V. Rudd
(Date) (Signature)



FIELD PARAMETER FORM (CC2)

Form 0002
Sample Management
12/89

ETC JOB # CA4756
Sample Point N 26W1H
Source Code _____ Sample Point I.D. _____

FIELD PROCEDURES

9.0.0.9.2.4 PURGE DATE (YY MM DD) 117318 START PURGE (2400 Hr Clock) 15 ELAPSED HRS _____ WATER VOL. IN CASING (Gallons) _____ VOLUME PURGED (Gallons)

SAMPLING METHOD: PRIVATE WELL PUMP PLUMBING

Sampler Type A-Submersible Pump D-Dipper/Bottle X-Other _____
 B-ISCO E-Bailer (SPECIFY OTHER) _____
 C-Bladder Pump F-Scoop/Shovel

Sampler Material A-Teflon C-PVC X-Other _____
 B-Metal D-Plastic (SPECIFY OTHER) _____

Tubing Material A-Teflon C-Polyethylene X-Other _____
 B-Tygon D-Silicon (SPECIFY OTHER) _____

Sample Compositing Y N Procedure/Proportions _____

FIELD MEASUREMENTS

Well Elevation (ft/msl) _____ Well Depth (ft) _____
Depth to Ground water (ft) _____ Sample Depth (non-well) (ft) _____
Groundwater Elevation (ft msl) _____

| | | | | | |
|------------------------------|----------------------------|----------------|-------------------------|-------------|-------------|
| 1st <u>7.07</u> (STD) pH | 1st <u>439</u> spec. cond. | um/cm at 25 °C | _____ (other parameter) | _____ value | _____ units |
| 2nd <u>7.07</u> (STD) pH | 2nd <u>437</u> spec. cond. | um/cm at 25 °C | _____ (other parameter) | _____ value | _____ units |
| 3rd <u>7.07</u> (STD) pH | 3rd <u>439</u> spec. cond. | um/cm at 25 °C | _____ (other parameter) | _____ value | _____ units |
| 4th <u>7.08</u> (STD) pH | 4th <u>438</u> spec. cond. | um/cm at 25 °C | _____ (other parameter) | _____ value | _____ units |
| <u>13.3</u> (°C) Sample Temp | _____ NTU | Turbidity | | | |

FIELD COMMENTS

Sample Appearance: CLEAR - NO ODR
Weather Conditions: PC - 60° F OUTSIDE
Other: SAMPLE TAKEN IN BASEMENT, APPROX TEMP 80° F

FILTERING: Use Chain of Custody (CC1) to indicate which bottles were filtered

Sampler: John Y Rudd (Print) Employer: John Y Rudd PFLA

I certify that sampling procedures were in accordance with applicable EPA state and corporate protocols.

9/20/90 (Date) John Y Rudd (Signature)

ORIGINAL

CHAIN OF CUSTODY FORM (CC1)

Company: WASTE MANAGEMENT, INC. Attn.: DEE BENCICH
 Facility/Site: _____ Phone: (708) 409 - 0700
 Address: , DANE COUNTY, WI

SAMPLE IDENTIFICATION

Facility: 40E (Facility/Site Code) _____ (Optional Sample Point Descriptions)
 Sample Point: 1X-2FIB (Source Code) 91009119 (Your Sample Point ID) _____ (Start Date) _____ (Start Time) _____ (Elapsed Hours)
 Source Codes:
 Well (W) _____ Outfall (O) _____ Bottom Sediment (B) _____ Surface Impoundment (I) _____ Leachate Collection Sys. (L) _____ Other (X) LAB
 Soil (S) _____ River/Stream (R) _____ Generation Point (G) _____ Treatment Facility (T) _____ Lake/Ocean (L) _____ Specify _____

SHUTTLE CONTENTS

| No | BOTTLE | | | ANALYSIS | SAMPLER | | LAB |
|----|--------|--------------------------------------|-------------|-----------------|------------|--------------|--------------|
| | Type | Size | Preserv. | | FIL. (Y/N) | Observations | Observations |
| 2 | VOA | 40 | HCL | VOLATILES | N | | ✓ |
| 1 | CONU | 125 | NONE | CHLORIDE | N | | ✓ |
| 1 | FL | 125 | NONE | FLUORIDE | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | TOC2 | N | | ✓ |
| 1 | CONS | 125 | H2SO4 | PHENOLS/NO2/NO3 | N | | ✓ |
| | MET | 1000 | HNO3 | METALS | N | | ✓ |
| 1 | CYAN | <u>1000 w. 9-30-90</u> <u>500</u> | <u>NaOH</u> | CY/T | N | | ✓ |

CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) DAN GREEN Date: 9-3-90 Time: 15:15
 Signature: [Signature] Seal #: 0185151 Intact: NO

2. I have received these materials in good condition from the above person.
 Name: John Y. Rudd Signature: [Signature]
 Date: 9/15/90 Time: 1000 Remarks: OK

3. I have received these materials in good condition from the above person.
 Name: _____ Signature: _____
 Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John Y. Rudd Date: 9/15/90 Time: 1205
 Signature: [Signature] Seal #: 0175392 Intact: OK

LAB USE ONLY Opened By: [Signature] Date: 9/20/90 Time: 920
 SHUTTLE # 500 TEMP. °C 6° SEAL # 175392 COND. [Signature]

ETC JOB # CA 4763
 Sample Point 21FR _____
Source Code Sample Point I.D.

FIELD PROCEDURES

PURGE DATE (YY MM DD) _____ START PURGE (2400 Hr Clock) _____ ELAPSED HRS _____ WATER VOL. IN CASING (Gallons) _____ VOLUME PURGED (Gallons) _____

SAMPLING METHOD: FIELD BLANK

Sampler Type A-Submersible Pump D-Dipper/Bottle
 B-ISCO E-Bailer
 C-Bladder Pump F-Scoop/Shovel X-Other _____ (SPECIFY OTHER)

Sampler Material A-Teflon C-PVC
 B-Metal D-Plastic X-Other _____ (SPECIFY OTHER)

Tubing Material A-Teflon C-Polyethylene
 B-Tygon D-Silicon X-Other _____ (SPECIFY OTHER)

Sample Compositd Y/N _____
Procedure/Proportions

FIELD MEASUREMENTS

Well Elevation (ft/msl) _____ Well Depth (ft) _____
 Depth to Ground water (ft) _____ Sample Depth (non-well) (ft) _____
 Groundwater Elevation (ft msl) _____

| | | | |
|--------------------|-----------------------|-----------------|---|
| 1st _____ (STD) ph | 1st _____ spec. cond. | uniform at 25°C | _____ (other parameter) _____ value _____ units |
| 2nd _____ (STD) ph | 2nd _____ spec. cond. | uniform at 25°C | _____ (other parameter) _____ value _____ units |
| 3rd _____ (STD) ph | 3rd _____ spec. cond. | uniform at 25°C | _____ (other parameter) _____ value _____ units |
| 4th _____ (STD) ph | 4th _____ spec. cond. | uniform at 25°C | _____ (other parameter) _____ value _____ units |

Sample Temp _____ (°C) Turbidity _____ NTU

FIELD COMMENTS

Sample Appearance: _____
 Weather Conditions: OVERCAST, COOL 50°F, WIND NW @ 6
 Other: _____

FILTERING: Use Chain of Custody (CC1) to indicate which bottles were filtered

Sampler: John Y. Rudd Employer: PELA
(Print)

I certify that sampling procedures were in accordance with applicable EPA state and corporate protocols.

5/15/90 John Y. Rudd
(Date) (Signature)



CHAIN OF CUSTODY FORM (CC1)

ORIGINAL

Date Sealed 9-5-90By: KJSCompany: WMTAttn.: Don GreenFacility/Site: Town of Swan

Phone: _____

Address: _____

SAMPLE IDENTIFICATION

Facility: _____

Facility/Site Code

41015

(Optional Sample Point Description)

Sample Point: X-2.FBSource Code
(from below)Your Sample Point ID
(left justify)Start Date
(YY/MM/DD)Start Time
(240 hr. clock)Elapsed Hours
(composite)9/09/90

Source Codes:

Well (W)

Outfall (O)

Bottom Sediment (B)

Surface Impoundment (I)

Leachate Collection Sys. (X)

Soil (S)

River/Stream (R)

Generation Point (G)

Treatment Facility (T)

Lake/Ocean (X)

LAB

SHUTTLE CONTENTS

| BOTTLE | | | | ANALYSIS | SAMPLER | | LAB |
|--------|------|------|----------|---------------------------|-----------|--------------|--------------|
| No | Type | Size | Preserv. | | FIL (Y/N) | Observations | Observations |
| 1 | P | 250 | Ø | TDS, Bicarbonate, Sulfate | | | full, clean |
| 1 | P | 250 | H2SO4 | NH3, TKN | | | ↓ |
| | | | | | | | |
| | | | | | | | |
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CHAIN OF CUSTODY CHRONICLE

1. Shuttle Opened By: (print) Don Green Date: 9-5-90 Time: 14:15
Signature: Don Green Seal #: 01376-01377 Intact: ✓

2. I have received these materials in good condition from the above person.
Name: John Y. Rudd Signature: John Y. Rudd
Date: 9/19/90 Time: 10:00 Remarks: OK

3. I have received these materials in good condition from the above person.
Name: _____ Signature: _____
Date: _____ Time: _____ Remarks: _____

4. Shuttle Sealed By: (print) John Y. Rudd Date: 9/19/90 Time: 18:15
Signature: John Y. Rudd Seal #: 01378 Intact: OK

LAB USE ONLY Opened By: Kelly Schell Date: 9-20-90 Time: 13:00
SHUTTLE # _____ TEMP. °C 40 SEAL # 1378 COND. Intact 140

WET CHEMISTRIES QC DATA

Batch: QW30597

Parameter: CHLORIDE

Date: 9-25-90

QC-Batch: QW60813
QW70264

Method Ref: 325.2

Time: 12:00 PM

Det Limit: 1.0

Instrument: GTPC

Analyst: H. Chaudhry

Verified: JLS 9/27/90

Matrix: H2O

INSTRUMENT CALIBRATION STANDARDS: Units = MG/L

| Standard | 200 | 100 | 75 | 50 | 5 |
|----------|-------|-------|------|------|------------|
| Observed | 200.9 | 100.5 | 76.0 | 49.6 | 4.75 (abs) |
| % Recov | 100 | 101 | 101 | 99 | 95 |

Limits: % Rec. = +/- 10.0 %, low std +/- 15% M_r

Comments:

Source Lot: _____
 y = mx + b
 Slope: NA
 Intep: _____
 Corr: _____
 Corr. Coef > 0.99500 []

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|--------|------------|---------|-------|-------|--------|-------|--|
| - | ICVER | +/- 15% | - | - | - | [] | Lot # _____ |
| - | ICBLK | < MDL | - | - | - | [] | _____ |
| 13, 14 | M. BLK - 1 | < MDL | 0.611 | - | - | [] | _____ |
| 15, 16 | S. BLK - 1 | +/- 15% | 49.8 | 50 | 99 | [] | Lot # _____ |
| 17, 18 | EXROC - 1 | +/- 20% | 97.9 | 97 | 101 | [] | Lot # ERA 9932 |
| 33 | CCV - 1 | +/- 15% | 104.9 | 100 | 105 | [] | _____ |
| 38 | CCB - 1 | < MDL | 0.797 | - | - | [] | _____ |
| 48 | CCV - 2 | +/- 15% | 105.2 | 100 | 105 | [] | _____ |
| 47 | CCB - 2 | < MDL | 1.3 | - | - | [] | Low level QO |
| 63 | CCV - 3 | +/- 15% | 106.4 | 100 | 106 | [] | _____ |
| 62 | CCB - 3 | < MDL | 1.9 | - | - | [] | Low level QO |
| - | M. BLK - 2 | < MDL | - | - | - | [] | _____ |
| - | S. BLK - 2 | +/- 15% | - | - | - | [] | Lot # _____ |
| - | EXROC - 2 | +/- 20% | - | - | - | [] | Lot # _____ |
| 78 | CCV - 4 | +/- 15% | 108.3 | 100 | 108 | [] | _____ |
| 77 | CCB - 4 | < MDL | 1.9 | - | - | [] | Low level QO |
| 93 | CCV - 5 | +/- 15% | 107.2 | 100 | 108 | [] | _____ |
| 92 | CCB - 5 | < MDL | 1.8 | - | - | [] | Low level QO |
| 108 | CCV - 6 | +/- 15% | 108.0 | 100 | 108 | [] | _____ |
| 107 | CCB - 6 | < MDL | 1.2 | - | - | [] | Low level QO |
| - | M. BLK - 3 | < MDL | - | - | - | [] | _____ |
| - | S. BLK - 3 | +/- 15% | - | - | - | [] | Lot # _____ |
| - | EXROC - 3 | +/- 20% | - | - | - | [] | Lot # _____ |
| 123 | CCV - 7 | +/- 15% | 106.5 | 100 | 107 | [] | _____ |
| 122 | CCB - 7 | < MDL | 1.5 | - | - | [] | Low level QO |
| 134 | CCV - 8 | +/- 15% | 106.5 | 100 | 107 | [] | _____ |
| 137 | CCB - 8 | < MDL | 2.710 | - | - | [] | 1/2 R 135 Cup #153 is little high on return |
| 153 | CCV - 9 | +/- 15% | 115.2 | 100 | 115 | [] | Sample from Cup #135-166 |
| 152 | CCB - 9 | < MDL | 1.2 | - | - | [] | _____ |

| QUALITY CONTROL | | | | Duplicate | | | Matrix Spikes | | | | |
|-----------------|-------|--------|------|-----------|-------|--------|---------------|-------|---------|------|--|
| Seq # | ETC # | ETC # | Orig | Dupe | % RPD | ETC # | Orig | Added | Recover | % R | |
| 20 | - | GB2974 | 9.3 | 9.5 | 0.2 | GB2974 | 9.3 | 50 | 58.0 | 98 | |
| 34 | - | GB2982 | 5.6 | 5.4 | 0.4 | GB2982 | 5.6 | 50 | 56.1 | 101 | |
| 49 | - | GB3528 | OVER | OVER | - | GB3528 | OVER | 50 | OVER | 97.2 | |
| 64 | - | GB2985 | 81.2 | 81.2 | 0.0 | GB2985 | 81.2 | 50 | 129.8 | 1-17 | |
| 79 | - | GB2527 | OVER | OVER | - | GB3527 | OVER | 50 | OVER | 1-17 | |
| 94 | - | GB3529 | OVER | OVER | - | GB3529 | OVER | 50 | OVER | 1-17 | |

C L C

WET CHEMISTRIES QC DATA

P.2/2.

S-Batch: QW 30597
QW 20264
QC-Batch: _____
Verified: _____

Parameter: CHLORIDE
Method Ref: 325.B
Det Limit: 1.0
Instrument: GTPC
Matrix: H₂O

Date: 9-25-90
Time: 12:00 PM
Analyst: H. Chandley

INSTRUMENT CALIBRATION STANDARDS: Units = MG/L
Standard _____
Observed See Page 1/2. _____ (abs)
% Recov _____
Limits: % Rec. = +/- 10.0 %, low std +/- 15% [],
Comments: _____

Source Lot: _____
y = mx + b
Slope: _____
Intep: _____
Corr: _____
Corr. Coef > 0.99500 []

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|-------|------------|---------|--------|-------|--------|-------|---------------|
| --- | ICVER | +/- 15% | --- | --- | --- | { } | Lot # _____ |
| --- | ICBLK | < MDL | --- | --- | --- | { } | _____ |
| --- | M. BLK - 1 | < MDL | --- | --- | --- | { } | _____ |
| --- | S. BLK - 1 | +/- 15% | --- | --- | --- | { } | Lot # _____ |
| --- | EXROC - 1 | +/- 20% | --- | --- | --- | { } | Lot # _____ |
| 168 | CCV - 10 | +/- 15% | 101.0 | 100 | 101 | { } | _____ |
| 167 | CCB - 10 | < MDL | 0.710 | --- | --- | { } | _____ |
| 183 | CCV - 8 11 | +/- 15% | 101.6 | 100 | 102 | { } | _____ |
| 182 | CCB - 8 11 | < MDL | 0.797 | --- | --- | { } | _____ |
| 198 | CCV - 8 12 | +/- 15% | 102.5 | 100 | 103 | { } | _____ |
| 197 | CCB - 8 12 | < MDL | -0.160 | --- | --- | { } | _____ |
| --- | M. BLK - 2 | < MDL | --- | --- | --- | { } | _____ |
| --- | S. BLK - 2 | +/- 15% | --- | --- | --- | { } | Lot # _____ |
| --- | EXROC - 2 | +/- 20% | --- | --- | --- | { } | Lot # _____ |
| 213 | CCV - 4 13 | +/- 15% | 102.3 | 100 | 102 | { } | _____ |
| 212 | CCB - 4 13 | < MDL | 2.6 | --- | --- | { } | Low level Pb. |
| 238 | CCV - 8 14 | +/- 15% | 103.6 | 100 | 104 | { } | _____ |
| 237 | CCB - 8 14 | < MDL | 0.291 | --- | --- | { } | _____ |
| 259 | CCV - 8 15 | +/- 15% | 103.7 | 100 | 104 | { } | _____ |
| 258 | CCB - 8 15 | < MDL | 1.2 | --- | --- | { } | Low level Cl. |
| --- | M. BLK - 3 | < MDL | --- | --- | --- | { } | _____ |
| --- | S. BLK - 3 | +/- 15% | --- | --- | --- | { } | Lot # _____ |
| 281 | EXROC - 3 | +/- 20% | --- | --- | --- | { } | Lot # _____ |
| 286 | CCV - 7 16 | +/- 15% | 104.0 | 100 | 104 | { } | _____ |
| 287 | CCB - 7 16 | < MDL | 0.3 | --- | --- | { } | _____ |
| 288 | CCV - 8 17 | +/- 15% | --- | --- | --- | { } | _____ |
| --- | CCB - 8 17 | < MDL | --- | --- | --- | { } | _____ |
| --- | CCV - 8 18 | +/- 15% | --- | --- | --- | { } | _____ |
| --- | CCB - 8 18 | < MDL | --- | --- | --- | { } | _____ |

M-C
9/26/90

| QUALITY CONTROL | | | | | Matrix Spikes | | | | | |
|-----------------|------|--------|------|------|---------------|--------|------|-----|-------|-----|
| Seq # | QC # | ETC # | Orig | Dupe | % MPD | ETC # | Orig | Add | Recon | % R |
| 109 | | FB2614 | 873 | 876 | 00 | FB2614 | 873 | 50 | 1364 | 98 |
| 124 | | FB1951 | 1.0 | 0.6 | --- | FB1951 | 1.0 | 50 | 50.6 | 99 |
| 139 | | FB2606 | 1.5 | 1.0 | --- | FB2606 | 1.5 | 50 | 1.0 | 100 |
| 154 | | CA4753 | 315 | 309 | --- | CA4753 | 315 | 50 | 314 | 100 |

Do not
also FB2606
CA4753
for QC.

ETC

P.2/2 POK
9/25/90

CHLORIDE

QW-Batch: QW30597
QW60813
QC-Batch: QW70264
QW05234
Verified: 3/25/90
Instrument: AAT

Method Ref: EPA 325.2, SW-846 9251
(Colorimetric/Automated Ferri Cyanide)
MDL: 1.0 mg/l
Matrix: Aqueous

Page 1 of 7
Date: 9-25-90
Time: 12:00 PM
Analyst: M. Chandley

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|---------------|-----------------------|
| 20 | GB2974 | 9.31 | - | 9.3 | |
| 24 | GB2975 | 5.09 | - | 5.1 | |
| 25 | GB2980 | 55.2 | - | 55.2 | |
| 26 | GB2983 | 4.73 | - | 4.7 | |
| 27 | GB4103 | 0.421 | - | BMDL | |
| 28 | GB3506 | OVER. | 1:10 | | |
| 30 | GB3510 | OVER | 1:10 | | |
| 34 | GB2982 | 5.57 | - | 5.6 | |
| 38 | GB3511 | OVER | 1:10 | | |
| 40 | GB3514 | OVER | 1:10 | | |
| 42 | GB3519 | 76.0 | 1:10 | 760 | |
| 44 | GB3520 | OVER | 1:10 | | |
| 46 | GB3538 | 87.3 | 1:10 | 873 | |
| 49 | GB3518 | OVER | - | | |
| 54 | GB4075 | CARRY OVER. | - | | |
| 55 | GB3505 | OVER | 1:10 | | |
| 56 | GB3508. | OVER | 1:10 | | |
| 57 | GB3509 | OVER | 1:10 | | |
| 58 | GB2976 | C. OVER. | - | | |
| 59 | GB2981 | 19.9 | - | 19.9 | |
| 60 | GB2995 | OVER | 1:10 | | |
| 61 | GB2984 | C. OVER. | - | | |
| 64 | GB2985 | 81.2 | - | 81.2 | |
| 68 | GB2986 | 2.11 | - | 2.1 | |

ETC

CHLORIDE

QW-Batch: QW30597
QW20213

QC-Batch: QW30597
QW20213

Verified: 9/27/90

Instrument: AA II

Method Ref: EPA 325.2, SW-846 9251

(Colorimetric/Automated Ferri Cyanide)

MDL: 1.0 mg/l

Matrix: Aqueous

Page 2 of 7

Date: 9-25-90

Time: 12:00 PM

Analyst: H. Chaudhry

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|---------------|-----------------------|
| 69 | GB3400 | OVER | - | | |
| 70 | GB3579 | C. OVER | - | | |
| 71 | GB3408 | C. OVER | - | | |
| 72 | GB3058 | 46.6 | - | 46.6 | |
| 73 | GB3701 | 21.6 | - | 21.6 | |
| 74 | GB2997 | OVER | - | | |
| 75 | GB3517 | OVER | - | | |
| 76 | GB3524 | OVER | - | | |
| 79 | GB3527 | OVER | - | | |
| 83 | GB3502 | OVER | 1:10 | | |
| 84 | GB3522 | OVER | - | | |
| 85 | GB3582 | C-OVER | 1:10 | | |
| 86 | GB3529 | OVER | - | | |
| 87 | GB3531 | OVER | - | | |
| 88 | GB3537 | OVER | - | | |
| 89 | GB3540 | OVER | - | | |
| 90 | GB3525. | OVER | - | | |
| 91 | GB3526 | OVER | - | | |
| 94 | GB3528 | OVER | - | | |
| 98 | GB3533 | OVER | - | | |
| 99 | GB3534 | OVER | - | | |
| 100 | GB3536 | OVER | - | | |
| 101 | GB2968 | C. OVER | - | | |
| 102 | GB2969 | 12.7 | - | 12.7 | |

CHLORIDE

QW-Batch: QW30597
QW60813
 QC-Batch: QW70254
QW00289
 Verified: RA 9/27/90
 Instrument: AATI.

Method Ref: EPA 325.2, SW-846 9251
 (Colorimetric/Automated Ferri Cyanide)
 MDL: 1.0 mg/l
 Matrix: Aqueous

Page 3 of 7
 Date: 9-25-90
 Time: 12:00 PM.
 Analyst: H. Chaudhry

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|--------------------|-----------------------|
| 103 | GB2970 | 14.9 | - | 14.9 | |
| 104 | FB2604 | 129.9 | - | 130.299 | M.C 9/26 |
| 105 | FB2605 | 115.8 | - | 116 | |
| 106 | FB2613 | 128.4 | - | 128 | |
| 109 | FB2614 | 87.3 | - | 87.3 | |
| 113 | FB2615 | 95.7 | - | 95.7 | |
| 114 | FB2616 | 152.4 | - | 152 | |
| 115 | FB2618 | 142.2 | - | 142 | |
| 116 | FB2623 | 172.5 | - | 173 | |
| 117 | FB2625 | 95.6 | - | 95.6 | |
| 118 | FB2634 | 119.2 | - | 119 | |
| 119 | FB2637 | 1.15 | - | 1.2 | |
| 120 | FB2655 | 20.6 | 1:10 | 206 | |
| 121 | FB1849 | OVER | - | | |
| 124 | FB1851 | 1.04 | - | 1.0 | |
| 128 | FB0782 | 2.77 | 1:20 | | |
| 129 | FB2645 | 42.3 | - | 42.3 | |
| 130 | FB2646 | 42.6 | - | 42.6 | |
| 131 | FB2647 | 44.6 | - | 44.6 | |
| 132 | FB2648 | 43.6 | - | 43.6 | |
| 133 | FB2649 | 11.6 | - | 11.6 | |
| 134 | FB2650 | 42.5 | - | 42.5 | |
| 135 | FB2607 | 12.2 | - | 12.2 | |
| 136 | FB2602 | 10.7 | - | 10.7 | |

ETC

CHLORIDE

QW-Batch: QW30597
QW60813

Method Ref: EPA 325.2, SW-846 9251

Page 4 of 7

QC-Batch: QW70264
QW05284

(Colorimetric/Automated Ferri Cyanide)

Date: 9-25-90

Verified: RAS 9/21/90

MDL: 1.0 mg/l

Time: 12:00 PM.

Instrument: AATL

Matrix: Aqueous

Analyst: M. Chaudhry

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|--------------------|-----------------|---------------|-----------------------|
| 139 | FB2603 | 10.3 | - | / | |
| 143 | FB2606 | 20.8 | - | / | |
| 144 | FB2608 | 10.3 | - | / | |
| 145 | FB2609 | 2.93 | - | / | |
| 146 | FB2610 | 47.0 | - | / | %R of ISS Cup #153 |
| 147 | FB2611 | 21.3 | - | / | out of Range - Return |
| 148 | FB2622 | 24.8 | - | / | Samples from |
| 149 | FB2636 | 1.22 | - | / | Cup #139-166. |
| 150 | HA2842 | OVER. | - | / | |
| 151 | HA2843 | 13.3 | - | / | |
| 154 | CA4753 | 12 31.3 | 1:26 | / | |
| 158 | CA4754 | 27.9 | - | / | |
| 159 | CA4755 | 14.5 | - | / | |
| 160 | CA4756 | 9.00 | - | / | |
| 161 | CA4763 | 1.45 | - | / | |
| 163 | GB3506 | OVER. | 1:50 | / | |
| 165 | GB3506 | 90.6 | 1:100 | / | |
| 166 | GB3506 | 42.4 | 1:200 | / | |
| 169 | GB3510 | OVER | 1:50 | / | |
| 171 | GB3510 | 162.3 | 1:100 | 16200 | |
| 172 | GB3510 | 76.4 | 1:200 | / | |
| 174 | GB3511 | 187.3 | 1:50 | 9370 | |
| 176 | GB3511 | 87.2 | 1:100 | / | |
| 177 | GB3511 | 43.1 | 1:200 | / | |

ETC

CHLORIDE

QW-Batch: QW30597
QW50813

Method Ref: EPA 325.2, SW-846 9251

Page 5 of 7

QC-Batch: QW302649
BT05549

(Colorimetric/Automated Ferri Cyanide)

Date: 9-25-90

Verified: PAH 9/27/90

MDL: 1.0 mg/l

Time: 12:00 PM

Instrument: AAII

Matrix: Aqueous

Analyst: M. Chaudhry

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|--|-----------------|---------------|--|
| 179 | GB3514 | OVER | 1:50 | | |
| 181 | GB3514 | 115.2 | 1:100 | 11500 | |
| 185 | GB3514 | ^{MC} 916 Lot 6 49.6 | 1:200 | | |
| 186 | GB3520 | 195.5 | 1:50 | | |
| 188 | GB3520 | 91.7 | 1:100 | 9200 | |
| 189 | GB3520 | 45.4 | 1:200 | | |
| 191 | GB3518 | 79.4 | 1:5 | 397 | |
| 193 | GB3518 | 39.1 | 1:10 | | |
| 194 | GB3518 | 20.6 | 1:20 | | |
| 196 | GB4075 | -0.0541 | - | BMDL | |
| 199 | GB3505 | OVER | 1:50 | | |
| 201 | GB3505 | 187.1 | 1:100 | 18700 | |
| 202 | GB3505 | 84.2 | 1:200 | | |
| 204 | GB3508 | OVER | 1:50 | | |
| 206 | GB3508 | OVER | 1:100 | | |
| 207 | GB3508 | OVER | 1:200 | | Return on 9/26/90 with batch # QW30597A |
| 209 | GB3509 | OVER | 1:50 | | |
| 211 | GB3509 | OVER | 1:100 | | |
| 214 | GB3509 | 156.2 | 1:200 | 31200 | |
| 216 | GB2976 | 4.37 | STR | 4.4 | |
| 217 | GB8995 | OVER | 1:50 | | |
| 219 | GB8995 | OVER | 1:100 | | |
| 220 | GB8995 | 121.6 | 1:200 | 24300 | |
| 222 | GB2984 | 4.12 | STR | 4.1 | |

ETC

CHLORIDE

QW-Batch: QW30597

Method Ref: EPA 325.2, SW-846 9251

Page 6 of 7QC-Batch: QW60813
QW70264

(Colorimetric/Automated Ferri Cyanide)

Date: 9-25-90Verified: 291 5/2/90

MDL: 1.0 mg/l

Time: 12:00 PMInstrument: AATI

Matrix: Aqueous

Analyst: M. Chandley

QTC05289

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|---------------|-----------------------|
| 223 | GB3400 | 189.1 | 1:5 | | |
| 225 | GB3400 | 44.5 | 1:20 | | |
| 226 | GB3400 | 91.8 | 1:10 | 918 | |
| 228 | GB3579 | 86.1 | 1:2 | 172 | |
| 230 | GB3408 | 99.9 | STR | 99.9 | |
| 231 | GB2997 | 99.7 | 1:2 | 199 | |
| 233 | GB3517 | 115.3 | 1:2 | 231 | |
| 235 | GB3524 | 42.9 | 1:5 | 215 | |
| 239 | GB3527 | 86.0 | 1:5 | 430 | |
| 241 | GB3527 | 20.7 | 1:20 | | |
| 242 | GB3527 | 40.3 | 1:10 | | |
| 244 | GB3502 | OVER | 1:50 | | |
| 246 | GB3502 | 62.0 | 1:200 | | |
| 247 | GB3502 | 144.2 | 1:100 | 14400 | |
| 249 | GB3522 | 39.0 | 1:10 | | |
| 251 | GB3522 | 19.3 | 1:20 | | |
| 252 | GB3522 | 80.7 | 1:5 | 404 | |
| 254 | GB3582 | 30.8 | 1:10 | 308 | |
| 255 | GB3582 | OVER. | STR | | |
| 260 | GB3529 | 12.3 | 1:20 | | |
| 261 | GB3529 | 24.7 | 1:10 | 247 | |
| 263 | GB3531 | 15.3 | 1:20 | | |
| 264 | GB3531 | 32.9 | 1:10 | 329 | |
| 266 | GB3537 | 103.2 | 1:2 | 206 | |

ETC

CHLORIDE

OW-Batch: QW30597
QW60813
QC-Batch: QW702640
QW005289
Verified: BAK 9/27/90
Instrument: AAII

Method Ref: EPA 325.2, SW-846 9251
(Colorimetric/Automated Ferri Cyanide)
MDL: 1.0 mg/l
Matrix: Aqueous

Page 7 of 7
Date: 9-25-90
Time: 12:00 PM.
Analyst: M. Chandley

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|---------------|-----------------------|
| 267 | GB3540 | 11.9 | 1:20 | | |
| 268 | GB3540 | 24.3 | 1:10 | | |
| 269 | GB3540 | 46.5 | 1:5 | 233 | |
| 271 | GB3525 | 20.1 | 1:20 | | |
| 272 | GB3525 | 41.0 | 1:10 | | |
| 273 | GB3525 | 85.7 | 1:5 | 429 | |
| 275 | GB3526 | 15.6 | 1:20 | | |
| 276 | GB3526 | 32.4 | 1:10 | | |
| 277 | GB3526 | 65.1 | 1:5 | 326 | |
| | | | | | Return sample after |
| | | | | | cup # 281 - Computer |
| | | | | | took cup # upto |
| | | | | | 269. |
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

WET CHEMISTRIES QC DATA

P.1/1.

REC 9/26/90

S-Batch: QW 30597A

Parameter: CHLORIDE⁻

Date: 9-26-90

QC-Batch: QW 70264A

Method Ref: 325-2

Time: 9:00 AM

Verified: PHS 9/21/90

Det Limit: 1.0

Instrument: GTFC

Matrix: H₂O

Analyst: H. Chandley

INSTRUMENT CALIBRATION STANDARDS: Units = MG/L

Standard 900 100 75 50 5

Observed 2006 100.4 75.2 50.1 5.07 (abs)

% Recovr 100 100 100 100 101

Limits: % Rec. = +/- 10.0 %, low std +/- 15% []

Comments:

Source Lot: _____

y = mx + b _____

Slope: N/A

Intcp: _____

Corr: _____

Corr. Coef > 0.99500 []

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|---------------|-------------------|-----------------|--------------|-------------|------------|-------|-------------------------|
| | ICVER | +/- 15% | | | | [] | Lot # _____ |
| | ICBLK | < MDL | | | | [] | |
| <u>13, 14</u> | <u>M. BLK - 1</u> | <u>< MDL</u> | <u>0.133</u> | | | [✓] | |
| <u>15, 16</u> | <u>S. BLK - 1</u> | <u>+/- 15%</u> | <u>50.0</u> | <u>50</u> | <u>100</u> | [✓] | Lot # _____ |
| <u>17, 18</u> | <u>EXROC - 1</u> | <u>+/- 20%</u> | <u>99.1</u> | <u>97.0</u> | <u>102</u> | [✓] | Lot # <u>9932 - ERA</u> |
| <u>33</u> | <u>CCV - 1</u> | <u>+/- 15%</u> | <u>100.4</u> | <u>100</u> | <u>100</u> | [✓] | |
| <u>32</u> | <u>CCB - 1</u> | <u>< MDL</u> | <u>0.438</u> | | | [✓] | |
| <u>48</u> | <u>CCV - 2</u> | <u>+/- 15%</u> | <u>101.1</u> | <u>100</u> | <u>101</u> | [✓] | |
| <u>47</u> | <u>CCB - 2</u> | <u>< MDL</u> | <u>0.850</u> | | | [✓] | |
| <u>63</u> | <u>CCV - 3</u> | <u>+/- 15%</u> | <u>101.9</u> | <u>100</u> | <u>102</u> | [✓] | |
| <u>62</u> | <u>CCB - 3</u> | <u>< MDL</u> | <u>0.4</u> | | | [✓] | |
| | <u>M. BLK - 2</u> | <u>< MDL</u> | | | | [] | |
| | <u>S. BLK - 2</u> | <u>+/- 15%</u> | | | | [] | Lot # _____ |
| | <u>EXROC - 2</u> | <u>+/- 20%</u> | | | | [] | Lot # _____ |
| <u>78</u> | <u>CCV - 4</u> | <u>+/- 15%</u> | <u>101.5</u> | <u>100</u> | <u>102</u> | [✓] | |
| <u>77</u> | <u>CCB - 4</u> | <u>< MDL</u> | <u>0.238</u> | | | [✓] | |
| | <u>CCV - 5</u> | <u>+/- 15%</u> | | | | [] | |
| | <u>CCB - 5</u> | <u>< MDL</u> | | | | [] | |
| | <u>CCV - 6</u> | <u>+/- 15%</u> | | | | [] | |
| | <u>CCB - 6</u> | <u>< MDL</u> | | | | [] | |
| | <u>M. BLK - 3</u> | <u>< MDL</u> | | | | [] | |
| | <u>S. BLK - 3</u> | <u>+/- 15%</u> | | | | [] | Lot # _____ |
| | <u>EXROC - 3</u> | <u>+/- 20%</u> | | | | [] | Lot # _____ |
| | <u>CCV - 7</u> | <u>+/- 15%</u> | | | | [] | |
| | <u>CCB - 7</u> | <u>< MDL</u> | | | | [] | |
| | <u>CCV - 8</u> | <u>+/- 15%</u> | | | | [] | |
| | <u>CCB - 8</u> | <u>< MDL</u> | | | | [] | |
| | <u>CCV - 9</u> | <u>+/- 15%</u> | | | | [] | |
| | <u>CCB - 9</u> | <u>< MDL</u> | | | | [] | |

| QUALITY CONTROL | | Duplicates | | | Matrix Spikes | | | | | |
|-----------------|-------|---------------|-------------|-------------|---------------|---------------|-------------|-----------|-------------|-----------|
| Seq # | ETC # | ETC # | Orig | Dupe | % RPD | ETC # | Orig | Added | Recur | % R |
| <u>20 - 23</u> | | <u>FB2603</u> | <u>9.2</u> | <u>9.3</u> | <u>01</u> | <u>FB2603</u> | <u>9.2</u> | <u>50</u> | <u>58.6</u> | <u>99</u> |
| <u>34 - 37</u> | | <u>CA4753</u> | <u>30.7</u> | <u>30.7</u> | <u>00</u> | <u>CA4753</u> | <u>30.7</u> | <u>50</u> | <u>76.4</u> | <u>91</u> |

WET CHEMISTRIES QC DATA

QW30597
QW60813
S-Batch: QW70264

Parameter: CHLORIDE
Method Ref: EPA 825.2
Det Limit: 1.0 mg/L
Instrument: GTPC
Matrix: H₂O

Date: 9/25/90

QC-Batch: _____

Time: 12:00pm

Verified: JHS 9/27/90

Analyst: M. Chaudry

COMMENTS/CONFORMANCE SUMMARY

CCV failure @ 115.2% - Analyst judgment to
re-analyze affected samples. Some low level
contaminant on CCAs however this is not Method
paper criteria. All other Method paper
criteria were acceptable.

Verification: [Signature]
Date: 9/27/90

Analyst Signature: [Signature] for ME
Date: 9/27/90

ETC

CHLORIDE

QW-Batch: QW30597A
QW6083A

Method Ref: EPA 325.2, SW-846 9251

Page 2 of 92 2nd of 92

QC-Batch: Q470364A
QTC05388

(Colorimetric/Automated Ferri Cyanide)

Date: 9-26-90

Verified: PHS 9/27/90

MDL: 1.0 mg/l

Time: 9:00 AM

Instrument: CTPC

Matrix: Aqueous

Analyst: M. Chavelly

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|---------------------|-----------------------|
| 20 | FB 2603 | 9.18 | - | 9.2 | |
| 24 | FB 2606 | 19.8 | - | 19.8 | |
| 25 | FB 2608 | 9.87 | - | 9.9 | |
| 26 | FB 2609 | 2.66 | - | 2.7 | |
| 27 | FB 2610 | 44.8 | - | 44.8 | |
| 28 | FB 2611 | 19.9 | - | 19.9 | |
| 29 | FB 2622 | 22.4 | - | 22.4 | |
| 30 | FB 2636 | 0.537 | - | BMDL | |
| 31 | HA 2843 | 11.6 | - | 11.6 | |
| 34 | CA 4753 | 30.7 | - | 30.7 | |
| 38 | CA 4754 | 26.3 | - | 26.3 | |
| 39 | CA 4755 | 13.6 | - | 13.6 | |
| 40 | CA 4756 | 8.27 | - | 8.3 | |
| 41 | CA 4763 | 0.447 | - | BMDL | |
| 42 | HA 2842 | 83.6 | 1:10 | 836 | |
| 44 | HA 2842 | 41.9 | 1:20 | | |
| 46 | GB 3506. | 89.3 | 1:100 | 8930 | |
| 50 | GB 3528 | 43.7 | 1:5 | | |
| 51 | GB 3528 | 116.3 | 1:2 | 233 | |
| 53 | GB 3533 | 82.6 | 1:5 | 413 | |
| 54 | GB 3533 | over. | 1:2 | | |
| 57 | GB 3534 | 91.4 | 1:10 | 914 | |
| 59 | GB 3536 | 140.7 | 1:20 | 140 1410 | M.C 9/27/90 |
| 60 | GB 3536 | 67.5 | 1:20 | | |

M.C

ETC

FLUORIDE

QW-Batch: QW70261 30621

Method Ref: SM 413 E

Page 1 of 2

QC-Batch: QTC05299

(Colorimetric, Automated Complexone)

Date: 9-28-90

Verified: 8/6/90

MDL: 0.1 mg/l

Time: 8:00am

Instrument: GTPL

Matrix: Aqueous

Analyst: OB

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|---------------|-----------------------|
| 20 | CA4753 | .261 | / | 0.3 | |
| 24 | CA4763 | .148 | / | 0.2 | See Cup # 145 c/o |
| 25 | CA4755 | .245 | / | 0.2 | |
| 26 | CA4754 | .242 | / | 0.2 | |
| 27 | CA4756 | .194 | / | 0.2 | |
| 31 | 6B3529 | 1.23 | / | 1.2 | |
| 35 | 6B3540 | .800 | / | 0.8 | |
| 36 | 6B3537 | 1.18 | / | 1.2 | |
| 37 | 6B3531 | .951 | / | 1.0 | |
| 38 | 6B3528 | 1.30 | / | 1.3 | |
| 39 | 6B3533 | 1.21 | / | 1.2 | |
| 40 | 6B3525 | 1.02 | / | 1.0 | |
| 41 | 6B3526 | 1.55 | / | 1.6 | |
| 42 | 6B4069 | .419 | / | 0.5 | |
| 43 | 6B4067 | 1.226 | / | 0.2 | |
| 44 | 6B4073 | .0706 | / | BMDL | |
| 48 | 6B4070 | .197 | / | 0.2 | |
| 53 | 6B3531 | .992 | / | 1.0 | confirmed ✓ |
| 54 | 6B3534 | 1.17 | / | 1.2 | |
| 55 | 6B4075 | .0860 | / | BMDL | |
| 56 | 6B4072 | .0307 | / | BMDL | |
| 57 | 6B4068 | .165 | / | 0.2 | |
| 58 | 6B4071 | .187 | / | 0.2 | |
| 59 | 6B4074 | 1.0138 | / | BMDL | |

ETC

FLUORIDE

QW-Batch: QW7026130621 Method Ref: SM 413 E Page 2 of 2
 QC-Batch: GC05299 (Colorimetric, Automated Complexone) Date: 9-28-90
 Verified: 26/2/90 MDL: 0.1 mg/l Time: 8:00am
 Instrument: GTPC Matrix: Aqueous Analyst: AR

| Seq # | ETC Job # | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------------|-----------------|------------------|-----------------------|
| 60 | 6B4076 | 0.0109 | — | BMDL | |
| 64 | 6B3518 | 1.08 | — | 1.1 | |
| 69 | 6B3527 | 1.12 | — | 1.1 | |
| 70 | 6B3536 | OVER | Redo 1:10 | — | — |
| 71 | 6B2999 | C/O | Redo STR | — | — |
| 72 | 6B4066 | 1.292 | — | 1.292 | Redo str C/O |
| 73 | 6B3535 | Redo | 1:2 | OVER | — |
| 82 | 6B3000 | Redo | 1:5 | ISS FAILED | C/O |
| 84 | 6B3503 | Redo | STR | ↓ | |
| 98 | 6B3501 | Redo | 1:5 | ↓ | |
| 101 | 6B3509 | Redo | 1:100 1:200 | ↓ | |
| 113 | 6B3502 | C/O on Wash | ISS 1:20 | Redo | |
| 116 | 6B3508 | Redo | 1:2, 1:5 | ISS FAILED | C/O |
| 123 | 6B2995 | Redo | 1:10 | ↓ | |
| 128 | 6B3505 | Redo | 1:10 | ↓ | |
| 133 | 6B3504 | Redo | 1:50 | ↓ | |
| 138 | 6B3515 | Redo | 1:50 | ↓ | |
| 145 | CA4763 | 1.0815 | — | BADK | Redo str |
| 153 | 6B2991 | 1.237 | — | 1.237 | Redo str |
| 155 | 6B3522 | 1.982 | — | 1.0 | Redo str |
| | | | | | |
| | | | | | |
| | | | | | |
| | | | | | |

ETC

WET CHEMISTRIES QC DATA

S-Batch: QW70261

Parameter: Amides

Date: 10-17-93

QC-Batch: QTC05342

Method Ref: 4139

Time: 8:00am

Verified: mc 10/16/93

Det Limit: 0.1212

Instrument: 6782

Matrix: 15203

Analyst: DB

INSTRUMENT CALIBRATION STANDARDS: Units = mg/l

Standard 20 15 10 0.5 0

Source Lot: _____

y = mx + b _____

Observed 1.99 1.49 1.04 0.476 — (abs)

Slope: _____

Intcp: _____

% Recov 99% 99% 104% 95% _____

Corr: _____

Limits: % Rec. = +/- 10.0 %, low std +/- 15% []

Corr. Coef > 0.99500 []

Comments: _____

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|-----------|-------------------|-----------------|--------------|-------------|-------------|-------|-------------|
| <u>18</u> | <u>ICVER</u> | <u>+/- 15%</u> | <u>.493</u> | <u>.500</u> | <u>98%</u> | [] | Lot # _____ |
| <u>16</u> | <u>ICBLK</u> | <u>< MDL</u> | <u>Lmdl</u> | | | [] | _____ |
| <u>15</u> | <u>M. BLK - 1</u> | <u>< MDL</u> | <u>Lmdl</u> | | | [] | _____ |
| <u>17</u> | <u>S. BLK - 1</u> | <u>+/- 15%</u> | <u>.476</u> | <u>.500</u> | <u>95%</u> | [] | Lot # _____ |
| <u>19</u> | <u>EXROC - 1</u> | <u>+/- 20%</u> | <u>1.18</u> | <u>.97</u> | <u>120%</u> | [] | Lot # _____ |
| <u>27</u> | <u>CCV - 1</u> | <u>+/- 15%</u> | <u>4.58</u> | <u>1.00</u> | <u>96%</u> | [] | _____ |
| <u>28</u> | <u>CCB - 1</u> | <u>< MDL</u> | <u>Lmdl</u> | | | [] | _____ |
| <u>31</u> | <u>CCV - 2</u> | <u>+/- 15%</u> | <u>1.026</u> | <u>1.00</u> | <u>103%</u> | [] | _____ |
| <u>32</u> | <u>CCB - 2</u> | <u>< MDL</u> | <u>Lmdl</u> | | | [] | _____ |
| _____ | <u>CCV - 3</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCB - 3</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>M. BLK - 2</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>S. BLK - 2</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | Lot # _____ |
| _____ | <u>EXROC - 2</u> | <u>+/- 20%</u> | _____ | _____ | _____ | [] | Lot # _____ |
| _____ | <u>CCV - 4</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCB - 4</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCV - 5</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCB - 5</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCV - 6</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCB - 6</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>M. BLK - 3</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>S. BLK - 3</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | Lot # _____ |
| _____ | <u>EXROC - 3</u> | <u>+/- 20%</u> | _____ | _____ | _____ | [] | Lot # _____ |
| _____ | <u>CCV - 7</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCB - 7</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCV - 8</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCB - 8</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCV - 9</u> | <u>+/- 15%</u> | _____ | _____ | _____ | [] | _____ |
| _____ | <u>CCB - 9</u> | <u>< MDL</u> | _____ | _____ | _____ | [] | _____ |

| QUALITY CONTROL | | | Duplicates | | | Matrix Spikes | | | | |
|-----------------|-------|---------------|-------------|-------------|----------|---------------|-------------|------------|--------------|-------------|
| Seq # | QTC # | ETC # | Orig | Dupe | % RPD | ETC # | Orig | Added | Result | % R |
| _____ | _____ | <u>Q29763</u> | <u>Lmdl</u> | <u>Lmdl</u> | <u>0</u> | <u>Q29763</u> | <u>Lmdl</u> | <u>1.0</u> | <u>1.078</u> | <u>107%</u> |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |

ETC

WET CHEMISTRIES QC DATA

S-Batch: QW60812

Parameter: N03-N

Date: 9/27/90

QC-Batch: _____

Method Ref: 353.2

Time: _____

Verified: JLS 9/28/90

Det Limit: 0.1 mg/L

Instrument: ETC

Matrix: H2O

Analyst: C.O.

COMMENTS/CONFORMANCE SUMMARY

ISS (OCU) failure @ Seawater # 217. ALL
analyses to be re-analyzed from Seawater
211 → 266. All other QC
Criteria were acceptable.

Verification: JLS

Date: 9/28/90

Analyst Signature: C.O.

Date: 9/28/90

ETC

WET CHEMISTRIES QC DATA

Batch: 06260812

Parameter: NO3-N

P. 1 of 3.

Date: 9/27/90

QC-Batch: _____

Method Ref: 353.2

Time: _____

Verified: RJS 9/28/90

Det Limit: 0.1 mg/L

Instrument: ETC

Matrix: H2O's

Analyst: CO

INSTRUMENT CALIBRATION STANDARDS: Units = mg/L

Standard 2.00 1.00 0.50 0.10

Observed 2.01 1.01 0.504 0.107

% Recov 100 101 101 107

Limits: % Rec. = +/- 10.0 %, low std +/- 15% []

Comments: _____

Source Lot: _____

y = mx + b

Slope: N/A

Intcp: _____

Corr: _____

Corr. Coef > 0.9950 []

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|----------------|------------|---------|--------------|--------------|-------------|-------|-----------------------|
| - | ICVER | +/- 15% | | | | [] | Lot # _____ |
| - | ICBLK | < MDL | | | | [] | |
| <u>19,20</u> | M. BLK - 1 | < MDL | <u>0.013</u> | | | [] | |
| <u>21,22</u> | S. BLK - 1 | +/- 15% | <u>0.506</u> | <u>0.500</u> | <u>101</u> | [] | Lot # _____ |
| <u>25,26</u> | EDROC - 1 | +/- 20% | <u>0.964</u> | <u>0.98</u> | <u>98.4</u> | [] | Lot # <u>ERA 9920</u> |
| <u>41</u> | CCV - 1 | +/- 15% | <u>0.991</u> | <u>1.00</u> | <u>99.1</u> | [] | |
| <u>40</u> | CCB - 1 | < MDL | <u>0.018</u> | | | [] | |
| <u>57</u> | CCV - 2 | +/- 15% | <u>1.02</u> | <u>1.00</u> | <u>102</u> | [] | |
| <u>56</u> | CCB - 2 | < MDL | <u>0.015</u> | | | [] | |
| <u>73</u> | CCV - 3 | +/- 15% | <u>1.03</u> | <u>1.00</u> | <u>103</u> | [] | |
| <u>72</u> | CCB - 3 | < MDL | <u>0.018</u> | | | [] | |
| - | M. BLK - 2 | < MDL | | | | [] | |
| - | S. BLK - 2 | +/- 15% | | | | [] | Lot # _____ |
| - | EDROC - 2 | +/- 20% | | | | [] | Lot # _____ |
| <u>89</u> | CCV - 4 | +/- 15% | <u>1.04</u> | <u>1.00</u> | <u>104</u> | [] | |
| <u>88</u> | CCB - 4 | < MDL | <u>0.019</u> | | | [] | |
| <u>105</u> | CCV - 6 | +/- 15% | <u>1.05</u> | <u>1.00</u> | <u>105</u> | [] | |
| <u>104</u> | CCB - 6 | < MDL | <u>0.020</u> | | | [] | |
| <u>121</u> | CCV - 6 | +/- 15% | <u>1.05</u> | <u>1.00</u> | <u>105</u> | [] | |
| <u>120</u> | CCB - 6 | < MDL | <u>0.021</u> | | | [] | |
| - | M. BLK - 3 | < MDL | | | | [] | |
| - | S. BLK - 3 | +/- 15% | | | | [] | Lot # _____ |
| - | EDROC - 3 | +/- 20% | | | | [] | Lot # _____ |
| <u>137</u> | CCV - 7 | +/- 15% | <u>1.03</u> | <u>1.00</u> | <u>103</u> | [] | |
| <u>136</u> | CCB - 7 | < MDL | <u>0.025</u> | | | [] | |
| <u>148</u> | CCV - 8 | +/- 15% | <u>1.02</u> | <u>1.00</u> | <u>102</u> | [] | |
| <u>147</u> | CCB - 8 | < MDL | <u>0.023</u> | | | [] | |
| <u>153,144</u> | CCV - 9 | +/- 15% | <u>1.04</u> | <u>1.00</u> | <u>104</u> | [] | |
| <u>152</u> | CCB - 9 | < MDL | <u>0.023</u> | | | [] | |

RJS
9/28/90

| QUALITY CONTROL | | Duplicates | | | Matrix Spikes | | | | | |
|-----------------|-------|---------------|-------------|-------------|---------------|---------------|-------------|-------------|--------------|------------|
| Seq # | ETC # | ETC # | Orig | Dups | % RPD | ETC # | Orig | Added | Reconst | % R |
| <u>32-35</u> | | <u>FB2604</u> | <u>0.03</u> | <u>0.03</u> | <u>0</u> | <u>FB2604</u> | <u>0.03</u> | <u>0.50</u> | <u>0.524</u> | <u>99</u> |
| <u>43-46</u> | | <u>FB2616</u> | <u>0.02</u> | <u>0.02</u> | <u>0</u> | <u>FB2616</u> | <u>0.02</u> | <u>0.50</u> | <u>0.522</u> | <u>100</u> |
| <u>59-62</u> | | <u>FB2647</u> | <u>0.05</u> | <u>0.05</u> | <u>0</u> | <u>FB2647</u> | <u>0.05</u> | <u>0.50</u> | <u>0.538</u> | <u>98</u> |
| <u>76-79</u> | | <u>FB2610</u> | <u>0.04</u> | <u>0.03</u> | <u>29</u> | <u>FB2610</u> | <u>0.04</u> | <u>0.50</u> | <u>0.532</u> | <u>98</u> |
| <u>97-100</u> | | <u>FB3525</u> | <u>0.05</u> | <u>0.05</u> | <u>0</u> | <u>FB3525</u> | <u>0.05</u> | <u>0.50</u> | <u>0.554</u> | <u>101</u> |
| <u>107-110</u> | | <u>FB3533</u> | <u>0.03</u> | <u>0.03</u> | <u>0</u> | <u>FB3533</u> | <u>0.03</u> | <u>0.50</u> | <u>0.528</u> | <u>100</u> |

CV-16 5/90 Rev. 0

ETC

WET CHEMISTRIES QC DATA

S-Batch: QW60812

Parameter: N03-N
Method Ref: 353.2
Det Limit: 0.109/L
Instrument: ETC
Matrix: H2O

P. 2 of 3.

Date: 9/27/90

QC-Batch: _____

Time: _____

Verified: RJK 9/28/90

Analyst: CO.

INSTRUMENT CALIBRATION STANDARDS: Units = _____

Source Lot: _____

Standard _____

y = mx + b

Observed See p 1/1 (abs)

Slope: _____

Intercept: _____

Corr: _____

% Recovr _____

Corr. Coef > 0.99500 []

Limits: % Rec. = +/- 10.0 %, low std +/- 15% []

Comments: _____

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|---------|------------|---------|--------|-------|--------|-------|-------------|
| - | ICVER | +/- 15% | | | | ++ | Lot # _____ |
| - | ICBLK | < MDL | | | | ++ | |
| - | M. BLK - 1 | < MDL | | | | ++ | |
| - | S. BLK - 1 | +/- 15% | | | | ++ | Lot # _____ |
| - | EXROC - 1 | +/- 20% | | | | ++ | Lot # _____ |
| 163 | CCV - 10 | +/- 15% | 1.03 | 1.00 | 103 | ++ | |
| 162 | CCB - 10 | < MDL | 0.024 | | | ++ | |
| 169 | CCV - 211 | +/- 15% | 1.06 | 1.00 | 106 | ++ | |
| 168 | CCB - 211 | < MDL | 0.020 | | | ++ | |
| 185 | CCV - 312 | +/- 15% | 1.07 | 1.00 | 107 | ++ | |
| 184 | CCB - 312 | < MDL | 0.032 | | | ++ | |
| - | M. BLK - 2 | < MDL | | | | ++ | |
| - | S. BLK - 2 | +/- 15% | | | | ++ | Lot # _____ |
| - | EXROC - 2 | +/- 20% | | | | ++ | Lot # _____ |
| 193 | CCV - 413 | +/- 15% | 1.07 | 1.00 | 107 | ++ | |
| 192 | CCB - 413 | < MDL | 0.035 | | | ++ | |
| 201 | CCV - 514 | +/- 15% | 1.08 | 1.00 | 108 | ++ | |
| 200 | CCB - 514 | < MDL | 0.034 | | | ++ | |
| 209 | CCV - 515 | +/- 15% | 0.945 | 1.00 | 94.5 | ++ | |
| 208 | CCB - 515 | < MDL | 0.041 | | | ++ | |
| - | M. BLK - 3 | < MDL | | | | ++ | |
| - | S. BLK - 3 | +/- 15% | | | | ++ | Lot # _____ |
| - | EXROC - 3 | +/- 20% | | | | ++ | Lot # _____ |
| 216-218 | CCV - 716 | +/- 15% | 0.421 | 1.00 | 42.1 | ++ | FAILED |
| 216-218 | CCB - 716 | < MDL | | | | ++ | |
| 225 | CCV - 817 | +/- 15% | FAILED | | | ++ | |
| 224 | CCB - 817 | < MDL | | | | ++ | |
| 233 | CCV - 818 | +/- 15% | FAILED | | | ++ | |
| 232 | CCB - 818 | < MDL | | | | ++ | |

RJK
9/28/90

| QUALITY CONTROL | | | Duplicates | | Matrix Spikes | | | | | |
|-----------------|------|--------|-------------|-------|---------------|--------|-------|-------|-------|-----|
| Seq # | QC # | ETC # | Orig | Dups | % RPD | ETC # | Orig | Added | Recur | % R |
| 124 - 127 | | CA4753 | - OFF SCALE | | | | | | | |
| 132 - 135 | | CB2781 | 0.099 | 0.099 | 0 | CB2781 | 0.099 | 0.50 | 0.593 | 99 |

CV-16 5/90 Rev. 0

ETC

P. 3 of 3

WET CHEMISTRIES QC DATA

S-Batch: QW60812

Parameter: NO3-N
Method Ref: 353.2
Det Limit: 2.1 mg/L
Instrument: ETC
Matrix: H2O

Date: 9/27/90

QC-Batch: _____

Time: _____

Verified: PKS 9/28/90

Analyst: C.O.

INSTRUMENT CALIBRATION STANDARDS: Units = _____

Source Lot: _____

Standard _____

y = mx + b

Observed _____

See P 19 (abs)

Slope: _____

Interp: _____

Corr: _____

% Recov _____

Corr. Coef > 0.99500 []

Limits: % Rec. = +/- 10.0 %, low std +/- 15% []

Comments: _____

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obev. | Known | % Rec. | [Ver] | Comments |
|-------|------------|---------|--------|-------|--------|-------|----------|
| — | ICVER | +/- 15% | — | — | — | — | Lot # |
| — | ICBLK | < MDL | — | — | — | — | — |
| — | M. BLK - 1 | < MDL | — | — | — | — | Lot # |
| — | S. BLK - 1 | +/- 15% | — | — | — | — | Lot # |
| — | EDROC - 1 | +/- 20% | — | — | — | — | — |
| 241 | CCV - 19 | +/- 15% | FAILED | — | — | — | — |
| 240 | CCB - 19 | < MDL | — | — | — | — | — |
| 249 | CCV - 20 | +/- 15% | FAILED | — | — | — | — |
| 248 | CCB - 20 | < MDL | — | — | — | — | — |
| 257 | CCV - 21 | +/- 15% | FAILED | — | — | — | — |
| 256 | CCB - 21 | < MDL | — | — | — | — | — |
| — | M. BLK - 2 | < MDL | — | — | — | — | Lot # |
| — | S. BLK - 2 | +/- 15% | — | — | — | — | Lot # |
| — | EDROC - 2 | +/- 20% | — | — | — | — | — |
| 265 | CCV - 22 | +/- 15% | FAILED | — | — | — | — |
| 264 | CCB - 22 | < MDL | — | — | — | — | — |
| — | CCV - 23 | +/- 15% | — | — | — | — | — |
| — | CCB - 23 | < MDL | — | — | — | — | — |
| — | CCV - 24 | +/- 15% | — | — | — | — | — |
| — | CCB - 24 | < MDL | — | — | — | — | — |
| — | M. BLK - 3 | < MDL | — | — | — | — | Lot # |
| — | S. BLK - 3 | +/- 15% | — | — | — | — | Lot # |
| — | EDROC - 3 | +/- 20% | — | — | — | — | — |
| — | CCV - 7 | +/- 15% | — | — | — | — | — |
| — | CCB - 7 | < MDL | — | — | — | — | — |
| — | CCV - 8 | +/- 15% | — | — | — | — | — |
| — | CCB - 8 | < MDL | — | — | — | — | — |
| — | CCV - 9 | +/- 15% | — | — | — | — | — |
| — | CCB - 9 | < MDL | — | — | — | — | — |

| QUALITY CONTROL | | | | Matrix Spikes | | | | | | |
|-----------------|-------|-------|------|---------------|-------|-------|------|-------|-------|-----|
| Seq # | ETC # | ETC # | Orig | Dups | % RPD | ETC # | Orig | Added | Recur | % R |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |
| — | — | — | — | — | — | — | — | — | — | — |

ETC

WET CHEMISTRIES QC DATA

S-Batch: QW 70262

Parameter: NO2-N
 Method Ref: 253.2
 Det Limit: 0.1 mg/L
 Instrument: ETC
 Matrix: H2O3

P. 1 of 3

Date: 9/27/90

QC-Batch: _____

Time: _____

Verified: JHS 9/29/90

Analyst: CO.

INSTRUMENT CALIBRATION STANDARDS: Units = mg/L

Standard 1.00 0.500 0.200 0.100

Source Lot: _____

y = mx + b

Observed 1.00 0.499 0.198 0.100 mg/L
 (obs)

Slope: NA

Interp: _____

% Recov 100 99.8 99 100

Corr: _____

Limits: % Rec. = +/- 10.0 %, low std +/- 15% U

Corr. Coef > 0.99500 ++

Comments: _____

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|-------|------------|---------|-------|-------|--------|-------|-------------|
| - | ICVER | +/- 15% | | | | | Lot # _____ |
| - | ICBLK | < MDL | | | | | |
| 19,20 | M. BLK - 1 | < MDL | 0.006 | | | | |
| 23,24 | S. BLK - 1 | +/- 15% | 0.198 | 0.200 | 99 | | Lot # _____ |
| 24,25 | EXROC - 1 | +/- 20% | NA | | | | Lot # _____ |
| 40,42 | CCV - 1 | +/- 15% | 0.503 | 0.500 | 100.6 | | |
| 42,40 | CCB - 1 | < MDL | 0.005 | | | | |
| 58 | CCV - 2 | +/- 15% | 0.504 | 0.500 | 100.8 | | |
| 56 | CCB - 2 | < MDL | 0.002 | | | | |
| 74 | CCV - 3 | +/- 15% | 0.507 | 0.500 | 101 | | |
| 72 | CCB - 3 | < MDL | 0.002 | | | | |
| - | M. BLK - 2 | < MDL | | | | | |
| - | S. BLK - 2 | +/- 15% | | | | | Lot # _____ |
| - | EXROC - 2 | +/- 20% | | | | | Lot # _____ |
| 90 | CCV - 4 | +/- 15% | 0.508 | 0.500 | 102 | | |
| 88 | CCB - 4 | < MDL | 0.003 | | | | |
| 106 | CCV - 5 | +/- 15% | 0.510 | 0.500 | 102 | | |
| 104 | CCB - 5 | < MDL | 0.003 | | | | |
| 122 | CCV - 6 | +/- 15% | 0.518 | 0.500 | 104 | | |
| 120 | CCB - 6 | < MDL | 0.005 | | | | |
| - | M. BLK - 3 | < MDL | | | | | |
| - | S. BLK - 3 | +/- 15% | | | | | Lot # _____ |
| - | EXROC - 3 | +/- 20% | | | | | Lot # _____ |
| 138 | CCV - 7 | +/- 15% | 0.518 | 0.500 | 104 | | |
| 136 | CCB - 7 | < MDL | 0.005 | | | | |
| 149 | CCV - 8 | +/- 15% | 0.487 | 0.500 | 97.4 | | |
| 147 | CCB - 8 | < MDL | 0.012 | | | | |
| 154 | CCV - 9 | +/- 15% | 0.519 | 0.500 | 104 | | |
| 152 | CCB - 9 | < MDL | 0.005 | | | | |

| QUALITY CONTROL | | | | | Matrix Spikes | | | | | |
|-----------------|-------|-------|------|------|---------------|-------|------|-------|-------|-----|
| Seq # | ETC # | ETC # | Orig | Dupe | % RPD | ETC # | Orig | Added | Recur | % R |
| | | | | | | | | | | |
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ETC

WET CHEMISTRIES OC DATA

S-Batch: DU70262

Parameter: NO2-N

P. 2 of 3

Date: 9/27/90

OC-Batch: _____

Method Ref: 3532

Time: _____

Verified: RHS 9/28/90

Det Limit: 0.100

Analyst: CO

Instrument: ETC

Source Lot: _____

Units = _____

Y = mx + b

Observed See P 1/1 (abs)

Slope: _____

% Recov _____

Intercept: _____

Limits: % Rec. = +/- 10.0 % low std +/- 15% []

Corr. Coef > 0.99500 []

Comments: _____

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Unit | Obsv. | Known | % Rec. | [Var] | Comments |
|-------|------------|---------|-------|-------|--------|-------|-------------|
| 1 | ICVEN | +/- 15% | | | | | Lot # _____ |
| 2 | ICBLK | < MDL | | | | | Lot # _____ |
| 3 | M. BLK - 1 | < MDL | | | | | Lot # _____ |
| 4 | S. BLK - 1 | +/- 15% | | | | | Lot # _____ |
| 5 | DNOC - 1 | +/- 20% | | | | | Lot # _____ |
| 6 | CCV - 10 | +/- 15% | 0.526 | 0.500 | 105 | | Lot # _____ |
| 7 | CCB - 10 | < MDL | 0.002 | | | | Lot # _____ |
| 8 | CCV - 11 | +/- 15% | 0.518 | 0.500 | 104 | | Lot # _____ |
| 9 | CCB - 11 | < MDL | 0.002 | | | | Lot # _____ |
| 10 | CCV - 12 | +/- 15% | 0.525 | 0.500 | 105 | | Lot # _____ |
| 11 | CCB - 12 | < MDL | 0.003 | | | | Lot # _____ |
| 12 | M. BLK - 2 | < MDL | | | | | Lot # _____ |
| 13 | S. BLK - 2 | +/- 15% | | | | | Lot # _____ |
| 14 | DNOC - 2 | +/- 20% | | | | | Lot # _____ |
| 15 | CCV - 13 | +/- 15% | 0.519 | 0.500 | 104 | | Lot # _____ |
| 16 | CCB - 13 | < MDL | 0.003 | | | | Lot # _____ |
| 17 | CCV - 14 | +/- 15% | 0.530 | 0.500 | 106 | | Lot # _____ |
| 18 | CCB - 14 | < MDL | 0.005 | | | | Lot # _____ |
| 19 | CCV - 15 | +/- 15% | 0.528 | 0.500 | 106 | | Lot # _____ |
| 20 | CCB - 15 | < MDL | 0.005 | | | | Lot # _____ |
| 21 | M. BLK - 3 | < MDL | | | | | Lot # _____ |
| 22 | S. BLK - 3 | +/- 15% | | | | | Lot # _____ |
| 23 | DNOC - 3 | +/- 20% | | | | | Lot # _____ |
| 24 | CCV - 16 | +/- 15% | 0.528 | 0.500 | 106 | | Lot # _____ |
| 25 | CCB - 16 | < MDL | 0.008 | | | | Lot # _____ |
| 26 | CCV - 17 | +/- 15% | 0.529 | 0.500 | 106 | | Lot # _____ |
| 27 | CCB - 17 | < MDL | 0.005 | | | | Lot # _____ |
| 28 | CCV - 18 | +/- 15% | 0.536 | 0.500 | 107 | | Lot # _____ |
| 29 | CCB - 18 | < MDL | 0.009 | | | | Lot # _____ |

| Seq # | ETC # | Duplicate | Days | % RSD | ETC # | Spikes | Added | Recur | % R |
|-------|-------|-----------|------|-------|-------|--------|-------|-------|-----|
| 1 | | | | | | | | | |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
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| 30 | | | | | | | | | |

ETC

WET CHEMISTRIES QC DATA

S-Batch: 0W70262

Parameter: NO2-N

P.3 of 3

Date: 9/27/90

QC-Batch: _____

Method Ref: 352.2

Time: _____

Verified: [Signature] 9/27/90

Det Limit: 0.1

Instrument: ETA

Analyst: CO.

Matrix: H2O5

INSTRUMENT CALIBRATION STANDARDS: Units = _____

Source Lot: _____

Standard _____

y = mx + b

Observed See P11 (abc)

Slope: _____

Intcp: _____

Corr: _____

% Recover _____

Corr. Coef > 0.99800 []

Limits: % Rec. = +/- 10.0 %, low std +/- 15% []

Comments: _____

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|--------------------|------------|---------|--------------------|--------------|------------|-------|-------------|
| — | ICVER | +/- 15% | — | — | — | [] | Lot # _____ |
| — | ICBLK | < MDL | — | — | — | [] | Lot # _____ |
| — | M. BLK - 1 | < MDL | — | — | — | [] | Lot # _____ |
| — | S. BLK - 1 | +/- 15% | — | — | — | [] | Lot # _____ |
| — | EDROC - 1 | +/- 20% | — | — | — | [] | Lot # _____ |
| <u>242</u> | CCV - 19 | +/- 15% | <u>0.540</u> | <u>0.500</u> | <u>108</u> | [] | Lot # _____ |
| <u>240</u> | CCB - 19 | < MDL | <u>0.010</u> | — | — | [] | Lot # _____ |
| <u>250</u> | CCV - 20 | +/- 15% | <u>0.534</u> | <u>0.500</u> | <u>107</u> | [] | Lot # _____ |
| <u>248</u> | CCB - 20 | < MDL | <u>0.008</u> | — | — | [] | Lot # _____ |
| <u>258</u> | CCV - 21 | +/- 15% | <u>0.524</u> | <u>0.500</u> | <u>105</u> | [] | Lot # _____ |
| <u>256</u> | CCB - 21 | < MDL | <u>0.017</u> | — | — | [] | Lot # _____ |
| — | M. BLK - 2 | < MDL | — | — | — | [] | Lot # _____ |
| — | S. BLK - 2 | +/- 15% | — | — | — | [] | Lot # _____ |
| — | EDROC - 2 | +/- 20% | — | — | — | [] | Lot # _____ |
| <u>266</u> | CCV - 22 | +/- 15% | <u>0.526</u> | <u>0.500</u> | <u>105</u> | [] | Lot # _____ |
| <u>264</u> | CCB - 22 | < MDL | <u>0.004</u> | — | — | [] | Lot # _____ |
| <u>[Signature]</u> | CCV - 23 | +/- 15% | <u>[Signature]</u> | — | — | [] | Lot # _____ |
| <u>[Signature]</u> | CCB - 23 | < MDL | <u>[Signature]</u> | — | — | [] | Lot # _____ |
| <u>[Signature]</u> | CCV - 24 | +/- 15% | <u>[Signature]</u> | — | — | [] | Lot # _____ |
| <u>[Signature]</u> | CCB - 24 | < MDL | <u>[Signature]</u> | — | — | [] | Lot # _____ |
| — | M. BLK - 3 | < MDL | — | — | — | [] | Lot # _____ |
| — | S. BLK - 3 | +/- 15% | — | — | — | [] | Lot # _____ |
| — | EDROC - 3 | +/- 20% | — | — | — | [] | Lot # _____ |
| — | CCV - 7 | +/- 15% | — | — | — | [] | Lot # _____ |
| — | CCB - 7 | < MDL | — | — | — | [] | Lot # _____ |
| — | CCV - 8 | +/- 15% | — | — | — | [] | Lot # _____ |
| — | CCB - 8 | < MDL | — | — | — | [] | Lot # _____ |
| — | CCV - 9 | +/- 15% | — | — | — | [] | Lot # _____ |
| — | CCB - 9 | < MDL | — | — | — | [] | Lot # _____ |

| QUALITY CONTROL | | | | | Matrix Spikes | | | | | |
|-----------------|-------|-------|------|------|---------------|-------|------|-------|---------|-----|
| Seq # | ETC # | ETC # | Orig | Dupe | % RPD | ETC # | Orig | Added | Reconst | % R |
| | | | | | | | | | | |
| | | | | | | | | | | |
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ETC

NITRATE-NITRITE

QW-Batch: QW60812

Method Ref: EPA 353.2

Page 1 of 5

QC-Batch: QTC 05286

(Colorimetric, Automated Cd-Reduction)

Date: 9/27/90

Verified: R. H. 9/27/90

MDL: 0.1 mg/l

Time: -

Instrument: GTPC

Matrix: Aqueous

Analyst: C.O.

| Seq # | ETC Job # | pH Int. | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|------------|---------|---------------|-----------------|---------------|-----------------------|
| 28 | FB 2637 | 42 | 0.0271 | - | <0.1 | |
| 29 | FB 1851 | 42 | 0.0468 | - | <0.1 | |
| 30 | FB 2636 | 42 | 0.0212 | - | <0.1 | |
| 31 | FB 2638 | 42 | 0.0476 | - | <0.1 | |
| 32 | FB 2604 | 42 | 0.0292 | - | <0.1 | |
| 36 | FB 2605 | 42 | 0.116 | - | 0.12 | .1 c.o. 10/17 |
| 37 | FB 2613 | 42 | 0.0185 | - | <0.1 | |
| 38 | FB 2614 | 42 | 0.0298 | - | <0.1 | |
| 39 | FB 2615 | 42 | 0.0512 | - | <0.1 | |
| 43 | FB 2616 | 42 | 0.0213 | - | <0.1 | |
| 47 | FB 2618 | 42 | 0.103 | - | 0.10 | |
| 48 | FB 2618-23 | 42 | 0.0565 | - | <0.1 | |
| 49 | FB 2625 | 42 | 0.164 | - | 0.16 | .2 c.o. 10/17 |
| 50 | FB 2625 | 42 | 0.0231 | - | <0.1 | |
| 52 | FB 2637 | 1849 | OFFSCALE | - | | See Seq 123 |
| 53 | FB 2782 | 42 | 0.0346 | - | <0.1 | |
| 54 | FB 2645 | 42 | 0.0941 | - | <0.1 | |
| 55 | FB 2646 | 42 | 0.0841 | - | <0.1 | |
| 59 | FB 2647 | 42 | 0.0470 | - | <0.1 | |
| 63 | FB 2648 | 42 | 0.0686 | - | <0.1 | |
| 64 | FB 2649 | 42 | 0.0212 | - | <0.1 | |
| 65 | FB 2650 | 42 | 0.0884 | - | <0.1 | |
| 66 | FB 2607 | 42 | 0.609 | - | 0.61 | .6 c.o. 10/17 |
| 67 | FB 2603 | 42 | 0.0293 | - | <0.1 | |

ETC

NITRATE-NITRITE

QW-Batch: QW60812

Method Ref: EPA 383.2

Page 2 of 5

QC-Batch: WTC05286

(Colorimetric, Automated Cd-Reduction)

Date: 9/27/90

Verified: PLS 9/28/90

MDL: 0.1 mg/l

Time:

Instrument: GTPC

Matrix: Aqueous

Analyst: CO.

| Seq # | ETC Job # | pH Int. | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------|---------------|-----------------|-----------------|-----------------------|
| 68 | FB 2606 | <2 | 0.262 | - | 0.26 | .3 c.o. 10/1/90 |
| 69 | FB 2607 | <2 | 0.0933 | - | <0.1 | |
| 70 | FB 2608 | <2 | 1.70 | - | 1.7 | |
| 71 | FB 2609 | <2 | 0.0347 | - | <0.1 | |
| 76 | FB 2610 | <2 | 0.0352 | - | <0.1 | |
| 80 | FB 2611 | <2 | 0.268 | - | 0.27 | .3 c.o. 10/1/90 |
| 81 | FB 2622 | <2 | 0.0256 | - | <0.1 | |
| 82 | FB 2617 | <2 | 0.0453 | - | <0.1 | |
| 83 | FB 2619 | <2 | 0.129 | - | 0.13 | .1 c.o. 10/1/90 |
| 84 | FB 2620 | <2 | 0.0595 | - | <0.1 | |
| 85 | FB 2621 | <2 | 0.0316 | - | <0.1 | |
| 86 | FB 2624 | <2 | 0.0401 | - | <0.1 | |
| 87 | FB 2626 | <2 | 0.551 | - | 0.55 | .6 c.o. 10/1/90 |
| 91 | FB 2627 | <2 | 0.0714 | - | <0.1 | |
| 92 | FB 2628 | <2 | 0.642 | - | 0.64 | .6 c.o. 10/1/90 |
| 93 | FB 2629 | <2 | 0.0996 | - | <0.1 | |
| 94 | FB 2630 | <2 | 0.205 | - | 0.20 | .2 c.o. 10/1/90 |
| 95 | FB 2632 | <2 | 0.0975 | - | <0.1 | |
| 96 | FB 2633 | <2 | 0.0344 | - | <0.1 | |
| 97 | GB 3525 | <2 | 0.0529 | - | <0.1 | |
| 101 | GB 3526 | <2 | 1.76 | - | 1.76 | 1.8 c.o. 10/1/90 |
| 102 | GB 3528 | <2 | 0.0338 | - | <0.1 | |
| 107 | GB 3533 | <2 | 0.0338 | - | <0.1 | |
| 112 | GB 3534 | <2 | 0.278 | - | 0.28 | c.o. 10/1/90 |

ETC

NITRATE-NITRITE

QW-Batch: DW60812

Method Ref: EPA 353.2

Page 3 of 5

QC-Batch: QTC05286

(Colorimetric, Automated Cd-Reduction)

Date: 9/27/90

Verified: 9/28/90

MDL: 0.1 mg/l

Time:

Instrument: GTPC

Matrix: Aqueous

Analyst: C.O.

| Seq # | ETC Job # | pH Int. | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------|---------------------|-----------------|-----------------|-----------------------|
| 113 | GB 3536 | | 0.0467 | - | LO.1 | C.O. 10/1/90 |
| 114 | GB 2968 | | 1.62 | - | 1.62 | 1.6 10/1/90 |
| 115 | GB 2969 | | 0.113 | - | 0.11 | .1 10/1/90 |
| 116 | GB 2970 | | 0.335 | - | 0.34 | .3 10/1/90 |
| 117 | GB 2980 | | 0.0554 | - | LO.1 | |
| 118 | GB 2983 | | 0.0634 | - | LO.1 | |
| 123 | FB 1849 | | 0.15 1:2 | | 2.3 | |
| 124 | CA 4753 | | OFFSCALE | - | | |
| 128 | CA 4754 | | 0.127 | - | 0.13 | C/Over C.O. 10/1/90 |
| 129 | CA 4755 | | OFFSCALE | - | | |
| 130 | CA 4756 | | 0.0662 | - | LO.1 | |
| 131 | CA 4763 | | 0.0546 | - | LO.1 | |
| 132 | GB 2981 | | 0.0978 | - | LO.1 | |
| 139 | GB 2982 | | 0.0346 | - | LO.1 | |
| 140 | GB 2984 | | 0.0403 | - | LO.1 | |
| 141 | GB 2985 | | 0.0304 | - | LO.1 | |
| 142 | GB 2986 | | 0.0327 | - | LO.1 | |
| 143 | GB 2991 | | 1.43 | - | 1.43 | 1.4 C.O. 10/1/90 |
| 144 | GB 2999 | | 0.0598 | - | LO.1 | |
| 146 | GB 3523 | | 0.0325 | - | LO.1 | |
| 150 | FB 2655 | | 0.0542 | 1:2 | | |
| 151 | FB 2655 | | 0.0224 | STR | LO.1 | |
| 155 | GB 3534 | | 0.281 | - | 0.28 | .3 10/1/90 |
| 156 | GB 3536 | | 0.0616 | - | LO.1 | |

ETC

NITRATE-NITRITE

QW-Batch: QW60812

Method Ref: EPA 353.2

Page 4 of 5

QC-Batch: QTC05286

(Colorimetric, Automated Cd-Reduction)

Date: 9/27/90

Verified: 9/29/90

MDL: 0.1 mg/l

Time: -

Instrument: GTPC

Matrix: Aqueous

Analyst: C.O.

| Seq # | ETC Job # | pH Int. | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------|------------------------|-----------------|-----------------|-----------------------|
| 157 | GB 2968 | | 1.60 | - | 1.60 | C.O. 10/1/90 |
| 158 | GB 2969 | | 0.122 | - | 0.12 | C.O. 10/1/90 |
| 159 | GB 2970 | | 0.329 | - | 0.33 | C.O. 10/1/90 |
| 160 | GB 2980 | | 0.0569 | - | 0.1 | C.O. 10/1/90 |
| 161 | GB 2983 | | 0.0536 | - | 0.1 | C.O. 10/1/90 |
| 165 | CA 4753 | | 1.07 | 1:10 | 10.7 | |
| 166 | CA 4754 | | 0.0961 | STR | <0.1 | |
| 167 | CA 4755 | | 0.865 | 1:10 | 8.65 | 8.7 10/1/90 |
| 171 | GB 3503 | | 0.0288 | 1:20 | | |
| 172 | GB 3000 | | 0.0435 | 1:20 | | |
| 173 | GB 3520 | | 0.0374 | 1:20 | | REDO@ 1:2 |
| 174 | GB 3538 | | 0.0582 | STR | <0.1 | |
| 175 | GB 3501 | | 0.0488 | 1:20 | | |
| 176 | GB 3511 | | 0.178 | 1:20 | | SEE SEQ# 179 |
| 177 | GB 3509 | | 0.0300 | 1:20 | | |
| 178 | GB 3519 | | 0.0441 | STR | <0.1 | |
| 179 | GB 3535 | | 0.0330 | STR | <0.1 | |
| 180 | GB 3506 | | 0.815 | 1:20 | 16.3 | C.O. 10/1/90 |
| 181 | GB 3510 | | 0.788 ^{0.643} | 1:20 | 12.9 | |
| 182 | GB 3514 | | 0.119 | 1:20 | | |
| 183 | GB 3502 | | 1.303 | 1:20 | 26.1 | |
| 187 | GB 3515 | | 0.997 | 1:20 | 19.9 | |
| 188 | GB 3504 | | 0.555 | 1:20 | 11.1 | |
| 189 | GB 3508 | | 0.0614 | 1:20 | | REDO@ 1:5 |

ETC

1052

WET CHEMISTRIES QC DATA

S-Batch: 8W 30613 Parameter: PHENOLS.
8W 702667 Method Ref: 420.2
 QC-Batch: OTC05304 Det Limit: 0.05
 Verified: 8/6/90 MC 1974 Instrument: AAII
 Matrix: H2O

Date: 9-29-90
 Time: 6:50 AM

Analyst: H. Chaudhry

INSTRUMENT CALIBRATION STANDARDS: Units = ug/L

Standard 0.300 0.250 0.200 0.100 0.050
 Observed 0.300 0.251 0.203 0.105 0.0523 (std)
 % Recover 100 100 102 105 105
 Limits: % Rec. = +/- 10.0 % low std +/- 15% []
 Comments: Corr. Coef > 0.99900 []

Source Lot: _____
 Y = mx + b
 Slope: _____
 Intercept: _____
 Corr: _____

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Units

| Seq # | ETC ID | Unit | Obsv. | Known | % Rec. | Ver | Comments |
|-------|------------|---------|----------|-------|--------|-----|-------------------|
| 13,14 | M. BLK - 1 | < MDL | 0.00203 | 0.100 | 97 | { } | Lot # _____ |
| 15,16 | S. BLK - 1 | +/- 15% | 0.0498 | 0.208 | 113 | { } | Lot # <u>1927</u> |
| 17,18 | EXROC - 1 | +/- 20% | 0.235 | 0.200 | 102 | { } | |
| 35 | CCV - 1 | +/- 15% | 0.203 | 0.200 | 102 | { } | |
| 32 | CCB - 1 | < MDL | 0.00203 | 0.200 | 100 | { } | |
| 48 | CCV - 2 | +/- 15% | 0.200 | 0.200 | 100 | { } | |
| 47 | CCB - 2 | < MDL | 0.00419 | 0.200 | 102 | { } | |
| 63 | CCV - 3 | +/- 15% | 0.205 | 0.200 | 102 | { } | |
| 62 | CCB - 3 | < MDL | 0.00621 | 0.200 | 102 | { } | |
| - | M. BLK - 2 | < MDL | - | - | - | { } | Lot # _____ |
| - | S. BLK - 2 | +/- 15% | - | - | - | { } | Lot # _____ |
| - | EXROC - 2 | +/- 20% | - | - | - | { } | |
| 78 | CCV - 4 | +/- 15% | 0.204 | 0.200 | 102 | { } | |
| 77 | CCB - 4 | < MDL | 0.00851 | 0.200 | 101 | { } | |
| 93 | CCV - 5 | +/- 15% | 0.202 | 0.200 | 101 | { } | |
| 92 | CCB - 5 | < MDL | 0.00909 | 0.200 | 93 | { } | |
| 108 | CCV - 6 | +/- 15% | 0.196 | 0.200 | 93 | { } | |
| 107 | CCB - 6 | < MDL | 0.00371 | 0.200 | 93 | { } | |
| - | M. BLK - 3 | < MDL | - | - | - | { } | Lot # _____ |
| - | S. BLK - 3 | +/- 15% | - | - | - | { } | Lot # _____ |
| - | EXROC - 3 | +/- 20% | - | - | - | { } | |
| 123 | CCV - 7 | +/- 15% | 0.193 | 0.200 | 97 | { } | |
| 122 | CCB - 7 | < MDL | 0.00179 | 0.200 | 92 | { } | |
| 122 | CCV - 8 | +/- 15% | 0.194 | 0.200 | 92 | { } | |
| 133 | CCB - 8 | < MDL | -0.00871 | 0.200 | 73 | { } | |
| 133 | CCV - 9 | +/- 15% | 0.145 | 0.200 | 73 | { } | |
| 152 | CCB - 9 | < MDL | -0.00894 | 0.200 | 73 | { } | |

100 up to 153 Forward
 5 Retire samples from
 6/19/90 - 167

QUALITY CONTROL

| Seq # | ETC ID | Disposition | Dupe | % MPD | ETC ID | Con | Agree | Recur | % R |
|-------|--------|-------------|------|-------|--------|------|-------|-------|-----|
| 34 | 68370 | SMX | SMXL | 00 | 68370 | SMXL | 0100 | 00980 | 93 |
| 49 | 68375 | SMXL | SMXL | 00 | 68375 | SMXL | 0100 | 0101 | 101 |
| 64 | 68402 | SMXL | SMXL | 00 | 68402 | SMXL | 0100 | 0100 | 100 |
| 79 | 68475 | SMXL | SMXL | 00 | 68475 | SMXL | 0100 | 00992 | 100 |
| 94 | 68481 | SMXL | SMXL | 00 | 68481 | SMXL | 0100 | 00982 | 97 |
| 109 | 68489 | SMXL | SMXL | 00 | 68489 | SMXL | 0100 | 00948 | 97 |

WET CHEMISTRIES QC DATA

S-Batch: QW30613
QW 60851

Parameter: PHENOLS

Date: 9-29-90

QC-Batch: QW 70267
QW05304 H-2

Method Ref: 420.2

Time: 6:30 AM

Verified: S. Contreras 10/4

Det Limit: 0.05

Instrument: AAIT

Analyst: H. Chambliss

Matrix: H₂O

INSTRUMENT CALIBRATION STANDARDS: Units = _____

Source Lot: _____

Standard _____

y = mx + b

Observed _____ (abs)

Slope: _____

% Recov _____

Interp: _____

Limits: % Rec. = +/- 10.0 %, low std +/- 15% []

Corr: _____

Comments: _____

Corr. Coef > 0.99500 []

QUALITY CONTROL DATA SUMMARY: Batch and Instrumental QC Results/Limits

| Seq # | ETC ID | Limit | Obsv. | Known | % Rec. | [Ver] | Comments |
|-------|------------|---------|----------|---------|--------|---------|-------------|
| | ICVER | +/- 15% | | | | [] | Lot # _____ |
| | ICBLK | < MDL | | | | [] | |
| | M. BLK - 1 | < MDL | | | | { } | |
| | S. BLK - 1 | +/- 15% | | | | { } | Lot # _____ |
| | EXROC - 1 | +/- 20% | | | | { } | Lot # _____ |
| 168 | CCV - 10 | +/- 15% | 0.190 | 0.200 | 95 | [] | |
| 167 | CCB - 10 | < MDL | -0.00232 | | | [] | |
| 183 | CCV - 2 | +/- 15% | 0.191 | 0.200 | 96 | [] | |
| 192 | CCB - 2 | < MDL | 0 | | | [] | |
| 201 | CCV - 3 | +/- 15% | 0.189 | 0.200 | 94 | [] | |
| 200 | CCB - 3 | < MDL | -0.00201 | | | [] | |
| | M. BLK - 2 | < MDL | | | | { } | |
| | S. BLK - 2 | +/- 15% | | | | { } | Lot # _____ |
| | EXROC - 2 | +/- 20% | | | | { } | Lot # _____ |
| 218 | CCV - 4 | +/- 15% | 0.193 | 0.200 | 97 | [] | |
| 217 | CCB - 4 | < MDL | -0.00134 | 0.00134 | | [] | |
| | CCV - 5 | +/- 15% | | | | [] | |
| | CCB - 5 | < MDL | | | | [] | |
| | CCV - 6 | +/- 15% | | | | [] | |
| | CCB - 6 | < MDL | | | | [] | |
| | M. BLK - 3 | < MDL | | | | { } | |
| | S. BLK - 3 | +/- 15% | | | | { } | Lot # _____ |
| | EXROC - 3 | +/- 20% | | | | { } | Lot # _____ |
| | CCV - 7 | +/- 15% | | | | [] | |
| | CCB - 7 | < MDL | | | | [] | |
| | CCV - 8 | +/- 15% | | | | [] | |
| | CCB - 8 | < MDL | | | | [] | |
| | CCV - 9 | +/- 15% | | | | [] | |
| | CCB - 9 | < MDL | | | | [] | |

| QUALITY CONTROL | | | | | Matrix Spikes | | | | | |
|-----------------|------|-------|------|------|---------------|-------|------|------|--------|-----|
| Seq # | QC # | ETC # | Orig | Dupe | % RPD | ETC # | Orig | Adds | Result | % R |
| | | | | | | | | | | |
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ETC

TOTAL PHENOLICS

QW60831

QW-Batch: QW 301613

QC-Batch: QW 70267 QTC 05304

Verified: S. Grant 10/10/00

Instrument: _____

Method Ref: EPA 420.2, SW-846 9086

Colorimetric, Auto-4AAP, w/Distillation

MDL: 0.050 mg/l

Matrix: Aqueous

Page 1 of 4

Date: 9/29/90

Time: 6:30 AM

Analyst: M.C.

| Seq # | ETC Job # | pH Int. | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------|---------------|-----------------|---------------|-----------------------|
| | FB2565 | | .0165 | str. | BMDL | |
| | GB2968 | | .00739 | str. | BMDL | |
| | GB2969 | | .00329 | str. | BMDL | |
| | GB2970 | | .00748 | str. | BMDL | |
| | GB3582 | | .0276 | str. | BMDL | |
| | GB2980 | | .00388 | str. | BMDL | |
| | GB2983 | | .00266 | str. | BMDL | |
| | GB3918 | | .00225 | str. | BMDL | |
| | GB4075 | | .00176 | str. | BMDL | |
| | GB2981 | | .00207 | str. | BMDL | |
| | GB2982 | | .00221 | str. | BMDL | |
| | GB2984 | | .00189 | str. | BMDL | |
| | GB2985 | | .0109 | str. | BMDL | |
| | GB2986 | | .00698 | str. | BMDL | |
| | GB4068 | | .0102 | str. | BMDL | |
| | GB4003 | | MX | 2x | | |
| | | | mx | str | | |
| | | | mx | 5x | | |
| | | | .00421 | 10x | <.500 | |
| | | | .00246 | 20x | | |
| | GB4069 | | .00856 | str. | BMDL | |
| | GB4071 | | .0163 | str. | BMDL | |
| | GB4072 | | .00396 | str. | BMDL | |
| | GB4074 | | .00311 | str. | BMDL | |

TOTAL PHENOLICS

QW-Batch: QW60831 ^{Mc} Method Ref: EPA 420.2, SW-846 9088
 QC-Batch: QW30613 ⁷⁴⁴ Colorimetric, Auto-4AAP, w/Distillation
 Verified: S. G. 10/1/90 MDL: 0.050 mg/l
 Instrument: _____ Matrix: Aqueous

Page 2 of 4
 Date: 9/29/90
 Time: 6:30 AM
 Analyst: M.C.

| Seq # | ETC Job # | pH Int. | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------|---------------|-----------------|---------------|-----------------------|
| | GB4076 | | .00784 | str. | BMDL | |
| | GB4066 | | .0116 | str. | BMDL | |
| | GB4067 | | .0154 | str. | BMDL | |
| | GB4070 | | .0288 | str. | BMDL | |
| | GB4073 | | .00545 | str. | BMDL | |
| | CA4753 | | .00738 | str. | BMDL | |
| | CA4754 | | .00743 | str. | BMDL | |
| | CA4755 | | .00531 | str. | BMDL | |
| | CA4763 | | .00635 | str. | BMDL | |
| | HA 2998 | | MX | str. | / | |
| | | | .00168 | 2x | < .100 | |
| | | | .00435 | 5x | / | |
| | | | .00553 | 10x | / | |
| | CA4756 | | .00860 | str. | BMDL | |
| | FB2492 | | .0469 | str. | BMDL | |
| | FB2495 | | .00661 | str. | BMDL | |
| | FB2422 | | .00612 | str. | BMDL | |
| | FB2425 | | .0124 | str. | BMDL | |
| | FB2426 | | .00639 | str. | BMDL | |
| | FB2431 | | .00798 | str. | BMDL | |
| | FB2423 | | .00742 | str. | BMDL | |
| | FB2427 | | .00562 | str. | BMDL | |
| | FB2421 | | .00999 | str. | BMDL | |
| | FB2424 | | .00603 | str. | BMDL | |

ETC

TOTAL PHENOLICS

QW-Batch: ^{GW60831} GW 30613

Method Ref: EPA 420.2, SW-846 9088

Page 3 of 4

QC-Batch: GW 70267 ETC 05304

Colorimetric, Auto-4AAP, w/Distillation

Date: 9/29/90

Verified: E. L. [signature] 10/1/90

MDL: 0.050 mg/l

Time: 6:30 AM

Instrument: _____

Matrix: Aqueous

Analyst: M.C.

| Seq # | ETC Job # | pH Int. | Analyzed mg/l | Dilution Factor | Reported mg/l | Comments Calculations |
|-------|-----------|---------|---------------|-----------------|---------------|-----------------------|
| | FB2928 | | .109 | str. | .109 | |
| | FB2929 | | .0153 | str. | BMDL | |
| | FB2930 | | .00733 | str. | BMDL | |
| | FB2932 | | .00648 | str. | BMDL | |
| | FB2933 | | .00454 | str. | BMDL | |
| | FB2934 | | .00467 | str. | BMDL | |
| | FB2935 | | .00346 | str. | BMDL | |
| | FB2936 | | .00170 | str. | BMDL | |
| | FB2939 | | .00490 | str. | BMDL | |
| | FB2782 | | .0176 | str. | BMDL | |
| | FB1851 | | .00296 | str. | BMDL | |
| | FB1849 | | .00679 | str. | BMDL | |
| | GB4116 | | .115 | 10x | 1.15 | |
| | | | over | 5x | | |
| | | | mx/over | 2x | | |
| | | | mx/over | str | | |
| | GB3753 | | .0253 | 200x | | |
| | | | .0540 | 100x | | |
| | | | .116 | 50x | 5.80 | |
| | GB4119 | | .0147 | 200x | | |
| | | | .0438 | 100x | | ISS FAILED |
| | | | .101 | 50x | | ISS FAILED |
| | | | .117 | 50x | 5.85 | RERUN -OK |
| | GB3094 | | .0103 | 10x | | ISS FAILED |

WET CHEMISTRIES OC DATA

S-Date: 0130615 Parameter: TOC Date: 10/6/90
 OC-Serial: DT05324 Method Ref: EPA4151, SW346, 9060 Time: William
 Verifier: Blair Smith/klh Instrument: HC-1289 Meter: H2O Analyte: TOC

INSTRUMENT CALIBRATION STANDARDS: Units = mg/L Source Lot: _____
 Standard 0 20 50 110 400 Y = mx + b _____
 Observed 107 218 507 939 399.8 (lbs) Slope: _____
 % Recov 107 109 101 99 100 Intercept: _____
 Limits: % Rec. = +/- 10.0 %, low end +/- 15% [] Corr. Corr. Coef > 0.9990 []
 Comments: _____

QUALITY CONTROL DATA SUMMARY: Each and Instrumental QC Results/Units

| Seq # | ETC ID | Unit | Obsv. | Known | % Rec. | [Var] | Comments |
|------------|------------|---------|-------------|------------|------------|-------|-------------|
| | ICVER | +/- 15% | | | | | Lot # _____ |
| | ICBLK | < MDL | | | | | |
| <u>12</u> | M. BLK - 1 | < MDL | <u>0.6</u> | | | | Lot # _____ |
| <u>34</u> | S. BLK - 1 | +/- 15% | <u>107</u> | <u>10</u> | <u>107</u> | { } | Lot # _____ |
| <u>112</u> | DNOC - 1 | +/- 20% | <u>47</u> | <u>46</u> | <u>102</u> | { } | Lot # _____ |
| <u>41</u> | CCV - 1 | +/- 15% | <u>1013</u> | <u>100</u> | <u>101</u> | { } | |
| <u>69</u> | CCB - 1 | < MDL | <u>1032</u> | <u>100</u> | <u>103</u> | { } | |
| <u>90</u> | CCV - 2 | +/- 15% | <u>1044</u> | <u>100</u> | <u>104</u> | { } | |
| | CCB - 2 | < MDL | | | | | |
| | CCV - 3 | +/- 15% | | | | | |
| | CCB - 3 | < MDL | | | | | |
| | M. BLK - 2 | < MDL | | | | | Lot # _____ |
| | S. BLK - 2 | +/- 15% | | | | | Lot # _____ |
| | DNOC - 2 | +/- 20% | | | | | |
| | CCV - 4 | +/- 15% | | | | | |
| | CCB - 4 | < MDL | | | | | |
| | CCV - 6 | +/- 15% | | | | | |
| | CCB - 6 | < MDL | | | | | |
| | CCV - 8 | +/- 15% | | | | | |
| | CCB - 8 | < MDL | | | | | |
| | M. BLK - 3 | < MDL | | | | | Lot # _____ |
| | S. BLK - 3 | +/- 15% | | | | | Lot # _____ |
| | DNOC - 3 | +/- 20% | | | | | |
| | CCV - 7 | +/- 15% | | | | | |
| | CCB - 7 | < MDL | | | | | |
| | CCV - 9 | +/- 15% | | | | | |
| | CCB - 9 | < MDL | | | | | |
| | CCV - 9 | +/- 15% | | | | | |
| | CCB - 9 | < MDL | | | | | |

| QUALITY CONTROL | | Durations | | Matrix Spikes | | | | | |
|-----------------|-----------|---------------|------------|---------------|------------|---------------|------------|-----------|-------------|
| Seq # | ETC # | Obsv | Known | ETC # | Obsv | Added | Recur | % R | |
| <u>17</u> | <u>18</u> | <u>HA242</u> | <u>7.8</u> | <u>0</u> | <u>0.2</u> | <u>HA292</u> | <u>7.7</u> | <u>30</u> | <u>25.5</u> |
| <u>44</u> | <u>44</u> | <u>CA4754</u> | <u>6.0</u> | <u>6.0</u> | <u>0</u> | <u>CA4754</u> | <u>6.6</u> | <u>30</u> | <u>21.3</u> |
| <u>22</u> | <u>22</u> | <u>HA2159</u> | <u>6.7</u> | <u>6.6</u> | <u>0</u> | <u>HA2159</u> | <u>6.7</u> | <u>30</u> | <u>27.6</u> |

CV-16 5/90 Rev. 0

Total Organic Carbon (Water)

QW-Batch: 19W30615 Method Ref: EPA 415.1 SW846. 9060
 QC-Batch: QTC05328 (UV-Persulfate/IR)
 Verified: RC 1/10/10 MDL: 1.0 mg/l
 Instrument: HG1299 Matrix: Aqueous

Page: 1 of 6
 Date: 10/1/10
 Time: 11:00 am
 Analyst: Lew

| Log Link or Bottle # | ETC Job # STD or known: | pH Initial | Inject # | Analyzed Conc. | Dilution Factor | Reported Conc. | Comments % Rec. Calcs |
|-------------------------|----------------------------|---------------|-------------|-------------------|--------------------|-------------------|--------------------------------|
| | 400ppm | | 1 | 399.9 | | | |
| | | | 2 | 398.5 | | | |
| | | | 3 | 398.6 | | | |
| | | | 4 | 44.3 | | | |
| | Cal | | | 200.44 | | | |
| | Cal Rec. | | | 400.3 | | | |
| | Cal Adj. | | | 399.8 | | 399.8 | 100% rec |
| | 70.0 | | 1 | 0.459 | | 8.14 | |
| | | | 2 | 0.637 | | 8.14 | |
| | 10 ppm | | 3 | 10.73 | | 10.7 | 90R=1.7 |
| | | | 4 | 13.74 | | 10.7 | |
| | 20 ppm | | 5 | 24.62 | | | Cancelled Re-sample 10-1-10 |
| | | | 6 | 21.49 | | 21.5 | 70R=1.09 |
| | | | 7 | 21.97 | | 22.0 | |
| | 50 ppm | | 8 | 49.54 | | 50.0 | 100R=1.1 |
| | | | 9 | 51.34 | | 51.3 | |
| | | | 10 | 49.30 | | | Cancelled Re-sample |

ETC

Total Organic Carbon (Water)

OW-Batch: QW30615

Method Ref: EPA 415.1 SW846. 9060

Page 2 of 6

QC-Batch: DTCA 1324

(UV-Persulfate/IR)

Date: 10/6/90

Verified: RL 10/11/90

MDL: 1.0 mg/l

Time: 11:00 am

Instrument: HC1299

Matrix: Aqueous

Analyst: LW

| Log Link or Bottle # | ETC Job # STD or known | pH Initial | Inject # | Analyzed Conc. | Dilution Factor | Reported Conc. | Comments % Rec. Calcs |
|-------------------------|---------------------------|---------------|-------------|-------------------|--------------------|-------------------|--------------------------|
| | EPA "41" | | 11 | 47.40 | | 47.4 | 7.0 R = 1.24 |
| | | | 12 | 48.59 | | 48.6 | |
| | 100 ppm | | 13 | 97.14 | | 97.2 | 7.0 R = 99 |
| | | | 14 | 100.6 | | 100.6 | |
| 100626 | HA2992 | <2 | 15 | 7.652 | | 7.7 | |
| | | | 16 | 7.759 | | 7.8 | |
| | Dup | | 17 | 7.973 | | | |
| | | | 18 | 8.086 | | | |
| | SPK | | 19 | 22.27 | | | Cancelled |
| | | | 20 | 25.59 | | | Cancelled |
| | | | 21 | 25.28 | | | 7.0 R = 99 10/11/90 |
| 100626 | HA2993 | <2 | 22 | 28.62 | | 28.6 | |
| | | | 23 | 28.57 | | 28.6 | |
| 100626 | HA2994 | <2 | 24 | 10.55 | | 10.6 | |
| | | | 25 | 10.67 | | 10.7 | |
| 100626 | HA2995 | <2 | 26 | 11.89 | | 11.9 | |
| | | | 27 | 11.85 | | 11.9 | |

Total Organic Carbon (Water)

QW-Batch: QW30615

Method Ref: EPA 415.1 SW846. 9060

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QC-Batch: QTC053 > 9

(UV-Persulfate/IR)

Date: 10/6/90

Verified: SC + 10/11/90

MDL: 1.0 mg/l

Time: 11:00am

Instrument: HC1299

Matrix: Aqueous

Analyst: SW

| Log Link or Bottle # | ETC Job # STD or known | pH Initial | Inject # | Analyzed Conc. | Dilution Factor | Reported Conc. | Comments % Rec. Calc |
|-------------------------|---------------------------|---------------|-------------|-------------------|--------------------|-------------------|----------------------------------|
| 100626 | HA2996 | 4.2 | 28 | 11.47 | | 11.5 | |
| | | | 29 | 11.67 | | 11.7 | |
| 100626 | HA2998 | 4.2 | 30 | 4.269 | | 4.3 | |
| | | | 31 | 3.961 | | 4.0 | |
| 100625 | CA4753 | 4.2 | 32 | 1.189 | | 1.2 | |
| | | | 33 | 1.104 | | 1.1 | |
| 100625 | CA4754 | 4.2 | 34 | 0.854 | | | Cancelled Run extra for check |
| | | | 35 | 1.127 | | 1.1 | |
| | | | 36 | 1.175 | | 1.2 | |
| 100625 | CA4755 | 4.2 | 37 | 1.945 | | 1.9 | |
| | | | 38 | 2.111 | | 2.1 | |
| 100625 | CA4763 | 4.2 | 39 | 1.597 | | 1.6 | |
| | | | 40 | 1.801 | | 1.8 | |
| | 100 ppm | | 41 | 10.3 | | 10.3 | %R = 101 |
| 100631 | CA4754 | 4.2 | 42 | 0.592 | | BMDt | |
| | | | 43 | 0.599 | | BMDt | |

ETC

Total Organic Carbon (Water)

QW-Batch: RU30615

Method Ref: EPA 815.1 SW846. 9060

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QC-Batch: QTC05324

(UV-Persulfate/IR)

Date: 10/6/90

Verified: R. Cantor 10/19/90

MDL: 1.0 mg/l

Time: 11:00am

Instrument: HG1299

Matrix: Aqueous

Analyst: LW

| Log Link or Bottle # | ETC Job # STD or known | pH Initial | Inject # | Analyzed Conc. | Dilution Factor | Reported Conc. | Comments % Rec. Calcs |
|-------------------------|---------------------------|---------------|-------------|-------------------|--------------------|-------------------|--------------------------|
| | Dup | | 44 | 0.721 | | | |
| | | | 45 | 0.578 | | | |
| | SPK | | 46 | 11.68 | | | Cancelled |
| | | | 47 | 11.71 | | | Re-pte. Re-pte |
| | | | 48 | 41.44 | | | Cancelled |
| | | | 49 | 21.52 | | | Re-pte. Re-pte |
| | | | 50 | 21.07 | | | |
| 302236 | 482976 | 42 | 51 | 0.572 | | 8MDL | |
| | | | 52 | 0.698 | | 8MDL | |
| 302236 | 482981 | 42 | 53 | 5.531 | | 5.5 | |
| | | | 54 | 5.804 | | 5.8 | |
| 302236 | 482982 | 42 | 55 | 7.439 | | 7.4 | |
| | | | 56 | 7.245 | | 7.2 | |
| 302236 | 482984 | 42 | 57 | 3.549 | | 3.5 | |
| | | | 58 | 3.613 | | 3.6 | |
| 302236 | 482985 | 42 | 59 | 19.64 | | 19.6 | |
| | | | 60 | 19.61 | | 19.6 | |

ETC

Total Organic Carbon (Water)

QW-Batch: RW30615

Method Ref: EPA 415.1 SW846. 9060

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QC-Batch: OTC053-4

(UV-Persulfate/IR)

Date: 10/6/70

Verified: Blank dilution

MDL: 1.0 mg/l

Time: 11:00 am

Instrument: HG1299

Matrix: Aqueous

Analyst: LW

| Log Link or Bottle # | ETC Job # STD or known | pH Initial | Inject # | Analyzed Conc. | Dilution Factor | Reported Conc. | Comments % Rec. Calcs |
|-------------------------|---------------------------|---------------|-------------|-------------------|--------------------|-------------------|--------------------------|
| 302236 | UB2986 | <2 | 61 | 2.229 | | 2.2 | |
| | | | 62 | 2.261 | | 2.3 | |
| 302250 | UB2991 | <2 | 63 | 1.885 | | 1.9 | |
| | | | 64 | 1.769 | | 1.8 | |
| 302250 | UB3523 | <2 | 65 | 9.381 | | 9.4 | |
| | | | 66 | 9.127 | | 9.1 | |
| 302274 | UB3010 | <2 | 67 | 39.25 | | 39.3 | |
| | | | 68 | 39.35 | | 39.4 | |
| | 100 ppm | | 69 | 103.4 | | 103.4 | %R = 103 |
| 302271 | UB2658 | <2 | 70 | 6.527 | | 6.5 | |
| | | | 71 | 6.938 | | 6.9 | |
| | UB | | 72 | 6.532 | | / | |
| | | | 73 | 6.752 | | / | |
| | SPK | | 74 | 27.10 | | / | %R = 1.2 |
| | | | 75 | 26.81 | | / | |

ETC

Total Organic Carbon (Water)

QW-Batch: QW30615

Method Ref: EPA 415.1 SW846. 9060

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QC-Batch: QTC053

(UV-Persulfate/IR)

Date: 10/6/90

Verified: S. C. ...

MDL: 1.0 mg/l

Time: 11:00 am

Instrument: HG1299

Matrix: Aqueous

Analyst: Lu

| Log Link or Bottle # | ETC Job # STD or known | pH Initial | Inject # | Analyzed Conc. | Dilution Factor | Reported Conc. | Comments % Rec. Calcs |
|-------------------------|---------------------------|---------------|-------------|-------------------|--------------------|----------------------|---|
| 302250 | 4B2999 | ~2 | 76 | 4.216 | 1:20 | | Cancelled needs dilution |
| | | | 77 | 41.84 | 1:2 | 83.6 | |
| | | | 78 | 43.40 | 1:2 | 86.8 | |
| | | | 79 | 43.05 | 1:2 | 86.2 86.1 | 95% no |
| | | | 80 | 43.40 | 1:2 | 86.8 | |
| 302250 | 4B3000 | ~2 | 81 | 24.68 | 1:20 | 494 | 6/10/6/90 |
| | | | 82 | 24.67 | 1:20 | 494 494 | |
| | | | 83 | 24.36 | 1:20 | 488 | |
| | | | 84 | 24.73 | 1:20 | 494 | |
| 302251 | 4B4003 | ~2 | 85 | 850.3 | 1:2 | | Cancelled needs dilution |
| | | | 86 | 85.66 | 1:20 | 1914 | 1710 |
| | | | 87 | 86.62 | 1:20 | 1914 1932 | 1730 |
| 302243 | | | | | | | |
| 302251 | 4B3501 | ~2 | 88 | 24.40 | 1:20 | 488 | |
| | | | 89 | 24.98 | 1:20 | 500 | |
| | | | 90 | 104.4 | | 1044 | 7.8 = 1.14 |

for
10/6/90