

**Technical Report**  
for  
**ENVIRONMENTAL STRATEGIES CORP.**  
**WASHINGTON HARBOUR**  
**3050 KK ST. SUITE 325**  
**WASHINGTON, DC 20007**

*Chain of Custody Data Required for ETC Data Management Summary Reports*

P4850	ENVIRONMENTAL STRATEGIES CORP.	CHAMPION	WCSP-1	861110	1100	
<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>



**John J. Fitzgerald**  
*Vice President*  
*Research and Operations*

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## Introduction

This report contains the analytical results on your sample. It is designed to include comprehensive data from the entire analytical process in order to satisfy the needs of various levels of review.

The results obtained from your sample are presented in tabular format immediately following this introduction. Quality assurance data is tabulated along with the appropriate sample results for verification. Depending on the analyses ordered, the quality assurance data may include results from blank, spiked blank, spiked sample (i.e. matrix spike) and replicate sample as well as results from surrogate compound analyses. Quality assurance data for verification of proper instrument performance is also included where appropriate. The report appendices include the chain of custody record for your sample and, where appropriate, the gas chromatograms and mass spectra.

The procedures used in the analysis of the sample are described in this report's methodology section. All analytical procedures within our laboratory are performed within a strictly enforced Quality Assurance Protocol. A description of this Protocol is included in the report.

## Results

Sample results, and associated quality assurance data, are always tabulated in one or more of this report's Quantitative Results Tables. The format of each table varies with the class of analysis.

### *Priority Pollutants*

The priority pollutant compounds and elements are listed with their NPDES (National Pollution Discharge Elimination System) numbers, and the Method Detection Limit (MDL) published in the Federal Register. When a compound or element is present below its published MDL it is reported as BMDL (Below Method Detection Limit). When a compound or element is not present at any detectable concentrations it is reported as ND (Not Detected). MDL's for non-aqueous matrices are based on USEPA published MDL's but are adjusted as per sample weight. Matrix spike and replicate analyses, where included, were performed on samples randomly chosen within each quality assurance batch and are therefore not necessarily spikes and replicates of this report's sample. Surrogate compound recovery data and instrument calibration data are included in the Method Performance Data Tables.

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Volatile Compounds - GC/MS Analysis Data (QR01)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4850 ENVIRONMENTAL STRATEGIES CORP CHAMPION WCSP-1 861110 1100  
 ETC Sample No. Company Sample Point Date Time Flashed Hours

NPDES Number	Compound <small>Acetone and Methylene chloride values are screen only</small>	Results		QC Replicate		QC Blank and Spiked Blank		QC Matrix Spike			
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concn. Added ug/l	% Recov	Unspiked Sample ug/l	Concn. Added ug/l	% Recov
1V	Acrolein	ND	100	ND	ND	ND	800	88	ND	800	98
2V	Acrylonitrile	ND	100	ND	ND	ND	80.0	94	ND	80.0	84
3V	Benzene	ND	4.4	ND	ND	ND	18.0	99	ND	18.0	100
4V	bis(Chloromethyl)ether	ND	10	ND	ND	ND	18.0	-	ND	18.0	-
5V	Bromoform	ND	4.7	ND	ND	ND	18.0	92	ND	18.0	90
6V	Carbon tetrachloride	ND	2.8	ND	ND	ND	18.0	103	ND	18.0	98
7V	Chlorobenzene	ND	6.0	ND	ND	ND	18.0	99	ND	18.0	99
8V	Chlorodibromomethane	ND	3.1	ND	ND	ND	18.0	99	ND	18.0	95
9V	Chloroethane	ND	10	ND	ND	ND	18.0	103	ND	18.0	96
10V	2-Chloroethylvinyl ether	ND	10	ND	ND	ND	18.0	0 <sup>a</sup>	ND	18.0	0 <sup>a</sup>
11V	Chloroform	ND	1.6	ND	ND	ND	18.0	103	ND	18.0	98
12V	Dichlorobromomethane	ND	2.2	ND	ND	ND	18.0	100	ND	18.0	96
13V	Dichlorodifluoromethane	ND	10	ND	ND	ND	18.0	109	ND	18.0	104
14V	1,1-Dichloroethane	ND	4.7	ND	ND	ND	18.0	100	ND	18.0	99
15V	1,2-Dichloroethane	ND	2.8	ND	ND	ND	18.0	104	ND	18.0	96
16V	1,1-Dichloroethylene	ND	2.8	ND	ND	ND	18.0	101	ND	18.0	97
17V	1,2-Dichloropropane	ND	6.0	ND	ND	ND	18.0	98	ND	18.0	101
18V	cis-1,3-Dichloropropylene	ND	5.0	ND	ND	ND	18.0	97	ND	18.0	100
19V	Ethylbenzene	ND	7.2	ND	ND	ND	18.0	100	ND	18.0	100
20V	Methyl bromide	ND	10	ND	ND	ND	18.0	103	ND	18.0	109
21V	Methyl chloride	ND	10	ND	ND	ND	18.0	107	ND	18.0	109
22V	Methylene chloride	ND	2.8	209	206	ND	18.0	98	45.0	18.0	89
23V	1,1,2,2-Tetrachloroethane	ND	6.9	ND	ND	ND	18.0	101	ND	18.0	106
24V	Tetrachloroethylene	ND	4.1	ND	ND	ND	18.0	98	ND	18.0	103
25V	Toluene	ND	6.0	ND	ND	ND	18.0	100	ND	18.0	102
26V	1,2-Trans-dichloroethylene	ND	1.6	ND	ND	ND	18.0	97	ND	18.0	99
27V	1,1,1-Trichloroethane	ND	3.8	ND	ND	ND	18.0	104	ND	18.0	97
28V	1,1,1,2-Trichloroethane	ND	5.0	ND	ND	ND	18.0	104	ND	18.0	100
29V	Trichloroethylene	ND	1.9	ND	ND	ND	18.0	93	ND	18.0	87
30V	Trichlorofluoromethane	ND	10	ND	ND	ND	18.0	105	ND	18.0	99
31V	Vinyl chloride	ND	10	ND	ND	ND	18.0	110	ND	18.0	99
18V	trans-1,3-Dichloropropylene	ND	10	ND	ND	ND	18.0	94	ND	18.0	96

<sup>a</sup> Recovery normally variable using established methodology.

305110



805111

**ETC** ENVIRONMENTAL TESTING and CERTIFICATION

NOV 29, 1986

### TABLE 1: QUANTITATIVE RESULTS Metals - Analysis Data (QR52)

Chain of Custody Data Required for ETC Data Management Summary Reports

P4850 ENVIRONMENTAL STRATEGIES CORP CHAMPION WCSP-1 861110 1100

ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

Compound	Results	
	Sample Concn. ug/l	MDL ug/l
Barium	32	3.2
Chromium	ND	26
Copper	BMDL	13
Iron	BMDL	140
Manganese	BMDL	6.4
Nickel	ND	11
Sodium	9110	72
Zinc	29	2.9

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Conventional Analysis Data (QR10)**

Chain of Custody Data Required for ETC Data Management Summary Reports

P4850	ENVIRONMENTAL STRATEGIES CORP. CHAMPION	WCSF-1	861110	1100
ETC Sample No.	Company	Facility	Sample Point	Date Time Elapsed Hours

NPDES Number	Results		Sample Concn. mg/l	MDL mg/l	Facility	Sample Point	Date	Time	Elapsed Hours
	Sample Concn. mg/l	MDL mg/l							
Fluoride Nitrate as N Sulfate as SO4 Phenolics, Total Cyanide, Total Solids, total suspended	16	.10							
	5.50	.10							
	85	2							
	<.050	.050							
	490	.025							

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ENVIRONMENTAL TESTING and CERTIFICATION

NOV 25, 1986

**TABLE 2: METHOD PERFORMANCE DATA**  
**Surrogate Recovery - Aqueous Matrices (QR20)**

Chain of Custody Data Required for ETC Data Management Summary Reports

P4850	ENVIRONMENTAL STRATEGIES CORP.	CHAMPION	WCSP-1	86111	1100	0
ETC Sample No.	Company	Facility	Sample Point	Date	Time	Elapsed Hours

Compound	Amount added	% Recovery	Control Limits	
			Lower	Upper
<b>VOLATILE FRACTION (GC/MS)</b>				
Toluene-D8	.250	100	88	110
p-Bromofluorobenzene	.250	104	86	115
1,2-Dichloroethane-D4	.250	108	76	114
<b>ACID FRACTION (GC/MS)</b>				
Phenol-D5	-	-	10	94
2-Fluorophenol	-	-	21	100
2,4,6-Tribromophenol	-	-	10	123
<b>BASE/NEUTRAL FRACTION (GC/MS)</b>				
Nitrobenzene-D5	-	-	35	114
2-Fluorobiphenyl	-	-	43	116
Terphenyl-D14	-	-	33	141
<b>PESTICIDE/PCB FRACTION (GC)</b>				
Dibutylchloride	-	-	24**	154**

\* 17B EPA Control Limits  
 \*\* Mercury Limits Only



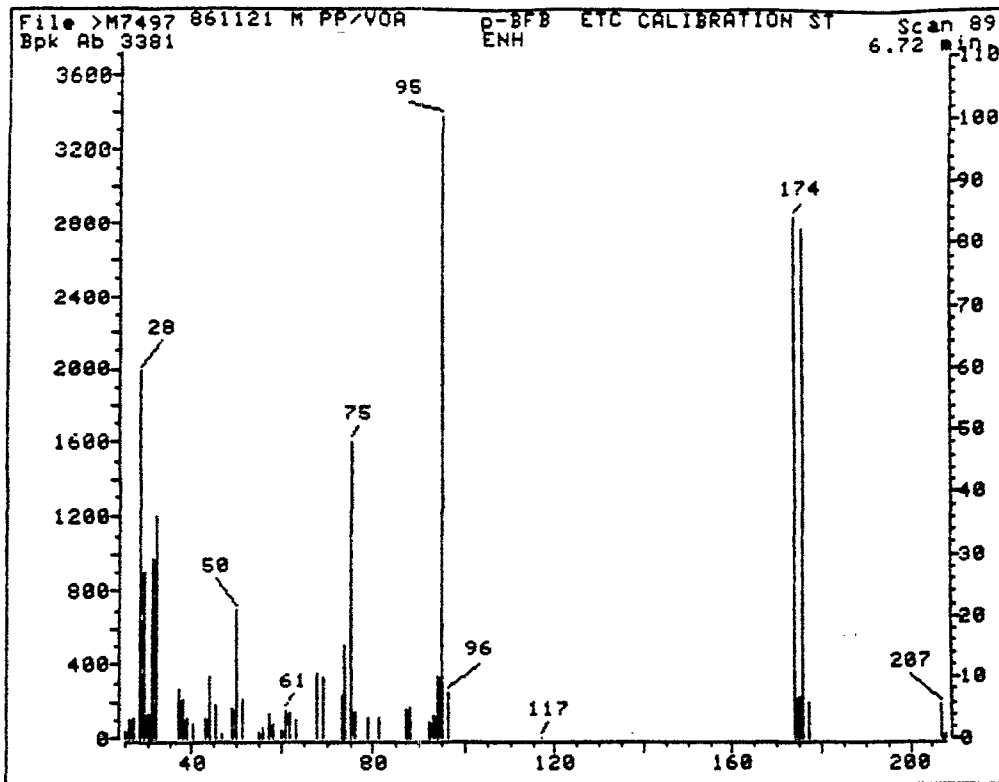


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	20.80	20.80	Ok
75	30-60% of mass 95	47.52	47.52	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.18	7.18	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	83.99	83.99	Ok
175	5-9% of mass 174	6.60	7.85	Ok
176	95-101% of mass 174	82.16	97.82	Ok
177	5-9% of mass 176	5.86	7.13	Ok

Injection Date: 11/21/86  
 Injection Time: 05:57  
 Run No: >M7497  
 Spectrum No: 89

Analyst: Joseph O...  
 Processor: (unreadable)  
 QC Batch: QV5931  
 Samples: P4240, P4241, P4242, P4243, P4236, P4237, P4239, P4235, P5415, P5414, P4850, P4851, P...

31720  
 305115

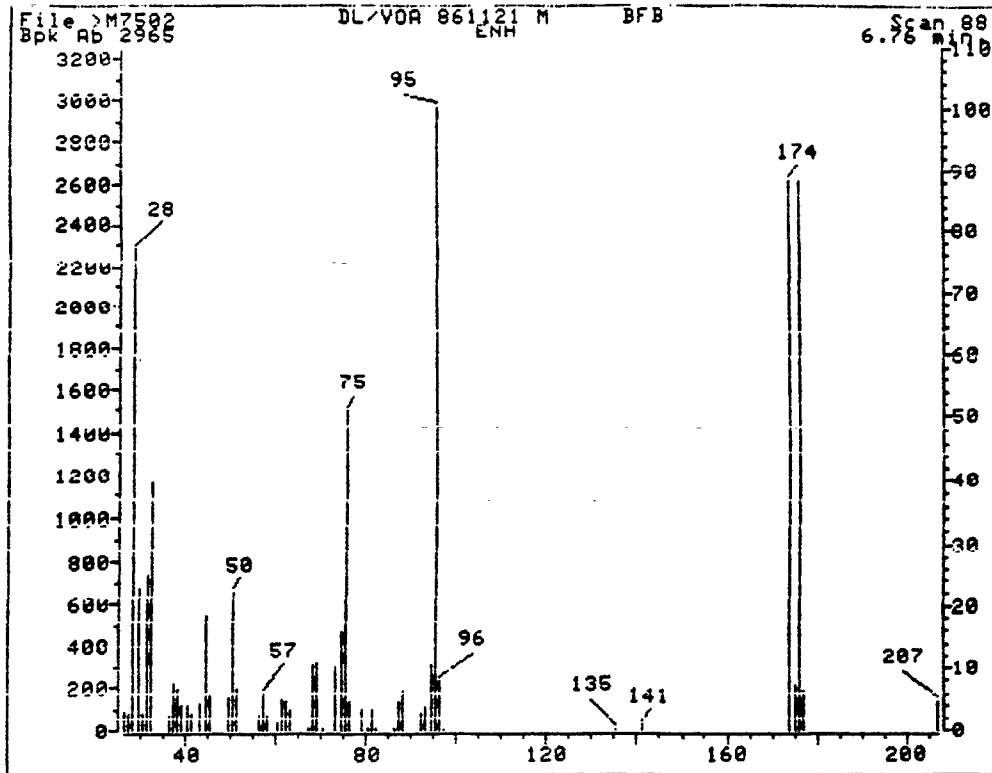


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	22.00	22.00	Ok
75	30-60% of mass 95	50.62	50.62	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.63	7.63	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	88.52	88.52	Ok
175	5-9% of mass 174	7.03	7.95	Ok
176	95-101% of mass 174	88.43	99.89	Ok
177	5-9% of mass 176	6.31	7.75	Ok

Injection Date: 11/21/86  
 Injection Time: 09:32  
 Run No: >M7502  
 Spectrun No: 88

Analyst: Nancy Albert  
 Processor: Quinn Christopher  
 QC Batch: QV 5931  
 Samples: P4240, P4241, P4242, P4243, P4244  
P4236, P4237, P4238, P4239, P533  
P5415, P5414, P4850, P4851, P4852

## Methodology for GC/MS Analysis of Priority Pollutant Compounds

The methods employed in the GC/MS analysis for priority pollutants are established EPA methods. Rigid compliance with the instrument parameters and performance criteria of the published methods was achieved. In some cases, the precise amounts of sample used and the sample handling procedures vary with the complexity of the sample matrix. Qualitative identification of the priority pollutants was performed using the relative retention times, the relative abundance of three characteristic ions and the abundance ratios. The entire mass spectrum was reviewed to confirm each identification. Quantitative analysis of detected compounds was performed by using a response factor generated by a major characteristic ion of the specific compound and an internal standard.

Compounds, in addition to those on the priority pollutant list, were identified through a computer-aided search of the NBS-EPA spectra library. After review the identifications are included in a separate tabulation and labelled "tentatively identified".

### *Volatile Priority Pollutant Compounds*

For the analysis of Volatile priority pollutants, EPA Method 624 was used. The method can be summarized as follows: Helium is bubbled through a 5 ml water sample contained in a specially designed purging chamber at ambient temperature. The purgeable volatile organic compounds are transferred from the aqueous phase to the vapor phase. The vapor is swept through a sorbent column where the organic components are trapped. After the purge cycle is complete, the sorbent column is heated and backflushed with helium to desorb the organic purgeables onto a gas chromatographic column. The gas chromatograph is temperature programmed to separate the purgeable mixture. The separated purgeable components are then identified and quantitated using a computerized mass spectrometer.

### *Acid, Base/Neutral and Pesticide Priority Pollutant Compounds*

For the analysis of the Acid, Base/Neutral and Pesticide priority pollutants in an aqueous liquid matrix, EPA Method 625 was used. The method can be summarized as follows: A measured volume of sample, approximately 1 liter, is adjusted to a pH greater than 11 and extracted with methylene chloride. The pH of the sample is adjusted to a value less than 2 and extracted with an aliquot of fresh methylene chloride. A separatory funnel or continuous extractor is used to perform the extractions. The two extracts are dried and concentrated to a 1 ml final volume. Each extract is injected into a GC/MS instrument specifically configured for the correct fraction.

## Methodology for Analysis of Metals

### **AQUEOUS**

The determination of metals in aqueous samples is performed according to the methods published by EPA in "Methods for Chemical Analysis of Water and Wastes," EPA-600/4-79-020, March, 1983, and the Federal Register, October 26, 1983. Arsenic, selenium and thallium are determined by furnace AA; silver, aluminum, barium, beryllium, boron, cadmium, calcium, chromium, copper, cobalt, iron, magnesium, manganese, molybdenum, nickel, lead, sodium, antimony, tin, titanium, vanadium, and zinc are determined by ICP emission spectrometry, except where lower levels of detection are required; in these cases (e.g. lead in groundwater monitoring samples) furnace AA is used. All furnace AA parameters are run by method of standard additions. The determination of mercury is performed by cold vapor AA.

### **EP TOXICITY**

The determination of metals in aqueous EP Toxicity leachates is performed according to the methods published by EPA in "Test Methods for Evaluating Solid Waste" EPA SW-846, revised April, 1984 and the Federal Register, Oct. 26, 1983, 1979. Silver, arsenic, barium, cadmium, chromium, lead and selenium are determined by ICP emission spectrometry. Mercury is determined using cold vapor AA. For leachates that are organic in nature, the analyses are performed according to the methods described under **OIL/SLUDGE** below.

### **SOIL/SEDIMENT**

The determination of silver, beryllium, cadmium, chromium, copper, nickel, antimony, lead, and zinc in sediment samples is performed according to methods published by EPA in "Interim Methods for the Sampling and Analysis of Priority Pollutants in Sediments and Fish Tissue", EPA 600/4-81-055, October 1980. Mercury is determined according to the sediment method published by EPA in "Method for Chemical Analysis of Water and Wastes", EPA 600/4-79-020, March 1983. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion.

### **OIL/SLUDGE**

The determination of silver, aluminum, boron, barium, beryllium, calcium, cadmium, copper, chromium, cobalt, iron, magnesium, manganese, molybdenum, sodium, nickel, lead, antimony, tin, titanium, vanadium, and zinc in sludge/petroleum-based samples is performed by ICP emission spectrometry using a magnesium nitrate dry ashing digestion technique. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion. Mercury is determined by cold vapor AA using the same digestion technique.



## Summary of Quality Assurance/Quality Control Procedures (QA/QC)

ETC bases its quality assurance protocols on the following government guidelines:

- . "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-330/9/79/001-R, October 1979;
- . the recommended guidelines for EPA Methods 624 and 625. (Federal Register, December 3, 1979, updated on October 26, 1984);
- . "Manual of Analytical Methods for the Analysis of Pesticides in Humans and Environmental Samples," EPA 600/8-80-038, June 1980;
- . "Determination of 2,3,7,8-TCDD in Soil and Sediment" EPA, Region VII, Kansas City, September 1983;
- . Organic Analysis: Multi-media, Multi Concentration-IFB WA84-A267; and
- . Dioxin Analysis: Soil/Sediment Matrix; Multi-Concentration; Selected Ion Monitoring with Jar Extration Procedure-IFB WA84-A002

However, we have modified our protocols to provide a higher level of QA/QC than the guidelines require. For example, we analyze a higher than required number of quality control samples and we pay especially careful attention to the certification of the "reference standard" compounds we use in analysis. Below are listed the key QA/QC elements for the methods we used.

### Analysis of Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry

- Each batch of 13 samples consists of 9 customer samples (at a maximum), one blank sample, one spiked blank, one spiked sample and one replicate sample. This amounts to a 30% quality control factor.
- Three surrogate compounds are added to each sample in the batch of 13.
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and retuned, if necessary, at the beginning of each day to ensure that its performance on bromofluorobenzene (BFB) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of Volatile Organic Priority Pollutant "standards" at a minimum of 3 different concentrations and using a mixture of 3 internal standards at a constant concentration.
- The calibration curve is verified with a mixture of priority pollutant standards every day. If the response factors vary greater than 25%, the instrument must be recalibrated.

### Analysis of Organic Compounds Extracted in Acid or Base/Neutral Solutions by Gas Chromatography/Mass Spectrometry

- Each batch of 20 samples consists of 16 customer samples (at a maximum), one blank sample, one spiked blank (for water matrices), one sample spiked with the priority pollutant standard mixture and a duplicate customer sample. This amounts to a 20% quality control factor.

- Three surrogate compounds are added to each sample in the batch for Base/Neutral analysis.
- Three surrogate compounds are added to each sample in the batch for Acid analysis.
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and retuned, if necessary, at the beginning of each day to ensure that its performance on decafluorotriphenylphosphine (DFTPP) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of standards composed of either the Organic Acid or Base/Neutral Extractable Compounds at a minimum of 3 concentrations and using five internal standards for quantitation.

### Analysis of Metals

#### All Samples

- New standards are prepared for each batch of samples
- Normal calibration is performed using a blank sample and four standards that have been through the sample preparation procedure. A regression analysis is used to construct the calibration curve.
- All EP Toxicity samples and all samples determined by furnace atomic absorption are calculated by the "method of additions".
- For each sample analysis that requires the use of the "method of additions" technique, a three point calibration is performed using U.S. EPA "Methods for Chemical Analysis of Water and Wastes, 1979". Results are obtained using linear regression analysis. Any regression with a coefficient of correlation below 0.990 is considered suspect, necessitating review of calibration data or sample re-analysis.
- In constructing the normal calibration curves the lowest concentration levels we use are values greater than or equal to 5 times the Instrumental Detection Limit (IDL).
- All calibration standards are analyzed in duplicate, at a minimum.
- Independent reference standards are used to check the accuracy of calibration standards.
- A check standard is analyzed every ten samples to validate the normal calibration curve.
- One customer sample out of every ten is analyzed in triplicate.

#### Homogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are homogeneous, the QC program is a minimum of 25% and consists of analyzing:

- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 1 independent reference standard;
- 4 Calibration standards (processed using the sample preparation method);
- 4 Calibration standards (w/ out sample preparation); and
- 1 Reagent Blank.

Heterogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are heterogeneous, the QC program is a minimum of 35% and consists of analyzing

- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 1 Replicate independent reference standards;
- 4 Calibration standards (processed using the sample preparation method);
- 1 Procedural Blank;
- 4 Calibration standards (without sample preparation); and
- 1 Reagent Blank.

Analysis of Mercury

To analyze samples for mercury we group them by matrix in batches of 30 or less. Our QC program is a minimum of 30% and consists of analyzing:

- each of the 30 customer samples in duplicate;
- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 2 Replicate independent reference standards;
- 10 Calibration standards (processed using the sample preparation method); and
- 2 Procedural Blanks.

Analysis of Pesticides, Herbicides and PCB's by Gas Chromatography

Pesticide, herbicide and PCB samples are grouped in batches of 16 customer samples or less according to the type of analysis to be performed. The QC program for each of these three types of analyses is a minimum of 20% and consists of analyzing:

- 1 procedural blank sample (a reagent blank is analyzed in the case of non-water matrices);
- 1 spiked blank sample (the spiked blank is eliminated in the case of non-water matrices);
- 1 replicate sample;
- 1 replicate spiked sample; and
- 1 known reference QC sample for at least each 100 samples analyzed.

The instrument is calibrated each run with three standards, and checked every 10 samples.

Analysis of Cyanides, Phenols, Fluoride, Chloride, Nitrate and Nitrite

- All parameters are analyzed using a Technicon Autoanalyzer II GT.
- 3 calibration standards are analyzed at the beginning and end of each batch.

- Each batch (up to 80 samples) consists of analyzing one blank, one spiked blank, one duplicate and spiked sample every 20 samples, and an EPA known reference sample.

#### Analysis of Total Organic Carbon (TOC)

TOC samples are analyzed on a daily basis with the number of samples analyzed per day dependent on the request for duplicate or quadruplicate analyses. The quality control program is designed to maintain the appropriate amount of QC and consists of the following elements:

- Daily instrument calibration
- One blank
- Standard recalibration every 10 samples
- Spiked samples at a low and high level
- Every sample is run in duplicate at a minimum

#### Analysis of Total Organic Halide (TOX)

- Blank reagent water for absolute carbon background must contain less than 5 ug/l of halide (as chloride).
- Using a trichlorophenol standard, the mean adsorption efficiency must be within +/- 15% of the standard value.
- Calibration standards are run every 10 samples.
- Every sample is run in duplicate at a minimum.

#### Analysis of 2,3,7,8-TCDD (Dioxin) by GC/MS (SIM)

- Each sample is dosed with a known quantity of  $^{13}\text{C}_{12}$ -2,3,7,8-TCDD as internal standard and  $^{37}\text{Cl}_4$ -TCDD as surrogate standard. The action limits for surrogate standard results is +/- 40% of the true value. Samples showing surrogate standard results outside of these limits are reextracted and reanalyzed.
- Two laboratory "method blanks" are run along with each set of 24 or fewer samples. The method blank is also dosed with the internal standard and surrogate standard
- At least one per set of 24 samples is run in duplicate to determine intralaboratory precision.
- Qualitative Requirements. The following are met in order to confirm the presence of native 2,3,7,8-TCDD:
  - a. Isomer specificity must be demonstrated initially and verified once per 8-hour work shift. The verification consists of injecting a mixture containing TCDD isomers which elute close to 2,3,7,8-TCDD. The 2,3,7,8-TCDD must be separated from interfering isomers, with no more than 25% valley relative to the 2,3,7,8-TCDD peak.
  - b. The 320/322 ratio is within the range of 0.67 to 0.87.
  - c. Ions 320, 322, and 257 are all present and maximize together the signal to mean noise ratio must be 2.5 to 1 or better for all 3 ions.
  - d. The retention time is equal (within 3 seconds) the retention time for the isotopically labeled 2,3,7,8-TCDD.
  - e. At least one of the positives can be confirmed by obtaining partial scan spectra from mass 150 to mass 350. The partial scan guidelines are as follows:

- . the 320/324 ratio should be 1.58 +/- 0.16
- . the 257/259 ratio should be 1.03 +/- 0.10
- . the 194/196 ratio should be 1.54 +/- 0.15
- One sample is spiked with native 2,3,7,8-TCDD at a level of 1.0 PPB (for soil) for each set of 24 or fewer samples.
- In cases where no native 2,3,7,8-TCDD is detected, the actual detection limit is estimated and reported based on a signal to noise ratio of 2.5 to 1 at ions 320 and 322.
- For each sample, the internal standard is present with at least a 10 to 1 signal to noise ratio for both mass 332 and mass 334. Also, the internal standard 332/334 ratio must be within the range of 0.67 to 0.87.

#### Subcontractor QA/QC

Each subcontractor is required to maintain an appropriate level of quality control. To insure this, each subcontractor is required to submit to ETC the quality control data for all analyses it performs. This data is kept on file at ETC. In general, the amount of quality control required is one duplicate sample with one spiked sample for every ten analyses.

#### Chain-of-Custody

The chain-of-custody procedure is part of our quality assurance protocol. We believe our chain-of-custody record fully complies with the legal requirements of federal, state and local government agencies and of the courts of law. The record covers:

- labeling of sample bottles, packing the Sample Shuttle and transferring the Shuttle under seal to the custody of a shipper;
- outgoing shipping manifests;
- the chain-of-custody form completed by the person(s) breaking the Shuttle seal, taking the sample, resealing the Shuttle and transferring custody to a shipper;
- incoming shipping manifests;
- breaking the Shuttle's reseat;
- storing each labeled sample bottle in a secured area;
- disposition of each sample to an analyst or technician; and
- the use of the sample in each bottle in a testing procedure appropriate to the intended purpose of the sample.

The records show for each link in this process:

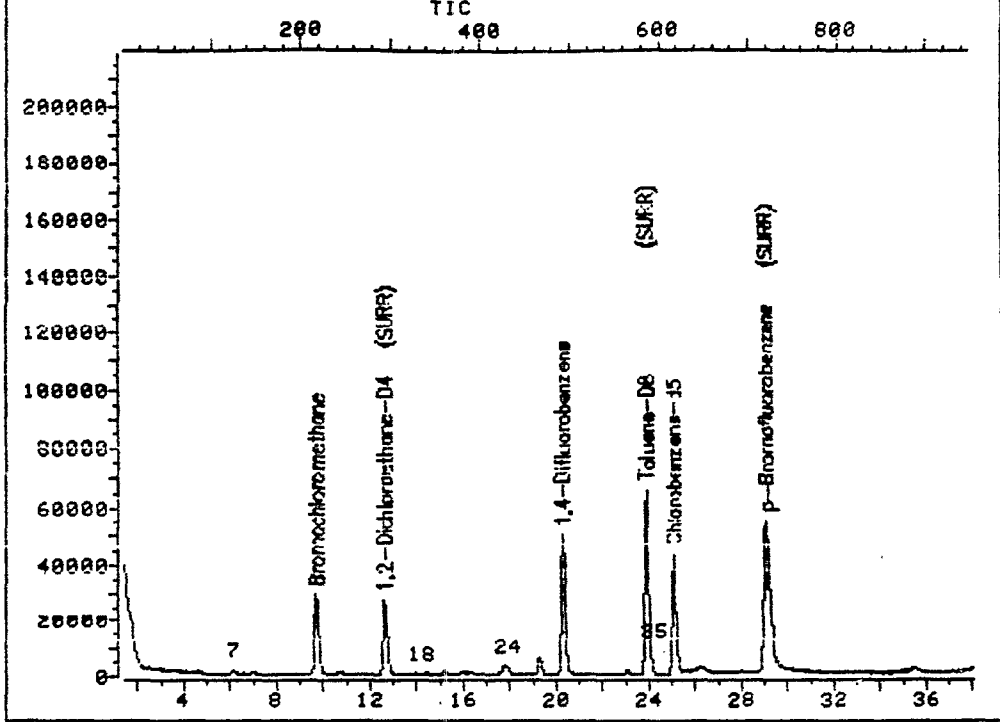
- the person with custody; and
- the time and date each person accepted or relinquished custody.

**Appendix A**  
**Mass Spectral Data**  
**for**  
**Quantitated Compounds**

- 1) A total ion chromatogram for each sample analyzed by a GC/MS instrument.
- 2) A Quant. report used by the analyst to determine qualitative and quantitative results of the compounds present.
- 3) A mass spectrum and a reference spectrum for each priority pollutant compound detected in the sample.

TOTAL ION CHROMATOGRAM

File >M7520 40.0-265.0 amu. PP/VOA 861121 M P4850V, QV5931, L, 5, 5,



Data File: >M7520::U4  
Name: PP/VOA 861121 M  
Misc: P4850V, QV5931, L, 5, 5,

Quant Output File: ^M7520::AQ

Id File: MVOA::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861121 12:01

Operator ID: DG1199  
Quant Time: 861122 04:05  
Injected at: 861122 03:26

305125

QUANT REPORT

Operator ID: DG1199  
 Output File: ^M7520::AQ  
 Data File: >M7520::U4  
 Name: PP/VOA 861121 M  
 Misc: P4850V,QU5931,L,5,5,

Quant Rev: 6      Quant Time: 861122 04:05  
 Injected at: 861122 03:26  
 Dilution Factor: 1.00000

ID File: MVOA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861121 12:01

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *Bromochloromethane	9.73	216	61033	250.00	NG	98
<del>7) Methylene chloride</del>	<del>6.13</del>	<del>123</del>	<del>3022</del>	<del>22.80</del>	<del>NG</del>	<del>96</del>
16) *1,4-Difluorobenzene	20.34	490	225753	250.00	NG	81
17) 1,2-Dichloroethane-D4 (SUHR)	12.67	292	53629	269.48	NG	94
<del>18) 1,1,1-Trichloroethane</del>	<del>14.83</del>	<del>327</del>	<del>1214</del>	<del>2.95</del>	<del>NG</del>	<del>98</del>
<del>24) Benzene</del>	<del>17.78</del>	<del>424</del>	<del>8248</del>	<del>9.87</del>	<del>NG</del>	<del>85</del>
30) *Chlorobenzene-d5	25.14	614	152293	250.00	NG	95
34) Toluene-D8 (SUHR)	23.98	584	275620	249.60	NG	94
<del>35) Toluene</del>	<del>24.13</del>	<del>588</del>	<del>3027</del>	<del>2.78</del>	<del>NG</del>	<del>91</del>
38) p-Bromofluorobenzene (SUHR)	29.17	718	163579	258.87	NG	98

\* Compound is ISTD

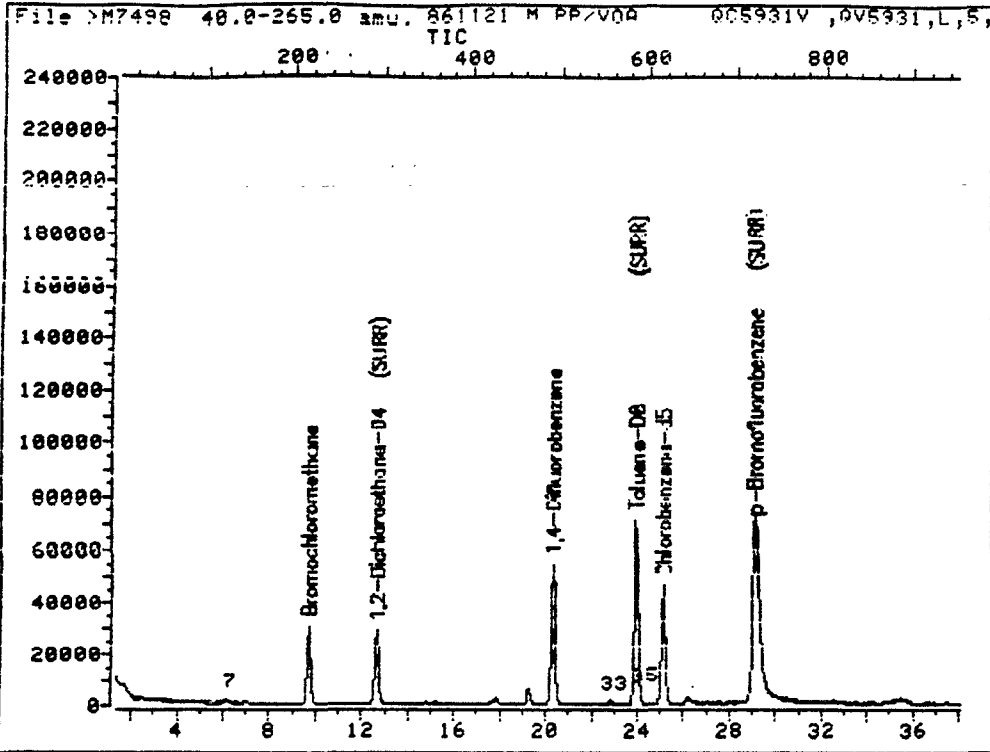
SNC 11/24/86



**Appendix C1**  
**GC/MS Subsidiary Data**

305127

TOTAL ION CHROMATOGRAM



Data File: >M7498::U3  
Name: 861121 M PP/VDA  
Misc: Q05931V ,Q05931,L,5,5,

Quant Output File: ^M7498::AU

Id File: M00A::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861121 12:01

Operator ID: JQ6275  
Quant Time: 861121 12:03  
Injected at: 861121 06:18

QUANT REPORT

Operator ID: JW6275  
 Output File: ^M/498::AQ  
 Data File: >M/498::U3  
 Name: 861121 M PP/VOA  
 Misc: QC5931U ,WV5931,L,5,5,

Quant Rev: 6      Quant Time: 861121 12:03  
 Injected at: 861121 06:18  
 Dilution Factor: 1.00000

ID File: MVUA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861121 12:01

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *Bromochloromethane	9.70	217	66376	250.00	NG	99
7) Methylene chloride	6.18	126	2829	19.63	NG	86
16) *1,4-Difluorobenzene	20.31	491	249275	250.00	NG	82
17) 1,2-Dichloroethane-D4 (SURR)	12.65	293	57076	259.74	NG	94
30) *Chlorobenzene-d5	25.11	615	170859	250.00	NG	99
<del>33) Tetrachloroethylene</del>	<del>22.83</del>	<del>556</del>	<del>2829</del>	<del>6.88</del>	<del>NG</del>	<del>94</del>
34) Toluene-D8 (SURR)	23.95	585	307632	248.31	NG	91
<del>35) Toluene</del>	<del>24.15</del>	<del>598</del>	<del>1931</del>	<del>1.58</del>	<del>NG</del>	<del>92</del>
38) p-Bromofluorobenzene (SURR)	29.10	718	181687	256.29	NG	97

\* Compound is ISTD

ENC 1124186

**Appendix D**  
**Subcontractor's Data**

- 1) A copy of the originating subcontractor's report is included for all data not generated within ETC's laboratory.



## Appendix E

### Chain-of Custody Forms

- 1) A field Chain-of-Custody form (CC1) is included for all samples shipped by ETC shuttle.
- 2) An in-house sample Chain-of-Custody form is included for all samples not shipped by ETC shuttle.
- 3) Any additional Chain-of-Custody material provided by a client or by a client's sampling agent is also included.
- 4) A subcontractor's Chain-of-Custody form is included for any analytical work not performed within ETC's laboratory.
- 5) Analysis and Extraction Custody forms are included for the period the sample was in ETC's possession.

**CHAIN OF CUSTODY FORM (CC1)**

Seal No. 67753 ETC Job # P4850

Date Sealed 6 NOV. 1986 By: M. I. /

Company: HELLERTOWN MANUFACTURING

Attn.: MARK SCHULTZ

Facility/Site: C/O GUARD OFFICE-SILVER ROAD

Phone: \_\_\_\_\_

Address: HELLERTOWN, PA. 18055

**SAMPLE IDENTIFICATION**

Facility: CHAMPAIGN (Optional Sample Point Descriptions)

Sample Point: M-CSP-1 11110186 1100 1100  
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.	Filt. (Y/N)	Observations
1	M	500 ML	HN03	METALS		/
1	PN	125 ML	H2S04	PHENOLS		VERY LOW SAMPLE VOLUME
1	B	125 ML	H2S04	NITRATE		/
1	CN	125 ML	NADH	CYANIDE		/
1	A	500 ML	BAKED DI	SOLIDS, TD/SULFATE		/
1	D	500 ML	RINSE SOD.	FLUORIDE		/
2	U	40 ML	THIOL GC/MS	VOA		/
1	TB	40 ML	H2O	TRIP BLANK		/

**CHAIN OF CUSTODY CHRONICLE**

1. Shuttle Opened By: (print) M. SCHULTZ Date: 11-10-86 Time: \_\_\_\_\_  
 Signature: [Signature] Seal #: \_\_\_\_\_ Intact: /

2. I have received these materials in good condition from the above person.  
 Name: \_\_\_\_\_ Signature: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_ Remarks: \_\_\_\_\_

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

4. Shuttle Sealed By: (print) M. SCHULTZ Date: 11-10-86 Time: 16:30  
 Signature: [Signature] Seal #: 67754 Intact: \_\_\_\_\_

LAB USE ONLY Opened By: Mark Schadt Date: 11/11/86 Time: 5:22 PM

SHUTTLE # 381 TEMP. °C 12 SEAL # 67754 COND INTACT

# 805133









SUBCONTRACT  
REQUEST FOR ANALYSIS  
and SAMPLE CHAIN OF CUSTODY

LOGLINK: 17866  
MATRIX: WATER

NAME OF SUBCONTRACT LAB: CHYUN

ETC JOB NUMBERS: P4850 P4853  
P4851  
P4852

TURNAROUND IN DAYS: NORMAL DATE DATA REQUIRED: 11/22  
(If deadline cannot be met, contact ETC Subcontract Group).

Send invoice, bill and reports to : ETC Subcontract Group  
(201)225-6786

-----  
Please perform the analyses requested below:

<input type="checkbox"/> Acidity	<input type="checkbox"/> Alkalinity	<input type="checkbox"/> Ammonia(probe)
<input type="checkbox"/> Ammonia(dist)	<input type="checkbox"/> Bicarbonate	<input type="checkbox"/> BOD(5 day)
<input type="checkbox"/> Bromide	<input type="checkbox"/> Carbonate	<input type="checkbox"/> Chloride
<input type="checkbox"/> Chem.Oxygen Dem.	<input type="checkbox"/> Chromium+6	<input type="checkbox"/> Fecal Coliform
<input type="checkbox"/> Total Coliform	<input type="checkbox"/> Color,apparent	<input type="checkbox"/> Cyanide,total
<input type="checkbox"/> Fluoride	<input type="checkbox"/> Formaldehyde(UV)	<input type="checkbox"/> Hardness
<input type="checkbox"/> Nitrate(NO3)	<input type="checkbox"/> Nitrite(NO2)	<input type="checkbox"/> TKN
<input type="checkbox"/> Odor	<input type="checkbox"/> Oil & Grease(grav)	<input type="checkbox"/> T.Organic Carb.
<input type="checkbox"/> Pet.Hydro(IR)	<input type="checkbox"/> Pet.Hydro(grav)	<input type="checkbox"/> Phenolics,tot.
<input type="checkbox"/> Phenolics(5ug/l)	<input type="checkbox"/> Phosphate(ortho)	<input type="checkbox"/> Phosphate(total)
<input type="checkbox"/> Phosphorus(tot.)	<input type="checkbox"/> Silica(dissolved)	<input checked="" type="checkbox"/> Sulfate(SO4)
<input type="checkbox"/> Sulfide(S)	<input type="checkbox"/> Sulfite(SO3)	<input type="checkbox"/> Surfactant(MBAS)
<input type="checkbox"/> Solids,Total	<input type="checkbox"/> Solids,Tot. Diss.	<input type="checkbox"/> Solids,Tot.Set.
<input checked="" type="checkbox"/> Solids,Tot. Sus.	<input type="checkbox"/> Solids,Tot. Vol.	<input type="checkbox"/> Turbidity
<input type="checkbox"/> Gross Alpha,Beta	<input type="checkbox"/> Radium 226	<input type="checkbox"/> Radium 228

Others: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

-----  
Chain of Custody Section for Courier Pick up at ETC

Sample(s) Relinquished by:(ETC) Mark Schmitt  
Time: 4:30pm Date: 11/2/86  
Sample(s) Received by: Mark Kelly  
Time: 4:30pm Date: 11/2/86

Chain of Custody Section for Sending Sample from ETC

Sample Shuttle sealed by : \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_  
Sample Shuttle opened by : \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_  
Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

Chain of Custody Section for Returning Sample to ETC

Sample Shuttle sealed by : \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_  
Sample Shuttle opened by : \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_  
Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

305137



TEST: NITRATE UNITS: mg/L  
 Meth. REF. # 353-2 MDL: 0.10  
 Instrument # #1 (Channel 1)

BY: CH  
 BOOK # CV-12 P. 97  
 FROM PAGE # 96

Log/Lab	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	Log/Lab	Job #	ANALYZED CONC.	DILN	REPORTED CONC.
7514	N6032	.334	5x	1.67	17866	P4851	.793	10x	7.93
7514	N6033	.0134	5x	Redo	17866	P4852	.0487	10x	
7562	N6122	.804	5x	4.02	17866	P4853	.442	10x	4.42
7503	P2683	.0213	5x	Redo	17866	P4854	.0482	10x	
7798	P2469	.276	10x	2.76	17866	P3117	.1153	10x	1.153
7798	P2477	1.54	10x	15.4	17866	P3120	.0154	10x	
17798	P2524	.517	10x	5.17	17866	P3122	.0211	10x	
17798	P2538	.583	10x	5.83	17866	P3124	.0224	10x	
17796	P2539	.856	10x	8.56	17866	P3126	.0176	10x	
17798	P2540	.618	10x	6.18	17866	P2543	.528	10x	5.28
17814	P3555	.101	10x		17866	P2544	.0253	10x	
17822	P3107	.417	10x	4.17	17866	P2545	1.33	10x	13.3
17822	P3112	.0139	10x		17866	P2547	1.44	10x	14.4
17828	P2473	.442	10x	4.42	17866	P2549	.924	10x	9.24
17828	P2472	.547	10x	5.47	17866	P2550	.401	10x	4.01
17828	P2522	.417	10x	4.17	17866	P2551	1.03	10x	10.3
17828	P2536	.435	10x	4.35	17866	P2552	.717	10x	7.17
17828	P2548	.525	10x	5.25					
17864	P2481	.0142	10x						
17864	P2521	.0194	10x						
17864	P2541	1.60	10x	1.60					
17864	P2546	.694	10x	6.94					
17864	P2557	.198	10x	1.98					
17866	P4850	.550	10x	5.50					

STANDARDS		
STANDARD mg/L	VALUE OBTAINED	% of Theoretical
PREPARED ON _____		
BY _____		
FROM _____ mg/L		
STOCK # _____		
SLOPE _____		
Intercept _____		
Corr. Coef. _____		

**ADDITIONAL COMMENTS**  
 Sample Red Line Through  
 Samples Done on Channel  
 2. No Dilution Page 98  
 Ref N6033 } Sample  
 P2683 }

Analyses QA/QC Data Report

QC Batch # N3003 & N3004

Verified: 11/24/86 *[Signature]*

METHOD BLANK		PRECISION	
Spiked Blank (Known)		Job # Dup (ETC)	
Spiked Blank Result		Original Result	
% Recovery		Duplicate Result	
		R Range	
		<b>ACCURACY</b>	
		Job # Spiked (ETC)	
		Original Result	
		Amount Spiked	
		Spiked S Result	
		% Recovery	

**EXTERNAL REF. QC**  
 Known Value \_\_\_\_\_  
 Observed Result \_\_\_\_\_  
 % Recovery \_\_\_\_\_

TEST: NITRATE UNITS: mg/l  
 METH. REF. # 353.2 MDL: 0.10  
 INSTRUMENT # #1 (Channel 2)

BY: CH  
 BOOK # CV-12 P. 98  
 FROM PAGE # 3

Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	STANDARDS															
7798	P2469	OVER		*	17886	P3117	1.15		1.15	PREPARED ON _____ BY _____ FROM _____ mg/L STOCK # _____															
7798	P2471	OVER		*	17866	P3120	.0207		BMDL																
7798	P2524	OVER		*	17866	P3122	.0511		BMDL																
7798	P2538	OVER		*	17866	P3124	.0173		BMDL																
7798	P2539	OVER		*	17866	P3126	.0127		BMDL																
788	P2540	OVER		*	17882	P2543	OVER		*	<table border="1"> <thead> <tr> <th>STANDARD mg/L</th> <th>VALUE OBTAINED</th> <th>% of Threshold</th> </tr> </thead> <tbody> <tr> <td>2.00</td> <td>1.98</td> <td></td> </tr> <tr> <td>1.00</td> <td>1.04</td> <td></td> </tr> <tr> <td>.50</td> <td>.497</td> <td></td> </tr> <tr> <td>.20</td> <td>.0934</td> <td></td> </tr> </tbody> </table>	STANDARD mg/L	VALUE OBTAINED	% of Threshold	2.00	1.98		1.00	1.04		.50	.497		.20	.0934	
STANDARD mg/L	VALUE OBTAINED	% of Threshold																							
2.00	1.98																								
1.00	1.04																								
.50	.497																								
.20	.0934																								
7814	P3555	.487		.49	17882	P2544	.0250		BMDL																
7822	P3107	OVER		*	17882	P2545	OVER		*																
17822	P3112	.007		BMDL	17882	P2547	OVER		*																
17828	P2473	OVER		*	17882	P2549	OVER		*																
7828	P2472	OVER		*	17882	P2550	OVER		*																
17828	P2522	OVER		*	17882	P2551	OVER		*																
17828	P2536	OVER		*	17882	P2552	OVER		*																
17828	P2548	OVER		*																					
17864	P2481	.006		BMDL																					
17864	P2531	.0245		BMDL																					
17864	P2541	1.73		1.73																					
17864	P2546	OVER		*																					
17864	P2554	2.06		2.06																					
17866	P4850	OVER		*																					
17866	P4851	OVER		*																					
7866	P4852	.398		.40																					
7866	P4853	OVER		*																					
7878	P2770	.0462		BMDL																					

ADDITIONAL COMMENTS  
 \* RPT on Channel 1  
 complete Page 97

Analyses QA/QC Data Report

QC Batch # NCO03

Verified: 11/28/86

METHOD BLANK	.003	.0177
Spiked Blank (Known)	.5	.5
Spiked Blank Result	.801	.532
% Recovery	160%	1.06

INTERNAL REF. QC	EPA	
Known Value	.93	
Observed Result	.953	
% Recovery	1.02	

PRECISION	P2473	P4850	P3126
Job # Dup (ETC)	P2473	P4850	P3126
Original Result	OVER	OVER	.0217
Duplicate Result	↓	↓	.0194
A Range			
ACCURACY	P2473	P4850	P3126
Job # Spiked (ETC)	P2473	P4850	P3126
Original Result	.050	OVER	.021
Amount Spiked			.5
Spiked B Result	↓	↓	.806
% Recovery			97%

To Page No. \_\_\_\_\_

TEST: NITRITE, UNITS: mg/L  
 Meth. REF. # 553.2, MDL: 0.10  
 Instrument # #1 (Channel 2)

BY: S.T.  
 BOOK # CV-12 P. 100  
 FROM PAGE # 39

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.
17524	N6110	0		BMDL	17866	P4850	.00151		BMDL
17524	N6120	.0028		BMDL	17866	P4851	.0166		BMDL
17524	N6126	.0040		BMDL	17866	P4852	.0490		BMDL
17531	P2710	.0024		BMDL	17866	P4853	.0446		BMDL
17531	P2725	0		BMDL	17878	P2776	.00181		BMDL
17539	P2117	.0043		BMDL	17866	P3120	.00114		BMDL
17539	P2116	.0954		BMDL	17866	P3122	0		BMDL
17539	P2114	.0571		BMDL	17866	P3124	.0164		BMDL
17524	P1711	.0043		BMDL	17866	P3126	0		BMDL
17524	P1713	.0040		BMDL	17882	P2543	.00471		BMDL
17524	P1714	0		BMDL	17882	P2544	.00213		BMDL
17524	P1717	.00114		BMDL	17882	P2545	.00173		BMDL
17562	N6128	0		BMDL	17882	P2547	.00784		BMDL
17562	N6148	.00119		BMDL	17882	P2549	.00119		BMDL
17573	P2717	0		BMDL	17886	P3117	.0066		BMDL
17579	P2726	.118		.12					
17584	P3068	.100		.10					
17503	P2683	.121		.12					
17562	N6122	.00357		BMDL					
17524	P1715	.00213		BMDL					
17864	P2551	.00119		BMDL					
17864	P2541	.0611		BMDL					
17864	P2546	.0691		BMDL					
17864	P2557	.00286		BMDL					

STANDARDS

PREPARED ON \_\_\_\_\_  
 BY \_\_\_\_\_  
 FROM \_\_\_\_\_ mg/L  
 STOCK # \_\_\_\_\_

STANDARD mg/L	VALUE OBTAINED	% of Threshold
1.00	.99	99%
0.50	.51	102%
0.20	.197	98.5%
0.05	.046	92%

SLOPE	Intercept	Corr. Coef.
.0120	-.0005	.9998

ADDITIONAL COMMENTS

All Calculations  
 By  $X = MY + B$ .

Analyses QA/QC Data Report GC Batch # (N3004 N02) Verified: 11/24/86

PRECISION		ACCURACY	
METHOD BLANK	.00327	Job # Spiked (ETC)	P1714
Spiked Blank (Known)	.20	Original Result	0
Spiked Blank Result	.198	Duplicate Result	0
% Recovery	99	R Time	0.0
EXTERNAL REF. QC	N/A	Job # Spiked (ETC)	P1714
Known Value	.5	Original Result	0
Observed Result	-.532	Amount Spiked	.0282
% Recovery		Spiked S' Result	
		% Recovery	1.31





TEST: Cyanide, UNITS: mg/L  
 Meth. REF. # 335.3, MDL: 0.025  
 Instrument # #1 (citron 1)

BY: C.H.  
 BOOK # CV-12 P. 119  
 FROM PAGE #     

Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	STANDARDS		
17837	N4709	0		BNDL	17697	P4748	.005		BNDL	PREP ON <u>12-2</u>		
17837	N4714	0		BNDL	17697	P4748	.004		BNDL	BY <u>D.B.</u>		
17863	P0489	.007		BNDL	17697	P4748	.004		BNDL	FROM <u>    </u> mg/L		
17863	P0493	.009		BNDL	17698	P1710	.003		BNDL	STOCK # <u>    </u>		
17863	P0494	.005		BNDL	17701	N4710	.004		BNDL	STANDARD	VALUE	% of
17863	P0495	.002		BNDL	17701	N4715	.004		BNDL	mg/L	OBTAINED	Threshold
17866	P4850	.539		.539	17866	P4851	over		Redo	.500	.493	98.6%
17868	P4510	*		Redo	17866	P4852	over		Redo	.400	.405	101.0%
17878	P2776	.007		BNDL	17866	P4853	X		Redo	.200	.205	102.5%
17881	P1865	.002		BNDL	17868	P4510	X		Redo	.100	.102	102%
17883	N4708	.002		BNDL	17868	P4511	X		Redo	SLOPE Intercept Corr. Coef.		
17883	N4711	.003		BNDL	17868	P4512	X		Redo	ADDITIONAL COMMENTS		
17833	P4241	.0197		BNDL	17868	P4513	X		Redo	* RPT DUE TO		
17833	P4242	.0169		BNDL	17883	N4712	X		Redo	INSTRUMENTATION Problem		
17833	P4243	.0211		BNDL	17883	N4713	X		Redo	P4851 10x		
17833	P4244	.0186		BNDL	17889	P2759	X		Redo	P4852 20x		
17834	P4248	.0185		BNDL	17889	P2761	X		Redo			
17834	P4249	.0188		BNDL	17889	P2774	X		Redo			
17834	P4250	.0231		BNDL	17889	P2775	X		Redo			
17834	P4251	.0189		BNDL	17833	P4235	X		Redo			
17834	P4254	.026		BNDL	17833	P4236	X		Redo			
17693	P2916	.0185		BNDL	17833	P4237	X		Redo			
17693	P2917	.0212		BNDL	17833	P4238	X		Redo			
17697	P4748	.006		BNDL	17423	P4239	X		Redo			

Analyses QA/QC Data Report      QC Batch # CY003      Verified: 12/3/86

PRECISION		ACCURACY	
METHOD BLANK	.008	Job # Dup (ETC)	N4709 P4748
Spiked Blank (Known)	.100	Original Result	0 0.006
Spiked Blank Result	.103	Duplicate Result	0 0.005
% Recovery	103%	R Range	0.0 0.001
EXTERNAL REF. QC	97%	Job # Spiked (ETC)	N4709
Known Value	.224	Original Result	0.0
Observed Result	.228	Amount Spiked	.200
% Recovery	101.7%	Spiked S' Result	.201
		% Recovery	99%

305143



305145

305146

1000  
1000

305147

**Technical Report**

**for**

**ENVIRONMENTAL STRATEGIES CORP.**

**WASHINGTON HARBOUR**

**3050 KK ST. SUITE 325**

**WASHINGTON, DC 20007**

*Chain of Custody Data Required for ETC Data Management Summary Reports*

<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>
P4851	ENVIRONMENTAL STRATEGIES CORP.	CHAMPION	WCSP-2	B61110	1300	



**John J. Fitzgerald**

*Vice President  
Research and Operations*

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## Introduction

This report contains the analytical results on your sample. It is designed to include comprehensive data from the entire analytical process in order to satisfy the needs of various levels of review.

The results obtained from your sample are presented in tabular format immediately following this introduction. Quality assurance data is tabulated along with the appropriate sample results for verification. Depending on the analyses ordered, the quality assurance data may include results from blank, spiked blank, spiked sample (i.e. matrix spike) and replicate sample as well as results from surrogate compound analyses. Quality assurance data for verification of proper instrument performance is also included where appropriate. The report appendices include the chain of custody record for your sample and, where appropriate, the gas chromatograms and mass spectra.

The procedures used in the analysis of the sample are described in this report's methodology section. All analytical procedures within our laboratory are performed within a strictly enforced Quality Assurance Protocol. A description of this Protocol is included in the report.

## Results

Sample results, and associated quality assurance data, are always tabulated in one or more of this report's Quantitative Results Tables. The format of each table varies with the class of analysis.

### *Priority Pollutants*

The priority pollutant compounds and elements are listed with their NPDES (National Pollution Discharge Elimination System) numbers, and the Method Detection Limit (MDL) published in the Federal Register. When a compound or element is present below its published MDL it is reported as BMDL (Below Method Detection Limit). When a compound or element is not present at any detectable concentrations it is reported as ND (Not Detected). MDL's for non-aqueous matrices are based on USEPA published MDL's but are adjusted as per sample weight. Matrix spike and replicate analyses, where included, were performed on samples randomly chosen within each quality assurance batch and are therefore not necessarily spikes and replicates of this report's sample. Surrogate compound recovery data and instrument calibration data are included in the Method Performance Data Tables.



**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Volatile Compounds - GC/MS Analysis Data (QR01)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4851 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WOSP-2 861110 1300  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

NPDES Number	Compound <small>Acrolein and Acrylonitrile values are screen only.</small>	Results			QC Replicate		QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concn. Added ug/l	% Recov	Unspiked Sample ug/l	Concn. Added ug/l	% Recov	
1V	Acrolein	ND	100	ND	ND	ND	800	88	ND	800	98	
2V	Acrylonitrile	ND	100	ND	ND	ND	80.0	94	ND	80.0	84	
3V	Benzene	ND	4.4	ND	ND	ND	18.0	99	ND	18.0	100	
4V	bis(Chloromethyl) ether	ND	10	ND	ND	ND	0	-	ND	0	-	
5V	Bromoform	ND	4.7	ND	ND	ND	18.0	92	ND	18.0	90	
6V	Carbon tetrachloride	ND	2.8	ND	ND	ND	18.0	103	ND	18.0	98	
7V	Chlorobenzene	ND	6.0	ND	ND	ND	18.0	99	ND	18.0	99	
8V	Chlorodibromomethane	ND	3.1	ND	ND	ND	18.0	99	ND	18.0	95	
9V	Chloroethane	ND	10	ND	ND	ND	18.0	103	ND	18.0	96	
10V	2-Chloroethylvinyl ether	ND	10	ND	ND	ND	18.0	0 <sup>a</sup>	ND	18.0	0 <sup>a</sup>	
11V	Chloroform	ND	1.6	ND	ND	ND	18.0	103	ND	18.0	98	
12V	Dichlorobromomethane	ND	2.2	ND	ND	ND	18.0	100	ND	18.0	96	
13V	Dichlorodifluoromethane	ND	10	ND	ND	ND	18.0	109	ND	18.0	104	
14V	1,1-Dichloroethane	ND	4.7	ND	ND	ND	18.0	100	ND	18.0	99	
15V	1,1,2-Dichloroethane	ND	2.8	ND	ND	ND	18.0	104	ND	18.0	96	
16V	1,1-Dichloroethylene	ND	2.8	ND	ND	ND	18.0	101	ND	18.0	97	
17V	1,2-Dichloropropane	ND	6.0	ND	ND	ND	18.0	98	ND	18.0	101	
18V	cis-1,3-Dichloropropylene	ND	5.0	ND	ND	ND	18.0	97	ND	18.0	96	
19V	Ethylbenzene	ND	7.2	ND	ND	ND	18.0	100	ND	18.0	100	
20V	Methyl bromide	ND	10	ND	ND	ND	18.0	103	ND	18.0	109	
21V	Methyl chloride	ND	10	209	206	ND	18.0	107	ND	18.0	109	
22V	Methylene chloride	ND	2.8	ND	ND	ND	18.0	98	45.0	18.0	89	
23V	1,1,2,2-Tetrachloroethane	ND	6.9	ND	ND	ND	18.0	101	ND	18.0	106	
24V	Tetrachloroethylene	ND	4.1	ND	ND	ND	18.0	98	ND	18.0	103	
25V	Toluene	ND	6.0	ND	ND	ND	18.0	100	ND	18.0	102	
26V	1,2-Trans-dichloroethylene	ND	1.6	ND	ND	ND	18.0	97	ND	18.0	99	
27V	1,1,1-Trichloroethane	ND	3.8	ND	ND	ND	18.0	104	ND	18.0	97	
28V	1,1,1,2-Trichloroethane	ND	5.0	ND	ND	ND	18.0	104	ND	18.0	100	
29V	Trichloroethylene	ND	1.9	ND	ND	ND	18.0	93	ND	18.0	87	
30V	Trichlorofluoromethane	3.28	10	ND	ND	ND	18.0	105	ND	18.0	99	
31V	Vinyl chloride	ND	10	ND	ND	ND	18.0	110	ND	18.0	99	
18V	trans-1,3-Dichloropropylene	ND	10	ND	ND	ND	18.0	94	ND	18.0	96	

<sup>a</sup> Recovery normally variable using established methodology.

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**ETC** ENVIRONMENTAL TESTING and CERTIFICATION

NOV 29, 1986

**TABLE 1: QUANTITATIVE RESULTS**

**Metals - Analysis Data (QR52)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4851 ENVIRONMENTAL STRATEGIES CORP. CHAMPION HCSP-2 861110 1300  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

Compound	Results	
	Sample Concn. ug/l	MDL ug/l
Barium	61	3.2
Chromium	BMDL	26
Copper	13	13
Iron	BMDL	140
Manganese	110	6.4
Nickel	ND	11
Sodium	12900	72
Zinc	67	2.9

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Conventional Analysis Data (QR10)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4851** ENVIRONMENTAL STRATEGIES CORP. CHAMPION MCSP-2 861110 1300  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

NPDES Number	Results	
	Sample Concen. mg/l	MDL mg/l
Fluoride	.21	.10
Nitrate as N	7.93	.10
Sulfate as SO4	94	2
Phenolics, Total	<.050	.050
Cyanide, Total	.029	.025
Solids, total suspended	890	2

**TABLE 2: METHOD PERFORMANCE DATA**  
**Surrogate Recovery - Aqueous Matrices (QR 20)**

Chain of Custody Data Required for ETC Data Management Summary Reports

P4851 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WOSP-2 86111 1300 0  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

Compound	Amount added ug	% Recovery	Control Limits	
			Lower	Upper
VOLATILE FRACTION (GC/MS)				
Toluene-D8	.250	100	88	110
p-Bromofluorobenzene	.250	103	86	115
1,2-Dichloroethane-D4	.250	104	76	114
ACID FRACTION (GC/MS)				
Phenol-D5	-	-	10	94
2-Fluorophenol	-	-	21	100
2,4,6-Tribromophenol	-	-	10	123
BASE/NEUTRAL FRACTION (GC/MS)				
Nitrobenzene-D5	-	-	35	114
2-Fluorobiphenyl	-	-	43	116
Terphenyl-D14	-	-	33	141
PESTICIDE/PCB FRACTION (GC)				
Dibutylchlorendate	-	-	24**	154**

\* IFB EPA Control Limits  
 \*\* Recovery Limits Only

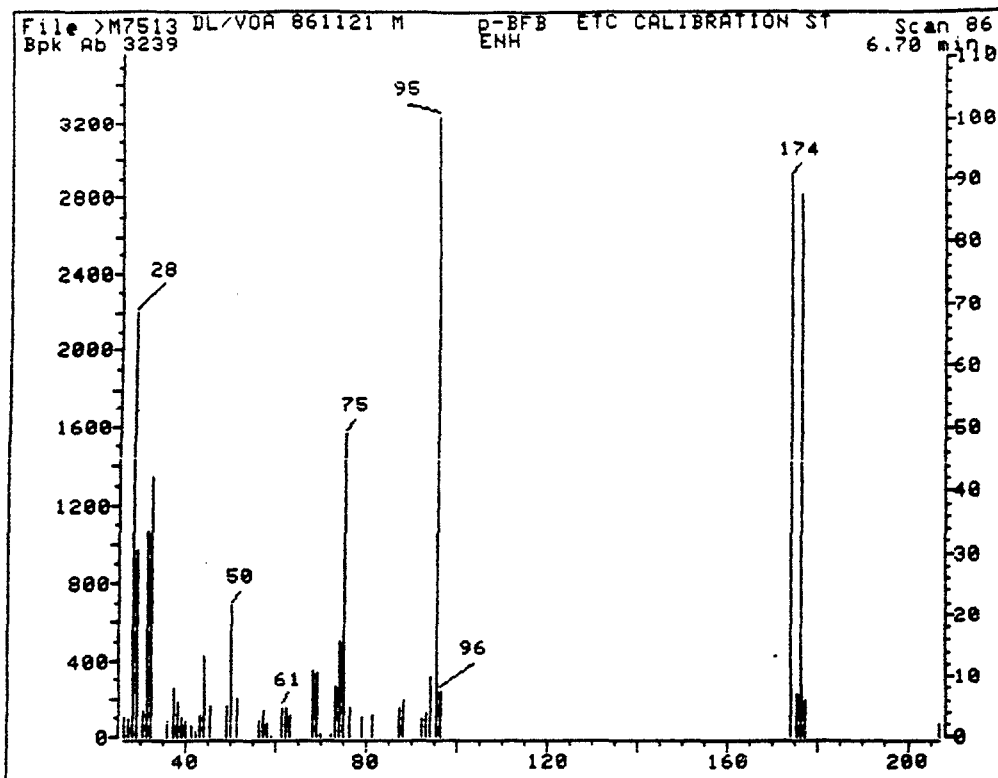


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	21.09	21.09	Ok
75	30-60% of mass 95	48.39	48.39	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.22	7.22	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	89.98	89.98	Ok
175	5-9% of mass 174	6.74	7.49	Ok
176	95-101% of mass 174	87.27	96.99	Ok
177	5-9% of mass 176	5.98	6.86	Ok

Injection Date: 11/21/86  
 Injection Time: 20:42  
 Run No: >M7513  
 Spectrun No: 86

Analyst: J. Quinn  
 Processor: QUANTALAB  
 QC Batch: QV 5431  
 Samples: P4240, P4241, P4242, P4243, P4244, P4236, P4237, P4239, P4235, P533, P5415, P5414, P4850, P4851, P485

*Handwritten signature and date:* 11/25/86

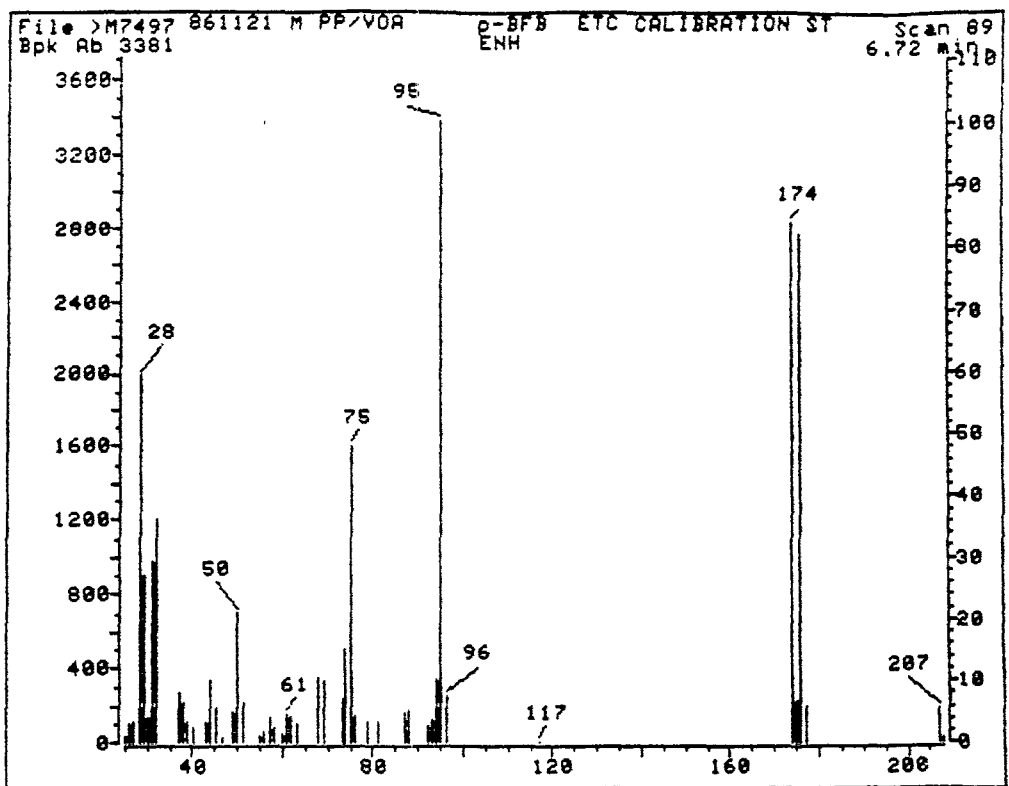


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance		Status
		Base Peak	Appropriate Peak	
50	15-40% of mass 95	20.80	20.80	Ok
75	30-60% of mass 95	47.52	47.52	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.18	7.18	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	83.99	83.99	Ok
175	5-9% of mass 174	6.60	7.85	Ok
176	95-101% of mass 174	82.16	97.82	Ok
177	5-9% of mass 176	5.86	7.13	Ok

Injection Date: 11/21/86  
 Injection Time: 05:57  
 Run No: >M7497  
 Spectrun No: 89

Analyst: Joseph Quinn  
 Processor: BLUON CHRISTOPHER  
 QC Batch: QV 8931  
 Samples: P4240, P4241, P4242, P4243, P4236, P4237, P4239, P4235, P5415, P5414, P4850, P4851, P

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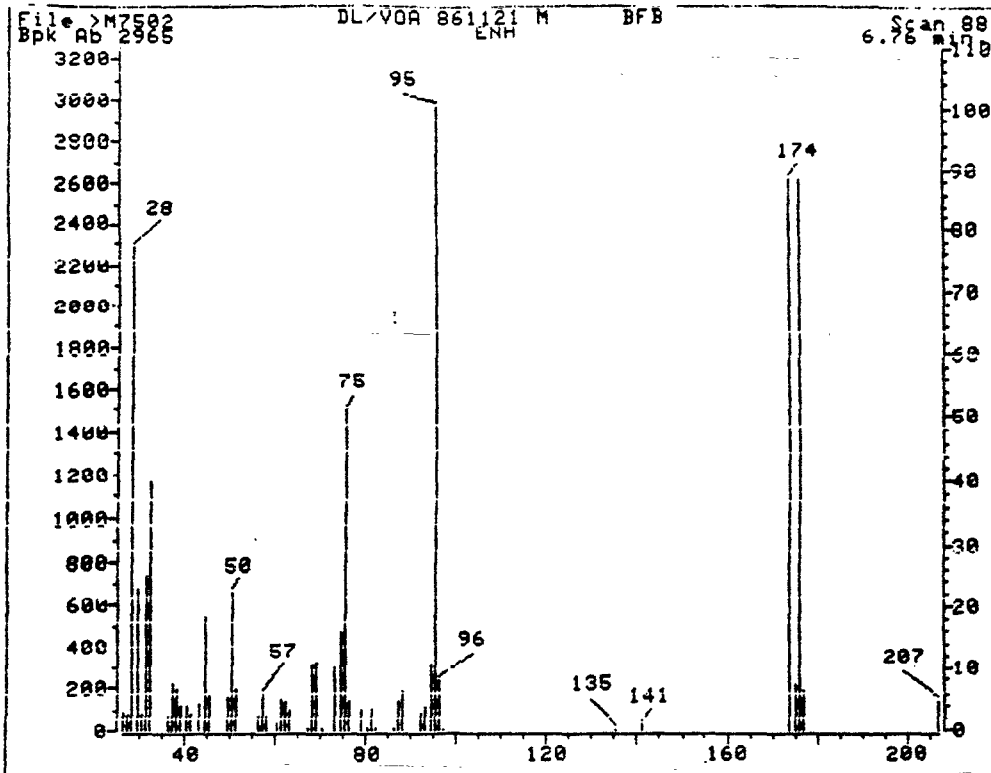


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	22.00	22.00	Ok
75	30-60% of mass 95	50.62	50.62	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.63	7.63	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	88.52	88.52	Ok
175	5-9% of mass 174	7.03	7.95	Ok
176	95-101% of mass 174	88.43	99.89	Ok
177	5-9% of mass 176	6.31	7.15	Ok

Injection Date: 11/21/86  
 Injection Time: 09:32  
 Run No: >M7502  
 Spectrum No: 88

Analyst: *Nancy Albert*  
 Processor: *Quinn Christopher*  
 QC Batch: *QV 5931*  
 Samples: *P4240, P4241, P4242, P4243, P4244, P4236, P4237, P4238, P4235, P5334, P5415, P5414, P4850, P4851, P4852*

## Methodology for GC/MS Analysis of Priority Pollutant Compounds

The methods employed in the GC/MS analysis for priority pollutants are established EPA methods. Rigid compliance with the instrument parameters and performance criteria of the published methods was achieved. In some cases, the precise amounts of sample used and the sample handling procedures vary with the complexity of the sample matrix. Qualitative identification of the priority pollutants was performed using the relative retention times, the relative abundance of three characteristic ions and the abundance ratios. The entire mass spectrum was reviewed to confirm each identification. Quantitative analysis of detected compounds was performed by using a response factor generated by a major characteristic ion of the specific compound and an internal standard.

Compounds, in addition to those on the priority pollutant list, were identified through a computer-aided search of the NBS-EPA spectra library. After review the identifications are included in a separate tabulation and labelled "tentatively identified".

### ***Volatile Priority Pollutant Compounds***

For the analysis of Volatile priority pollutants, EPA Method 624 was used. The method can be summarized as follows: Helium is bubbled through a 5 ml water sample contained in a specially designed purging chamber at ambient temperature. The purgeable volatile organic compounds are transferred from the aqueous phase to the vapor phase. The vapor is swept through a sorbent column where the organic components are trapped. After the purge cycle is complete, the sorbent column is heated and backflushed with helium to desorb the organic purgeables onto a gas chromatographic column. The gas chromatograph is temperature programmed to separate the purgeable mixture. The separated purgeable components are then identified and quantitated using a computerized mass spectrometer.

### ***Acid, Base/Neutral and Pesticide Priority Pollutant Compounds***

For the analysis of the Acid, Base/Neutral and Pesticide priority pollutants in an aqueous liquid matrix, EPA Method 625 was used. The method can be summarized as follows: A measured volume of sample, approximately 1 liter, is adjusted to a pH greater than 11 and extracted with methylene chloride. The pH of the sample is adjusted to a value less than 2 and extracted with an aliquot of fresh methylene chloride. A separatory funnel or continuous extractor is used to perform the extractions. The two extracts are dried and concentrated to a 1 ml final volume. Each extract is injected into a GC/MS instrument specifically configured for the correct fraction.



## Methodology for Analysis of Metals

### **AQUEOUS**

The determination of metals in aqueous samples is performed according to the methods published by EPA in "Methods for Chemical Analysis of Water and Wastes," EPA-600/4-79-020, March, 1983, and the Federal Register, October 26, 1983. Arsenic, selenium and thallium are determined by furnace AA; silver, aluminum, barium, beryllium, boron, cadmium, calcium, chromium, copper, cobalt, iron, magnesium, manganese, molybdenum, nickel, lead, sodium, antimony, tin, titanium, vanadium, and zinc are determined by ICP emission spectrometry, except where lower levels of detection are required; in these cases (e.g. lead in groundwater monitoring samples) furnace AA is used. All furnace AA parameters are run by method of standard additions. The determination of mercury is performed by cold vapor AA.

### **EP TOXICITY**

The determination of metals in aqueous EP Toxicity leachates is performed according to the methods published by EPA in "Test Methods for Evaluating Solid Waste" EPA SW-846, revised April, 1984 and the Federal Register, Oct. 26, 1983, 1979. Silver, arsenic, barium, cadmium, chromium, lead and selenium are determined by ICP emission spectrometry. Mercury is determined using cold vapor AA. For leachates that are organic in nature, the analyses are performed according to the methods described under **OIL/SLUDGE** below.

### **SOIL/SEDIMENT**

The determination of silver, beryllium, cadmium, chromium, copper, nickel, antimony, lead, and zinc in sediment samples is performed according to methods published by EPA in "Interim Methods for the Sampling and Analysis of Priority Pollutants in Sediments and Fish Tissue", EPA 600/4-81-055, October 1980. Mercury is determined according to the sediment method published by EPA in "Method for Chemical Analysis of Water and Wastes", EPA 600/4-79-020, March 1983. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion.

### **OIL/SLUDGE**

The determination of silver, aluminum, boron, barium, beryllium, calcium, cadmium, copper, chromium, cobalt, iron, magnesium, manganese, molybdenum, sodium, nickel, lead, antimony, tin, titanium, vanadium, and zinc in sludge/petroleum-based samples is performed by ICP emission spectrometry using a magnesium nitrate dry ashing digestion technique. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion. Mercury is determined by cold vapor AA using the same digestion technique.

## Summary of Quality Assurance/Quality Control Procedures (QA/QC)

ETC bases its quality assurance protocols on the following government guidelines:

- "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-330/9/79/001-R, October 1979;
- the recommended guidelines for EPA Methods 624 and 625. (Federal Register, December 3, 1979, updated on October 26, 1984);
- "Manual of Analytical Methods for the Analysis of Pesticides in Humans and Environmental Samples," EPA 600/8-80-038, June 1980;
- "Determination of 2,3,7,8-TCDD in Soil and Sediment" EPA, Region VII, Kansas City, September 1963;
- Organic Analysis: Multi-media Multi Concentration-IFB WA84-A267; and
- Dioxin Analysis: Soil/Sediment Matrix; Multi-Concentration; Selected Ion Monitoring with Jar Extraction Procedure-IFB WA84-A002

However, we have modified our protocols to provide a higher level of QA/QC than the guidelines require. For example, we analyze a higher than required number of quality control samples and we pay especially careful attention to the certification of the "reference standard" compounds we use in analysis. Below are listed the key QA/QC elements for the methods we used.

### Analysis of Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry

- Each batch of 13 samples consists of 9 customer samples (at a maximum), one blank sample, one spiked blank, one spiked sample and one replicate sample. This amounts to a 30% quality control factor.
- Three surrogate compounds are added to each sample in the batch of 13.
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and retuned, if necessary, at the beginning of each day to ensure that its performance on bromofluorobenzene (BFB) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of Volatile Organic Priority Pollutant "standards" at a minimum of 3 different concentrations and using a mixture of 3 internal standards at a constant concentration.
- The calibration curve is verified with a mixture of priority pollutant standards every day. If the response factors vary greater than 25%, the instrument must be recalibrated.

### Analysis of Organic Compounds Extracted in Acid or Base/Neutral Solutions by Gas Chromatography/Mass Spectrometry

- Each batch of 20 samples consists of 16 customer samples (at a maximum), one blank sample, one spiked blank (for water matrices), one sample spiked with the priority pollutant standard mixture and a duplicate customer sample. This amounts to a 20% quality control factor.

- Three surrogate compounds are added to each sample in the batch for Base/Neutral analysis.
- Three surrogate compounds are added to each sample in the batch for Acid analysis.
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and retuned, if necessary, at the beginning of each day to ensure that its performance on decafluorotriphenylphosphine (DFTPP) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of standards composed of either the Organic Acid or Base/Neutral Extractable Compounds at a minimum of 3 concentrations and using five internal standards for quantitation.

### Analysis of Metals

#### All Samples

- New standards are prepared for each batch of samples
- Normal calibration is performed using a blank sample and four standards that have been through the sample preparation procedure. A regression analysis is used to construct the calibration curve.
- All EP Toxicity samples and all samples determined by furnace atomic absorption are calculated by the "method of additions".
- For each sample analysis that requires the use of the "method of additions" technique, a three point calibration is performed using U.S. EPA "Methods for Chemical Analysis of Water and Wastes, 1979". Results are obtained using linear regression analysis. Any regression with a coefficient of correlation below 0.990 is considered suspect, necessitating review of calibration data or sample re-analysis.
- In constructing the normal calibration curves the lowest concentration levels we use are values greater than or equal to 5 times the Instrumental Detection Limit (IDL).
- All calibration standards are analyzed in duplicate, at a minimum.
- Independent reference standards are used to check the accuracy of calibration standards.
- A check standard is analyzed every ten samples to validate the normal calibration curve.
- One customer sample out of every ten is analyzed in triplicate.

#### Homogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are homogeneous, the QC program is a minimum of 25% and consists of analyzing:

- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 1 independent reference standard;
- 4 Calibration standards (processed using the sample preparation method);
- 4 Calibration standards (without sample preparation); and
- 1 Reagent Blank.

Heterogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are heterogeneous, the QC program is a minimum of 35% and consists of analyzing

- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 1 Replicate independent reference standards;
- 4 Calibration standards (processed using the sample preparation method);
- 1 Procedural Blank;
- 4 Calibration standards (without sample preparation); and
- 1 Reagent Blank.

Analysis of Mercury

To analyze samples for mercury we group them by matrix in batches of 30 or less. Our QC program is a minimum of 30% and consists of analyzing:

- each of the 30 customer samples in duplicate;
- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 2 Replicate independent reference standards;
- 10 Calibration standards (processed using the sample preparation method); and
- 2 Procedural Blanks.

Analysis of Pesticides, Herbicides and PCB's by Gas Chromatography

Pesticide, herbicide and PCB samples are grouped in batches of 16 customer samples or less according to the type of analysis to be performed. The QC program for each of these three types of analyses is a minimum of 20% and consists of analyzing:

- 1 procedural blank sample (a reagent blank is analyzed in the case of non-water matrices);
- 1 spiked blank sample (the spiked blank is eliminated in the case of non-water matrices);
- 1 replicate sample;
- 1 replicate spiked sample; and
- 1 known reference QC sample for at least each 100 samples analyzed.

The instrument is calibrated each run with three standards, and checked every 10 samples.

Analysis of Cyanides, Phenols, Fluoride, Chloride, Nitrate and Nitrite

- All parameters are analyzed using a Technicon Autoanalyzer II GT.
- 3 calibration standards are analyzed at the beginning and end of each batch.

- Each batch (up to 80 samples) consists of analyzing one blank, one spiked blank, one duplicate and spiked sample every 20 samples, and an EPA known reference sample.

#### Analysis of Total Organic Carbon (TOC)

TOC samples are analyzed on a daily basis with the number of samples analyzed per day dependent on the request for duplicate or quadruplicate analyses. The quality control program is designed to maintain the appropriate amount of QC and consists of the following elements:

- Daily instrument calibration
- One blank
- Standard recalibration every 10 samples
- Spiked samples at a low and high level
- Every sample is run in duplicate at a minimum

#### Analysis of Total Organic Halide (TOX)

- Blank reagent water for absolute carbon background must contain less than 5 ug/l of halide (as chloride)
- Using a trichlorophenol standard, the mean adsorption efficiency must be within +/- 15% of the standard value.
- Calibration standards are run every 10 samples.
- Every sample is run in duplicate at a minimum.

#### Analysis of 2,3,7,8-TCDD (Dioxin) by GC/MS (SIM)

- Each sample is dosed with a known quantity of  $^{13}\text{C}_{12}$ -2,3,7,8-TCDD as internal standard and  $^{37}\text{Cl}_4$ -TCDD as surrogate standard. The action limits for surrogate standard results is +/- 40% of the true value. Samples showing surrogate standard results outside of these limits are reextracted and reanalyzed.
- Two laboratory "method blanks" are run along with each set of 24 or fewer samples. The method blank is also dosed with the internal standard and surrogate standard.
- At least one method blank sample is run in duplicate to determine intralaboratory precision.
- Qualitative Requirements. The following are met in order to confirm the presence of native 2,3,7,8-TCDD:
  - a. Isomer specificity must be demonstrated initially and verified once per 8-hour work shift. The verification consists of injecting a mixture containing TCDD isomers which elute close to 2,3,7,8-TCDD. The 2,3,7,8-TCDD must be separated from interfering isomers, with no more than 25% valley relative to the 2,3,7,8-TCDD peak.
  - b. The 320/322 ratio is within the range of 0.67 to 0.67.
  - c. Ions 320, 322, and 257 are all present and maximize together the signal to mean noise ratio must be 2.5 to 1 or better for all 3 ions.
  - d. The retention time is equal (within 3 seconds) the retention time for the isotopically labeled 2,3,7,8-TCDD.
  - e. At least one of the positives can be confirmed by obtaining partial scan spectra from mass 150 to mass 350. The partial scan guidelines are as follows:

- . the 320/324 ratio should be 1.58 +/- 0.16
- . the 257/259 ratio should be 1.03 +/- 0.10
- . the 194/196 ratio should be 1.54 +/- 0.15
- One sample is spiked with native 2,3,7,8-TCDD at a level of 1.0 PPB (for soil) for each set of 24 or fewer samples.
- In cases where no native 2,3,7,8-TCDD is detected, the actual detection limit is estimated and reported based on a signal to noise ratio of 2.5 to 1 at ions 320 and 322.
- For each sample, the internal standard is present with at least a 10 to 1 signal to noise ratio for both mass 332 and mass 334. Also, the internal standard 332/334 ratio must be within the range of 0.67 to 0.87.

#### Subcontractor QA/QC

Each subcontractor is required to maintain an appropriate level of quality control. To insure this, each subcontractor is required to submit to ETC the quality control data for all analyses it performs. This data is kept on file at ETC. In general, the amount of quality control required is one duplicate sample with one spiked sample for every ten analyses.

#### Chain-of-Custody

The chain-of-custody procedure is part of our quality assurance protocol. We believe our chain-of-custody record fully complies with the legal requirements of federal, state and local government agencies and of the courts of law. The record covers:

- labeling of sample bottles, packing the Sample Shuttle and transferring the Shuttle under seal to the custody of a shipper;
- outgoing shipping manifests;
- the chain-of-custody form completed by the person(s) breaking the Shuttle seal, taking the sample, resealing the Shuttle and transferring custody to a shipper;
- incoming shipping manifests;
- breaking the Shuttle's reseal;
- storing each labeled sample bottle in a secured area;
- disposition of each sample to an analyst or technician; and
- the use of the sample in each bottle in a testing procedure appropriate to the intended purpose of the sample.

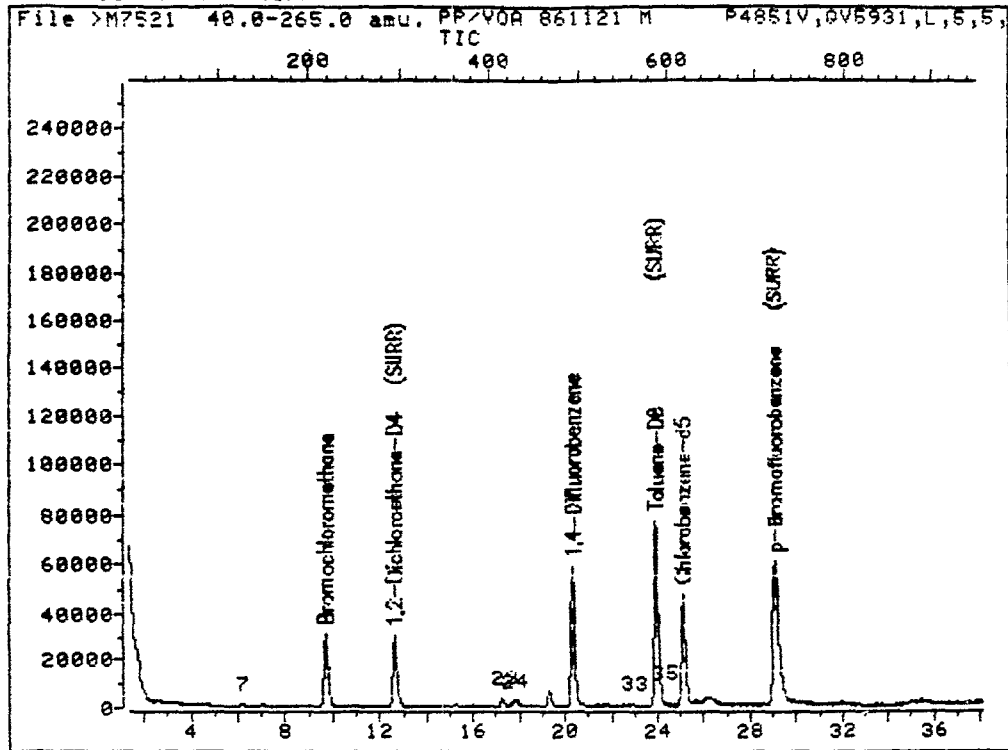
The records show for each link in this process:

- the person with custody; and
- the time and date each person accepted or relinquished custody.

**Appendix A**  
**Mass Spectral Data**  
**for**  
**Quantitated Compounds**

- 1) A total ion chromatogram for each sample analyzed by a GC/MS instrument.
- 2) A Quant. report used by the analyst to determine qualitative and quantitative results of the compounds present.
- 3) A mass spectrum and a reference spectrum for each priority pollutant compound detected in the sample.

TOTAL ION CHROMATOGRAM



Data File: >M7521::U4  
Name: PP/VQA 861121 M  
Misc: P4851V,QV5931,L,5,5,

Quant Output File: ^M7521::AQ

Id File: MVQA::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861121 12:01

Operator ID: DG1199  
Quant Time: 861122 04:51  
Injected at: 861122 04:12

305166



QUANT REPORT

Operator ID: DG1199  
 Output File: <M7521::AQ  
 Data File: >M7521::U4  
 Name: PP/UQA 861121 M  
 Misc: P4851U, QV5931, L, 5, 5,

Quant Rev: 6      Quant Time: 861122 04:51  
 Injected at: 861122 04:12  
 Dilution Factor: 1.00000

ID File: MVDA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861121 12:01

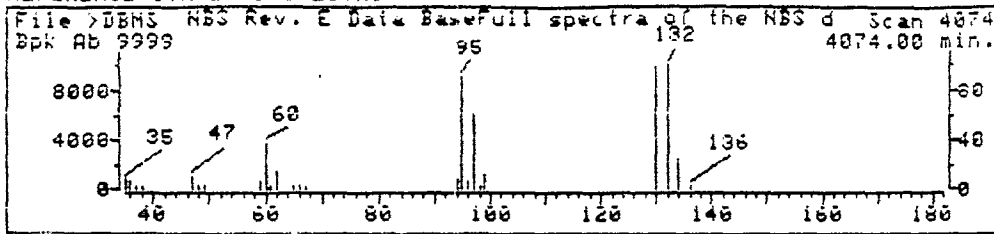
Compound	R.T.	Scan#	Area	Conc	Units	q
1) *Bromochloromethane	9.70	217	65287	250.00	NG	94
<del>7) Methylene chloride</del>	<del>6.14</del>	<del>125</del>	<del>2874</del>	<del>28.17</del>	<del>NG</del>	<del>92</del>
16) *1,4-Difluorobenzene	20.35	492	257366	250.00	NG	82
17) 1,2-Dichloroethane-D4 (SURR)	12.64	293	59194	260.91	NG	96
23) Trichloroethylene	17.25	412	6765	16.40	NG	95
<del>24) Benzene</del>	<del>17.72</del>	<del>424</del>	<del>4623</del>	<del>4.46</del>	<del>NG</del>	<del>95</del>
30) *Chlorobenzene-d5	25.11	615	178018	250.00	NG	97
<del>33) Tetrachloroethylene</del>	<del>22.87</del>	<del>557</del>	<del>1737</del>	<del>3.58</del>	<del>NG</del>	<del>83</del>
34) Toluene-DB (SURR)	23.95	585	321390	248.99	NG	94
<del>35) Toluene</del>	<del>24.14</del>	<del>598</del>	<del>2561</del>	<del>2.81</del>	<del>NG</del>	<del>96</del>
38) p-Bromofluorobenzene (SURR)	29.14	719	190591	258.04	NG	95

\* Compound is ISTD

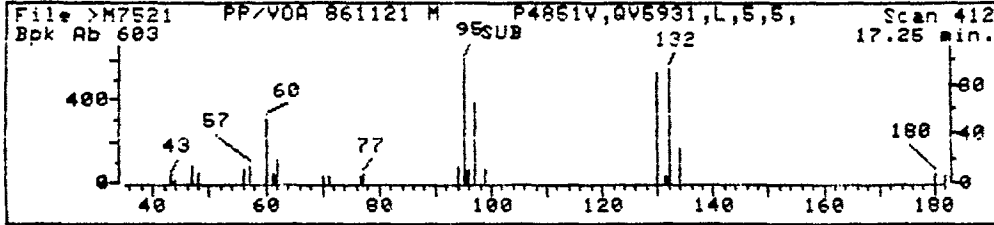
SNC 11/24/86

905167

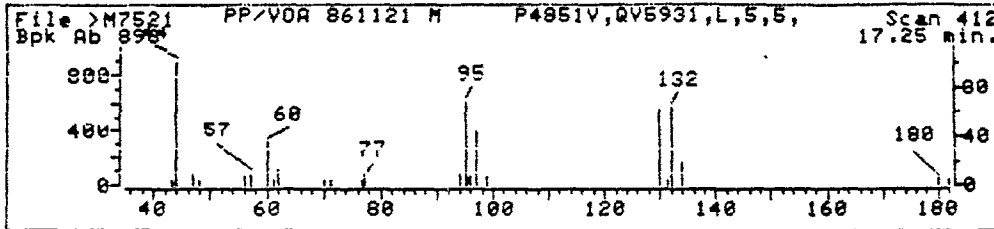
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >M7521::U4  
Name: PP/VOA 861121 M  
Misc: P4851V,QV5931,L,5,5,  
Quant Time: 861122 04:51  
Injected at: 861122 04:12

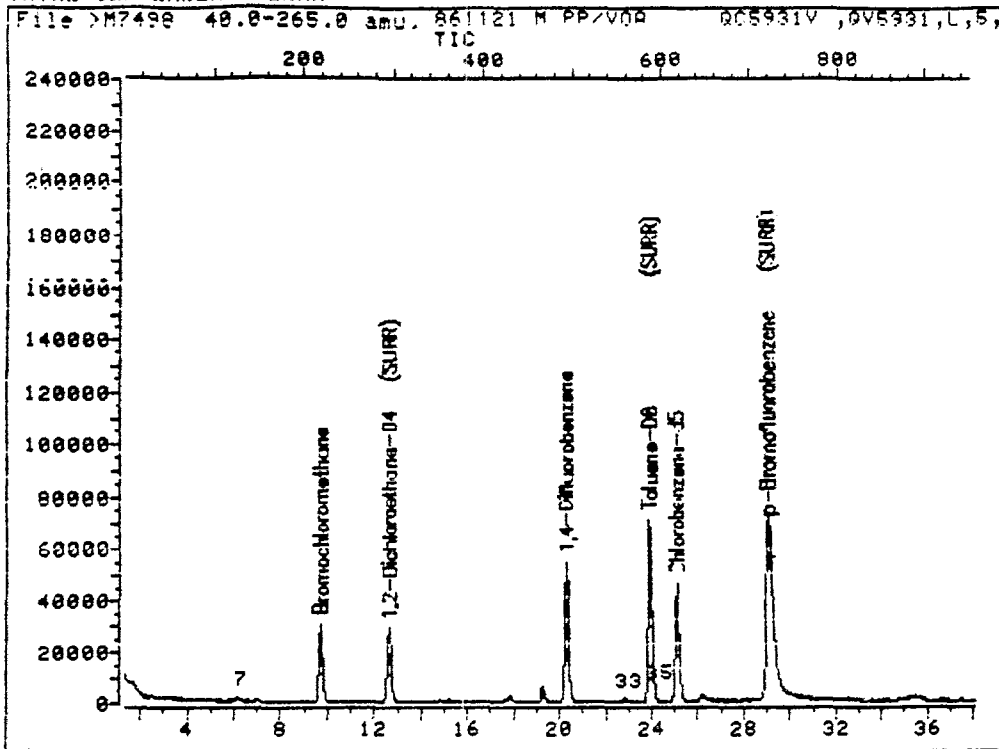
Quant Output File: ^M7521::AQ  
Quant ID File: MVDA::US  
Last Calibration: 861121 12:01

Compound No: 23  
Compound Name: Trichloroethylene  
Scan Number: 412  
Retention Time: 17.25 min.  
Quant Ion: 95.0  
Area: 6765  
Concentration: 16.40 NG  
q-value: 95

305168  
nd 10/11

**Appendix C1**  
**GC/MS Subsidiary Data**

TOTAL ION CHROMATOGRAM



Data File: >M7498::U3  
Name: 861121 M PP/VDA  
Misc: Q05931V ,Q05931,L,5,5,

Quant Output File: ^M7498::AQ

Id File: MVUA::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861121 12:01

Operator ID: JQ6275  
Quant Time: 861121 12:03  
Injected at: 861121 06:10

305170

QUANT REPORT

Operator ID: JW6275  
 Output File: ^M/498::AQ  
 Data File: >M/498::U3  
 Name: 861121 M PP/VOA  
 Misc: QC5931U ,QV5931,L,5,5,

Quant Rev: 6      Quant Time: 861121 12:03  
 Injected at: 861121 06:18  
 Dilution Factor: 1.00000

ID File: MVDA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861121 12:01

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *Bromochloromethane	9.70	217	66376	250.00	NG	99
7) Methylene chloride	6.18	126	2829	19.63	NG	86
16) *1,4-Difluorobenzene	20.31	491	249275	250.00	NG	82
17) 1,2-Dichloroethane-D4 (SURR)	12.65	293	57076	259.74	NG	94
30) *Chlorobenzene-d5	25.11	615	170859	250.00	NG	99
<del>33) Tetrachloroethylene</del>	<del>22.83</del>	<del>556</del>	<del>2829</del>	<del>6.88</del>	<del>NG</del>	<del>94</del>
34) Toluene-D8 (SURR)	23.95	585	307632	248.31	NG	91
<del>35) Toluene</del>	<del>24.15</del>	<del>598</del>	<del>1931</del>	<del>1.58</del>	<del>NG</del>	<del>92</del>
38) p-Bromofluorobenzene (SURR)	29.10	718	181687	256.29	NG	97

\* Compound is ISTD

ENC 11/24/86

**Appendix D**  
**Subcontractor's Data**

- 1) A copy of the originating subcontractor's report is included for all data not generated within ETC's laboratory.



## Appendix E

### Chain-of Custody Forms

1. A field Chain-of-Custody form (CC1) is included for all samples shipped by ETC shuttle.
2. An in-house sample Chain-of-Custody form is included for all samples not shipped by ETC shuttle.
3. Any additional Chain-of-Custody material provided by a client or by a client's sampling agent is also included.
4. A subcontractor's Chain-of-Custody form is included for any analytical work not performed within ETC's laboratory.
5. Analysis and Extraction Custody forms are included for the period the sample was in ETC's possession.



**CHAIN OF CUSTODY FORM (CC1)**

Company: HELLERTOWN MANUFACTURING Attn.: MARK SCHULTZ  
 Facility/Site: C/O GUARD OFFICE-SILVER ROAD Phone: \_\_\_\_\_  
 Address: HELLERTOWN, PA. 18055

**SAMPLE IDENTIFICATION**

Facility: CHAMPION (Optional Sample Point Descriptions)  
 Sample Point: W-1 C-SIP-2 111086 1300  
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	M	500 ML	HN03	METALS			/
1	PN	125 ML	H2S04	PHENOLS			/
1	B	125 ML	H2S04	NITRATE			/
1	CN	125 ML	NAOH	CYANIDE			/
1	A	500 ML	BAKED	SOLIDS, TD/SULFATE			/
1	D	500 ML	RINSE	FLUORIDE			/
2	U	40 ML	THIOL	VOA			VI-BUB

**CHAIN OF CUSTODY CHRONICLE**

1. Shuttle Opened By: (print) M. SCHULTZ Date: 11-10-86 Time: \_\_\_\_\_  
 Signature: Mark Schultz Seal #: \_\_\_\_\_ Intact: \_\_\_\_\_  
 2. I have received these materials in good condition from the above person.  
 Name: \_\_\_\_\_ Signature: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_ Remarks: \_\_\_\_\_  
 3. I have received these materials in good condition from the above person.  
 Name: \_\_\_\_\_ Signature: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_ Remarks: \_\_\_\_\_  
 Shuttle Sealed By: (print) M. SCHULTZ Date: 11-10-86 Time: 16:30  
 Signature: Mark Schultz Seal #: 6067754 Intact: \_\_\_\_\_

LAB USE ONLY Opened By: Mark Schultzt Date: 11/11/86 Time: 5:22 PM  
 SHUTTLE # 381 TEMP. °C 12 SEAL # 67754 COND. INTACT

305175







ANALYZED ON: 11/26/86

TEST: Fluoride, UNITS: mg/L  
 Meth. REF. # STD METHOD 413E, MDL: 0.10  
 Instrument # A.A. Channel 7

BY: C. K. ...  
 BOOK # CV-12 P. 108  
 FROM PAGE # 107

Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	STANDARDS
17455	m0831	0.07	-	BMDL	17828	P2472	0.03	-	BMDL	PREPARED ON _____ BY _____ FROM _____ mg/L STOCK # _____
17455	m0703	0.26	10x	2.60	17828	P2473	0.03	-	BMDL	
17455	m0704	0.27	10x	2.70	17828	P2522	0.08	-	BMDL	
17455	m0705	0.23	10x	2.30	17828	P2536	0.03	-	BMDL	
17455	m0706	0.28	10x	2.80	17828	P2548	0.02	-	BMDL	
17455	m0707	0.23	10x	2.30	17866	P4850	0.16	-	0.16	
17455	m0708	0.03	-	BMDL	17866	P4851	0.21	-	0.21	
17514	<del>m0708</del> FL003	0.44	-	0.44	17866	P4852	2.24	-	RPT	
17514	<del>m0708</del> FL003	0.54	-	0.54	17866	P4853	0.98	-	0.98	
17749	P2500	0.01	-	BMDL	17250	N4520	1.64	-	1.64	
17749	P2513	0.07	-	BMDL	17441	P2237	0.07	-	BMDL	STANDARD mg/L    VALUE OBTAINED    % of Theoretical
17749	P2514	0.03	-	BMDL	17878	P2776	0.087	-	BMDL	
17749	P2515	0.07	-	BMDL	17864	P2481	0.06	-	BMDL	SLOPE    Intercept    Corr. Coef.
17749	P2526	0.04	-	BMDL	17864	P2531	0.02	-	BMDL	
17778	P2497	0.03	-	BMDL	17864	P2541	0.04	-	BMDL	ADDITIONAL COMMENTS RPT: N4519 - Lake 7 m. 1 P4852 - and Lake 5 site in treatment raise Dilute 1:10 and check
17778	P2499	0.03	-	BMDL	17864	P2546	0.02	-	BMDL	
17778	P2507	0.13	-	0.13	17864	P2554	0.01	-	BMDL	
17778	P2517	0.12	-	0.12						
17748	P2469	0.03	-	BMDL						
17748	P2477	0.05	-	BMDL						
17748	P2524	0.08	-	BMDL						
17748	P2538	0.05	-	BMDL						
17748	P2539	0.04	-	BMDL						
17748	P2540	0.02	-	BMDL						

Analyses QA/QC Data Report  
 See CV-12, P. 107.

QC Batch # FL001 + FL003

Verified: [Signature] 11/28/86

METHOD BLANK	PRECISION
Spiked Blank (Known)	Job # Dup (ETC) <u>P2472</u>
Spiked Blank Result	Original Result <u>.0297</u>
% Recovery	Duplicate Result <u>.0339</u>
	R Range <u>.0042</u>
EXTERNAL REF. QC	ACCURACY
Known Value	Job # Spiked (ETC) <u>P2472</u>
Observed Result	Original Result <u>.0313</u>
% Recovery	Amount Spiked <u>1.00</u>
	Spiked S' Result <u>1.07</u>
	% Recovery <u>104.40</u>

TEST: NITRATE UNITS: mg/L  
 METH. REF. # 353.2 MDL: 0.10  
 INSTRUMENT # #1 (Channel 1)

BY: CH  
 BOOK # CV-12 P. 97  
 FROM PAGE # 96

Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.
7514	N6032	.334	5X	1.67	17866	P4851	.793	10X	7.93
7514	N6033	.0134	5X	ReLo	17866	P4852	.0487	10X	
7562	N6122	.804	5X	4.02	17866	P4853	.442	10X	4.42
7503	P2683	.0213	5X	ReLo	17866	P4854	.0192	10X	
7798	P2469	.276	10X	2.76	17866	P4857	.153	10X	1.53
7798	P2477	1.54	10X	15.4	17866	P4858	.0198	10X	
17798	P2524	.517	10X	5.17	17866	P4859	.0211	10X	
17798	P2538	.583	10X	5.83	17866	P4861	.0227	10X	
17796	P2539	.856	10X	8.56	17866	P4862	.0178	10X	
17798	P2540	.618	10X	6.18	17866	P2543	.528	10X	5.28
17814	P3555	1.01	10X		17866	P2544	.0253	10X	
17822	P3107	.417	10X	4.17	17866	P2548	1.33	10X	13.3
17822	P3112	.0139	10X		17866	P2547	1.44	10X	14.4
17828	P2473	.442	10X	4.42	17866	P2549	.924	10X	9.24
17828	P2472	.547	10X	5.47	17866	P2550	.401	10X	4.01
17828	P2522	.417	10X	4.17	17866	P2551	1.03	10X	10.3
17828	P2536	.435	10X	4.35	17866	P2552	.717	10X	7.17
17828	P2548	.525	10X	5.25					
17864	P2481	.0182	10X						
17864	P2521	.0198	10X						
17864	P2541	.160	10X	1.60					
17864	P2546	.697	10X	6.97					
17864	P2537	.196	10X	1.96					
17866	P4850	.550	10X	5.50					

STANDARDS		
STANDARD mg/L	VALUE OBTAINED	% of Theoretical
PREP ON _____		
BY _____		
FROM _____ mg/L		
STOCK# _____		
SLOPE _____		
Intercept _____		
Corr. Coef. _____		

**ADDITIONAL COMMENTS**  
 STRAIGHT Red Line Through  
 Samples Done on Channel  
 2. No dilution Page 98  
 RPT N6033 } straight  
 P2683 }

**Analyses QA/QC Data Report**

QC Batch # N3003 & N3004

Verified: 11/24/91 *[Signature]*

PRECISION			
METHOD BLANK		Job # Dup (ETC)	
Spiked Blank (Known)		Original Result	
Spiked Blank Result		Duplicate Result	
% Recovery		R RANGE	
		ACCURACY	
EXTERNAL REF. QC		Job # Spiked (ETC)	
Known Value		Original Result	
Observed Result		Amount Spiked	
% Recovery		Spiked S' Result	
		% Recovery	

305180

TEST: NITRITE, UNITS: mg/L  
 METH. REF. # 553.2, MDL: 0.10  
 INSTRUMENT # #1 (Channel 2)

BY: S.T.  
 BOOK # CV-12 P. 100  
 FROM PAGE # 39

Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	STANDARDS
17524	N6110	0		BMDL	17866	P4850	.00151		BMDL	PREPARED ON _____ BY _____ FROM _____ mg/L STOCK # _____
17529	N6120	.0028		BMDL	17866	P4851	.0166		BMDL	
17529	N6126	.0040		BMDL	17866	P4852	.0490		BMDL	
17531	P2710	.0024		BMDL	17866	P4853	.0446		BMDL	
17531	P2725	0		BMDL	17878	P2776	.00181		BMDL	
17539	P2117	.0043		BMDL	17866	P3120	.00114		BMDL	STANDARD mg/L
17539	P2116	.0954		BMDL	17866	P3122	0		BMDL	VALUE OBTAINED
17539	P2119	.0571		BMDL	17866	P3124	.0164		BMDL	% of theoretical
17524	P1711	.0043		BMDL	17866	P3126	0		BMDL	1.00
17524	P1713	.0040		BMDL	17882	P2543	.00472		BMDL	0.50
17524	P1714	0		BMDL	17882	P2544	.00213		BMDL	0.20
17524	P1717	.00114		BMDL	17882	P2545	.00473		BMDL	0.05
17562	N6128	0		BMDL	17882	P2547	.00784		BMDL	SLOPE
17562	N6148	.00119		BMDL	17882	P2549	.00119		BMDL	Intercept
17573	P2717	0		BMDL	17886	P3117	.0066		BMDL	Corr. Coef.
17579	P2726	.114		.12						ADDITIONAL COMMENTS <u>All calculations by X=MY+B.</u>
17584	P3068	.100		.10						
17503	P2683	.121		.12						
17562	N6124	.00357		BMDL						
17524	P1715	.00213		BMDL						
17864	P2531	.00119		BMDL						
17864	P2541	.0611		BMDL						
17864	P2546	.0691		BMDL						
17864	P2557	.00286		BMDL						

Analyses QA/QC Data Report

QC Batch # (N3004 N02)

Verified: 11/24/86 *[Signature]*

METHOD BLANK		PRECISION	
Spiked Blank (Known)	.20	Job # Dup (ETC)	N6110 P1714
Spiked Blank Result	198	Original Result	0 0
% Recovery	99	Duplicate Result	0 .0015
		R Range	0.0 .0015
EXTERNAL REF. QC	N/A	ACCURACY	
Known Value	.5	Job # Spiked (ETC)	P1714
Observed Result	-.537	Original Result	0
% Recovery		Amount Spiked	.0282
		Spiked S' Result	
		% Recovery	1.31

TEST: NITRATE UNITS: mg/L  
 Meth. REF. # 353.2 MDL: 0.10  
 Instrument # #1 (Channel 2)

BY: CH  
 BOOK # CV-12 P. 98  
 FROM PAGE #     

Lab #	Job #	ANALYZED CONC.	Dim	REPORTED CONC.	LogLink	Job #	ANALYZED CONC.	Dim	REPORTED CONC.	STANDARDS
7798	P2469	over		*	17866	P3117	1.15		1.15	PREPARED ON _____ BY _____ FROM _____ mg/L STOCK# _____
7798	P2471	over		*	17866	P3120	.0207		BMDL	
7798	P2524	over		*	17866	P3122	.0511		BMDL	
7798	P2538	over		*	17866	P3124	.0173		BMDL	
7798	P2539	over		*	17866	P3126	.0127		BMDL	
7814	P2540	over		*	17882	P2543	over		*	STANDARD mg/L    VALUE OBTAINED    % of theoretical
7814	P3555	.487		.49	17882	P2544	.0250		BMDL	
7822	P3107	over		*	17882	P2545	over		*	
17822	P3112	.007		BMDL	17882	P2547	over		*	
17828	P2473	over		*	17882	P2549	over		K	
7828	P2472	over		*	17882	P2550	over		K	SLOPE    Intercept    Corr. Coef.
17828	P2522	over		*	17882	P2551	over		*	
17828	P2536	over		*	17882	P2552	over		K	
17828	P2548	over		*						
7864	P2481	.006		BMDL						
17864	P2531	.0245		BMDL						ADDITIONAL COMMENTS * RPT on Channel 1 Complete Page 97
17864	P2541	1.73		1.73						
17864	P2546	over		*						
17864	P2554	2.06		2.06						
17866	P4850	over		*						
17866	P4851	over		*						
7866	P4852	.398		.40						
7866	P4853	over		*						
17878	P2770	.0462		BMDL						

Analyses QA/QC Data Report      QC Batch # N C 003      Verified: 11/24/86 [Signature]

PRECISION			ACCURACY			
METHOD BLANK	.003	.0177	Job # Dup (ETC)	P2473	P4850	P3126
Spiked Blank (Known)	.5	.5	Original Result	over	over	.0217
Spiked Blank Result	.501	.532	Duplicate Result	↓	↓	.0194
% Recovery	160%	1.06	A RANGE	↓	↓	
EXTERNAL REF. QC	EPA		Job # Spiked (ETC)	P2473	P4850	P3126
Known Value	93		Original Result	over	over	.021
Observed Result	93		Amount Spiked	↓	↓	.5
% Recovery	1.02		Spiked B' Result	↓	↓	.506
			% Recovery			97%

305182



TEST: CYANIDE UNITS: mg/L  
 Meth. REF. # 335.3 MDL: 0.025  
 Instrument # #1 (Channel 1)

BY: C.H.  
 BOOK # CV-12 P. 119  
 FROM PAGE #     

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.
17837	N4709	0		B=DL	17697	P4748	.005		B=DL
17837	N4714	0		B=DL	17697	P4748	.004		B=DL
17863	P0489	.007		B=DL	17697	P4748	.004		B=DL
17863	P0493	.009		B=DL	17698	P1710	.003		B=DL
17863	P0494	.005		B=DL	17701	N4710	.004		B=DL
17863	P0495	.002		B=DL	17701	N4715	.004		B=DL
17866	P4850	.539		.539	17866	P4851	over		Redo
17868	P4510	x		Redo	17866	P4852	over		Redo
17878	P2776	.007		B=DL	17866	P4853	x		Redo
17881	P1865	.002		B=DL	17868	P4510	x		Redo
17883	N4708	.002		B=DL	17868	P4511	x		Redo
17883	N4711	.003		B=DL	17868	P4512	x		Redo
17833	P4241	.0197		B=DL	17868	P4513	x		Redo
17833	P4242	.0169		B=DL	17883	N4712	x		Redo
17833	P4243	.0211		B=DL	17883	N4715	x		Redo
17833	P4244	.0186		B=DL	17889	P2754	x		Redo
17834	P4248	.0185		B=DL	17889	P2761	x		Redo
17834	P4249	.0188		B=DL	17889	P2774	x		Redo
17834	P4250	.0231		B=DL	17889	P2775	x		Redo
17834	P4251	.0229		B=DL	17833	P4235	x		Redo
<del>17834</del>	<del>P4224</del>	<del>.006</del>		<del>B=DL</del>	17833	P4226	x		Redo
17693	P2916	.0185		B=DL	17833	P4227	x		Redo
17693	P2917	.0312		B=DL	17833	P4228	x		Redo
17697	P4748	.006		B=DL	17443	P4239	x		Redo

PREPARED ON 12-2  
 BY P.B.  
 FROM      mg/L  
 STOCK #     

STANDARD mg/L	VALUE OBTAINED	% of Theoretical
.500	.493	98.6%
.400	.405	101.2%
.200	.205	102.5%
.100	.102	102%

SLOPE	Intercept	Corr. Coef.

**ADDITIONAL COMMENTS**  
 \* RPT DUE TO INSTRUMENTATION Problem  
 P4851 10x  
 P4852 20x

QC Batch # CY003

Verified: 12/3/86 *[Signature]*

Analyses QA/QC Data Report

METHOD BLANK	.008
Spiked Blank (Known)	.100
Spiked Blank Result	.103
% Recovery	103%
EXTERNAL REF. QC	CPA
Known Value	.224
Observed Result	.228
% Recovery	101.7%

PRECISION		Job # Dup (ETC)	N4709	P4748
Original Result	0		0.005	
Duplicate Result	0		0.005	
R TIME	0.0		0.001	
ACCURACY		Job # Spiked (ETC)	N4709	
Original Result	0.0			
Amount Spiked	.200			
Spiked S' Result	.201			
% Recovery	99%			

To Page No.

ANALYZED ON: 11/29/86  
 BY: P. Newman / R. Hill  
 BOOK # CV-12 P. 115  
 FROM PAGE # 114

TEST: Phenol, UNITS: mg/L  
 METH. REF. # 420.2, MDL: 0.050  
 INSTRUMENT # AA Channel #1

Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	STANDARDS
17639	P2121	0.037	X10	.375	1749	P2514	TH		MATRIX RPT STR	PREPARED ON _____ BY _____ FROM _____ mg/L STOCK # _____
↓	P2123	0.043	STR	BMDL						
17343	P4131	TH	X10	RPT						
17584	P3067	0.013	-	BMDL						
17639	N9882	TH	X5	RPT @ 1:100						
17681	N9884	TH	X5	RPT @ 1:100						
↓	N9880	TH	X5	RPT @ 1:100						
↓	P2614	TH	X5	STR/TH RPT						
17096	P0582	TH	-	STR RPT						
17824	P4180	TH	X5	RPT @ 1:10						
17828	P2472	TH	X5	RPT						
17844	P2531	TH	X5	RPT						
17868	P4512	TH	X5	RPT						
17866	P4851	0.067	X5	0.335						
17864	P2554	0.017	-	BMDL						
17866	P4851	0.013	-	BMDL						
↓	P4852	0.017	-	BMDL						
↓	P4853	0.022	-	BMDL						
17886	P5124	0.128	X10	1.28						
17707	N8672	0.048	-	BMDL						
↓	N8675	0.027	-	BMDL						
17721	N8674	0.131	X10	1.31						
1749	P2500	0.031	-	BMDL						
↓	P2513	0.096	X10	.964						

STANDARD mg/L	VALUE OBTAINED	# of Titrations
300		
200		
100		
050		

SLOPE	Intercept	Corr. Coef

ADDITIONAL COMMENT

Analyses QA/QC Data Report  
 SEE CV-12, P 114

QC Batch # PN003

Verified: [Signature] 11/29/86

METHOD BLANK		PRECISION	
Spiked Blank (Known)		Job # Dup (ETC)	
Spiked Blank Result		Original Result	
% Recovery		Duplicate Result	
		R Range	
		ACCURACY	
EXTERNAL REF. QC		Job # Spiked (ETC)	
Known Value		Original Result	
Observed Result		Amount Spiked	
% Recovery		Spiked S' Result	
		% Recovery	

305184

ANALYZED ON: 11/20/86

TEST: Phenol, UNITS: mg/L  
 METH. REF. # 420.2, MDL: 0.050  
 INSTRUMENT # AA Channel #1

BY: L. Venturini/R. Hillit  
 BOOK # CV-12 P. 115  
 FROM PAGE # 114

Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.
17639	P2121	0.037	X10	.375	1749	P2514	TH		MATRIX RPT STR
↓	P2123	0.043	STR	BMDL					
17343	P4131	TH	X10	RPT <sup>STR</sup>					
17584	P3067	0.013	-	BMDL					
17639	N9882	TH	X5	RPT <sup>1:100</sup>					
17681	N9884	TH	X5	RPT <sup>1:100</sup>					
↓	N9880	TH	X5	RPT <sup>1:100</sup>					
↓	P2614	TH	X5	RPT <sup>STR/TH</sup>					
17696	P0582	TH	-	RPT <sup>STR</sup>					
17824	P4480	TH	X5	RPT <sup>1:10</sup>					
17828	P2472	TH	X5	RPT					
17844	P253	TH	X5	RPT					
17868	P4512	TH	X5	RPT					
17868	P4851	0.067	X5	0.335					
17868	P2554	0.017	-	BMDL					
17868	P4851	0.013	-	BMDL					
↓	P4852	0.017	-	BMDL					
↓	P4853	0.022	-	BMDL					
17886	P5124	0.128	X10	1.28					
17707	N8672	0.048	-	BMDL					
↓	N8675	0.027	-	BMDL					
17721	N8674	0.131	X10	1.31					
1749	P2510	0.031	-	BMDL					
↓	P2513	0.096	X10	.964					

STANDARDS		
PREPARED ON _____		
BY _____		
FROM _____ mg/L		
STOCK # _____		
STANDARD mg/L	VALUE OBTAINED	% of Theoretical
300		
200		
100		
050		
SLOPE	Intercept	Corr. Coef.

ADDITIONAL COMMENTS

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Analyses QA/QC Data Report  
 SEE CV-12, P 114

QC Batch # PN003

Verified: [Signature] 11/29/86

METHOD BLANK				PRECISION			
Spiked Blank (Known)				Job # Dup (ETC)			
Spiked Blank Result				Original Result			
% Recovery				Duplicate Result			
				R Range			
EXTERNAL REF. QC				ACCURACY			
Known Value				Job # Spiked (ETC)			
Observed Result				Original Result			
% Recovery				Amount Spiked			
				Spiked S' Result			
				% Recovery			

Analyzed on: 11/29/86

TEST: Phenol, UNITS: mg/L  
 Meth. REF. # AA/EP4202, MDL: 0.50  
 Instrument # AA Channel 2

By: J. Hernandez/R. H. H.  
 Book # CV-12 P. 11  
 From Page #     

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	STANDARDS
17898	P2469	0.030	-	BMDL	17866	P4851	~	-	RPT	PREP ON _____ BY _____ FROM _____ mg/L STOCK # _____
	P2471	0.014	-	BMDL		P4852	~	-	RPT	
	P2524	0.012	-	BMDL		P4853	~	-	RPT	
	P2538	0.012	-	BMDL	17868	P4510	0.076	-	0.76	
	P2539	0.012	-	BMDL		P4511	0.054	-	0.54	STANDARD mg/L   VALUE OBTAINED   % of Theoretical
	P2540	0.014	-	BMDL		P4512	0.056	-	0.56	
17899	P3193	0.038	-	BMDL		P4513	0.017	-	BMDL	
	P3194	0.028	-	BMDL	17878	P2776	0.015	-	BMDL	
17801	P4447	0.004	-	BMDL	17886	P3120	0.027	-	BMDL	PKH 360 85.5 85.5%
17823	P4724	0.002	-	BMDL		P3122	0.010	-	BMDL	200 63.5 63.5%
17824	P4480	~	-	RPT		P3124	~	-	RPT	100 33.5 33.5%
17828	P2472	0.028	-	BMDL		P3126	0.084	-	0.84	50 18.5 18.5%
	P2473	0.017	-	BMDL	17882	P2543	0.036	-	BMDL	OFF SCALE
	P2522	0.017	-	BMDL		P2544	0.060	-	0.60	SLP
	P2536	0.025	-	BMDL		P2545	0.051	-	0.51	INT
	P2548	0.015	-	BMDL		P2547	0.017	-	BMDL	Y
17837	P4709	0.015	-	BMDL		P2549	0.015	-	BMDL	INTER
	P4714	0.023	-	BMDL		P2550	0.021	-	BMDL	REFERENCE
17864	P2481	0.025	-	BMDL		P2551	0.031	-	BMDL	P4850 - RPT for dilution
	P2531	0.006	-	BMDL		P2552	0.021	-	BMDL	P4851 -
	P2541	0.030	-	BMDL						P4852 -
	P2546	0.017	-	BMDL						P4853 -
	P2554	~	-	RPT						P3124 -
17866	P4850	0.321	-	RPT						ALL CALCS BY X=MY+B.

Analyses QA/QC Data Report

QC Batch # P1004

Verified: 11/29/86 JPH/Hee

PRECISION			ACCURACY				
METHOD BLANK	0.019		Job # Dup (ETC)	P2469	P4480	P3120	P2551
Spiked Blank (Known)	0.50	1.00	Original Result	0.030	MATRIX	0.028	0.036
Spiked Blank Result	0.045	0.095	Duplicate Result	0.030		0.026	0.026
% Recovery	90%	95%	R Range		INTER-	0.002	0.010
EXTERNAL REF. QC	N/A	N/A	Job # Spiked (ETC)	P2469	PKH	P3120	P2551
Known Value			Original Result	0.030	(CH)	0.027	0.031
Observed Result			Amount Spiked	0.100		100	100
% Recovery			Spiked S' Result	0.123		0.117	0.130
			% Recovery	91.2		90%	99%



REQUEST FOR ANALYSIS  
and SAMPLE CHAIN OF CUSTODY

LOGLINK: 17566

MATRIX: WATER

NAME OF SUBCONTRACT LAB: CHYUN

ETC JOB NUMBERS: P4850 P4853  
P4851  
P4852

TURNAROUND IN DAYS: NORMAL DATE DATA REQUIRED: 11/22  
(If deadline cannot be met, contact ETC Subcontract Group).

Send invoice, bill and reports to : ETC Subcontract Group  
(201)225-6786

Please perform the analyses requested below:

<input type="checkbox"/> Acidity	<input type="checkbox"/> Alkalinity	<input type="checkbox"/> Ammonia(probe)
<input type="checkbox"/> Ammonia(dist)	<input type="checkbox"/> Bicarbonate	<input type="checkbox"/> BOD(5 day)
<input type="checkbox"/> Bromide	<input type="checkbox"/> Carbonate	<input type="checkbox"/> Chloride
<input type="checkbox"/> Chem.Oxygen Dem.	<input type="checkbox"/> Chromium+6	<input type="checkbox"/> Fecal Coliform
<input type="checkbox"/> Total Coliform	<input type="checkbox"/> Color, apparent	<input type="checkbox"/> Cyanide, total
<input type="checkbox"/> Fluoride	<input type="checkbox"/> Formaldehyde(UV)	<input type="checkbox"/> Hardness
<input type="checkbox"/> Nitrate(NO3)	<input type="checkbox"/> Nitrite(NO2)	<input type="checkbox"/> TKN
<input type="checkbox"/> Odor	<input type="checkbox"/> Oil & Grease(grav)	<input type="checkbox"/> T.Organic Carb.
<input type="checkbox"/> Pet.Hydro(IR)	<input type="checkbox"/> Pet.Hydro(grav)	<input type="checkbox"/> Phenolics, tot.
<input type="checkbox"/> Phenolics(5ug/l)	<input type="checkbox"/> Phosphate(ortho)	<input type="checkbox"/> Phosphate(total)
<input type="checkbox"/> Phosphorus(tot.)	<input type="checkbox"/> Silica(dissolved)	<input checked="" type="checkbox"/> Sulfate(SO4)
<input type="checkbox"/> Sulfide(S)	<input type="checkbox"/> Sulfite(SO3)	<input type="checkbox"/> Surfactant(MBAS)
<input type="checkbox"/> Solids, Total	<input type="checkbox"/> Solids, Tot. Diss.	<input type="checkbox"/> Solids, Tot. Set.
<input checked="" type="checkbox"/> Solids, Tot. Sus.	<input type="checkbox"/> Solids, Tot. Vol.	<input type="checkbox"/> Turbidity
<input type="checkbox"/> Gross Alpha, Beta	<input type="checkbox"/> Radium 226	<input type="checkbox"/> Radium 228

Others: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

Chain of Custody Section for Courier Pick up at ETC

Sample(s) Relinquished by: (ETC) Mark Jekstadt  
Time: 4:30pm Date: 11/12/86

Sample(s) Received by: Mark Kelly  
Time: 4:30pm Date: 11/12/86

Chain of Custody Section for Sending Sample from ETC

Sample Shuttle sealed by : \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Sample Shuttle opened by : \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

Chain of Custody Section for Returning Sample to ETC

Sample Shuttle sealed by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Sample Shuttle opened by : \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

305188



**Technical Report**

for

**ENVIRONMENTAL STRATEGIES CORP.**

**WASHINGTON HARBOUR**

**3050 KK ST. SUITE 325**

**WASHINGTON, DC 20007**

*Chain of Custody Data Required for ETC Data Management Summary Reports*

<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>
P4852	ENVIRONMENTAL STRATEGIES CORP.	CHAMPION	WCSP-3	861110	1400	



**John J. Fitzgerald**

*Vice President  
Research and Operations*

This Technical Report is an INSITE<sup>SM</sup> service generated by LODESTAR<sup>SM</sup> Data Management Software.



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## Introduction

This report contains the analytical results on your sample. It is designed to include comprehensive data from the entire analytical process in order to satisfy the needs of various levels of review.

The results obtained from your sample are presented in tabular format immediately following this introduction. Quality assurance data is tabulated along with the appropriate sample results for verification. Depending on the analyses ordered, the quality assurance data may include results from blank, spiked blank, spiked sample (i.e. matrix spike) and replicate sample as well as results from surrogate compound analyses. Quality assurance data for verification of proper instrument performance is also included where appropriate. The report appendices include the chain of custody record for your sample and, where appropriate, the gas chromatograms and mass spectra.

The procedures used in the analysis of the sample are described in this report's methodology section. All analytical procedures within our laboratory are performed within a strictly enforced Quality Assurance Protocol. A description of this Protocol is included in the report.

## Results

Sample results, and associated quality assurance data, are always tabulated in one or more of this report's Quantitative Results Tables. The format of each table varies with the class of analysis.

### *Priority Pollutants*

The priority pollutant compounds and elements are listed with their NPDES (National Pollution Discharge Elimination System) numbers, and the Method Detection Limit (MDL) published in the Federal Register. When a compound or element is present below its published MDL it is reported as BMDL (Below Method Detection Limit). When a compound or element is not present at any detectable concentrations it is reported as ND (Not Detected). MDL's for non-aqueous matrices are based on USEPA published MDL's but are adjusted as per sample weight. Matrix spike and replicate analyses, where included, were performed on samples randomly chosen within each quality assurance batch and are therefore not necessarily spikes and replicates of this report's sample. Surrogate compound recovery data and instrument calibration data are included in the Method Performance Data Tables.

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Volatile Compounds - GC/MS Analysis Data (QR01)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4852 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WOSP-3 861110 1400  
 ETC Sample No. Company Sample Point Date Time Elapsed hours

NPDES Number	Compound <small>Acetone and hexamethylenetetramine values are screen only.</small>	Results			QC Replicate		QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concn. Added ug/l	% Recov	Unspiked Sample ug/l	Concn. Added ug/l	% Recov	
1V	Acrolein	ND	100	ND	ND	ND	800	88	ND	800	98	
2V	Acrylonitrile	ND	100	ND	ND	ND	80.0	94	ND	80.0	84	
3V	Benzene	ND	4.4	ND	ND	ND	18.0	99	ND	18.0	100	
4V	bis(Chloromethyl) ether	ND	10	ND	ND	ND	18.0	92	ND	18.0	90	
5V	Bromoform	ND	4.7	ND	ND	ND	18.0	103	ND	18.0	98	
6V	Carbon tetrachloride	ND	2.8	ND	ND	ND	18.0	99	ND	18.0	99	
7V	Chlorobenzene	ND	6.0	ND	ND	ND	18.0	99	ND	18.0	95	
8V	Chlorodibromomethane	ND	3.1	ND	ND	ND	18.0	103	ND	18.0	96	
9V	Chloroethane	ND	10	ND	ND	ND	18.0	0 <sup>a</sup>	ND	18.0	0 <sup>a</sup>	
10V	2-Chloroethylvinyl ether	ND	10	ND	ND	ND	18.0	103	ND	18.0	98	
11V	Chloroform	ND	1.6	ND	ND	ND	18.0	100	ND	18.0	96	
12V	Dichlorobromomethane	ND	2.2	ND	ND	ND	18.0	109	ND	18.0	104	
13V	Dichlorodifluoromethane	ND	10	ND	ND	ND	18.0	100	ND	18.0	99	
14V	1,1-Dichloroethane	5.88	4.7	ND	ND	ND	18.0	100	ND	18.0	96	
15V	1,2-Dichloroethane	3.44	2.8	ND	ND	ND	18.0	104	ND	18.0	99	
16V	1,1-Dichloroethylene	ND	6.0	ND	ND	ND	18.0	98	ND	18.0	97	
17V	1,2-Dichloropropane	ND	5.0	ND	ND	ND	18.0	97	ND	18.0	101	
18V	cis-1,3-Dichloropropylene	ND	7.2	ND	ND	ND	18.0	100	ND	18.0	100	
19V	Ethylbenzene	ND	10	ND	ND	ND	18.0	103	ND	18.0	109	
20V	Methyl bromide	ND	10	ND	ND	ND	18.0	107	ND	18.0	109	
21V	Methyl chloride	ND	10	209	206	ND	18.0	98	45.0	18.0	89	
22V	Methylene chloride	ND	2.8	ND	ND	ND	18.0	101	ND	18.0	106	
23V	1,1,2,2-Tetrachloroethane	ND	6.9	ND	ND	ND	18.0	98	ND	18.0	103	
24V	Tetrachloroethylene	ND	4.1	ND	ND	ND	18.0	100	ND	18.0	102	
25V	Toluene	1040	6.0	ND	ND	ND	18.0	97	ND	18.0	99	
26V	1,2-Trans-dichloroethylene	5.33	1.6	ND	ND	ND	18.0	104	ND	18.0	97	
27V	1,1-Trichloroethane	ND	3.8	ND	ND	ND	18.0	104	ND	18.0	100	
28V	1,1,2-Trichloroethane	167	5.0	ND	ND	ND	18.0	93	ND	18.0	87	
29V	Trichloroethylene	513	1.9	ND	ND	ND	18.0	105	ND	18.0	99	
30V	Trichlorofluoromethane	ND	10	ND	ND	ND	18.0	110	ND	18.0	99	
31V	Vinyl chloride	ND	10	ND	ND	ND	18.0	94	ND	18.0	96	
18V	trans-1,3-Dichloropropylene	ND	10	ND	ND	ND	18.0		ND	18.0		

<sup>a</sup> Recovery normally variable using established methodology.

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**ETC** ENVIRONMENTAL TESTING and CERTIFICATION

NOV 29, 1986

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Acid Compounds - GC/MS Analysis Data (QR02)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4852** ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-3 861110 1400  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results		QC Replicate		QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concen. Added ug/l	% Recov	Unspiked Sample ug/l	Concen. Added ug/l	% Recov
1A	2-Chlorophenol	ND	3.5	ND	ND	ND	100	89	ND	102	107
2A	2,4-Dichlorophenol	ND	2.9	ND	ND	ND	100	92	ND	102	100
3A	2,4-Dimethylphenol	ND	2.9	ND	ND	ND	100	18 <sup>a</sup>	ND	102	100
4A	4,6-Dinitro-o-cresol	ND	26	ND	ND	ND	100	92	ND	102	90
5A	2,4-Dinitrophenol	ND	45	ND	ND	ND	100	77	ND	102	88
6A	2-Nitrophenol	ND	9	ND	ND	ND	100	80	ND	102	89
7A	4-Nitrophenol	ND	3.6	ND	ND	ND	100	85	ND	102	97
8A	p-Chloro-m-cresol	ND	3.2	ND	ND	ND	100	77	ND	102	68
9A	Pentachlorophenol	ND	3.9	ND	ND	ND	100	74	ND	102	82
10A	Phenol	ND	1.6	ND	ND	ND	100	89	ND	102	109
11A	2,4,6-Trichlorophenol	ND	2.9	ND	ND	ND	100	89	ND	102	94

<sup>a</sup> Recovery normally variable using established methodology.

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA  
BASE/NEUTRAL COMPOUNDS - GC/MS ANALYSIS DATA (QR03)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4852** ENVIRONMENTAL STRATEGIES CORP. CHAMPION WOSP-3 861110 1400  
 ETC Sample No. Company Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results			QC Replicate		QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concn. Added ug/l	% Recov	Unspiked Sample ug/l	Concn. Added ug/l	% Recov	
1B	Acenaphthene	ND	2.0	ND	ND	ND	100	87	ND	102	87	
2B	Acenaphthylene	ND	3.8	ND	ND	ND	100	90	ND	102	92	
3B	Anthracene	ND	2.0	ND	ND	ND	100	84	ND	102	86	
4B	Benzo(a)anthracene	ND	47	ND	ND	ND	100	0 <sup>a</sup>	ND	102	3 <sup>a</sup>	
5B	Benzo(a)pyrene	ND	8.4	ND	ND	ND	100	89	ND	102	90	
6B	Benzo(b)fluoranthene	ND	2.7	ND	ND	ND	100	98	ND	102	99	
7B	Benzo(k)fluoranthene	ND	11	ND	ND	ND	100	109	ND	102	105	
8B	Benzo(g)h)perylene	ND	4.4	ND	ND	ND	0	-	ND	0	-	
9B	Benzo(k)fluoranthene	ND	3.8	ND	ND	ND	100	78	ND	102	87	
10B	bis(2-Chloroethoxy)methane	ND	5.7	ND	ND	ND	100	80	ND	102	83	
11B	bis(2-Chloroethyl) ether	ND	6.1	ND	ND	ND	100	87	ND	102	89	
12B	bis(2-Chloroisopropyl) ether	ND	6.1	ND	ND	ND	100	88	ND	102	92	
13B	bis(2-Ethylhexyl)phthalate	ND	11	ND	ND	ND	100	92	ND	102	93	
14B	4-Bromophenyl phenyl ether	ND	2.0	ND	ND	ND	100	99	ND	102	95	
15B	Butyl benzyl phthalate	ND	11	ND	ND	ND	100	75	ND	102	78	
16B	2-Chloronaphthalene	ND	2.0	ND	ND	ND	100	87	ND	102	85	
17B	4-Chlorophenyl phenyl ether	ND	4.5	ND	ND	ND	100	87	ND	102	91	
18B	Chrysene	ND	2.7	ND	ND	ND	100	88	ND	102	88	
19B	Dibenzo(a,h)anthracene	ND	11	ND	ND	ND	0	-	ND	0	-	
20B	1,2-Dichlorobenzene	ND	2.0	ND	ND	ND	100	76	ND	102	76	
21B	1,3-Dichlorobenzene	ND	2.0	ND	ND	ND	100	67	ND	102	72	
22B	1,4-Dichlorobenzene	ND	4.7	ND	ND	ND	100	71	ND	102	75	
23B	3,3'-Dichlorobenzidine	ND	18	ND	ND	ND	100	52	ND	102	54	
24B	Diethyl phthalate	ND	11	ND	ND	ND	100	45	ND	102	48	
25B	Dimethyl phthalate	ND	11	ND	ND	ND	100	22 <sup>a</sup>	ND	102	18 <sup>a</sup>	
26B	Di-n-butyl phthalate	ND	11	ND	ND	ND	100	63	ND	102	61	
27B	2,4-Dinitrotoluene	ND	6.1	ND	ND	ND	100	81	ND	102	110	
28B	2,6-Dinitrotoluene	ND	2.0	ND	ND	ND	100	94	ND	102	106	
29B	Di-n-octyl phthalate	ND	11	ND	ND	ND	100	101	ND	102	103	
30B	1,2-Diphenylhydrazine	ND	11	ND	ND	ND	100	88	ND	102	96	
31B	Fluoranthene	ND	2.4	ND	ND	ND	100	70	ND	102	58	
32B	Fluorene	ND	2.0	ND	ND	ND	100	88	ND	102	95	

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA  
BASE/NEUTRAL COMPOUNDS - GC/MS ANALYSIS DATA (QR03)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4852** ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-3 861110 1400  
 ETC Sample No. Company Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results			QC Replicate			QC Blank and Spiked Blank			QC Matrix Spike			
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concen. Added ug/l	% Recov	Unspiked Sample ug/l	Concen. Added ug/l	% Recov			
33B	Hexachlorobenzene	ND	2.0	ND	ND	ND	100	93	ND	102	94	ND	102	94
34B	Hexachlorobutadiene	ND	1.97	ND	ND	ND	100	71	ND	102	75	ND	102	75
35B	Hexachlorocyclopentadiene	ND	1.7	ND	ND	ND	0	-	ND	0	-	ND	0	-
36B	Hexachloroethane	ND	5.1	ND	ND	ND	100	62	ND	102	63	ND	102	63
37B	Indeno(1,2,3-c,d)pyrene	ND	2.4	ND	ND	ND	0	-	ND	0	-	ND	0	-
38B	Isophorone	ND	1.7	ND	ND	ND	100	76	ND	102	80	ND	102	80
39B	Naphthalene	ND	2.0	ND	ND	ND	100	80	ND	102	84	ND	102	84
40B	Nitrobenzene	ND	1.7	ND	ND	ND	100	89	ND	102	88	ND	102	88
41B	N-Nitrosodimethylamine	ND	11	ND	ND	ND	0	-	ND	0	-	ND	0	-
42B	N-Nitrosodi-n-propylamine	ND	11	ND	ND	ND	100	65	ND	102	85	ND	102	85
43B	N-Nitrosodiphenylamine	ND	2.0	ND	ND	ND	100	83	ND	102	110	ND	102	110
44B	Phenanthrene	ND	5.8	ND	ND	ND	100	89	ND	102	89	ND	102	89
45B	Pyrene	ND	2.0	ND	ND	ND	100	69	ND	102	55	ND	102	55
46B	1,2,4-Trichlorobenzene	ND	2.0	ND	ND	ND	100	158 <sup>a</sup>	ND	100	152 <sup>a</sup>	ND	102	152 <sup>a</sup>

<sup>a</sup> Recovery normally variable using established methodology.

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**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Pesticide/PCB Compounds - GC/MS Analysis Data (QR04)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4852 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-3 861110 1400  
 ETC Sample No. Company Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results		QC Replicate		QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concen. Added ug/l	% Retov	Unspiked Sample ug/l	Concen. Added ug/l	% Recov
1P	Aldrin	ND	2.0	ND	ND	ND	100	89	ND	102	55
2P	Alpha-BHC	ND	11	ND	ND	ND	100	39	ND	102	34
3P	Beta-BHC	ND	4.7	ND	ND	ND	100	57	ND	102	67
4P	Gamma-BHC	ND	11	ND	ND	ND	100	34	ND	102	32
5P	Delta-BHC	ND	3.3	ND	ND	ND	100	46	ND	102	44
6P	Chlordane	ND	11	ND	ND	ND	200	69	ND	204	56
7P	4,4'-DDT	ND	3.0	ND	ND	ND	100	72	ND	102	36
8P	4,4'-DDE	ND	6.0	ND	ND	ND	100	84	ND	102	47
9P	4,4'-DDD	ND	5.1	ND	ND	ND	100	71	ND	102	46
10P	Dieldrin	ND	2.7	ND	ND	ND	100	83	ND	102	82
11P	Endosulfan I	ND	11	ND	ND	ND	100	29	ND	102	11 <sup>a</sup>
12P	Endosulfan II	ND	11	ND	ND	ND	100	30	ND	102	8 <sup>a</sup>
13P	Endosulfan sulfate	ND	6.0	ND	ND	ND	100	44	ND	102	44
14P	Endrin	ND	11	ND	ND	ND	100	74	ND	102	72
15P	Endrin aldehyde	ND	11	ND	ND	ND	100	29	ND	102	30
16P	Heptachlor	ND	2.0	ND	ND	ND	100	97	ND	102	68
17P	Heptachlor epoxide	ND	2.4	ND	ND	ND	100	86	ND	102	106
18P	PCB-1242	ND	39	ND	ND	ND	0	-	ND	0	-
19P	PCB-1254	ND	39	ND	ND	ND	0	-	ND	0	-
20P	PCB-1221	ND	39	ND	ND	ND	0	-	ND	0	-
21P	PCB-1232	ND	39	ND	ND	ND	0	-	ND	0	-
22P	PCB-1248	ND	39	ND	ND	ND	0	-	ND	0	-
23P	PCB-1260	ND	39	ND	ND	ND	0	-	ND	0	-
24P	PCB-1016	ND	39	ND	ND	ND	100	93	ND	102	49
25P	Toxaphene	ND	11	ND	ND	ND	0	-	ND	0	-

<sup>a</sup> Recovery normally variable using established methodology.

**TABLE 1: QUANTITATIVE RESULTS**  
**Metals - Analysis Data (QR52)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4852 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-3 861110 1400  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

Compound	Results	
	Sample Concn. ug/l	MDL ug/l
Barium	52	3.2
Chromium	ND	26
Copper	BMDL	13
Iron	BMDL	140
Manganese	500	6.4
Nickel	18	11
Sodium	191000	72
Zinc	150	2.9

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**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Conventional Analysis Data (QR10)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4852 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-3 861110 1400  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

NPDES Number	Results	
	Sample Concn. mg/l	MDL mg/l
Fluoride	2.53	.10
Nitrate as N	.40	.10
Sulfate as SO4	180	2
Phenolics, Total	< .050	.050
Cyanide, Total	1.08	.025
Solids, total suspended	17300	2

**TABLE 2: METHOD PERFORMANCE DATA**  
**Surrogate Recovery - Aqueous Matrices (QR20)**

**Chain of Custody Data Required for ETC Data Management Summary Reports**

P4852

ETC Sample No.      Company      Facility      Sample Point      Date      Time      Elapsed Hours

Compound	Amount Added ug	% Recovery	Control Limits %	
			Lower	Upper
<b>VOLATILE FRACTION (GC/MS)</b>				
Toluene-D8	.250	95	88	110
Bromofluorobenzene	.250	97	86	115
1,2-Dichloroethane-D4	.250	100	76	114
<b>ACID FRACTION (GC/MS)</b>				
Phenol-D5	100	98	10	94
2-Fluorophenol	100	102	21	100
2,4,6-Tribromophenol	100	78	10	123
<b>BASE/NEUTRAL FRACTION (GC/MS)</b>				
Nitrobenzene-D5	50	99	35	114
2-Fluorobiphenyl	50	80	43	116
Terphenyl-D14	50	24***	33	141
<b>PESTICIDE/PCB FRACTION (GC/MS)</b>				
Dibutylchloroendate	-	-	24**	154**

\* IFB EPA Control Limits.  
 \*\* Advisory Limits Only.  
 \*\*\* Recovery manually verified.

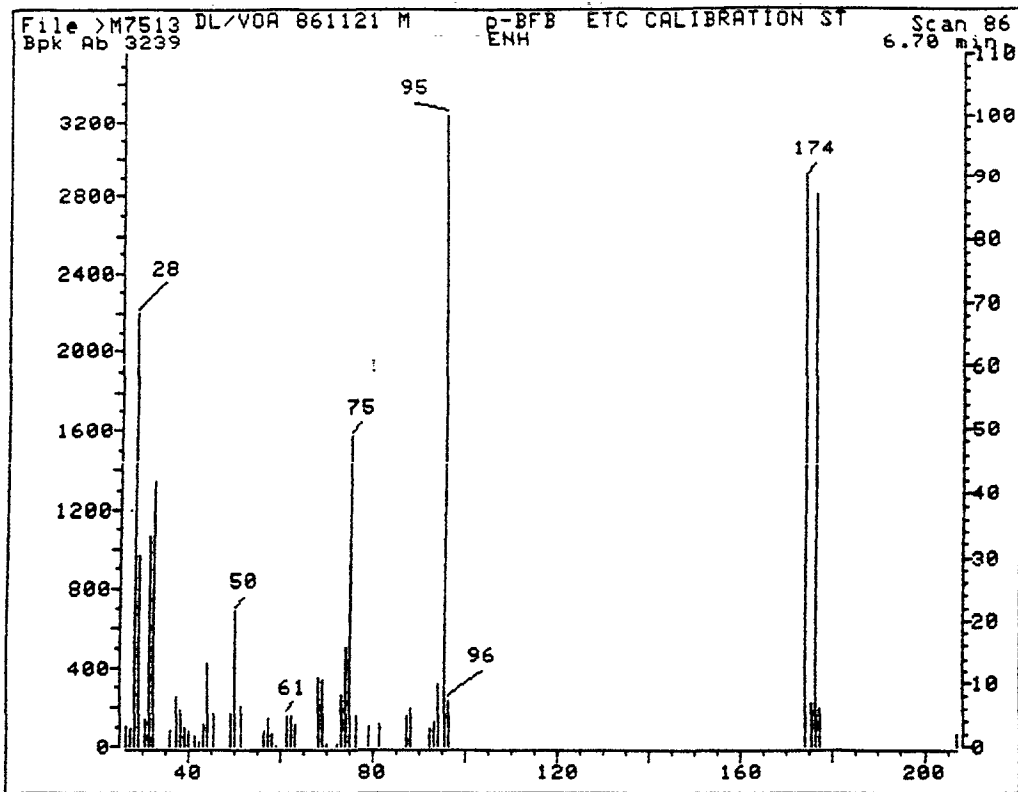


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance		Status
		Base Peak	Appropriate Peak	
50	15-40% of mass 95	21.09	21.09	Ok
75	30-60% of mass 95	48.39	48.39	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.22	7.22	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	89.98	89.98	Ok
175	5-9% of mass 174	6.74	7.49	Ok
176	95-101% of mass 174	87.27	96.99	Ok
177	5-9% of mass 176	5.98	6.86	Ok

Injection Date: 11/21/86  
 Injection Time: 20:42  
 Run No: >M7513  
 Spectrun No: 86

Analyst: J. Owen  
 Processor: ELUON COORDINATOR  
 QC Batch: QV 5431  
 Samples: P4240, P4241, P4242, P4243, P4244  
P4236, P4237, P4239, P4235, P533  
P5415, P5414, P4850, P4851, P485

*Handwritten signature and date:* 11/25/86

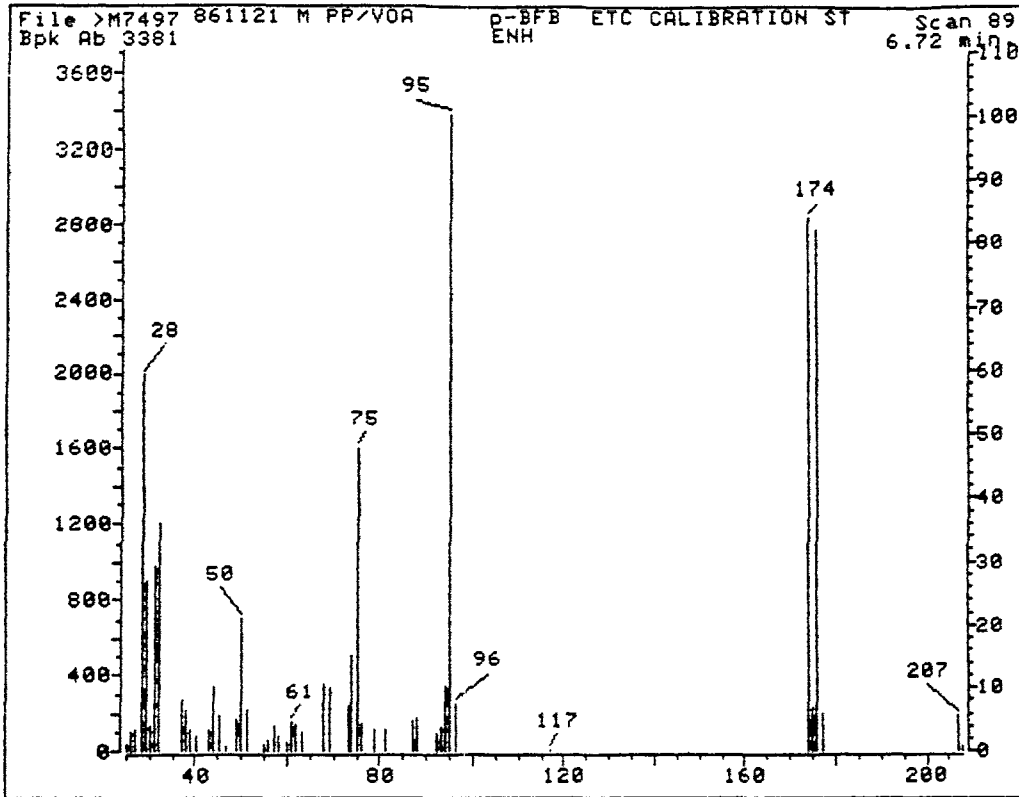


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance		Status
		Base Peak	Appropriate Peak	
50	15-40% of mass 95	20.80	20.80	Ok
75	30-60% of mass 95	47.52	47.52	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.18	7.18	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	83.99	83.99	Ok
175	5-9% of mass 174	6.60	7.85	Ok
176	95-101% of mass 174	82.16	97.82	Ok
177	5-9% of mass 176	5.86	7.13	Ok

Injection Date: 11/21/86  
 Injection Time: 05:57  
 Run No: >M7497  
 Spectrun No: 89

Analyst: *Joseph Quinn*  
 Processor: *BUDD CHRISTOPHER*  
 QC Batch: *QV5931*  
 Samples: *P4240, P4241, P4242, P4243, P4244, P4236, P4237, P4239, P4235, P533, P5415, P5414, P4850, P4851, P4852*

*JQ*  
 005202

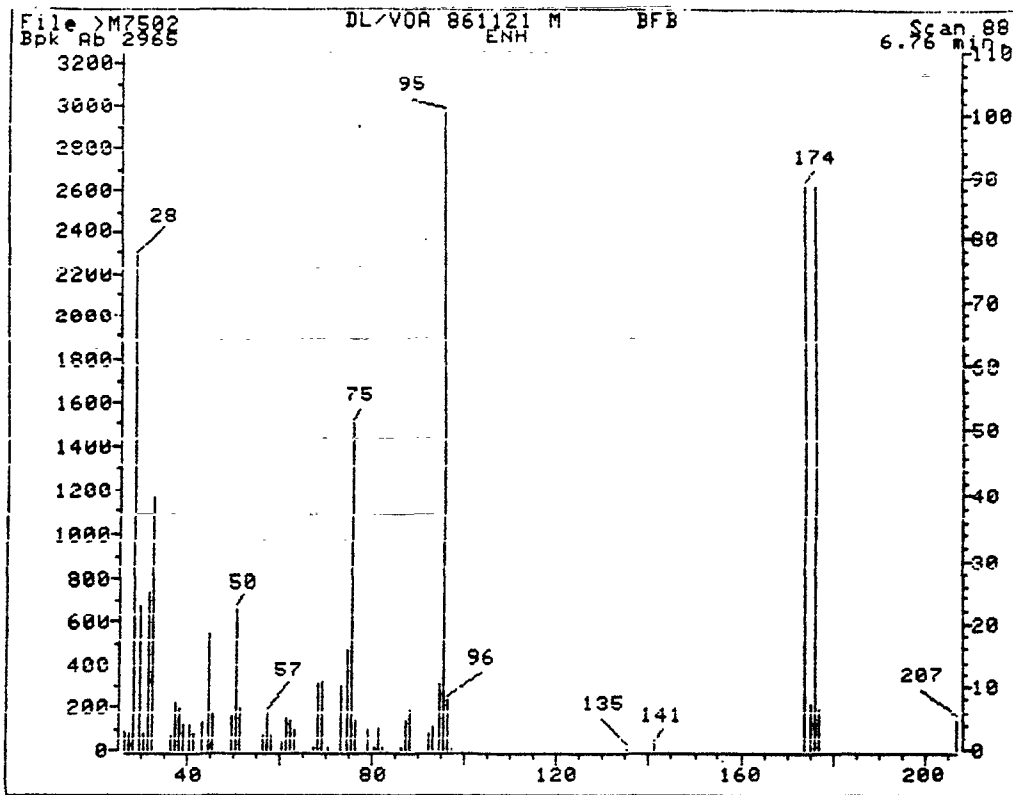


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	22.00	22.00	Ok
75	30-60% of mass 95	50.62	50.62	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.63	7.63	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	88.52	88.52	Ok
175	5-9% of mass 174	7.03	7.95	Ok
176	95-101% of mass 174	88.43	99.89	Ok
177	5-9% of mass 176	6.31	7.15	Ok

Injection Date: 11/21/86  
 Injection Time: 09:32  
 Run No: >M7502  
 Spectrum No: 88

Analyst: *Nancy Albert*  
 Processor: *Quinn Christopher*  
 QC Batch: *QV 5931*  
 Samples: *P4240, P4241, P4242, P4243, P4244, P4236, P4237, P4238, P4235, P5331, P5415, P5414, P4850, P4851, P4852*

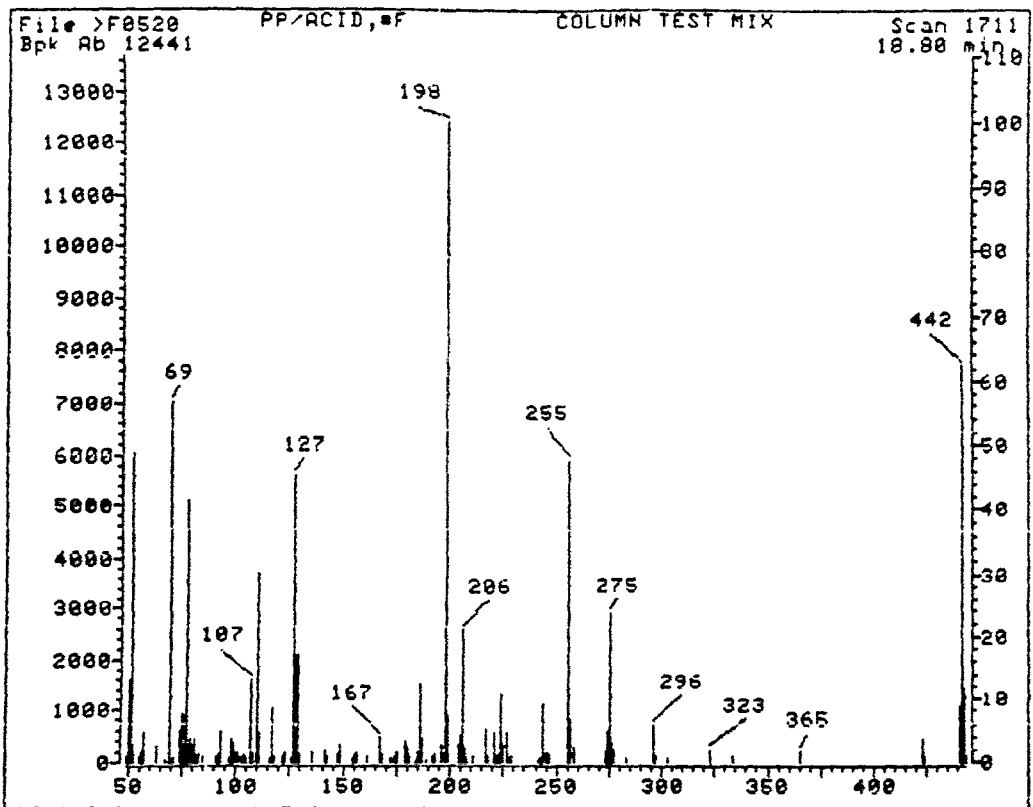


TABLE 2: METHOD PERFORMANCE DATA (QR22)

GC/MS Tuning Data - Decafluorotriphenylphospine (DFTPP) for Acids Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	Appropriate Peak	Status
51	30-60% of mass 198	48.52	48.52	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	56.51	56.51	Ok
70	Less than 2% of mass 69	.42	.74	Ok
127	40-60% of mass 198	45.35	45.35	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	7.19	7.19	Ok
275	10-30% of mass 198	23.50	23.50	Ok
365	Greater than 1% of mass 198	1.89	1.89	Ok
441	0-100% of mass 443	8.79	76.34	Ok
442	Greater than 40% of mass 198	62.49	62.49	Ok
443	17-23% of mass 442	11.52	18.43	Ok

Injection Date: 11/26/86  
 Injection Time: 07:29  
 Run No: >F0520  
 Spectrun No: 1711

Analyst: J. AO  
 Processor: Michael Trank  
 QC Batch: 046013  
 Samples: 00447-00450, 00457, 00458,  
00467, 00468, 00471, 00472,  
00489, 00493-00495, 04852,  
04853, 00472

*KS*  
*11/29/86*

305204

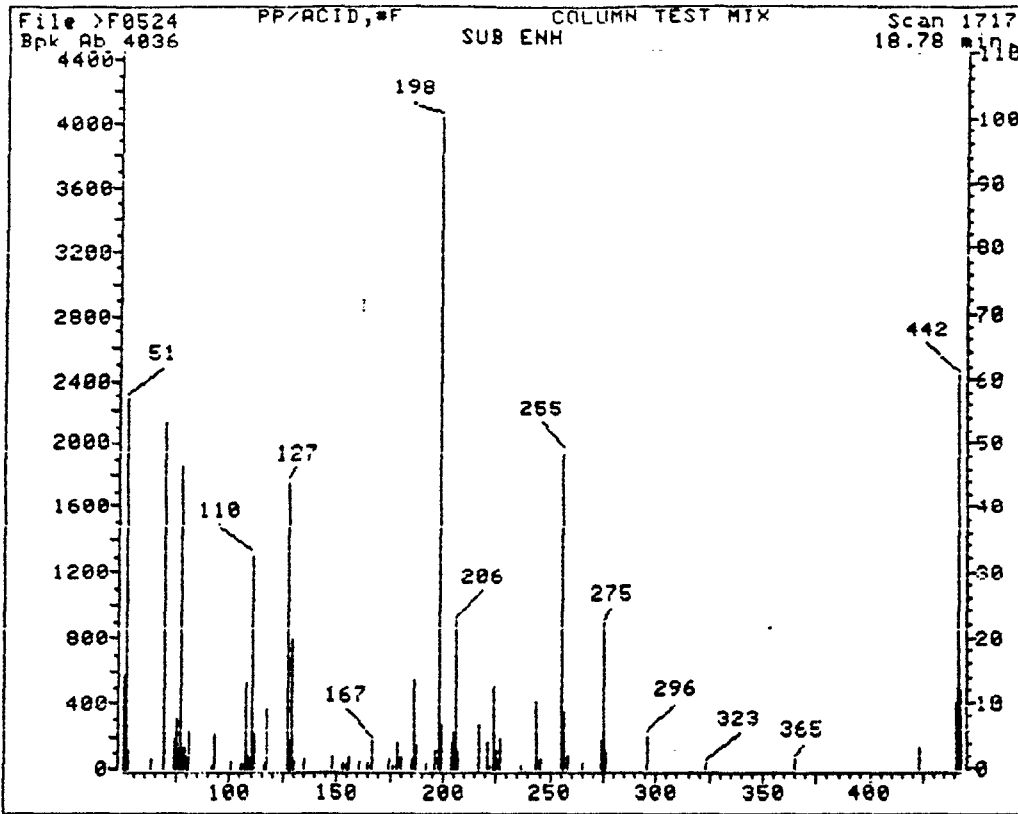


TABLE 2: METHOD PERFORMANCE DATA (QR22)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFIPP) for Acids Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	56.57	56.57	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	52.49	52.49	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	43.39	43.39	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	6.54	6.54	Ok
275	10-30% of mass 198	21.76	21.76	Ok
365	Greater than 1% of mass 198	1.26	1.26	Ok
441	0-100% of mass 443	10.00	84.54	Ok
442	Greater than 40% of mass 198	60.51	60.51	Ok
443	17-23% of mass 442	11.83	19.55	Ok

Injection Date: 11/26/86  
 Injection Time: 11:04  
 Run No: >F0524  
 Spectrun No: 1717

Analyst: J. A. O.  
 Processor: Paul Trant  
 QC Batch: QA6013  
 Samples: 80447-80450, 80452, 80453,  
80457, 80458, 80467, 80468,  
80471, 80472.

*KS*  
*11/25/86*

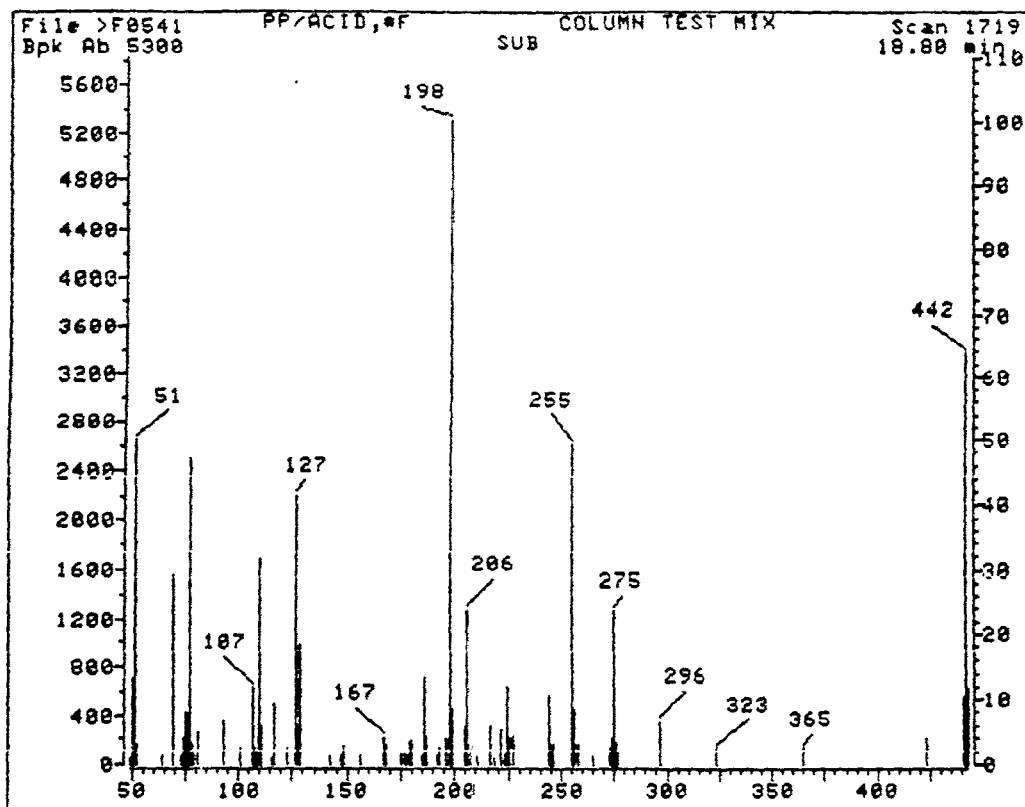


TABLE 2: METHOD PERFORMANCE DATA (QR22)

GC/MS Tuning Data - Decafluorotriphenylphospine (DFTPP) for Acids Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	49.92	49.92	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	29.34	29.34	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	41.43	41.43	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	8.38	8.38	Ok
275	10-30% of mass 198	23.66	23.66	Ok
365	Greater than 1% of mass 198	2.28	2.28	Ok
441	0-100% of mass 443	10.34	88.96	Ok
442	Greater than 40% of mass 198	63.66	63.66	Ok
443	17-23% of mass 442	11.62	18.26	Ok

KS  
11/29/86

Injection Date: 11/26/86  
 Injection Time: 22:39  
 Run No: >F0541  
 Spectrun No: 1719

Analyst: Stephen AS  
 Processor: Phil Frank  
 QC Batch: 0A0013  
 Samples: 00409, 00403, 00404, 01495

305206



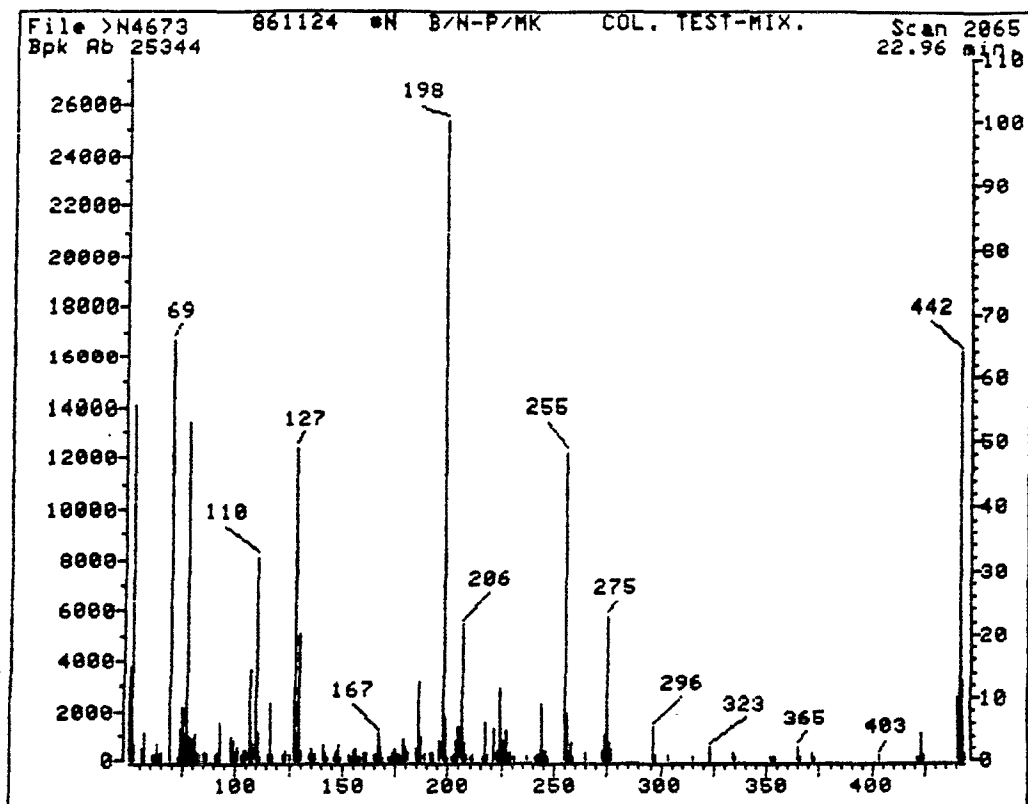


TABLE 2: METHOD PERFORMANCE DATA (QR23)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFTPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	55.50	55.50	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	65.88	65.88	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	48.83	48.83	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	6.67	6.67	Ok
275	10-30% of mass 198	22.60	22.60	Ok
365	Greater than 1% of mass 198	2.14	2.14	Ok
441	0-100% of mass 443	9.87	78.53	Ok
442	Greater than 40% of mass 198	64.00	64.00	Ok
443	17-23% of mass 442	12.57	19.64	Ok

Injection Date: 11/25/86  
Injection Time: 14:37  
Run No: >N4673  
Spectrum No: 2065

Analyst: J. Hendell  
Processor: W. J. C. C.  
QC Batch: QB6013  
Samples: 20447-20450, 20457, 20458  
20457, 20458, 20471, 20472, 20489  
20493-20495, 20852, 20853

LAN 11-30

305207

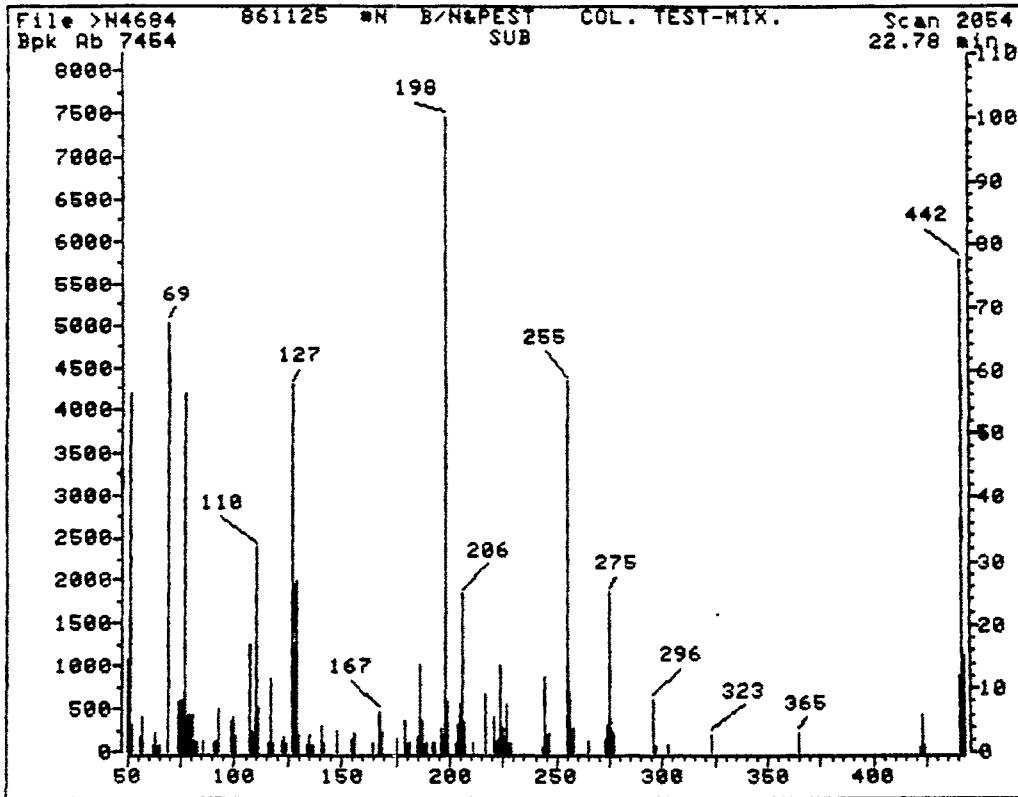


TABLE 2: METHOD PERFORMANCE DATA (QR23)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFTPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	56.12	56.12	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	67.41	67.41	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	57.59	57.59	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	7.90	7.90	Ok
275	10-30% of mass 198	25.03	25.03	Ok
365	Greater than 1% of mass 198	3.03	3.03	Ok
441	0-100% of mass 443	12.03	79.24	Ok
442	Greater than 40% of mass 198	77.68	77.68	Ok
443	17-23% of mass 442	15.19	19.55	Ok

Injection Date: 11/26/86  
 Injection Time: 03:17  
 Run No: >N4684  
 Spectrum No: 2054

Analyst: *[Signature]*  
 Processor: W.A.C.L.  
 QC Batch: QB6013  
 Samples: P0447, P0450, P0450, P0451, P0451, P0467, P0468, P0471, P0472, P0472, P0493-P0495, P4852, P4853

305208

LAH  
11-30

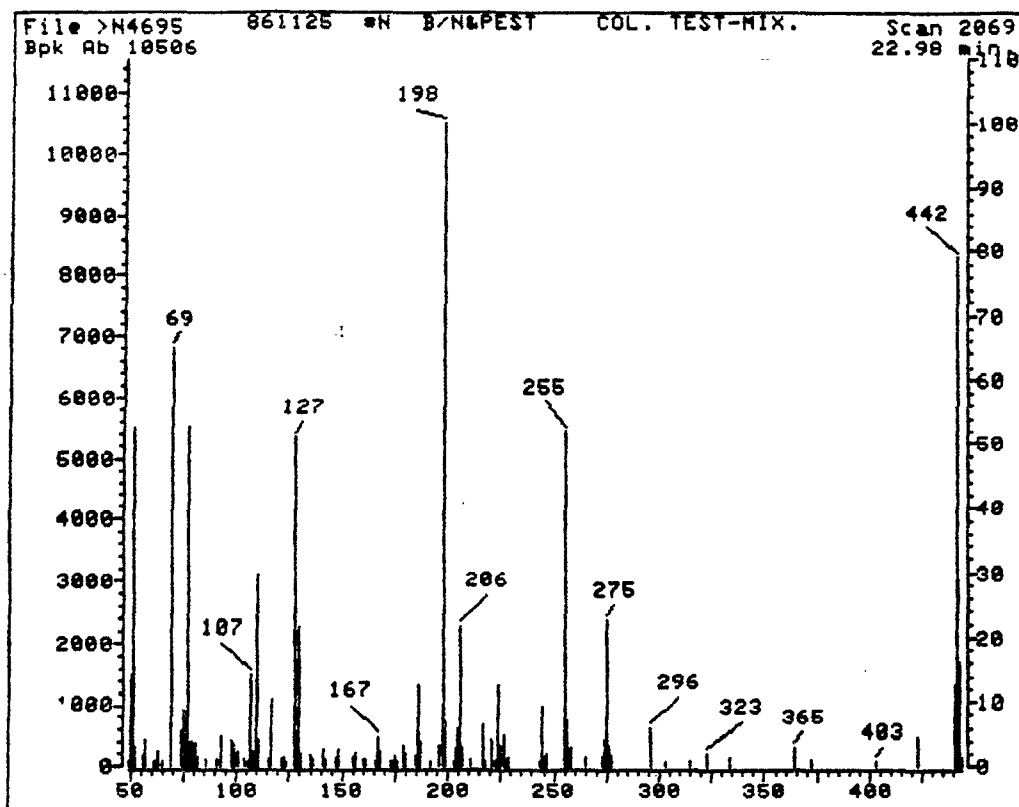


TABLE 2: METHOD PERFORMANCE DATA (QR23)

GC/MS Tuning Data - Decafluorotriphenylphospine (DFTPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	52.28	52.28	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	64.83	64.83	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	50.89	50.89	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	6.94	6.94	Ok
275	10-30% of mass 198	22.61	22.61	Ok
365	Greater than 1% of mass 198	3.04	3.04	Ok
441	0-100% of mass 443	12.55	77.32	Ok
442	Greater than 40% of mass 198	79.16	79.16	Ok
443	17-23% of mass 442	16.24	20.51	Ok

Injection Date: 11/26/86  
 Injection Time: 12:15  
 Run No: >N4695  
 Spectrun No: 2069

Analyst: V. [Signature]  
 Processor: W. D. C. C.  
 QC Batch: QB6013  
 Samples: P0447-P0450, P0457, P0458  
P0467, P0468.

LAA  
 11-30

305209

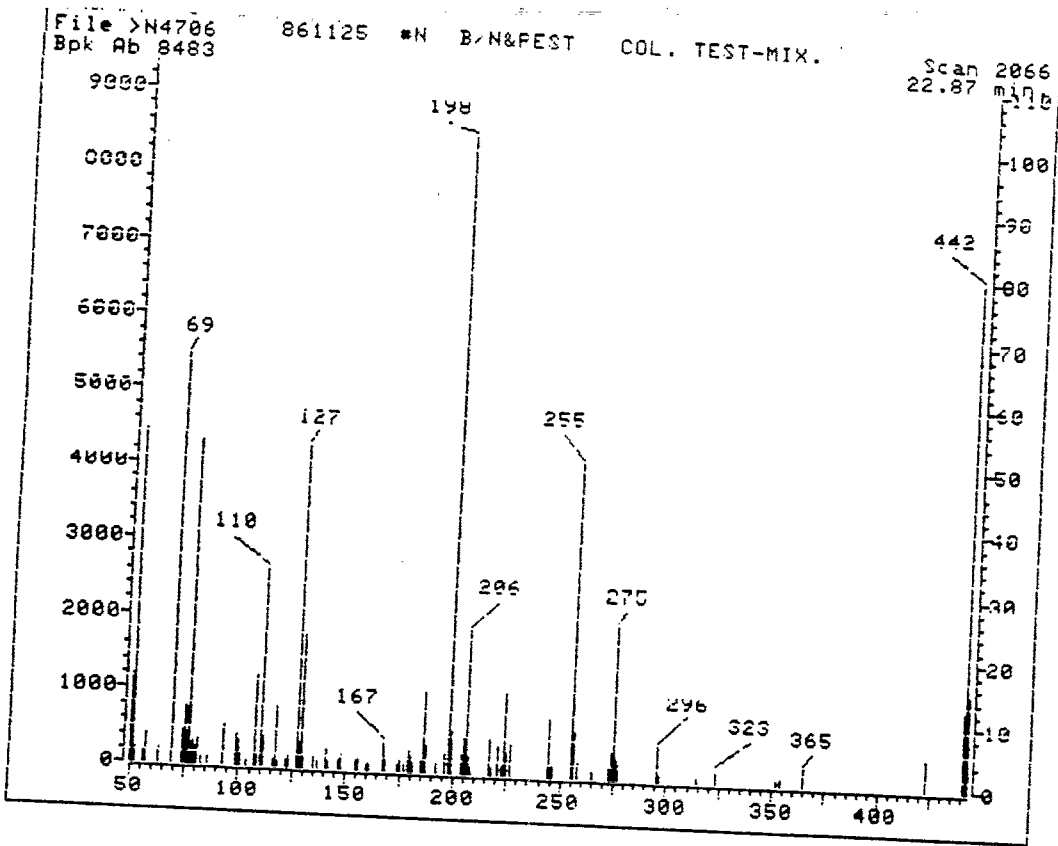


TABLE 2: METHOD PERFORMANCE DATA (WR23)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFPPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	52.22	52.22	UK
68	Less than 2% of mass 69	0.00	0.00	UK
69	(reference only)	64.29	64.29	UK
70	Less than 2% of mass 69	0.00	0.00	UK
127	40-60% of mass 198	50.49	50.49	UK
197	Less than 1% of mass 198	0.00	0.00	UK
198	Base peak, 100% relative abundance	100.00	100.00	UK
199	5-9% of mass 198	6.64	6.64	UK
275	10-30% of mass 198	24.07	24.07	UK
365	Greater than 1% of mass 198	2.91	2.91	UK
441	0-100% of mass 443	12.58	82.33	UK
442	Greater than 40% of mass 198	19.88	19.88	UK
443	1/-23% of mass 442	15.28	19.13	UK

Injection Date: 11/26/86  
Injection Time: 21:29  
Run No: >N4/U6  
Spectrum No: 2066

Analyst: Sharon Malachuk  
Processor: A. N. C.  
QC Batch: 086013  
Samples: P0471, P0472, P0489, P0493, P0494, P0495, P0496, P0497, P0498, P0499

LAM  
11-30

305210

## Methodology for GC/MS Analysis of Priority Pollutant Compounds

The methods employed in the GC/MS analysis for priority pollutants are established EPA methods. Rigid compliance with the instrument parameters and performance criteria of the published methods was achieved. In some cases, the precise amounts of sample used and the sample handling procedures vary with the complexity of the sample matrix. Qualitative identification of the priority pollutants was performed using the relative retention times, the relative abundance of three characteristic ions and the abundance ratios. The entire mass spectrum was reviewed to confirm each identification. Quantitative analysis of detected compounds was performed by using a response factor generated by a major characteristic ion of the specific compound and an internal standard.

Compounds, in addition to those on the priority pollutant list, were identified through a computer-aided search of the NBS-EPA spectra library. After review the identifications are included in a separate tabulation and labelled "tentatively identified".

### ***Volatile Priority Pollutant Compounds***

For the analysis of Volatile priority pollutants, EPA Method 624 was used. The method can be summarized as follows: Helium is bubbled through a 5 ml water sample contained in a specially designed purging chamber at ambient temperature. The purgeable volatile organic compounds are transferred from the aqueous phase to the vapor phase. The vapor is swept through a sorbent column where the organic components are trapped. After the purge cycle is complete, the sorbent column is heated and backflushed with helium to desorb the organic purgeables onto a gas chromatographic column. The gas chromatograph is temperature programmed to separate the purgeable mixture. The separated purgeable components are then identified and quantitated using a computerized mass spectrometer.

### ***Acid, Base/Neutral and Pesticide Priority Pollutant Compounds***

For the analysis of the Acid, Base/Neutral and Pesticide priority pollutants in an aqueous liquid matrix, EPA Method 625 was used. The method can be summarized as follows: A measured volume of sample, approximately 1 liter, is adjusted to a pH greater than 11 and extracted with methylene chloride. The pH of the sample is adjusted to a value less than 2 and extracted with an aliquot of fresh methylene chloride. A separatory funnel or continuous extractor is used to perform the extractions. The two extracts are dried and concentrated to a 1 ml final volume. Each extract is injected into a GC/MS instrument specifically configured for the correct fraction.

## Methodology for Analysis of Metals

### **AQUEOUS**

The determination of metals in aqueous samples is performed according to the methods published by EPA in "Methods for Chemical Analysis of Water and Wastes," EPA-600/4-79-020, March, 1983, and the Federal Register, October 26, 1983. Arsenic, selenium and thallium are determined by furnace AA; silver, aluminum, barium, beryllium, boron, cadmium, calcium, chromium, copper, cobalt, iron, magnesium, manganese, molybdenum, nickel, lead, sodium, antimony, tin, titanium, vanadium, and zinc are determined by ICP emission spectrometry, except where lower levels of detection are required; in these cases (e.g. lead in groundwater monitoring samples) furnace AA is used. All furnace AA parameters are run by method of standard additions. The determination of mercury is performed by cold vapor AA.

### **EP TOXICITY**

The determination of metals in aqueous EP Toxicity leachates is performed according to the methods published by EPA in "Test Methods for Evaluating Solid Waste" EPA SW-846, revised April, 1984 and the Federal Register, Oct. 26, 1983, 1979. Silver, arsenic, barium, cadmium, chromium, lead and selenium are determined by ICP emission spectrometry. Mercury is determined using cold vapor AA. For leachates that are organic in nature, the analyses are performed according to the methods described under **OIL/SLUDGE** below.

### **SOIL/SEDIMENT**

The determination of silver, beryllium, cadmium, chromium, copper, nickel, antimony, lead, and zinc in sediment samples is performed according to methods published by EPA in "Interim Methods for the Sampling and Analysis of Priority Pollutants in Sediments and Fish Tissue", EPA 600/4-81-055, October 1980. Mercury is determined according to the sediment method published by EPA in "Method for Chemical Analysis of Water and Wastes", EPA 600/4-79-020, March 1983. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion.

### **OIL/SLUDGE**

The determination of silver, aluminum, boron, barium, beryllium, calcium, cadmium, copper, chromium, cobalt, iron, magnesium, manganese, molybdenum, sodium, nickel, lead, antimony, tin, titanium, vanadium, and zinc in sludge/petroleum-based samples is performed by ICP emission spectrometry using a magnesium nitrate dry ashing digestion technique. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion. Mercury is determined by cold vapor AA using the same digestion technique.

## Summary of Quality Assurance/Quality Control Procedures (QA/QC)

ETC bases its quality assurance protocols on the following government guidelines:

- "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-330/9/79/001-R, October 1979;
- the recommended guidelines for EPA Methods 624 and 625. (Federal Register, December 3, 1979, updated on October 26, 1984);
- "Manual of Analytical Methods for the Analysis of Pesticides in Humans and Environmental Samples," EPA 600/8-80-038, June 1980;
- "Determination of 2,3,7,8-TCDD in Soil and Sediment" EPA, Region VII, Kansas City, September 1983;
- Organic Analysis: Multi-media, Multi Concentration-IFB WA84-A267; and
- Dioxin Analysis: Soil/Sediment Matrix; Multi-Concentration; Selected Ion Monitoring with Jar Extraction Procedure-IFB WA84-A002

However, we have modified our protocols to provide a higher level of QA/QC than the guidelines require. For example, we analyze a higher than required number of quality control samples and we pay especially careful attention to the certification of the "reference standard" compounds we use in analysis. Below are listed the key QA/QC elements for the methods we used.

### Analysis of Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry

- Each batch of 13 samples consists of 9 customer samples (at a maximum), one blank sample, one spiked blank, one spiked sample and one replicate sample. This amounts to a 30% quality control factor.
- Three surrogate compounds are added to each sample in the batch of 13.
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and returned, if necessary, at the beginning of each day to ensure that its performance on bromofluorobenzene (BFB) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of Volatile Organic Priority Pollutant "standards" at a minimum of 3 different concentrations and using a mixture of 3 internal standards at a constant concentration.
- The calibration curve is verified with a mixture of priority pollutant standards every day. If the response factors vary greater than 25%, the instrument must be recalibrated.

### Analysis of Organic Compounds Extracted in Acid or Base/Neutral Solutions by Gas Chromatography/Mass Spectrometry

- Each batch of 20 samples consists of 16 customer samples (at a maximum), one blank sample, one spiked blank (for water matrices), one sample spiked with the priority pollutant standard mixture and a duplicate customer sample. This amounts to a 20% quality control factor.

- Three surrogate compounds are added to each sample in the batch for Base/Neutral analysis
- Three surrogate compounds are added to each sample in the batch for Acid analysis.
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and retuned, if necessary, at the beginning of each day to ensure that its performance on decafluorotriphenylphosphine (DFTPP) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of standards composed of either the Organic Acid or Base/Neutral Extractable Compounds at a minimum of 3 concentrations and using five internal standards for quantitation.

### Analysis of Metals

#### All Samples

- New standards are prepared for each batch of samples
- Normal calibration is performed using a blank sample and four standards that have been through the sample preparation procedure. A regression analysis is used to construct the calibration curve.
- All EP Toxicity samples and all samples determined by furnace atomic absorption are calculated by the "method of additions".
- For each sample analysis that requires the use of the "method of additions" technique, a three point calibration is performed using U.S. EPA "Methods for Chemical Analysis of Water and Wastes, 1979". Results are obtained using linear regression analysis. Any regression with a coefficient of correlation below 0.990 is considered suspect, necessitating review of calibration data or sample re-analysis.
- In constructing the normal calibration curves the lowest concentration levels we use are values greater than or equal to 5 times the Instrumental Detection Limit (IDL).
- All calibration standards are analyzed in duplicate, at a minimum.
- Independent reference standards are used to check the accuracy of calibration standards.
- A check standard is analyzed every ten samples to validate the normal calibration curve.
- One customer sample out of every ten is analyzed in triplicate.

#### Homogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are homogeneous, the QC program is a minimum of 25% and consists of analyzing:

- 3 sets of triplicate analyses;
- 2 Replicate spikes.
- 1 independent reference standard;
- 4 Calibration standards (processed using the sample preparation method);
- 4 Calibration standards (without sample preparation); and
- 1 Reagent Blank.



Heterogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are heterogeneous, the QC program is a minimum of 35% and consists of analyzing:

- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 1 Replicate independent reference standards;
- 4 Calibration standards (processed using the sample preparation method);
- 1 Procedural Blank;
- 4 Calibration standards (without sample preparation); and
- 1 Reagent Blank.

Analysis of Mercury

To analyze samples for mercury we group them by matrix in batches of 30 or less. Our QC program is a minimum of 30% and consists of analyzing:

- each of the 30 customer samples in duplicate;
- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 2 Replicate independent reference standards;
- 10 Calibration standards (processed using the sample preparation method); and
- 2 Procedural Blanks.

Analysis of Pesticides, Herbicides and PCB's by Gas Chromatography

Pesticide, herbicide and PCB samples are grouped in batches of 16 customer samples or less according to the type of analysis to be performed. The QC program for each of these three types of analysis is a minimum of 20% and consists of analyzing:

- 1 procedural blank sample (a reagent blank is analyzed in the case of non-water matrices);
- 1 spiked blank sample (the spiked blank is eliminated in the case of non-water matrices);
- 1 replicate sample;
- 1 replicate spiked sample; and
- 1 known reference QC sample for at least each 100 samples analyzed.

The instrument is calibrated each run with three standards, and checked every 10 samples.

Analysis of Cyanides, Phenols, Fluoride, Chloride, Nitrate and Nitrite

- All parameters are analyzed using a Technicon Autoanalyzer II GT.
- 3 calibration standards are analyzed at the beginning and end of each batch.

- Each batch (up to 80 samples) consists of analyzing one blank, one spiked blank, one duplicate and spiked sample every 20 samples, and an EPA known reference sample.

#### Analysis of Total Organic Carbon (TOC)

TOC samples are analyzed on a daily basis with the number of samples analyzed per day dependent on the request for duplicate or quadruplicate analyses. The quality control program is designed to maintain the appropriate amount of QC and consists of the following elements:

- Daily instrument calibration
- One blank
- Standard recalibration every 10 samples
- Spiked samples at a low and high level
- Every sample is run in duplicate at a minimum

#### Analysis of Total Organic Halide (TOX)

- Blank reagent water for absolute carbon background must contain less than 5 ug/l of halide (as chloride).
- Using a trichlorophenol standard, the mean adsorption efficiency must be within +/- 15% of the standard value.
- Calibration standards are run every 10 samples.
- Every sample is run in duplicate at a minimum.

#### Analysis of 2,3,7,8-TCDD (Dioxin) by GC/MS (SIM)

- Each sample is dosed with a known quantity of  $^{13}\text{C}_{12}$ -2,3,7,8-TCDD as internal standard and  $^{37}\text{Cl}_4$ -TCDD as surrogate standard. The action limits for surrogate standard results is +/- 40% of the true value. Samples showing surrogate standard results outside of these limits are reextracted and reanalyzed.
- Two laboratory "method blanks" are run along with each set of 24 or fewer samples. The method blank is also dosed with the internal standard and surrogate standard.
- At least one per set of 24 samples is run in duplicate to determine intralaboratory precision.
- Qualitative Requirements. The following are met in order to confirm the presence of native 2,3,7,8-TCDD:
  - a. Isomer specificity must be demonstrated initially and verified once per 8-hour work shift. The verification consists of injecting a mixture containing TCDD isomers which elute close to 2,3,7,8-TCDD. The 2,3,7,8-TCDD must be separated from interfering isomers, with no more than 25% valley relative to the 2,3,7,8-TCDD peak.
  - b. The 320/322 ratio is within the range of 0.67 to 0.87.
  - c. Ions 320, 322, and 257 are all present and maximize together the signal to mean noise ratio must be 2.5 to 1 or better for all 3 ions.
  - d. The retention time is equal (within 3 seconds) the retention time for the isotopically labeled 2,3,7,8-TCDD.
  - e. At least one of the positives can be confirmed by obtaining partial scan spectra from mass 150 to mass 350. The partial scan guidelines are as follows:

- . the 320/324 ratio should be 1.58 +/- 0.16
- . the 257/259 ratio should be 1.03 +/- 0.10
- . the 194/196 ratio should be 1.54 +/- 0.15
- One sample is spiked with native 2,3,7,8-TCDD at a level of 1.0 PPB (for soil) for each set of 24 or fewer samples.
- In cases where no native 2,3,7,8-TCDD is detected, the actual detection limit is estimated and reported based on a signal to noise ratio of 2.5 to 1 at ions 320 and 322.
- For each sample, the internal standard is present with at least a 10 to 1 signal to noise ratio for both mass 332 and mass 334. Also, the internal standard 332/334 ratio must be within the range of 0.67 to 0.87.

#### Subcontractor QA/QC

Each subcontractor is required to maintain an appropriate level of quality control. To insure this, each subcontractor is required to submit to ETC the quality control data for all analyses it performs. This data is kept on file at ETC. In general, the amount of quality control required is one duplicate sample with one spiked sample for every ten analyses.

#### Chain-of-Custody

The chain-of-custody procedure is part of our quality assurance protocol. We believe our chain-of-custody record fully complies with the legal requirements of federal, state and local government agencies and of the courts of law. The record covers:

- labeling of sample bottles, packing the Sample Shuttle and transferring the Shuttle under seal to the custody of a shipper;
- outgoing shipping manifests;
- the chain-of-custody form completed by the person(s) breaking the Shuttle seal, taking the sample, resealing the Shuttle and transferring custody to a shipper;
- incoming shipping manifests;
- breaking the Shuttle's resealed seal;
- storing each labeled sample bottle in a secured area;
- disposition of each sample to an analyst or technician; and
- the use of the sample in each bottle in a testing procedure appropriate to the intended purpose of the sample.

The records show for each link in this process:

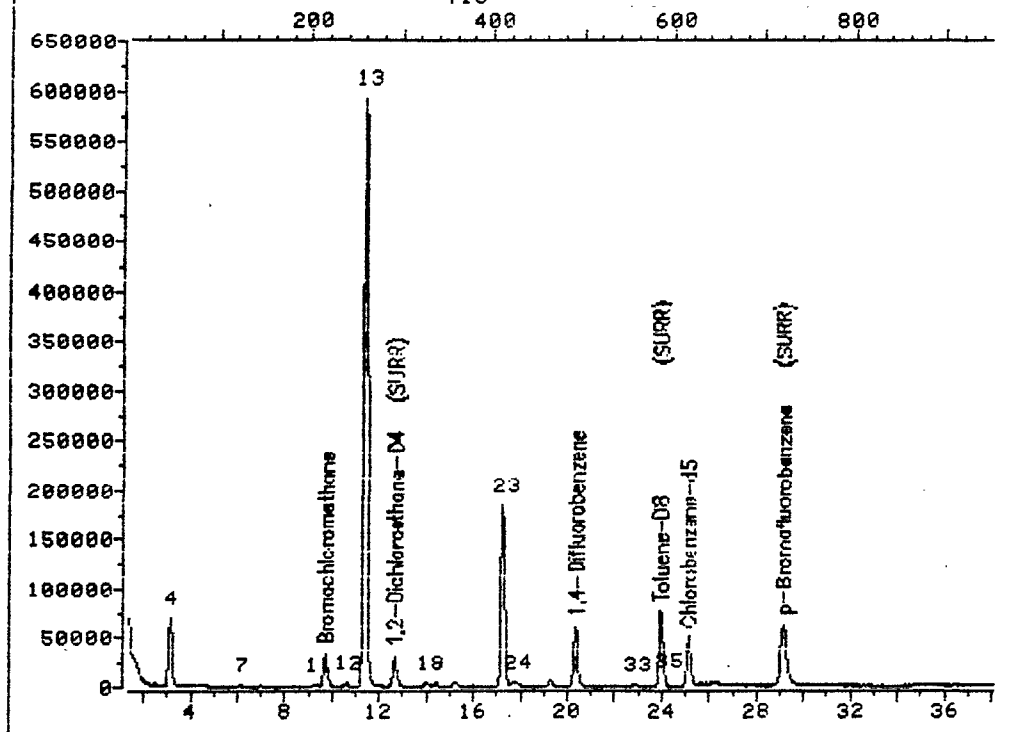
- the person with custody; and
- the time and date each person accepted or relinquished custody.

**Appendix A**  
**Mass Spectral Data**  
**for**  
**Quantitated Compounds**

- 1) A total ion chromatogram for each sample analyzed by a GC/MS instrument.
- 2) A Quant. report used by the analyst to determine qualitative and quantitative results of the compounds present.
- 3) A mass spectrum and a reference spectrum for each priority pollutant compound detected in the sample.

TOTAL ION CHROMATOGRAM

File >M7522 40.0-265.0 amu. PP/VDA 861121 M P4852V, QV5931, L, 5, 5, TIC



Data File: >M7522::U4  
Name: PP/VDA 861121 M  
Misc: P4852V, QV5931, L, 5, 5,

Quant Output File: ^M7522::AQ

Id File: MVOA::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861121 12:01

Operator ID: DG1199  
Quant Time: 861122 05:37  
Injected at: 861122 04:57

QUANT REPORT

Operator ID: DG1199  
 Output File: ^M7522::AQ  
 Data File: >M7522::U4  
 Name: PP/VOA 861121 M  
 Misc: P4852V,QU5931,L,5,5,

Quant Rev: 6      Quant Time: 861122 05:37  
 Injected at: 861122 04:57  
 Dilution Factor: 1.0000U

ID File: MVOA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861121 12:01

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *Bromochloromethane	9.71	217	68147	250.00	NG	98
4) Vinyl chloride	3.13	47	413329	2566.75	NG	97
<del>7) Methylene chloride</del>	<del>6.15</del>	<del>125</del>	<del>2913</del>	<del>19.69</del>	<del>NG</del>	<del>98</del>
11) 1,1-Dichloroethylene	9.28	206	7173	17.19	NG	97
12) 1,1-Dichloroethane	10.60	240	16448	29.38	NG	97
13) 1,2-Trans-dichloroethylene	11.37	260	2251711	5177.55	NG	97
16) *1,4-Difluorobenzene	20.36	492	271475	250.00	NG	82
17) 1,2-Dichloroethane-D4 (SURR)	12.65	293	60126	251.24	NG	92
18) 1,1,1-Trichloroethane	14.01	328	13202	26.64	NG	88
<del>19) Carbon tetrachloride</del>	<del>14.01</del>	<del>328</del>	<del>1814</del>	<del>3.63</del>	<del>NG</del>	<del>98</del>
23) Trichloroethylene	17.26	412	364275	837.31	NG	99
<del>24) Benzene</del>	<del>17.76</del>	<del>425</del>	<del>5477</del>	<del>5.01</del>	<del>NG</del>	<del>97</del>
30) *Chlorobenzene-d5	25.12	615	190573	250.00	NG	97
<del>33) Tetrachloroethylene</del>	<del>22.88</del>	<del>557</del>	<del>2952</del>	<del>5.69</del>	<del>NG</del>	<del>89</del>
34) Toluene-D8 (SURR)	23.96	585	327007	236.65	NG	91
<del>35) Toluene</del>	<del>24.15</del>	<del>590</del>	<del>3720</del>	<del>2.73</del>	<del>NG</del>	<del>87</del>
38) p-Bromofluorobenzene (SURR)	29.15	719	191071	241.64	NG	99

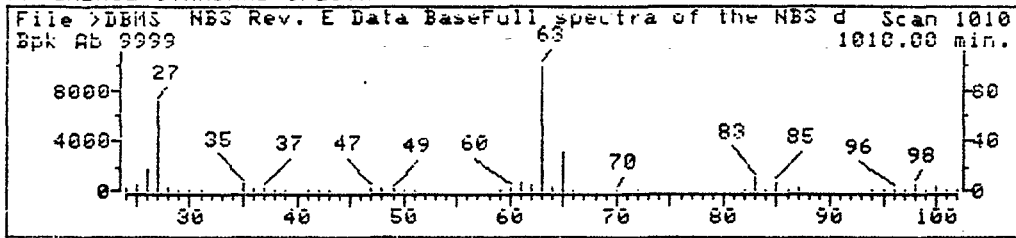
\* Compound is ISTD

SIC 11124186

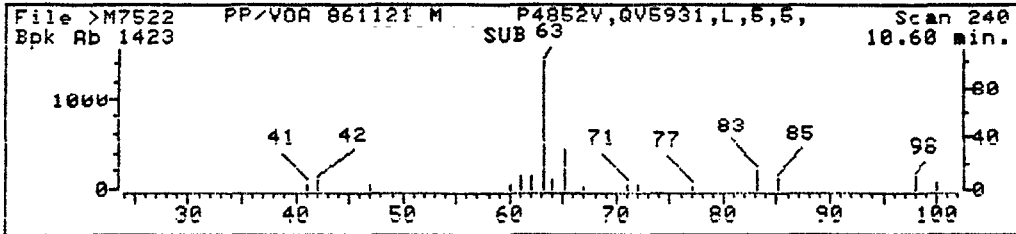
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1986

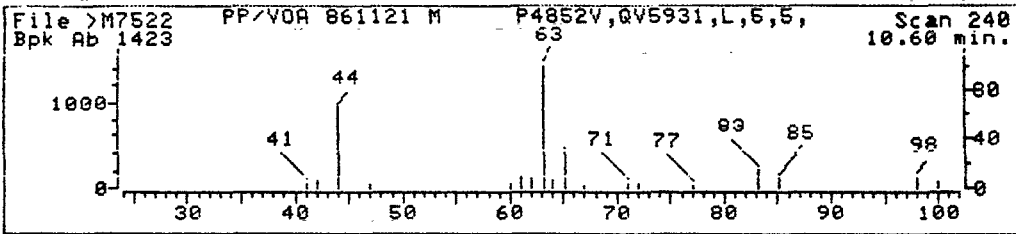
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

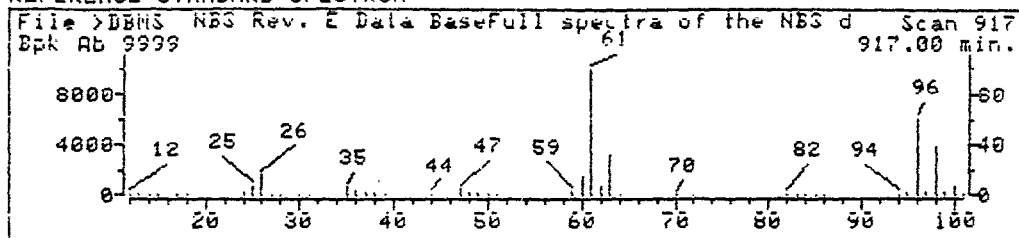


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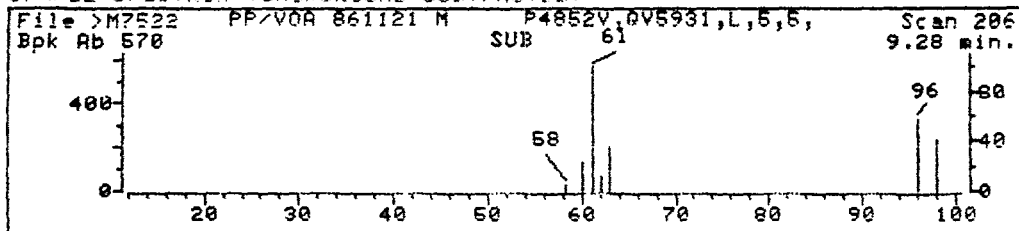
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Quant ID File: MUVA::US  
Last Calibration: 861121 12:01

Compound No: 12  
Compound Name: 1,1-Dichloroethane  
Scan Number: 240  
Retention Time: 10.60 min.  
Quant Ion: 63.0  
Area: 16448  
Concentration: 29.38 NG  
q-value: 97

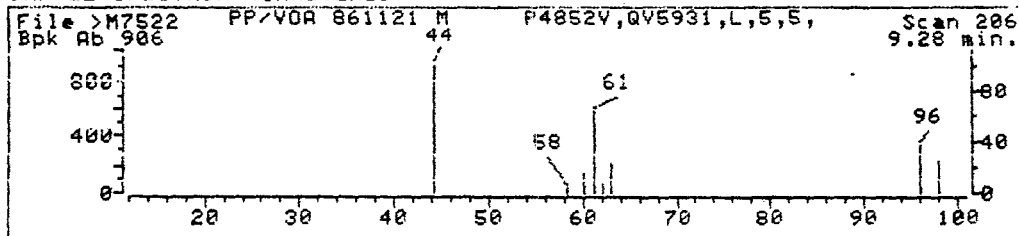
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



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Quant Time: 861122 05:37  
Injected at: 861122 04:57

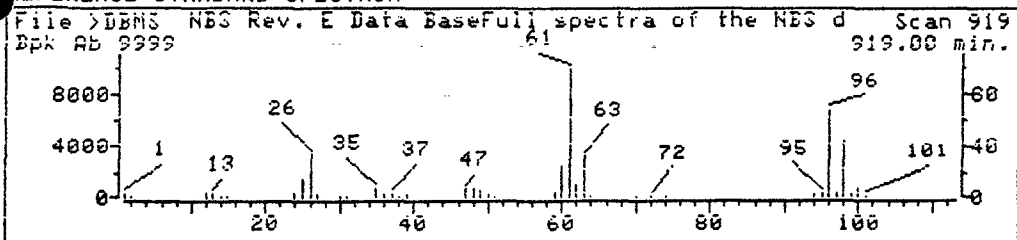
Quant Output File: ^M7522::AQ  
Quant ID File: MVUA::US  
Last Calibration: 861121 12:01

Compound No: 11  
Compound Name: 1,1-Dichloroethylene  
Scan Number: 206  
Retention Time: 9.28 min.  
Quant Ion: 61.0  
Area: 7173  
Concentration: 17.19 NG  
q-value: 97

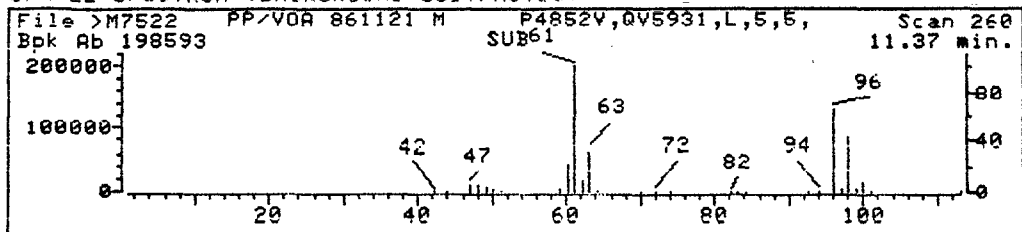
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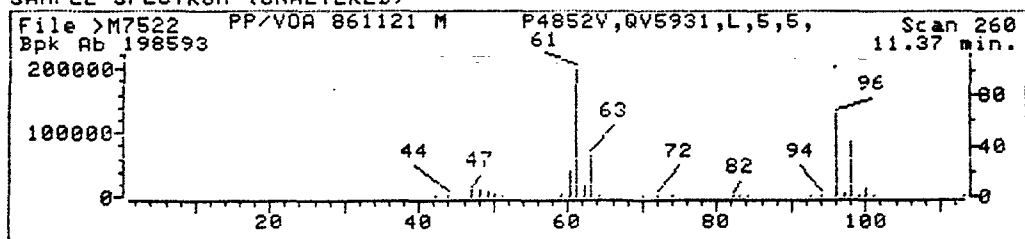
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SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



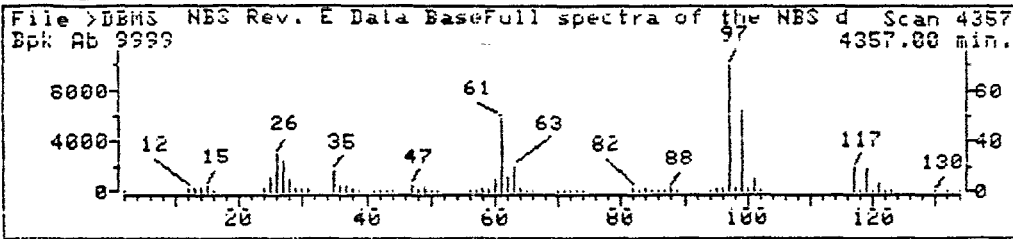
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Quant Time: 861122 05:37  
Injected at: 861122 04:57

Quant Output File: ^M7522::AW

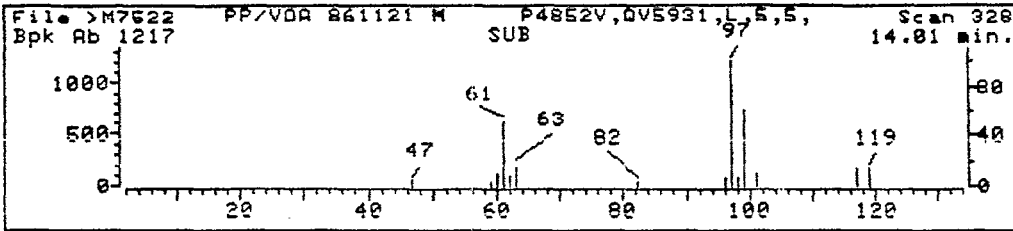
Quant ID File: MVOA::US  
Last Calibration: 861121 12:01

Compound No: 13  
Compound Name: 1,2-Trans-dichloroethylene  
Scan Number: 260  
Retention Time: 11.37 min.  
Quant Ion: 61.0  
Area: 2251711  
Concentration: 5177.55 NG  
q-value: 97

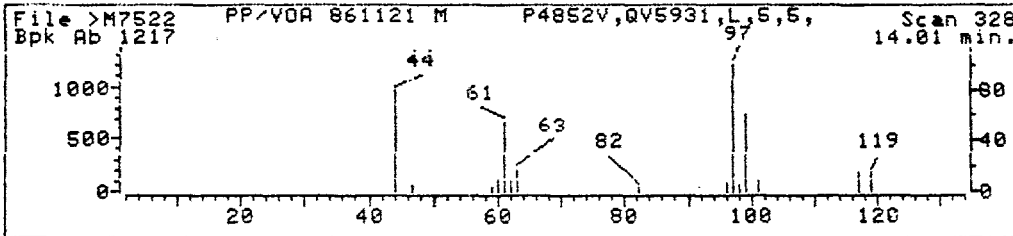
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



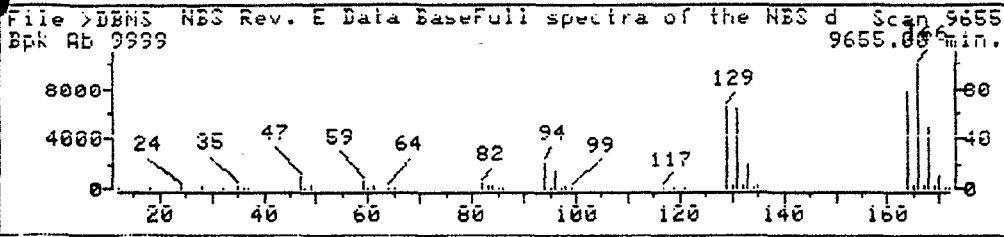
Data File: >M7522::U4  
Name: PP/VDA 861121 M  
Misc: P4852V,QV5931,L,5,5,  
Quant Time: 861122 05:37  
Injected at: 861122 04:57

Quant Output File: ^M7522::AQ  
Quant ID File: MVUA::US  
Last Calibration: 861121 12:01

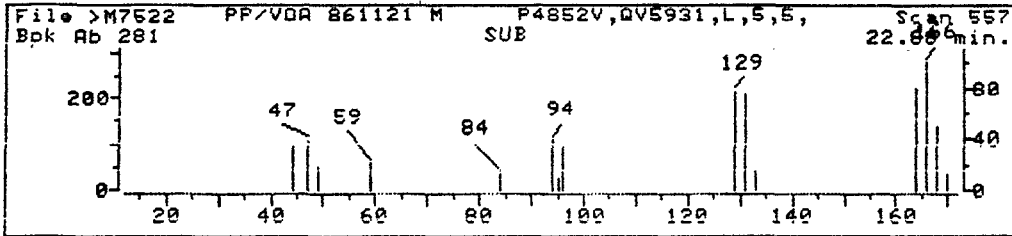
Compound No: 18  
Compound Name: 1,1,1-Trichloroethane  
Scan Number: 328  
Retention Time: 14.01 min.  
Quant Ion: 97.0  
Area: 13202  
Concentration: 26.64 NG  
q-value: 88

305224

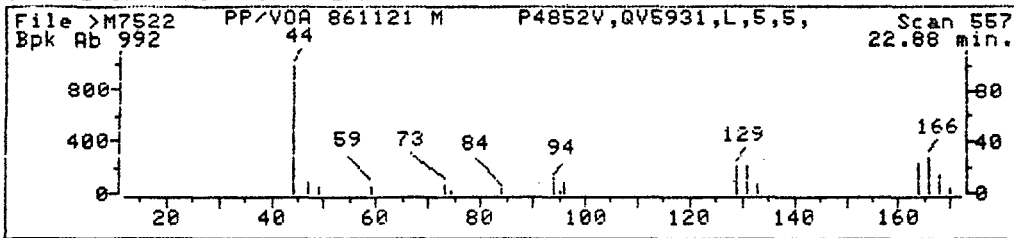
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

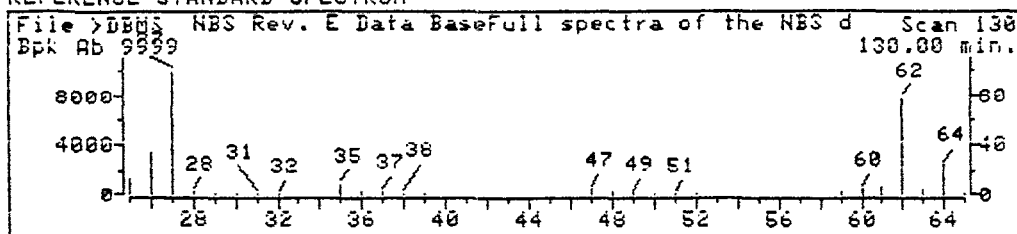


Data File: >M7522::U4  
Name: PP/VDA 861121 M  
Misc: P4852V,QV5931,L,5,5,  
Quant Time: 861122 05:37  
Injected at: 861122 04:57

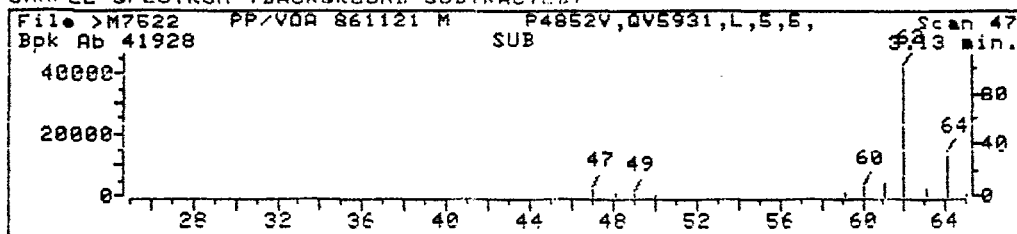
Quant Output File: ^M7522::AQ  
Quant ID File: MVDA::US  
Last Calibration: 861121 12:01

Compound No: 33  
Compound Name: Tetrachloroethylene  
Scan Number: 557  
Retention Time: 22.88 min.  
Quant Ion: 166.0  
Area: 2952  
Concentration: 5.69 NG  
q-value: 89

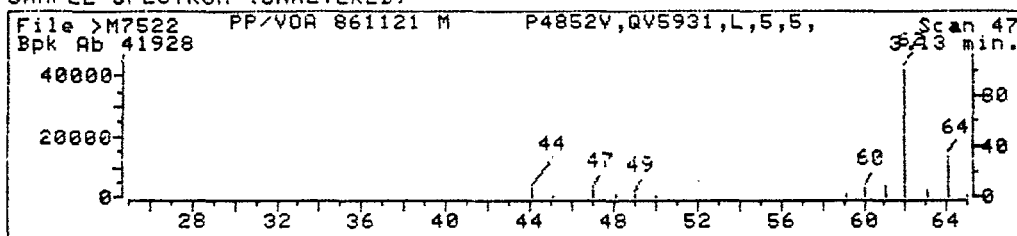
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >M7522::U4  
Name: PP/VOA 861121 M  
Misc: P4852V, QV5931, L, 5, 5,  
Quant Time: 861122 05:37  
Injected at: 861122 04:57

Quant Output File: ^M7522::AQ

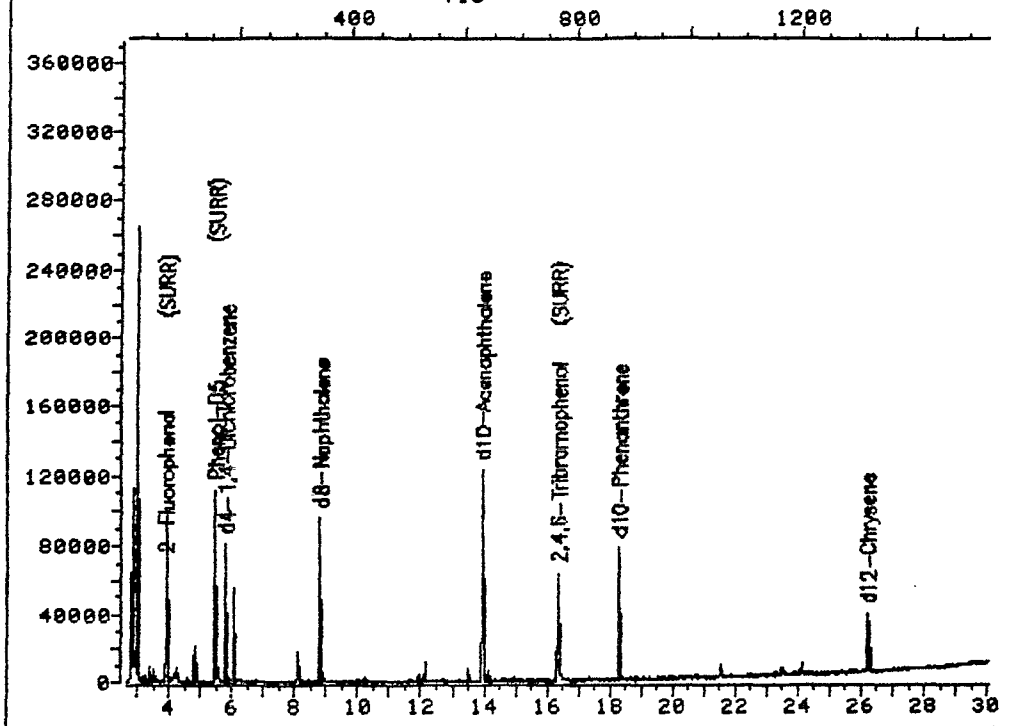
Quant ID File: MVUA::US  
Last Calibration: 861121 12:01

Compound No: 4  
Compound Name: Vinyl chloride  
Scan Number: 47  
Retention Time: 3.13 min.  
Quant Ion: 62.0  
Area: 413329  
Concentration: 2566.75 NG  
q-value: 97

305226

TOTAL ION CHROMATOGRAM

File >F0533 45.0-450.0 amu. PP/ACID,#F P4852A ,QA6013,L, 9  
TIC



Data File: >F0533::U3  
Name: PP/ACID,#F  
Misc: P4852A ,QA6013,L, 930,1

Quant Output File: ^F0533::AQ

BTL# 9

Id File: FACID::US  
Title: PP/ACID IDFILE  
Last Calibration: 861126 11:03

Operator ID: JA4996  
Quant Time: 861126 17:48  
Injected at: 861126 17:15

305227

QUANT REPORT

Operator ID: JA4996  
 Output File: ^F0533::AQ  
 Data File: >F0533::U3  
 Name: PP/ACID,#F  
 Misc: P4852A ,QA6013,L, 930,1

Quant Rev: 6      Quant Time: 861126 17:48  
                   Injected at: 861126 17:15  
                   Dilution Factor: 1.00000

BTL# 9

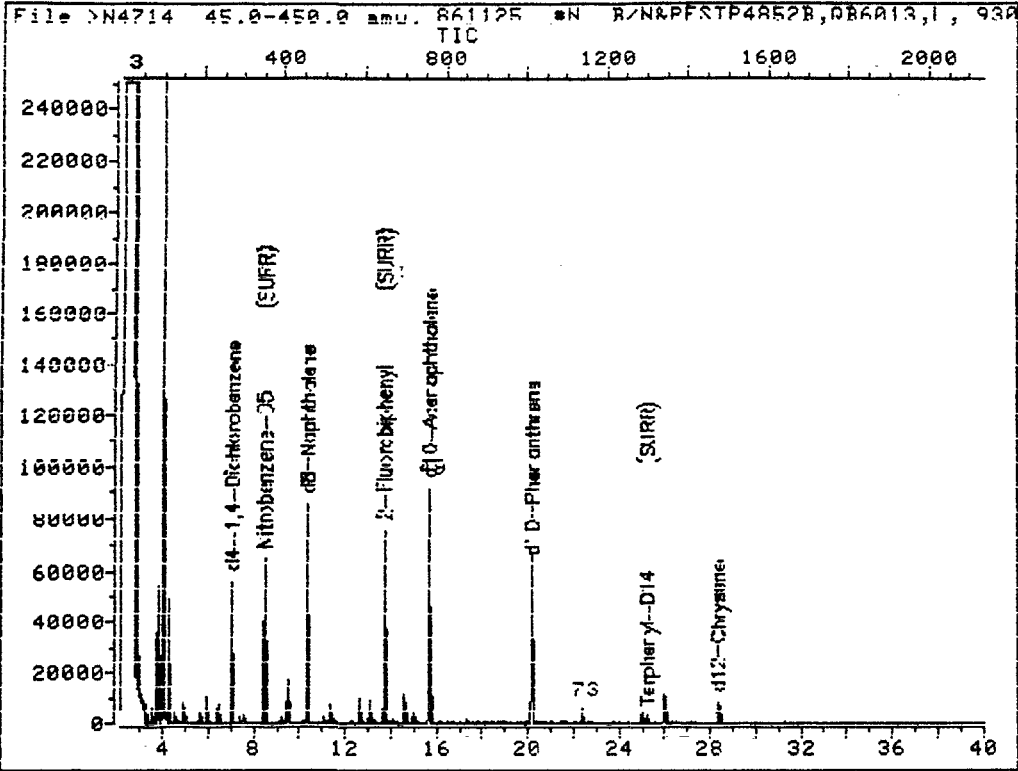
ID File: FACID::US  
 Title: PP/ACID IDFILE  
 Last Calibration: 861126 11:03

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *d4-1,4-Dichlorobenzene	5.80	177	33098	40.00	UG/ML	92
3) 2-Fluorophenol (SURR)	3.89	70	60236	101.66	UG/ML	71
9) Phenol-D5 (SURR)	5.46	158	88911	98.17	UG/ML	96
12) *d8-Naphthalene	8.81	345	94177	40.00	UG/ML	96
18) *d10-Acenaphthalene	13.94	632	57335	40.00	UG/ML	96
25) *d10-Phenanthrene	18.29	875	77353	40.00	UG/ML	99
26) 2,4,6-Tribromophenol (SURR)	16.36	767	22922	78.43	UG/ML	98
29) *d12-Chrysene	26.21	1317	34356	40.00	UG/ML	100

\* Compound is ISTD

305228

TOTAL ION CHROMATOGRAM



Data File: >N4/14::U4  
Name: 861125 #N B/N&PEST  
Misc: P4852B,QB6013,L, 930,1

Quant Output File: ^N4/14::AQ

BIL#41

Id File: NBNP::U5  
Title: B/N+PEST ID FILE  
Last Calibration: 861126 12:44

Operator ID: KUU/86  
Quant Time: 861127 08:16  
Injected at: 861127 03:57

305229

QUANT REPORT

Operator ID: KVV786  
 Output File: ^N4714::AQ  
 Data File: >N4714::U4  
 Name: 861125 #N B/N&PESI  
 Misc: P4852B, Q86013, L, 930, 1

Quant Rev: 6 Quant time: 861127 08:16  
 Injected at: 861127 03:57  
 Dilution Factor: 1.00000

BIL#41

ID File: NBNP::US  
 Title: B/N+PESI ID FILE  
 Last Calibration: 861126 12:44

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *d4-1,4-Dichlorobenzene	7.00	272	17048	40.00	UG/ML	98
3) N-Nitrosodimethylamine	2.60	24	1625	<del>3.67</del>	UG/ML	100
3) N-Nitrosodimethylamine	2.69	29	5717	<del>12.90</del>	UG/ML	100
12) Nitrobenzene-D5 (SURK)	8.40	351	50089	49.62	UG/ML	80
21) *d8-Naphthalene	10.30	458	93843	40.00	UG/ML	99
22) 2-Fluorobiphenyl (SURK)	13.65	647	68194	39.78	UG/ML	96
42) *d10-Acenaphthalene	15.65	760	57536	40.00	UG/ML	98
49) Dimethyl phthalate	15.65	760	15168	<del>8.00</del>	UG/ML	59
66) *d10-Phenanthrene	20.14	1013	79402	40.00	UG/ML	97
73) Di-n-butyl phthalate	22.35	1138	7308	<del>2.84</del>	UG/ML	95
85) *d12-Chrysene	28.33	1475	9719	40.00	UG/ML	100
98) Terphenyl-D14 (SURK)	25.21	1299	3448	11.90	UG/ML	96

\* Compound is ISTD

WJ  
 11-9-86

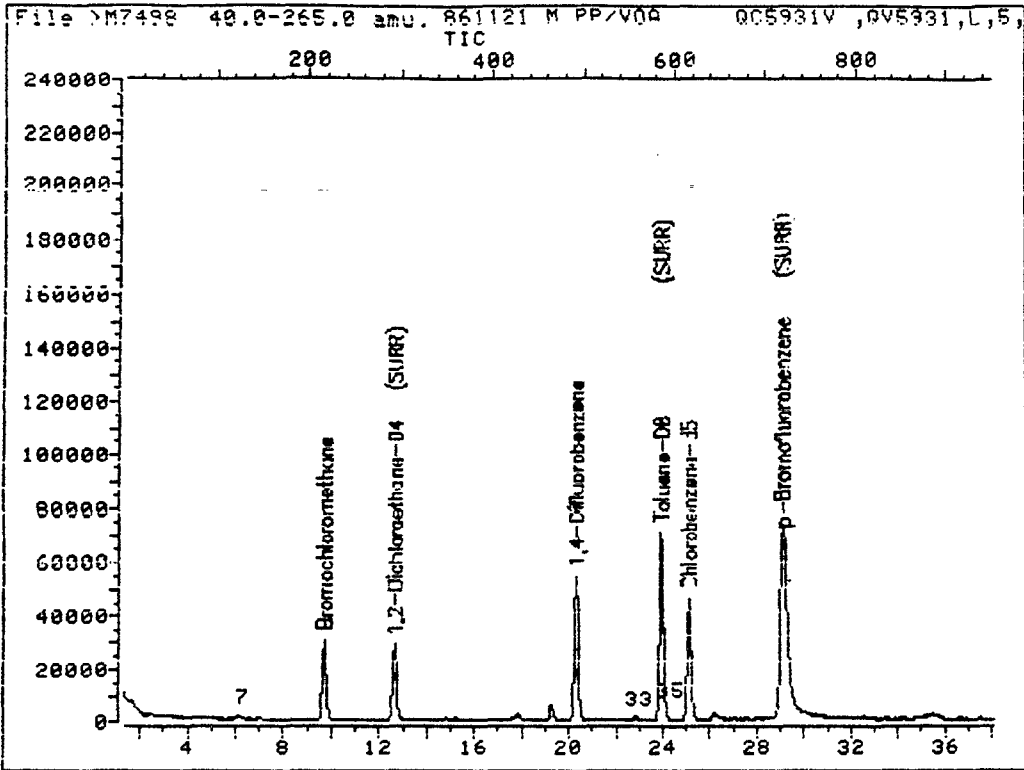
305230



**Appendix C1**  
**GC/MS Subsidiary Data**

305231

TOTAL ION CHROMATOGRAM



Data File: >M7498::U3  
Name: 861121 M PP/V0A  
Misc: QC5931V ,QV5931,L,5,5,

Quant Output File: >M7498::AQ

Id File: MV0A::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861121 12:01

Operator ID: JQ6275  
Quant Time: 861121 12:03  
Injected at: 861121 06:18

305232

QUANT REPORT

Operator ID: JQ6275                      Quant Rev: 6      Quant Time: 861121 12:03  
 Output File: ^M/498::AQ                      Injected at: 861121 06:18  
 Data File: ^M/498::U3                      Dilution Factor: 1.00000  
 Name: 861121 M PP/VOA  
 Misc: QC5931U ,WV5931,L,5,5,

ID File: MVOA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861121 12:01

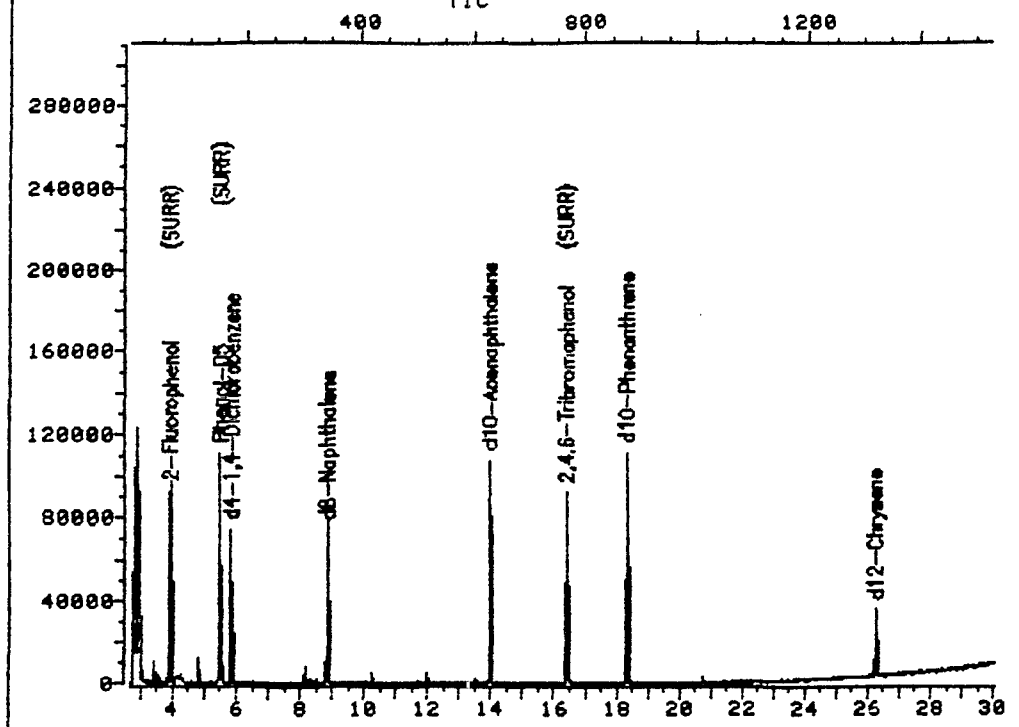
Compound	R.T.	Scan#	Area	Conc	Units	q
1) *Bromochloromethane	9.70	217	66376	250.00	NG	99
7) Methylene chloride	6.18	126	2829	19.63	NG	86
16) *1,4-Difluorobenzene	20.31	491	249275	250.00	NG	82
17) 1,2-Dichloroethane-D4 (SURR)	12.65	293	57076	259.74	NG	94
30) *Chlorobenzene-d5	25.11	615	170859	250.00	NG	99
<del>33) Tetrachloroethylene</del>	<del>22.83</del>	<del>556</del>	<del>2829</del>	<del>6.08</del>	<del>NG</del>	<del>94</del>
34) Toluene-D8 (SURR)	23.95	585	307632	248.31	NG	91
<del>35) Toluene</del>	<del>24.15</del>	<del>598</del>	<del>1931</del>	<del>1.58</del>	<del>NG</del>	<del>92</del>
38) p-Bromofluorobenzene (SURR)	29.10	718	181687	256.29	NG	97

\* Compound is ISTD

ENC 11/24/86

TOTAL ION CHROMATOGRAM

File >F0526 45.0-450.0 amu. PP/ACID,#F QC6013A ,QA6013,L,10  
TIC



Data File: >F0526::U3  
Name: PP/ACID,#F  
Misc: QC6013A ,QA6013,L,1000,1

Quant Output File: ^F0526::AQ

BTL# 2

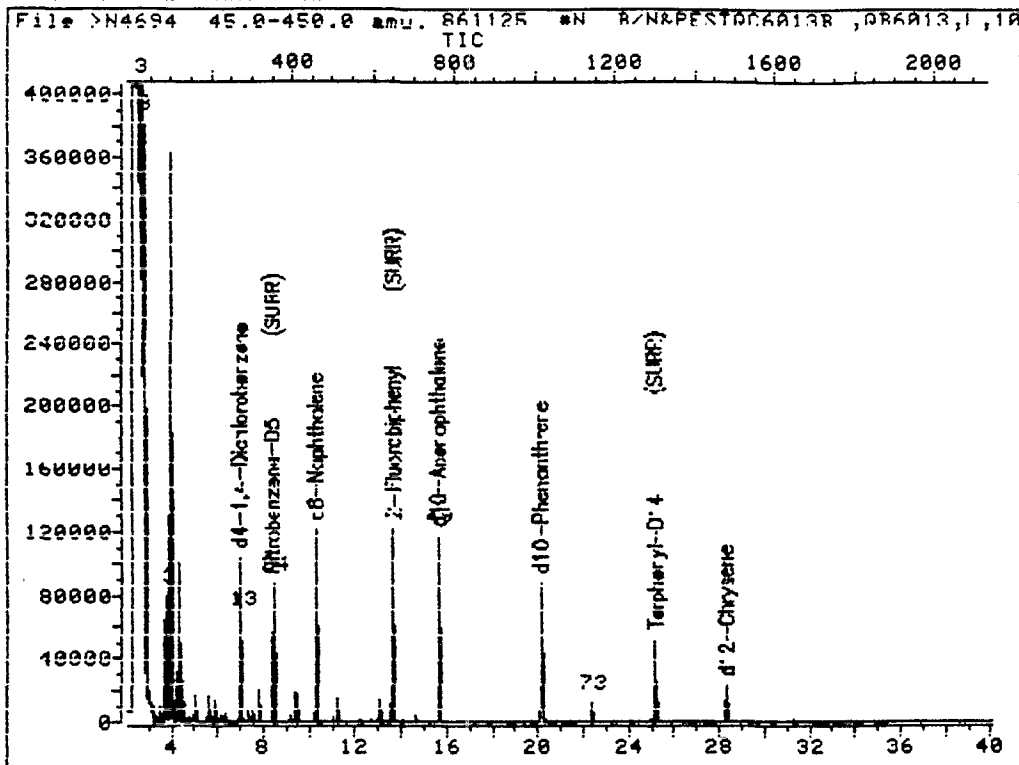
Id File: FACID::US  
Title: PP/ACID IDFILE  
Last Calibration: 861126 11:03

Operator ID: JA4996  
Quant Time: 861126 13:17  
Injected at: 861126 12:45

305234



TOTAL ION CHROMATOGRAM



Data File: >N4694::U4  
Name: 861126 #N B/N&PEST  
Misc: QC6U13B ,QB6U13,L,1000,1

Quant Output File: ^N4694::AQ

BIL#21

Id File: NBNP::US  
Title: B/N+PEST ID FILE  
Last Calibration: 861126 12:44

Operator ID: KUU/86  
Quant time: 861126 16:26  
Injected at: 861126 11:25

305236

QUANT REPORT

Operator ID: KVV786      Quant Rev: 6      Quant time: 861126 16:26  
 Output File: ^N4694:IAQ      Injected at: 861126 11:25  
 Data File: >N4694:U4      Dilution Factor: 1.00000  
 Name: 861125 #N B/N&PEST  
 Misc: QC6U13B ,QB6U13,L,1000,1      BIL#21

ID File: NBNP::US  
 Title: B/N+PEST ID FILE:  
 Last Calibration: 861126 12:44

Compound	R.I.	Scan#	Area	Conc	Units	q
1) *d4-1,4-Dichlorobenzene	6.95	267	34441	40.00	UG/ML	98
3) N-Nitrosodimethylamine	2.66	25	10063	<del>11.24</del>	UG/ML	100
3) N-Nitrosodimethylamine	2.76	31	1340	<del>1.50</del>	UG/ML	100
3) N-Nitrosodimethylamine	3.78	88	2107	<del>2.35</del>	UG/ML	100
12) Nitrobenzene-D5 (SURR)	8.35	346	84430	41.40	UG/ML	83
13) bis(2-Chloroisopropyl)ether	6.97	268	1836	<del>6.80</del>	UG/ML	74
21) *d8-Naphthalene	10.25	453	144769	40.00	UG/ML	98
22) 2-Fluorobiphenyl (SURR)	13.63	643	107010	40.47	UG/ML	97
24) N-Nitrosodi-n-propylamine	8.35	346	12555	<del>18.77</del>	UG/ML	66
42) *d10-Acenaphthalene	15.61	755	73999	40.00	UG/ML	97
49) Dimethyl phthalate	15.61	755	18965	<del>7.78</del>	UG/ML	56
66) *d10-Phenanthrene	20.11	1009	102784	40.00	UG/ML	96
53) Di-n-butyl phthalate	22.31	1133	17098	5.13	UG/ML	97
7) *d12-Chrysene	28.30	1471	25929	40.00	UG/ML	100
8) Terphenyl-D14 (SURR)	25.18	1295	49419	63.93	UG/ML	94

\* Compound is ISTD

WD  
11/29/88

**Appendix D**  
**Subcontractor's Data**

- 1) A copy of the originating subcontractor's report is included for all data not generated within ETC's laboratory.

305238





## Appendix E

### Chain-of Custody Forms

1. A field Chain-of-Custody form (CC1) is included for all samples shipped by ETC shuttle.
2. An in-house sample Chain-of-Custody form is included for all samples not shipped by ETC shuttle.
3. Any additional Chain-of-Custody material provided by a client or by a client's sampling agent is also included.
4. A subcontractor's Chain-of-Custody form is included for any analytical work not performed within ETC's laboratory.
5. Analysis and Extraction Custody forms are included for the period the sample was in ETC's possession.

**ETC ENVIRONMENTAL TESTING and CERTIFICATION**  
**CHAIN OF CUSTODY FORM (CC1)**

Seal No. 67751 ETC Job # P4852  
 Date Sealed 6 NOV. 1986 By: M. I. /

Company: HELLERTOWN MANUFACTURING Attn.: MARK SCHULTZ  
 Facility/Site: C/O GUARD OFFICE-SILVER ROAD Phone: \_\_\_\_\_  
 Address: HELLERTOWN, PA. 18055

**SAMPLE IDENTIFICATION**

Facility: CHAMPION (Facility/Site Code) \_\_\_\_\_ (Optional Sample Point Descriptions)  
 Sample Point: M-CSP-3 \_\_\_\_\_ 411086 1400 \_\_\_\_\_  
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: 86/11/10 MB861113  
 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	Type	BOTTLE		ANALYSIS	SAMPLER		LAB
		Size	Preserv.		Fill. (Y/N)	Observations	Observations
1	M	500 ML	HN03	METALS			/
1	PN	125 ML	H2S04	PHENOLS			/
1	B	125 ML	H2S04	NITRATE			/
1	CN	125 ML	NAOH	CYANIDE			/
1	A	500 ML	BAKED DI	SOLIDS, TD/SULFATE			/
1	D	500 ML	RINSE SOD.	FLUORIDE			/
2	V	40 ML	THIOL	VDA			/
3	E	1 L	BAKED GC/MS	EXTRACTABLES			/
1	TB	40 ML	H2O	TRIP BLANK			/

**CHAIN OF CUSTODY CHRONICLE**

1. Shuttle Opened By: (print) M. SCHULTZ Date: 11-10-86 Time: \_\_\_\_\_  
 Signature: Mark Schultz Seal #: \_\_\_\_\_ Intact: \_\_\_\_\_  
 I have received these materials in good condition from the above person.  
 Name: \_\_\_\_\_ Signature: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_ Remarks: \_\_\_\_\_  
 I have received these materials in good condition from the above person.  
 Name: \_\_\_\_\_ Signature: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_ Remarks: \_\_\_\_\_  
 Shuttle Sealed By: (print) M. SCHULTZ Date: 11-10-86 Time: 16:30  
 Signature: Mark Schultz Seal #: 0067752 Intact: \_\_\_\_\_

LAB USE ONLY Opened By: Mark Schultz Date: 11/11/86 Time: 5:20 PM  
 SHUTTLE # 1165 TEMP. °C 12 SEAL # 67752 COND. T/INTACT







Sample Number	Log Link	Sample Vol. (ml)	Extract Vol. (ml)		Comments
			BN	ACID	
P0447	17862	1000	1.0	1.0	pp/acid/ps/peis
P0448		1000	1.0	1.0	
P0449		1000	1.0	1.0	
P0450		1000	1.0	1.0	
P0457		1000	1.0	1.0	
P0458		1000	1.0	1.0	
P0467		1000	1.0	1.0	
P0468		1000	1.0	1.0	
P0471		1000	1.0	1.0	
<del>P0472</del>		<del>995</del>	<del>1.0</del>	<del>1.0</del>	
P0489	17863	1000	1.0	1.0	pp/T
P0493		980	1.0	1.0	
P0494		1000	1.0	1.0	
P0495		1000	1.0	1.0	
P4852	17866	930	1.0	1.0	pp/acid/ps/peis
P4853		930	1.0	1.0	
P0472	17862	1000	1.0	1.0	
QC 6013		1000	1.0	1.0	
QC 6013	S	1000	1.0	1.0	
P0447	S	980	1.0	1.0	
P0448	R	995	1.0	1.0	

QC Batch # 6013

Analysis \_\_\_\_\_  
SEE COMMENTS

Matrix H<sub>2</sub>O

Turnaround NORM

Date 11/17/86

**Extraction Method:**

Sep. Funnel L.B. 11-17-86  
BN ON: 11:30A 11-17-86  
 Continuous OFF BN 11:30A 11-18-86  
ACID ON: 1:00PM 11/18/86  
 Soxhlet OFF: 1:00PM 11/19/86

Other \_\_\_\_\_

**COMMENTS**

⊕ P0472 - Acid fraction lost during conc. repeated using a separate aliquot

FRACTION	SPIKE		
	Amt (ml)	(ug/ml) Conc.	Lot #
PP/BN	1.0	100	13,929
PP/ACID	1.0	100	13,829
Dioxin 1260	1.0	100	14,404
Heptachlor	1.0	100/200	14,631
SCMI-VDA			

SURROGATE		
Amt. (ml)	Conc. (ug/ml)	Lot #
1.0	BN's - 50	14,597
	Acid's - 100	

Set-up: L. Boyer 11-17-86 Aug Nazarek 11-17-86  
 Conc.: L. Boyer 11-24-86

UPD/Supervisor: Raymond [Signature] 11-25-86  
 Spike/Surr. Verified: Kenneth Boyer 11-17-86  
305745

DATE 11-20-82 SHIFT \_\_\_\_\_  
 FRACTION PA/ACID  
 INSTRUMENT F  
 TUNE FILE MTF001  
 SEQUENCE FILE SAFI/JAF  
 METHOD FILE AUDF  
 ID FILE TALID  
 ANALYST(S) J. AO  
 SUPERVISOR Stapenrich  
 BATCH #'s QA 6013  
QA 6056  
 (PLEASE INITIAL)

STANDARD	CONC PPM	LOT NO.	LOT VOL.
PP/ACID STD	300	14429	
↓	300	14428	
↓	60	14427	
Int. STD Test MIX	400	14585	
		14616	

*page 193*

CURRENT CS05 STATUS		STANDARDS UPDATED	
ACQ		DATE	
WIP		BY	

NAME	DATA FILE	UL INJ	ALS #	DIL	TAPE #	SPECIALS (WRITE A-TYPE)	PLUS Y/N
Column Test mix	> F0520					OK	
300 ppm PP/ACID STD	> F0521					← CRUPP	
100 ppm PP/ACID STD	> 0522					← CRCAE	
60 ppm PP/ACID STD	> 0523					←	
Column Test MIX	> 0524				11:04	OK	
QC 6013 AS	> 0525		1				
QC 6013 A	> 0526		2				
P0447 AS	> 0527		3				
P0447 A	> 0528		4				
P0448 A	> 0529		5				
P0448 AR	> F0530		6				
P0449 A	> 0531		7				
P0450 A	> 0532		8				
P0452 A	> 0533		9				
P0453 A	> 0534		10				
P0457 A	> 0535		11				
P0458 A	> 0536		12				
P0467 A	> 0537		13				
P0468 A	> 0538		14				
P0471 A	> 0539		15				
P0472 A	> F0540		16				
Column Test MIX	> 0541		17		22:39		
100 ppm PP/ACID STD	> 0542		18			OK	
P0489 A	> 0543		19				
P0493 A	> F0544		20				

3052





11-30

DATE Nov. 26th 86 SHIFT \_\_\_\_\_  
 FRACTION B/N  
 INSTRUMENT "N"  
 TUNE FILE \_\_\_\_\_  
 SEQUENCE FILE XV25  
 METHOD FILE BNPN; ACIDN  
 ID FILE N3UP  
 ANALYST(S) K. Valach  
 SUPERVISOR (Signature)  
 BATCH #'s Q 6057; Q 6013

STANDARD	CONC PPM	LOT NO.	LOT VOL
Pest. I	150	14433	
B/N I	200	14432	
B/N II	100	14713	
R/N I	60	14430	
Aniline	300	13468	
Aniline	100	13469	
Aniline	60	12470	
2-Picoline	300	14411	
2-Picoline	100	14412	
2-Picoline	60	14413	
Methoxy chlo	300	14352	
Methoxy chlo	100	14353	
Methoxy chlo	60	14354	
1st. std.	400	14585	
Col. Test Mix	25	14616	

(PLEASE INITIAL)

CURRENT CS05 STATUS	STANDARDS UPDATED
ACQ	DATE <u>Nov. 26th 1986</u>
WIP	BY <u>KV</u>

NAME	DATA FILE	uL INJ	ALS #	DIL	TAPE #	SPECIALS (WRITE A-TYPE)	PLUS YIN
Col. Test-Mix	>N4673					OK 14:37	
Pest. col. std. IV; 150ppm	>N4674		1				
B/N col. std. II; 200ppm	>N4675		2				
B/N col. std. II; 100ppm	>N4676		3				
B/N col. std. I; 60ppm	>N4677		4				
Aniline std.; 300ppm	>N4678		5				
Aniline std.; 100ppm	>N4679		6				
Aniline std.; 60ppm	>N4680		7				
2-Picoline Methoxy chlo std.; 300ppm	>N4681		8				
2-Picoline Methoxy chlo std.; 100ppm	>N4682		9				
2-Picoline Methoxy chlo std.; 60ppm	>N4683		10			OK	
Col. Test-Mix	>N4684		11			OK 03:17	
QC 6057 BS	>N4685		12			AN1260	
P 2712 BS	>N4686		13			AN1260	
QC 6057 B	>N4687		14			PCBSCN	S Y
P 2711 B	>N4688		15				S Y
P 2711 BR	>N4689		16				
P 2712 B	>N4690		17				S Y
N 6321 B	>N4691		18			V	S Y
QC 6013 BS	>N4692		19			AN1260	
P 0447 BS	>N4693		20			AN1260	
QC 6013 B	>N4694		21			PCBSCN	
Col. Test-Mix	>N4695		22			OK 12:15	
B/N col. std. II; 100ppm	>N4696		23			AN1260; OK	
P 2712 B	>N4697		24			PCBSCN	

305218





TEST: NITRITE, UNITS: mg/L  
 METH. REF. # 353.2, MDL: 0.10  
 INSTRUMENT # #1 (Channel 2)

BY: S.T.  
 BOOK # CV-12 P. 100  
 FROM PAGE # 39

Log Link	JOB #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	JOB #	ANALYZED CONC.	Dilu	REPORTED CONC.
17524	N6110	0		BMDL	17866	P4850	.00157		BMDL
17529	N6120	.0028		BMDL	17866	P4851	.0166		BMDL
17529	N6126	.0040		BMDL	17866	P4852	.0490		BMDL
17531	P2710	.0024		BMDL	17866	P4853	.0446		BMDL
17531	P2725	0		BMDL	17878	P2774	.00181		BMDL
17539	P2117	.0043		BMDL	17866	P3120	.00114		BMDL
17539	P2116	.0954		BMDL	17866	P3122	0		BMDL
17539	P2119	.0571		BMDL	17866	P3124	.0167		BMDL
17524	P1711	.0043		BMDL	17866	P3126	0		BMDL
17524	P1713	.0040		BMDL	17882	P2543	.00471		BMDL
17524	P1714	0		BMDL	17882	P2544	.00213		BMDL
17524	P1717	.00119		BMDL	17882	P2545	.00473		BMDL
17562	N6128	0		BMDL	17882	P2547	.00784		BMDL
17562	N6148	.00119		BMDL	17882	P2549	.00119		BMDL
17573	P2717	0		BMDL	17886	P3117	.0066		BMDL
17579	P2724	.114		.12					
17584	P3068	.100		.10					
17503	P2683	.121		.12					
17562	N6122	.00357		BMDL					
17524	P1715	.00213		BMDL					
17864	P2531	.00119		BMDL					
17864	P2541	.0611		BMDL					
17864	P2546	.0691		BMDL					
17864	P2557	.00282		BMDL					

STANDARDS		
PREPARED ON _____		
BY _____		
FROM _____ mg/L		
STOCK # _____		
STANDARD mg/L	VALUE OBTAINED	% of Threshold
1.00	.99	99%
0.50	.51	102%
0.20	.197	98.5%
0.05	.046	92%
SLOPE	Intercept	Corr. Coef.
.0120	-.0005	.9998

**ADDITIONAL COMMENTS**  
 ALL CALCULATIONS  
 BY  $X = MY + B$

Analyses QA/QC Data Report      QC Batch # (103004 NOL)      Verified: 11/24/86 *[Signature]*

PRECISION	
METHOD BLANK	.00387
Spiked Blank (Known)	.20
Spiked Blank Result	.198
% Recovery	99
EXTERNAL REF. QC	N/A
Known Value	.5
X Observed Result	-.532
X % Recovery	

ACCURACY	
Job # Dup (ETC)	N6110 P1714
Original Result	0 0
Duplicate Result	0 .0015
RANGE	0.0 .0015
Job # Spiked (ETC)	P1714
Original Result	0
Amount Spiked	.0282
Spiked S' Result	
% Recovery	1.31

TEST: NITRATE, UNITS: mg/L  
 METH. REF. # 353.2, MDL: 0.10  
 INSTRUMENT # #1 (Channel 2)

BY: CH  
 BOOK # CV-12 P. 98  
 FROM PAGE #     

Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	DILN	REPORTED CONC.	STANDARDS	
7798	P2469	over		*	17886	P3117	1.15		1.15	PREP'D ON _____ BY _____ FROM _____ mg/L STOCK# _____	
7798	P2471	over		*	17866	P3120	.0207		BMDL		
7798	P2524	over		*	17866	P3122	.0511		BMDL		
7798	P2538	over		*	17866	P3124	.0173		BMDL		
7798	P2539	over		*	17866	P3126	.0127		BMDL		
7798	P2540	over		*	17882	P2543	over		*	STANDARD mg/L   VALUE OBTAINED   % of theoretical	
7814	P3535	.487		.49	17882	P2544	.0250		BMDL		2.00   1.98
7822	P3107	over		*	17882	P2545	over		*		1.00   1.04
17822	P3112	.007		BMDL	17882	P2547	over		*		.50   .497
17828	P2473	over		*	17882	P2549	over		*		.20   .0934
7828	P2472	over		*	17882	P2550	over		*		
17828	P2522	over		*	17882	P2551	over		*		SLOPE   Intercept   Corr. Coef.
17828	P2536	over		*	17882	P2552	over		*		
17828	P2548	over		*							
17864	P2481	.006		BMDL							
17864	P2531	.0245		BMDL							
17864	P2541	1.73		1.73							
17864	P2546	over		*							
17864	P2554	2.06		2.06							
17866	P4850	over		*							
17866	P4851	over		*							
7866	P4852	.398		.40							
7866	P4853	over		*							
17878	P2778	.0462		BMDL							

ADDITIONAL COMMENTS  
 \* RPT on Channel 1  
 complete Page 97

**Analyses QA/QC Data Report**

QC Batch # NCO03

Verified: 11/24/86 [Signature]

METHOD BLANK	.003	.0177
Spiked Blank (Known)	.5	.5
Spiked Blank Result	.501	.532
% Recovery	160%	1.06

**PRECISION**

Job # Dup (ETC)	P2473	P4850	P3126
Original Result	over	over	.0217
Duplicate Result			.0194
R Range	↓	↓	

Job # Spiked (ETC)	P2473	P4850	P3126
Original Result	over	over	.021
Amount Spiked			.5
Spiked S' Result	↓	↓	.506
% Recovery			97%

205256

TEST: CYANIDE, UNITS: mg/L  
 METH. REF. # 335.3, MDL: 0.025  
 INSTRUMENT # #1 (citanna 1)

BY: C.H.  
 BOOK # CY-12 P. 119  
 FROM PAGE #     

Log Link	JOB #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	JOB #	ANALYZED CONC.	Dilu	REPORTED CONC.
17837	N4709	0		BMDL	17697	P4748	.005		BMDL
17837	N4714	0		BMDL	17697	P4748	.004		BMDL
17862	P0489	.007		BMDL	17697	P4748	.004		BMDL
17863	P0493	.009		BMDL	17698	P1710	.003		BMDL
17863	P0494	.005		BMDL	17701	N4710	.004		BMDL
17863	P0495	.002		BMDL	17701	N4715	.004		BMDL
17866	P4850	.539		.539	17866	P4851	over		Redo
17868	P4510	*		Redo	17866	P4852	over		Redo
17878	P2776	.007		BMDL	17866	P4853	X		Redo
17881	P1865	.002		BMDL	17868	P4510	*		Redo
17883	N4708	.002		BMDL	17868	P4511	K		Redo
17883	N4711	.003		BMDL	17868	P4512	K		Redo
17833	P4241	.0197		BMDL	17868	P4513	K		Redo
17833	P4242	.0169		BMDL	17883	N4712	X		Redo
17833	P4243	.0211		BMDL	17883	N4713	K		Redo
17833	P4244	.0186		BMDL	17889	P2754	X		Redo
17834	P4248	.0185		BMDL	17889	P2761	*		Redo
17834	P4249	.0189		BMDL	17889	P2774	X		Redo
17834	P4250	.0231		BMDL	17889	P2775	K		Redo
17834	P4251	.0189		BMDL	17833	P4235	X		Redo
<del>17834</del>	<del>P2724</del>	<del>.026</del>		<del>BMDL</del>	17833	P4236	N		Redo
17693	P2916	.0185		BMDL	17833	P4237	L		Redo
17693	P2917	.0312		BMDL	17833	P4238	L		Redo
17697	P4748	.006		BMDL	17833	P4239	F		Redo

STANDARDS

PREPARED ON 12-3  
 BY D.B.  
 FROM      mg/L  
 STOCK #     

STANDARD mg/L	VALUE OBTAINED	% of theoretical
.500	.493	98.6%
.400	.405	101.0%
.200	.205	102.5%
.100	.102	102.0%

SLOPE	Intercept	Corr. Coef.

ADDITIONAL COMMENTS

\* RPT DUE TO INSTRUMENTATION Problem  
 P4851 10x  
 P4852 20x

Analyses QA/QC Data Report

QC Batch # CY003

Verified: 12/3/86

METHOD BLANK	Result
METHOD BLANK	.008
Spiked Blank (Known)	.100
Spiked Blank Result	.103
% Recovery	103%

INTERNAL REF. QC	Result
INTERNAL REF. QC	CYA
Known Value	.224
Observed Result	.228
% Recovery	101.7%

PRECISION			
Job # Dup (ETC)	N4709	P4748	
Original Result	0	0.006	
Duplicate Result	0	0.005	
R Range	0.0	0.001	

ACCURACY			
Job # Spiked (ETC)	N4709		
Original Result	0.0		
Amount Spiked	.200		
Spiked S' Result	.201		
% Recovery	99%		

ANALYZED ON: 11/29/86

TEST: Phenol, UNITS: mg/L  
 METH. REF. # 420.2, MDL: 0.050  
 INSTRUMENT # AA Channel #1

BY: P. Johnson/R. Hill  
 BOOK # CV-12 P. 115  
 FROM PAGE # 114

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	STANDARDS
17639	P2121	0.037	X10	.375	1749	P2514	TH		MATRIX RPT STR	PREP'D ON _____ BY _____ FROM _____ mg/L STOCK # _____
↓	P2123	0.043	STR	BMDL						
17343	P4131	TH	X10	RPT <sup>STR</sup>						
17584	P3067	0.013	-	BMDL						
17639	N9882	TH	X5	RPT <sup>1:100</sup>						
17681	N9884	TH	X5	RPT <sup>1:100</sup>						
↓	N9880	TH	X5	RPT <sup>1:150</sup>						
↓	P2614	TH	X5	RPT <sup>STR/TH</sup>						
17096	P0582	TH	-	STR. RPT						
17824	P4480	TH	X5	RPT <sup>1:10</sup>						
17828	P2472	TH	X5	RPT						
17844	P253	TH	X5	RPT						
17808	P4512	TH	X5	RPT						
17810	P4851	0.067	X5	0.335						
17804	P2554	0.017	-	BMDL						
17804	P4851	0.013	-	BMDL						
↓	P4852	0.017	-	BMDL						
↓	P4853	0.022	-	BMDL						
17886	P5124	0.128	X10	1.28						
17707	N8072	0.048	-	BMDL						
↓	N8075	0.027	-	BMDL						
17721	N8074	0.131	X10	1.31						
1749	P2510	0.031	-	BMDL						
↓	P2513	0.096	X10	.964						

STANDARD mg/L	VALUE OBTAINED	% of Theoretical
300		
200		
100		
050		

SLOPE	Intercept	Corr. Coef

ADDITIONAL COMMENT

Analyses QA/QC Data Report  
 SEE CV-12, P 114

QC Batch # PN003

Verified: [Signature] 11/29/86

METHOD BLANK	PRECISION
Spiked Blank (Known)	Job # Dup (ETC)
Spiked Blank Result	Original Result
% Recovery	Duplicate Result
	R Time
EXTERNAL REF. QC	ACCURACY
Known Value	Job # Spiked (ETC)
Observed Result	Original Result
% Recovery	Amount Spiked
	Spiked S' Result
	% Recovery



ANALYZED ON: 11/29/86

TEST: Phenol, UNITS: mg/L  
 Meth. REF. # AA/EPA420.2, MDL: 0.50  
 INSTRUMENT # AA Channel #2

BY: J. Hernandez/R. Hult  
 BOOK # CV-12 P. 113  
 FROM PAGE #

Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	STANDARDS
1798	P2469	0.030	-	BMDL	17866	P4851	-	-	RPT	PREP'D ON _____ BY _____ FROM _____ mg/L STOCK# _____
	P2471	0.014	-	BMDL		P4852	-	-	RPT	
	P2524	0.012	-	BMDL		P4853	-	-	RPT	
	P2538	0.012	-	BMDL	17868	P4510	0.076	-	0.76	
	P2539	0.012	-	BMDL		P4511	0.054	-	0.54	STANDARD mg/L   VALUE OBTAINED   % of theoretical
	P2540	0.014	-	BMDL		P4512	0.056	-	0.56	
	P3993	0.038	-	BMDL		P4513	0.017	-	BMDL	
	P3994	0.028	-	BMDL	17878	P2776	0.015	-	BMDL	
17801	P4447	0.004	-	BMDL	17886	P3120	0.027	-	BMDL	.300   PKH   85.5% .200   63.5   63.5% .100   33.5   33.5% .050   18.5   18.5%
17823	P4724	0.002	-	BMDL		P3122	0.010	-	BMDL	SLOPE   Intercept   Corr. Coef. 270.169   6.3474   0.99719
17824	P4480	-	-	RPT		P3124	-	-	RPT	
17828	P2472	0.028	-	BMDL		P3126	0.084	-	0.84	ADDITIONAL COMMENTS P4480 - matrix inter-ference P2534 - " " " P4850 - RPT for dilution P4851 - P4852 - P4853 - P3124 - ALL CALLS BY X=MY+B.
	P2473	0.017	-	BMDL	17882	P2543	0.036	-	BMDL	
	P2522	0.017	-	BMDL		P2544	0.060	-	0.60	
	P2536	0.025	-	BMDL		P2545	0.051	-	0.51	
	P2548	0.015	-	BMDL		P2547	0.017	-	BMDL	
17837	P4709	0.015	-	BMDL		P2549	0.015	-	BMDL	
	P4714	0.023	-	BMDL		P2550	0.021	-	BMDL	
17864	P2481	0.025	-	BMDL		P2551	0.031	-	BMDL	
	P2531	0.006	-	BMDL		P2552	0.021	-	BMDL	
	P2541	0.030	-	BMDL						
	P2546	0.017	-	BMDL						
	P2554	-	-	RPT						
17866	P4850	0.321	-	RPT						

Analyses QA/QC Data Report

QC Batch # PH004

Verified: 11/29/86 JPH/Hult

PRECISION			ACCURACY				
METHOD BLANK	0.019		Job # Dup (ETC)	P2469	P4480	P3120	P2551
Spiked Blank (Known)	0.050	0.100	Original Result	0.030	MATRIX	0.028	0.036
Spiked Blank Result	0.045	0.095	Duplicate Result	0.030	Inter-	0.026	0.026
% Recovery	90%	95%	A RANGE			0.002	0.010
EXTERNAL REF. QC	N/A	At 100	Job # Spiked (ETC)	P2469	Inter-	P3120	P2551
Known Value			Original Result	0.030	(CH)	0.027	0.031
Observed Result			Amount Spiked	0.100		0.100	0.100
% Recovery			Spiked S' Result	0.123		0.117	0.130
			% Recovery	91.2		90%	99%

LABORATORY CHRONICLE: Metals Department

Samples P4850 - P4853

	Chemist	Date
Hg Prep	<u>—</u>	<u>—</u>
AA Prep	<u>—</u>	<u>—</u>
ICAP Prep	<u>Kathy Lundberg</u>	<u>11/24/86</u>

Lab Supervisor Joan Komarek Date 11/28/86

SUBCONTRACT  
REQUEST FOR ANALYSIS  
and SAMPLE CHAIN OF CUSTODY

LOGLINK: 17866

MATRIX: WATER

NAME OF SUBCONTRACT LAB: CHYUN

ETC JOB NUMBERS: P4850 P4853  
P4851  
P4852

TURNAROUND IN DAYS: NORMAL DATE DATA REQUIRED: 11/22  
(If deadline cannot be met, contact ETC Subcontract Group).

Send invoice, bill and reports to : ETC Subcontract Group  
(201)225-6786

-----  
Please perform the analyses requested below:

<input type="checkbox"/> Acidity	<input type="checkbox"/> Alkalinity	<input type="checkbox"/> Ammonia (probe)
<input type="checkbox"/> Ammonia (dist)	<input type="checkbox"/> Bicarbonate	<input type="checkbox"/> BOD (5 day)
<input type="checkbox"/> Bromide	<input type="checkbox"/> Carbonate	<input type="checkbox"/> Chloride
<input type="checkbox"/> Chem. Oxygen Dem.	<input type="checkbox"/> Chromium+6	<input type="checkbox"/> Fecal Coliform
<input type="checkbox"/> Total Coliform	<input type="checkbox"/> Color, apparent	<input type="checkbox"/> Cyanide, total
<input type="checkbox"/> Fluoride	<input type="checkbox"/> Formaldehyde (UV)	<input type="checkbox"/> Hardness
<input type="checkbox"/> Nitrate (NO3)	<input type="checkbox"/> Nitrite (NO2)	<input type="checkbox"/> TKN
<input type="checkbox"/> Odor	<input type="checkbox"/> Oil & Grease (grav)	<input type="checkbox"/> T. Organic Carb.
<input type="checkbox"/> Pet. Hydro (IR)	<input type="checkbox"/> Pet. Hydro (grav)	<input type="checkbox"/> Phenolics, tot.
<input type="checkbox"/> Phenolics (5ug/l)	<input type="checkbox"/> Phosphate (ortho)	<input type="checkbox"/> Phosphate (total)
<input type="checkbox"/> Phosphorus (tot.)	<input type="checkbox"/> Silica (dissolved)	<input checked="" type="checkbox"/> Sulfate (SO4)
<input type="checkbox"/> Sulfide (S)	<input type="checkbox"/> Sulfite (SO3)	<input type="checkbox"/> Surfactant (MBAS)
<input type="checkbox"/> Solids, Total	<input type="checkbox"/> Solids, Tot. Diss.	<input type="checkbox"/> Solids, Tot. Set.
<input checked="" type="checkbox"/> Solids, Tot. Sus.	<input type="checkbox"/> Solids, Tot. Vol.	<input type="checkbox"/> Turbidity
<input type="checkbox"/> Gross Alpha, Beta	<input type="checkbox"/> Radium 226	<input type="checkbox"/> Radium 228

Others: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

-----  
Chain of Custody Section for Courier Pick up at ETC

Sample(s) Relinquished by: (ETC) Mark Jekstadt  
Time: 4:30pm Date: 11/21/86

Sample(s) Received by: Mark Kelly  
Time: 4:30pm Date: 11/21/86

Chain of Custody Section for Sending Sample from ETC

Sample Shuttle sealed by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Sample Shuttle opened by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

-----  
Chain of Custody Section for Returning Sample to ETC

Sample Shuttle sealed by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Sample Shuttle opened by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

305257

305258



**Technical Report**

for

**ENVIRONMENTAL STRATEGIES CORP.**

**WASHINGTON HARBOUR**

**3050 KK ST. SUITE 325**

**WASHINGTON, DC 20007**

*Chain of Custody Data Required for ETC Data Management Summary Reports*

<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>
P4853	ENVIRONMENTAL STRATEGIES CORP.	CHAMPION	WCSP-4	861110	1500	



**John J. Fitzgerald**  
*Vice President*  
*Research and Operations*

This Technical Report is an INSITE<sup>SM</sup> service generated by LODESTAR<sup>SM</sup> Data Management Software.

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QA Protocol

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Appendix C - Mass Spectral Data for Tentatively Identified Compounds

Appendix E - Chain of Custody Forms

## Introduction

This report contains the analytical results on your sample. It is designed to include comprehensive data from the entire analytical process in order to satisfy the needs of various levels of review.

The results obtained from your sample are presented in tabular format immediately following this introduction. Quality assurance data is tabulated along with the appropriate sample results for verification. Depending on the analyses ordered, the quality assurance data may include results from blank, spiked blank, spiked sample (i.e. matrix spike) and replicate sample as well as results from surrogate compound analyses. Quality assurance data for verification of proper instrument performance is also included where appropriate. The report appendices include the chain of custody record for your sample and, where appropriate, the gas chromatograms and mass spectra.

The procedures used in the analysis of the sample are described in this report's methodology section. All analytical procedures within our laboratory are performed within a strictly enforced Quality Assurance Protocol. A description of this Protocol is included in the report.

## Results

Sample results, and associated quality assurance data, are always tabulated in one or more of this report's Quantitative Results Tables. The format of each table varies with the class of analysis.

### *Priority Pollutants*

The priority pollutant compounds and elements are listed with their NPDES (National Pollution Discharge Elimination System) numbers, and the Method Detection Limit (MDL) published in the Federal Register. When a compound or element is present below its published MDL it is reported as BMDL (Below Method Detection Limit). When a compound or element is not present at any detectable concentrations it is reported as ND (Not Detected). MDL's for non-aqueous matrices are based on USEPA published MDL's but are adjusted as per sample weight. Matrix spike and replicate analyses, where included, were performed on samples randomly chosen within each quality assurance batch and are therefore not necessarily spikes and replicates of this report's sample. Surrogate compound recovery data and instrument calibration data are included in the Method Performance Data Tables.



**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Volatile Compounds - GC/MS Analysis Data (QR01)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4853** ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-4 861110 1500  
 ETC Sample No. Company Sample Point Date Time Elapsed Hours

NPDES Number	Compound <small>Acrolean and Acrylonitrile values are screen only.</small>	Results			QC Replicate		QC Blank and Spiked Blank			QC Matrix Spike				
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concn. Added ug/l	% Recov	Unspiked Sample ug/l	Concn. Added ug/l	% Recov			
1V	Acrolean	ND	100	ND	ND	ND	800	119	ND	800	88	ND	800	88
2V	Acrylonitrile	ND	100	ND	ND	ND	80.0	91	ND	80.0	80	ND	80.0	80
3V	Benzene	13.5	4.4	ND	ND	ND	18.0	105	71.1	18.0	98	ND	18.0	98
4V	bis(Chloromethyl)ether	ND	10	ND	ND	ND	18.0	-	ND	18.0	-	ND	18.0	-
5V	Bromoform	ND	4.7	ND	ND	ND	18.0	85	ND	18.0	83	ND	18.0	83
6V	Carbon tetrachloride	ND	2.8	ND	ND	ND	18.0	100	ND	18.0	100	ND	18.0	100
7V	Chlorobenzene	ND	6.0	ND	ND	ND	18.0	100	ND	18.0	101	ND	18.0	101
8V	Chlorodibromomethane	ND	3.1	ND	ND	ND	18.0	92	ND	18.0	94	ND	18.0	94
9V	Chloroethane	ND	10	ND	ND	ND	18.0	96	ND	18.0	96	ND	18.0	96
10V	2-Chloroethylvinyl ether	ND	10	ND	ND	ND	18.0	0 <sup>a</sup>	ND	18.0	0 <sup>a</sup>	ND	18.0	0 <sup>a</sup>
11V	Chloroform	ND	1.6	ND	ND	ND	18.0	92	ND	18.0	104	ND	18.0	104
12V	Dichlorobromomethane	ND	2.2	ND	ND	ND	18.0	102	ND	18.0	101	ND	18.0	101
13V	Dichlorodifluoromethane	ND	10	ND	ND	ND	18.0	93	ND	18.0	91	ND	18.0	91
14V	1,1-Dichloroethane	ND	4.7	ND	ND	ND	18.0	95	ND	18.0	101	ND	18.0	101
15V	1,2-Dichloroethane	ND	2.8	ND	ND	ND	18.0	94	ND	18.0	98	ND	18.0	98
16V	1,1-Dichloroethylene	ND	2.8	ND	ND	ND	18.0	96	ND	18.0	97	ND	18.0	97
17V	1,2-Dichloropropane	ND	6.0	ND	ND	ND	18.0	99	ND	18.0	98	ND	18.0	98
18V	cis-1,3-Dichloropropylene	ND	5.0	ND	ND	ND	18.0	94	ND	18.0	95	ND	18.0	95
19V	Ethylbenzene	ND	7.2	ND	ND	ND	18.0	101	ND	18.0	110	ND	18.0	110
20V	Methyl bromide	ND	10	ND	ND	ND	18.0	118	ND	18.0	125	ND	18.0	125
21V	Methyl chloride	ND	10	ND	ND	ND	18.0	104	ND	18.0	92	ND	18.0	92
22V	Methylene chloride	13.9	2.8	ND	ND	ND	18.0	73	7.63	18.0	115	ND	18.0	115
23V	1,1,2,2-Tetrachloroethane	ND	6.9	ND	ND	ND	18.0	113	ND	18.0	103	ND	18.0	103
24V	1,1,2,2-Tetrachloroethylene	19.7	4.1	ND	ND	ND	18.0	97	ND	18.0	100	24.5	18.0	100
25V	Toluene	ND	6.0	ND	ND	ND	18.0	101	ND	18.0	100	ND	18.0	100
26V	1,2-Trans-dichloroethylene	150	1.6	ND	ND	ND	18.0	93	ND	18.0	95	ND	18.0	95
27V	1,1,1-Trichloroethane	ND	3.8	ND	ND	ND	18.0	106	ND	18.0	102	ND	18.0	102
28V	1,1,1,2-Trichloroethane	ND	5.0	ND	ND	ND	18.0	93	ND	18.0	95	ND	18.0	95
29V	Trichloroethylene	103	1.9	ND	ND	ND	18.0	93	ND	18.0	85	ND	18.0	85
30V	Trichlorofluoromethane	ND	10	ND	ND	ND	18.0	92	ND	18.0	97	ND	18.0	97
31V	Vinyl chloride	ND	10	ND	ND	ND	18.0	96	ND	18.0	92	ND	18.0	92
18V	trans-1,3-Dichloropropylene	ND	10	ND	ND	ND	18.0	103	ND	18.0	105	ND	18.0	105

<sup>a</sup> Recovery normally variable using established methodology.

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Acid Compounds - GC/MS Analysis Data (QR02)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4853 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-4 861110 1500  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results			QC Replicate			QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concen. Added ug/l	% Recov	Unspiked Sample ug/l	Concen. Added ug/l	% Recov		
1A	2-Chlorophenol	ND	3.5	ND	ND	ND	100	89	ND	102	107		
2A	2,4-Dichlorophenol	ND	2.9	ND	ND	ND	100	92	ND	102	100		
3A	2,4-Dimethylphenol	ND	2.9	ND	ND	ND	100	18 <sup>a</sup>	ND	102	100		
4A	4,6-Dinitro-o-cresol	ND	26	ND	ND	ND	100	92	ND	102	90		
5A	2,4-Dinitrophenol	ND	45	ND	ND	ND	100	77	ND	102	88		
6A	2-Nitrophenol	ND	3.9	ND	ND	ND	100	80	ND	102	89		
7A	4-Nitrophenol	ND	2.6	ND	ND	ND	100	85	ND	102	97		
8A	p-Chloro-m-cresol	ND	3.2	ND	ND	ND	100	77	ND	102	68		
9A	Pentachlorophenol	ND	3.9	ND	ND	ND	100	74	ND	102	82		
10A	Phenol	ND	1.6	ND	ND	ND	100	89	ND	102	109		
11A	2,4,6-Trichlorophenol	ND	2.9	ND	ND	ND	100	89	ND	102	94		

<sup>a</sup> Recovery normally variable using established methodology.

**ETC**

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**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA  
BASE/NEUTRAL COMPOUNDS - GC/MS ANALYSIS DATA (QR03)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4853 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WSCP-4 861110 1500**  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results		QC Replicate		QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concn. Added ug/l	% Recov	Unspiked Sample ug/l	Concn. Added ug/l	% Recov
1B	Acenaphthene	ND	2.0	ND	ND	ND	100	87	ND	102	87
2B	Acenaphthylene	ND	3.8	ND	ND	ND	100	90	ND	102	92
3B	Anthracene	ND	2.0	ND	ND	ND	100	84	ND	102	86
4B	Benztidine	ND	47	ND	ND	ND	100	0 <sup>a</sup>	ND	102	3 <sup>a</sup>
5B	Benzo(a)anthracene	ND	8.4	ND	ND	ND	100	89	ND	102	90
6B	Benzo(a)pyrene	ND	2.7	ND	ND	ND	100	98	ND	102	99
7B	Benzo(b)fluoranthene	ND	11	ND	ND	ND	100	109	ND	102	105
8B	Benzo(ghi)perylene	ND	4.4	ND	ND	ND	0	-	ND	0	-
9B	Benzo(k)fluoranthene	ND	3.8	ND	ND	ND	100	78	ND	102	87
10B	bis(2-Chloroethoxy)methane	ND	5.7	ND	ND	ND	100	80	ND	102	83
11B	bis(2-Chloroethyl) ether	ND	6.1	ND	ND	ND	100	87	ND	102	89
12B	bis(2-Chloroisopropyl) ether	ND	6.1	ND	ND	ND	100	88	ND	102	92
13B	bis(2-Ethylhexyl)phthalate	ND	11	ND	ND	ND	100	92	ND	102	93
14B	4-Bromophenyl phenyl ether	ND	2.0	ND	ND	ND	100	99	ND	102	95
15B	Butyl benzyl phthalate	ND	11	ND	ND	ND	100	75	ND	102	78
16B	2-Chloronaphthalene	ND	2.0	ND	ND	ND	100	87	ND	102	85
17B	4-Chlorophenyl phenyl ether	ND	4.5	ND	ND	ND	100	87	ND	102	91
18B	Chrysene	ND	2.7	ND	ND	ND	100	88	ND	102	88
19B	Dibenzo(a,h)anthracene	ND	11	ND	ND	ND	0	-	ND	0	-
20B	1,2-Dichlorobenzene	ND	2.0	ND	ND	ND	100	76	ND	102	76
21B	1,3-Dichlorobenzene	ND	2.0	ND	ND	ND	100	67	ND	102	72
22B	1,4-Dichlorobenzene	ND	4.7	ND	ND	ND	100	71	ND	102	75
23B	3,3'-Dichlorobenzidine	ND	18	ND	ND	ND	100	52	ND	102	54
24B	Diethyl phthalate	ND	11	ND	ND	ND	100	45	ND	102	48
25B	Dimethyl phthalate	ND	11	ND	ND	ND	100	22 <sup>a</sup>	ND	102	18 <sup>a</sup>
26B	Di-n-butyl phthalate	ND	11	ND	ND	ND	100	63	ND	102	61
27B	2,4-Dinitrotoluene	ND	6.1	ND	ND	ND	100	81	ND	102	110
28B	2,6-Dinitrotoluene	ND	2.0	ND	ND	ND	100	94	ND	102	106
29B	Di-n-octyl phthalate	ND	11	ND	ND	ND	100	101	ND	102	103
30B	1,2-Diphenylhydrazine	ND	11	ND	ND	ND	100	88	ND	102	96
31B	Fluoranthene	ND	2.4	ND	ND	ND	100	70	ND	102	58
32B	Fluorene	ND	2.0	ND	ND	ND	100	88	ND	102	95

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**ETC** ENVIRONMENTAL TESTING and CERTIFICATION

NOV 30, 1986

**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA  
BASE/NEUTRAL COMPOUNDS - GC/MS ANALYSIS DATA (QR03)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4853** ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-4 861110 1500  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results			QC Replicate			QC Blank and Spiked Blank			QC Matrix Spike		
		Sample Concn. ug/l	MDL ug/l		First ug/l	Second ug/l	Blank Data ug/l	Concn. Added ug/l	% Recov	Unspiked Sample ug/l	Concn. Added ug/l	% Recov	
33B	Hexachlorobenzene	ND	2.0		ND	ND	ND	100	93	ND	102	94	
34B	Hexachlorobutadiene	ND	1.97		ND	ND	ND	100	71	ND	102	75	
35B	Hexachlorocyclopentadiene	ND	1.7		ND	ND	ND	0	-	ND	0	-	
36B	Hexachloroethane	ND	1.7		ND	ND	ND	100	62	ND	102	63	
37B	Indeno(1,2,3-c,d)pyrene	ND	5.1		ND	ND	ND	0	-	ND	0	-	
38B	Isophorone	ND	2.4		ND	ND	ND	100	76	ND	102	80	
39B	Naphthalene	ND	1.7		ND	ND	ND	100	80	ND	102	84	
40B	Nitrobenzene	ND	2.0		ND	ND	ND	100	89	ND	102	88	
41B	N-Nitrosodimethylamine	ND	1.1		ND	ND	ND	0	-	ND	0	-	
42B	N-Nitrosodi-n-propylamine	ND	1.1		ND	ND	ND	100	65	ND	102	85	
43B	N-Nitrosodiphenylamine	ND	2.0		ND	ND	ND	100	83	ND	102	110	
44B	Phenanthrene	ND	5.8		ND	ND	ND	100	89	ND	102	89	
45B	Pyrene	ND	2.0		ND	ND	ND	100	69	ND	102	55	
46B	1,2,4-Trichlorobenzene	ND	2.0		ND	ND	ND	100	158 <sup>a</sup>	ND	102	152 <sup>a</sup>	

<sup>a</sup> Recovery normally variable using established methodology.

**ETC**

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**TABLE 1: QUANTITATIVE RESULTS and QUALITY ASSURANCE DATA**  
**Pesticide/PCB Compounds - GC/MS Analysis Data (QR04)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
**P4853 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WCSP-4 861110 1500**  
 ETC Sample No. Company Sample Point Date Time Elapsed Hours

NPDES Number	Compound	Results		QC Replicate		QC Blank and Spiked Blank		QC Matrix Spike			
		Sample Concn. ug/l	MDL ug/l	First ug/l	Second ug/l	Blank Data ug/l	Concen. Added ug/l	% Recov	Unspiked Sample ug/l	Concen. Added ug/l	% Recov
1P	Aldrin	ND	2.0	ND	ND	ND	100	89	ND	102	55
2P	Alpha-BHC	ND	11	ND	ND	ND	100	39	ND	102	34
3P	Beta-BHC	ND	4.7	ND	ND	ND	100	57	ND	102	67
4P	Gamma-BHC	ND	11	ND	ND	ND	100	34	ND	102	32
5P	Delta-BHC	ND	3.3	ND	ND	ND	100	46	ND	102	44
6P	Chlordane	ND	11	ND	ND	ND	200	69	ND	204	56
7P	4,4'-DDT	ND	3.0	ND	ND	ND	100	72	ND	102	36
8P	4,4'-DDE	ND	6.0	ND	ND	ND	100	84	ND	102	47
9P	4,4'-DDD	ND	5.1	ND	ND	ND	100	71	ND	102	46
10P	Dieldrin	ND	2.7	ND	ND	ND	100	83	ND	102	46
11P	Endosulfan I	ND	11	ND	ND	ND	100	29	ND	102	82
12P	Endosulfan II	ND	11	ND	ND	ND	100	30	ND	102	11 <sup>a</sup>
13P	Endosulfan sulfate	ND	6.0	ND	ND	ND	100	44	ND	102	8 <sup>a</sup>
14P	Endrin	ND	11	ND	ND	ND	100	74	ND	102	44
15P	Endrin aldehyde	ND	11	ND	ND	ND	100	29	ND	102	72
16P	Heptachlor epoxide	ND	2.4	ND	ND	ND	100	97	ND	102	30
17P	Heptachlor epoxide	ND	2.4	ND	ND	ND	100	86	ND	102	68
18P	PCB-1242	ND	39	ND	ND	ND	0	-	ND	0	106
19P	PCB-1254	ND	39	ND	ND	ND	0	-	ND	0	-
20P	PCB-1221	ND	39	ND	ND	ND	0	-	ND	0	-
21P	PCB-1232	ND	39	ND	ND	ND	0	-	ND	0	-
22P	PCB-1248	ND	39	ND	ND	ND	0	-	ND	0	-
23P	PCB-1260	ND	39	ND	ND	ND	100	93	ND	102	49
24P	PCB-1016	ND	39	ND	ND	ND	0	-	ND	0	-
25P	Toxaphene	ND	11	ND	ND	ND	0	-	ND	0	-

<sup>a</sup> Recovery normally variable using established methodology.

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**TABLE 1: QUANTITATIVE RESULTS**  
**Metals - Analysis Data (QR52)**

Chain of Custody Data Required for ETC Data Management Summary Reports  
 P4853 ENVIRONMENTAL STRATEGIES CORP. CHAMPION WOSP-4 861110 1500  
 ETC Sample No. Company Facility Sample Point Date Time Elapsed Hours

Compound	Results	
	Sample Concn. ug/l	MDL ug/l
Barium	66	3.2
Chromium	BMDL	26
Copper	BMDL	13
Iron	145000	180
Manganese	7360	3.7
Nickel	11	11
Sodium	1480000	87
Zinc	150	2.9



**TABLE 2: METHOD PERFORMANCE DATA**  
**Surrogate Recovery - Aqueous Matrices (QR20)**

**P4853**  
 Chain of Custody Data Required for ETC Data Management Summary Reports

ETC Sample No.	Company	Facility	Sample Point	Date	Time	Elapsed Hours

Compound	Amount Added ug	% Recovery	Control Limits *	
			Lower	Upper
<b>VOLATILE FRACTION (GC/MS)</b>				
Toluene-D8	250	98	88	110
Bromofluorobenzene	250	96	86	115
1,2-Dichloroethane-D4	250	97	76	114
<b>ACID FRACTION (GC/MS)</b>				
Phenol-D5	100	87	10	94
2-Fluorophenol	100	87	21	100
2,4,6-Tribromophenol	100	72	10	123
<b>BASE/NEUTRAL FRACTION (GC/MS)</b>				
Nitrobenzene-D5	50	88	35	114
2-Fluorobiphenyl	50	59	43	116
Terphenyl-D14	50	12***	33	141
<b>PESTICIDE/PCB FRACTION (GC/MS)</b>				
Dibutylchloroendate	-	-	24**	154**

\* IFB EPA Control Limits.  
 \*\* Advisory Limits Only.  
 \*\*\* Recovery manually verified.

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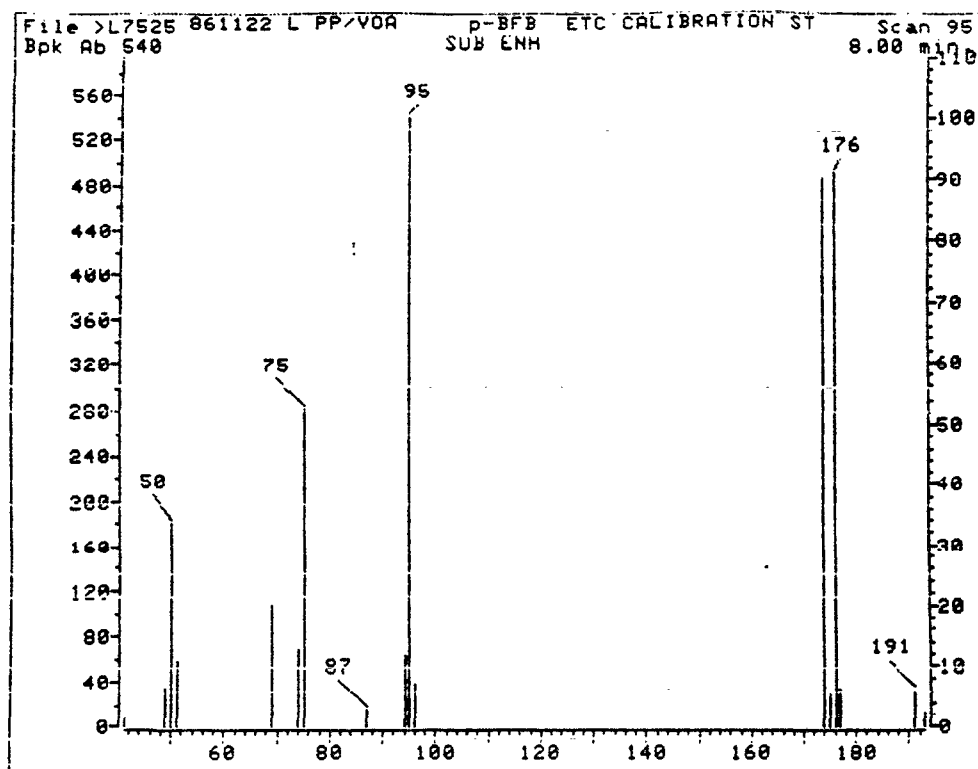


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	33.35	33.35	OK
75	30-60% of mass 95	52.25	52.25	UK
95	Base peak, 100% relative abundance	100.00	100.00	OK
96	5-9% of mass 95	7.13	7.13	UK
173	Less than 1% of mass 95	0.00	0.00	OK
174	Greater than 50% of mass 95	90.46	90.46	UK
175	5-9% of mass 174	5.56	6.14	OK
176	95-101% of mass 174	91.20	100.82	UK
177	5-9% of mass 176	6.16	6.75	OK

Injection Date: 11/22/86  
 Injection Time: 21:11  
 Run No: >L7525  
 Spectrun No: 95

Analyst: *J. Quinn*  
 Processor: *DAVID CHRISTOPHER*  
 QC Batch: *075936*  
 Samples: *P3328, P3329, P3330, P3331, P3332, P3333, P3334, P3335, P3336, P3337, P3338, P3339, P3340, P3341, P3342, P3343, P3344, P3345, P3346, P3347, P3348, P3349, P3350, P3351, P3352, P3353, P3354, P3355, P3356, P3357, P3358, P3359, P3360, P3361, P3362, P3363, P3364, P3365, P3366, P3367, P3368, P3369, P3370, P3371, P3372, P3373, P3374, P3375, P3376, P3377, P3378, P3379, P3380, P3381, P3382, P3383, P3384, P3385, P3386, P3387, P3388, P3389, P3390, P3391, P3392, P3393, P3394, P3395, P3396, P3397, P3398, P3399, P3400, P3401, P3402, P3403, P3404, P3405, P3406, P3407, P3408, P3409, P3410, P3411, P3412, P3413, P3414, P3415, P3416, P3417, P3418, P3419, P3420, P3421, P3422, P3423, P3424, P3425, P3426, P3427, P3428, P3429, P3430, P3431, P3432, P3433, P3434, P3435, P3436, P3437, P3438, P3439, P3440, P3441, P3442, P3443, P3444, P3445, P3446, P3447, P3448, P3449, P3450, P3451, P3452, P3453, P3454, P3455, P3456, P3457*

*WJ*  
*11/22/86*

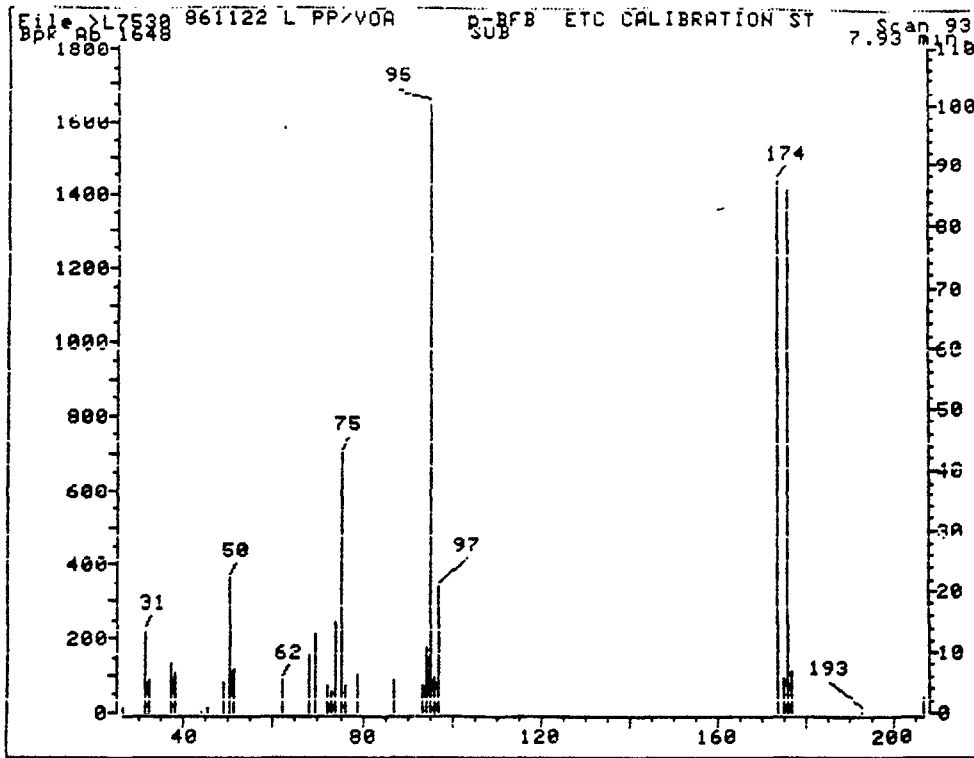


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	22.03	22.03	OK
75	30-60% of mass 95	42.60	42.60	OK
95	Base peak, 100% relative abundance	100.00	100.00	OK
96	5-9% of mass 95	6.01	6.01	OK
173	Less than 1% of mass 95	0.00	0.00	OK
174	Greater than 50% of mass 95	87.32	87.32	OK
175	5-9% of mass 174	5.58	6.39	OK
176	95-101% of mass 174	85.92	98.40	OK
177	5-9% of mass 176	6.80	7.91	OK

Injection Date: 11/23/86  
 Injection Time: 00:42  
 Run No: >L7550  
 Spectrum No: 93

Analyst: J. Quinn  
 Processor: (unreadable)  
 GC Batch: QV 5826  
 Samples: P2762, P2763, P2764, P2765, P2766, P2767, P2768, P2769, P2770, P2771, P2772, P2773, P2774, P2775, P2776, P2777, P2778, P2779, P2780, P2781, P2782, P2783, P2784, P2785, P2786, P2787, P2788, P2789, P2790, P2791, P2792, P2793, P2794, P2795, P2796, P2797, P2798, P2799, P2800, P2801, P2802, P2803, P2804, P2805, P2806, P2807, P2808, P2809, P2810, P2811, P2812, P2813, P2814, P2815, P2816, P2817, P2818, P2819, P2820, P2821, P2822, P2823, P2824, P2825, P2826, P2827, P2828, P2829, P2830, P2831, P2832, P2833, P1957

*Handwritten note:*  
 VS  
 11/29/86

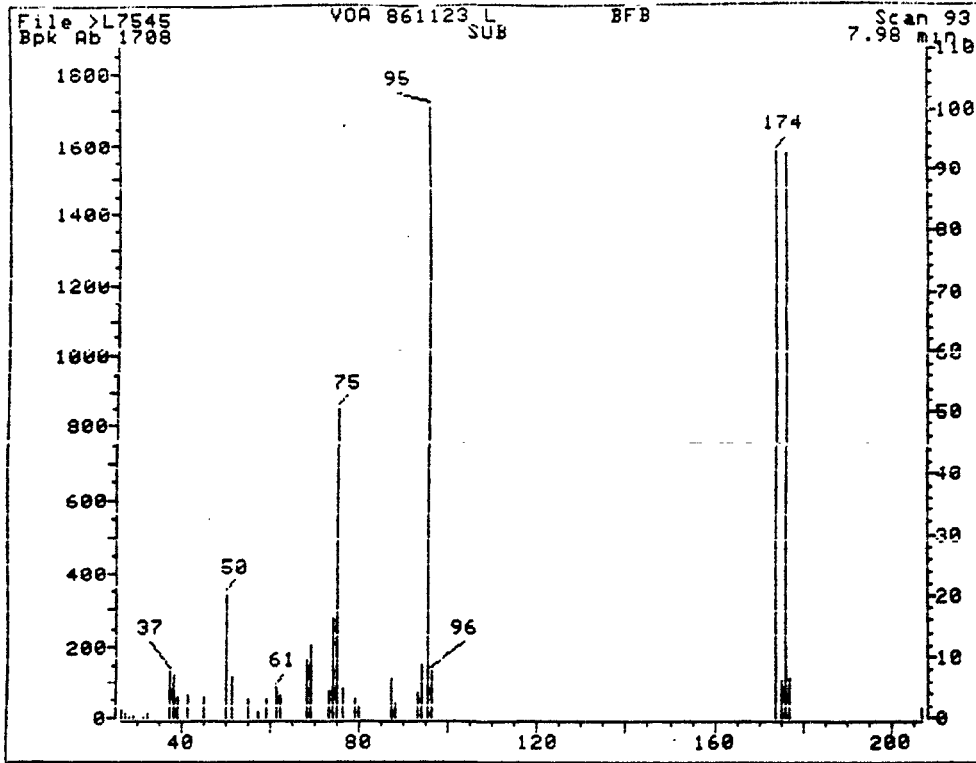


TABLE 2: METHOD PERFORMANCE DATA (QR21)

GC/MS Tuning Data - Bromofluorobenzene (BFB) for Volatiles Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
50	15-40% of mass 95	20.02	20.02	Ok
75	30-60% of mass 95	49.82	49.82	Ok
95	Base peak, 100% relative abundance	100.00	100.00	Ok
96	5-9% of mass 95	7.67	7.67	Ok
173	Less than 1% of mass 95	0.00	0.00	Ok
174	Greater than 50% of mass 95	92.51	92.51	Ok
175	5-9% of mass 174	6.32	6.84	Ok
176	95-101% of mass 174	92.21	99.68	Ok
177	5-9% of mass 176	6.38	6.92	Ok

Injection Date: 11/23/86  
 Injection Time: 13:45  
 Run No: >L7545  
 Spectrum No: 93

Analyst: *Paul Delato*  
 Processor: *DURON CHRISTOPHER*  
 QC Batch: *QV 5936*  
 Samples: *P3325, P3326, P3327, P3328, P3330, P3332, P3333, P1957*

*KS*  
*11/29/86*

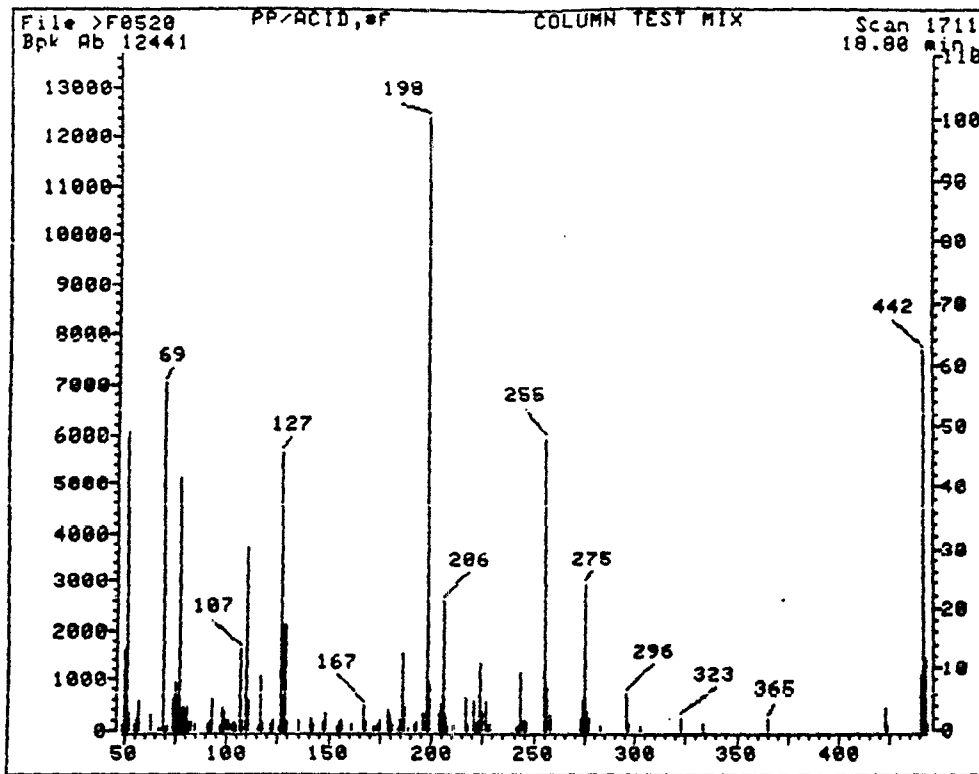


TABLE 2: METHOD PERFORMANCE DATA (QR22)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFTPP) for Acids Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	48.52	48.52	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	56.51	56.51	Ok
70	Less than 2% of mass 69	.42	.74	Ok
127	40-60% of mass 198	45.35	45.35	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	7.19	7.19	Ok
275	10-30% of mass 198	23.50	23.50	Ok
365	Greater than 1% of mass 198	1.89	1.89	Ok
441	0-100% of mass 443	8.79	76.34	Ok
442	Greater than 40% of mass 198	62.49	62.49	Ok
443	17-23% of mass 442	11.52	18.43	Ok

Injection Date: 11/26/86  
Injection Time: 07:29  
Run No: >F0520  
Spectrum No: 1711

Analyst: G. AO  
Processor: David T. ...  
QC Batch: NA6013  
Samples: G0447-G0450, G0457, G0458,  
G0467, G0468, G0471, G0472,  
G0489, G0493-G0495, G4952,  
G4953, G0472

WS  
11/29/86

305274

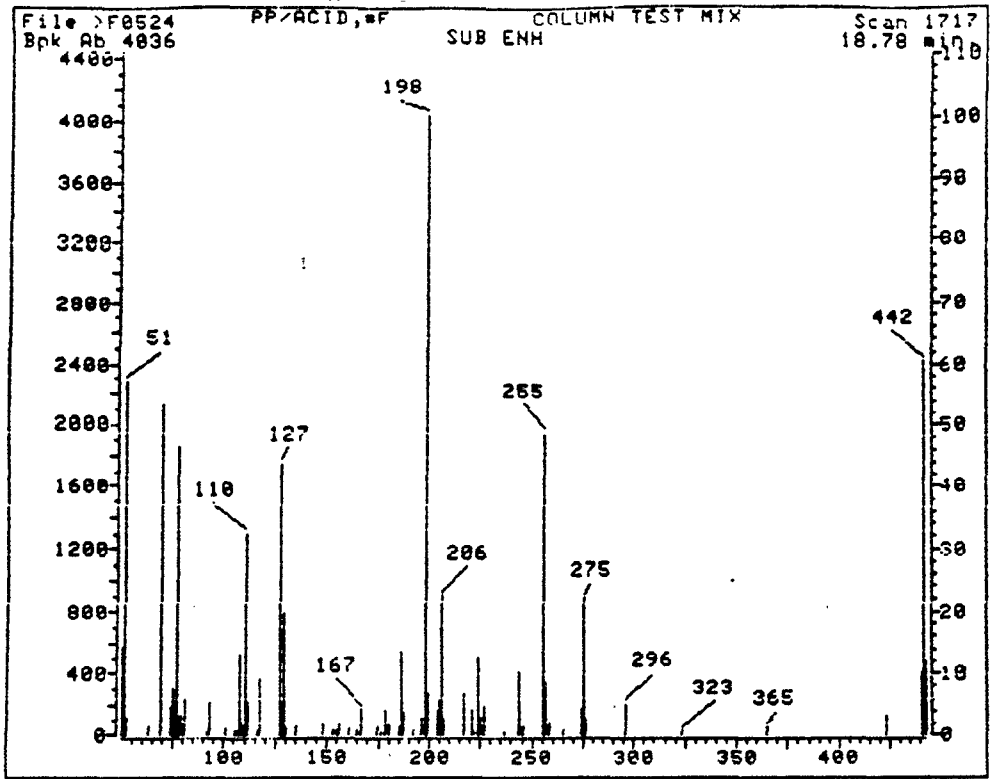


TABLE 2: METHOD PERFORMANCE DATA (QR22)

GC/MS Tuning Data - Decafluorotriphenylphospine (DFIPP) for Acids Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	56.57	56.57	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	52.49	52.49	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	43.39	43.39	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	6.54	6.54	Ok
275	10-30% of mass 198	21.76	21.76	Ok
365	Greater than 1% of mass 198	1.26	1.26	Ok
441	0-100% of mass 443	10.00	84.54	Ok
442	Greater than 40% of mass 198	60.51	60.51	Ok
443	17-23% of mass 442	11.83	19.55	Ok

*KS*  
*11/25/86*

Injection Date: 11/26/86  
 Injection Time: 11:04  
 Run No: >F0524  
 Spectrun No: 1717

Analyst: *g ad*  
 Processor: *Chad Trant*  
 QC Batch: *QA 6013*  
 Samples: *P0447 - P0450, P452, P453, P0457, P0458, P0467, P0468, P0471, P0472.*

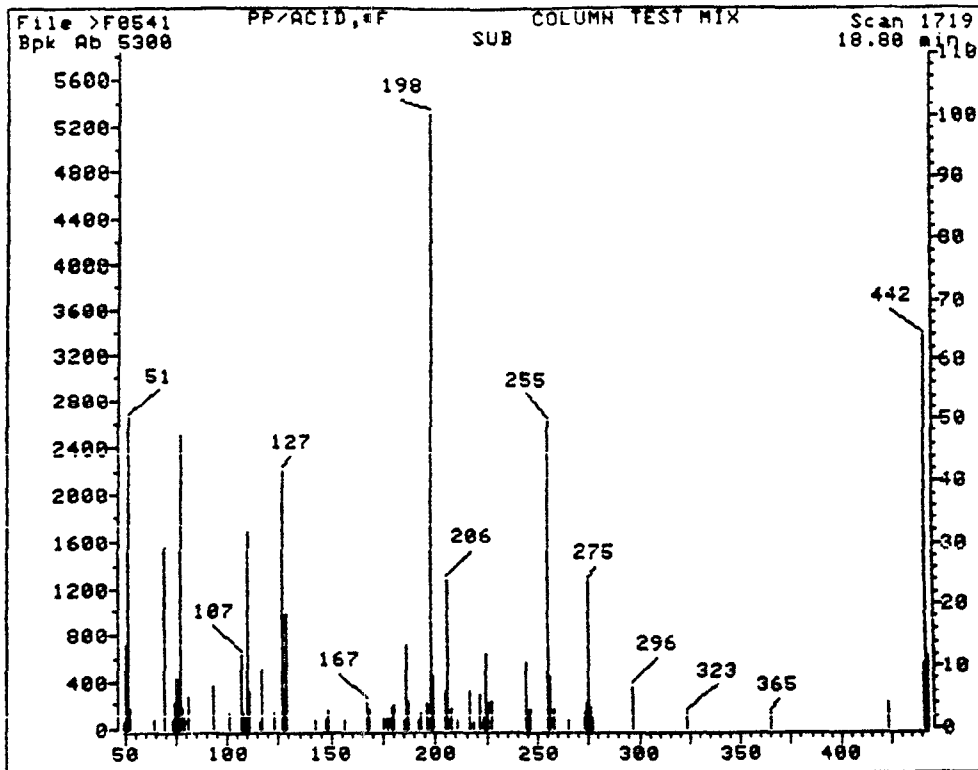


TABLE 2: METHOD PERFORMANCE DATA (QR22)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFTPP) for Acids Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	49.92	49.92	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	29.34	29.34	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	41.43	41.43	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	8.38	8.38	Ok
275	10-30% of mass 198	23.66	23.66	Ok
365	Greater than 1% of mass 198	2.28	2.28	Ok
441	0-100% of mass 443	10.34	88.96	Ok
442	Greater than 40% of mass 198	63.66	63.66	Ok
443	17-23% of mass 442	11.62	18.26	Ok

KS  
11/29/86

Injection Date: 11/26/86  
Injection Time: 22:39  
Run No: >F0541  
Spectrum No: 1719

Analyst:                       
Processor:                       
QC Batch:                       
Samples:                     

305276

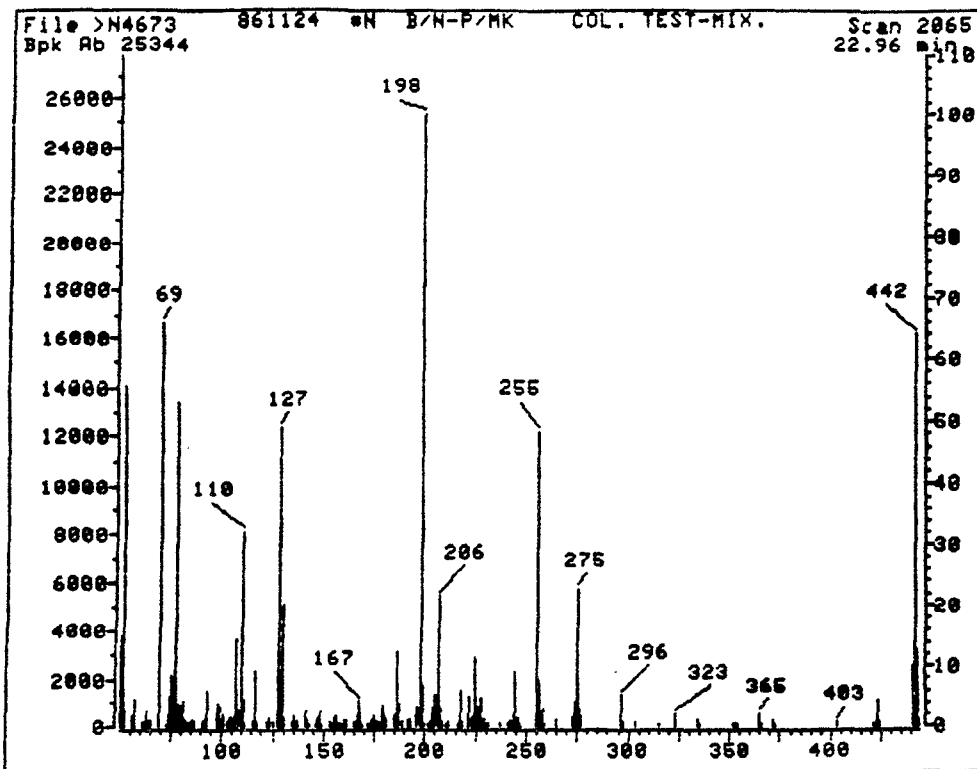


TABLE 2: METHOD PERFORMANCE DATA (QR23)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFTPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	55.50	55.50	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	65.88	65.88	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	48.83	48.83	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	6.67	6.67	Ok
275	10-30% of mass 198	22.60	22.60	Ok
365	Greater than 1% of mass 198	2.14	2.14	Ok
441	0-100% of mass 443	9.87	78.53	Ok
442	Greater than 40% of mass 198	64.00	64.00	Ok
443	17-23% of mass 442	12.57	19.64	Ok

Injection Date: 11/25/86  
Injection Time: 14:37  
Run No: >N4673  
Spectrum No: 2065

Analyst: J. Hendell  
Processor: J. W. C.  
QC Batch: QB6013  
Samples: 70447-70450, 70457, 70458  
70467, 70468, 70471, 70472, 70489  
70493-70495, 70852, 70453

LAM  
11-30

305277

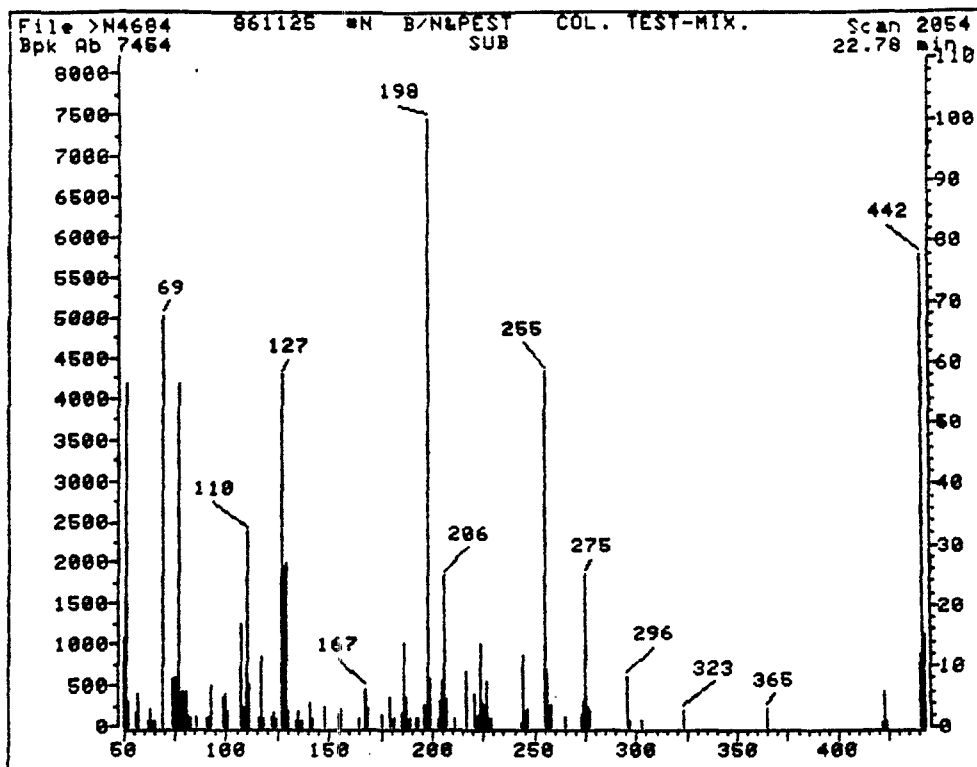


TABLE 2: METHOD PERFORMANCE DATA (QR23)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFTPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	56.12	56.12	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	67.41	67.41	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	57.59	57.59	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	7.90	7.90	Ok
275	10-30% of mass 198	25.03	25.03	Ok
365	Greater than 1% of mass 198	3.03	3.03	Ok
441	0-100% of mass 443	12.03	79.24	Ok
442	Greater than 40% of mass 198	77.68	77.68	Ok
443	17-23% of mass 442	15.19	19.55	Ok

Injection Date: 11/26/86  
 Injection Time: 03:17  
 Run No: >N4684  
 Spectrum No: 2054

Analyst: [Signature]  
 Processor: W.N.C.  
 QC Batch: 026013  
 Samples: P0447-50, P0450, P0451, P0455, P0467, P0468, P0471, P0472, P0493-P0495, P4852, P4853

LAN  
11-30

305278



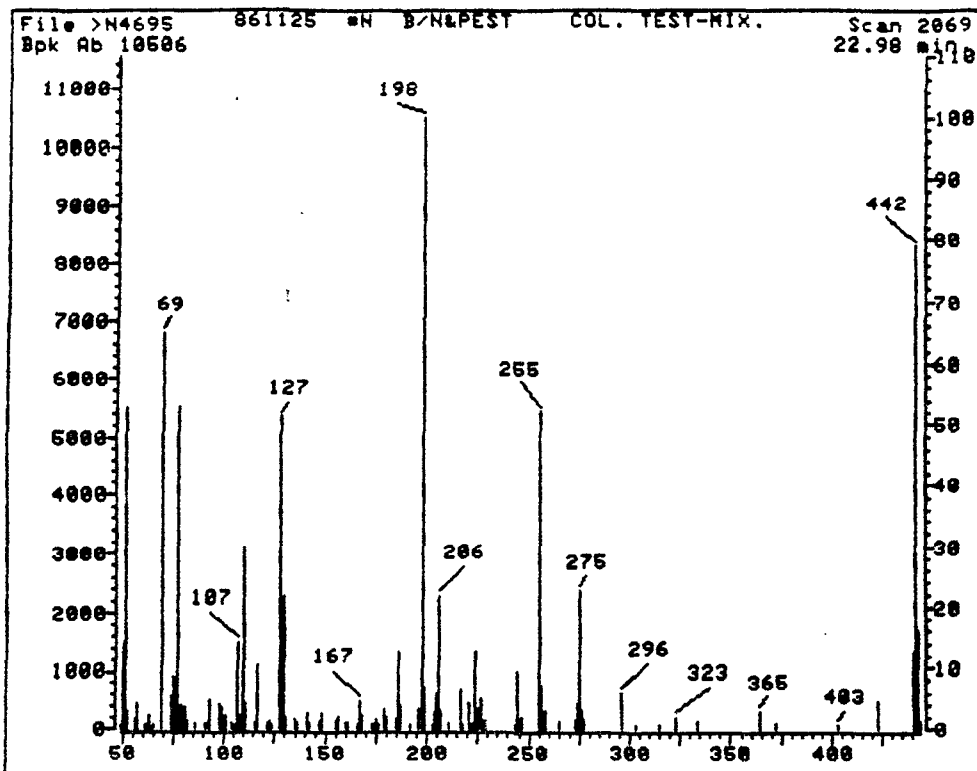


TABLE 2: METHOD PERFORMANCE DATA (QR23)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFTPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	30-60% of mass 198	52.28	52.28	Ok
68	Less than 2% of mass 69	0.00	0.00	Ok
69	(reference only)	64.83	64.83	Ok
70	Less than 2% of mass 69	0.00	0.00	Ok
127	40-60% of mass 198	50.89	50.89	Ok
197	Less than 1% of mass 198	0.00	0.00	Ok
198	Base peak, 100% relative abundance	100.00	100.00	Ok
199	5-9% of mass 198	6.94	6.94	Ok
275	10-30% of mass 198	22.61	22.61	Ok
365	Greater than 1% of mass 198	3.04	3.04	Ok
441	0-100% of mass 443	12.55	77.32	Ok
442	Greater than 40% of mass 198	79.16	79.16	Ok
443	17-23% of mass 442	16.24	20.51	Ok

Injection Date: 11/26/86  
Injection Time: 12:15  
Run No: >N4695  
Spectrum No: 2069

Analyst: [Signature]  
Processor: W. D. C. L.  
QC Batch: 006013  
Samples: P0447-P0450, P0457, P0458  
P0467, P0468.

LAH  
11-30

305279

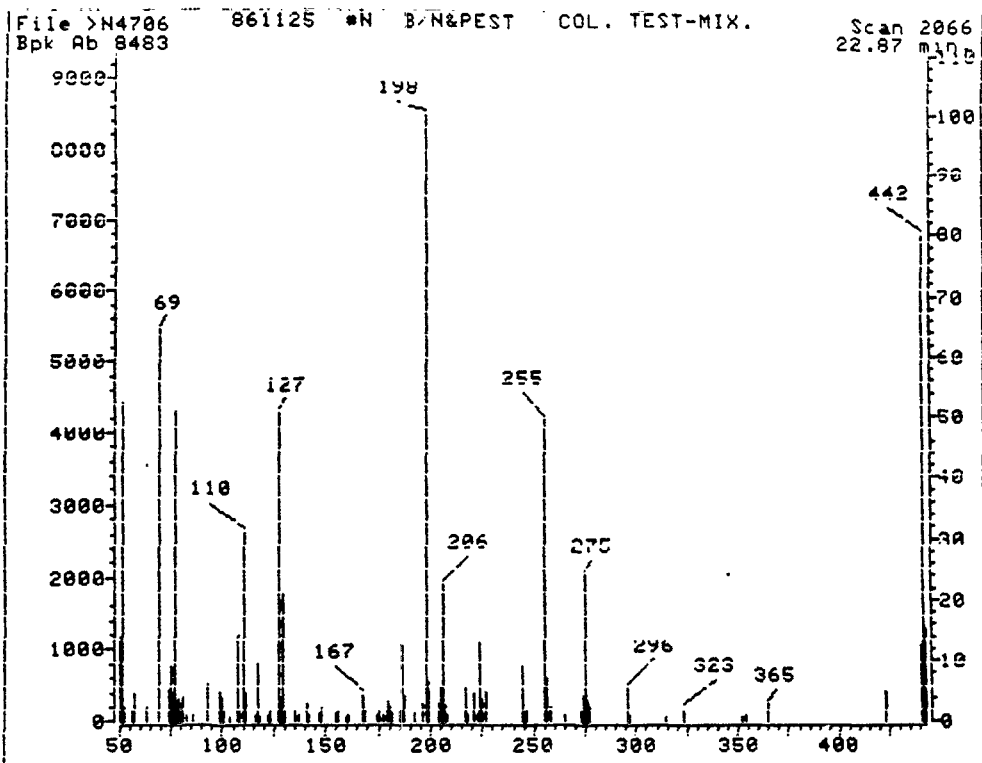


TABLE 2: METHOD PERFORMANCE DATA (WR25)

GC/MS Tuning Data - Decafluorotriphenylphosphine (DFIPP) for Base/Neutral Analysis

m/z	Ion Abundance Criteria	% Relative Abundance Base Peak	% Relative Abundance Appropriate Peak	Status
51	50-60% of mass 198	52.22	52.22	UK
68	Less than 2% of mass 69	0.00	0.00	UK
69	(reference only)	64.29	64.29	UK
70	Less than 2% of mass 69	0.00	0.00	UK
127	40-60% of mass 198	50.49	50.49	UK
197	Less than 1% of mass 198	0.00	0.00	UK
198	Base peak, 100% relative abundance	100.00	100.00	UK
199	5-9% of mass 198	6.64	6.64	UK
275	10-30% of mass 198	24.07	24.07	UK
365	Greater than 1% of mass 198	2.91	2.91	UK
441	0-100% of mass 442	12.58	82.33	UK
442	Greater than 40% of mass 198	19.88	19.88	UK
443	1/-25% of mass 442	15.28	19.13	UK

Injection Date: 11/26/86  
 Injection Time: 21:29  
 Run No: >N4/U6  
 Spectrum No: 2066

Analyst: Shawn Malachuk  
 Processor: A. N. C.  
 GC Batch: 086013  
 Samples: P0471, P0472, P0489, P0493, P0494, P0495, P0496, P0497, P0498, P0499

305280

LAH  
11-30

## Methodology for GC/MS Analysis of Priority Pollutant Compounds

The methods employed in the GC/MS analysis for priority pollutants are established EPA methods. Rigid compliance with the instrument parameters and performance criteria of the published methods was achieved. In some cases, the precise amounts of sample used and the sample handling procedures vary with the complexity of the sample matrix. Qualitative identification of the priority pollutants was performed using the relative retention times, the relative abundance of three characteristic ions and the abundance ratios. The entire mass spectrum was reviewed to confirm each identification. Quantitative analysis of detected compounds was performed by using a response factor generated by a major characteristic ion of the specific compound and an internal standard.

Compounds, in addition to those on the priority pollutant list, were identified through a computer-aided search of the NBS-EPA spectra library. After review the identifications are included in a separate tabulation and labelled "tentatively identified".

### ***Volatile Priority Pollutant Compounds***

For the analysis of Volatile priority pollutants, EPA Method 624 was used. The method can be summarized as follows: Helium is bubbled through a 5 ml water sample contained in a specially designed purging chamber at ambient temperature. The purgeable volatile organic compounds are transferred from the aqueous phase to the vapor phase. The vapor is swept through a sorbent column where the organic components are trapped. After the purge cycle is complete, the sorbent column is heated and backflushed with helium to desorb the organic purgeables onto a gas chromatographic column. The gas chromatograph is temperature programmed to separate the purgeable mixture. The separated purgeable components are then identified and quantitated using a computerized mass spectrometer.

### ***Acid, Base/Neutral and Pesticide Priority Pollutant Compounds***

For the analysis of the Acid, Base/Neutral and Pesticide priority pollutants in an aqueous liquid matrix, EPA Method 625 was used. The method can be summarized as follows: A measured volume of sample, approximately 1 liter, is adjusted to a pH greater than 11 and extracted with methylene chloride. The pH of the sample is adjusted to a value less than 2 and extracted with an aliquot of fresh methylene chloride. A separatory funnel or continuous extractor is used to perform the extractions. The two extracts are dried and concentrated to a 1 ml final volume. Each extract is injected into a GC/MS instrument specifically configured for the correct fraction.

## Methodology for Analysis of Metals

### **AQUEOUS**

The determination of metals in aqueous samples is performed according to the methods published by EPA in "Methods for Chemical Analysis of Water and Wastes," EPA-600/4-79-020, March, 1983, and the Federal Register, October 26, 1983. Arsenic, selenium and thallium are determined by furnace AA; silver, aluminum, barium, beryllium, boron, cadmium, calcium, chromium, copper, cobalt, iron, magnesium, manganese, molybdenum, nickel, lead, sodium, antimony, tin, titanium, vanadium, and zinc are determined by ICP emission spectrometry, except where lower levels of detection are required; in these cases (e.g. lead in groundwater monitoring samples) furnace AA is used. All furnace AA parameters are run by method of standard additions. The determination of mercury is performed by cold vapor AA.

### **EP TOXICITY**

The determination of metals in aqueous EP Toxicity leachates is performed according to the methods published by EPA in "Test Methods for Evaluating Solid Waste" EPA SW-846, revised April, 1984 and the Federal Register, Oct. 26, 1983, 1979. Silver, arsenic, barium, cadmium, chromium, lead and selenium are determined by ICP emission spectrometry. Mercury is determined using cold vapor AA. For leachates that are organic in nature, the analyses are performed according to the methods described under **OIL/SLUDGE** below.

### **SOIL/SEDIMENT**

The determination of silver, beryllium, cadmium, chromium, copper, nickel, antimony, lead, and zinc in sediment samples is performed according to methods published by EPA in "Interim Methods for the Sampling and Analysis of Priority Pollutants in Sediments and Fish Tissue", EPA 600/4-81-055, October 1980. Mercury is determined according to the sediment method published by EPA in "Method for Chemical Analysis of Water and Wastes", EPA 600/4-79-020, March 1983. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion.

### **OIL/SLUDGE**

The determination of silver, aluminum, boron, barium, beryllium, calcium, cadmium, copper, chromium, cobalt, iron, magnesium, manganese, molybdenum, sodium, nickel, lead, antimony, tin, titanium, vanadium, and zinc in sludge/petroleum-based samples is performed by ICP emission spectrometry using a magnesium nitrate dry ashing digestion technique. Arsenic, selenium and thallium are determined by furnace AA using nitric acid in a closed decomposition vessel for sample digestion. Mercury is determined by cold vapor AA using the same digestion technique.

## Summary of Quality Assurance/Quality Control Procedures (QA/QC)

ETC bases its quality assurance protocols on the following government guidelines:

- "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-330/9/79/001-R, October 1979;
- the recommended guidelines for EPA Methods 624 and 625. (Federal Register, December 3, 1979, updated on October 26, 1984);
- "Manual of Analytical Methods for the Analysis of Pesticides in Humans and Environmental Samples," EPA 600/8-80-038, June 1980;
- "Determination of 2,3,7,8-TCDD in Soil and Sediment" EPA, Region VII, Kansas City, September 1983;
- Organic Analysis: Multi-media, Multi Concentration-IFB WA84-A267; and
- Dioxin Analysis: Soil/Sediment Matrix; Multi-Concentration; Selected Ion Monitoring with Jar Extraction Procedure-IFB WA84-A002

However, we have modified our protocols to provide a higher level of QA/QC than the guidelines require. For example, we analyze a higher than required number of quality control samples and we pay especially careful attention to the certification of the "reference standard" compounds we use in analysis. Below are listed the key QA/QC elements for the methods we used.

### Analysis of Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry

- Each batch of 13 samples consists of 9 customer samples (at a maximum), one blank sample, one spiked blank, one spiked sample and one replicate sample. This amounts to a 30% quality control factor.
- Three surrogate compounds are added to each sample in the batch of 13.
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and retuned, if necessary, at the beginning of each day to ensure that its performance on bromofluorobenzene (BFB) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of Volatile Organic Priority Pollutant "standards" at a minimum of 3 different concentrations and using a mixture of 3 internal standards at a constant concentration.
- The calibration curve is verified with a mixture of priority pollutant standards every day. If the response factors vary greater than 25%, the instrument must be recalibrated.

### Analysis of Organic Compounds Extracted in Acid or Base/Neutral Solutions by Gas Chromatography/Mass Spectrometry

- Each batch of 20 samples consists of 16 customer samples (at a maximum), one blank sample, one spiked blank (for water matrices), one sample spiked with the priority pollutant standard mixture and a duplicate customer sample. This amounts to a 20% quality control factor.

- Three surrogate compounds are added to each sample in the batch for Base/Neutral analysis.
- Three surrogate compounds are added to each sample in the batch for Acid analysis
- A blind quality control sample is introduced to the laboratory for analysis on a weekly basis.
- Each GC/MS is checked and retuned, if necessary, at the beginning of each day to ensure that its performance on decafluorotriphenylphosphine (DFTPP) meets the EPA criteria.
- A calibration curve for quantitation is prepared using a mixture of standards composed of either the Organic Acid or Base/Neutral Extractable Compounds at a minimum of 3 concentrations and using five internal standards for quantitation.

### Analysis of Metals

#### All Samples

- New standards are prepared for each batch of samples
- Normal calibration is performed using a blank sample and four standards that have been through the sample preparation procedure. A regression analysis is used to construct the calibration curve.
- All EP Toxicity samples and all samples determined by furnace atomic absorption are calculated by the "method of additions".
- For each sample analysis that requires the use of the "method of additions" technique, a three point calibration is performed using U.S. EPA "Methods for Chemical Analysis of Water and Wastes, 1979". Results are obtained using linear regression analysis. Any regression with a coefficient of correlation below 0.990 is considered suspect, necessitating review of calibration data or sample re-analysis.
- In constructing the normal calibration curves the lowest concentration levels we use are values greater than or equal to 5 times the Instrumental Detection Limit (IDL).
- All calibration standards are analyzed in duplicate, at a minimum.
- Independent reference standards are used to check the accuracy of calibration standards.
- A check standard is analyzed every ten samples to validate the normal calibration curve.
- One customer sample out of every ten is analyzed in triplicate.

#### Homogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are homogeneous, the QC program is a minimum of 25% and consists of analyzing

- 3 sets of triplicate analyses,
- 2 Replicate spikes;
- 1 independent reference standard;
- 4 Calibration standards (processed using the sample preparation method);
- 4 Calibration standards (without sample preparation); and
- 1 Reagent Blank

Heterogeneous Samples (except for Mercury analysis)

Samples are analyzed in batches of 30 or less. For batches in which the sample matrices are heterogeneous, the QC program is a minimum of 35% and consists of analyzing:

- 3 sets of triplicate analyses;
- 2 Replicate spikes;
- 1 Replicate independent reference standards;
- 4 Calibration standards (processed using the sample preparation method);
- 1 Procedural Blank;
- 4 Calibration standards (without sample preparation); and
- 1 Reagent Blank.

Analysis of Mercury

To analyze samples for mercury we group them by matrix in batches of 30 or less. Our QC program is a minimum of 30% and consists of analyzing:

- each of the 30 customer samples in duplicate;
- 3 sets of triplicate analyses;
- 2 Replicate spikes,
- 2 Replicate independent reference standards;
- 10 Calibration standards (processed using the sample preparation method); and
- 2 Procedural Blanks

Analysis of Pesticides, Herbicides and PCB's by Gas Chromatography

Pesticide, herbicide and PCB samples are grouped in batches of 16 customer samples or less according to the type of analysis to be performed. The QC program for each of these three types of analyses is a minimum of 20% and consists of analyzing:

- 1 procedural blank sample (a reagent blank is analyzed in the case of non-water matrices);
- 1 spiked blank sample (the spiked blank is eliminated in the case of non-water matrices);
- 1 replicate sample;
- 1 replicate spiked sample; and
- 1 known reference QC sample for at least each 100 samples analyzed.

The instrument is calibrated each run with three standards, and checked every 10 samples.

Analysis of Cyanides, Phenols, Fluoride, Chloride, Nitrate and Nitrite

- All parameters are analyzed using a Technicon Autoanalyzer II GT.
- 3 calibration standards are analyzed at the beginning and end of each batch.

- Each batch (up to 80 samples) consists of analyzing one blank, one spiked blank, one duplicate and spiked sample every 20 samples, and an EPA known reference sample.

#### Analysis of Total Organic Carbon (TOC)

TOC samples are analyzed on a daily basis with the number of samples analyzed per day dependent on the request for duplicate or quadruplicate analyses. The quality control program is designed to maintain the appropriate amount of QC and consists of the following elements.

- Daily instrument calibration
- One blank
- Standard recalibration every 10 samples
- Spiked samples at a low and high level
- Every sample is run in duplicate at a minimum

#### Analysis of Total Organic Halide (TOX)

- Blank reagent water for absolute carbon background must contain less than 5 ug/l of halide (as chloride).
- Using a trichlorophenol standard, the mean adsorption efficiency must be within +/- 15% of the standard value.
- Calibration standards are run every 10 samples.
- Every sample is run in duplicate at a minimum.

#### Analysis of 2,3,7,8-TCDD (Dioxin) by GC/MS (SIM)

- Each sample is dosed with a known quantity of  $^{13}\text{C}_{12}$ -2,3,7,8-TCDD as internal standard and  $^{37}\text{Cl}_4$ -TCDD as surrogate standard. The action limits for surrogate standard results is +/- 40% of the true value. Samples showing surrogate standard results outside of these limits are reextracted and reanalyzed.
- Two laboratory "method blanks" are run along with each set of 24 or fewer samples. The method blank is also dosed with the internal standard and surrogate standard.
- At least one method blank of 24 samples is run in duplicate to determine intra-laboratory precision.
- Qualitative Requirements. The following are met in order to confirm the presence of native 2,3,7,8-TCDD.
  - a. Isomer specificity must be demonstrated initially and verified once per 8-hour work shift. The verification consists of injecting a mixture containing TCDD isomers which elute close to 2,3,7,8-TCDD. The 2,3,7,8-TCDD must be separated from interfering isomers, with no more than 25% valley relative to the 2,3,7,8-TCDD peak.
  - b. The 320/322 ratio is within the range of 0.67 to 0.87.
  - c. Ions 320, 322, and 257 are all present and maximize together the signal to mean noise ratio must be 2.5 to 1 or better for all 3 ions.
  - d. The retention time is equal (within 3 seconds) the retention time for the isotopically labeled 2,3,7,8-TCDD.
  - e. At least one of the positives can be confirmed by obtaining partial scan spectra from mass 150 to mass 350. The partial scan guidelines are as follows:



- . the 320/324 ratio should be 1.58 +/- 0.16
- . the 257/259 ratio should be 1.03 +/- 0.10
- . the 194/196 ratio should be 1.54 +/- 0.15
- One sample is spiked with native 2,3,7,8-TCDD at a level of 1.0 PPB (for soil) for each set of 24 or fewer samples.
- In cases where no native 2,3,7,8-TCDD is detected, the actual detection limit is estimated and reported based on a signal to noise ratio of 2.5 to 1 at ions 320 and 322.
- For each sample, the internal standard is present with at least a 10 to 1 signal to noise ratio for both mass 332 and mass 334. Also, the internal standard 332/334 ratio must be within the range of 0.67 to 0.87.

#### Subcontractor QA/QC

Each subcontractor is required to maintain an appropriate level of quality control. To insure this, each subcontractor is required to submit to ETC the quality control data for all analyses it performs. This data is kept on file at ETC. In general, the amount of quality control required is one duplicate sample with one spiked sample for every ten analyses.

#### Chain-of-Custody

The chain-of-custody procedure is part of our quality assurance protocol. We believe our chain-of-custody record fully complies with the legal requirements of federal, state and local government agencies and of the courts of law. The record covers:

- labeling of sample bottles, packing the Sample Shuttle and transferring the Shuttle under seal to the custody of a shipper;
- outgoing shipping manifests;
- the chain-of-custody form completed by the person(s) breaking the Shuttle seal, taking the sample, resealing the Shuttle and transferring custody to a shipper.
- incoming shipping manifests;
- breaking the Shuttle's reseal.
- storing each labeled sample bottle in a secured area;
- disposition of each sample to an analyst or technician; and
- the use of the sample in each bottle in a testing procedure appropriate to the intended purpose of the sample.

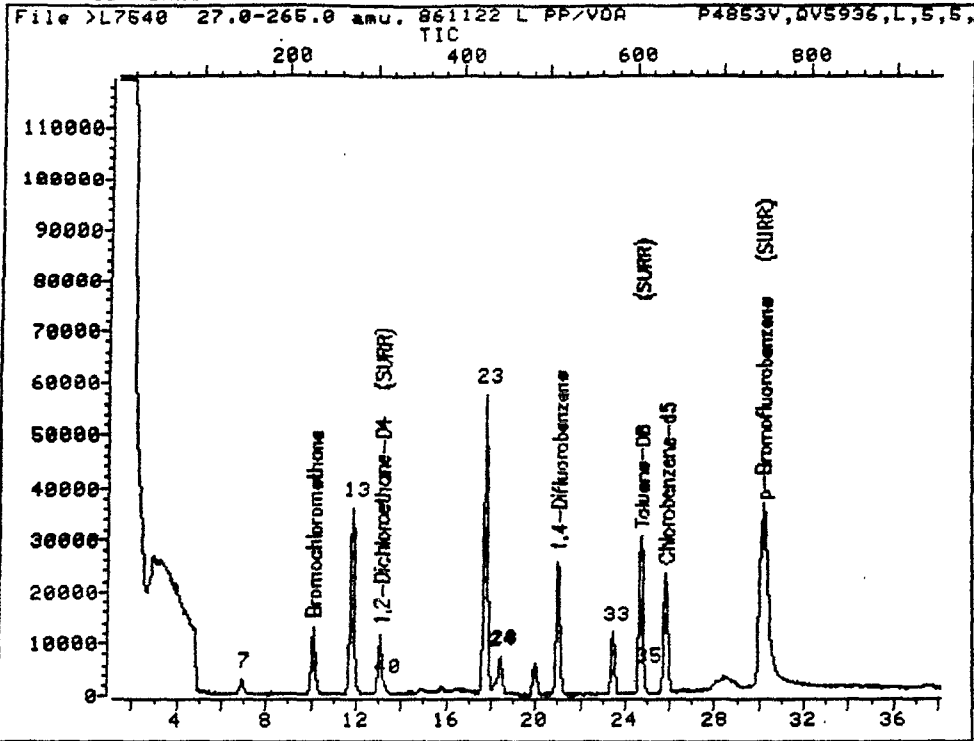
The records show for each link in this process:

- the person with custody; and
- the time and date each person accepted or relinquished custody.

**Appendix A**  
**Mass Spectral Data**  
**for**  
**Quantitated Compounds**

1. A total ion chromatogram for each sample analyzed by a GC/MS instrument.
2. A Quant. report used by the analyst to determine qualitative and quantitative results of the compounds present.
3. A mass spectrum and a reference spectrum for each priority pollutant compound detected in the sample.

TOTAL ION CHROMATOGRAM



Data File: >L7540::U2  
Name: 861122 L PP/VOA  
Misc: P4853U, QV5936, L, 5, 5,

Quant Output File: ^L7540::AQ

Id File: LVQA::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861123 05:33

Operator ID: JQ6275  
Quant Time: 861123 09:51  
Injected at: 861123 09:12

QUANT REPORT

Operator ID: JQ6275                      Quant Rev: 6            Quant Time: 861123 09:51  
 Output File: ^L7540::AQ                Injected at: 861123 09:12  
 Data File: >L7540::U2                 Dilution Factor: 1.00000  
 Name: 861122 L PP/VOA  
 Misc: P4853U,QU5936,L,5,5,

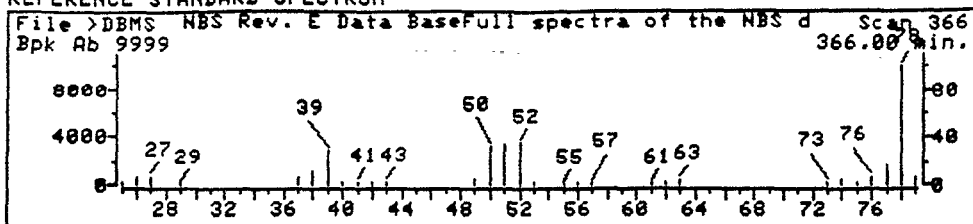
ID File: LVOA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861123 05:33

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *Bromochloromethane	10.09	231	27080	250.00	NG	92
7) Methylene chloride	6.94	150	6462	<del>107.81</del>	NG <del>94.5</del>	94
13) 1,2-Trans-dichloroethylene	11.84	276	148354	750.16	NG	99
16) *1,4-Difluorobenzene	21.01	512	123058	250.00	NG	94
17) 1,2-Dichloroethane-D4 (SURR)	13.08	308	24542	242.96	NG	82
23) Trichloroethylene	17.75	428	114888	512.54	NG	99
24) Benzene	18.41	445	7959	67.59	NG	83
<del>28) bis(Chloromethyl)ether</del>	<del>18.45</del>	<del>446</del>	<del>2571</del>	<del>41.76</del>	<del>NG</del>	<del>100</del>
30) *Chlorobenzene-d5	25.83	636	94499	250.00	NG	84
33) Tetrachloroethylene	23.46	575	23637	98.31	NG	89
34) Toluene-D8 (SURR)	24.71	607	140880	244.68	NG	87
<del>35) Toluene</del>	<del>24.90</del>	<del>612</del>	<del>2004</del>	<del>3.35</del>	<del>NG</del>	<del>87</del>
38) p-Bromofluorobenzene (SURR)	30.16	747	85124	240.04	NG	
40) Methyl ethyl ketone	13.31	314	11187	67.15	NG	

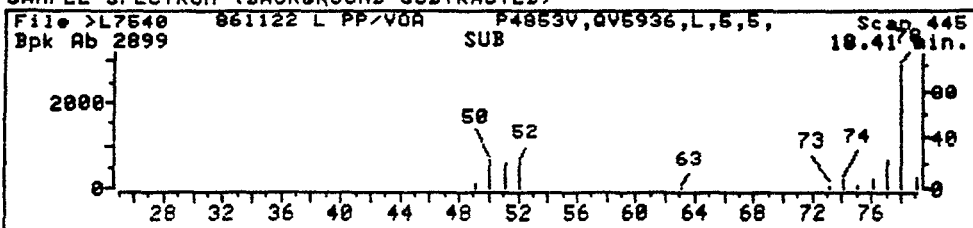
\* Compound is ISTD

src 11/26/86

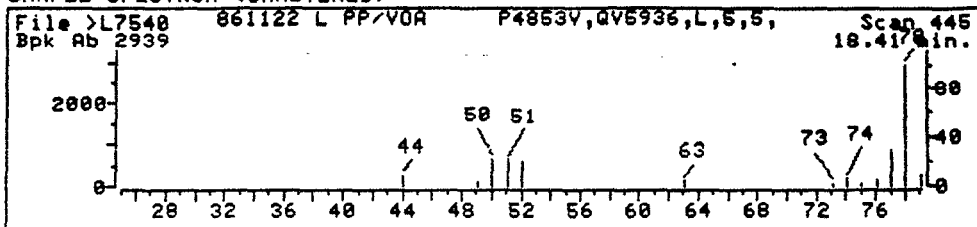
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



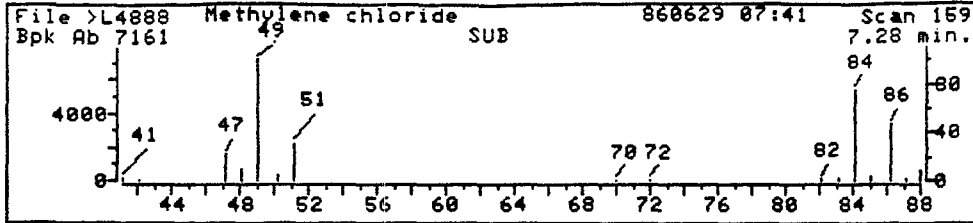
Data File: >L7540::U2  
Name: 861122 L PP/VOA  
Misc: P4853V,QU5936,L,5,5,  
Quant Time: 861123 09:51  
Injected at: 861123 09:12

Quant Output File: ^L7540::AQ

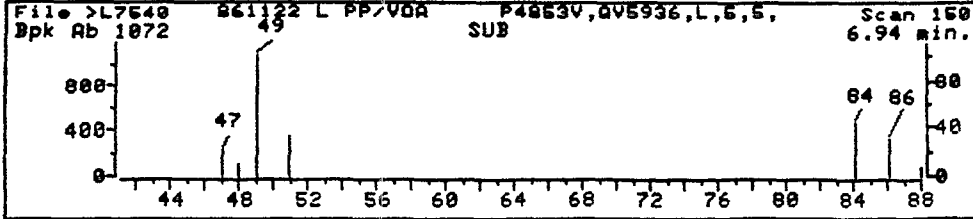
Quant ID File: LVQA::US  
Last Calibration: 861123 05:33

Compound No: 24  
Compound Name: Benzene  
Scan Number: 445  
Retention Time: 18.41 min.  
Quant Ion: 50.0  
Area: 7959  
Concentration: 67.59 NG  
q-value: 83

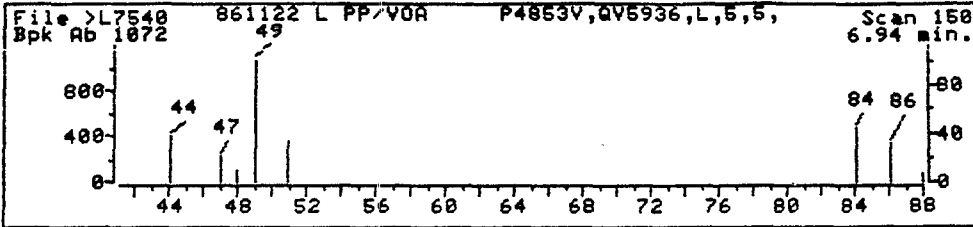
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >L7540::U2  
 Name: 861122 L PP/VOA  
 Misc: P4853V,QV5936,L,5,5,  
 Quant Time: 861123 09:51  
 Injected at: 861123 09:12

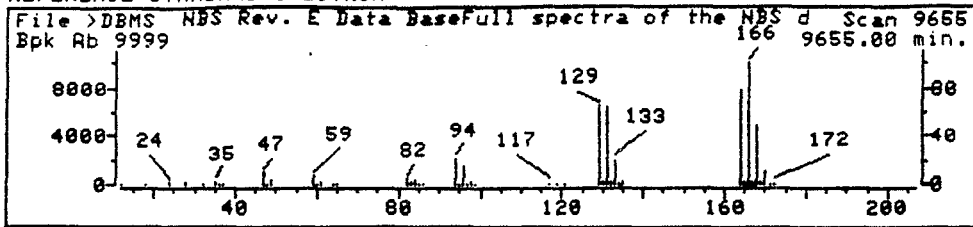
Quant Output File: ^L7540::AQ

Quant ID File: LVOA::US  
 Last Calibration: 861123 05:33

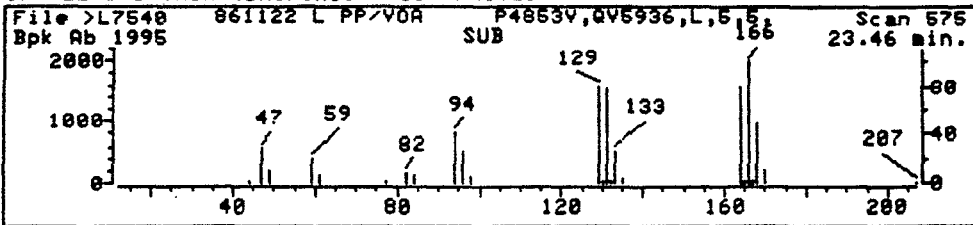
Compound No: 7  
 Compound Name: Methylene chloride  
 Scan Number: 150  
 Retention Time: 6.94 min.  
 Quant Ion: 84.0  
 Area: 6462  
 Concentration: 107.81 NG  
 q-value: 94

305292

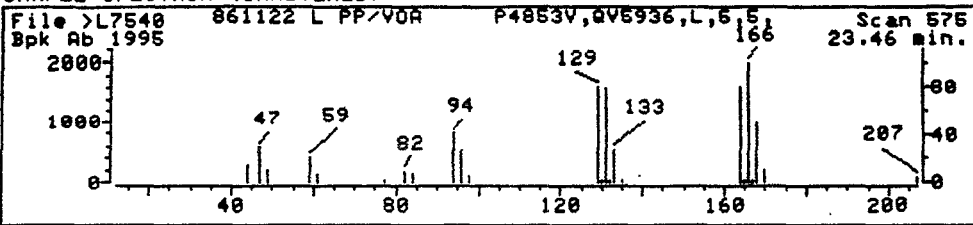
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

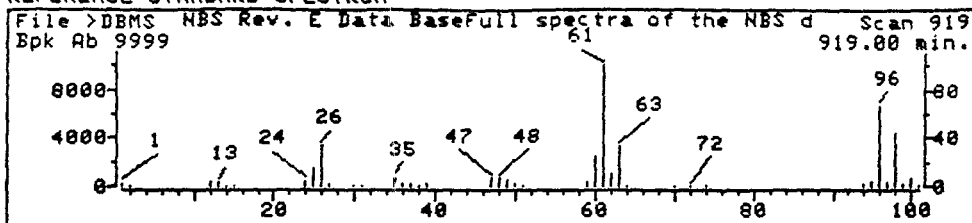


Data File: >L7540::U2  
Name: 861122 L PP/VDA  
Misc: P4853V,QU5936,L,5,5,  
Quant Time: 861123 09:51  
Injected at: 861123 09:12

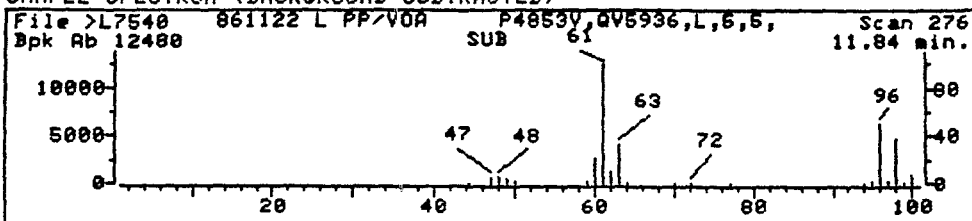
Quant Output File: ^L7540::AQ  
Quant ID File: LVDA::US  
Last Calibration: 861123 05:33

Compound No: 33  
Compound Name: Tetrachloroethylene  
Scan Number: 575  
Retention Time: 23.46 min.  
Quant Ion: 166.0  
Area: 23637  
Concentration: 98.31 NG  
q-value: 89

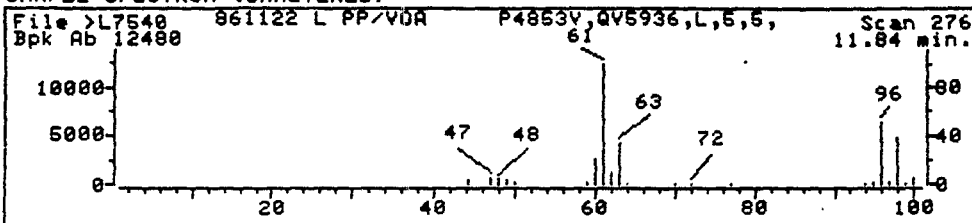
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >L7540::U2  
Name: 861122 L PP/VOA  
Misc: P4853U,QU5936,L,5,5,  
Quant Time: 861123 09:51  
Injected at: 861123 09:12

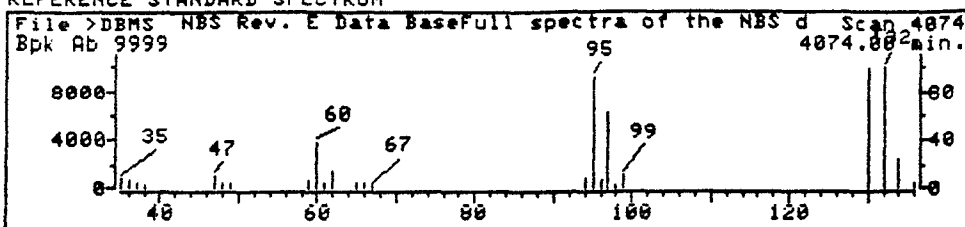
Quant Output File: ^L7540::AQ  
Quant ID File: LVDA::US  
Last Calibration: 861123 05:33

Compound No: 13  
Compound Name: 1,2-Trans-dichloroethylene  
Scan Number: 276  
Retention Time: 11.84 min.  
Quant Ion: 61.0  
Area: 148354  
Concentration: 750.16 NG  
q-value: 99

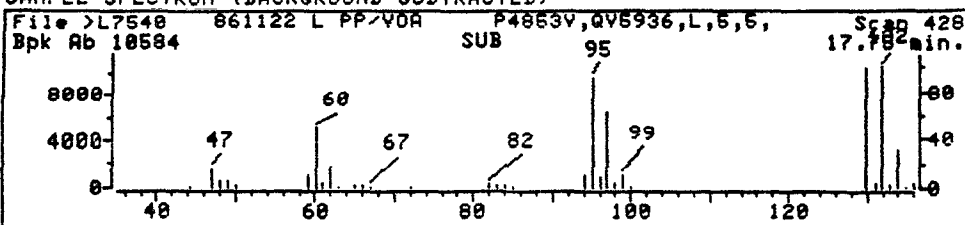
305294



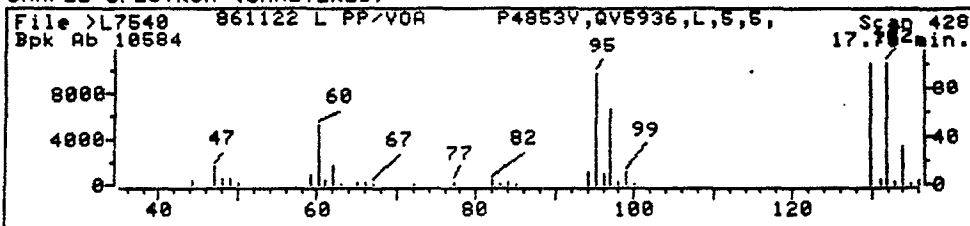
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >L7540::U2  
Name: 861122 L PP/VDA  
Misc: P4853U,QU5936,L,5,5,  
Quant Time: 861123 09:51  
Injected at: 861123 09:12

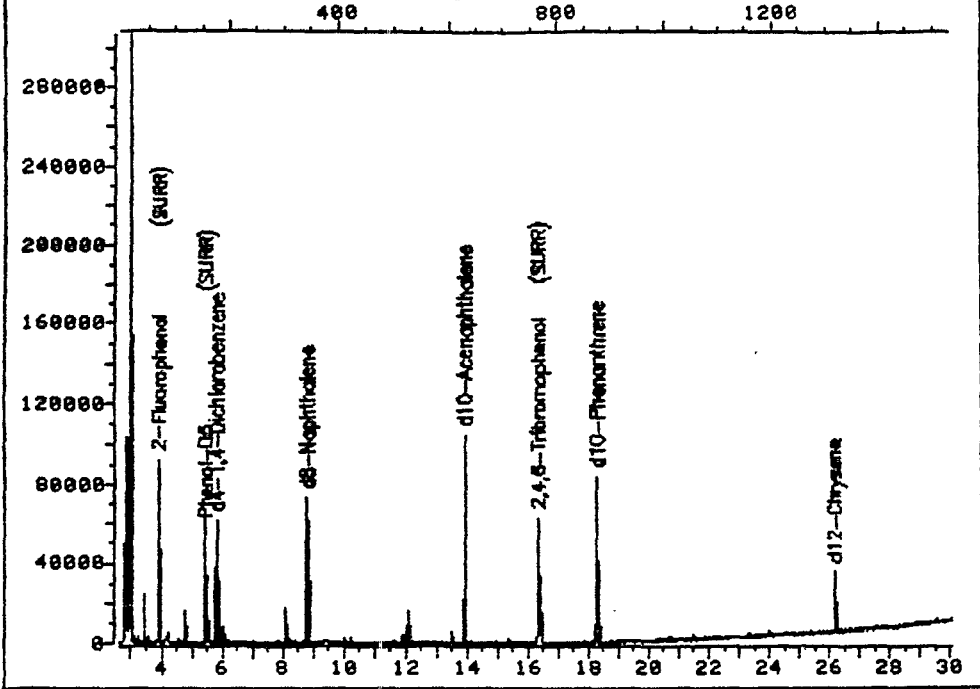
Quant Output File: ^L7540::AQ

Quant ID File: LVDA::US  
Last Calibration: 861123 05:33

Compound No: 23  
Compound Name: Trichloroethylene  
Scan Number: 428  
Retention Time: 17.75 min.  
Quant Ion: 95.0  
Area: 114888  
Concentration: 512.54 NG  
q-value: 99

TOTAL ION CHROMATOGRAM

File >F0534 45.0-450.0 amu. PP/ACID,#F P4853A ,QA6013,L, 9  
TIC



Data File: >F0534::U3  
Name: PP/ACID,#F  
Misc: P4853A ,QA6013,L, 930,1

Quant Output File: ^F0534::AQ

BTL#10

Id File: FACID::US  
Title: PP/ACID IDFILE  
Last Calibration: 861126 11:03

Operator ID: JA4996  
Quant Time: 861126 18:27  
Injected at: 861126 17:54

305296

QUANT REPORT

Operator ID: JA4996  
 Output File: ^F0534::AQ  
 Data File: >F0534::U3  
 Name: PP/ACID,#F  
 Misc: P4853A ,QA6013,L, 930,1

Quant Rev: 6      Quant Time: 861126 18:27  
 Injected at: 861126 17:54  
 Dilution Factor: 1.00000

BTL#10

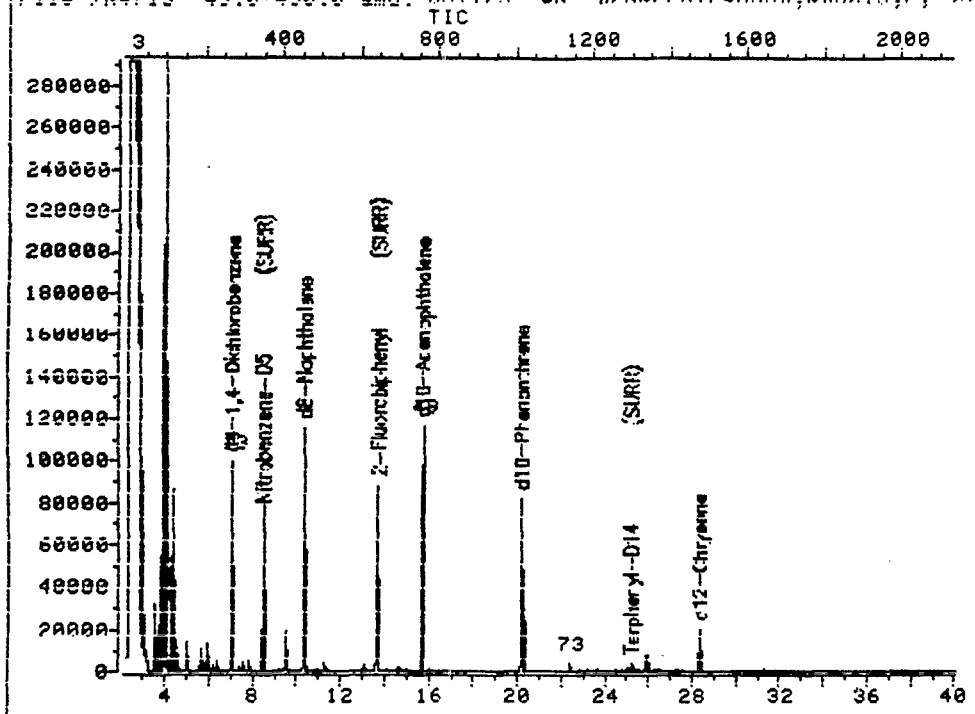
ID File: FACID::US  
 Title: PP/ACID IDFILE  
 Last Calibration: 861126 11:03

	Compound		R.T.	Scan#	Area	Conc	Units	q
1)	*d4-1,4-Dichlorobenzene		5.78	175	28442	40.00	UG/ML	95
3)	2-Fluorophenol	(SURR)	3.89	69	44528	87.45	UG/ML	80
9)	Phenol-D5	(SURR)	5.42	155	68076	87.47	UG/ML	91
12)	*d8-Naphthalene		8.77	342	77573	40.00	UG/ML	97
18)	*d10-Acenaphthalene		13.90	629	52966	40.00	UG/ML	96
25)	*d10-Phenanthrene		18.27	873	81641	40.00	UG/ML	99
26)	2,4,6-Tribromophenol	(SURR)	16.34	765	22297	72.28	UG/ML	99
29)	*d12-Chrysene		26.18	1315	30010	40.00	UG/ML	100

\* Compound is ISTD

TOTAL ION CHROMATOGRAM

File >N4715 45.0-450.0 amu. 861125 #N B/N&PESTP4853R,DR6013,1, 930



Data File: >N4/15::U4  
 Name: 861125 #N B/N&PEST  
 Misc: P4853B,UB6013,L, 930,1

Quant Output File: ^N4/15::AQ

BIL#42

Id File: NBNP::US  
 Title: B/N+PEST ID FILE  
 Last Calibration: 861126 12:44

Operator ID: KUU/86  
 Quant Time: 861127 08:22  
 Injected at: 861127 04:47

305298

QUANT REPORT

Operator ID: K00786  
 Output File: ^N4715::AQ  
 Data File: >N4715::U4  
 Name: 861125 #N B/N&PEST  
 Misc: P4853B,Q86U13,L, 93U,1

Quant Rev: 6 Quant Time: 861127 08:22  
 Injected at: 861127 04:47  
 Dilution Factor: 1.00000

BTL#42

ID File: NBNP::US  
 Title: B/N+PEST ID FILE  
 Last Calibration: 861126 12:44

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *d4-1,4-Dichlorobenzene	6.99	268	30663	40.00	UG/ML	95
3) N-Nitrosodimethylamine	2.64	23	8824	<del>11.07</del>	<del>UG/ML</del>	100
12) Nitrobenzene-D5 (SURR)	8.39	347	80015	44.07	UG/ML	80
13) bis(2-Chloroisopropyl)ether	6.99	268	1714	<del>7.14</del>	<del>UG/ML</del>	92
21) *d8-Naphthalene	10.29	454	148418	40.00	UG/ML	98
22) 2-Fluorobiphenyl (SURR)	13.64	643	79883	29.47	UG/ML	97
42) *d10-Acenaphthalene	15.64	756	76092	40.00	UG/ML	98
49) Dimethyl phthalate	15.64	756	20090	<del>8.01</del>	<del>UG/ML</del>	97
66) *d10-Phenanthrene	20.14	1010	97172	40.00	UG/ML	97
73) Di-n-butyl phthalate	22.34	1134	6344	<del>2.01</del>	<del>UG/ML</del>	93
85) *d12-Chrysene	28.34	1472	24567	40.00	UG/ML	100
98) Terphenyl-D14 (SURR)	25.20	1295	4444	6.07	UG/ML	93

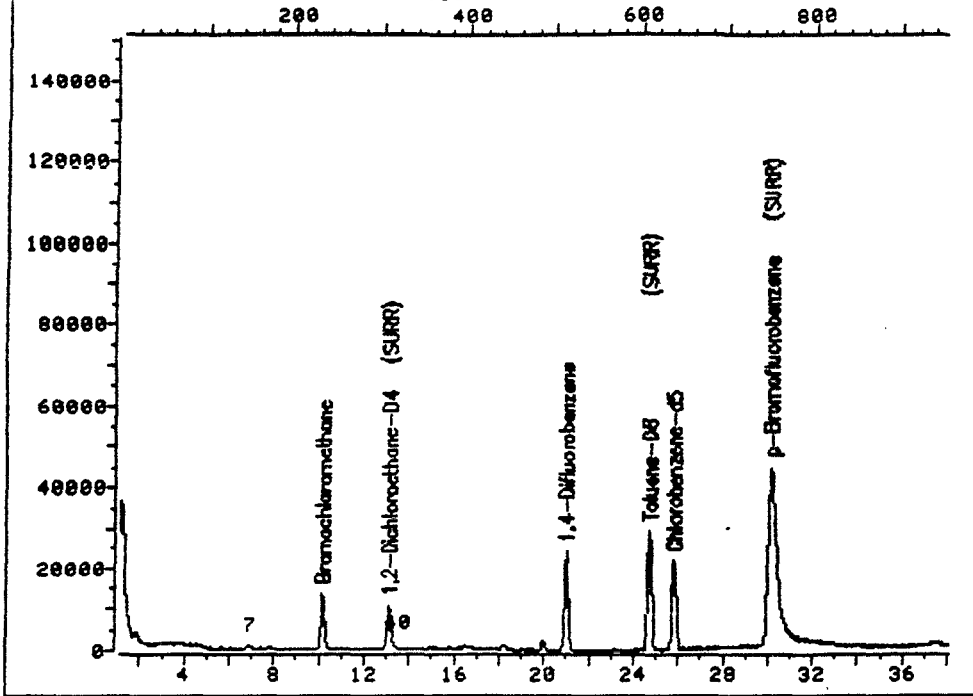
\* Compound is ISID

AD  
 11/9/88

**Appendix C1**  
**GC/MS Subsidiary Data**

TOTAL ION CHROMATOGRAM

File >L7526 27.0-265.0 amu. 861122 L PP/VDA QC5936V, QV5936, L, 5, 5  
TIC



Data File: >L7526::U1  
Name: 861122 L PP/VDA  
Misc: QC5936V, QV5936, L, 5, 5,

Quant Output File: ^L7526::AQ

Id File: LVDA::US  
Title: Volatile Priority Pollutant ID FILE  
Last Calibration: 861123 05:33

Operator ID: JQ6275  
Quant Time: 861123 05:56  
Injected at: 861122 21:45

305301

QUANT REPORT

Operator ID: JQ6275                      Quant Rev: 6      Quant Time: 861123 05:56  
 Output File: ^L7526::AQ                      Injected at: 861122 21:45  
 Data File: >L7526::U1                      Dilution Factor: 1.00000  
 Name: 861122 L PP/UDA  
 Misc: QC5936U,QU5936,L,5,5,

ID File: LUDA::US  
 Title: Volatile Priority Pollutant ID FILE  
 Last Calibration: 861123 05:33

	Compound	R.T.	Scan#	Area	Conc	Units	q
1)	*Bromochloromethane	10.16	233	27932	250.00	NG	99
7)	Methylene chloride	6.90	149	2359	38.16	NG	97
16)	*1,4-Difluorobenzene	21.01	512	115376	250.00	NG	96
17)	1,2-Dichloroethane-D4 (SURR)	13.12	309	22560	238.21	NG	82
30)	*Chlorobenzene-d5	25.83	636	90031	250.00	NG	84
34)	Toluene-D8 (SURR)	24.75	608	135260	246.57	NG	89
38)	p-Bromofluorobenzene (SURR)	30.20	748	86108	254.86	NG	97
40)	<del>Methyl ethyl ketone</del>	<del>13.35</del>	<del>315</del>	<del>11088</del>	<del>74.48</del>	<del>NG</del>	<del>92</del>

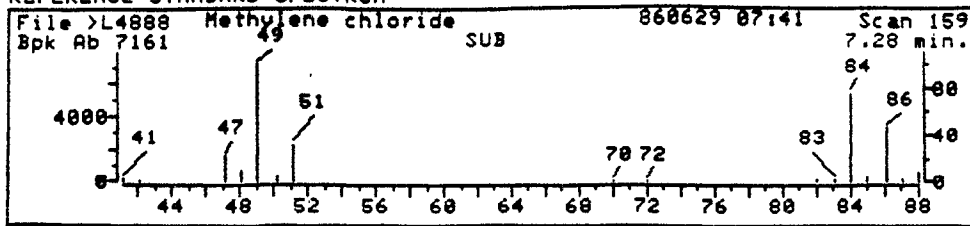
\* Compound is ISTD

SFC 11125184

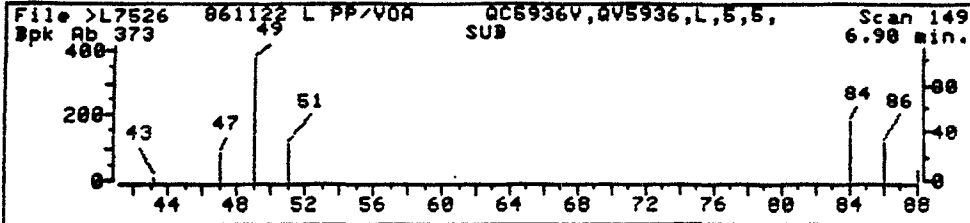
305302



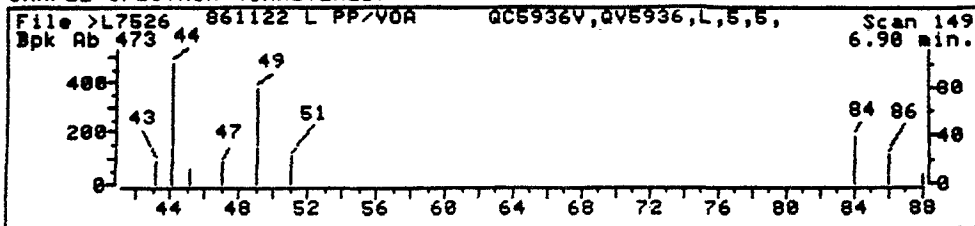
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



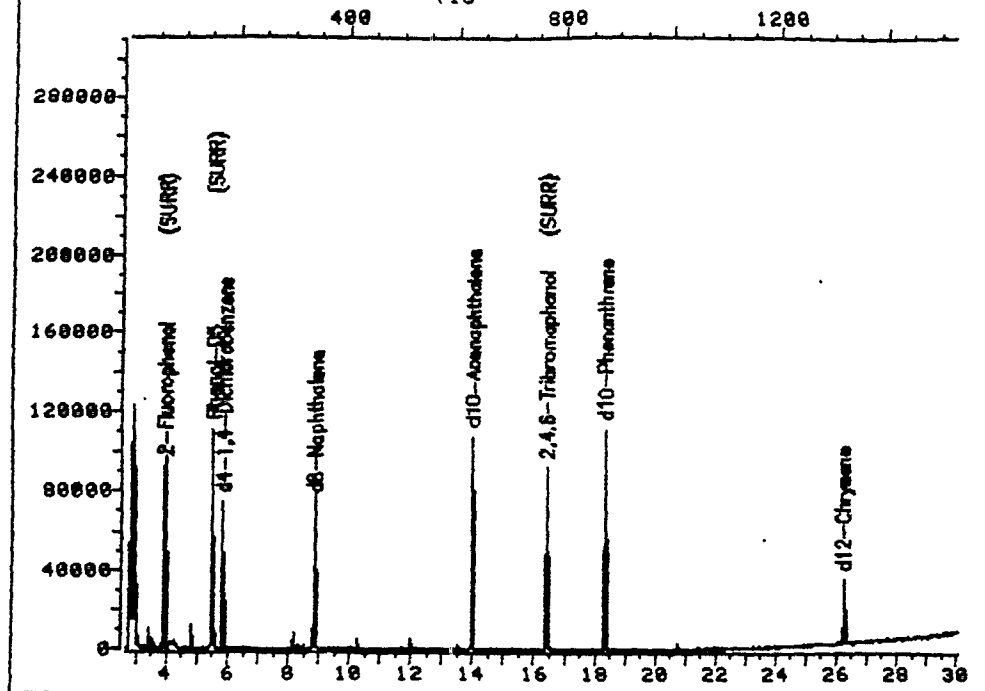
Data File: >L7526::U1  
 Name: 861122 L PP/VOA  
 Misc: QC5936V,QU5936,L,5,5,  
 Quant Time: 861123 05:56  
 Injected at: 861122 21:45

Quant Output File: ^L7526::AQ  
 Quant ID File: LVOA::US  
 Last Calibration: 861123 05:33

Compound No: 7  
 Compound Name: Methylene chloride  
 Scan Number: 149  
 Retention Time: 6.90 min.  
 Quant Ion: 84.0  
 Area: 2359  
 Concentration: 38.16 NG  
 q-value: 97

TOTAL ION CHROMATOGRAM

File >F0526 45.8-458.8 amu. PP/ACID,#F QC6013A ,QA6013,L,10  
TIC



Data File: >F0526::U3  
Name: PP/ACID,#F  
Misc: QC6013A ,QA6013,L,1000,1

Quant Output File: ^F0526::AQ

BTL# 2

Id File: FACID::US  
Title: PP/ACID IDFILE  
Last Calibration: 861126 11:03

Operator ID: JA4996  
Quant Time: 861126 13:17  
Injected at: 861126 12:45

305304

QUANT REPORT

Operator ID: JA4996  
 Output File: ^F0526::AQ  
 Data File: >F0526::U3  
 Name: PP/ACID,#F  
 Misc: QC6013A ,QA6013,L,1000,1

Quant Rev: 6      Quant Time: 861126 13:17  
 Injected at: 861126 12:45  
 Dilution Factor: 1.00000

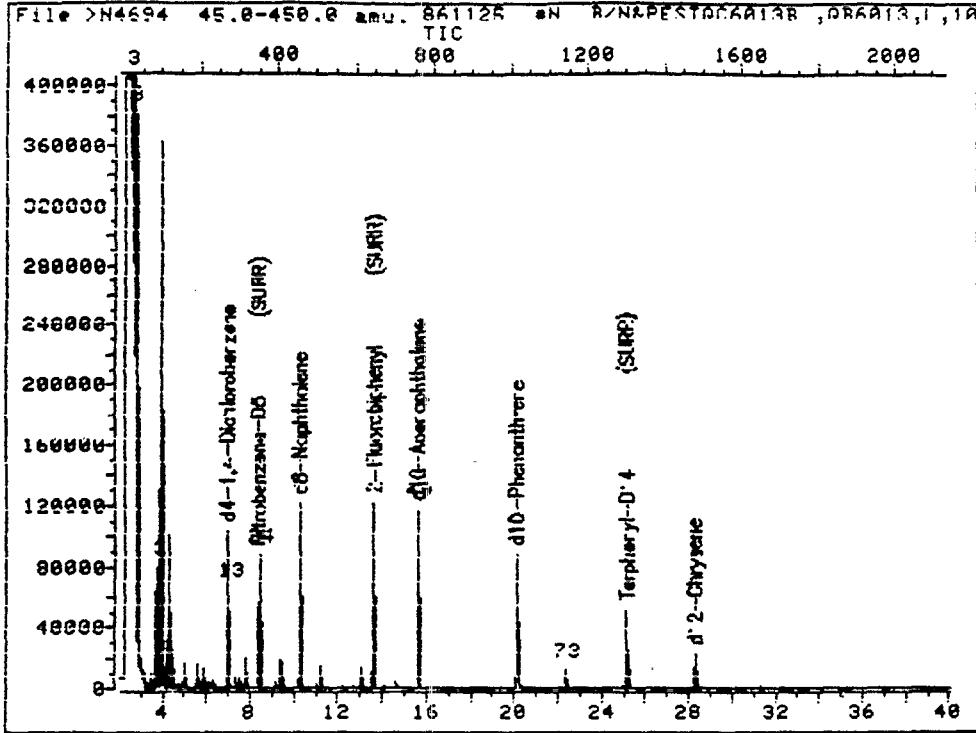
BTL# 2

ID File: FACID::US  
 Title: PP/ACID IDFILE  
 Last Calibration: 861126 11:03

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *d4-1,4-Dichlorobenzene	5.81	171	35658	40.00	UG/ML	89
3) 2-Fluorophenol (SURRE)	3.88	63	69006	108.10	UG/ML	76
9) Phenol-D5 (SURRE)	5.47	152	85528	87.66	UG/ML	96
12) *d8-Naphthalene	8.83	340	86861	40.00	UG/ML	98
18) *d10-Acenaphthalene	13.97	627	56444	40.00	UG/ML	97
25) *d10-Phenanthrene	18.33	871	102168	40.00	UG/ML	99
26) 2,4,6-Tribromophenol (SURRE)	16.40	763	32589	84.42	UG/ML	97
29) *d12-Chrysene	26.24	1313	29616	40.00	UG/ML	100

\* Compound is ISTD

TOTAL ION CHROMATOGRAM



Data File: >N4694::U4  
Name: 861126 #N B/N&PEST  
Misc: QB6013B ,QB6013,L,1000,1

Quant Output File: ^N4694::AQ

BIL#21

Id File: NBNP::US  
Title: B/N+PEST ID FILE  
Last Calibration: 861126 12:44

Operator ID: KUU/B6  
Quant time: 861126 16:26  
Injected at: 861126 11:25

305306

QUANT REPORT

Operator ID: K00786      Quant Rev: 6      Quant Time: 861126 16:26  
 Output File: ^N4694::AQ      Injected at: 861126 11:25  
 Data File: >N4694::U4      Dilution Factor: 1.00000  
 Name: 861125 #N B/N&PEST  
 Misc: QC6013B ,QB6013,L,1000,1      BIL#21

ID File: NBNP::US  
 Title: B/N&PEST ID FILE  
 Last Calibration: 861126 12:44

Compound	R.I.	Scan#	Area	Conc	Units	q
1) *d4-1,4-Dichlorobenzene	6.95	267	34441	40.00	UG/ML	98
3) N-Nitrosodimethylamine	2.66	25	10063	<del>11.24</del>	<del>UG/ML</del>	100
3) N-Nitrosodimethylamine	2.76	31	1340	<del>1.50</del>	<del>UG/ML</del>	100
3) N-Nitrosodimethylamine	3.78	88	2107	<del>2.35</del>	<del>UG/ML</del>	100
12) Nitrobenzene-D5 (SURR)	8.35	346	84430	41.40	UG/ML	83
13) bis(2-Chloroisopropyl)ether	6.97	268	1836	<del>6.80</del>	<del>UG/ML</del>	74
21) *d8-Naphthalene	10.25	453	144769	40.00	UG/ML	98
22) 2-Fluorobiphenyl (SURR)	13.63	643	107010	40.47	UG/ML	97
24) N-Nitrosodi-n-propylamine	8.35	346	12555	<del>18.77</del>	<del>UG/ML</del>	66
42) *d10-Acenaphthalene	15.61	755	73999	40.00	UG/ML	97
49) Dimethyl phthalate	15.61	755	18965	<del>7.78</del>	<del>UG/ML</del>	56
66) *d10-Phenanthrene	20.11	1009	102784	40.00	UG/ML	96
73) Di-n-butyl phthalate	22.31	1133	17098	5.13	UG/ML	97
85) *d12-Chrysene	28.30	1471	25929	40.00	UG/ML	100
98) Terphenyl-D14 (SURR)	25.18	1295	49419	63.93	UG/ML	94

\* Compound is ISTD

ND  
 11/29/83

**Appendix D**  
**Subcontractor's Data**

1) A copy of the originating subcontractor's report is included for all data not generated within ETC's laboratory.



## Appendix E

### Chain-of-Custody Forms

1. A field Chain-of-Custody form (CCT) is included for all samples shipped by ETC shuttle.
2. An in-house sample Chain-of-Custody form is included for all samples not shipped by ETC shuttle.
3. Any additional Chain-of-Custody material provided by a client or by a client's sampling agent is also included.
4. A subcontractor's Chain-of-Custody form is included for any analytical work not performed within ETC's laboratory.
5. Analysis and Extraction Custody forms are included for the period the sample was in ETC's possession.



**CHAIN OF CUSTODY FORM (CC1)**

Company: HELLERTOWN MANUFACTURING Attn.: MARK SCHULTZ

Facility/Site: C/O GUARD OFFICE-SILVER ROAD Phone: \_\_\_\_\_

Address: HELLERTOWN, PA. 18055

**SAMPLE IDENTIFICATION**

Facility: QHAMAIGN (Facility/Site Code) \_\_\_\_\_ (Optional Sample Point Descriptions)

Sample Point: W-1-C-S-P-4 (Source Code) 441096 (Your Sample Point ID) 1500 (Start Date) 1500 (Start Time) \_\_\_\_\_ (Elapsed Hours)

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	M	500 ML	HN03	METALS			/
1	PN	125 ML	H2S04	PHENOLS			/
1	B	125 ML	H2S04	NITRATE			/
1	CN	125 ML	NAOH	CYANIDE			/
1	A	500 ML	BAKED	SOLIDS, TD/SULFATE			/
1	D	500 ML	RINSE	FLUORIDE			/
2	U	40 ML	THIOL	VOA			/
3	E	1 L	BAKED	EXTRACTABLES			/
1	TB	40 ML	H2O	TRIP BLANK			/

**CHAIN OF CUSTODY CHRONICLE**

1. Shuttle Opened By: (print) M. SCHULTZ Date: 11-10-86 Time: \_\_\_\_\_  
Signature: Mark Schultz Seal #: \_\_\_\_\_ Intact: \_\_\_\_\_

2. I have received these materials in good condition from the above person.  
Name: \_\_\_\_\_ Signature: \_\_\_\_\_  
Date: \_\_\_\_\_ Time: \_\_\_\_\_ Remarks: \_\_\_\_\_

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

4. Shuttle Sealed By: (print) M. SCHULTZ Date: 11-10-86 Time: 16:30  
Signature: Mark Schultz Seal #: 0067756 Intact: \_\_\_\_\_

AB USE ONLY Opened By: Mark Schstadt Date: 11/11/86 Time: 5:10 PM  
SHUTTLE # 1318 TEMP. °C 12 SEAL # 67756 COND. INTACT

**305311**











Sample Number	Log Link	Sample Vol. (ml)	Extract Vol. (ml)		Comments
			BN	ACID	
P0447	17862	1000	1.0	1.0	ppb extract/prec
P0448		1000	1.0	1.0	
P0449		1000	1.0	1.0	
P0450		1000	1.0	1.0	
P0457		1000	1.0	1.0	
P0458		1000	1.0	1.0	
P0467		1000	1.0	1.0	
P0468		1000	1.0	1.0	
P0471		1000	1.0	1.0	
<del>P0472</del>	<del>995</del>	<del>1000</del>	<del>1.0</del>	<del>1.0</del>	<del></del>
P0489	17863	1000	1.0	1.0	pp/T
P0493		980	1.0	1.0	
P0494		1000	1.0	1.0	
P0495		1000	1.0	1.0	
P4852	17866	930	1.0	1.0	ppb extract/prec
P4853		930	1.0	1.0	
P0472	17862	1000	1.0	1.0	
QC 6013		1000	1.0	1.0	
QC 6013	S	1000	1.0	1.0	
P0447	S	980	1.0	1.0	
P0448	R	995	1.0	1.0	

QC Batch # 6013

Analysis \_\_\_\_\_  
SEE COMMENTS

Matrix H<sub>2</sub>O

Turnaround NORM

Date 11/17/86

**Extraction Method:**

Sep. Funnel K.B. 11-17-86  
BN 10:30 AM 11-17-86  
 Continuous OFF BN 11:30 AM 11-17-86  
ACID 1:50 PM 11/18/86  
 Soxhlet OFF: 1:20 PM 11/19/86

Other \_\_\_\_\_

**COMMENTS**

Ⓢ P0472 - Acid fraction lost during conc., repeated using a separate aliquot

FRACTION	SPIKE		
	Amt (ml)	Conc. (ug/ml)	Lot #
pp/BN	1.0	100	13,929
pp/ACID	1.0	100	13,829
Dioxibac-1260	1.0	100	14,404
pesticide	1.0	100/200	14,631
SEM-VDA			

SURROGATE		
Amt. (ml)	Conc. (ug/ml)	Lot #
1.0	BN's - 50	14,597
	ACID's - 100	

Set-up: K. Boyer 11-17-86 Aug. Dazarik 11-17-86  
 Conc.: 6/17/86 11-24-86

UPD/Supervisor: Ken - [Signature] 11-25-86  
 Spike/Surr. Verified: Kenneth Boyer 11-17-86







LAN  
11-30

DATE Nov 26 1986 SHIFT \_\_\_\_\_  
 FRACTION B/N  
 INSTRUMENT "N"  
 TUNE FILE \_\_\_\_\_  
 SEQUENCE FILE XV25  
 METHOD FILE BNPN-ACIDN  
 ID FILE N3UP  
 ANALYST(S) K. Valuk  
 SUPERVISOR Stapponich  
 BATCH #'s Q 6057, Q 6013

STANDARD	CONC PPM	LOT NO.	LOT VO
Test. IV	150	14433	
B/N II	200	14432	
B/N II	100	14713	
B/N I	60	14430	
Aniline	300	13468	
Aniline	100	13468	
Aniline	60	12470	
2-Picoline	300	14411	
2-Picoline	100	14412	
2-Picoline	60	14413	
Methoxy dln	300	14352	
Methoxy dln	100	14353	
Methoxy dln	60	14354	
Col. Test	400	14585	
Col Test Mix	25	14616	

(PLEASE INITIAL)

CURRENT CSOS STATUS	STANDARDS UPDATED
ACQ	DATE <u>Nov. 26 1986</u>
WIP	BY <u>KV</u>

NAME	DATA FILE	UL INJ	ALS #	DIL	TAPE #	SPECIALS (WRITE A-TYPE)	PLUS Y/N
22 Test-Mix	>N4673					OK 14:37	
Test. cal. sta. IV; 150ppm	>N4674		1				
B/N cal. sta. II; 200ppm	>N4675		2				
B/N cal. sta. II; 100ppm	>N4676		3				
B/N cal. sta. I; 60ppm	>N4677		4				
Aniline sta.; 300ppm	>N4678		5				
Aniline sta.; 100ppm	>N4679		6				
Aniline sta.; 60ppm	>N4680		7				
2-Picoline Methoxy dln sta.; 300ppm	>N4681		8				
2-Picoline Methoxy dln sta.; 100ppm	>N4682		9				
2-Picoline Methoxy dln sta.; 60ppm	>N4683		10			OK	
Col. Test-Mix	>N4684		11			OK 03:17	
QC 6057BS	>N4685		12			AN1260	
P 2712BS	>N4686		13			AN1260	
QC 6057B	>N4687		14			PCBSCN	S Y
P 2711B	>N4688		15				S Y
P 2711B2	>N4689		16				
P 2712B	>N4690		17				S Y
N 6321B	>N4691		18				S Y
QC 6013BS	>N4692		19			AN1260	
P 0447BS	>N4693		20			AN1260	
QC 6013B	>N4694		21			PCBSCN	
Col. Test-Mix	>N4695		22			OK 12:15	
B/N cal. sta. II; 100ppm	>N4696		23			AN1260; OK	
P 0447B	>N4697		24			PCBSCN	

305320



SUBCONTRACT  
REQUEST FOR ANALYSIS  
and SAMPLE CHAIN OF CUSTODY

LOGLINK: 17866

MATRIX: WATER

NAME OF SUBCONTRACT LAB: CHYUN

ETC JOB NUMBERS: P4850 P4853  
P4851  
P4852

TURNAROUND IN DAYS: NORMAL DATE DATA REQUIRED: 11/22  
(If deadline cannot be met, contact ETC Subcontract Group).

Send invoice, bill and reports to : ETC Subcontract Group  
(201)225-6786

Please perform the analyses requested below:

- |   |  |   |
|---|--|---|
| <input type="checkbox"/> Acidity                      | <input type="checkbox"/> Alkalinity          | <input type="checkbox"/> Ammonia (probe)          |
| <input type="checkbox"/> Ammonia (dist)               | <input type="checkbox"/> Bicarbonate         | <input type="checkbox"/> BOD (5 day)              |
| <input type="checkbox"/> Bromide                      | <input type="checkbox"/> Carbonate           | <input type="checkbox"/> Chloride                 |
| <input type="checkbox"/> Chem. Oxygen Dem.            | <input type="checkbox"/> Chromium+6          | <input type="checkbox"/> Fecal Coliform           |
| <input type="checkbox"/> Total Coliform               | <input type="checkbox"/> Color, apparent     | <input type="checkbox"/> Cyanide, total           |
| <input type="checkbox"/> Fluoride                     | <input type="checkbox"/> Formaldehyde (UV)   | <input type="checkbox"/> Hardness                 |
| <input type="checkbox"/> Nitrate (NO3)                | <input type="checkbox"/> Nitrite (NO2)       | <input type="checkbox"/> TKN                      |
| <input type="checkbox"/> Odor                         | <input type="checkbox"/> Oil & Grease (grav) | <input type="checkbox"/> T. Organic Carb.         |
| <input type="checkbox"/> Pet. Hydro (IR)              | <input type="checkbox"/> Pet. Hydro (grav)   | <input type="checkbox"/> Phenolics, tot.          |
| <input type="checkbox"/> Phenolics (5ug/l)            | <input type="checkbox"/> Phosphate (ortho)   | <input type="checkbox"/> Phosphate (total)        |
| <input type="checkbox"/> Phosphorus (tot.)            | <input type="checkbox"/> Silica (dissolved)  | <input checked="" type="checkbox"/> Sulfate (SO4) |
| <input type="checkbox"/> Sulfide (S)                  | <input type="checkbox"/> Sulfite (SO3)       | <input type="checkbox"/> Surfactant (MBAS)        |
| <input type="checkbox"/> Solids, Total                | <input type="checkbox"/> Solids, Tot. Diss.  | <input type="checkbox"/> Solids, Tot. Set.        |
| <input checked="" type="checkbox"/> Solids, Tot. Sus. | <input type="checkbox"/> Solids, Tot. Vol.   | <input type="checkbox"/> Turbidity                |
| <input type="checkbox"/> Gross Alpha, Beta            | <input type="checkbox"/> Radium 226          | <input type="checkbox"/> Radium 228               |

Others: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

Chain of Custody Section for Courier Pick up at ETC

Sample(s) Relinquished by: (ETC) Mark J. Eckardt  
Time: 4:30pm Date: 11/21/86  
Sample(s) Received by: Mark Kelly  
Time: 4:30pm Date: 11/21/86

Chain of Custody Section for Sending Sample from ETC

Sample Shuttle sealed by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Sample Shuttle opened by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_  
Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

Chain of Custody Section for Returning Sample to ETC

Sample Shuttle sealed by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_

Sample Shuttle opened by: \_\_\_\_\_  
Date: \_\_\_/\_\_\_/\_\_\_ Time: \_\_\_\_\_ Seal Number: \_\_\_\_\_  
Was the Seal intact? \_\_\_\_\_ Are the shuttle contents in good condition? \_\_\_\_\_

305322

ANALYZED ON: 11/26/86

TEST: Fluoride, UNITS: mg/L  
 METH. REF. # STD METHA 413E, MDL: 0.10  
 INSTRUMENT # A.A. CHANNEL 7

BY: CKERRAN  
 BOOK # CV-12 P. 108  
 FROM PAGE # 107

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	STANDARDS
17455	M0731	0.07	-	BMDL	17828	P2472	0.03	-	BMDL	PREP ON _____ BY _____ FROM _____ mg/L STOCK # _____
17455	M0703	0.26	10x	2.60	17828	P2473	0.03	-	BMDL	
17455	M0704	0.27	10x	2.70	17828	P2522	0.08	-	BMDL	
17455	M0705	0.23	10x	2.30	17828	P2536	0.03	-	BMDL	
17455	M0706	0.28	10x	2.80	17828	P2548	0.02	-	BMDL	
17455	M0707	0.23	10x	2.30	17806	P4850	0.16	-	0.16	
17455	M0708	0.03	-	BMDL	17806	P4851	0.21	-	0.21	STANDARD mg/L
17514	<del>P2532</del>	0.44	-	0.44	17806	P4852	2.24	-	RPT	
17514	<del>P2533</del>	0.54	-	0.54	17806	P4853	0.98	-	0.98	% of Theoretical
17749	P2500	0.01	-	BMDL	17250	N4520	1.64	-	1.64	SCOPE
17749	P2513	0.07	-	BMDL	17441	P2237	0.07	-	BMDL	
17749	P2514	0.03	-	BMDL	17878	P2776	0.087	-	BMDL	Corr. Coef.
17749	P2515	0.07	-	BMDL	17804	P2481	0.06	-	BMDL	<b>ADDITIONAL COMMENTS</b> RPT: IN 9319 - {over 7 mg/L P4852 - and excessive instrument noise Dilute 1:10 and check
17749	P2526	0.04	-	BMDL	17804	P2531	0.02	-	BMDL	
17778	P2497	0.03	-	BMDL	17804	P2541	0.04	-	BMDL	
17778	P2499	0.03	-	BMDL	17804	P2546	0.02	-	BMDL	
17778	P2507	0.13	-	0.13	17804	P2554	0.01	-	BMDL	
17778	P2517	0.12	-	0.12						
17748	P2469	0.03	-	BMDL						
17748	P2477	0.05	-	BMDL						
17748	P2524	0.08	-	BMDL						
17748	P2538	0.05	-	BMDL						
17748	P2539	0.04	-	BMDL						
17748	P2540	0.02	-	BMDL						

**Analyses QA/QC Data Report**  
 See CV-12, P. 107.

QC Batch # E1001 + FL003

Verified: [Signature] 11/28/86

METHOD BLANK		PRECISION	
Spiked Blank (Known)		Job # Dup (ETC)	P2472
Spiked Blank Result		Original Result	0.297
% Recovery		Duplicate Result	0.339
		A 7.4%	0.042
EXTERNAL REF. QC		ACCURACY	
Known Value		Job # Spiked (ETC)	P2472
Observed Result		Original Result	0.313
% Recovery		Amount Spiked	1.00
		Spiked S' Result	1.07
		% Recovery	104.4%

To Page No. \_\_\_\_\_

TEST: NITRATE, UNITS: mg/L  
 Meth. REF. # 353-2, MDL: 0.10  
 Instrument # #1 (Channel 1)

BY: CH  
 Book # CV-12 P. 97  
 FROM PAGE # 96

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.
7514	N6032	.334	5x	1.67	17866	P4851	.793	10x	7.93
7514	N6033	.0134	5x	Relo	<del>17866</del>	<del>P4852</del>	<del>.0487</del>	<del>10x</del>	
7562	N6122	.804	5x	4.02	17866	P4853	.442	10x	4.42
7503	P2683	.0213	5x	Relo	<del>17866</del>	<del>P4854</del>	<del>.0482</del>	<del>10x</del>	
7798	P2469	.276	10x	2.76	<del>17866</del>	<del>P3117</del>	<del>.153</del>	<del>10x</del>	<del>1.53</del>
7798	P2477	1.54	10x	15.4	<del>17866</del>	<del>P3120</del>	<del>.0175</del>	<del>10x</del>	
17798	P2524	.517	10x	5.17	<del>17866</del>	<del>P3123</del>	<del>.0211</del>	<del>10x</del>	
17798	P2538	.583	10x	5.83	<del>17866</del>	<del>P3124</del>	<del>.0229</del>	<del>10x</del>	
17798	P2539	.856	10x	8.56	<del>17866</del>	<del>P3126</del>	<del>.0176</del>	<del>10x</del>	
17798	P2540	.618	10x	6.18	17882	P2543	.528	10x	5.28
<del>7814</del>	<del>P3555</del>	<del>.101</del>	<del>10x</del>		<del>17882</del>	<del>P2544</del>	<del>.0253</del>	<del>10x</del>	
17822	P3107	.417	10x	4.17	17882	P2545	1.33	10x	13.3
<del>7822</del>	<del>P3112</del>	<del>.0139</del>	<del>10x</del>		17882	P2547	1.44	10x	14.4
17828	P2473	.442	10x	4.42	17882	P2549	.924	10x	9.24
17828	P2472	.547	10x	5.47	17882	P2550	.401	10x	4.01
17828	P2522	.417	10x	4.17	17882	P2551	1.03	10x	10.3
17828	P2536	.435	10x	4.35	17882	P2552	.717	10x	7.17
17828	P2548	.525	10x	5.25					
<del>7864</del>	<del>P2471</del>	<del>.042</del>	<del>10x</del>						
<del>7864</del>	<del>P2521</del>	<del>.0156</del>	<del>10x</del>						
<del>7864</del>	<del>P2471</del>	<del>.160</del>	<del>10x</del>	<del>1.60</del>					
17864	P2546	.697	10x	6.97					
<del>17864</del>	<del>P2537</del>	<del>.198</del>	<del>10x</del>	<del>1.98</del>					
17866	P4850	.550	10x	5.50					

PREPARED ON \_\_\_\_\_  
 BY \_\_\_\_\_  
 FROM \_\_\_\_\_ mg/L  
 STOCK # \_\_\_\_\_

STANDARD mg/L	VALUE OBTAINED	% of Theoretical
---------------	----------------	------------------

SLOPE	Intercept	Corr. Coef.
-------	-----------	-------------

ADDITIONAL COMMENTS

Straight Red Line Through  
Samples Done on Channel  
2, no dilution Page 98  
RET N6033 } Straight  
P2683 }

Analyses QA/QC Data Report

QC Batch # N3003 & N3004

Verified: 11/24/81 [Signature]

METHOD BLANK	
Spiked Blank (Known)	
Spiked Blank Result	
% Recovery	
EXTERNAL REF. QC	
Known Value	
Observed Result	
% Recovery	

PRECISION

Job # Dup (ETC)	Original Result	Duplicate Result	R %

ACCURACY

Job # Spiked (ETC)	Original Result	Amount Spiked	Spiked S' Result	% Recovery

305324

To Page No. \_\_\_\_\_

TEST: NITRATE, UNITS: mg/l  
 Meth. REF. # 353.2, MDL: 0.10  
 Instrument # #1 (Channel 2)

BY: CH  
 BOOK # CV-12 P. 98  
 FROM PAGE #     

Log Link	Job #	ANALYZED CONC.	DU	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	DU	REPORTED CONC.
1788	P2469	OVER		*	17886	P3117	1.15		1.15
1778	P2471	OVER		*	17866	P3120	.0207		BMDL
1778	P2524	OVER		*	17866	P3122	.0511		BMDL
1778	P2538	OVER		*	17866	P3124	.0173		BMDL
1778	P2539	OVER		*	17866	P3126	.0127		BMDL
1786	P2540	OVER		*	17882	P2543	OVER		*
1784	P3555	.487		.49	17882	P2544	.0250		BMDL
17822	P3107	OVER		*	17882	P2545	OVER		*
17822	P3112	.007		BMDL	17882	P2547	OVER		*
17828	P2473	OVER		*	17882	P2549	OVER		*
17828	P2472	OVER		*	17882	P2550	OVER		*
17828	P2522	OVER		*	17882	P2551	OVER		*
17828	P2530	OVER		*	17882	P2552	OVER		*
17828	P2548	OVER		*					
17864	P2481	.006		BMDL					
17864	P2531	.0245		BMDL					
17864	P2541	1.73		1.73					
17864	P2546	OVER		*					
17864	P2554	2.06		2.06					
17866	P4850	OVER		*					
17866	P4851	OVER		*					
17866	P4852	.398		.40					
17866	P4853	OVER		*					
17878	P2778	.0462		BMDL					

PREPARED ON \_\_\_\_\_  
 BY \_\_\_\_\_  
 FROM \_\_\_\_\_ mg/L  
 STOCK # \_\_\_\_\_

STANDARD mg/L	VALUE OBTAINED	% of Threshold
2.00	1.98	
1.00	1.04	
.50	.497	
.20	.0934	

SLOPE Intercept Corr. Coef.

ADDITIONAL COMMENTS  
 \* RPT on Channel 1  
 Complete Page 97

Analyses QA/QC Data Report      QC Batch # N003      Verified: 11/28/86 [Signature]

METHOD BLANK	.003	.0177
Spiked Blank (Known)	.5	.5
Spiked Blank Result	.501	.532
% Recovery	100%	1.06

EXTERNAL REF. OC	EPA
Known Value	95
Observed Result	953
% Recovery	1.02

Job # Dup (ETC)	P2473	P4850	P3126
Original Result	OVER	OVER	.0217
Duplicate Result			.0194
A RANGE	↓	↓	

Job # Spiked (ETC)	P2473	P4850	P3126
Original Result	.050	OVER	.021
Amount Spiked			.5
Spiked R Result	↓	↓	.506
% Recovery			97%

TEST: NITRITE, UNITS: mg/L  
 Meth. REF. # 353.2, MDL: 0.10  
 Instrument # #1 (Channel 2)

BY: S.T.  
 BOOK # CV-12 P. 100  
 FROM PAGE # 39

Log Link	Job #	ANALYZED CONC.	Dilut	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.
17524	N6110	0		B=DL	17866	P4850	.00151		B=DL
17524	N6120	.0028		B=DL	17866	P4851	.0166		B=DL
17524	N6126	.0040		B=DL	17866	P4852	.0490		B=DL
17531	P2710	.0024		B=DL	17866	P4853	.0446		B=DL
17531	P2725	0		B=DL	17878	P2776	.00181		B=DL
17539	P2117	.0043		B=DL	17866	P3120	.00114		B=DL
17539	P2116	.0954		B=DL	17866	P3122	0		B=DL
17539	P2114	.0571		B=DL	17866	P3124	.0164		B=DL
17524	P1711	.0043		B=DL	17866	P3126	0		B=DL
17524	P1713	.0040		B=DL	17882	P2543	.00475		B=DL
17524	P1714	0		B=DL	17882	P2544	.00213		B=DL
17524	P1717	.00114		B=DL	17882	P2545	.00473		B=DL
17562	N6128	0		B=DL	17882	P2547	.00784		B=DL
17562	N6148	.00119		B=DL	17882	P2549	.00119		B=DL
17573	P2717	0		B=DL	17886	P3117	.0066		B=DL
17579	P2726	.118		.12					
17584	P3068	.100		.10					
17503	P2683	.121		.12					
17562	N6122	.00357		B=DL					
17524	P1715	.00213		B=DL					
17864	P2551	.00119		B=DL					
17864	P2541	.0611		B=DL					
17864	P2546	.0691		B=DL					
17864	P2557	.00286		B=DL					

PREPARED ON \_\_\_\_\_  
 BY \_\_\_\_\_  
 FROM \_\_\_\_\_ mg/L  
 STOCK # \_\_\_\_\_

STANDARD mg/L	VALUE OBTAINED	% of theoretical
1.00	.99	99%
0.50	.51	102%
0.20	.197	98.5%
0.05	.046	92%

SLOPE	Intercept	Corr. Coef.
.0120	-.0005	.9998

**ADDITIONAL COMMENTS**

All calculations by  $X = mY + B$ .

**Analyses QA/QC Data Report**

QC Batch # (N3004 N02)

Verified: 11/24/86 *[Signature]*

METHOD BLANK		PRECISION	
Method Blank	.00337	Job # Dup (ETC)	N6110 P1714
Spiked Blank (Known)	.20	Original Result	0 0
Spiked Blank Result	.198	Duplicate Result	0 .0015
% Recovery	99	R Range	0.0 .0015
EXTERNAL REF. OC	N/A	ACCURACY	
Known Value	.5	Job # Spiked (ETC)	P1714
Observed Result	-.533	Original Result	0
% Recovery		Amount Spiked	.0382
		Spiked S' Result	
		% Recovery	1.31

305320



ANALYZED ON: 11/29/86  
 BY: L. V. ... / R. ...  
 BOOK # CV-12 P. 115  
 FROM PAGE # 114

TEST: Phenol, UNITS: mg/L  
 METH. REF. # 420.2, MDL: 0.050  
 INSTRUMENT # AA Channel #1

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.
17639	P2121	0.037	X10	.375	1749	P2514	TH		MATRIX RPT STR
↓	P2123	0.043	STR	BMDL					
17343	P4131	TH	X10	RPT <sup>STR</sup>					
17584	P3067	0.013	-	BMDL					
17639	N9882	TH	X5	RPT <sup>1:100</sup>					
17681	N9884	TH	X5	RPT <sup>1:100</sup>					
↓	N9880	TH	X5	RPT <sup>1:100</sup>					
↓	P2614	TH	X5	RPT <sup>STR</sup>					
17696	P0582	TH	-	RPT <sup>STR</sup>					
17824	P4180	TH	X5	RPT <sup>1:10</sup>					
17828	P2472	TH	X5	RPT					
17844	P253	TH	X5	RPT					
17868	P4512	TH	X5	RPT					
17866	P4851	0.067	X5	0.335					
17844	P2554	0.017	-	BMDL					
17866	P4851	0.013	-	BMDL					
↓	P4852	0.017	-	BMDL					
↓	P4853	0.022	-	BMDL					
17886	P5124	0.128	X10	1.28					
17707	N8672	0.048	-	BMDL					
↓	N8675	0.027	-	BMDL					
17721	N8674	0.131	X10	1.31					
1749	P2506	0.031	-	BMDL					
↓	P2513	0.096	X10	.964					

STANDARDS		
PREPARED ON	_____	
BY	_____	
FROM	_____	mg/L
STOCK #	_____	
STANDARD mg/L	VALUE OBTAINED	% of Theoretical
300		
200		
100		
0.050		
SLOPE	Intercept	Corr. Coef.

**ADDITIONAL COMMENTS**

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Analyses QA/QC Data Report QC Batch # PN003 Verified: [Signature] 11/29/86  
 SEE CV-12, P 114

METHOD BLANK		PRECISION	
Spiked Blank (Known)		Job # Dup (ETC)	
Spiked Blank Result		Original Result	
% Recovery		Duplicate Result	
		R 74-75	
		<b>ACCURACY</b>	
		Job # Spiked (ETC)	
		Original Result	
		Amount Spiked	
		Spiked S' Result	
		% Recovery	
<b>EXTERNAL REF. QC</b>			
Known Value			
X Observed Result			
X % Recovery			



TEST: Cyanide, UNITS: mg/L  
 METH. REF. # 335.3, MDL: 0.025  
 INSTRUMENT # #1 (citrate 1)

BY: C.H.  
 BOOK # CV-12 P. 119  
 FROM PAGE #     

Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	STANDARDS	
17837	N4709	0		BMDL	17697	P4748	.005		BMDL	PREPARED ON <u>12-2</u> BY <u>D.B.</u> FROM <u>    </u> mg/L STOCK # <u>    </u>	
17837	N4714	0		BMDL	17697	P4748	.004		BMDL		
17863	P0489	.007		BMDL	17697	P4748	.004		BMDL		
17863	P0493	.009		BMDL	17698	P1710	.003		BMDL		
17863	P0494	.005		BMDL	17701	N4710	.004		BMDL		
17863	P0495	.002		BMDL	17701	N4715	.004		BMDL	STANDARD mg/L	
17866	P4850	.539		539	17866	P4851	over		Redo	VALUE OBTAINED	
17868	P4510	*		Redo	17866	P4852	over		Redo	% of theoretical	
17878	P2776	.007		BMDL	17866	P4853	x		Redo	.500	
17881	P1865	.002		BMDL	17868	P4510	*		Redo	.400	
17883	N4708	.002		BMDL	17868	P4511	*		Redo	.200	
17883	N4711	.003		BMDL	17868	P4512	*		Redo	.100	
17833	P4241	.0197		BMDL	17868	P4513	*		Redo	SLOPE	
17833	P4242	.0169		BMDL	17883	N4712	*		Redo	Intercept	
17833	P4243	.0211		BMDL	17883	N4715	*		Redo	Corr. Coef.	
17833	P4244	.0186		BMDL	17889	P2754	*		Redo	ADDITIONAL COMMENTS	
17834	P4248	.0185		BMDL	17889	P2761	*		Redo	* RET. DUE TO INSTRUMENTATION Problem	
17834	P4249	.0189		BMDL	17889	P2774	*		Redo	P4851 10x	
17834	P4250	.0231		BMDL	17889	P2775	*		Redo	P4852 20x	
17834	P4251	.0189		BMDL	17833	P4235	*		Redo		
<del>17834</del>	<del>P2724</del>	<del>.026</del>		<del>BMDL</del>	17833	P4226	*		Redo		
17693	P2916	.0185		BMDL	17833	P4237	*		Redo		
17693	P2917	.0312		BMDL	17833	P4238	*		Redo		
17697	P4748	.006		BMDL	17833	P4239	*		Redo		

Analyses QA/QC Data Report

QC Batch # CY003

Verified: 12/3/86 *[Signature]*

PRECISION		ACCURACY	
METHOD BLANK	.008	Job # Dup (ETC)	N4709 P4748
Spiked Blank (Known)	.100	Original Result	0 0.006
Spiked Blank Result	.103	Duplicate Result	0 0.005
% Recovery	103%	R Range	0.0 0.001
EXTERNAL REF. QC	CPA	Job # Spikes (ETC)	N4709
Known Value	.224	Original Result	0.0
X Observed Result	.228	Amount Spiked	.200
X % Recovery	101.7%	Spiked S' Result	.201
		% Recovery	99%

To Page No.

**ETC ENVIRONMENTAL TESTING and CERTIFICATION**

**CONVENTIONAL CHEMISTRY LAB**

ANALYZED ON: 12-13-86

TEST: Cyanide

UNITS: mg/L

METH. REF. # 335.3

MDL: 0.025

INSTRUMENT # FI (Channel 1)

BY: C.B. / J.L.

BOOK # CV-12 P. 150

FROM PAGE #     

Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Diln	REPORTED CONC.
17866	P4853	.123	10x	1.23	17874	P4831	.006		BNDL
18065	P4032	.228	100	22.8	18236	P3873	.006		BNDL
18200	P6632	.0462		.046	18236	P3874	.006		BNDL
18203	P6634	.0246		.025	18236	P3881	.00958		BNDL
18203	P6641	.0252		.025	18236	P3882	.00777		BNDL
18207	P6665	.0183		BNDL	18236	P3883	.0059		BNDL
18207	P6666	.0270		.027	18236	P3884	.00589		BNDL
18207	P6667	.00568		BNDL	18236	P3885	.00587		BNDL
18207	P6668	.00391		BNDL	18236	P3886	.00586		BNDL
18203	P5794	.0667		.067	18236	P3889	.00584		BNDL
18205	P5796	.00509		BNDL	18236	P3890	.00583		BNDL
18210	P3809	.225	100	22.5	18234	P3867	.00581		BNDL
18210	P3813	interfer		Redo	18234	P3868	.0058		BNDL
18214	P6646	.006		BNDL	18234	P3877	.009		BNDL
18243	P6555	.00626		BNDL	18234	P3878	.00574		BNDL
18243	P6556	.00625		BNDL	18234	P3879	.00572		BNDL
18243	P6557	.00623		BNDL	18234	P3880	.00571		BNDL
18243	P6558	.006		BNDL	18234	P3893	.00569		BNDL
18243	P6559	.008		BNDL	18234	P3894	.00568		BNDL
18243	P6560	.006		BNDL	18234	P3895	.00566		BNDL
18243	P5792	.00972		BNDL	18234	P3896	.0113		BNDL
18253	P7918	.006		BNDL	18269	P3872	.007		BNDL
18264	P5795	.006		BNDL	18269	P3887	.00696		BNDL
18264	P6725	.0201		BNDL	18269	P3888	.0105		BNDL

PREPARED ON \_\_\_\_\_  
 BY \_\_\_\_\_  
 FROM \_\_\_\_\_ mg/L  
 STOCK # \_\_\_\_\_

STANDARD mg/L	VALUE OBTAINED	% of Theoretical
.500	.491	98.2%
.400	.406	101.5%
.200	.210	105%
.100	.0985	98.5%

SLOPE \_\_\_\_\_ Intercept \_\_\_\_\_ Cor \_\_\_\_\_

**ADDITIONAL COMMENTS**  
P3813 App Straight  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**Analyses QA/QC Data Report**

QC Batch # CY007

Verified: 12/16/86 J.L.

METHOD BLANK	Result
Spiked Blank (Known)	.10
Spiked Blank Result	.098
% Recovery	98%

EXTERNAL REF. QC	Result
Known Value	.224
Observed Result	.212
% Recovery	94.6%

PRECISION	P6667	P3881	P6558	P3888	P3877	P3881
Job # Dup (ETC)	P6667	P3881	P6558	P3888	P3877	P3881
Original Result	.00568	.009	.002	.0105	.106	.115
Duplicate Result	.00774	.006	.003	.007	.109	.118
R Range	.002	.003	.003	.009	.003	.003

ACCURACY	P6667	P3881	P6558	P3888	P3877	P3881
Job # Spiked (ETC)	P6667	P3881	P6558	P3888	P3877	P3881
Original Result	.0067	.008	.003	.009	.108	.117
Amount Spiked	.2	.2	.2	.2	.2	.2
Spiked S' Result	.228	.197	.215	.18	.304	.311
% Recovery	108.6%	98.5%	107.5%	92.5%	8%	97%

305330

TEST: CYANIDE, UNITS: mg/L  
 Meth. Ref. # EPA 335.3, MDL: 0.025 mg/L  
 Instrument # # 1 Chl # 1

By: MC/ST  
 Book # CV-12 p. 136  
 From Page # X

Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	Log Link	Job #	ANALYZED CONC.	Dilu	REPORTED CONC.	STANDARDS
17823	P4274	0.006	-	BMDL	18084	P4189	0.0102	-	BMDL	PREPARED ON _____ BY _____ FROM _____ mg/L STOCK # _____
17866	P4252	0.043	1:25	1.08	7	P4190	0.009	-	BMDL	
7	P4853	OFF SCALE	-	1:10	7	P4205	0.0098	-	BMDL	
7	P4851	0.0287	-	0.029	7	P4219	0.0104	-	BMDL	
17868	P4510	0.0024	-	BMDL	7	P4220	0.0104	-	BMDL	STANDARD mg/L    VALUE OBTAINED    % of Threshold
7	P4511	0.005	-	BMDL	7	P4223	0.0105	-	BMDL	
7	P4512	0.0718	-	0.072	7	P4224	0.0102	-	BMDL	
7	P4513	0.0081	-	BMDL	7	P5017	0.0139	-	BMDL	
17883	N4712	0.0075	-	BMDL	18121	P2983	0.0070	-	BMDL	0.500    0.204    100.8
7	N4713	0.0079	-	BMDL	18128	P5813	0.0087	-	BMDL	0.400    0.397    99.2
17889	P2759	0.0119	-	BMDL	7	P5814	0.0091	-	BMDL	0.200    0.198    99.0
7	P2761	0.0086	-	BMDL	18132	P2779	0.0088	-	BMDL	0.100    0.0952    95.2
7	P2774	0.006	-	BMDL	7	P2780	0.0064	-	BMDL	SLOPE    Intercept    Corr. Coef.
7	P2775	0.0093	-	BMDL	7	P2781	0.0101	-	BMDL	ADDITIONAL COMMENTS
17833	P4235	0.005	-	BMDL	18151	P5372	0.0108	-	BMDL	
7	P4236	0.007	-	BMDL	7	P5374	0.0116	-	BMDL	
7	P4237	0.008	-	BMDL	7	P5375	0.0108	-	BMDL	
7	P4238	0.009	-	BMDL	7	P5376	0.0124	-	BMDL	
7	P4239	0.006	-	BMDL	7	P5377	0.0141	-	BMDL	
7	P4240	0.009	-	BMDL	18195	P7549	0.0124	-	BMDL	
18065	P4030	0.154	1:100	15.4	7	P7550	0.0127	-	BMDL	
18068	N8922	0.006	-	BMDL	7	P7551	0.0127	-	BMDL	
18078	P5373	0.009	-	BMDL	7	P7552	0.0126	-	BMDL	
18084	P2987	0.0069	-	BMDL	7	P7553	0.0114	-	BMDL	

Analyses QA/QC Data Report    QC Batch # CY006    Verified: 12/9/86 FAK

METHOD BLANK		PRECISION					
Soaked Blank (Known)	0.200	Job # Dup (ETC)	P4274	N8922	P5017	P5376	P7557
Soaked Blank Result	0.182	Original Result	ND	0.007	ND	ND	ND
% Recovery	91.0	Duplicate Result	ND	0.017	ND	ND	ND
EXTERNAL REF. QC	EPA	R Range	0	0.010	0	0	0
Known Value	0.224	ACCURACY					
Observed Result	0.195	Job # Spiked (ETC)	P4274	N8922	P5017	P5376	P7557
% Recovery	87.0	Original Result	0	0.010	0	0	0
		Amount Spiked	0.20	0.100	0.200	0.200	0.300
		Spiked S' Result	0.182	0.121	0.264	0.26	0.293
		% Recovery	91	111.0	132.0	130.2	97.7

