

M E M O R A N D U M
Washington State Department of Ecology

April 2, 2013

TO: Weyerhaeuser Everett East File

FROM: David L. South

SUBJECT: Supplement to *Periodic Review, Weyerhaeuser Everett East Site, Facility Site ID# 11*, July 9, 2012.

Ecology conducted the subject periodic review to assess environmental conditions at the Weyerhaeuser Everett East cleanup site. Construction is complete for this site and the site is in long-term monitoring. This memorandum is a supplement to the periodic review. Please refer to the periodic review for additional detail.

The review found that ground water data collected by Sierra Pacific in 2005 found pentachlorophenol concentrations of 14,000 µg/L in the Upper Sand Aquifer and 15 µg/L in the Lower Sand Aquifer at location GP-10. These concentrations exceed the Consent Decree ground water cleanup level of 7.29 µg/L. GP-10 is located landward of the conditional point of compliance set at the Snohomish River.

Compliance monitoring data was not collected in the Lower Sand Aquifer during post-construction ground water monitoring. The Periodic Review recommended assessing pentachlorophenol concentrations in the vicinity of GP-10.

Ecology installed a number of wells to assess ground water conditions with respect to the Everett Smelter cleanup site. The Everett Smelter site includes Weyerhaeuser Everett East. Wells LLMW-17S (screened in the Upper Sand Aquifer) and LLMW-17D (screened in the Lower Sand Aquifer) were installed near the river bank downgradient from GP-10. The well locations are shown on Figure 1. Well information is as follows:

Location	Ecy_Well_Tag	Northing_Y	Easting_X	Surface_Elevation_feet	ToC_feet	momument	Screened_Interval_ft_bgs
LLMW-17S	BHU-039	371320.3207	1310602.283	15.32	18.27	Stickup	4-11
LLMW-17D	BHU-038	371317.6575	1310603.072	15.27	18.29	Stickup	15-25

Ecology's contractor, GeoEngineers, collected samples from these wells in January 2013 and sent them to Analytical Resources, Incorporated for analysis. GeoEngineers provided Weyerhaeuser with split samples for analysis by Weyerhaeuser Analytical & Testing Services. Ecology's results indicated pentachlorophenol was not detected in either well at a Method Reporting Limit of 0.25 µg/L. Weyerhaeuser's results indicated pentachlorophenol was not detected in either well at a Method Reporting Limit of 0.50 µg/L. Ecology's laboratory data sheets are included in Attachment A. Weyerhaeuser provided a table of their results, Attachment B.

As a result of this sampling Ecology considers the Weyerhaeuser Everett East site to currently be in compliance with the requirements of the Consent Decree. The next Periodic Review is to be performed by 2017. When performing this review, Ecology may require Weyerhaeuser to collect ground water quality data to evaluate whether protectiveness continues to be achieved at the ground water conditional point of compliance.

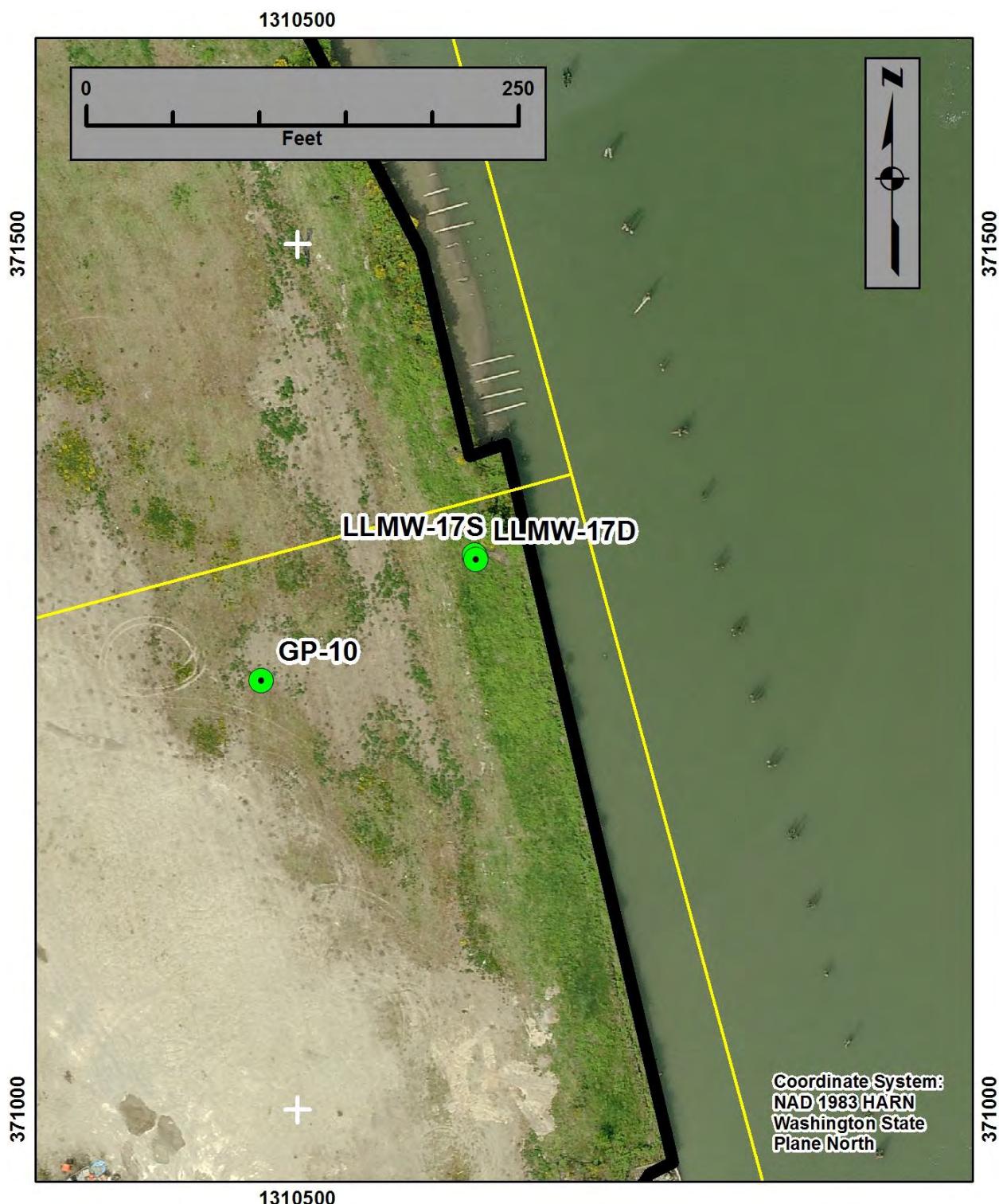


Figure 1: Location of GP-10, LLMW-17S, and LLMW-17D. Pentachlorophenol concentrations measured in LLMW-17S and LLMW-17D were < 0.25 µg/L. The cleanup level is 7.29 µg/L. See the Periodic Review for additional figures that show the entire site.

Attachment 1: Laboratory Data Sheets for Ecology Samples

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Client: GeoEngineers

Project: Everett

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bc
Signature

January-23-2013
Date



Analytical Resources, Incorporated
Analytical Chemists and Consultants

January 23, 2013

Garrett Leque
GeoEngineers, Inc.
1101 Fawcett, Suite 200
Tacoma, WA 98402

**RE: Everett
ARI Job No.: VZ72**

Dear Garrett:

Please find enclosed the Chain-of-Custody record (COC), sample receipt documentation, and the data package for samples from the project referenced above.

Sample receipt and details of these analyses are discussed in the Case Narrative.

An electronic copy of this package will remain on file with ARI. Should you have any questions or problems, please feel free to contact me at your convenience.

Sincerely,

ANALYTICAL RESOURCES, INC.

Cheronne Oreiro
Project Manager
(206) 695-6214
cheronneo@arilabs.com
www.arilabs.com

cc: eFile VZ72

Enclosures

Chain of Custody Documentation

ARI Job ID: VZ72

Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number:	Turn-around Requested:
✓772	NORMAL

ARI Client Company:	Phone:
GEOENGINEERS	253 383 4940

Client Contact:	GARRETT LEQUE
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Client Project Name:	EVERETT
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Client Project #:	NA	Samplers:	PAUL ROBICETTE
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Sample ID	Date	Time	Matrix	No. Containers
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LLMW175-130111-W **1/13** **1020** W **2** X

LLMW17D-130111-W **1/13** **1050** W **2** X

Page: **1** of **1**

Date: **1/11/13** Ice Present?

No. of Coolers: Cooler Temps:



Analytical Resources, Incorporated
Analytical Chemists and Consultants
4611 South 134th Place, Suite 100
Tukwila, WA 98168
206-695-6200 206-695-6201 (fax)

Penta EPA 2014*	Analysis Requested						Notes/Comments

Comments/Special Instructions * LDQ = 0.25 mg/L	Relinquished by (Signature) <i>Paul Robicette</i>	Received by (Signature) <i>Jennifer Milkay</i>	Relinquished by (Signature)	Received by (Signature)
	Printed Name: <i>Paul Robicette</i>	Printed Name: <i>Jennifer Milkay</i>	Printed Name	Printed Name
	Company: <i>GEOENGINEERS</i>	Company: <i>AKI</i>	Company	Company
	Date & Time: 1/14/13 1000	Date & Time: 1/14/13 1000	Date & Time	Date & Time

Limits of Liability: ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

Sample Retention Policy: All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.



ARI Client GeoEngineers

COC No(s). _____ NA

Assigned ARI Job No VZ72

Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES NO

Were custody papers included with the cooler? YES NO

Were custody papers properly filled out (ink, signed, etc) YES NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry) 18

If cooler temperature is out of compliance fill out form 00070F

Cooler Accepted by JM Date 1/14/13 Time 1600 Temp Gun ID# 132412024

Complete custody forms and attach all shipping documents

Log-In Phase:

Was a temperature blank included in the cooler? YES NO

What kind of packing material was used? ... Bubble Wrap Wet Ice Gel Packs Baggies Foam Block Paper Other: _____

Was sufficient ice used (if appropriate)? YES NO

Were all bottles sealed in individual plastic bags? YES NO

Did all bottles arrive in good condition (unbroken)? YES NO

Were all bottle labels complete and legible? YES NO

Did the number of containers listed on COC match with the number of containers received? YES NO

Did all bottle labels and tags agree with custody papers? YES NO

Were all bottles used correct for the requested analyses? YES NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs) YES NO

Were all VOC vials free of air bubbles? YES NO

Was sufficient amount of sample sent in each bottle? YES NO

Date VOC Trip Blank was made at ARI.. YES NO

Was Sample Split by ARI: YES Date/Time _____ Equipment: _____ Split by: _____

Samples Logged by _____ Date: 1/14/13 Time: 1601

**** Notify Project Manager of discrepancies or concerns ****

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

Additional Notes, Discrepancies, & Resolutions:

By _____ Date _____

Small Air Bubbles ~2mm • • •	Peabubbles' 2-4 mm • • •	LARGE Air Bubbles > 4 mm • • •	Small → "sm" Peabubbles → "pb" Large → "lg" Headspace → "hs"

Case Narrative, Data Qualifiers, Control Limits

ARI Job ID: VZ72

VZ72 : 00005



Case Narrative

Client: GeoEngineers

Project: Everett

ARI Job No.: VZ72

Sample Receipt

Two water samples were received on January 14, 2013 under ARI job VZ72. The cooler temperature measured by IR thermometer following ARI SOP was 0.8°C. For details regarding sample receipt, refer to the Cooler Receipt Forms.

Pentachlorophenol by SW8041

The samples were extracted and analyzed within the method recommended holding times.

Initial and continuing calibrations were within method requirements.

The surrogate percent recoveries were within control limits.

The method blank was clean at the reporting limit. The LCS and LCSD percent recoveries were within the control limits.

Sample ID Cross Reference ReportANALYTICAL
RESOURCES
INCORPORATED

ARI Job No: VZ72
Client: GeoEngineers
Project Event: N/A
Project Name: Everett

Sample ID	ARI Lab ID	ARI LIMS ID	Matrix	Sample Date/Time	VTSR
1. LLMW175-130111-W	VZ72A	13-907	Water	01/13/13 10:20	01/14/13 10:00
2. LLMW17D-130111-W	VZ72B	13-908	Water	01/13/13 10:50	01/14/13 10:00

Printed 01/14/13 Page 1 of 1

VZ72 : 00000 ?



Data Reporting Qualifiers

Effective 2/14/2011

Inorganic Data

- U Indicates that the target analyte was not detected at the reported concentration
- * Duplicate RPD is not within established control limits
- B Reported value is less than the CRDL but \geq the Reporting Limit
- N Matrix Spike recovery not within established control limits
- NA Not Applicable, analyte not spiked
- H The natural concentration of the spiked element is so much greater than the concentration spiked that an accurate determination of spike recovery is not possible
- L Analyte concentration is \leq 5 times the Reporting Limit and the replicate control limit defaults to ± 1 RL instead of the normal 20% RPD

Organic Data

-
- U Indicates that the target analyte was not detected at the reported concentration
 - * Flagged value is not within established control limits
 - B Analyte detected in an associated Method Blank at a concentration greater than one-half of ARI's Reporting Limit or 5% of the regulatory limit or 5% of the analyte concentration in the sample.
 - J Estimated concentration when the value is less than ARI's established reporting limits
 - D The spiked compound was not detected due to sample extract dilution
 - E Estimated concentration calculated for an analyte response above the valid instrument calibration range. A dilution is required to obtain an accurate quantification of the analyte
 - Q Indicates a detected analyte with an initial or continuing calibration that does not meet established acceptance criteria (<20%RSD, <20%Drift or minimum RRF).



Analytical Resources, Incorporated
Analytical Chemists and Consultants

- S Indicates an analyte response that has saturated the detector. **The calculated concentration is not valid; a dilution is required to obtain valid quantification of the analyte**
- NA The flagged analyte was not analyzed for
- NR Spiked compound recovery is not reported due to chromatographic interference
- NS The flagged analyte was not spiked into the sample
- M Estimated value for an analyte detected and confirmed by an analyst but with low spectral match parameters. This flag is used only for GC-MS analyses
- M2 The sample contains PCB congeners that do not match any standard Aroclor pattern. The PCBs are identified and quantified as the Aroclor whose pattern most closely matches that of the sample. The reported value is an estimate.
- N The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification"
- Y The analyte is not detected at or above the reported concentration. The reporting limit is raised due to chromatographic interference. The Y flag is equivalent to the U flag with a raised reporting limit
- EMPC Estimated Maximum Possible Concentration (EMPC) defined in EPA Statement of Work DLM02.2 as a value "calculated for 2,3,7,8-substituted isomers for which the quantitation and /or confirmation ion(s) has signal to noise in excess of 2.5, but does not meet identification criteria" **(Dioxin/Furan analysis only)**
- C The analyte was positively identified on only one of two chromatographic columns. Chromatographic interference prevented a positive identification on the second column
- P The analyte was detected on both chromatographic columns but the quantified values differ by ≥40% RPD with no obvious chromatographic interference
- X Analyte signal includes interference from polychlorinated diphenyl ethers. **(Dioxin/Furan analysis only)**
- Z Analyte signal includes interference from the sample matrix or perfluorokerosene ions **(Dioxin/Furan analysis only)**



Geotechnical Data

- A The total of all fines fractions. This flag is used to report total fines when only sieve analysis is requested and balances total grain size with sample weight.
- F Samples were frozen prior to particle size determination
- SM Sample matrix was not appropriate for the requested analysis. This normally refers to samples contaminated with an organic product that interferes with the sieving process and/or moisture content, porosity and saturation calculations
- SS Sample did not contain the proportion of "fines" required to perform the pipette portion of the grain size analysis
- W Weight of sample in some pipette aliquots was below the level required for accurate weighting



Chlorinated Phenols
DL, LOD, LOQ and Control Limit Summary ¹
EPA Method 8041A

Analyte	Aqueous Samples ^{2,3}				Solid Samples ^{2,4}				RPD ⁵
	DL µg/L	LOD µg/L	LOQ µg/L	LCS Recovery ⁶	DL µg/kg	LOD µg/kg	LOQ µg/kg	LCS Recovery ⁶	
Pentachlorophenol (PCP)	0.085	0.13	0.25	48 – 116%	0.83	3.13	6.25	56 – 111%	≤ 40
2,4,6-Trichlorophenol	0.106	0.13	0.25	30 – 160%	1.21	3.13	6.25	30 – 160%	≤ 40
2,3,6-Trichlorophenol	0.066	0.13	0.25	30 – 160%	2.96	3.13	6.25	30 – 160%	≤ 40
2,4,5-Trichlorophenol	0.077	0.13	0.25	30 – 160%	2.24	3.13	6.25	30 – 160%	≤ 40
2,3,4-Trichlorophenol	0.126	0.13	0.25	30 – 160%	1.67	3.13	6.25	30 – 160%	≤ 40
2,3,5,6-Tetrachlorophenol	0.091	0.13	0.25	30 – 160%	1.61	3.13	6.25	30 – 160%	≤ 40
2,3,4,5-Tetrachlorophenol	0.077	0.13	0.25	30 – 160%	1.26	3.13	6.25	30 – 160%	≤ 40
2,4-Dichlorophenol	1.510	3.0	6.0	30 – 160%	16.6	31.3	62.5	30 – 160%	≤ 40
Pentachlorophenol (PCP)(low level prep)	0.014 ⁷	0.013	0.025	36 – 159%					≤ 40
Surrogate % Recovery	MB / LCS	Sample			MB / LCS	Sample			
2,4,6-Tribromophenol	41 – 98	26 – 113			39 – 99	10 – 129 ⁸			≤ 40
2,4,6-Tribromophenol (low level preparation)	33 – 151	10 – 181 ⁸							≤ 40

(1) Detection Limit (DL), Limit of Detection (LOD) and Limit of Quantitation as defined in ARI SOP 1018S

(2) Control limits calculated using all data from 1/1/10 through 8/1/11.

(3) Separatory funnel extraction (EPA Method 3510C) 500 mL sample to 50 mL final volume (5 mL for low level Prep)

(4) Microwave assisted extraction (EPA Method 3546) 10 g sample to 25 mL final volume

(5) Relative Percent Difference between analytes in replicate analyzes. If C_O and C_D are the concentrations of the original and duplicate respectively then

$$RPD = \frac{|C_O - C_D|}{\frac{C_O + C_D}{2}} \times 100$$

(6) 30 – 160 are default values used when there is insufficient data to calculate historic control limits.

(7) MDL study QS97 (4/30/10)

(8) Highlighted control limits (**bold font**) are adjusted from the calculated values to reflect that ARI does not use control limits < 10 for the lower limit or < 100 for the upper limit.

**PCP/Chlorophenols Analysis
Report and Summary QC Forms**

ARI Job ID: VZ72

VZ72 : 00012

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
Extraction Method: SW3510C
 Page 1 of 1

Sample ID: LLMW175-130111-W
SAMPLE

Lab Sample ID: VZ72A
 LIMS ID: 13-907
 Matrix: Water
 Data Release Authorized: *MW*
 Reported: 01/23/13

QC Report No: VZ72-GeoEngineers
 Project: Everett

Date Sampled: 01/13/13
 Date Received: 01/14/13

Sample Amount: 500 mL
 Final Extract Volume: 50 mL
 Dilution Factor: 1.00

Date Extracted: 01/17/13
 Date Analyzed: 01/22/13 15:45
 Instrument/Analyst: ECD1/YZ

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	0.25	< 0.25 U
Reported in µg/L (ppb)			
Chlorophenol Surrogate Recovery			
	2, 4, 6-Tribromophenol		82.0%

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
Extraction Method: SW3510C
 Page 1 of 1

Sample ID: LLMW17D-130111-W
SAMPLE

Lab Sample ID: VZ72B
 LIMS ID: 13-908
 Matrix: Water
 Data Release Authorized: *MW*
 Reported: 01/23/13

QC Report No: VZ72-GeoEngineers
 Project: Everett

Date Sampled: 01/13/13
 Date Received: 01/14/13

Sample Amount: 500 mL
 Final Extract Volume: 50 mL
 Dilution Factor: 1.00

Date Extracted: 01/17/13
 Date Analyzed: 01/22/13 16:21
 Instrument/Analyst: ECD1/YZ

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	0.25	< 0.25 U

Reported in µg/L (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	80.8%
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SW8041 CHLOROPHENOLICS SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: VZ72-GeoEngineers
Project: Everett

<u>Client ID</u>	<u>TBP</u>	<u>TOT OUT</u>
LLMW175-130111-W	82.0%	0
MB-011713	76.4%	0
LCS-011713	91.4%	0
LCSD-011713	90.0%	0
LLMW17D-130111-W	80.8%	0

LCS/MB LIMITS QC LIMITS

(TBP) = 2,4,6-Tribromophenol (41-98) (26-113)

Prep Method: SW3510C
Log Number Range: 13-907 to 13-908

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LCS-011713
LCS/LCSD

Lab Sample ID: LCS-011713

QC Report No: VZ72-GeoEngineers
 Project: Everett

LIMS ID: 13-908

Matrix: Water

Data Release Authorized: *MW*

Reported: 01/23/13

Date Extracted LCS/LCSD: 01/17/13

Sample Amount LCS: 500 mL
 LCSD: 500 mL

Date Analyzed LCS: 01/22/13 13:56
 LCSD: 01/22/13 14:32

Final Extract Volume LCS: 50 mL
 LCSD: 50 mL

Instrument/Analyst LCS: ECD1/YZ
 LCSD: ECD1/YZ

Dilution Factor LCS: 1.00
 LCSD: 1.00

Analyte	LCS	Spike Added-LCS	LCS Recovery	LCSD	Spike Added-LCSD	LCSD Recovery	RPD
Pentachlorophenol	2.18	2.50	87.2%	2.19	2.50	87.6%	0.5%

Chlorophenols Surrogate Recovery

	LCS	LCSD
2,4,6-Tribromophenol	91.4%	90.0%

Results reported in $\mu\text{g}/\text{L}$

RPD calculated using sample concentrations per SW846.

4
CHLOROPHENOL METHOD BLANK SUMMARY

SAMPLE NO.

VZ72MBW1

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

Lab Sample ID: VZ72MBW1

Lab File ID: 0122A004

Matrix (soil/water) LIQUID

Extraction: (SepF/Cont/Sonc) SW3510C

Sulfur Cleanup (Y/N) Y

Date Extracted: 01/17/13

Date Analyzed (1): 01/22/13

Date Analyzed (2): 01/22/13

Time Analyzed (1): 1319

Time Analyzed (2): 1319

Instrument ID (1): ECD1

Instrument ID (2): ECD1

GC Column (1): STX CLP1 ID: 0.53 (mm)

GC Column (2): STX CLP2 ID: 0.53 (mm)

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

	CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED 1	DATE ANALYZED 2
01	VZ72LCSW1	VZ72LCSW1	01/22/13	01/22/13
02	VZ72LCSDW1	VZ72LCSDW1	01/22/13	01/22/13
03	LLMW175-1301	VZ72A	01/22/13	01/22/13
04	LLMW17D-1301	VZ72B	01/22/13	01/22/13

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
Extraction Method: SW3510C
Page 1 of 1

**ANALYTICAL
RESOURCES
INCORPORATED**

Sample ID: MB-011713
METHOD BLANK

Lab Sample ID: MB-011713
LIMS ID: 13-908
Matrix: Water
Data Release Authorized: *MMW*
Reported: 01/23/13

QC Report No: VZ72-GeoEngineers
Project: Everett

Date Sampled: NA
Date Received: NA

Sample Amount: 500 mL
Final Extract Volume: 50 mL
Dilution Factor: 1.00

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	0.25	< 0.25 U

Reported in µg/L (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	76.4%
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6D
CHLOROPHENOL INITIAL CALIBRATION
RETENTION TIME WINDOWS

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP1 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 09/21/12

COMPOUND	RT OF STANDARDS						MEAN	RT WINDOW	
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		RT	FROM
Pentachlorophenol	21.18	21.17	21.16	21.24	21.15	21.15	21.18	21.17	21.31
2,4,6-Trichloropheno	13.26	13.26	13.26	13.33	13.25	13.25	13.27	13.26	13.40
2,3,6-Trichloropheno	14.26	14.26	14.26	14.33	14.25	14.25	14.27	14.26	14.40
2,4,5-Trichloropheno	16.04	16.02	16.02	16.09	16.00	16.00	16.03	16.02	16.16
2,3,4-Trichloropheno	17.57	17.54	17.53	17.60	17.51	17.51	17.55	17.53	17.67
2,3,5,6-Tetrachlorop	17.33	17.32	17.32	17.39	17.31	17.31	17.33	17.32	17.46
2,3,4,5-Tetrachlorop	20.37	20.35	20.33	20.41	20.31	20.31	20.35	20.34	20.48
2,4-Dichlorophenol	12.78	12.78	12.78	12.78	12.78	12.78	12.78	12.71	12.85
2,4,6-Tribromophenol	18.79	18.78	18.77	18.84	18.76	18.75	18.79	18.77	18.91

6D
CHLOROPHENOL INITIAL CALIBRATION
RETENTION TIME WINDOWS

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP2 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 09/21/12

COMPOUND	RT OF STANDARDS						MEAN RT	RT WINDOW	
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		FROM	TO
Pentachlorophenol	22.91	22.90	22.89	22.96	22.88	22.88	22.91	22.89	23.03
2,4,6-Trichloropheno	14.27	14.27	14.27	14.34	14.26	14.26	14.28	14.27	14.41
2,3,6-Trichloropheno	15.51	15.51	15.51	15.58	15.50	15.50	15.52	15.51	15.65
2,4,5-Trichloropheno	17.45	17.43	17.43	17.50	17.42	17.41	17.45	17.43	17.57
2,3,4-Trichloropheno	19.00	18.98	18.97	19.05	18.96	18.96	18.99	18.98	19.12
2,3,5,6-Tetrachlorop	18.77	18.76	18.75	18.83	18.75	18.75	18.77	18.76	18.90
2,3,4,5-Tetrachlorop	22.04	22.03	22.02	22.09	22.01	22.00	22.04	22.02	22.16
2,4-Dichlorophenol	13.79	13.78	13.78	13.85	13.77	13.77	13.79	13.78	13.92
2,4,6-Tribromophenol	20.89	20.88	20.87	20.94	20.86	20.86	20.89	20.87	21.01

6E
CHLOROPHENOL INITIAL CALIBRATION
CALIBRATION FACTORS

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP1 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 09/21/12

COMPOUND	CALIBRATION FACTORS						R^2/ %RSD	CT
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		
Pentachlorophenol	77052	71102	62381	56971	51600	44551	0.9999	Q
2,4,6-Trichlorophenol	37800	36793	34435	30695	27572	23819	17.2	A
2,3,6-Trichlorophenol	39363	41432	38253	34053	27204	22897	0.9992	Q
2,4,5-Trichlorophenol	23290	24192	22777	19742	16572	13274	1.0000	Q
2,3,4-Trichlorophenol	35575	30905	28044	24569	21009	17260	1.0000	Q
2,3,5,6-Tetrachloroph	53827	53395	50427	46244	41856	36207	14.8	A
2,3,4,5-Tetrachloroph	50607	45516	41632	36089	31716	27069	0.9998	Q
2,4-Dichlorophenol	2363	1981	1759	1559	1390	1206	0.9998	Q
2,4,6-Tribromophenol	63570	57559	52456	46332	41843	35975	0.9999	Q
							AVE RSD	21.0

CT stands for Curve Types:

- A Indicates an Average Response Factor Curve
- L Indicates a Linear Curve
- Q Indicates a Quadratic Curve

CALIBRATION FILES

LVL 1: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A018.d
 LVL 2: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A019.d
 LVL 3: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A020.d
 LVL 4: /chem2/ecd1.i/PCP20120921.b/1205-1.b/1205A012.d
 LVL 5: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A021.d
 LVL 6: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A022.d

6E
CHLOROPHENOL INITIAL CALIBRATION
CALIBRATION FACTORS

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP2 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 09/21/12

COMPOUND	CALIBRATION FACTORS						R^2/ %RSD	CT
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		
Pentachlorophenol	42596	39717	37527	34681	31627	27985	15.0	A
2,4,6-Trichlorophenol	32528	33312	25170	22050	20938	17684	0.9992	Q
2,3,6-Trichlorophenol	25874	25680	24007	20973	17820	15188	0.9997	Q
2,4,5-Trichlorophenol	12584	13469	15051	12710	10198	7860	0.9996	Q
2,3,4-Trichlorophenol	23072	18508	18408	15761	13094	9728	0.9988	Q
2,3,5,6-Tetrachloroph	33035	31396	31059	28607	26066	23149	12.9	A
2,3,4,5-Tetrachloroph	25112	23550	21899	20040	18059	15675	17.0	A
2,4-Dichlorophenol	1292	1502	1138	970	877	712	0.9993	Q
2,4,6-Tribromophenol	33881	32521	32652	29584	27527	24749	11.7	A
							AVE RSD	19.8

CT stands for Curve Types:

- A Indicates an Average Response Factor Curve
- L Indicates a Linear Curve
- Q Indicates a Quadratic Curve

CALIBRATION FILES

LVL 1: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A018.d
 LVL 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A019.d
 LVL 3: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A020.d
 LVL 4: /chem2/ecd1.i/PCP20120921.b/1205-2.b/1205A012.d
 LVL 5: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A021.d
 LVL 6: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A022.d

7E
CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP1 ID: 0.53 (mm)

Init. Calib. Date(s): 09/21/12 12/05/12

Client Sample No. (PCP):

Date Analyzed : 01/22/13

Lab Sample ID (PCP): PCPCCAL

Time Analyzed : 1243

PCP MIX COMPOUND	RT	RT WINDOW FROM	TO	CALC AMOUNT	NOM AMOUNT	%D
Pentachlorophenol	21.27	21.17	21.31	26.3	25.0	5.2
2,4,6-Trichlorophenol	13.36	13.26	13.40	28.2	25.0	12.8
2,3,6-Trichlorophenol	14.36	14.26	14.40	25.3	25.0	1.2
2,4,5-Trichlorophenol	16.12	16.02	16.16	23.0	25.0	-8.0
2,3,4-Trichlorophenol	17.63	17.53	17.67	27.5	25.0	10.0
2,3,5,6-Tetrachlorophenol	17.42	17.32	17.46	26.3	25.0	5.2
2,3,4,5-Tetrachlorophenol	20.44	20.34	20.48	27.0	25.0	8.0
2,4-Dichlorophenol	12.81	12.71	12.85	271	250	8.4
2,4,6-Tribromophenol (surr)	18.87	18.77	18.91	26.3	25.0	5.2

AVERAGE %D = 7.1

7E
CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP2 ID: 0.53 (mm)

Init. Calib. Date(s): 09/21/12 12/05/12

Client Sample No. (PCP):

Date Analyzed :01/22/13

Lab Sample ID (PCP): PCPCCAL

Time Analyzed :1243

PCP MIX COMPOUND	RT	RT WINDOW FROM	TO	CALC AMOUNT	NOM AMOUNT	%D
Pentachlorophenol	22.99	22.89	23.03	26.7	25.0	6.8
2,4,6-Trichlorophenol	14.36	14.27	14.41	23.0	25.0	-8.0
2,3,6-Trichlorophenol	15.61	15.51	15.65	23.3	25.0	-6.8
2,4,5-Trichlorophenol	17.52	17.43	17.57	23.2	25.0	-7.2
2,3,4-Trichlorophenol	19.07	18.98	19.12	24.0	25.0	-4.0
2,3,5,6-Tetrachlorophenol	18.85	18.76	18.90	25.2	25.0	0.8
2,3,4,5-Tetrachlorophenol	22.12	22.02	22.16	26.2	25.0	4.8
2,4-Dichlorophenol	13.88	13.78	13.92	214	250	-14.4
2,4,6-Tribromophenol (surr)	20.97	20.87	21.01	25.5	25.0	2.0

AVERAGE %D = 6.1

7E
CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP1 ID: 0.53 (mm)

Init. Calib. Date(s): 09/21/12 12/05/12

Client Sample No.(PCP):

Date Analyzed :01/22/13

Lab Sample ID (PCP): PCPCCAL

Time Analyzed :1733

PCP MIX COMPOUND	RT	RT WINDOW FROM	TO	CALC AMOUNT	NOM AMOUNT	%D
Pentachlorophenol	21.27	21.17	21.31	26.7	25.0	6.8
2,4,6-Trichlorophenol	13.36	13.26	13.40	25.8	25.0	3.2
2,3,6-Trichlorophenol	14.36	14.26	14.40	22.6	25.0	-9.6
2,4,5-Trichlorophenol	16.12	16.02	16.16	23.5	25.0	-6.0
2,3,4-Trichlorophenol	17.63	17.53	17.67	28.0	25.0	12.0
2,3,5,6-Tetrachlorophenol	17.42	17.32	17.46	26.9	25.0	7.6
2,3,4,5-Tetrachlorophenol	20.44	20.34	20.48	27.9	25.0	11.6
2,4-Dichlorophenol	12.82	12.71	12.85	273	250	9.2
2,4,6-Tribromophenol (surr)	18.87	18.77	18.91	27.4	25.0	9.6

AVERAGE %D = 8.4

7E
CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP2 ID: 0.53 (mm)

Init. Calib. Date(s): 09/21/12 12/05/12

Client Sample No. (PCP):

Date Analyzed : 01/22/13

Lab Sample ID (PCP): PCPCCAL

Time Analyzed : 1733

PCP MIX COMPOUND	RT	RT WINDOW FROM	TO	CALC AMOUNT	NOM AMOUNT	%D
Pentachlorophenol	22.99	22.89	23.03	27.8	25.0	11.2
2,4,6-Trichlorophenol	14.37	14.27	14.41	23.9	25.0	-4.4
2,3,6-Trichlorophenol	15.61	15.51	15.65	27.8	25.0	11.2
2,4,5-Trichlorophenol	17.53	17.43	17.57	22.0	25.0	-12.0
2,3,4-Trichlorophenol	19.07	18.98	19.12	22.9	25.0	-8.4
2,3,5,6-Tetrachlorophenol	18.86	18.76	18.90	25.7	25.0	2.8
2,3,4,5-Tetrachlorophenol	22.12	22.02	22.16	26.8	25.0	7.2
2,4-Dichlorophenol	13.88	13.78	13.92	233	250	-6.8
2,4,6-Tribromophenol (surr)	20.97	20.87	21.01	26.0	25.0	4.0

AVERAGE %D = 7.6

CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP1 ID: 0.53 (mm) Instrument ID: ECD1

Init. Calib. Date(s): 09/21/12 12/05/12

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION S1 : 18.84				
CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	S1 RT #
01	PCP D	09/21/12	1928	18.76
02	PCP A	09/21/12	2004	18.79
03	PCP B	09/21/12	2041	18.78
04	PCP C	09/21/12	2117	18.77
05	PCP E	09/21/12	2153	18.76
06	PCP F	09/21/12	2230	18.75
07	PCPCCAL	01/22/13	1243	18.87
08	VZ72MBW1	01/22/13	1319	18.89
09	VZ72LCSW1	01/22/13	1356	18.89
10	VZ72LCSDW1	01/22/13	1432	18.89
11	ZZZZZ	01/22/13	1508	18.89
12	LLMW175-1301	VZ72A	01/22/13	1545
13	LLMW17D-1301	VZ72B	01/22/13	1621
14	PCPCCAL	01/22/13	1733	18.87

QC LIMITS
S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No.: VZ72

Project: EVERETT

GC Column: STX CLP2 ID: 0.53 (mm) Instrument ID: ECD1

Init. Calib. Date(s): 09/21/12 12/05/12

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION S1 : 20.94				
CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	S1 RT #
01	PCP D	09/21/12	1928	20.87
02	PCP A	09/21/12	2004	20.89
03	PCP B	09/21/12	2041	20.88
04	PCP C	09/21/12	2117	20.87
05	PCP E	09/21/12	2153	20.86
06	PCP F	09/21/12	2230	20.86
07	PCPCCAL	01/22/13	1243	20.97
08	VZ72MBW1	01/22/13	1319	20.98
09	VZ72LCSW1	01/22/13	1356	20.98
10	VZ72LCSDW1	01/22/13	1432	20.98
11	ZZZZZ	01/22/13	1508	20.98
12	LLMW175-1301	VZ72A	01/22/13	1545
13	LLMW17D-1301	VZ72B	01/22/13	1621
14		PCPCCAL	01/22/13	20.97
			1733	

QC LIMITS

S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

**PCP/Chlorophenols Raw Data
Extraction Bench Sheets and Notes**

ARI Job ID: VZ72

VZ72 : 00029



Analytical Resources,
Incorporated
Analytical Chemists and
Consultants

Organic Extractions Laboratory
Analyst Notes

ARI Job No.:

VZ72

Client ID:

Geo Engineers

Parameter: PCP

Client Project: Everett

Screens: Soil/Sediment/Solid/Other:	Analyst/Date
<input type="checkbox"/> No Anomalies (standard soil/wet sediment/sand/gravel)=	
<input type="checkbox"/> Standing Water Decanted (Not shared)=	
<input type="checkbox"/> Standing Water Homogenized (Shared samples)=	
<input type="checkbox"/> Clay/Clumps (Difficult to homogenize)=	
<input type="checkbox"/> Rocks (%+size)?	
<input type="checkbox"/> Organics (Leaves/sticks/grass)=	
<input type="checkbox"/> Oily, obvious fuel/sulfur odors=	
<input type="checkbox"/> Other (Details)=	
Aqueous:	
<input type="checkbox"/> No Anomalies	
<input checked="" type="checkbox"/> Turbid/Color= B ~ pale yellow.	SP 1/17/13
<input type="checkbox"/> Particulates(%)=(Note: >5% =Notify Supervisor/Lead)	
<input type="checkbox"/> Emulsions (%)=	
<input type="checkbox"/> Other (Details)=	
<input type="checkbox"/> Other Notes/Comments= (Note problems, concerns, corrective actions). (Centrifuge#1 used for all Centrifugations)	

**PCP/Chlorophenols Raw Data
Initial Calibration**

ARI Job ID: VZ72

VZ72 : 000032



Analytical Resources, Incorporated
Analytical Chemists and Consultants

GC Initial Calibration Notes

ARI SOP: 403S(PCB) 405S(Herb) 407S(TPH-D) 409S(HCID) 412S(PCP) 423S(Pest)
427S(Dir Inj) 428S(EPH) Other

Instrument: FID-3A FID-3B FID-4A FID-4B FID-5 FID-7 FID-8
FID-9 ECD-1 ECD-5 ECD-6 ECD-7 ECD-8

Curve Date(s): 9/21/2012 Internal Standard ID NA Expiration NA

Endrin/DDT Breakdown <15%? YES / NO / NA ICV Exceeding ±20%? YES / NO
ICal Meets %RSD & r^2 Criteria YES / NO ICV Exceeding ±30%? YES / NO
Manual Integrations for ICal? YES / NO Linear Fits Used? YES / NO
Minimum Response S/N Met YES / NO Quadratic Fits Used? YES / NO
Calibration Points Dropped? YES / NO

Primary Source	Standard #	Expiration	Secondary Source	Standard #	Expiration
<u>Arcu Standard</u>	<u>#4444-92712</u> <u>1919-3</u>	<u>10/15/2012</u>		<u>1951-2</u>	<u>6/8/2009</u>

Detail problems, corrective actions and/or other pertinent information below:

Analyst: / Date: 9/25/2012

Reviewer: BB Date: 9/26/12

GC LOG SUMMARY FOR DATABATCH - /chem2/ecd1.i/PCP20120921.b/ical-2.b

	Inject Date/Time	Filename	DF	LabID	ClientID
1	21-SEP-2012 19:28	0921A017.d	1	PCP D	
2	21-SEP-2012 20:04	0921A018.d	1	PCP A	
3	21-SEP-2012 20:41	0921A019.d	1	PCP B	
4	21-SEP-2012 21:17	0921A020.d	1	PCP C	
5	21-SEP-2012 21:53	0921A021.d	1	PCP E	
6	21-SEP-2012 22:30	0921A022.d	1	PCP F	
7	21-SEP-2012 23:06	0921A023.d	1	PCP ICV	

MANUAL INTEGRATION SUMMARY FOR DATABATCH - /chem2/ecd1.i/PCP20120921.b/ical-1.b

ARI Job No.: PCP Method: PCP.m Instrument: ecd1.i Date: 21-SEP-2012

Time	Filename	LabID	ClientId	DF	Manually Integrated Compounds
1928	0921A017.d	PCP D		1	NO MANUAL INTEGRATION
2004	0921A018.d	PCP A		1	NO MANUAL INTEGRATION
2041	0921A019.d	PCP B		1	NO MANUAL INTEGRATION
2117	0921A020.d	PCP C		1	NO MANUAL INTEGRATION
2153	0921A021.d	PCP E		1	NO MANUAL INTEGRATION
2230	0921A022.d	PCP F		1	NO MANUAL INTEGRATION

MANUAL INTEGRATION SUMMARY FOR DATABATCH - /chem2/ecd1.i/PCP20120921.b/ical-2.b

ARI Job No.: PCP Method: PCPB.m Instrument: ecd1.i Date: 21-SEP-2012

Time	Filename	LabID	ClientId	DP	Manually Integrated Compounds
1928	0921A017.d	PCP D		1	NO MANUAL INTEGRATION
2004	0921A018.d	PCP A		1	NO MANUAL INTEGRATION
2041	0921A019.d	PCP B		1	NO MANUAL INTEGRATION
2117	0921A020.d	PCP C		1	NO MANUAL INTEGRATION
2153	0921A021.d	PCP E		1	NO MANUAL INTEGRATION
2230	0921A022.d	PCP F		1	NO MANUAL INTEGRATION
2306	0921A023.d	PCP ICV		1	NO MANUAL INTEGRATION

Report Date : 26-Sep-2012 09:35

Page 1

Analytical Resources, Inc.
RETENTION TIME SUMMARY REPORT

Method File: /chem2/ecd1.i/PCP20120921.b/PCPB.m
Batch File: /chem2/ecd1.i/PCP20120921.b/ical-2.b
Inst ID: ecd1.i

ID:	RT01	RT02	RT03	RT04	RT05	RT06	RT07
FILENAME:	0921A017	0921A018	0921A019	0921A020	0921A021	0921A022	0921A023
INJ.DATE:	21-SEP-2012						
INJ.TIME:	19:28	20:04	20:41	21:17	21:53	22:30	23:06

Compound	RT01	RT02	RT03	RT04	RT05	RT06	RT07	EXPEC RT	RT WINDOW	Avg RT	STD DEV
1 2,4-Dichlorophenol	13.778	13.787	13.782	13.780	13.775	13.773	13.779	13.778	13.708-13.848	13.779	0.005
2 2,4,6-Trichlorophenol	14.264	14.268	14.266	14.265	14.261	14.259	14.264	14.264	14.194-14.334	14.264	0.003
3 2,3,6-Trichlorophenol	15.507	15.514	15.510	15.508	15.504	15.502	15.510	15.507	15.437-15.577	15.508	0.004
4 2,4,5-Trichlorophenol	17.423	17.448	17.434	17.427	17.418	17.414	17.433	17.423	17.353-17.493	17.428	0.012
5 2,3,5,6-Tetrachlorophenol	18.752	18.767	18.759	18.754	18.748	18.746	18.759	18.752	18.682-18.822	18.755	0.007
6 2,3,4-Trichlorophenol	18.968	19.001	18.984	18.974	18.961	18.957	18.982	18.968	18.898-19.038	18.975	0.015
7 2,4,6-Tribromophenol	20.867	20.887	20.877	20.870	20.863	20.861	20.878	20.867	20.797-20.937	20.872	0.009
8 2,3,4,5-Tetrachlorophenol	22.012	22.044	22.029	22.019	22.006	22.003	22.024	22.012	21.942-22.082	22.020	0.014
9 Pentachlorophenol	22.888	22.906	22.897	22.892	22.884	22.882	22.896	22.888	22.818-22.958	22.892	0.008

Reviewer 1

AF

Date: 9/26/2012

Reviewer 2

BB

Date: 9/26/12

Analytical Resources, Inc.
RETENTION TIME SUMMARY REPORT

Method File: /chem2/ecd1.i/PCP20120921.b/PCP.m
 Batch File: /chem2/ecd1.i/PCP20120921.b/ical-1.b
 Inst ID: ecd1.i

ID:	RT01	RT02	RT03	RT04	RT05	RT06
FILENAME:	0921A017	0921A018	0921A019	0921A020	0921A021	0921A022
INJ. DATE:	21-SEP-2012	21-SEP-2012	21-SEP-2012	21-SEP-2012	21-SEP-2012	21-SEP-2012
INJ. TIME:	19:28	20:04	20:41	21:17	21:53	22:30

Compound	RT01	RT02	RT03	RT04	RT05	RT06	EXPEC RT	RT WINDOW	Avg RT	STD DEV
1 2,4-Dichlorophenol	12.711	12.719	12.716	12.713	12.707	12.706	12.711	12.641-12.781	12.712	0.005
2 2,4,6-Trichlorophenol	13.256	13.260	13.257	13.256	13.254	13.252	13.256	13.186-13.326	13.256	0.003
3 2,3,6-Trichlorophenol	14.254	14.262	14.258	14.256	14.251	14.249	14.254	14.184-14.324	14.255	0.005
4 2,4,5-Trichlorophenol	16.010	16.044	16.025	16.015	16.003	15.999	16.010	15.940-16.080	16.016	0.016
5 2,3,5,6-Tetrachlorophenol	17.315	17.334	17.323	17.318	17.311	17.307	17.315	17.245-17.385	17.318	0.009
6 2,3,4-Trichlorophenol	17.519	17.566	17.541	17.527	17.511	17.506	17.519	17.449-17.589	17.528	0.022
7 2,4,6-Tribromophenol	18.763	18.793	18.776	18.768	18.757	18.753	18.763	18.693-18.833	18.768	0.014
8 2,3,4,5-Tetrachlorophenol	20.323	20.368	20.346	20.331	20.314	20.308	20.323	20.253-20.393	20.332	0.022
9 Pentachlorophenol	21.158	21.182	21.170	21.163	21.153	21.149	21.158	21.088-21.228	21.162	0.012

Reviewer 1 _____

AR Date: 9/26/2012

Reviewer 2 _____

BB Date: 9/26/12

Analytical Resources, Inc.

INITIAL CALIBRATION DATA

Start Cal Date : 21-SEP-2012 19:28
 End Cal Date : 21-SEP-2012 22:30
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem2/ecd1.i/PCP20120921.b/PCPB.m
 Cal Date : 25-Sep-2012 10:11 aron
 Curve Type : Average

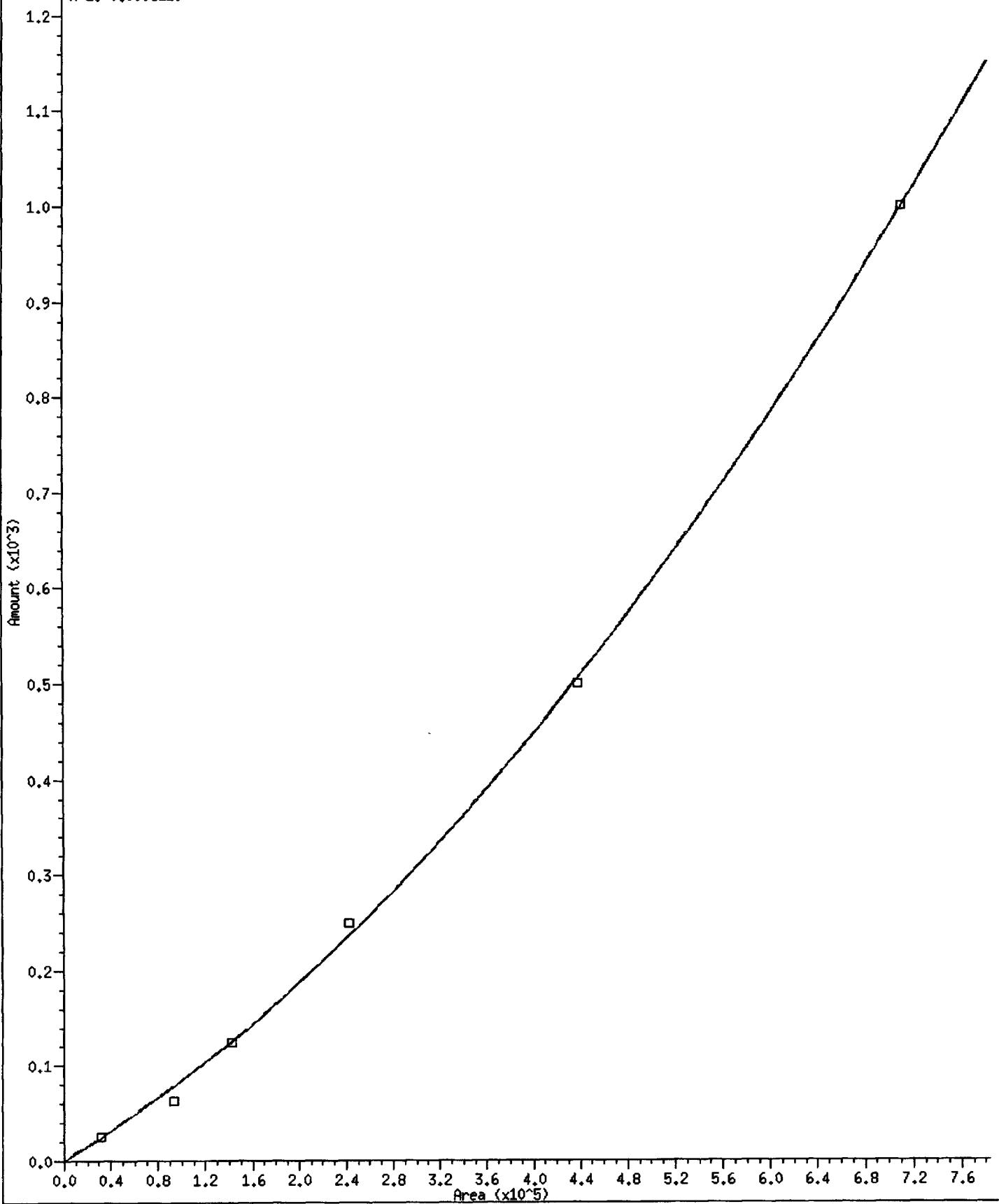
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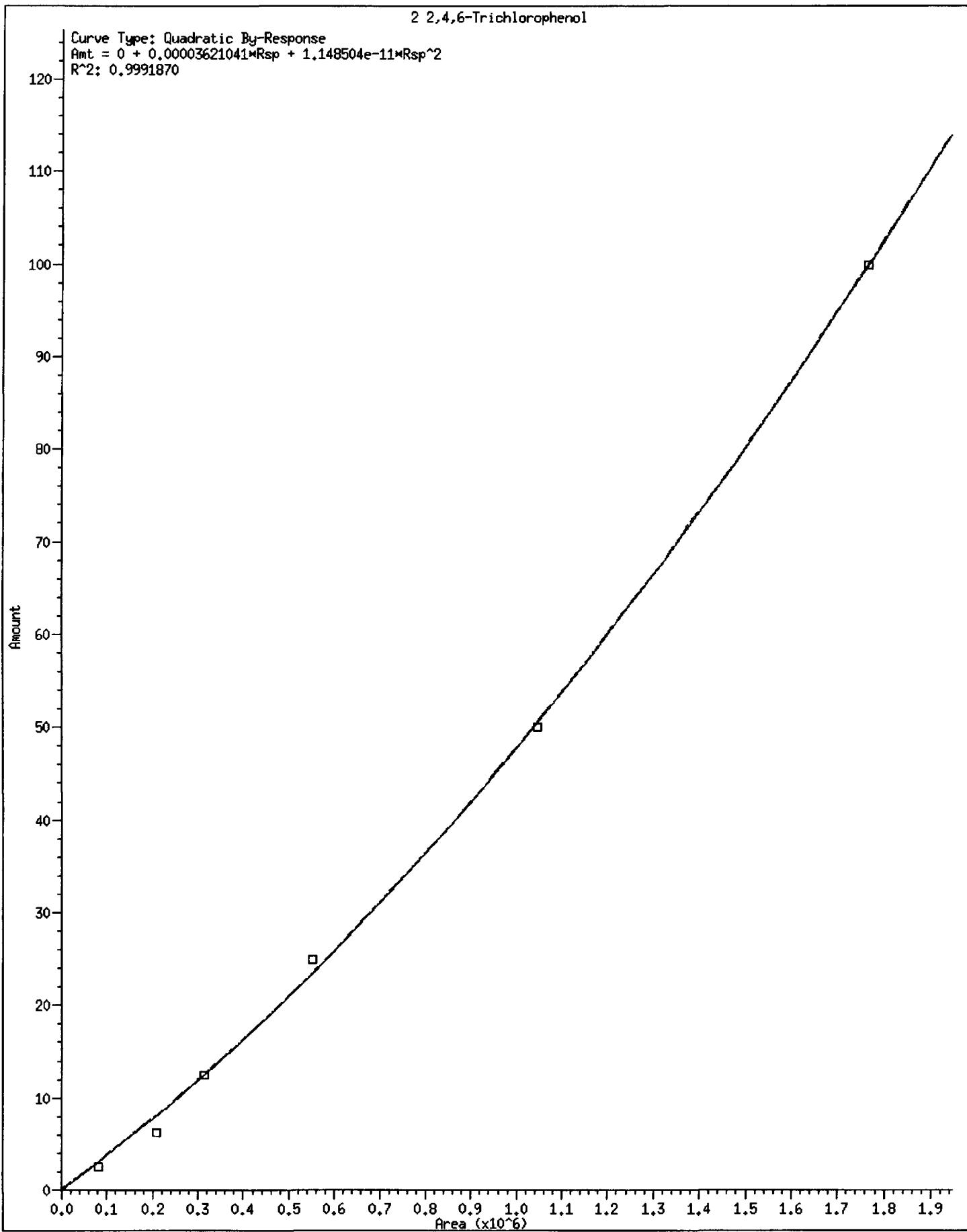
Level 1: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A018.d
 Level 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A019.d
 Level 3: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A020.d
 Level 4: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A017.d
 Level 5: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A021.d
 Level 6: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A022.d

Compound	2.500	6.250	12.500	25.000	50.000	100.000	RRF	% RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		
1 2,4-Dichlorophenol	1292	1502	1138	970	877	712	1082	26.637 <-
2 2,4,6-Trichlorophenol	32528	33312	25171	22050	20938	17684	25280	25.269 <-
3 2,3,6-Trichlorophenol	25874	25680	24007	20973	17820	15189	21590	20.347 <-
4 2,4,5-Trichlorophenol	12584	13469	15051	12710	10198	7860	11979	21.347 <-
5 2,3,5,6-Tetrachlorophenol	33035	31396	31059	28607	26066	23149	28886	12.877
6 2,3,4-Trichlorophenol	23072	18508	18408	15761	13094	9728	16428	28.400 <-
8 2,3,4,5-Tetrachlorophenol	25112	23550	21899	20040	18059	15675	20723	16.950
9 Pentachlorophenol	42596	39717	37527	34681	31627	27985	35689	15.046
\$ 7 2,4,6-Tribromophenol (surr)	33881	32521	32652	29584	27527	24749	30152	11.695

1 2,4-Dichlorophenol

Curve Type: Quadratic By-Response
Amt = 0 + 0.0007495282*Rsp + 9.19662e-10*Rsp^2
R^2: 0.9993119



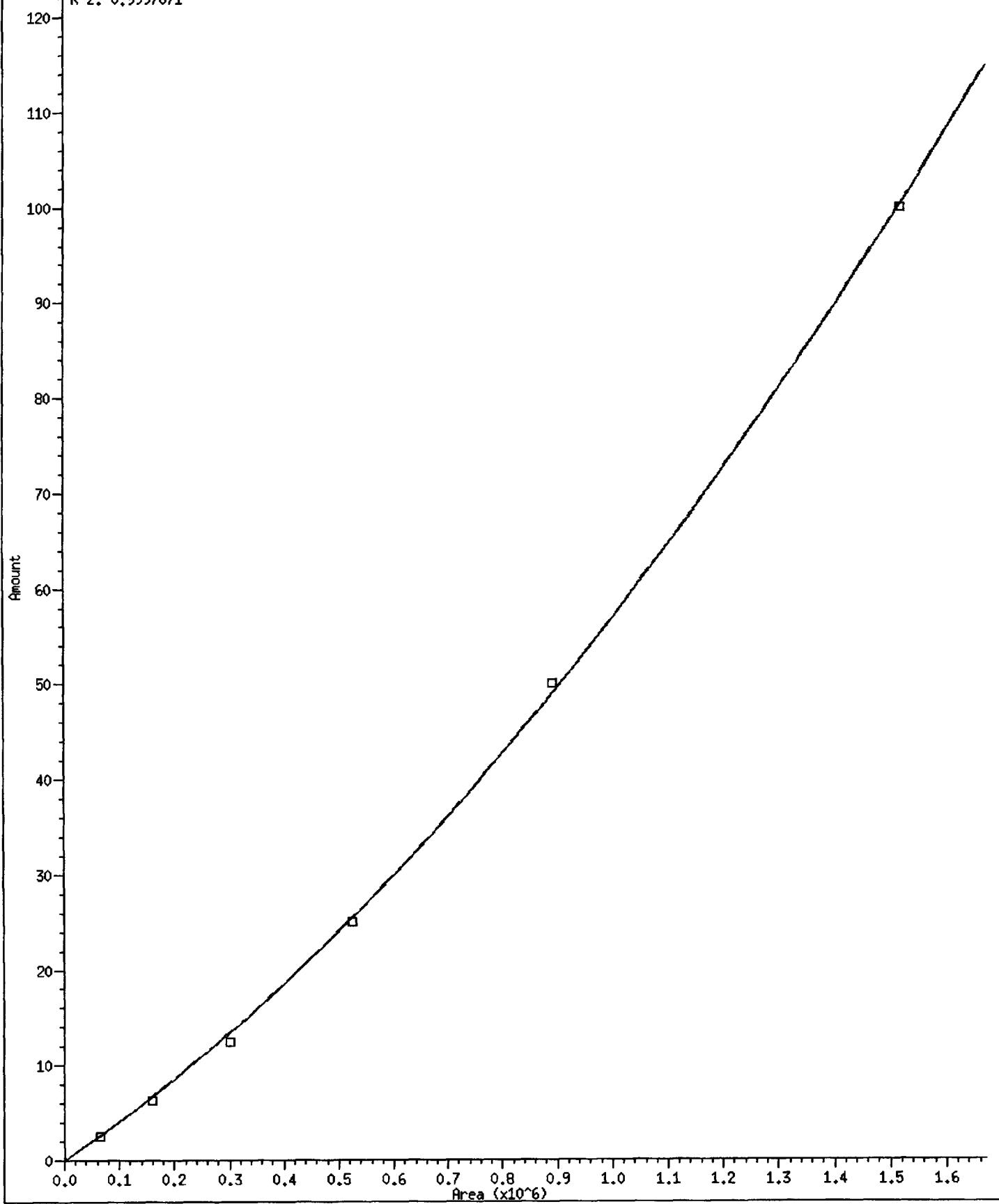


3 2,3,6-Trichlorophenol

Curve Type: Quadratic By-Response

Amt = 0 + 0.00003908481×Rsp + 1.774195e-11×Rsp²

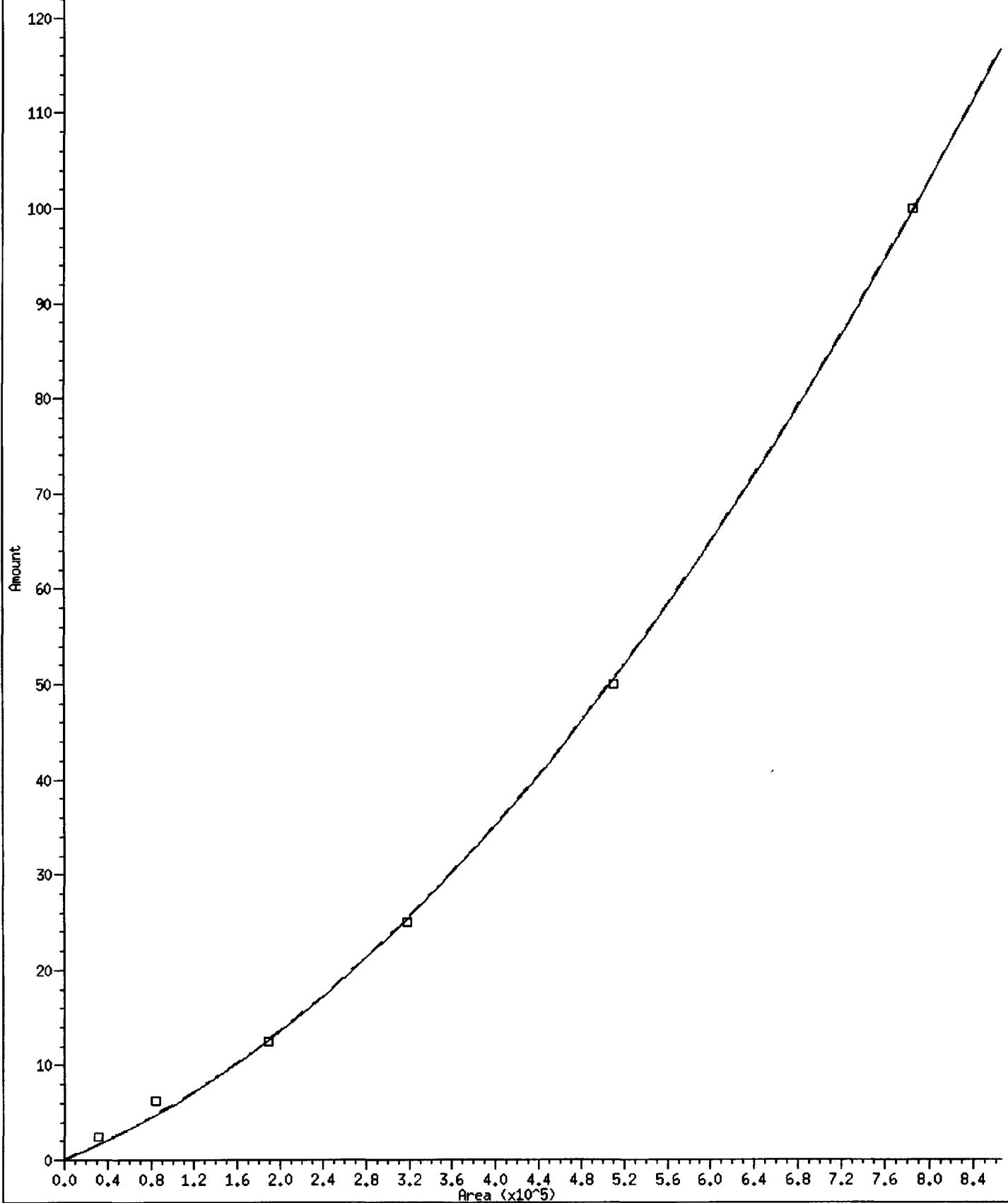
R²: 0.9997071



4 2,4,5-Trichlorophenol

Curve Type: Quadratic By-Response

$$\text{Amt} = 0 + 0.00004747477 \times \text{Rsp} + 1.010754 \times 10^{-10} \times \text{Rsp}^2$$

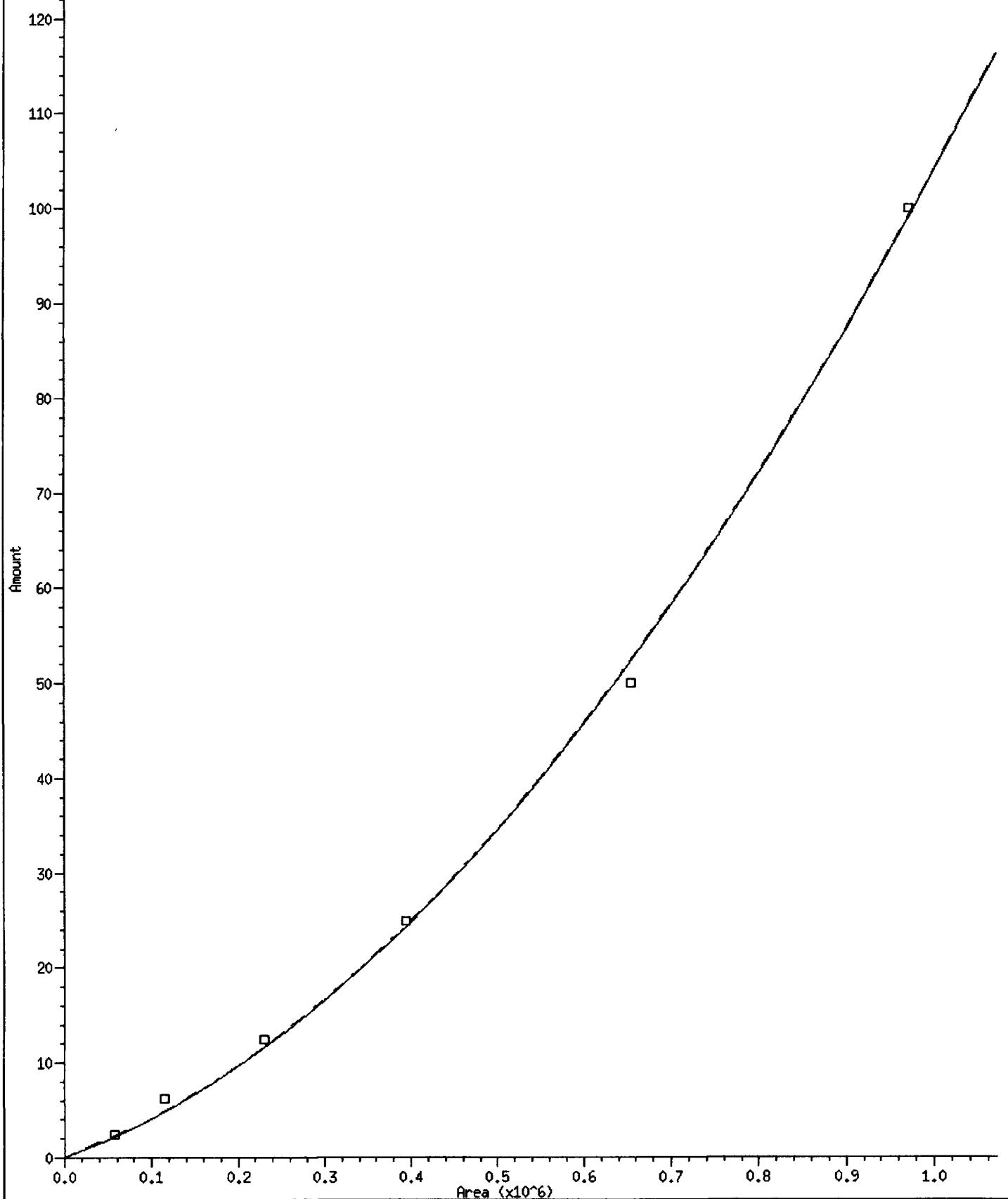
 $R^2: 0.9996490$ 

6 2,3,4-Trichlorophenol

Curve Type: Quadratic By-Response

Amt = 0 + 0.00003425395×Rsp + 6.958576e-11×Rsp²

R²: 0.9988476



Analytical Resources, Inc.
INITIAL CALIBRATION DATA

Start Cal Date : 21-SEP-2012 19:28
 End Cal Date : 21-SEP-2012 22:30
 Quant Method : ESTD
 Origin : Force
 Target Version : 3.50
 Integrator : HP Genie
 Method File : /chem2/ecd1.i/PCP20120921.b/PCPB.m
 Cal Date : 25-Sep-2012 10:11 aron

Calibration File Names:

Level 1: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A018.d
 Level 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A019.d
 Level 3: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A020.d
 Level 4: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A017.d
 Level 5: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A021.d
 Level 6: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A022.d

Compound	2	6	12	25	50	100	Curve	b	Coefficients	%RSD
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6		m1	m2	or R^2
1, 2, 4-Dichlorophenol	32292	93863	14227	242396	430585	711683	QUD	0.000e+00	0.00075	9.197e-10
2, 2, 4, 6-Trichlorophenol	81319	208197	314632	551249	1046878	1768391	QUD	0.000e+00	0.00004	1.143e-11
3, 2, 3, 6-Trichlorophenol	64685	160501	300088	524327	891000	1518833	QUD	0.000e+00	0.00004	1.774e-11
4, 2, 4, 5-Trichlorophenol	31460	84183	188136	317756	509875	786042	QUD	0.000e+00	0.00005	1.011e-10
5, 2, 3, 5, 6-Tetrachlorophenol	33035	31396	31039	28607	26066	23149	AVRG		28886	12.87707
6, 2, 3, 4-Trichlorophenol	57679	115678	23005	394022	654702	972757	QUD	0.000e+00	0.00003	6.959e-11
8, 2, 3, 4, 5-Tetrachlorophenol	25112	23550	21839	20040	18059	15675	AVRG		20723	16.95022
9 Pentachlorophenol	42596	39717	37527	34681	31627	27985	AVRG		35689	15.04595
\$ 7 2, 4, 6-Tribromophenol (surr)	33881	32521	32652	29584	27527	24749	AVRG		30152	11.69494

Analytical Resources, Inc.

INITIAL CALIBRATION DATA

Start Cal Date : 21-SEP-2012 19:28
 End Cal Date : 21-SEP-2012 22:30
 Quant Method : ESTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem2/ecd1.i/PCP20120921.b/PCP.m
 Cal Date : 25-Sep-2012 12:45 aron
 Curve Type : Average

Calibration File Names:

Level 1: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A018.d
 Level 2: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A019.d
 Level 3: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A020.d
 Level 4: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A017.d
 Level 5: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A021.d
 Level 6: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A022.d

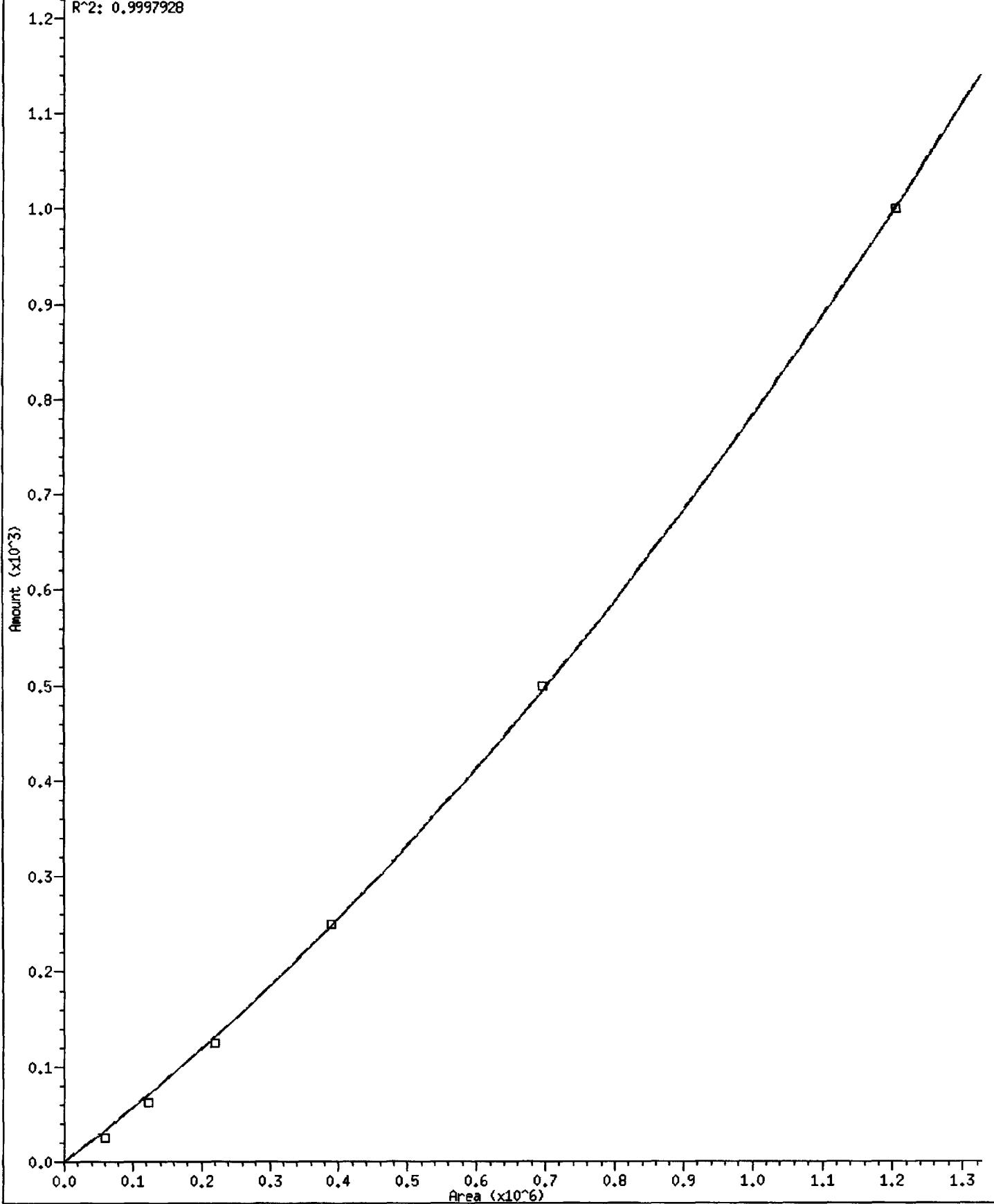
Compound	2.500	6.250	12.500	25.000	50.000	100.000	RRF	* RSD
1 2,4-Dichlorophenol /	2363	1981	1759	1559	1390	1206	1710	24.560 <-
2 2,4,6-Trichlorophenol /	37800	36793	34435	30695	27572	23819	31852	17.232
3 2,3,6-Trichlorophenol	39363	41432	38253	34053	27204	22897	33867	21.756 <-
4 2,4,5-Trichlorophenol	23290	24192	22777	19742	16572	13274	19974	21.611 <-
5 2,3,5,6-Tetrachlorophenol	53827	53395	50427	46244	41856	36207	46993	14.828
6 2,3,4-Trichlorophenol	35575	30905	28044	24569	21009	17260	26227	25.461 <-
8 2,3,4,5-Tetrachlorophenol	50607	45516	41632	36089	31716	27069	38772	22.711 <-
9 Pentachlorophenol	77052	71102	62381	56971	51600	44551	60610	20.016 <-
\$ 7 2,4,6-Tribromophenol (surr)	63570	57559	52456	46332	41843	35975	49623	20.614 <-

1 2,4-Dichlorophenol

Curve Type: Quadratic By-Response

Amt = 0 + 0.0005439634×Rsp + 2.380142e-10×Rsp²

R²: 0.9997928

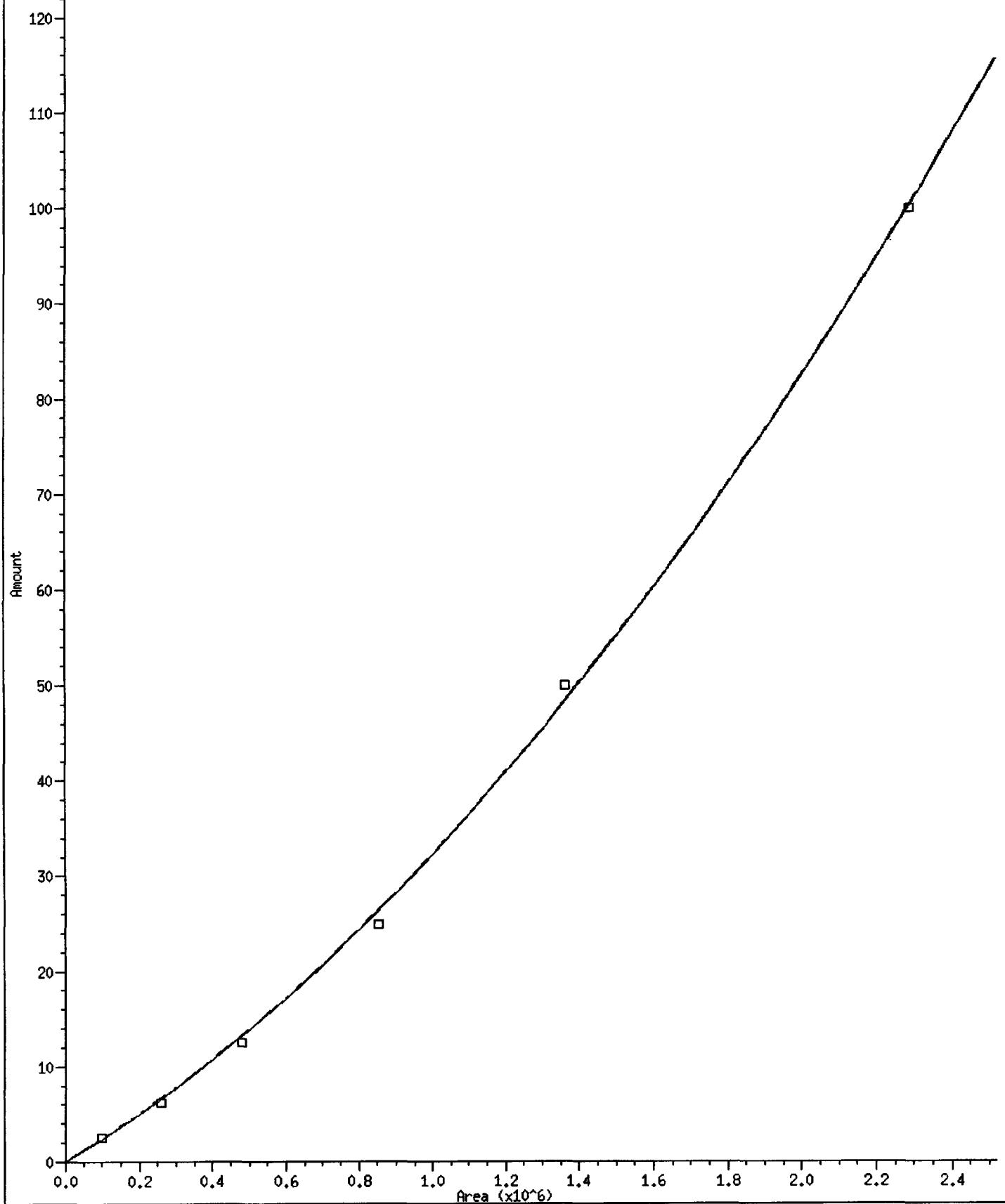


3 2,3,6-Trichlorophenol

Curve Type: Quadratic By-Response

Amt = 0 + 0.00002314655×Rsp + 9.045971e-12×Rsp²

R²: 0.9992332

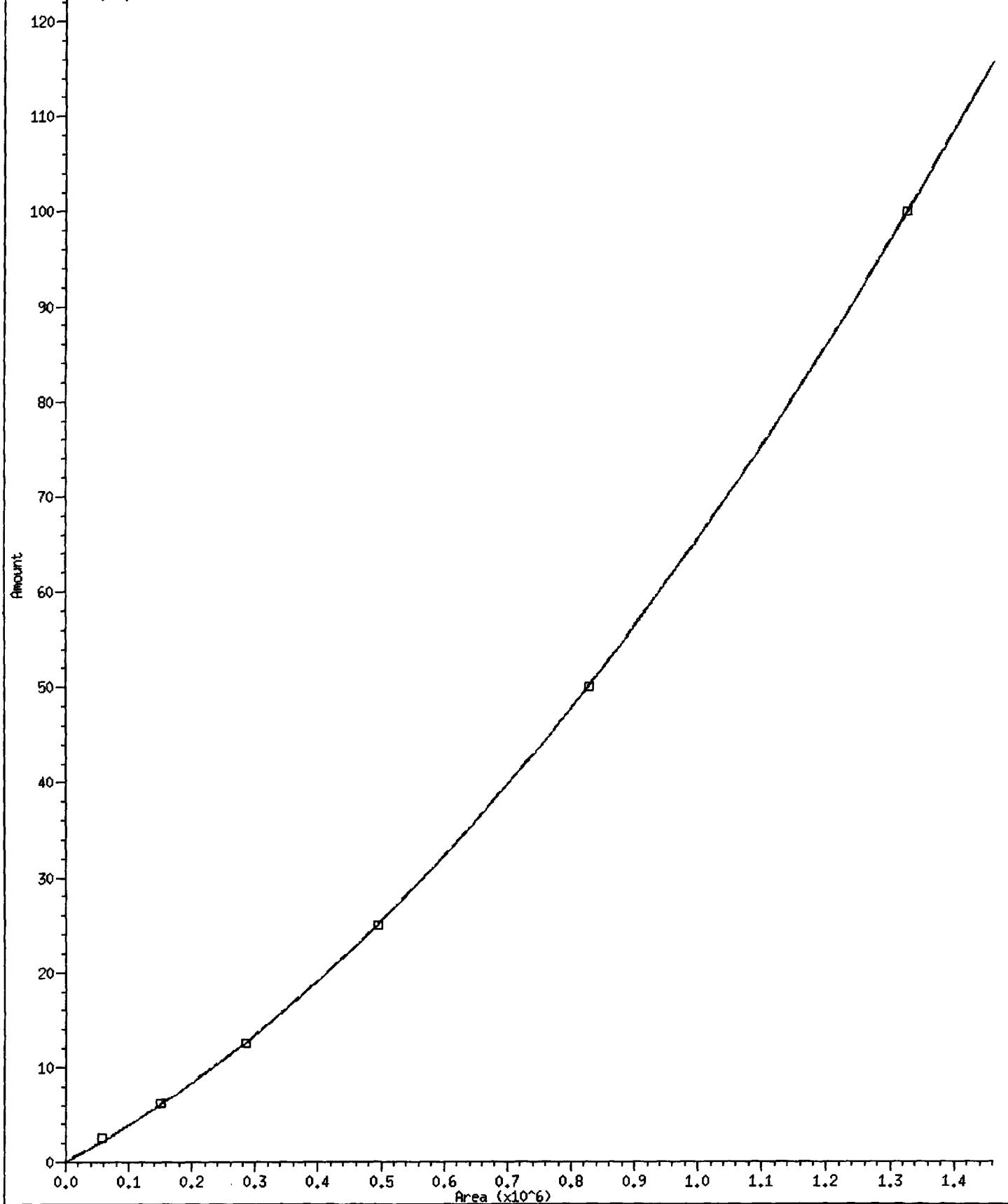


4 2,4,5-Trichlorophenol

Curve Type: Quadratic By-Response

Amt = 0 + 0.00003583434×Rsp + 2.973949e-11×Rsp²

R²: 0.9999836

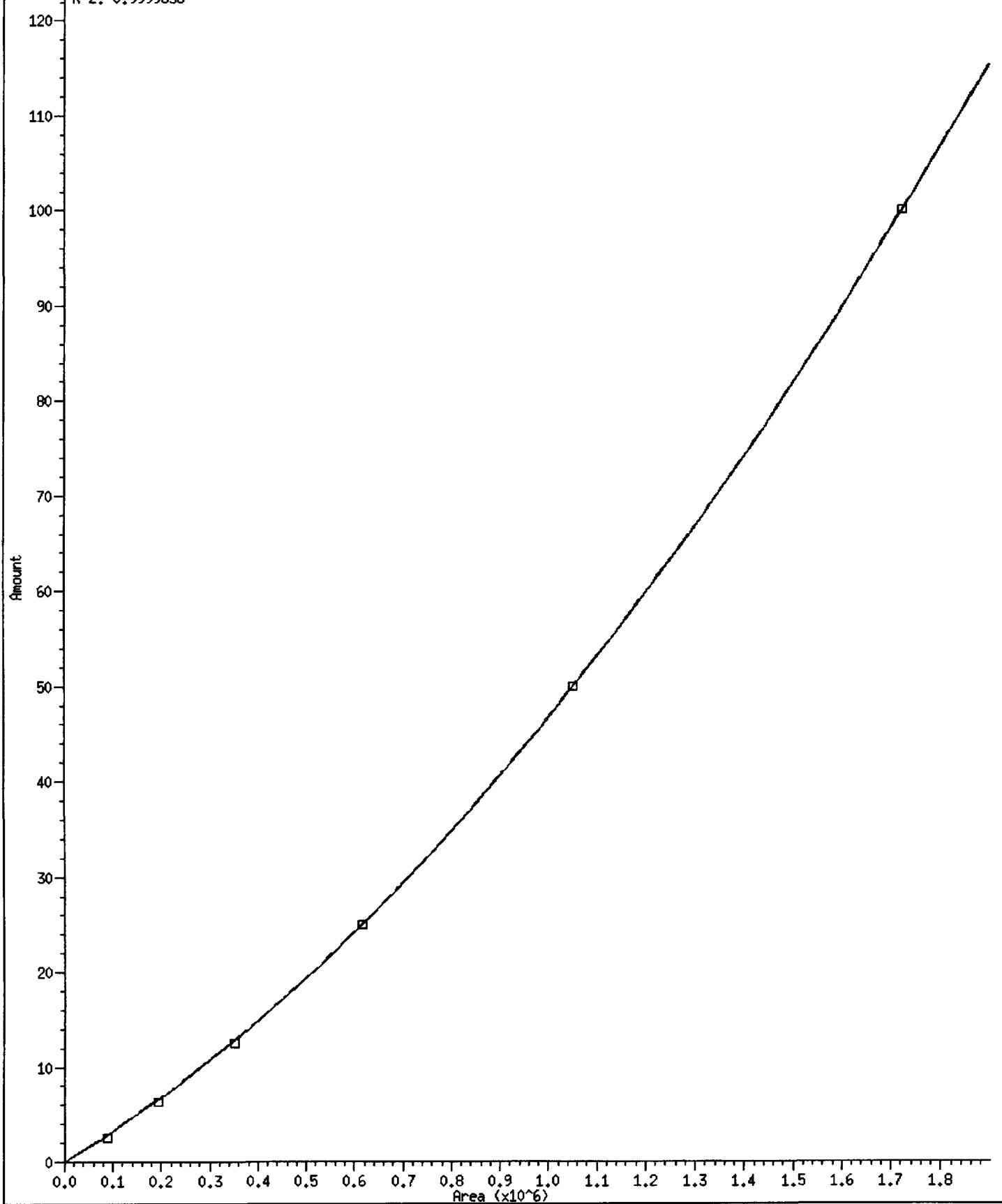


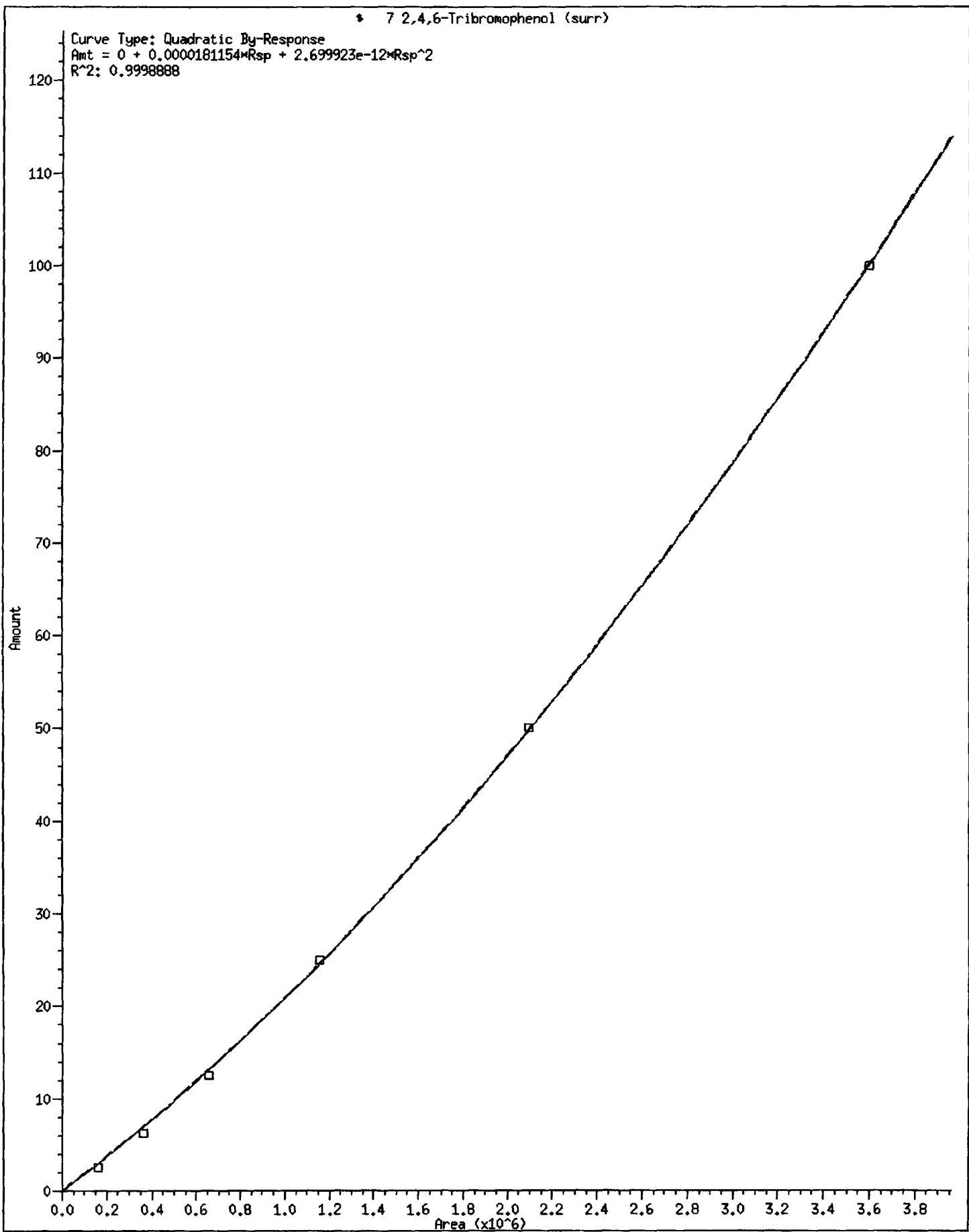
6 2,3,4-Trichlorophenol

Curve Type: Quadratic By-Response

Amt = 0 + 0.00003084114×Rsp + 1.573015e-11×Rsp²

R²: 0.9999656





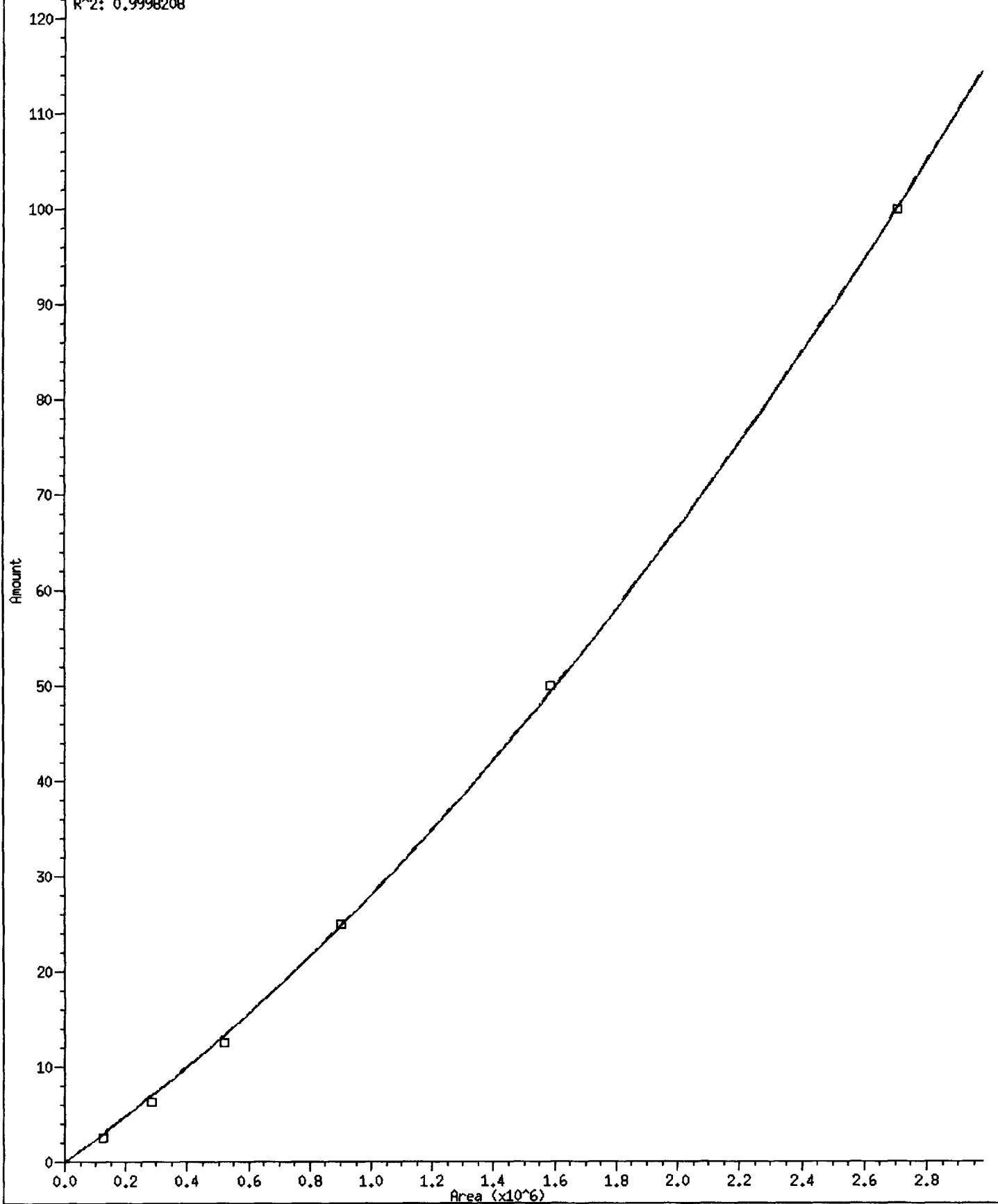
42721600051

8 2,3,4,5-Tetrachlorophenol

Curve Type: Quadratic By-Response

Amt = 0 + 0.00002275485×Rsp + 5.269859e-12×Rsp²

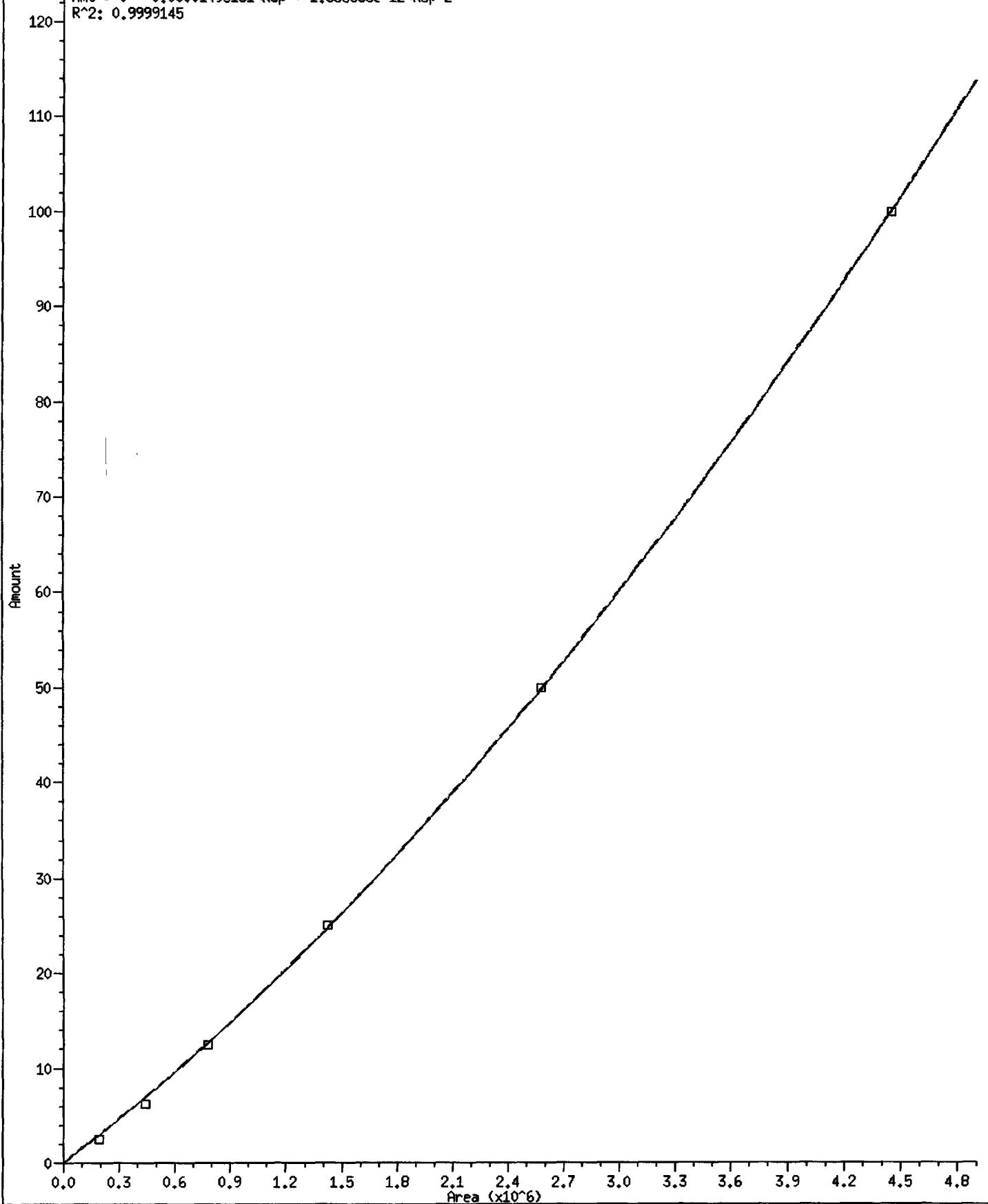
R²: 0.9998208



9 Pentachlorophenol

Curve Type: Quadratic By-Response

$$\text{Amt} = 0 + 0.00001495181 \times \text{Rsp} + 1.686386e-12 \times \text{Rsp}^2$$

 $R^2: 0.9999145$ 

Analytical Resources, Inc.

INITIAL CALIBRATION DATA

Start Cal Date : 21-SEP-2012 19:28
 End Cal Date : 21-SEP-2012 22:30
 Quant Method : ESTD
 Origin : Force
 Target Version : 3.50
 Integrator : HP Genie
 Method file : /chem2/ecd1.i/PCP20120921.b/PCP.m
 Cal Date : 25-Sep-2012 12:45 aron

Calibration File Names:

Level 1: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A018.d
 Level 2: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A019.d
 Level 3: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A020.d
 Level 4: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A017.d
 Level 5: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A021.d
 Level 6: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A022.d

Compound	2	6	12	25	50	100		Coefficients			%RSD or R^2
	Level 1	Level 2	Level 3	Level 4	Level 5	Level 6	Curve	b	m1	m2	
1 2,4-Dichlorophenol	59080	123823	219893	389692	695219	1205949	QUAD	0.000e+00	0.00054	2.380e-10	0.99979
2 2,4,6-Trichlorophenol	37800	36793	34435	30695	27572	23819	AVRG		31852		17.23219
3 2,3,6-Trichlorophenol	98407	258950	478158	851316	1360198	2289663	QUAD	0.000e+00	0.00002	9.046e-12	0.99923
4 2,4,5-Trichlorophenol	58226	151202	284711	493541	828590	1327374	QUAD	0.000e+00	0.00004	2.974e-11	0.99998
5 2,3,5,6-Tetrachlorophenol	53827	53395	50427	46244	41856	36207	AVRG		46993		14.82781
6 2,3,4-Trichlorophenol	88937	193157	350554	614215	1050439	1725987	QUAD	0.000e+00	0.00003	1.573e-11	0.99997
8 2,3,4,5-Tetrachlorophenol	126517	284477	520404	902216	1585782	2706943	QUAD	0.000e+00	0.00002	5.270e-12	0.99982
9 Pentachlorophenol	192630	444389	779765	1424284	2579984	4455069	QUAD	0.000e+00	0.00001	1.686e-12	0.99991
\$ 7 2,4,6-Tribromophenol (surr)	158926	359745	655694	1158311	2092155	3597466	QUAD	0.000e+00	0.00002	2.700e-12	0.99989

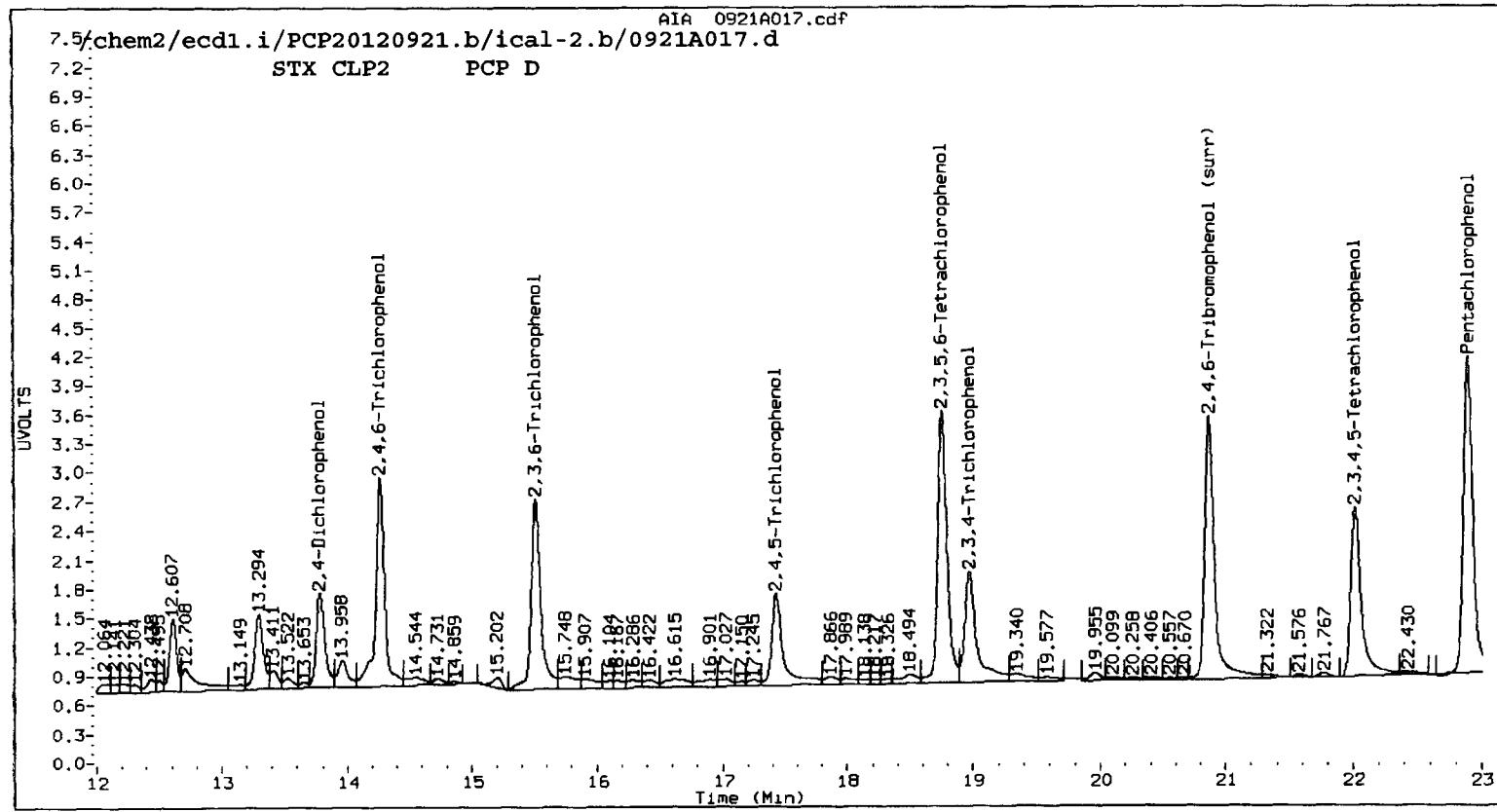
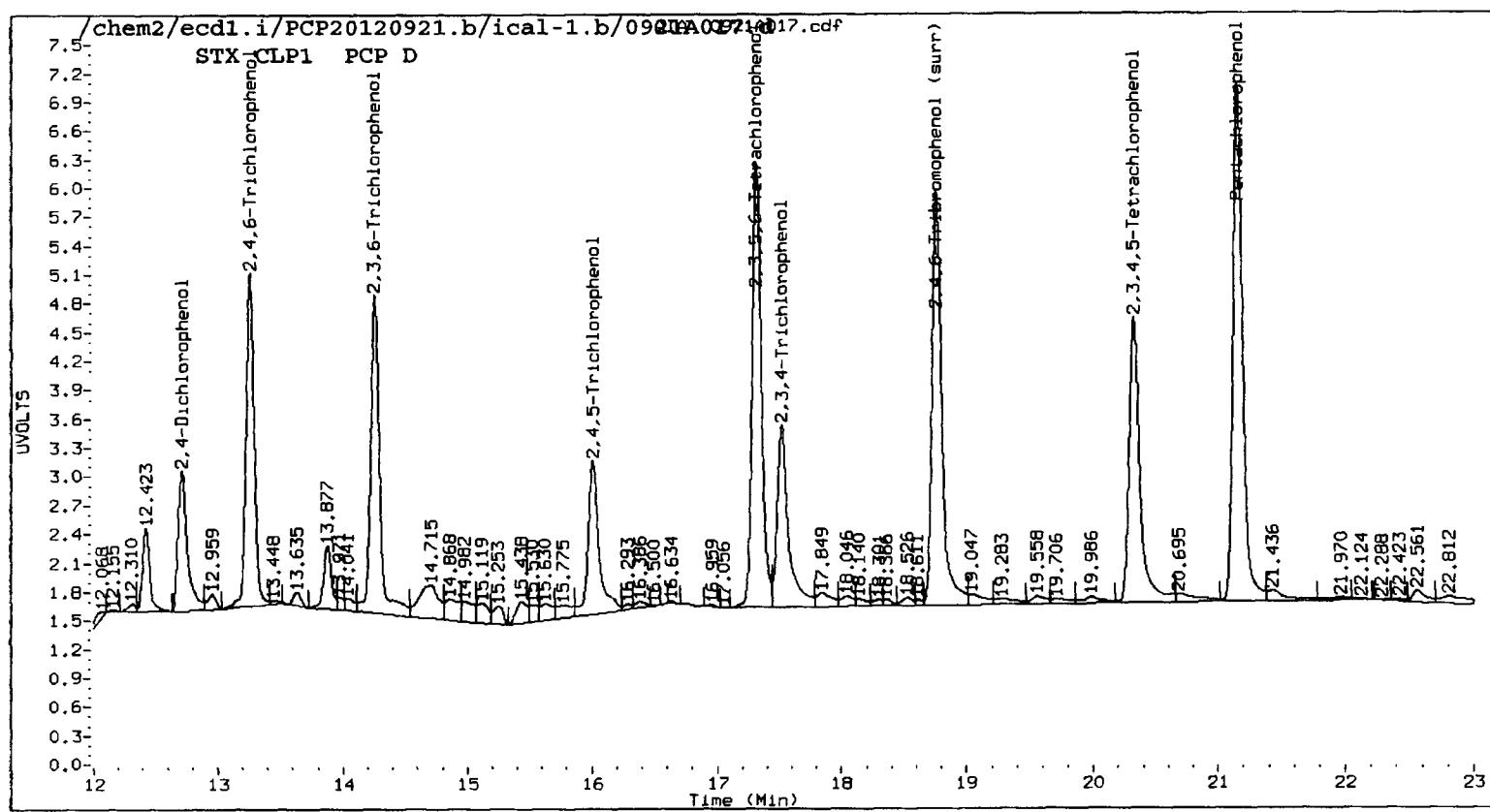
Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Data file 1: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A017.d ARI ID: PCP D
 Data file 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A017.d Client ID:
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 21-SEP-2012 19:28
 Compound Sublist: all Report Date: 09/25/2012 13:17
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

RT	STX CLP1 Col		STX CLP2 Col		on col	STX CLP1	STX CLP2	Compound
	Shift	Response	RT	Shift	Response			
21.158	0.000	1424284	22.888	0.000	867017	24.7166	24.2938	1.7 Pentachlorophenol
13.256	0.000	767382	14.264	0.000	551249	24.0918	23.4510	2.7 2,4,6-Trichlorophenol
14.254	0.000	851316	15.507	0.000	524327	26.2610	25.3708	3.4 2,3,6-Trichlorophenol
16.010	0.000	493541	17.423	0.000	317756	24.9297	25.2909	1.4 2,4,5-Trichlorophenol
17.519	0.000	614215	18.968	0.000	394022	24.8774	24.3002	2.3 2,3,4-Trichlorophenol
17.315	0.000	1156092	18.752	0.000	715183	24.6014	24.7592	0.6 2,3,5,6-Tetrachloroph
20.323	0.000	902216	22.012	0.000	501003	24.8194	24.1766	2.6 2,3,4,5-Tetrachlorophe
12.711	0.000	389692	13.778	0.000	242396	248.1230	235.7181	5.1 2,4-Dichlorophenol
18.763	0.000	1158311	20.867	0.000	739598	24.6	24.5	0.3 2,4,6-Tribromophenol (

PERCENT RECOVERY

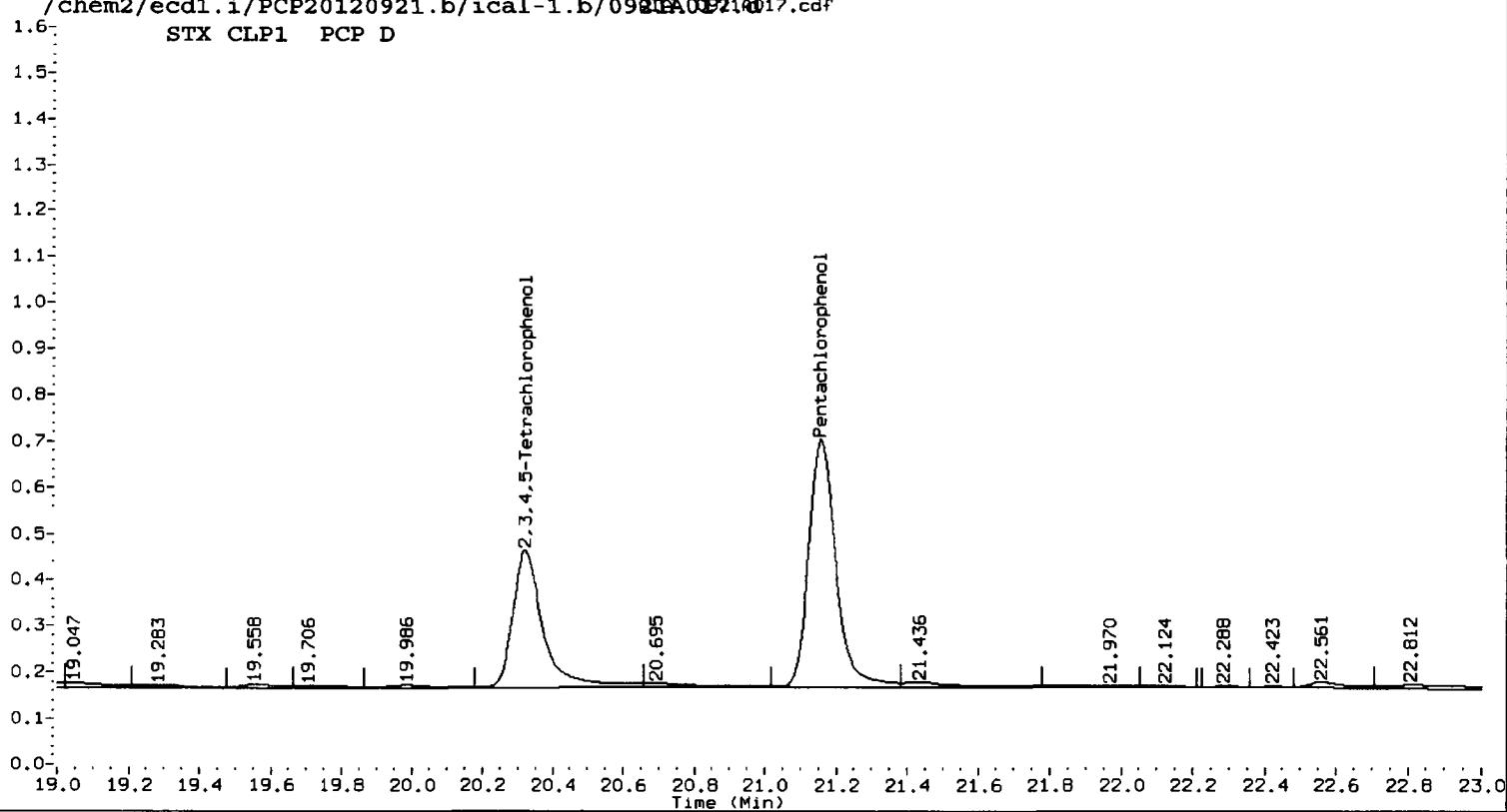
COMPOUND	Col1	Col2
2,4,6-TBP (surr)	98.4	98.1



/chem2/ecdl.i/PCP20120921.b/ical-1.b/0921A017.cdf

STX CLP1 PCP D

UVOLTS

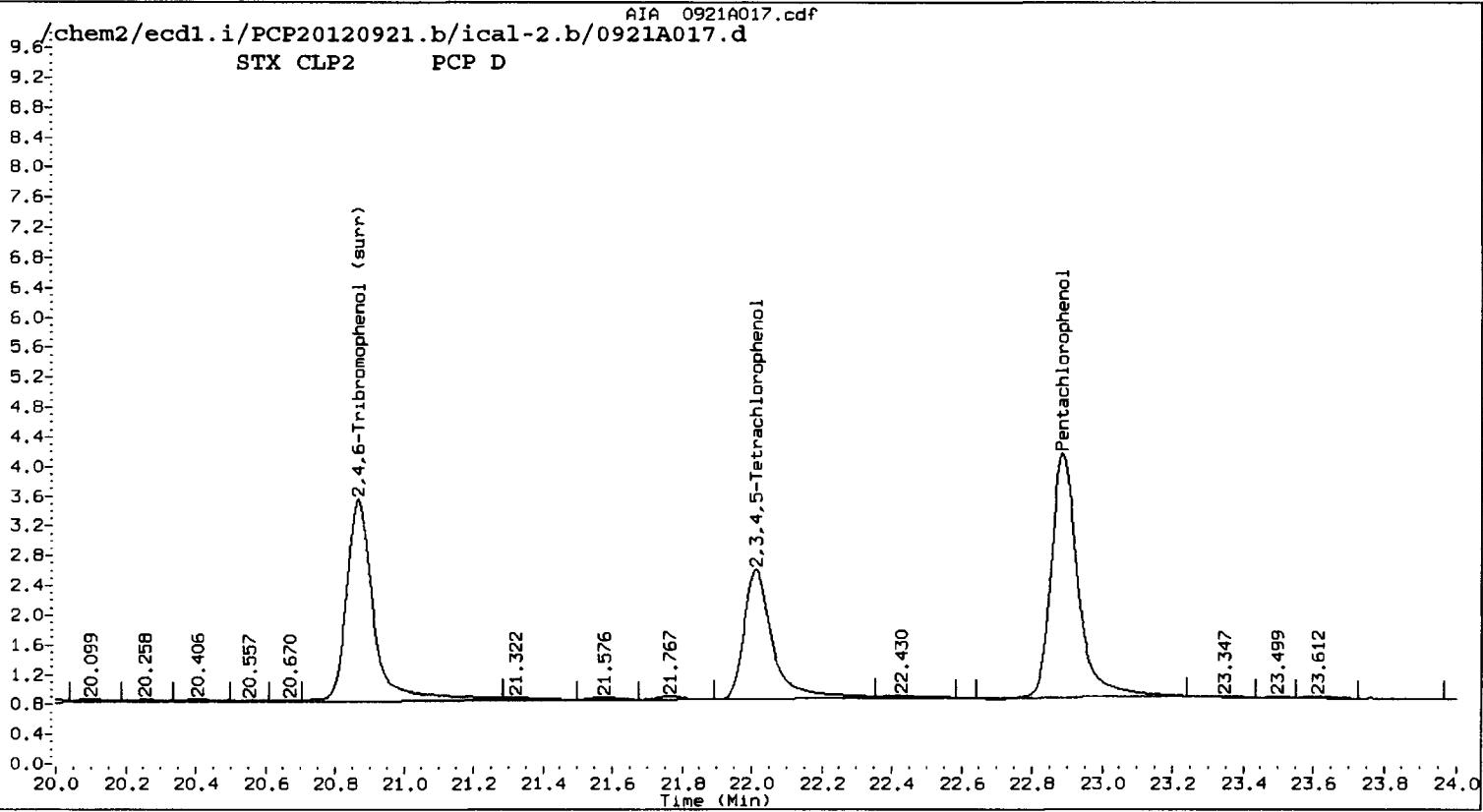


AIA 0921A017.cdf

/chem2/ecdl.i/PCP20120921.b/ical-2.b/0921A017.d

STX CLP2 PCP D

UVOLTS



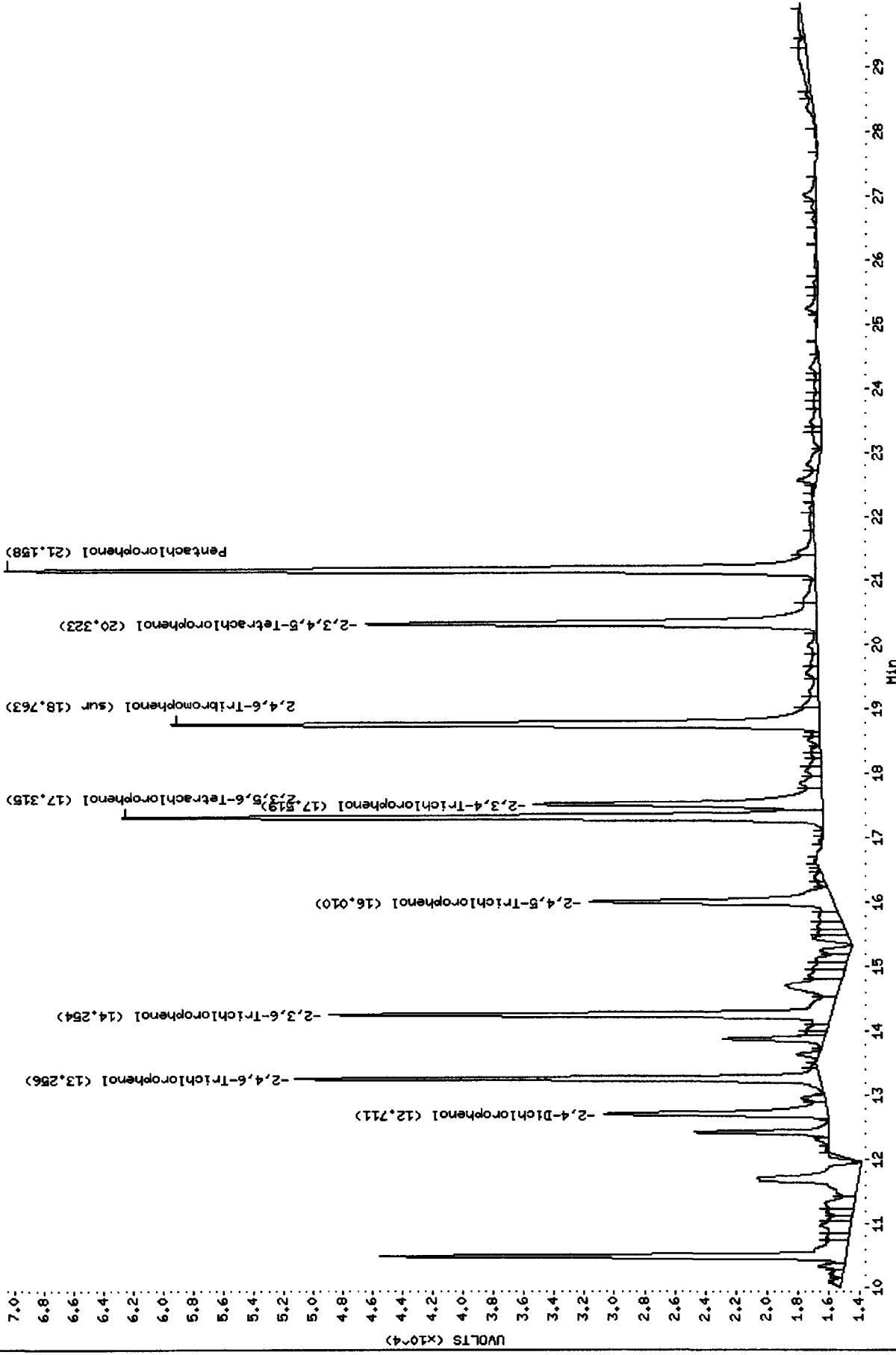
Data File: /chem2/ecd1.i /PCP20120921.b/ical-1.b/0921A017.d
Date : 21-SEP-2012 19:28

Client ID:
Sample Info: PCP D
Purge Volume: 500.0
Column Phase: STX CLP1

Instrument: ecd1.i

Operator: ar
Column diameter: 0.53

/chem2/ecd1.i /PCP20120921.b/ical-1.b/0921A017.d/0921A017.cdf



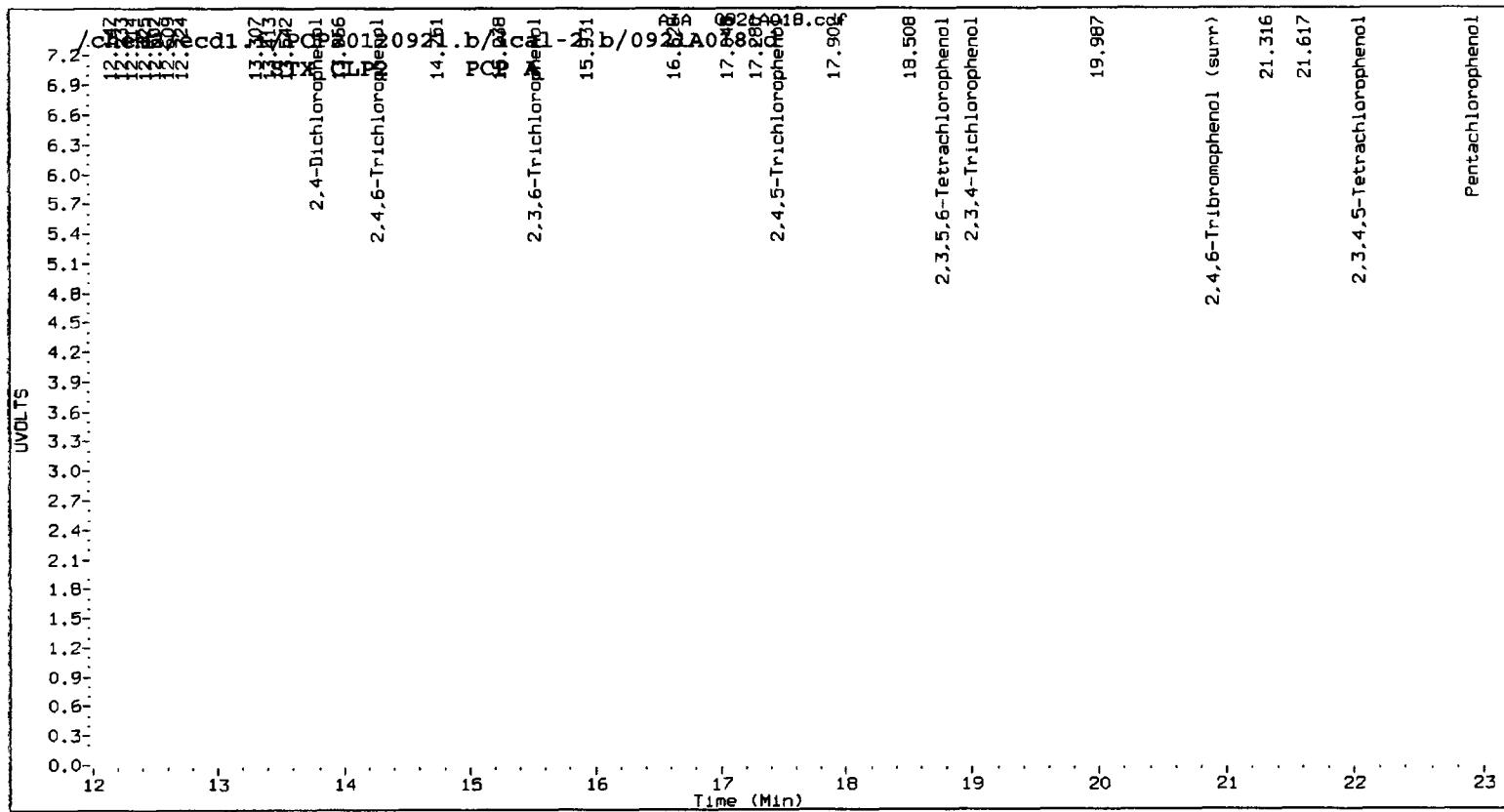
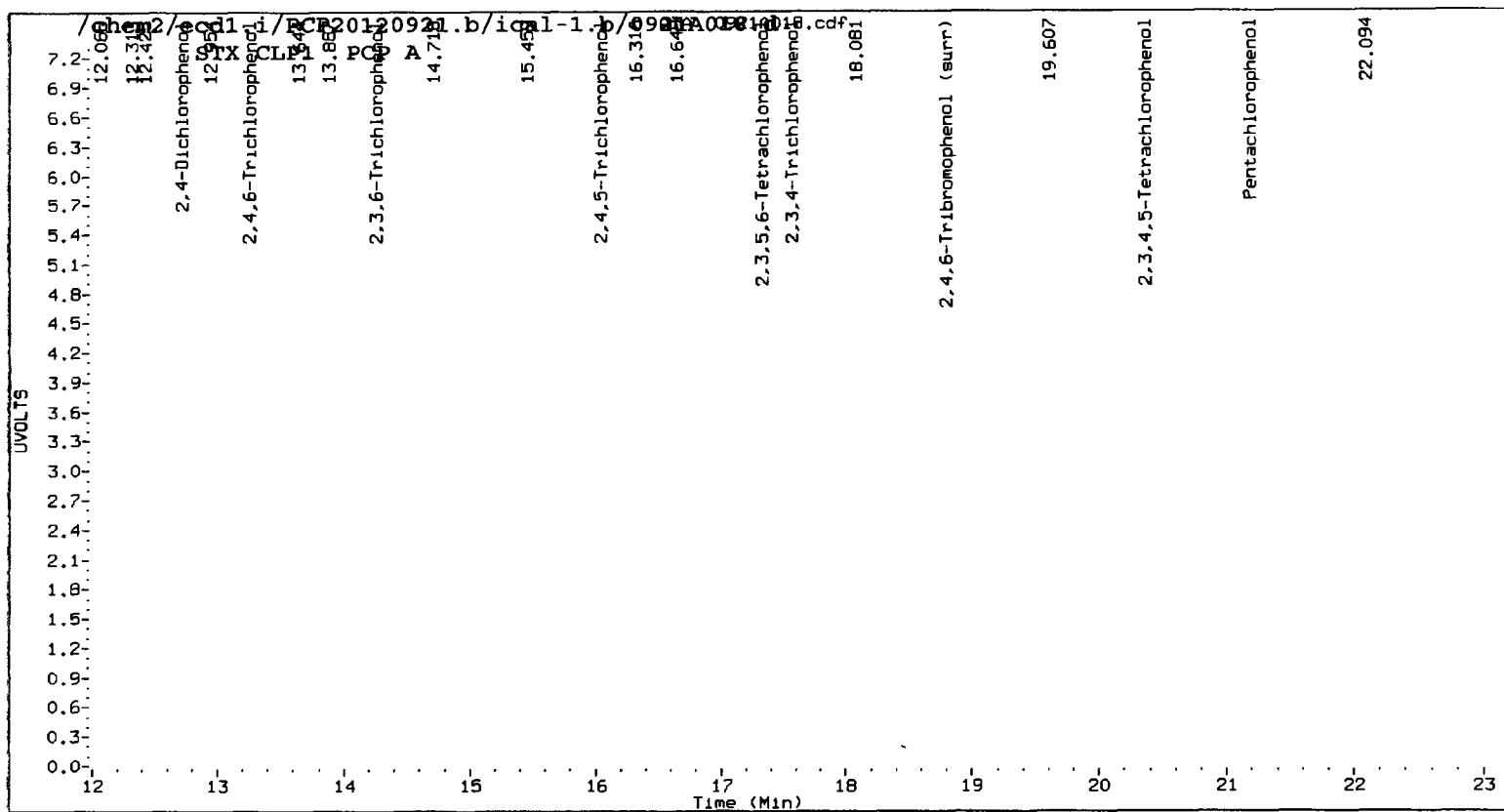
Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

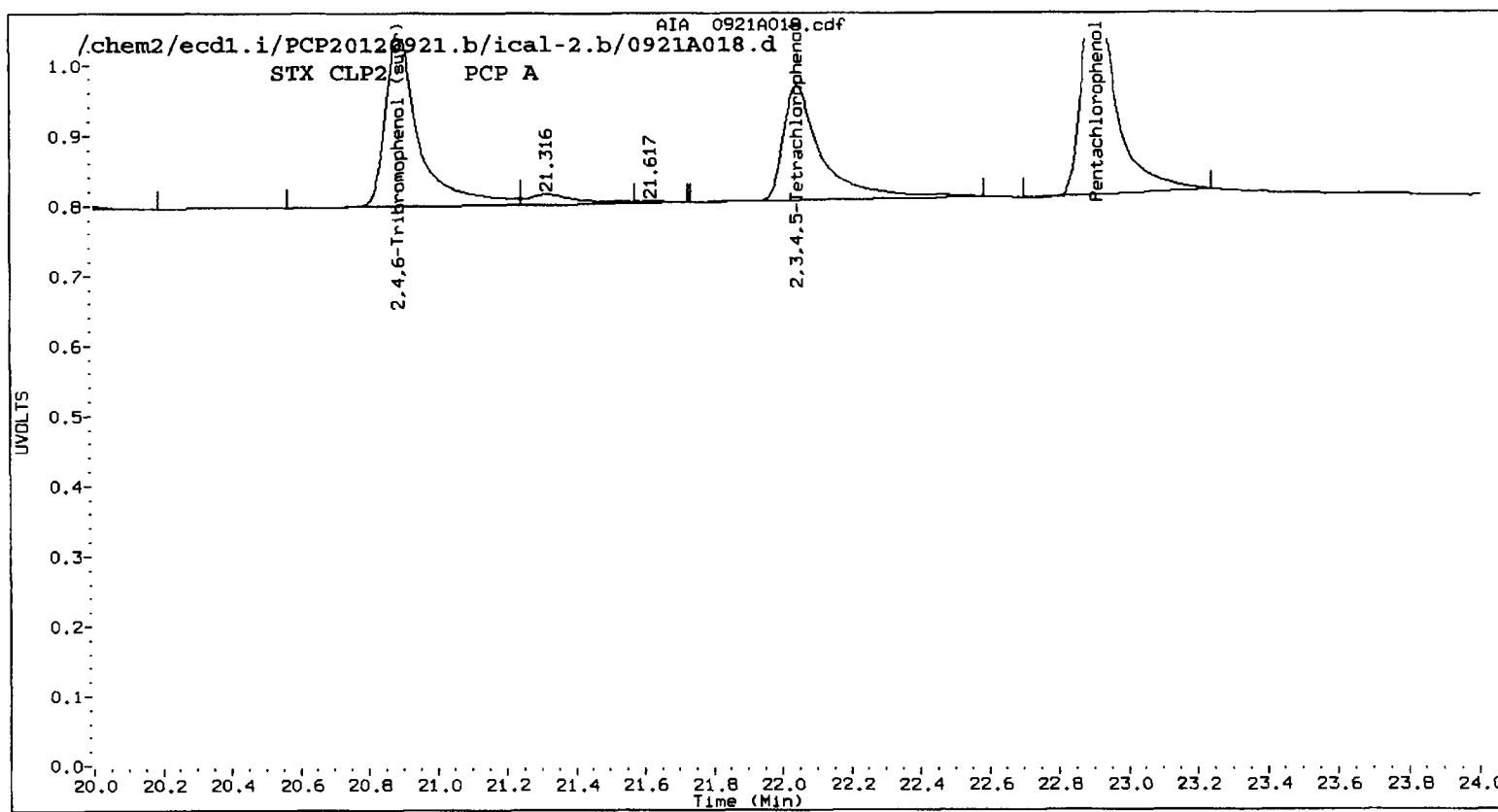
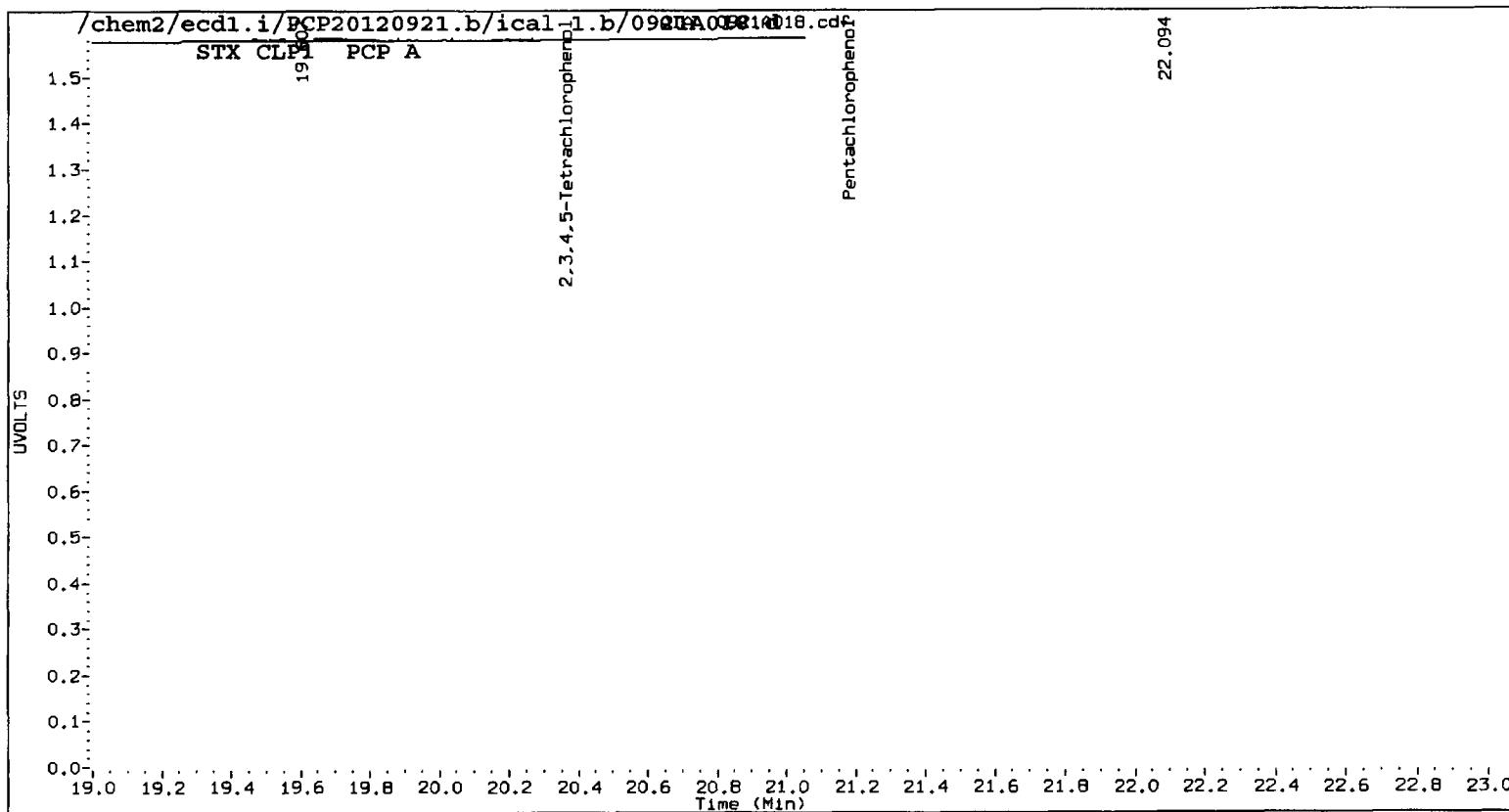
Data file 1: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A018.d ARI ID: PCP A
 Data file 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A018.d Client ID:
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 21-SEP-2012 20:04
 Compound Sublist: all Report Date: 09/25/2012 13:17
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

RT	STX CLP1 Col			STX CLP2 Col			STX CLP1 on col	STX CLP2 on col	RPD	Compound
	Shift	Response	RT	Shift	Response	RT				
21.182	0.024	192630	22.906	0.018	106490	2.9427	2.9838	1.4	Pentachlorophenol	
13.260	0.004	94500	14.268	0.004	81319	2.9668	3.0205	1.8	2,4,6-Trichlorophenol	
14.262	0.008	98407	15.514	0.007	64685	2.3654	2.6024	9.5	2,3,6-Trichlorophenol	
16.044	0.034	58226	17.448	0.025	31460	2.1873	1.5936	31.4	2,4,5-Trichlorophenol	
17.566	0.047	88937	19.001	0.033	57679	2.8673	2.2072	26.0	2,3,4-Trichlorophenol	
17.334	0.019	134568	18.767	0.015	82588	2.8636	2.8591	0.2	2,3,5,6-Tetrachlorophene	
20.368	0.045	126517	22.044	0.032	62781	2.9632	3.0296	2.2	2,3,4,5-Tetrachlorophene	
12.719	0.008	59080	13.787	0.009	32292	32.9681	25.1628	26.9	2,4-Dichlorophenol	
18.793	0.030	158926	20.887	0.020	84702	2.9	2.8	4.8	2,4,6-Tribromophenol (s)	

PERCENT RECOVERY

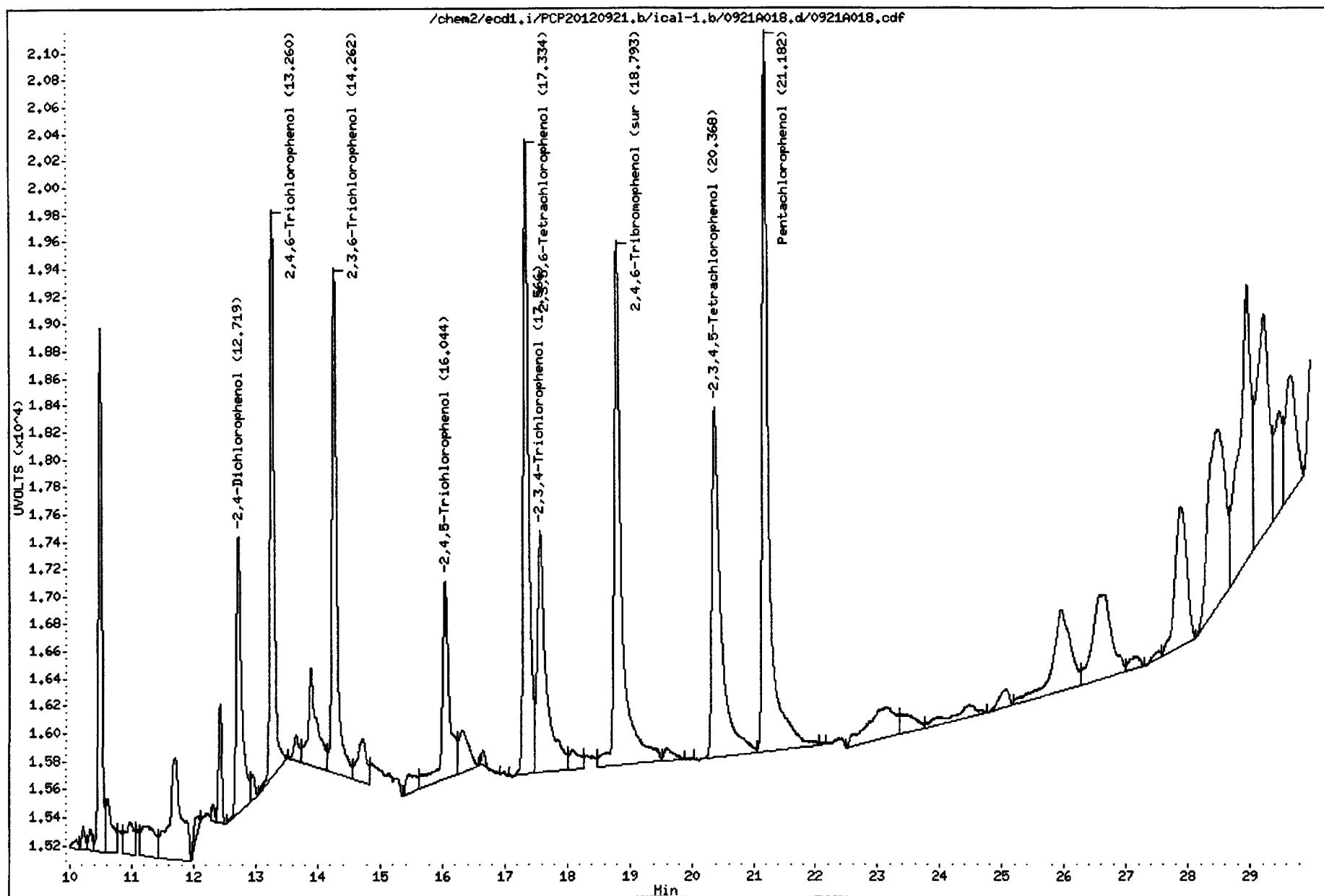
COMPOUND	Col1	Col2
2,4,6-TBP (surr)	11.8	11.2





Data File: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A018.d
Date : 21-SEP-2012 20:04
Client ID:
Sample Info: PCP A
Purge Volume: 500.0
Column phase: STX CLP1

Instrument: ecd1.i
Operator: ar
Column diameter: 0.53
Page 1



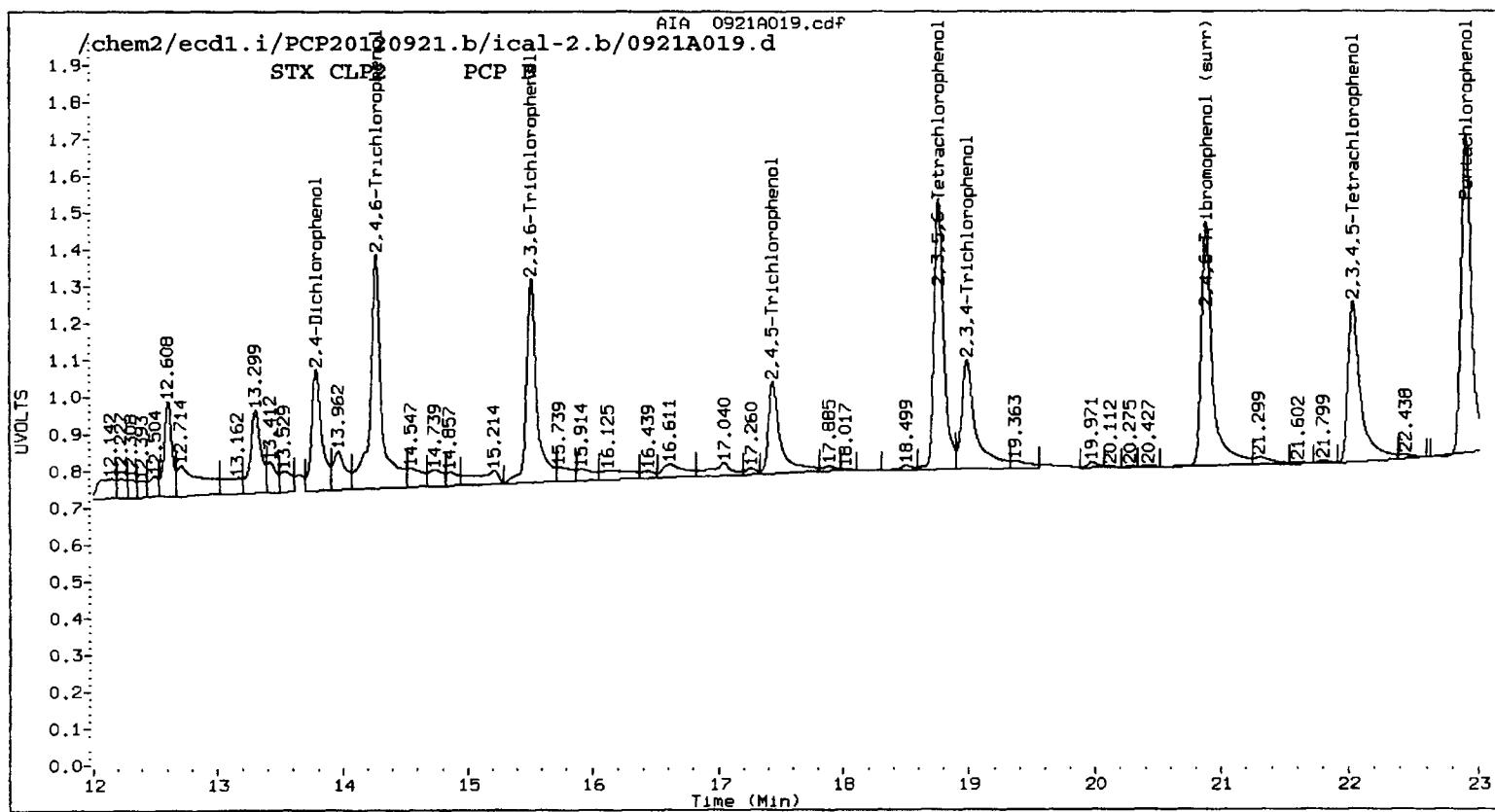
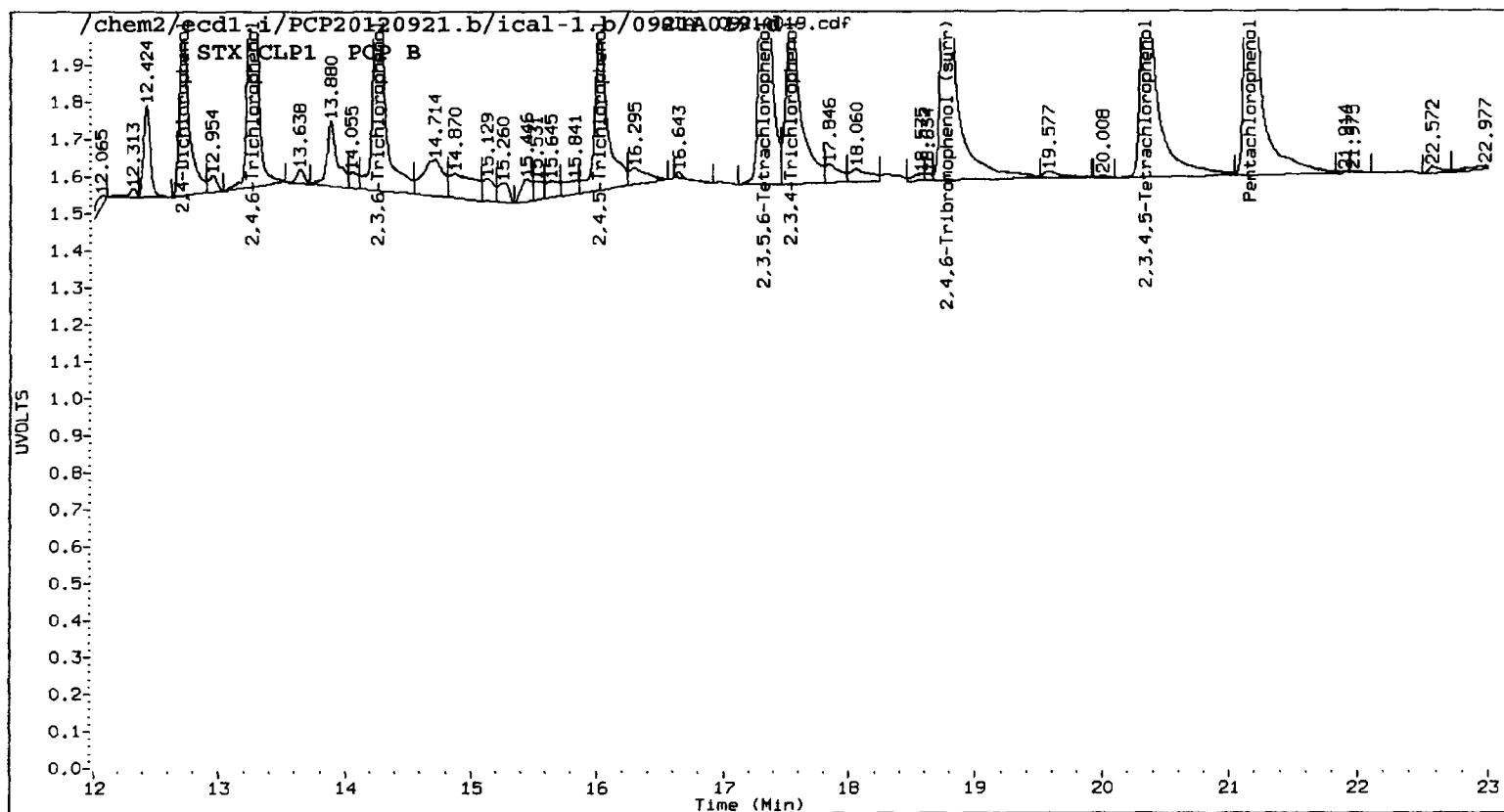
Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

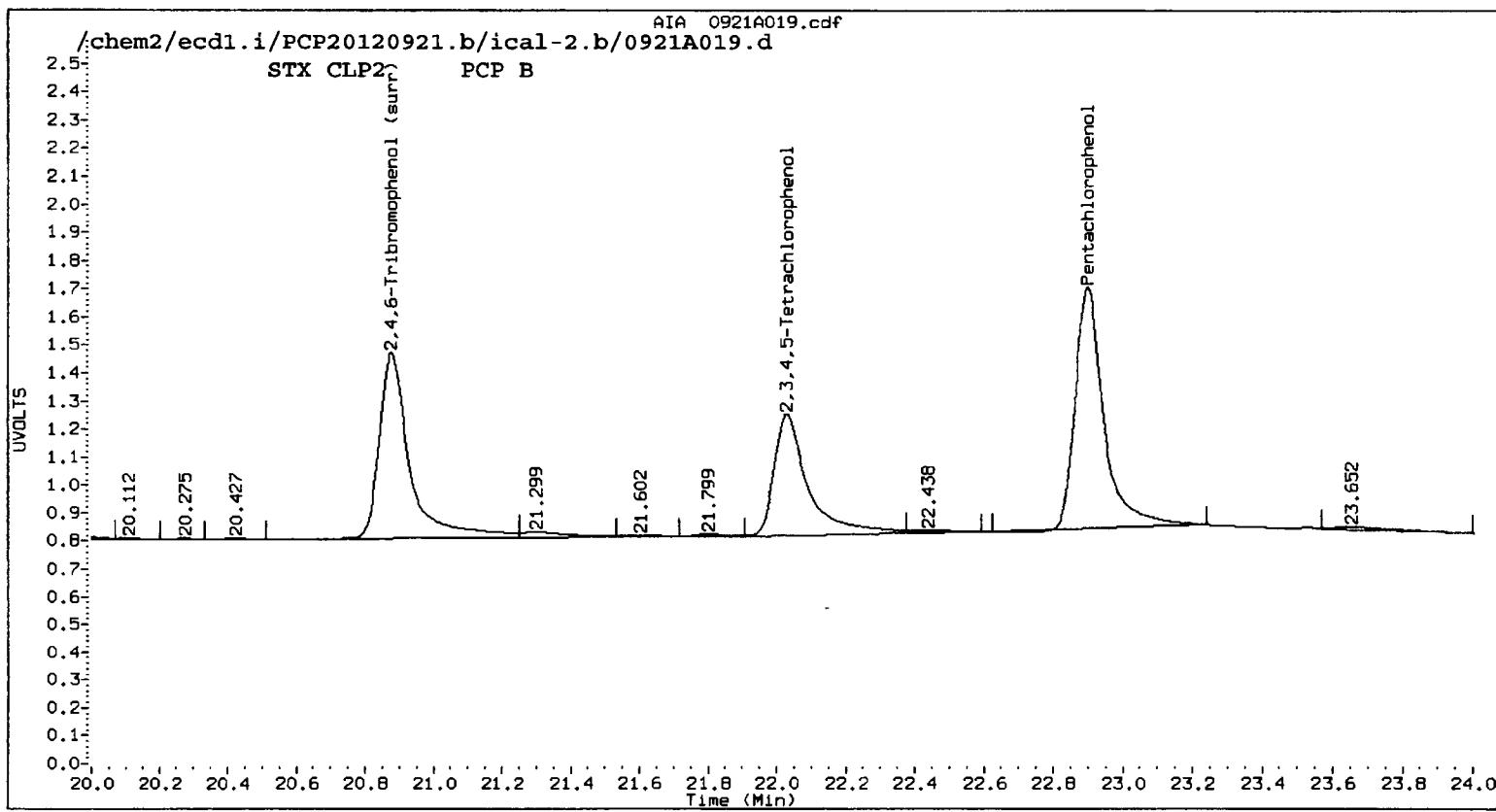
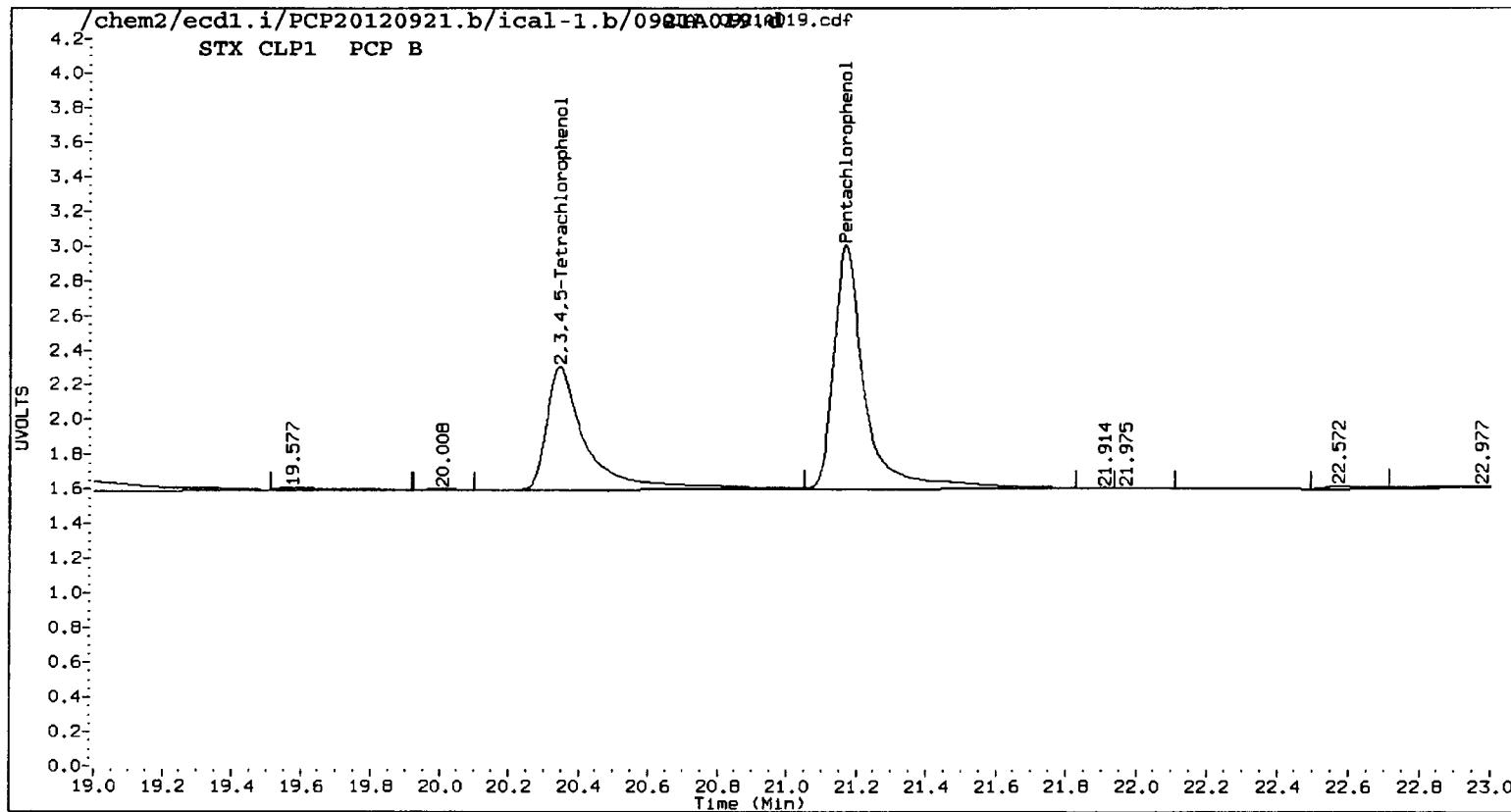
Data file 1: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A019.d ARI ID: PCP B
 Data file 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A019.d Client ID:
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 21-SEP-2012 20:41
 Compound Sublist: all Report Date: 09/25/2012 13:17
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

RT	STX CLP1 Col		STX CLP2 Col		on col	on col	RPD	STX CLP2 Compound
	Shift	Response	RT	Shift	Response			
21.170	0.012	444389	22.897	0.009	248232	6.9775	6.9555	0.3 Pentachlorophenol
13.257	0.001	229955	14.266	0.002	208197	7.2194	8.0367	10.7 2,4,6-Trichlorophenol
14.258	0.004	258950	15.510	0.003	160501	6.6004	6.7302	1.9 2,3,6-Trichlorophenol
16.025	0.015	151202	17.434	0.011	84183	6.0981	4.7129	25.6 2,4,5-Trichlorophenol
17.541	0.022	193157	18.984	0.016	115678	6.5441	4.8936	28.9 2,3,4-Trichlorophenol
17.323	0.008	333721	18.759	0.007	196228	7.1015	6.7933	4.4 2,3,5,6-Tetrachlorophene
20.346	0.023	284477	22.029	0.017	147190	6.8997	7.1029	2.9 2,3,4,5-Tetrachlorophene
12.716	0.005	123823	13.782	0.004	93863	71.0044	78.4554	10.0 2,4-Dichlorophenol
18.776	0.013	359745	20.877	0.010	203256	6.9	6.7	1.8 2,4,6-Tribromophenol (s)

PERCENT RECOVERY

COMPOUND	Col1	Col2
2,4,6-TBP (surr)	27.5	27.0





Date : 21-SEP-2012 20:41

Client ID:

Sample Info: PCP_B

Purge Volume: 500.0

Column Phase: STX CLP1

Instrument: eddi.i

Operator: ar

Column diameter: 0.53

/chem2/eddi.i/PCP20120921.b/ical-1.b/0924A019.d/0924A019.cdf

Permethropheanol (21.170)

2,3,4,5-Tetrachlorophenol (20.346)

2,4,6-Tribromophenol (sur (18.776))

2,3,4-Trichlorophenol (17.323)

2,3,6-Tetrachlorophenol (17.343)

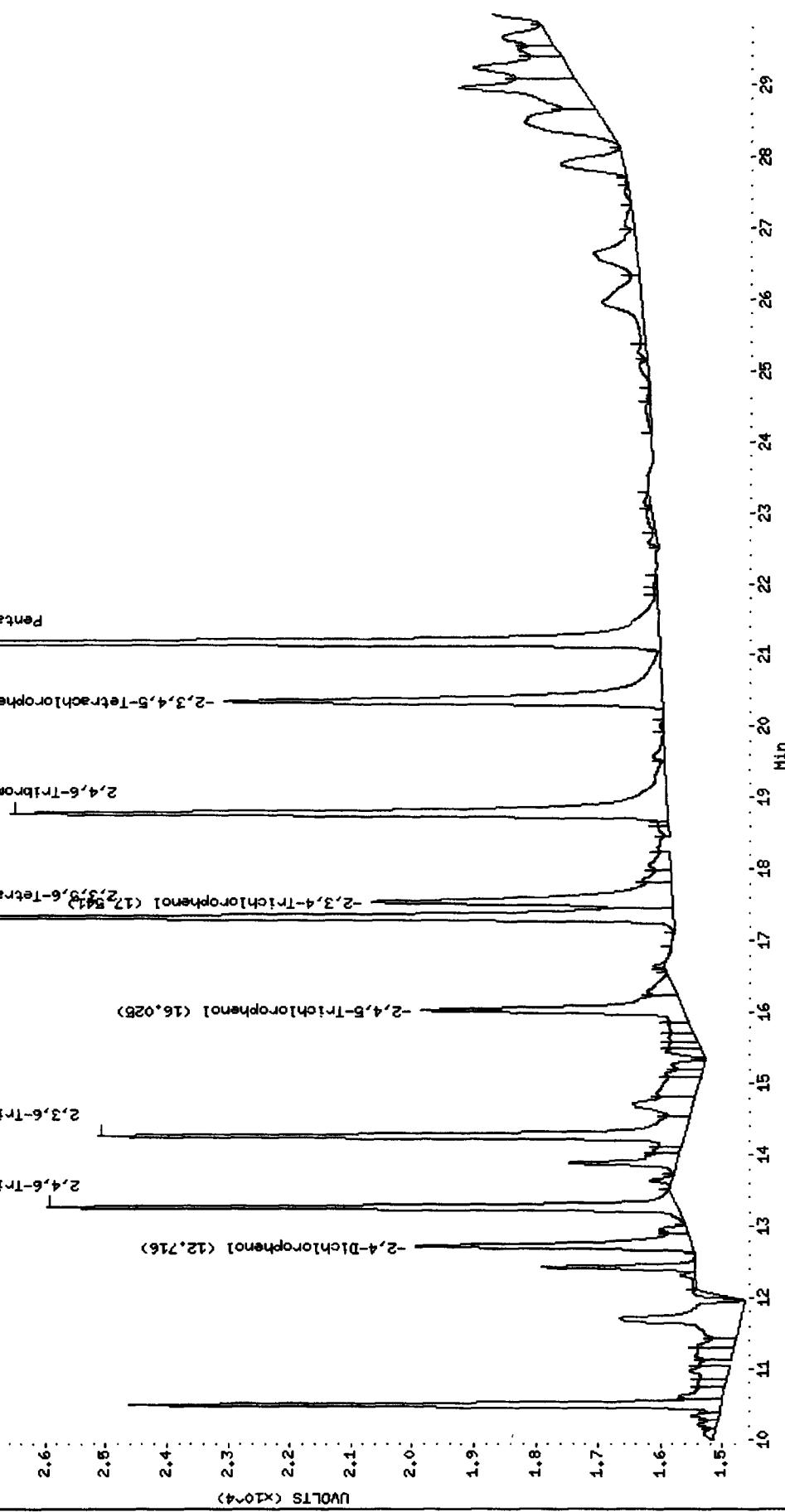
2,4,5-Trichlorophenol (16.025)

2,3,6-Trichlorophenol (14.258)

2,4,6-Trichlorophenol (13.257)

2,4-Dichlorophenol (12.716)

UVOLTS (x10⁻⁴)



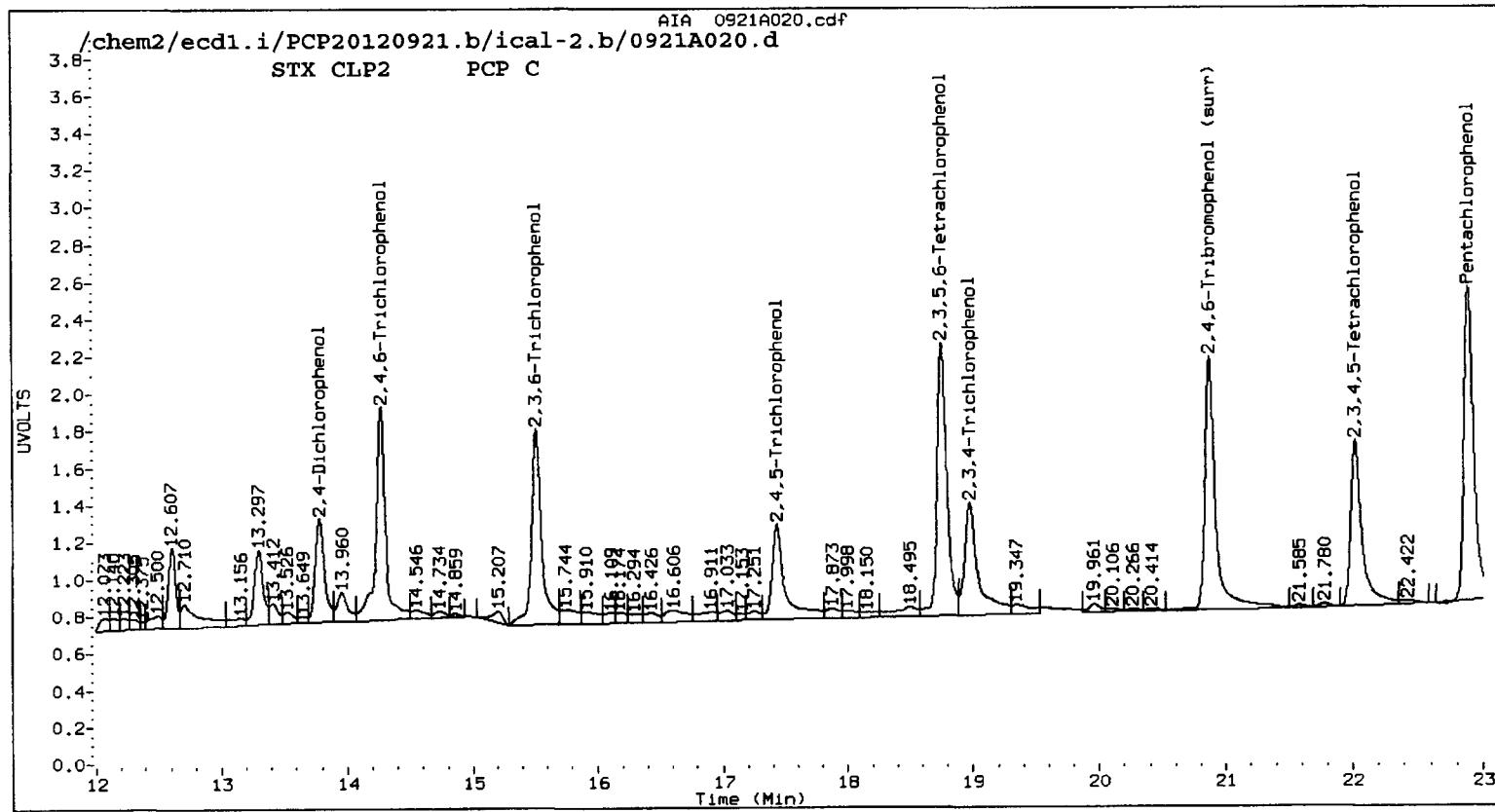
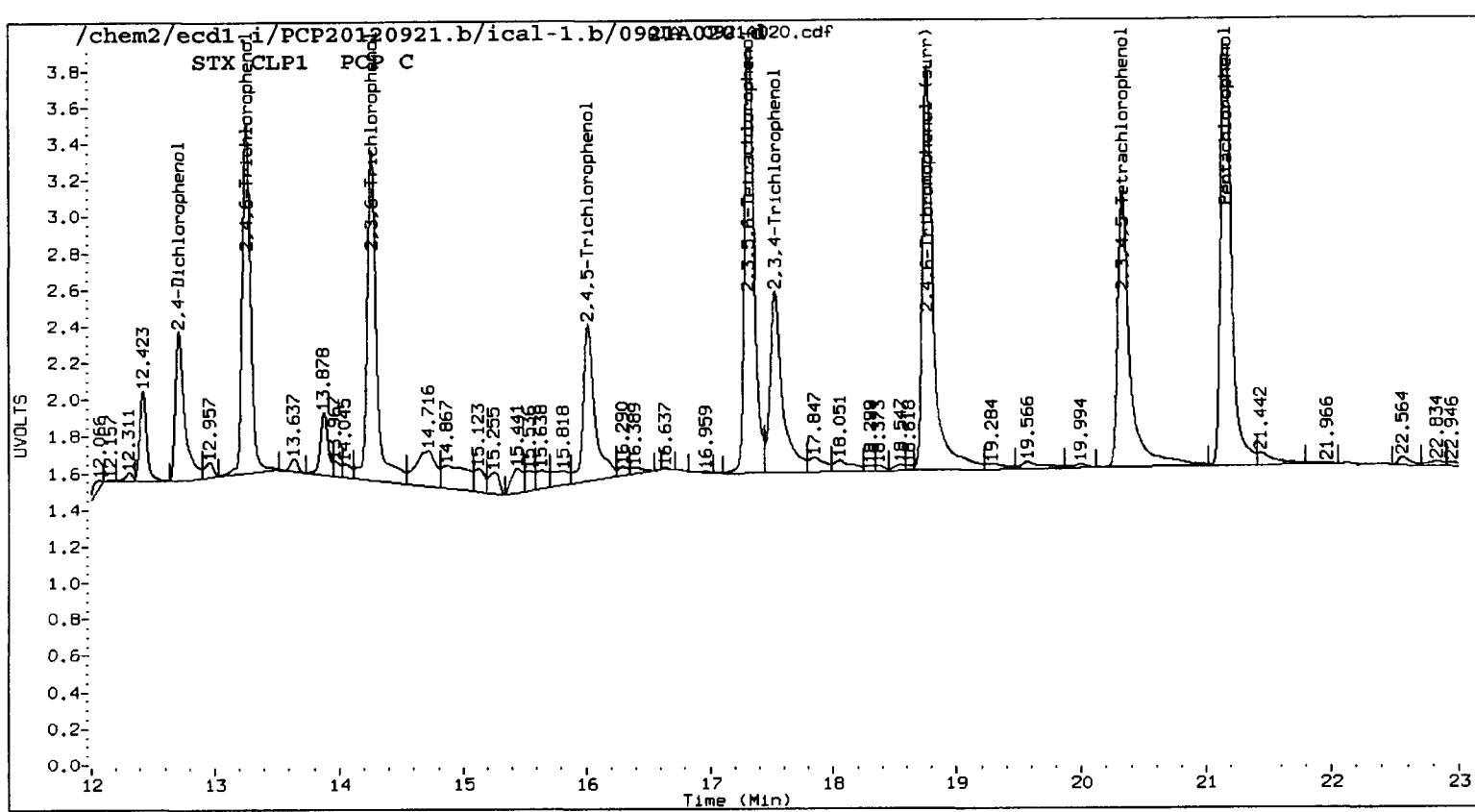
Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

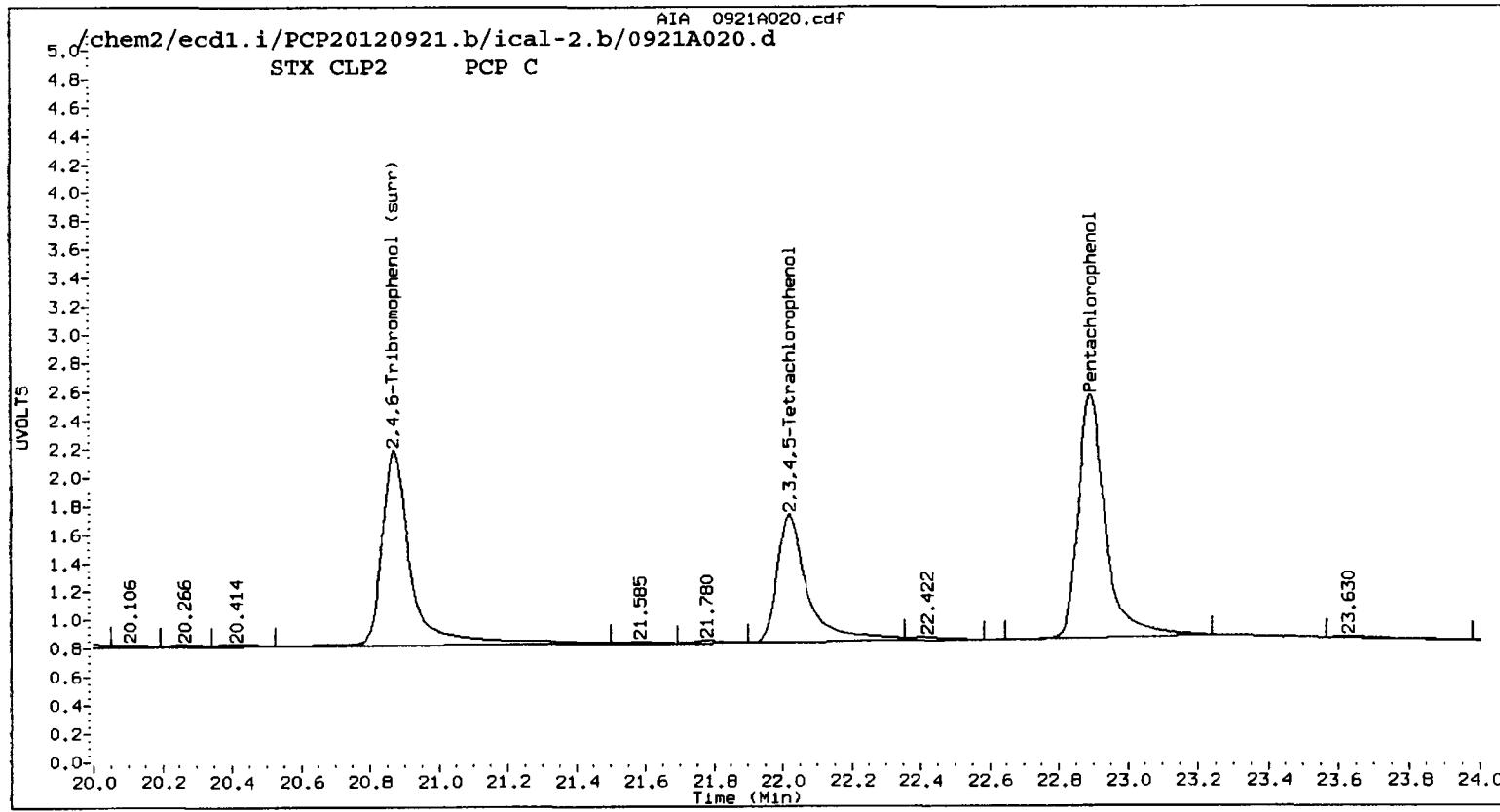
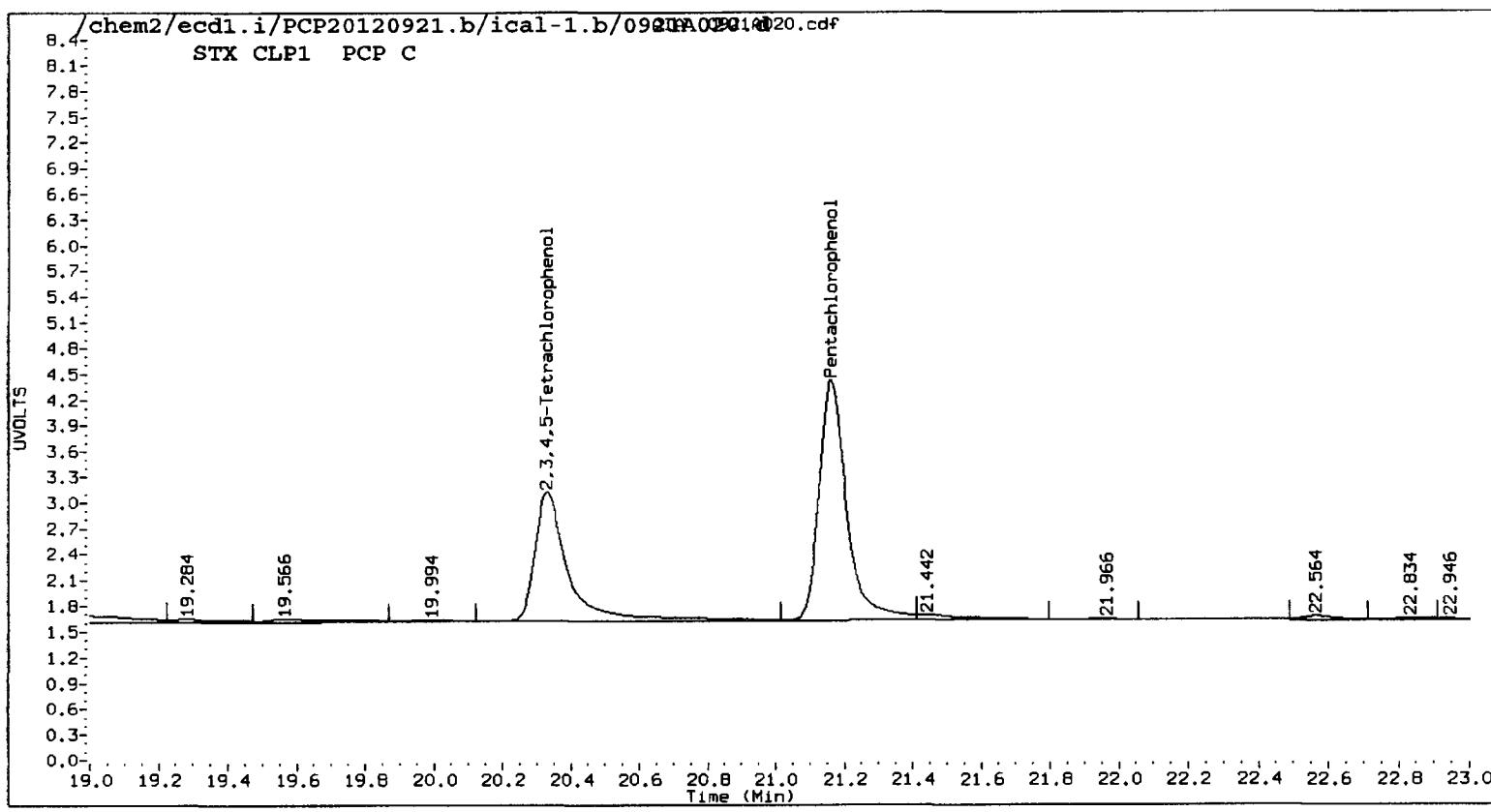
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 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 21-SEP-2012 21:17
 Compound Sublist: all Report Date: 09/25/2012 13:17
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

STX CLP1 Col			STX CLP2 Col			STX CLP1	STX CLP2		
RT	Shift	Response	RT	Shift	Response	on col	on col	RPD	Compound
21.163	0.005	779765	22.892	0.004	469083	12.6843	13.1437	3.6	Pentachlorophenol
13.256	0.000	430433	14.265	0.001	314632	13.5134	12.5299	7.6	2,4,6-Trichlorophenol
14.256	0.002	478158	15.508	0.001	300088	13.1359	13.3266	1.4	2,3,6-Trichlorophenol
16.015	0.005	284711	17.427	0.004	188136	12.6131	12.5093	0.8	2,4,5-Trichlorophenol
17.527	0.008	350554	18.974	0.006	230095	12.7445	11.5658	9.7	2,3,4-Trichlorophenol
17.318	0.003	630342	18.754	0.002	388236	13.4136	13.4405	0.2	2,3,5,6-Tetrachlorophenol
20.331	0.008	520404	22.019	0.007	273735	13.2689	13.2095	0.4	2,3,4,5-Tetrachlorophenol
12.713	0.002	219893	13.780	0.002	142277	131.1224	125.2571	4.6	2,4-Dichlorophenol
18.768	0.005	655694	20.870	0.003	408156	13.0	13.5	3.7	2,4,6-Tribromophenol (s)

PERCENT RECOVERY

COMPOUND	Col1	Col2
2,4,6-TBP (surr)	52.2	54.1

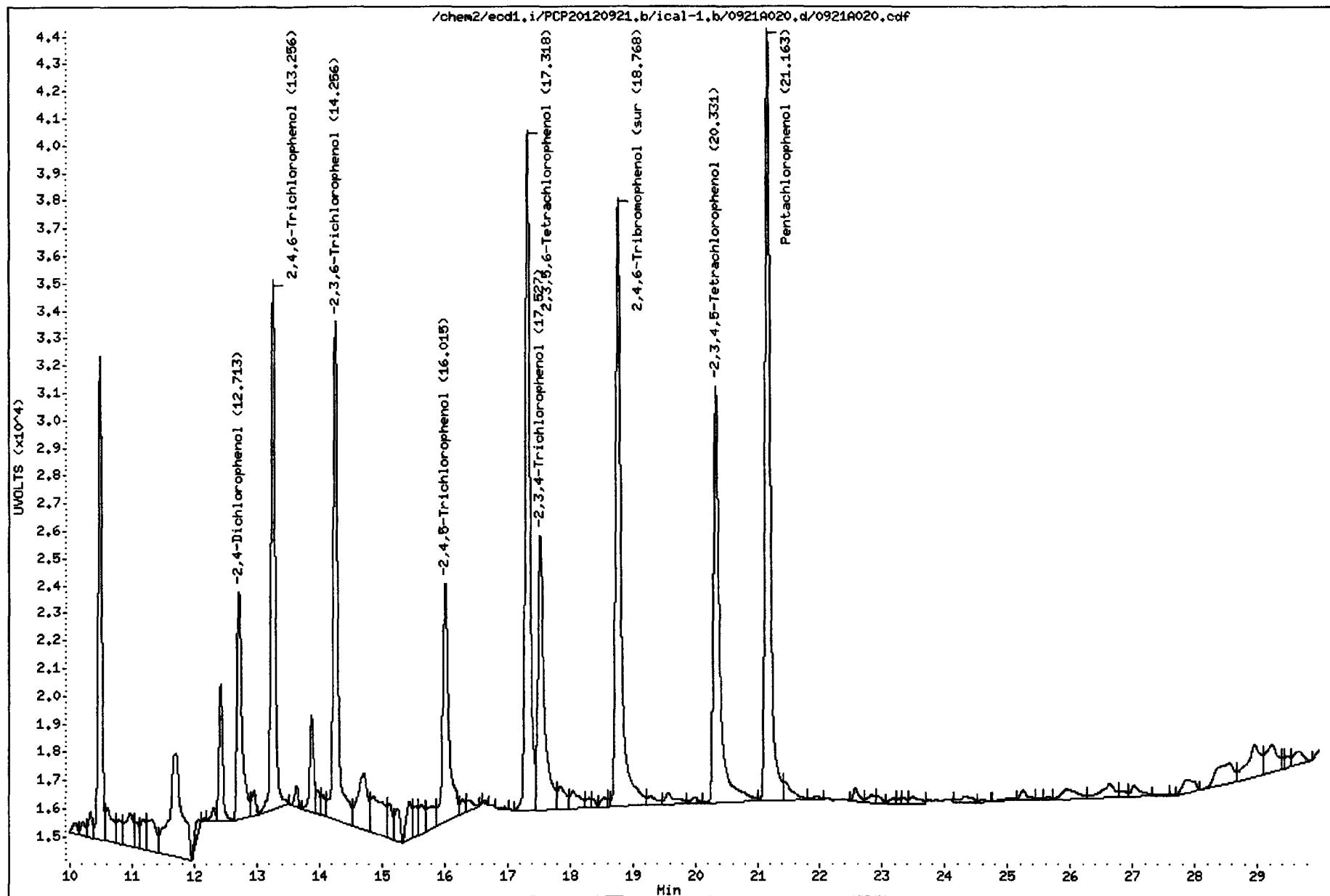




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Date : 21-SEP-2012 21:17
Client ID:
Sample Info: PCP C
Purge Volume: 500.0
Column phase: STX CLP1

Page 1

Instrument: ecd1.i
Operator: ar
Column diameter: 0.53



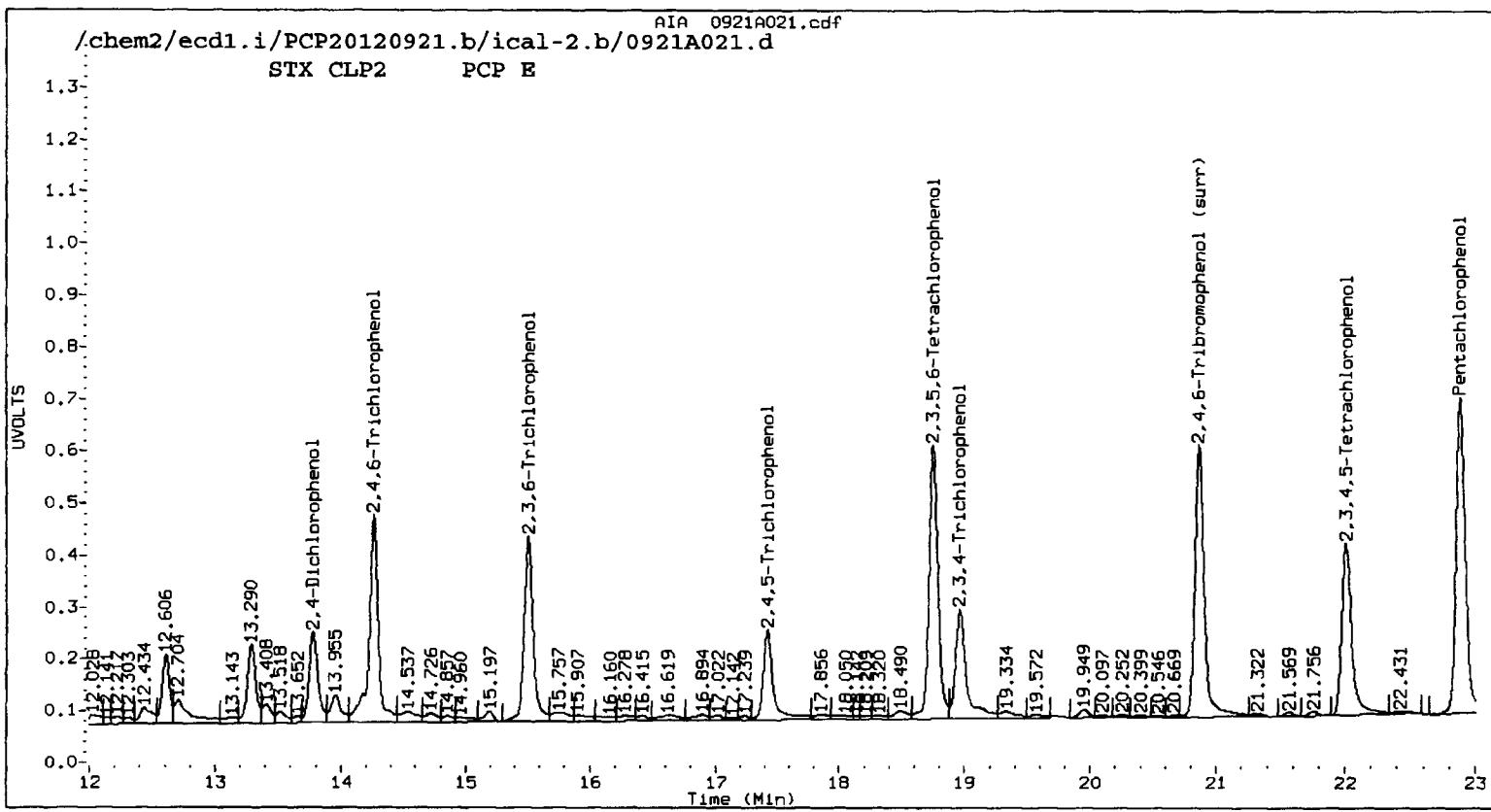
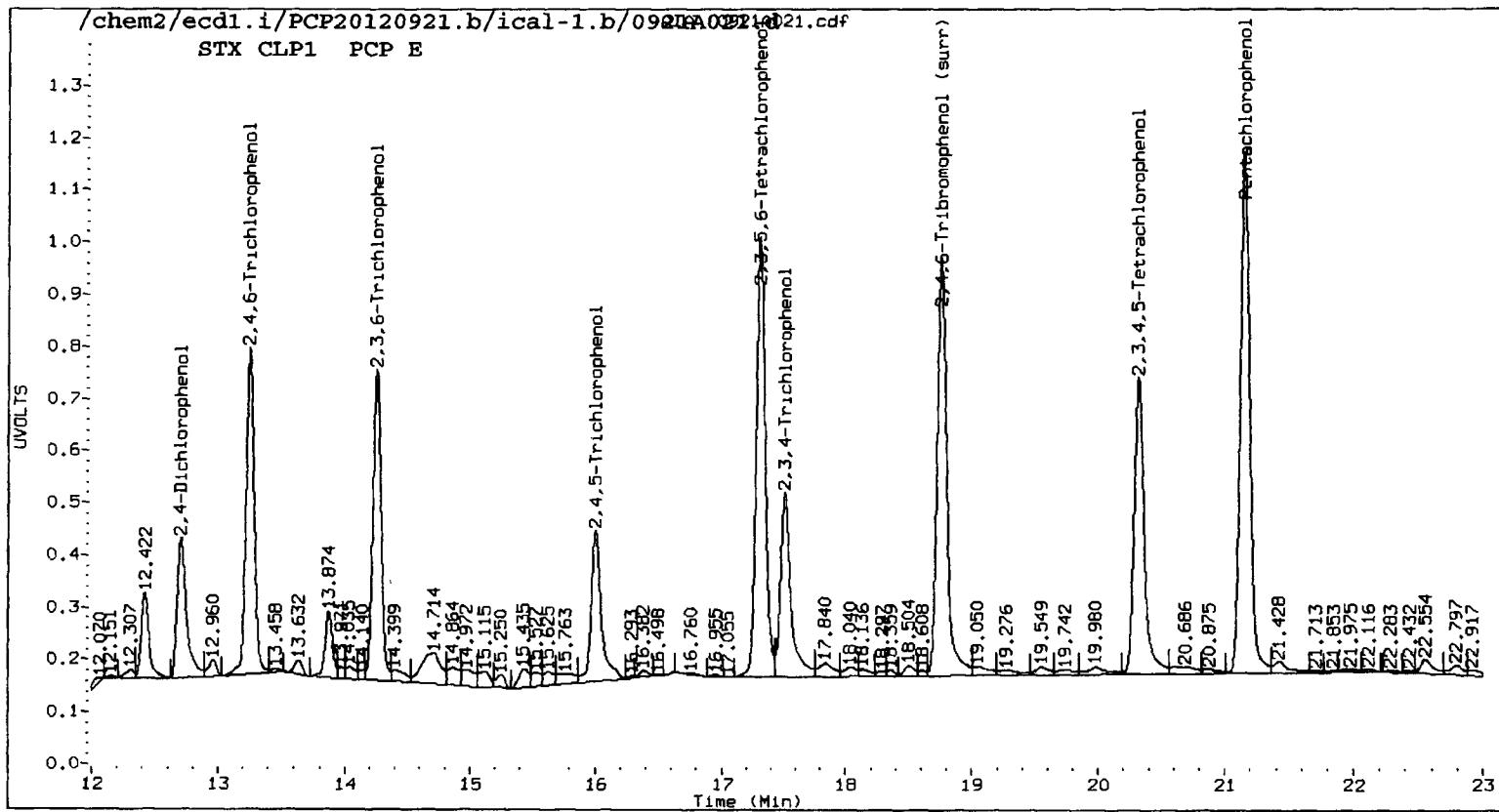
Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

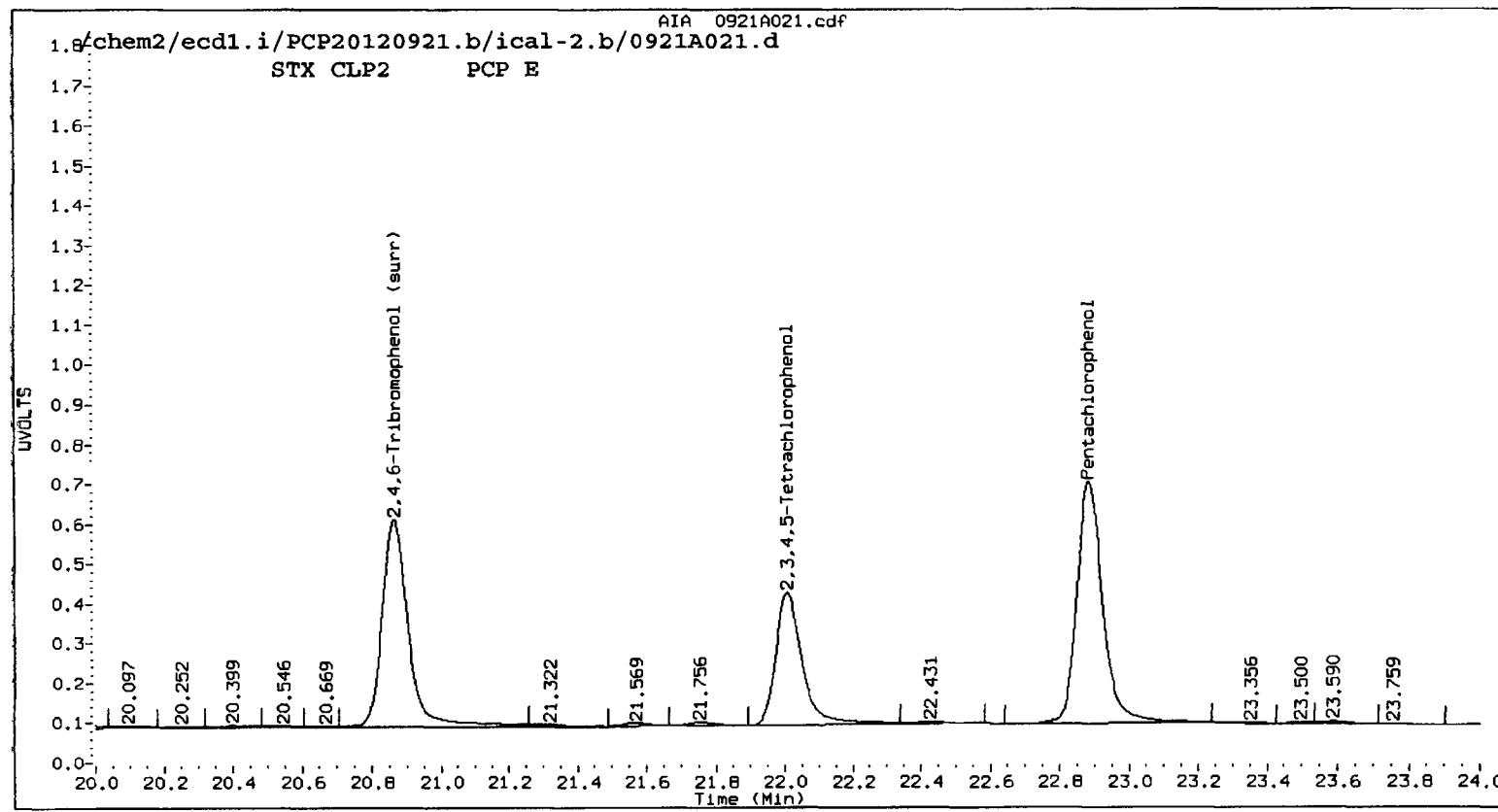
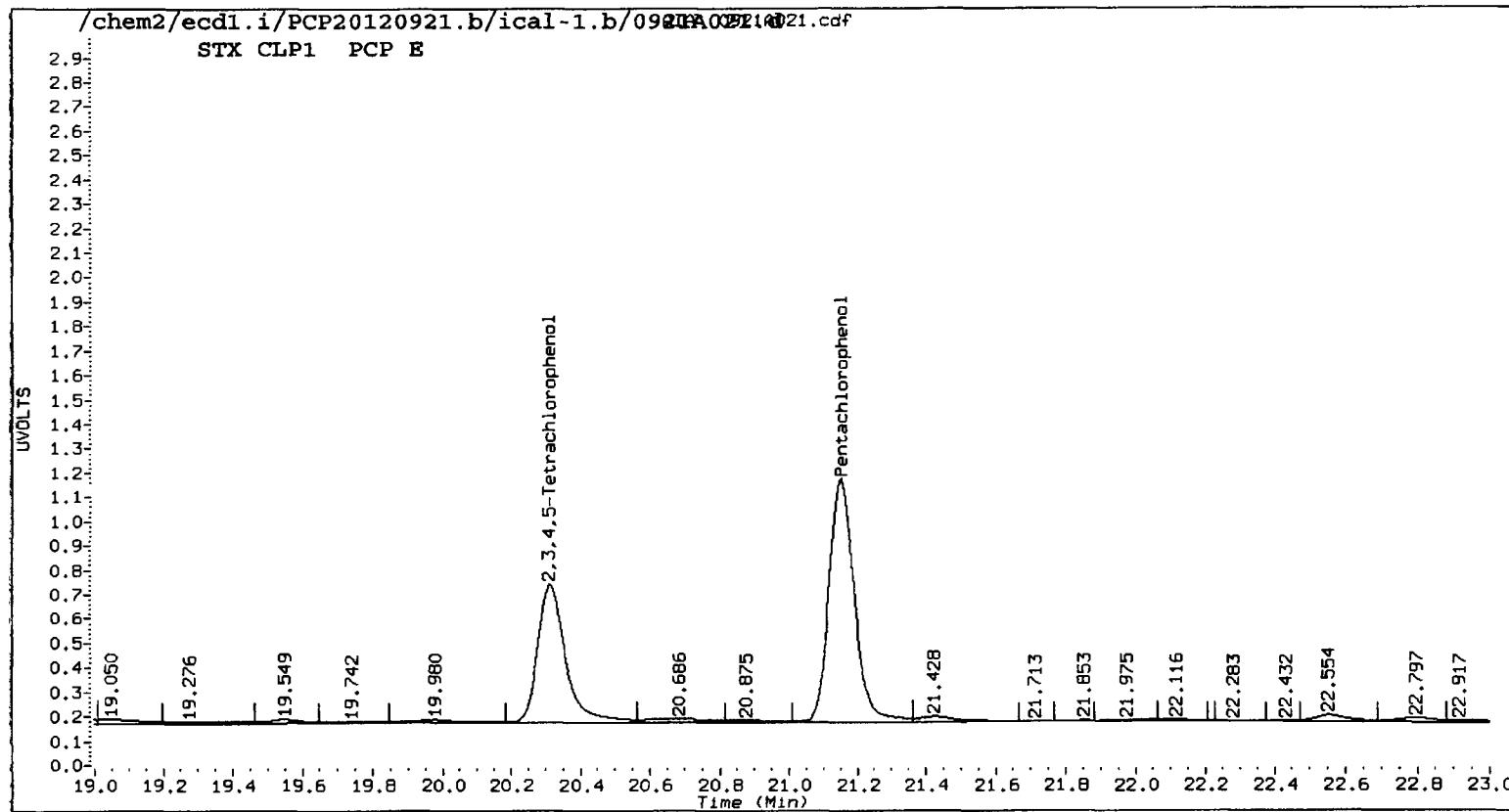
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Data file 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A021.d Client ID:
Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 21-SEP-2012 21:53
Compound Sublist: all Report Date: 09/25/2012 13:17
Instrument: ecd1.i Matrix: WATER
Operator: ar Dilution Factor: 1.000

RT	STX CLP1 Col Shift Response	RT	STX CLP2 Col Shift Response	on col	STX CLP1 on col	STX CLP2 RPD	Compound
21.153	-0.005 2579984	22.884 -0.004	1581355	49.8006 44.3095	11.7	Pentachlorophenol	
13.254	-0.002 1378613	14.261 -0.003	1046878	43.2813 50.4949	15.4	2,4,6-Trichloropheno	
14.251	-0.003 1360198	15.504 -0.003	891000	48.2202 48.9096	1.4	2,3,6-Trichlorophenol	
16.003	-0.007 828590	17.418 -0.005	509875	50.1100 50.4830	0.7	2,4,5-Trichlorophenol	
17.511	-0.008 1050439	18.961 -0.007	654702	49.7537 52.2530	4.9	2,3,4-Trichlorophenol	
17.311	-0.004 2092823	18.748 -0.004	1303311	44.5349 45.1198	1.3	2,3,5,6-Tetrachlorop	
20.314	-0.009 1585782	22.006 -0.006	902939	49.3364 43.5727	12.4	2,3,4,5-Tetrachloroph	
12.707	-0.004 695219	13.775 -0.003	438585	493.2130 505.6351	2.5	2,4-Dichlorophenol	
18.757	-0.006 2092155	20.863 -0.004	1376359	49.7 45.6	8.5	2,4,6-Tribromophenol	

PERCENT RECOVERY

COMPOUND	Col1	Col2
2,4,6-TBP (surr)	198.9	182.6

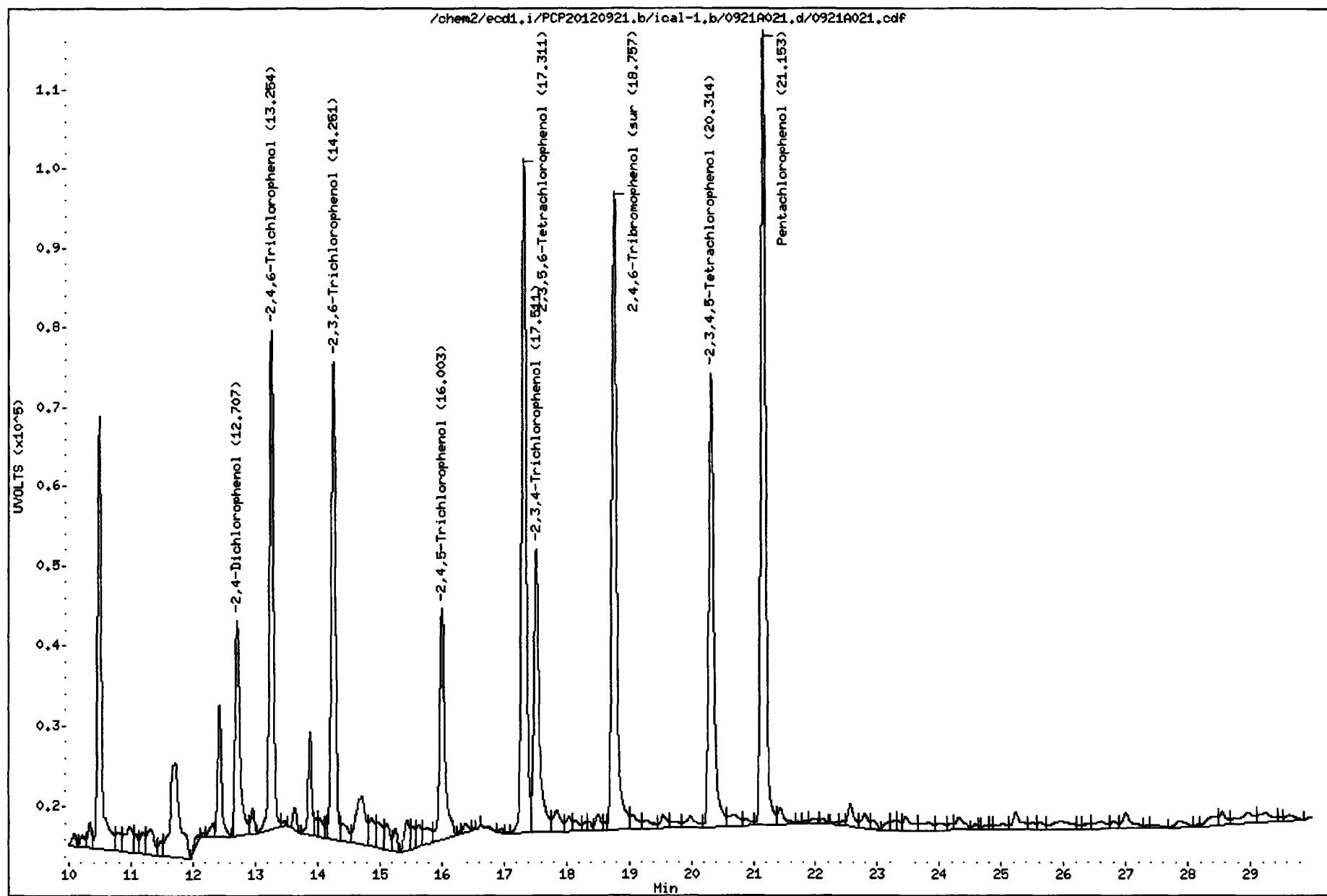




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Date : 21-SEP-2012 21:53
Client ID:
Sample Info: PCP E
Purge Volume: 500.0
Column phase: STX CLP1

Page 1

Instrument: ecd1.i
Operator: ar
Column diameter: 0.53



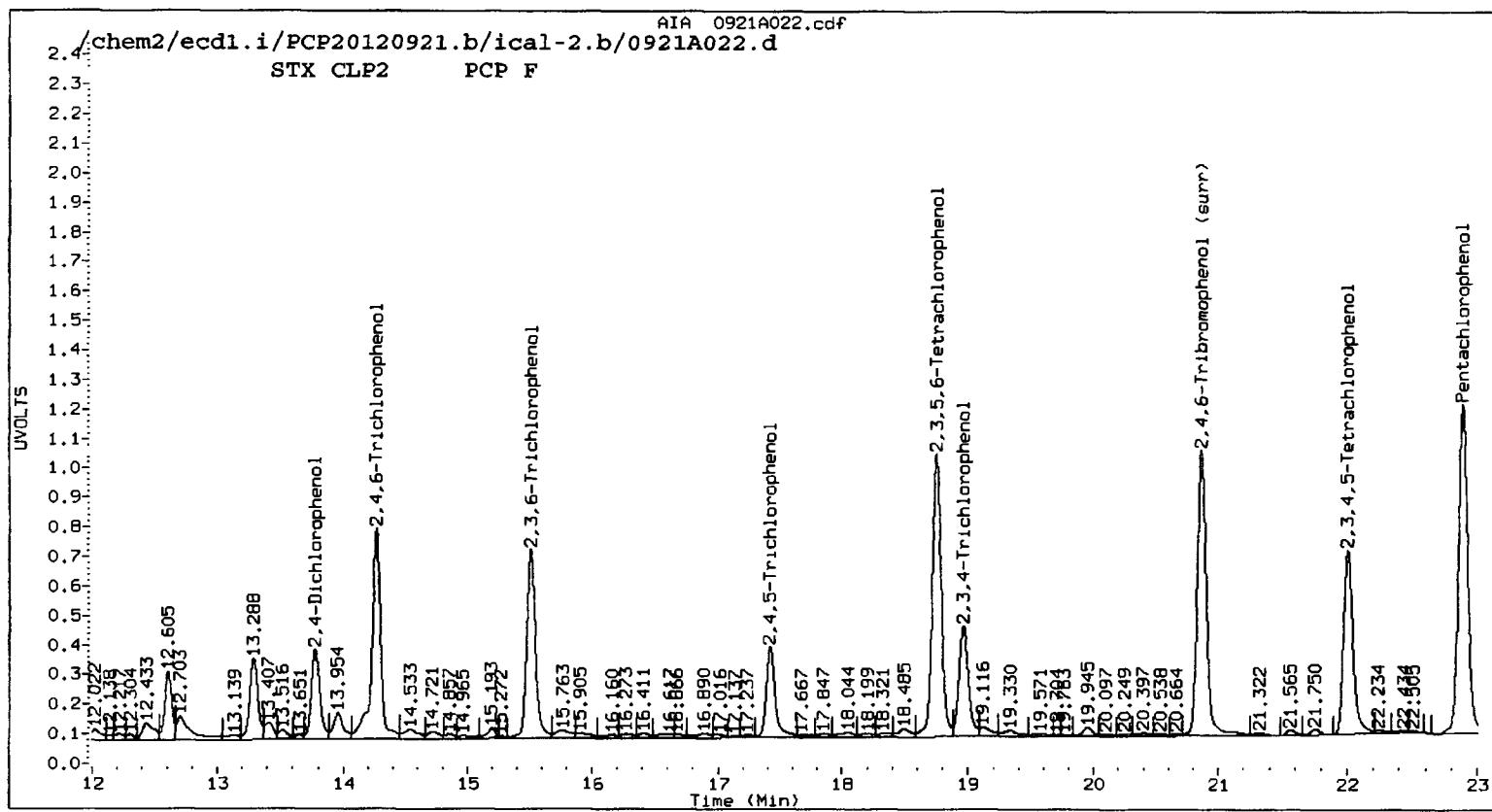
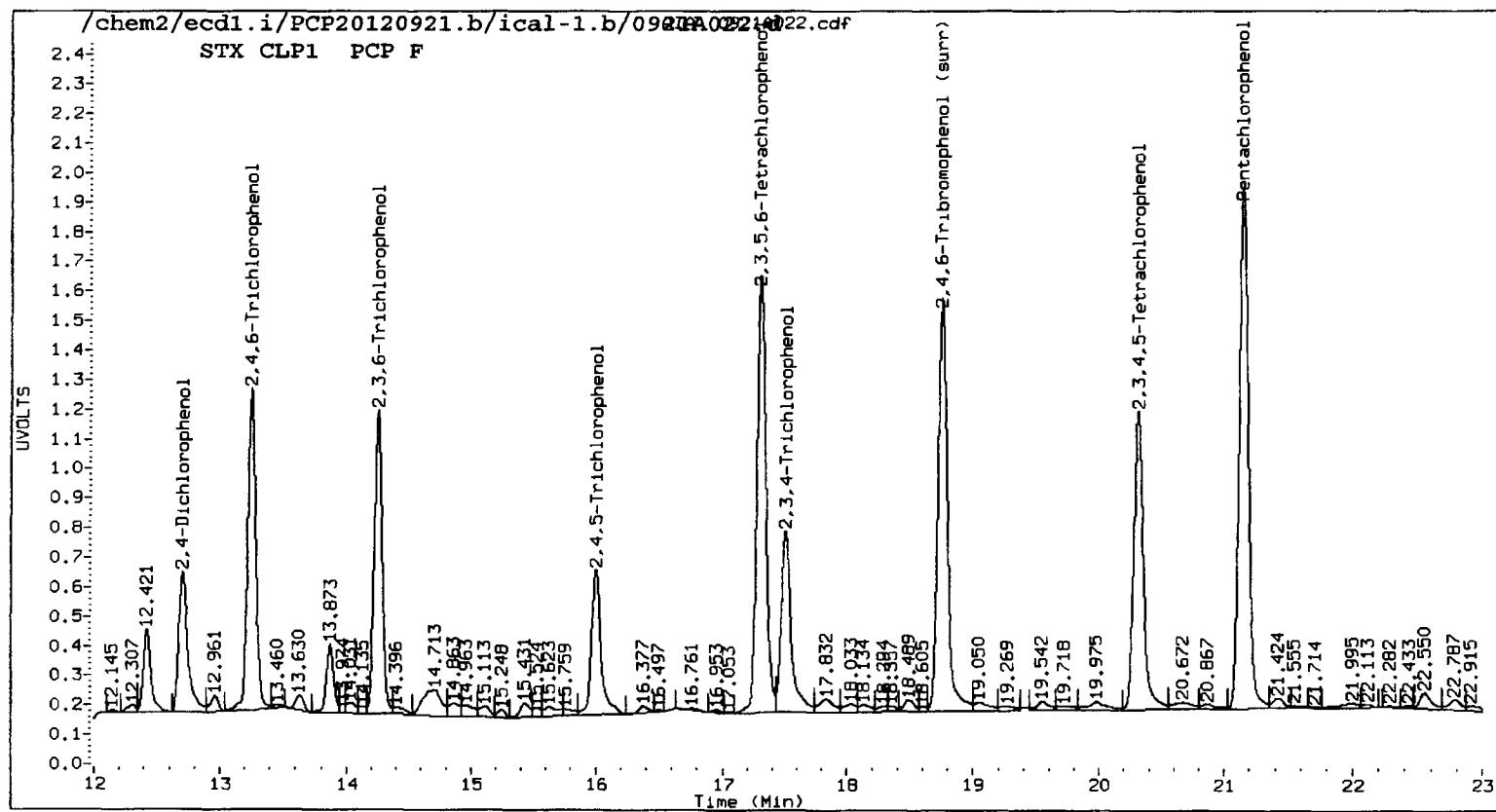
Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

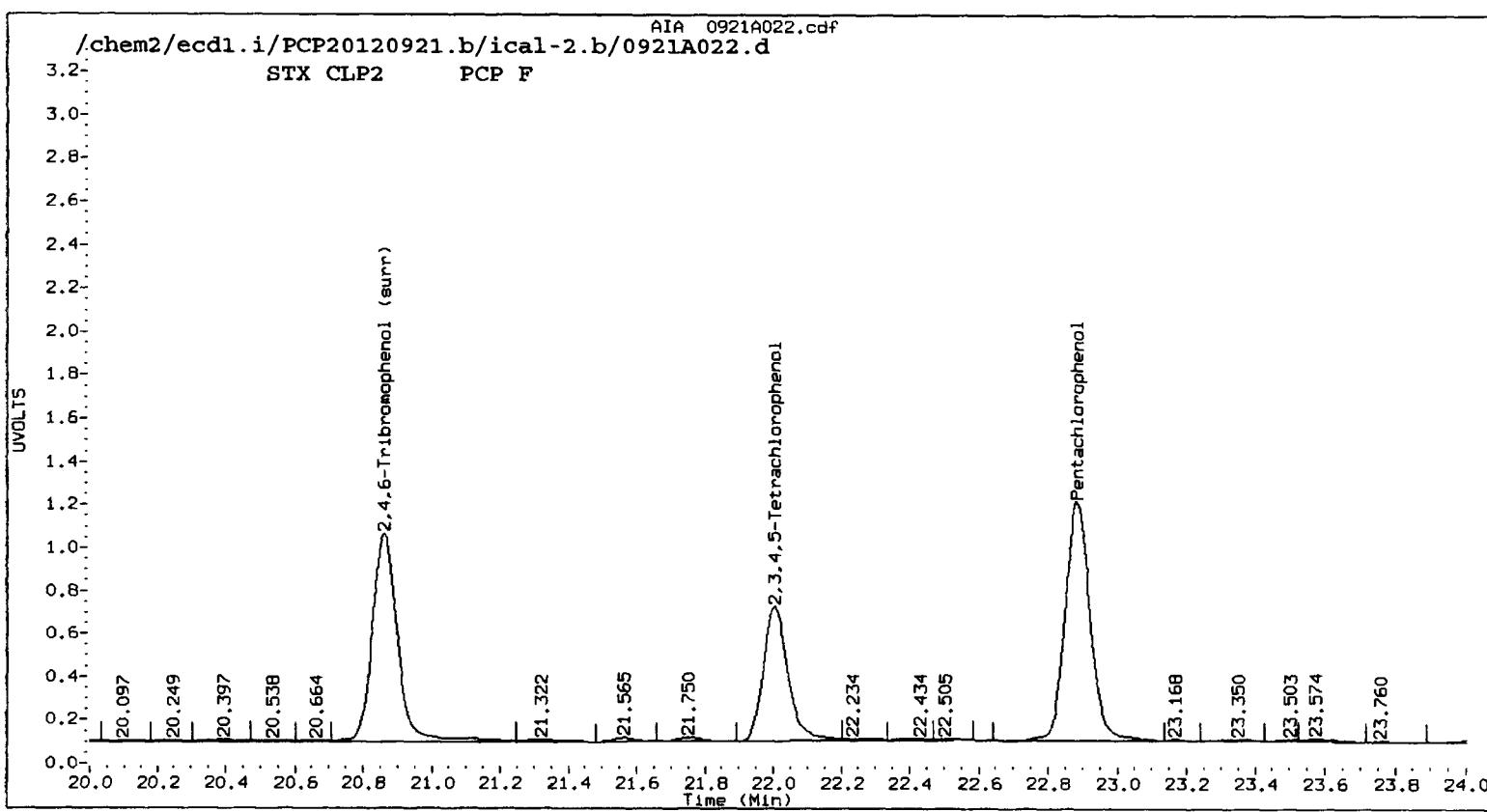
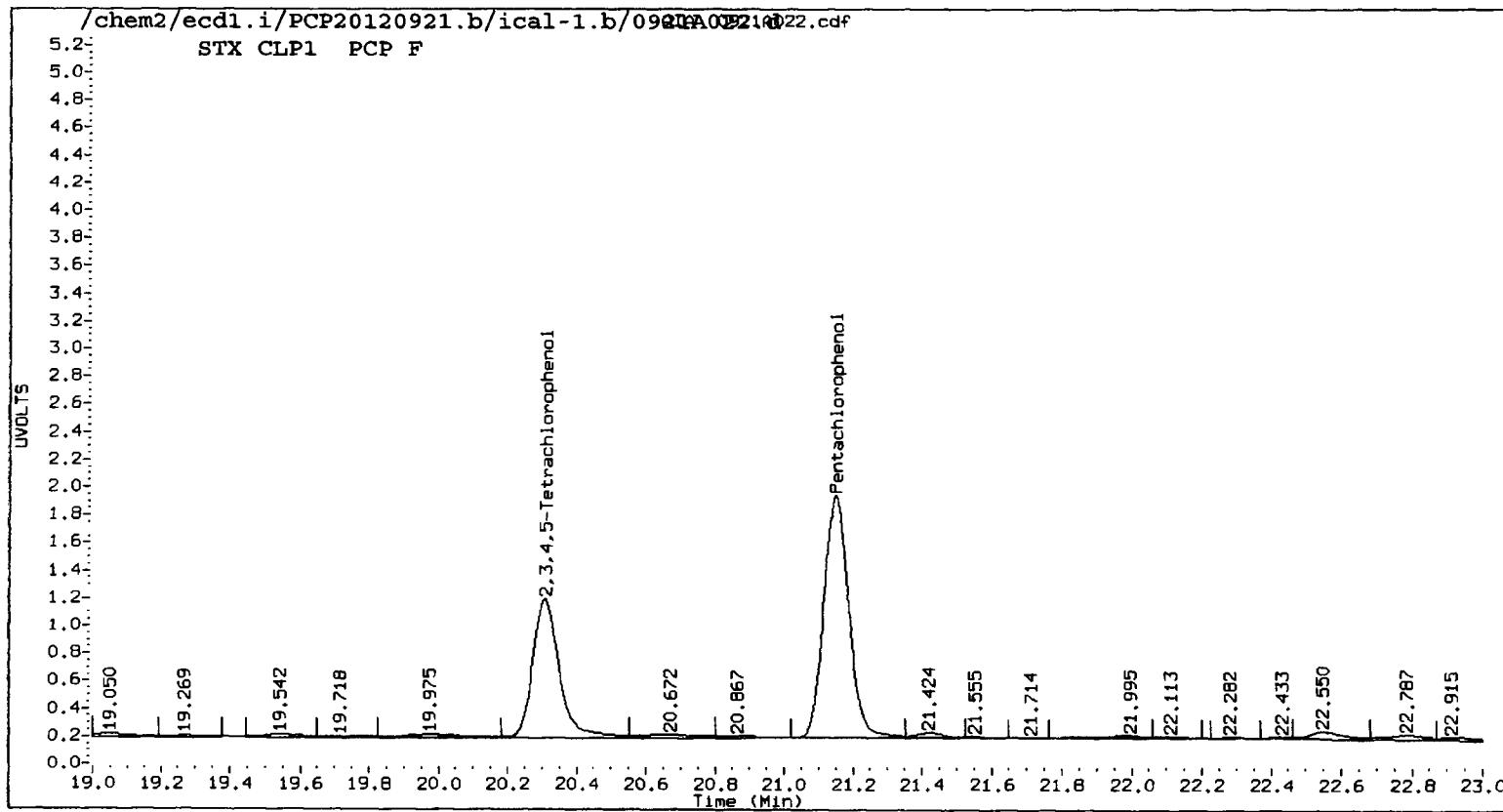
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 Data file 2: /chem2/ecdl.i/PCP20120921.b/ical-2.b/0921A022.d Client ID:
 Method: /chem2/ecdl.i/PCP20120921.b/PCP.m Injection Date: 21-SEP-2012 22:30
 Compound Sublist: all Report Date: 09/25/2012 13:17
 Instrument: ecdl.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

RT	STX CLP1 Col		STX CLP2 Col		on col	on col	RPD	STX CLP2 Compound
	Shift	Response	RT	Shift	Response			
21.149	-0.009	4455069	22.882	-0.006	2798545	100.0821	78.4151	24.3 Pentachlorophenol
13.252	-0.004	2381933	14.259	-0.005	1768391	74.7804	99.9503	28.8 2,4,6-Trichloropheno
14.249	-0.005	2289663	15.502	-0.005	1518853	100.4218	100.2932	0.1 2,3,6-Trichlorophen
15.999	-0.011	1327374	17.414	-0.009	786042	99.9642	99.7678	0.2 2,4,5-Trichlorophenol
17.506	-0.013	1725987	18.957	-0.011	972757	100.0920	99.1667	0.9 2,3,4-Trichlorophenol
17.307	-0.008	3620738	18.746	-0.006	2314935	77.0486	80.1416	3.9 2,3,5,6-Tetrachlorop
20.308	-0.015	2706943	22.003	-0.009	1567510	100.2112	75.6425	27.9 2,3,4,5-Tetrachlorop
12.706	-0.005	1205949	13.773	-0.005	711683	1002.1393	999.2285	0.3 2,4-Dichlorophenol
18.753	-0.010	3597466	20.861	-0.006	2474892	100.1	82.1	19.8 2,4,6-Tribromophenol

PERCENT RECOVERY

COMPOUND	Col1	Col2
2,4,6-TBP (surr)	400.4	328.3

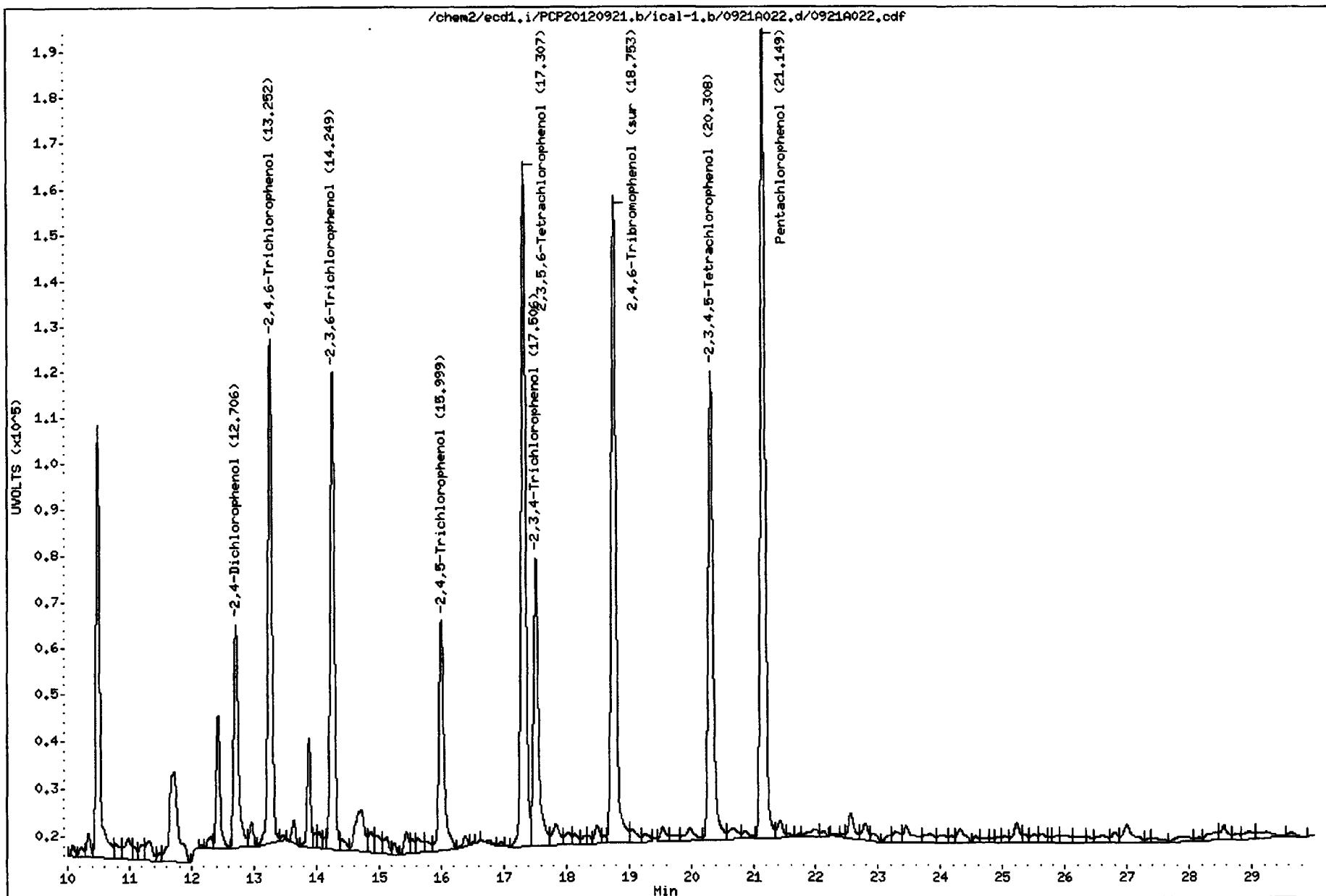




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Date : 21-SEP-2012 22:30
Client ID:
Sample Info: PCP F
Purge Volume: 500.0
Column phase: STX CLP1

Page 1

Instrument: ecd1.i
Operator: ar
Column diameter: 0.53



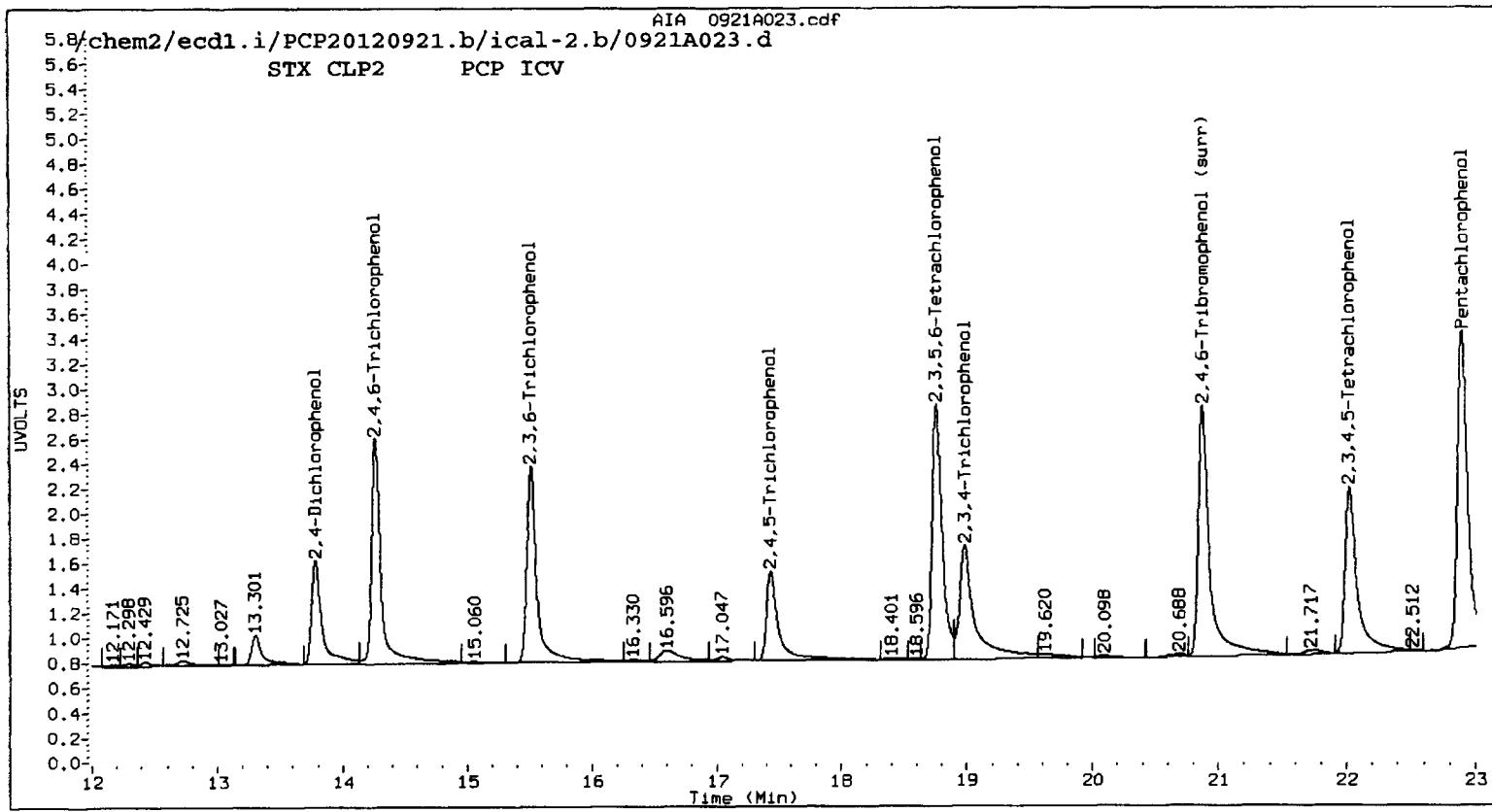
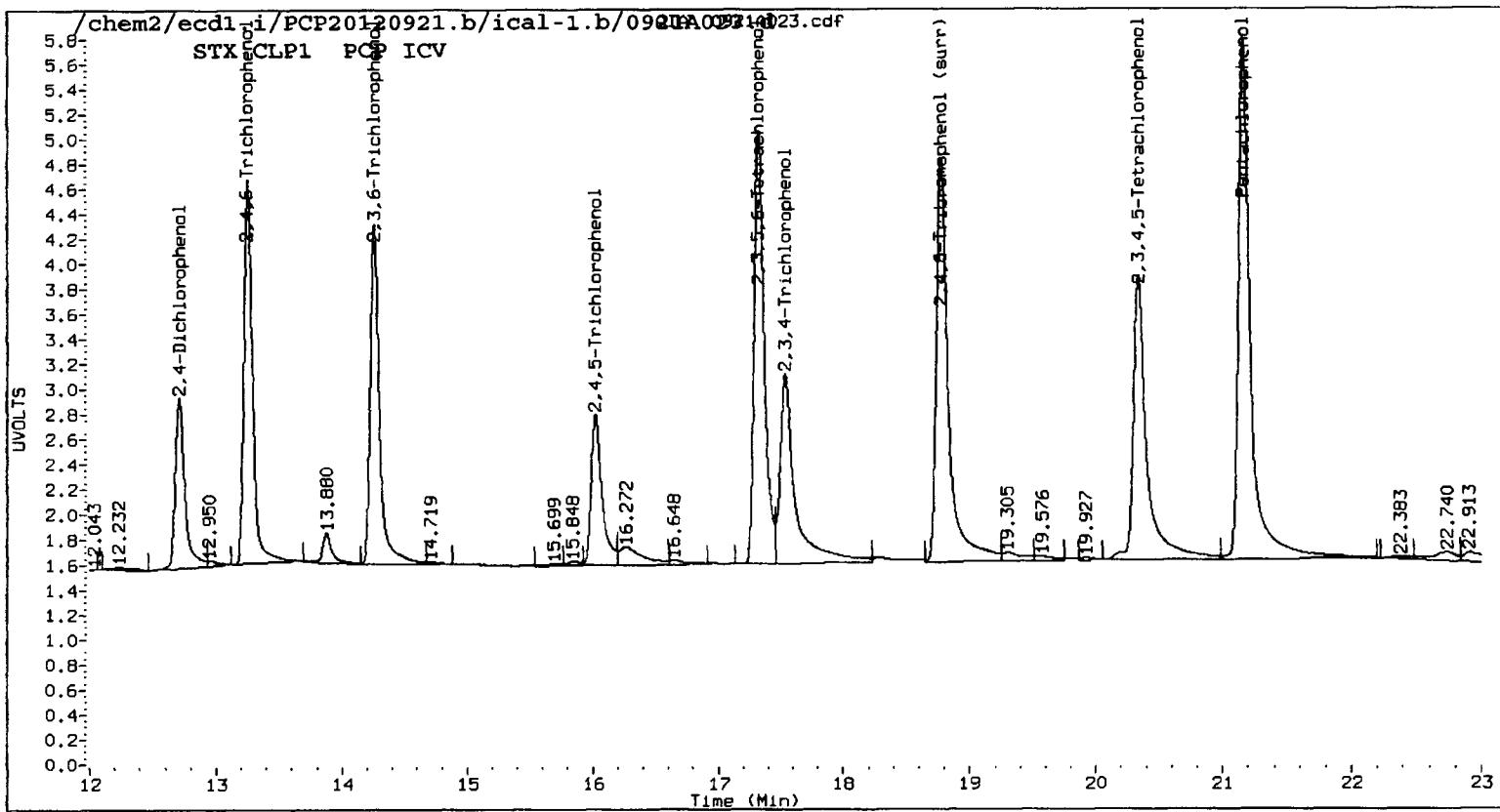
Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Data file 1: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A023.d ARI ID: PCP ICV
 Data file 2: /chem2/ecd1.i/PCP20120921.b/ical-2.b/0921A023.d Client ID:
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 21-SEP-2012 23:06
 Compound Sublist: all Report Date: 09/26/2012 09:48
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

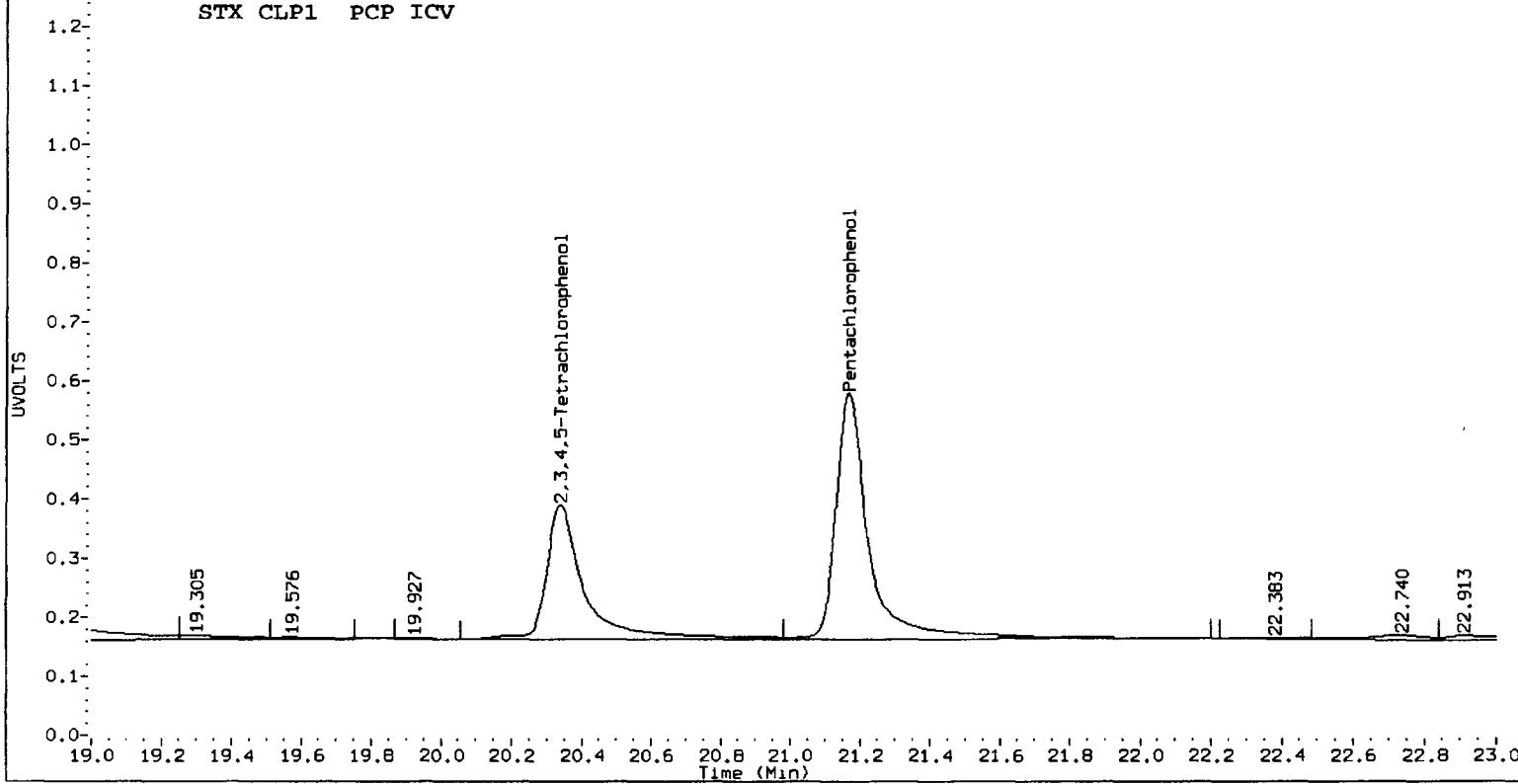
RT	STX CLP1 Col		STX CLP2 Col		on col	on col	RPD	STX CLP2 Compound
	Shift	Response	RT	Shift	Response			
21.169	0.011	1357525	22.896	0.008	737814	23.4053	20.6735	12.4 Pentachlorophenol
13.257	0.001	719668	14.264	0.000	486197	22.5938	20.3203	10.6 2,4,6-Trichlorophenol
14.257	0.003	693796	15.510	0.003	448159	20.4133	21.0796	3.2 2,3,6-Trichlorophenol
16.024	0.014	375368	17.433	0.010	252070	17.6414	18.3892	4.2 2,4,5-Trichlorophenol
17.540	0.021	645984	18.982	0.014	387901	26.4870	23.7575	10.9 2,3,4-Trichlorophenol
17.324	0.009	943248	18.759	0.007	557301	20.0721	19.2934	4.0 2,3,5,6-Tetrachlorophene
20.340	0.017	850177	22.024	0.012	442020	23.1547	21.3303	8.2 2,3,4,5-Tetrachlorophene
12.710	-0.002	362932	13.779	0.001	244804	228.7729	238.6019	4.2 2,4-Dichlorophenol
18.780	0.017	1048704	20.878	0.011	642772	22.0	21.3	3.0 2,4,6-Tribromophenol (

PERCENT RECOVERY

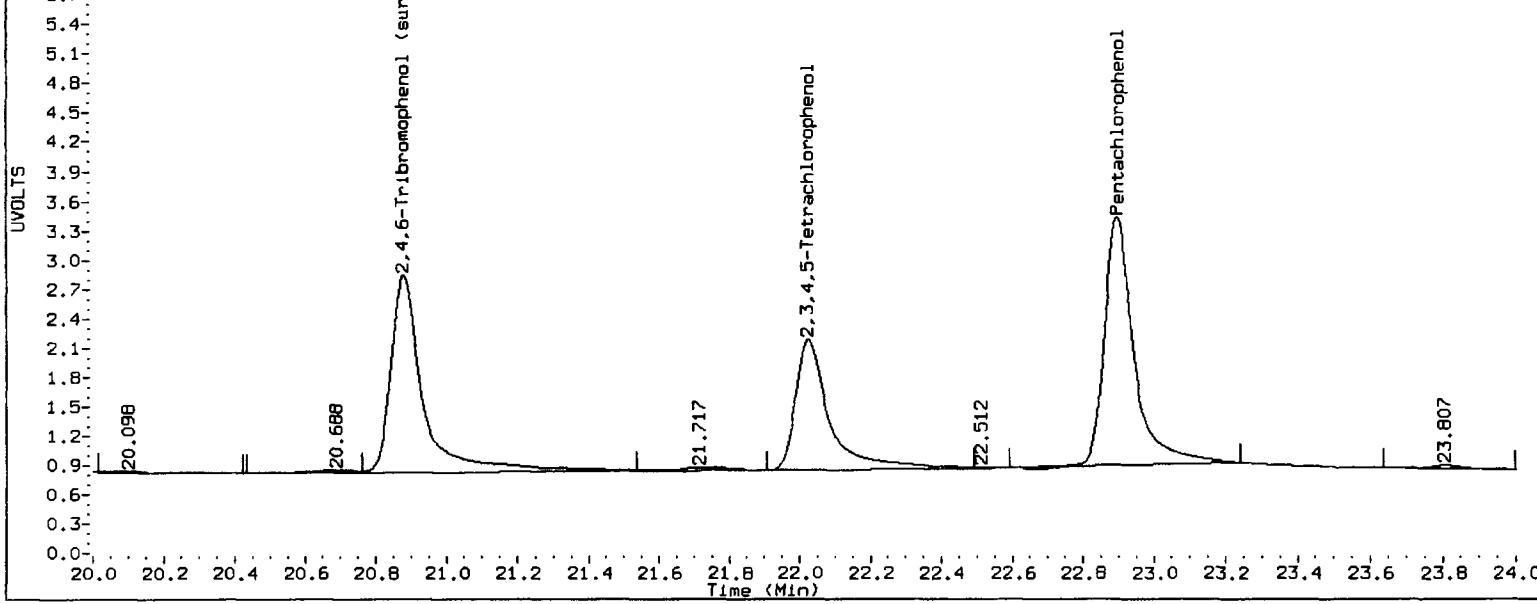
COMPOUND	Col1	Col2
Pentachlorophenol	93.6	82.7
2,4,6-Trichlorophenol	90.4	81.3
2,3,6-Trichlorophenol	81.7	84.3
2,4,5-Trichlorophenol	70.6	73.6
2,3,4-Trichlorophenol	105.9	95.0
2,3,5,6-Tetrachlorophenol	80.3	77.2
2,3,4,5-Tetrachlorophenol	92.6	85.3
2,4-Dichlorophenol	91.5	95.4
2,4,6-TBP (surr)	43.9	42.6
	87.8	85.2



/chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A023.cdf
STX CLP1 PCP ICV



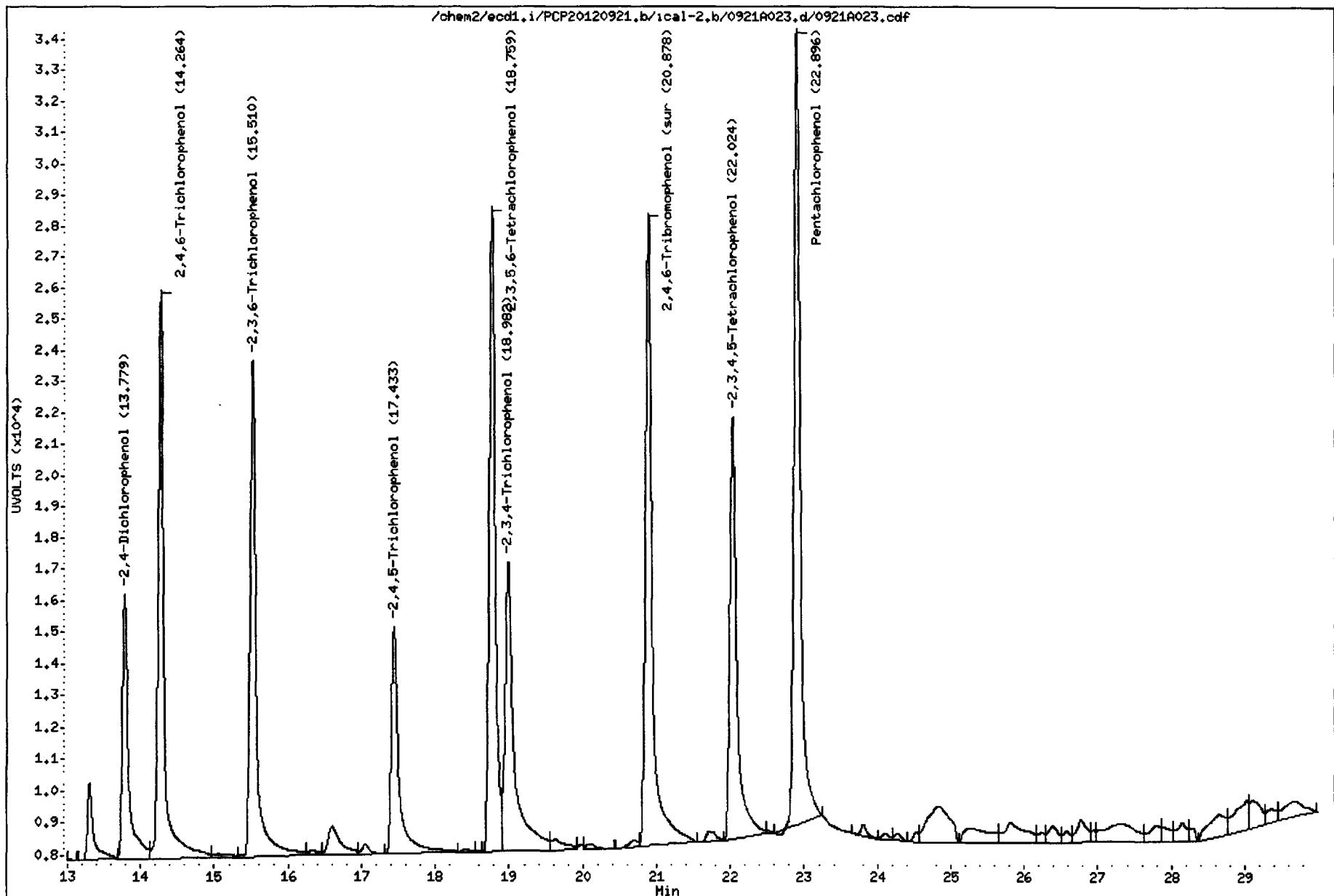
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STX CLP2 PCP ICV



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Date : 21-SEP-2012 23:06
Client ID:
Sample Info: PCP ICV
Purge Volume: 500.0
Column phase: STX CLP2

Page 1

Instrument: ecd1.i
Operator: ar
Column diameter: 0.53



Data File: /chem2/ecd1.i/PCP20120921.b/ical-1.b/0921A023.d

Page 1

Date : 21-SEP-2012 23:06

Client ID:

Instrument: ecd1.i

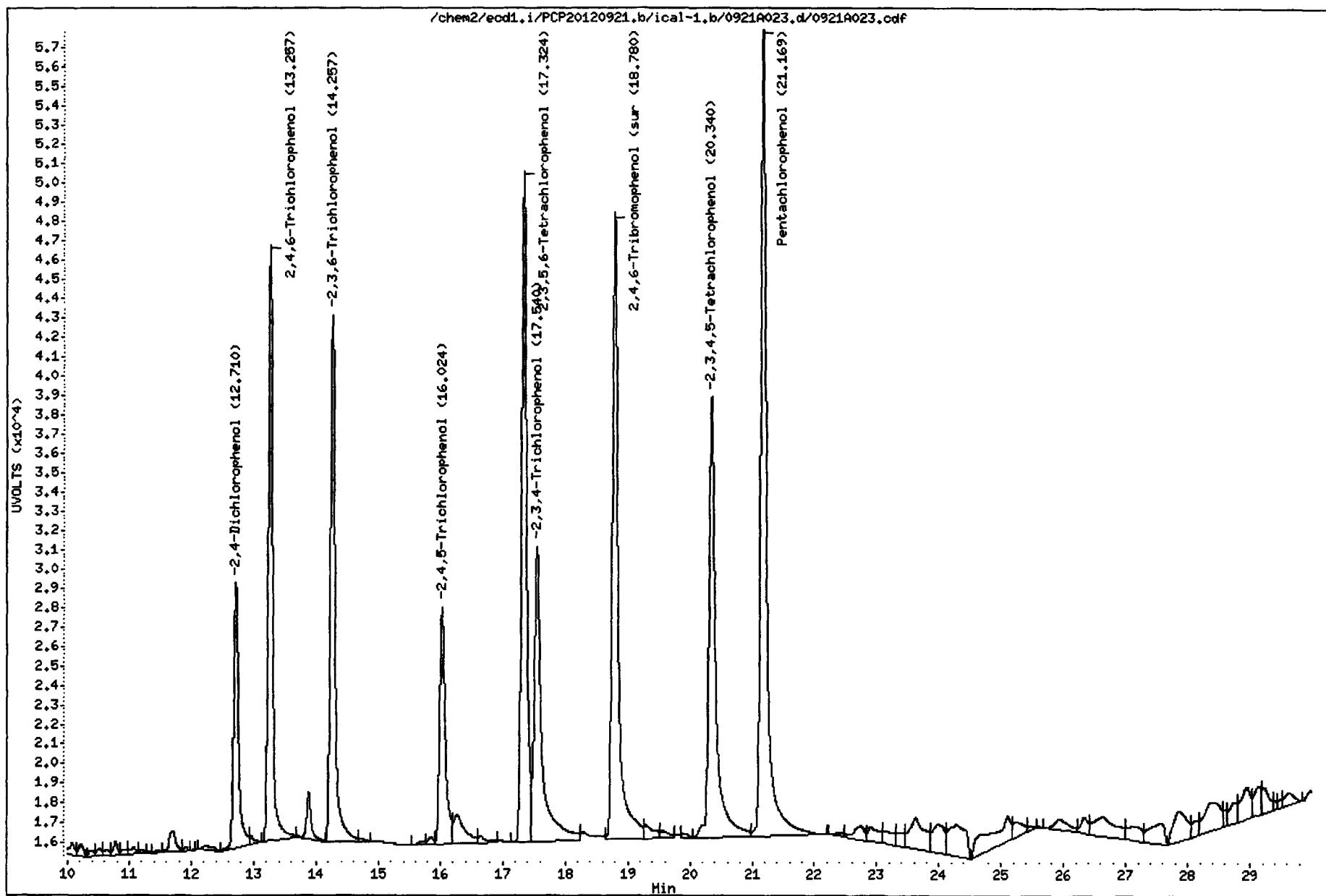
Sample Info: PCP ICV

Operator: ar

Purge Volume: 500.0

Column diameter: 0.53

Column phase: STX CLP1



**PCP/Chlorophenols Raw Data
Run Logs, Continuing Calibrations, and Raw Data**

ARI Job ID: VZ72

VZ72 : 000084

**GC Analyst Notes / Corrective Action Log**RI Project ID: 1372Client ID: bed Engineer 2RI SOP: **403S(PCB)** **405S(Herb)** **407S(TPH-D)** **409S(HCID)** **412S(PCP)** **423S(Pest)**
427S(Dir Inj) **428S(EPH)** **432S(EDB)** **Other**Parameter(s): PCPInstrument: FID-3A FID-3B FID-4A FID-4B FID-5 FID-7 FID-8
FID-9 ECD-1 ECD-5 ECD-6 ECD-7Dates: Curve: 09/25/13 Analysis Start: 09/25/13

Endrin/DDT Breakdown <15%? YES / NO / NA Method Blank In Control? YES / NO

Cal Meets RF & %RSD Criteria? YES / NO LCS/LCSD Recovery In Control? YES / NO

ICal Meets RF & %RSD Criteria? YES / NO Surrogate Recovery In Control? YES / NO

Manual Integrations for ICal? YES / NO Manual Integrations for Samples? YES / NO

Internal Standard Meets Criteria? YES / NO / NA Special Analysis Criteria Met? YES / NO / NA

Detail problems, corrective actions and/or other pertinent information below (use reverse side when necessary):**Additional Details on Reverse: Yes / No**Analyst: YZ Date: 09/23/13Reviewer: MW Date: 1/23

Analytical Resources Inc.: Organics Instrument Log
ECD1 Serial No.: 3410A39690

Date: 01/23/13

GC Program: Herbm

Calibration File: PCP20120921

Analysis: PCP

Column No: 922995/801642

Curve Date: 01/23/13

Analyst: YB

Column Type: STX CCP / CCR

Injection Vol.: 2 µl

IS/SS	Ical/Ccal	LCS/ICV
	<u>1034-3</u>	

Document All Maintenance Tasks In StarLIMS

GC LOG SUMMARY FOR DATABATCH - /chem2/ecd1.i/PCP20120921.b/0122-1.b

	Inject Date/Time	Filename	DF	LabID	ClientID
1	22-JAN-2013 12:43	0122A003.d	1	PCPCCAL	
2	22-JAN-2013 13:19	0122A004.d	1	VZ72MBW1	
3	22-JAN-2013 13:56	0122A005.d	1	VZ72LCSW1	
4	22-JAN-2013 14:32	0122A006.d	1	VZ72LCSDW1	
5	22-JAN-2013 15:08	0122A007.d	1	VZ72QLS1	
6	22-JAN-2013 15:45	0122A008.d	1	VZ72A	
7	22-JAN-2013 16:21	0122A009.d	1	VZ72B	
8	22-JAN-2013 16:57	0122A010.d	1	PCP	
9	22-JAN-2013 17:33	0122A011.d	1	PCPCCAL	

YB 01/23/13

Every line must contain information or be lined out. Make all entries legible.
Start a new page for each QC period. Document All Maintenance Tasks In StarLIMS

Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Data file 1: /chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A003.d ARI ID: PCPCCAL *Y2 01/23/13*
 Data file 2: /chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A003.d Client ID:
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 22-JAN-2013 12:43
 Compound Sublist: pcpca1 Report Date: 01/23/2013 11:16
 Instrument: ecd1.i Matrix: NONE
 Operator: ar Dilution Factor: 1.000

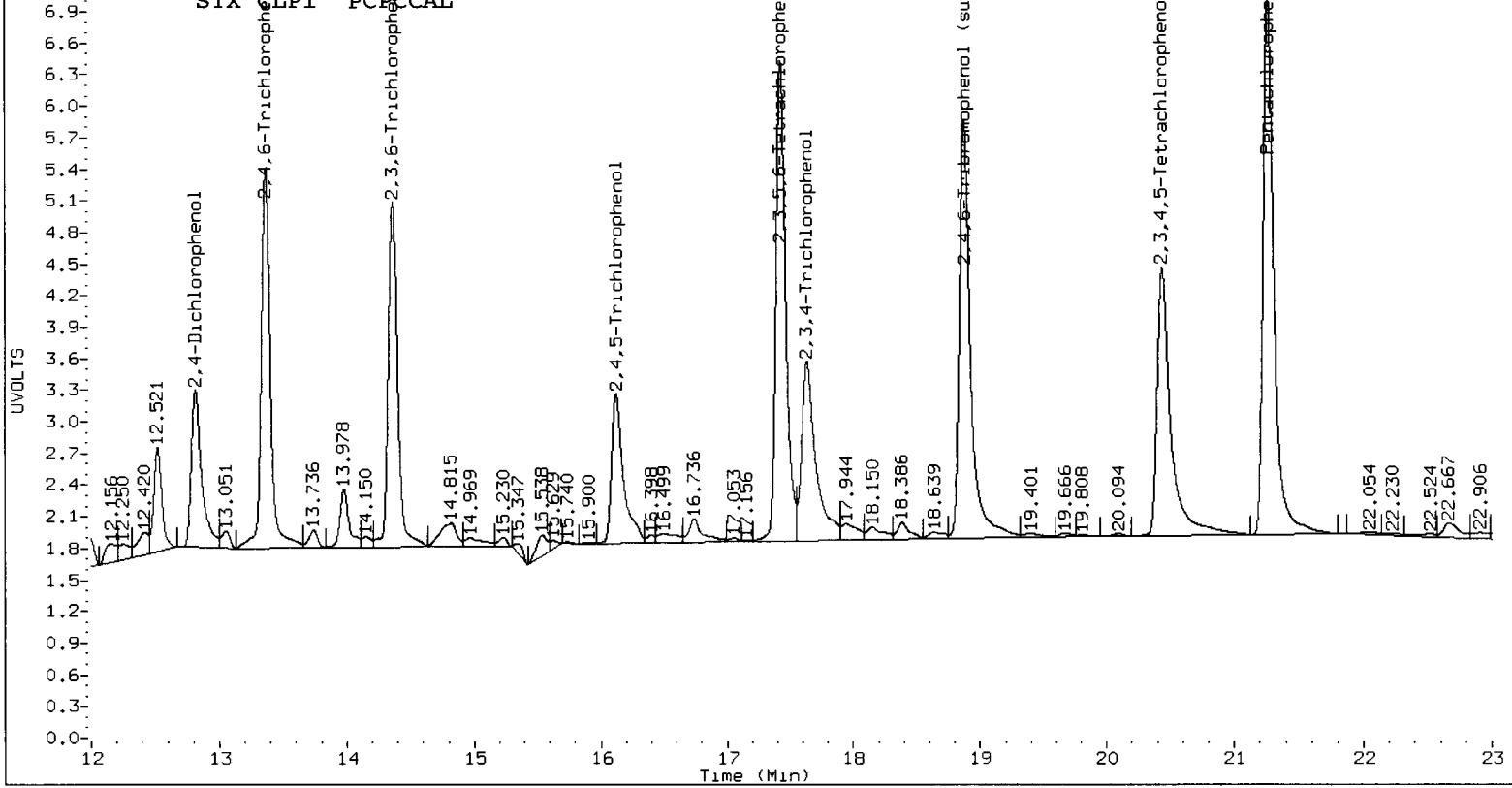
RT	STX CLP1 Col			STX CLP2 Col			on col	on col	RPD	STX CLP2 Compound
	Shift	Response	RT	Shift	Response					
21.266	0.030	1506094	22.989	0.026	951556	26.3441	26.6626	1.2	Pentachlorophenol	
13.359	0.029	898122	14.364	0.026	542354	28.1964	23.0172	20.2	2,4,6-Trichlorophenol	
14.360	0.030	826461	15.608	0.026	487392	25.3085	23.2643	8.4	2,3,6-Trichlorophenol	
16.120	0.032	463005	17.525	0.026	298393	22.9668	23.1658	0.9	2,4,5-Trichlorophenol	
17.633	0.031	666139	19.072	0.026	391274	27.5246	24.0560	13.4	2,3,4-Trichlorophenol	
17.422	0.030	1234608	18.853	0.026	728997	26.2722	25.2374	4.0	2,3,5,6-Tetrachloroph	
20.437	0.030	970317	22.117	0.026	542974	27.0411	26.2020	3.2	2,3,4,5-Tetrachlorophe	
12.814	0.029	421308	13.877	0.026	224236	271.4237	214.3137	23.5	2,4-Dichlorophenol	
18.874	0.031	1228819	20.970	0.026	768313	26.3	25.5	3.3	2,4,6-Tribromophenol (

PERCENT RECOVERY

COMPOUND	Col1	Col2
<hr/>		
Pentachlorophenol	105.4	106.7
2,4,6-Trichlorophenol	112.8	92.1
2,3,6-Trichlorophenol	101.2	93.1
2,4,5-Trichlorophenol	91.9	92.7
2,3,4-Trichlorophenol	110.1	96.2
2,3,5,6-Tetrachlorophenol	105.1	100.9
2,3,4,5-Tetrachlorophenol	108.2	104.8
2,4-Dichlorophenol	108.6	85.7
2,4,6-TBP (surr)	105.3	101.9

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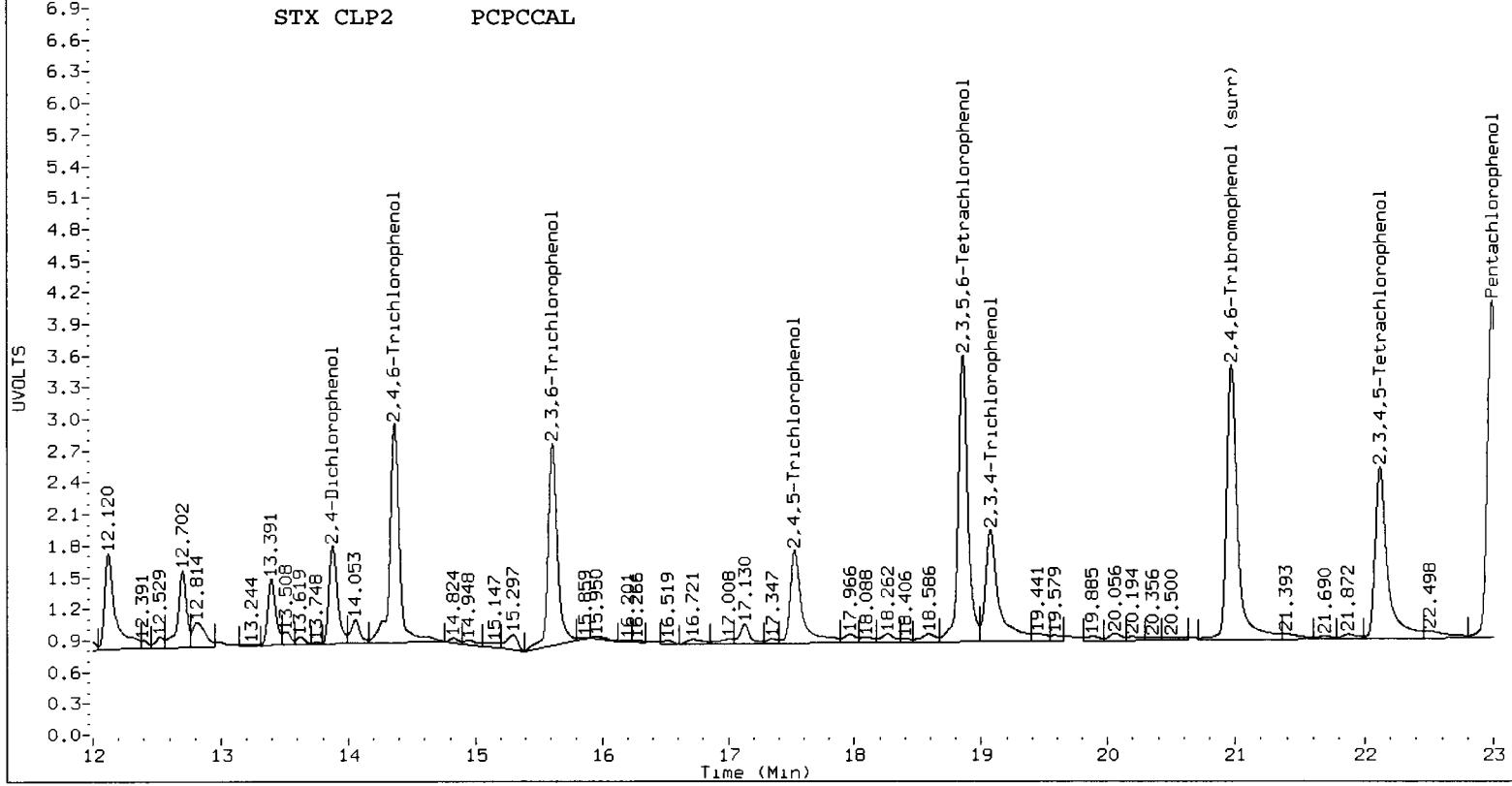
STX CLP1 PCPCCAL



AIA 0122A003.cdf

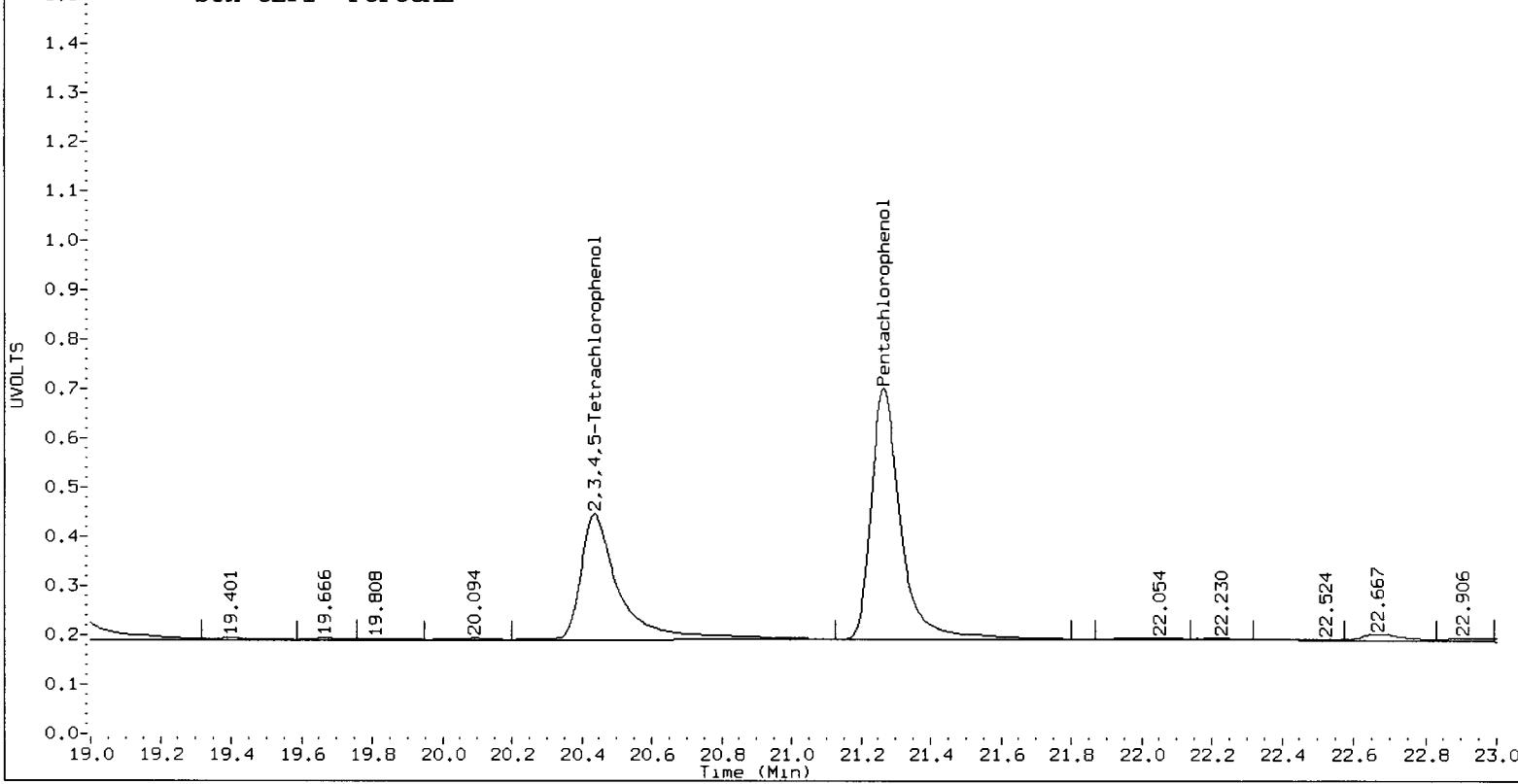
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STX CLP2 PCPCCAL



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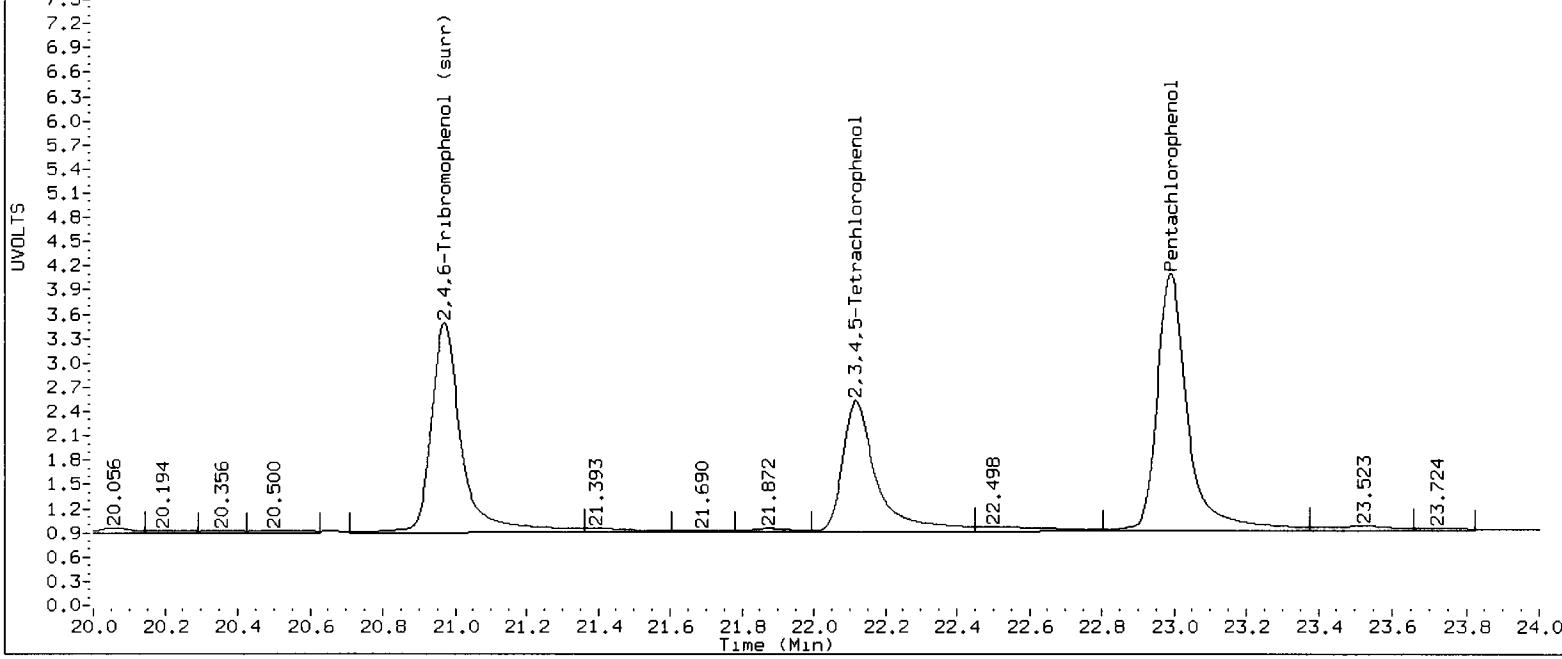
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AIA 0122A003.cdf

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STX CLP2 PCPCCAL



Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Y2 01/23/13

Data file 1: /chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A004.d ARI ID: VZ72MBW1
 Data file 2: /chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A004.d Client ID: VZ72MBW1
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 22-JAN-2013 13:19
 Compound Sublist: pcpcal Report Date: 01/23/2013 10:56
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

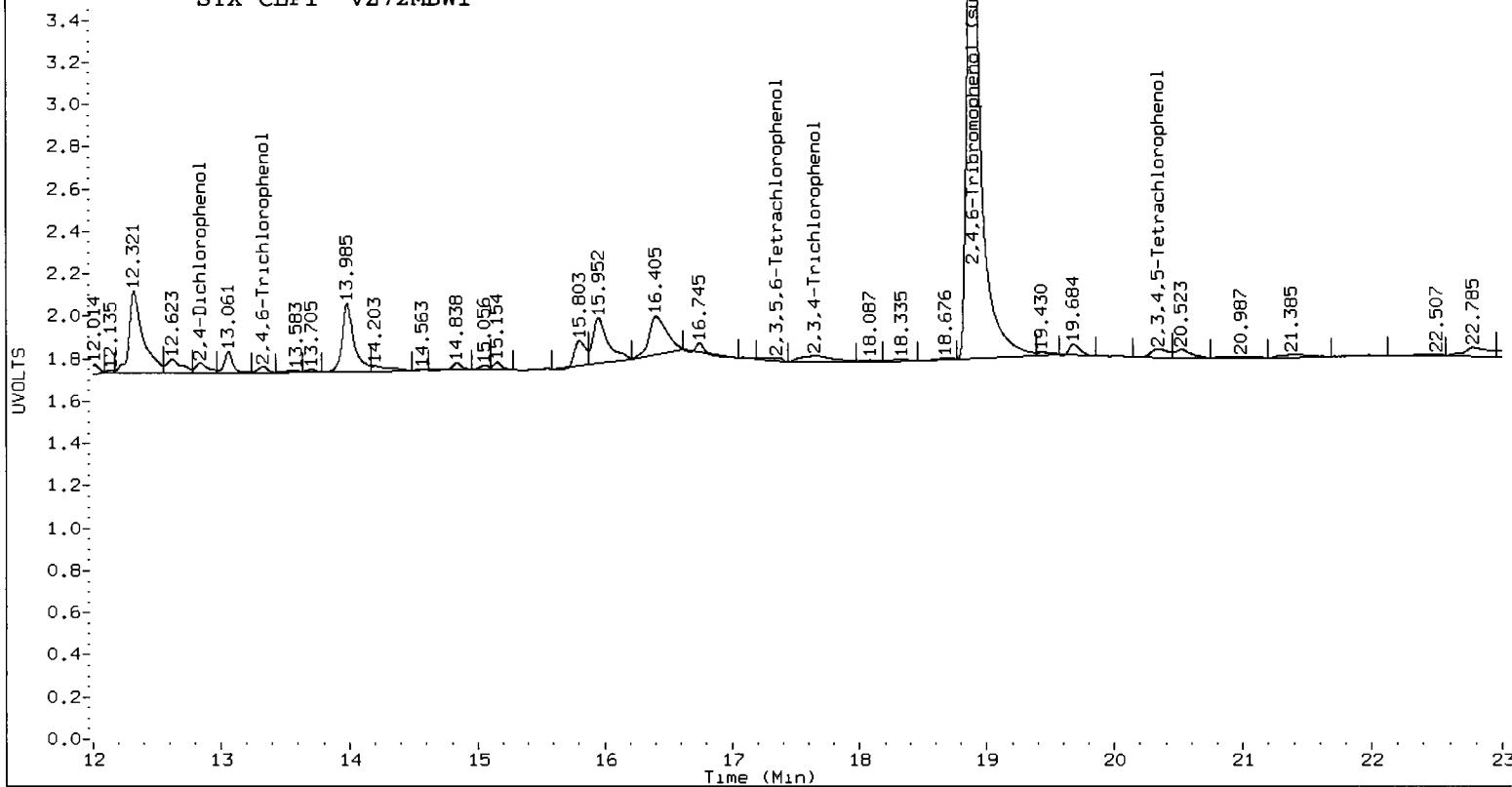
RT	STX CLP1 Col			STX CLP2 Col			STX CLP1 on col	STX CLP2 on col	RPD	Compound
	Shift	Response	RT	Shift	Response					
----			23.004	0.041	4032		0.0000	0.1130	---	Pentachlorophenol
13.330	-0.001	6935	14.385	0.047	3716		0.2177	0.1347	47.1*	2,4,6-Trichlorophenol
----			15.553	-0.029	4844		0.0000	0.1898	---	2,3,6-Trichlorophenol
----			----				0.0000	0.0000	---	2,4,5-Trichlorophenol
17.646	0.044	25911	----				0.8097	0.0000	---	2,3,4-Trichlorophenol
17.347	-0.045	7826	18.861	0.035	3194		0.1665	0.1106	40.4*	2,3,5,6-Tetrachlorophene
20.345	-0.062	19633	22.121	0.030	4365		0.4488	0.2106	72.2*	2,3,4,5-Tetrachlorophene
12.843	0.058	15696	----				< 8.5967	0.0000	---	2,4-Dichlorophenol
18.893	0.051	924370	20.982	0.038	560296		19.1	18.6	2.5	2,4,6-Tribromophenol (s)

PERCENT RECOVERY

COMPOUND	Col1	Col2
2,4,6-TBP (surr)	76.2	74.3

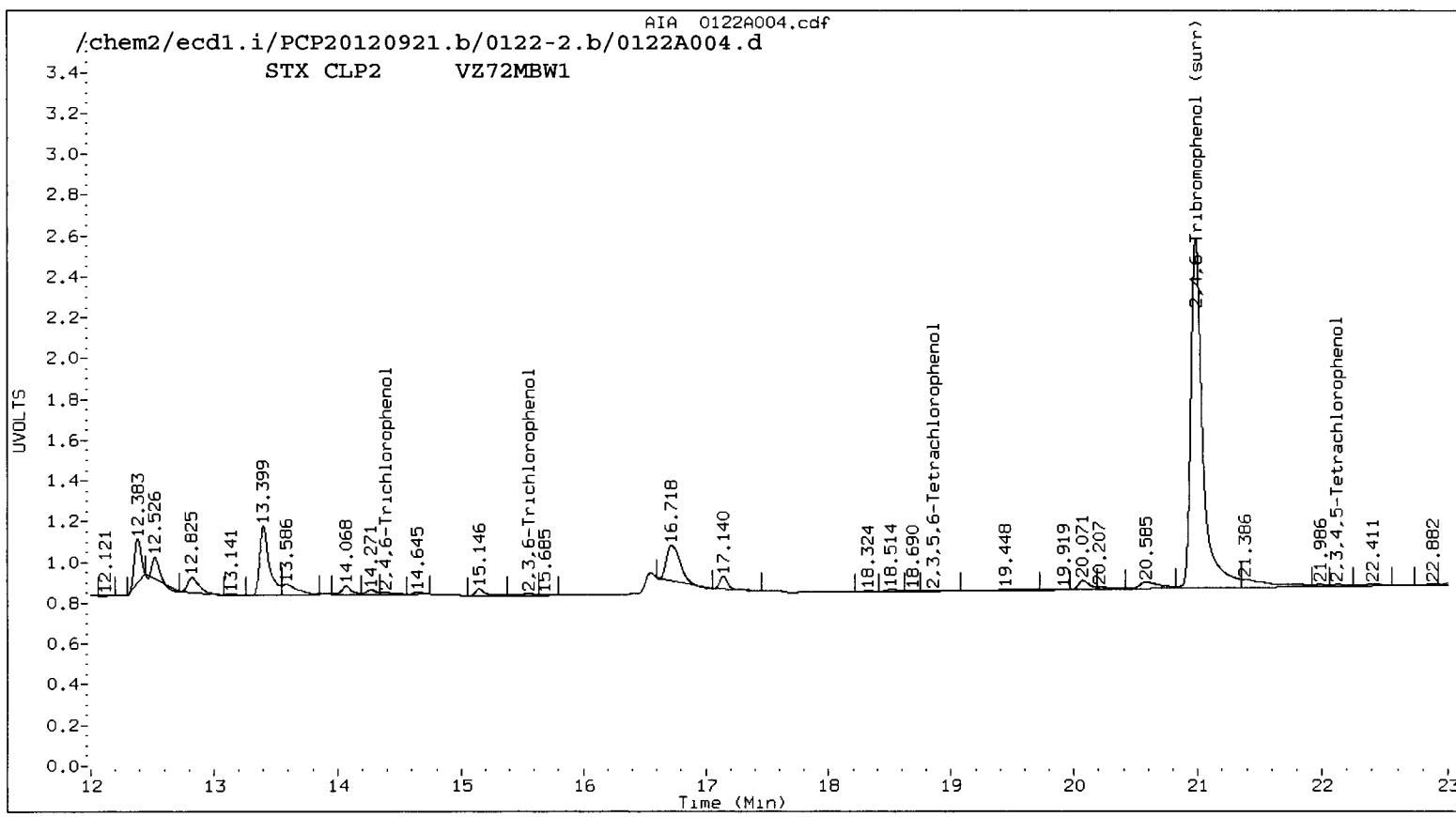
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STX CLP1 VZ72MBW1



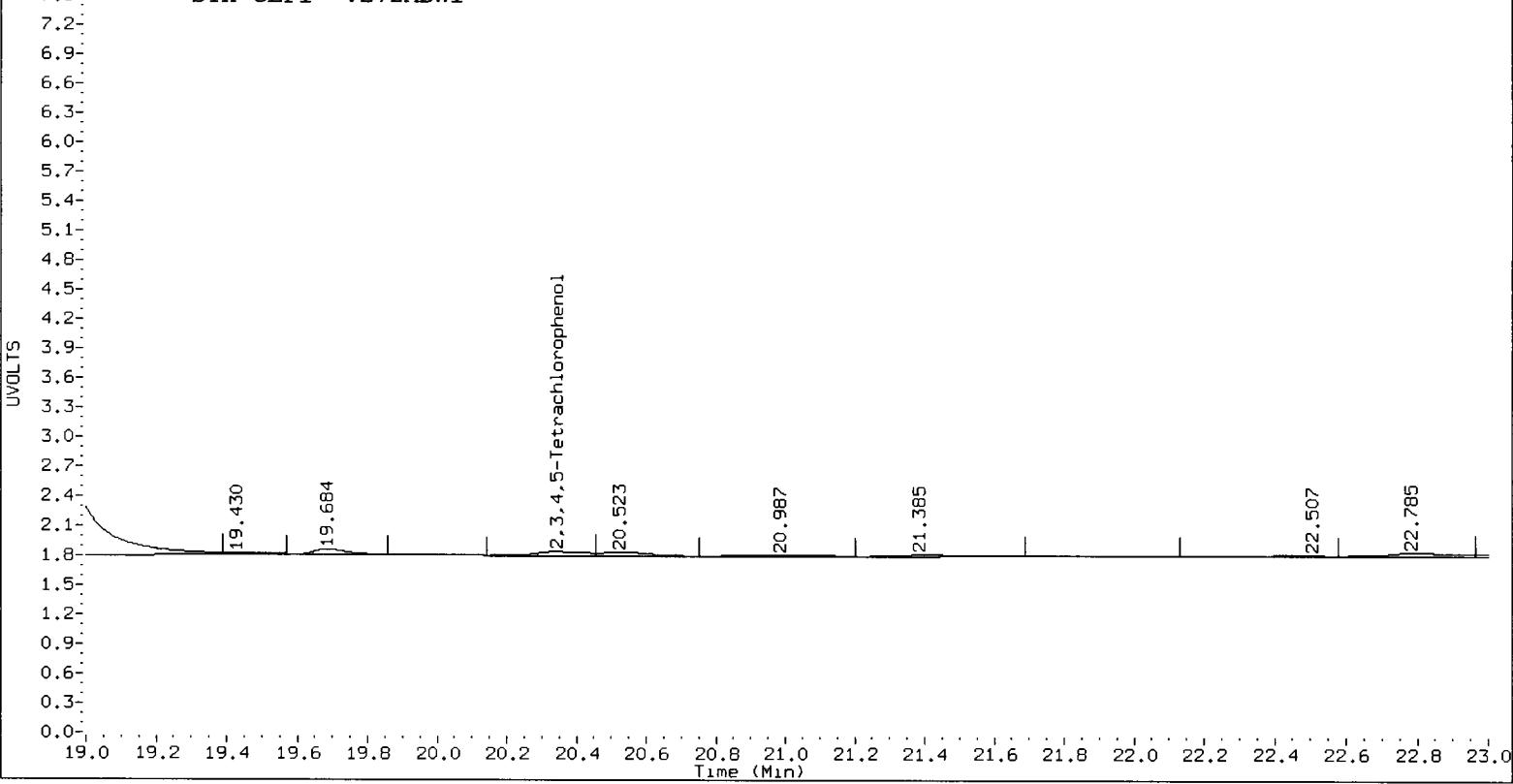
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STX CLP2 VZ72MBW1



/chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A004.cdf

STX CLP1 VZ72MBW1

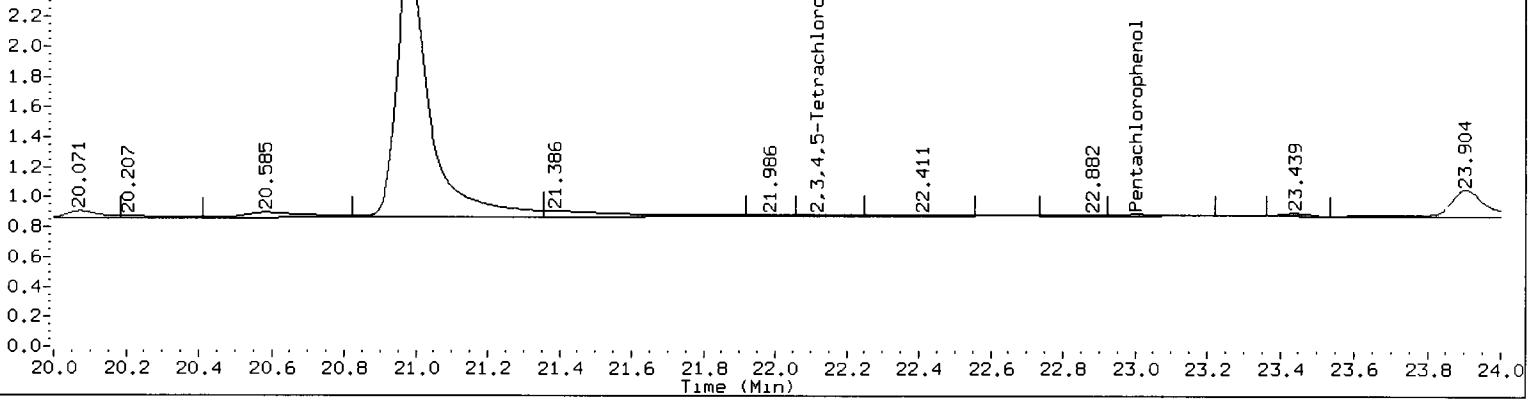


AIA 0122A004.cdf

/chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A004.d
STX CLP2 VZ72MBW1

UVOLTS

>>>2,4,6-Tribromophenol (surv)



VZ72 : 000042

Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Data file 1: /chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A005.d ARI ID: VZ72LCSW1 1/2 01/23/13
 Data file 2: /chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A005.d Client ID: VZ72LCSW1
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 22-JAN-2013 13:56
 Compound Sublist: pcpcal Report Date: 01/23/2013 10:56
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

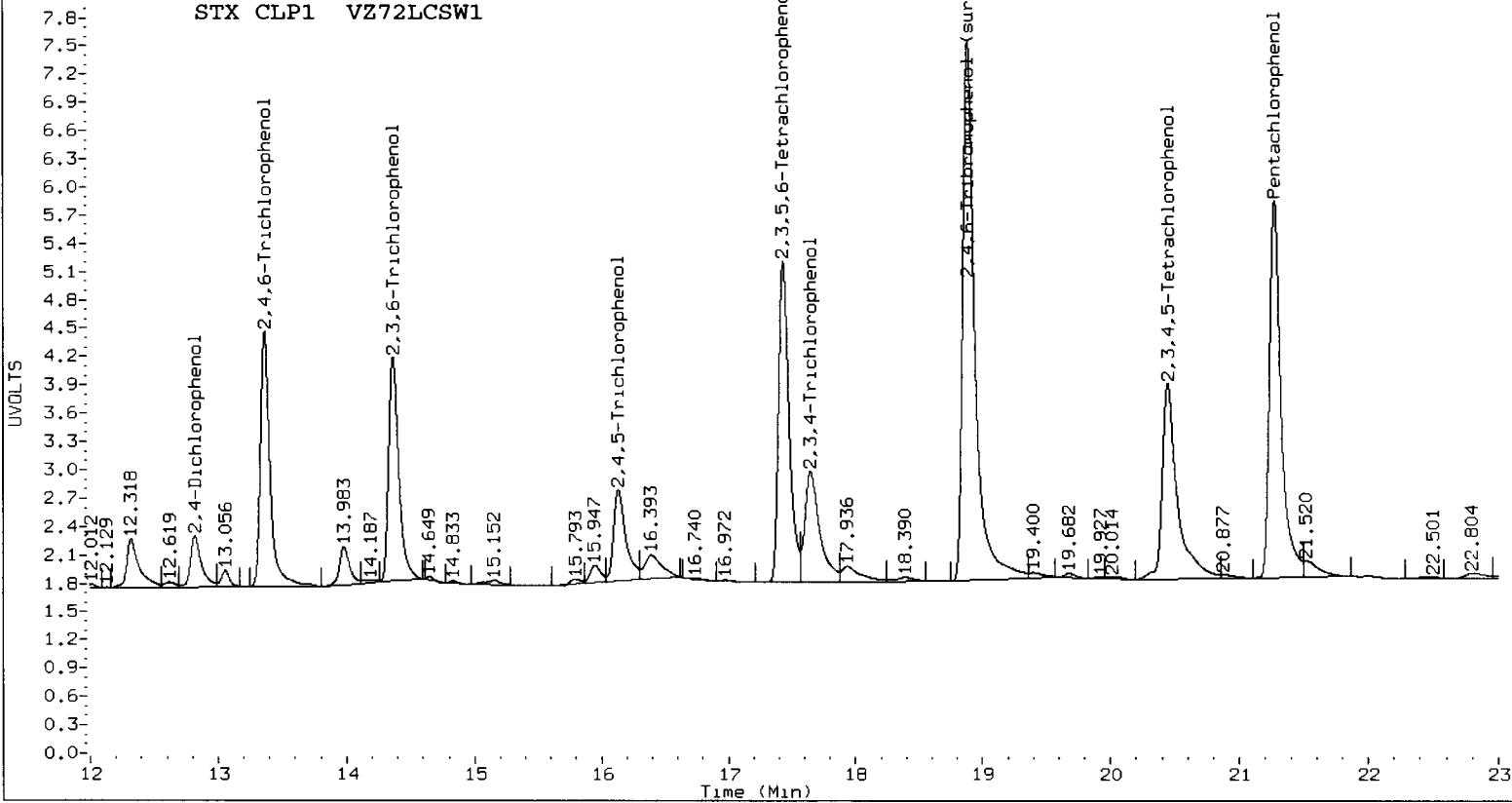
RT	STX CLP1 Col		STX CLP2 Col		on col	on col	RPD	Compound
	Shift	Response	RT	Shift	Response			
21.276	0.039	1271897	22.995	0.032	756141	21.7453	21.1870	2.6 Pentachlorophenol
13.362	0.032	703943	14.367	0.029	396504	22.1002	16.1632	31.0 2,4,6-Trichlorophenol
14.365	0.035	635410	15.613	0.031	380844	18.3598	17.4586	5.0 2,3,6-Trichlorophenol
16.134	0.045	327820	17.533	0.035	184162	14.9432	12.1712	20.4 2,4,5-Trichlorophenol
17.649	0.047	468252	19.082	0.036	318871	17.8904	17.9980	0.6 2,3,4-Trichlorophenol
17.433	0.040	993330	18.860	0.033	579551	21.1379	20.0637	5.2 2,3,5,6-Tetrachlorophene
20.449	0.042	854152	22.124	0.033	440724	23.2809	21.2678	9.0 2,3,4,5-Tetrachlorophene
12.820	0.035	161572	13.882	0.031	87876	94.1027	72.9674	25.3 2,4-Dichlorophenol
18.886	0.043	1952276	20.977	0.034	1194685	45.7	39.6	14.2 2,4,6-Tribromophenol

PERCENT RECOVERY

COMPOUND	Col1	Col2
Pentachlorophenol	87.0	84.7
2,4,6-Trichlorophenol	88.4	64.7
2,3,6-Trichlorophenol	73.4	69.8
2,4,5-Trichlorophenol	59.8	48.7
2,3,4-Trichlorophenol	71.6	72.0
2,3,5,6-Tetrachlorophenol	84.6	80.3
2,3,4,5-Tetrachlorophenol	93.1	85.1
2,4-Dichlorophenol	37.6	29.2
2,4,6-TBP (surr)	91.3	79.2

/chem2/ecdl.i/PCP20120921.b/0122-1.b/0122A005.cdf

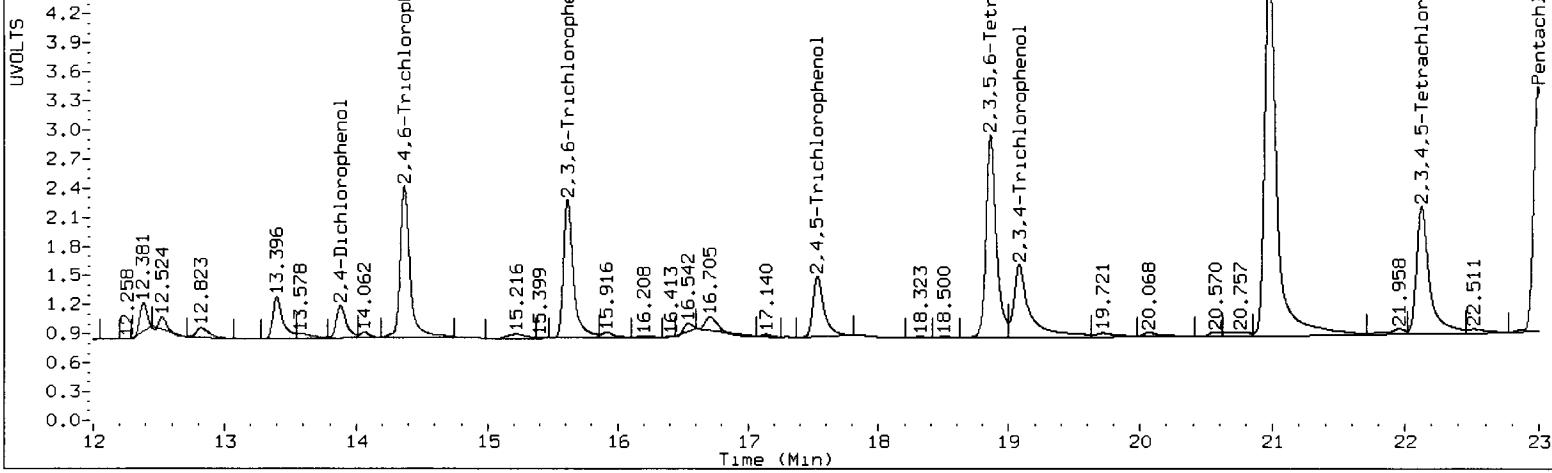
STX CLP1 VZ72LCSW1



AIA 0122A005.cdf

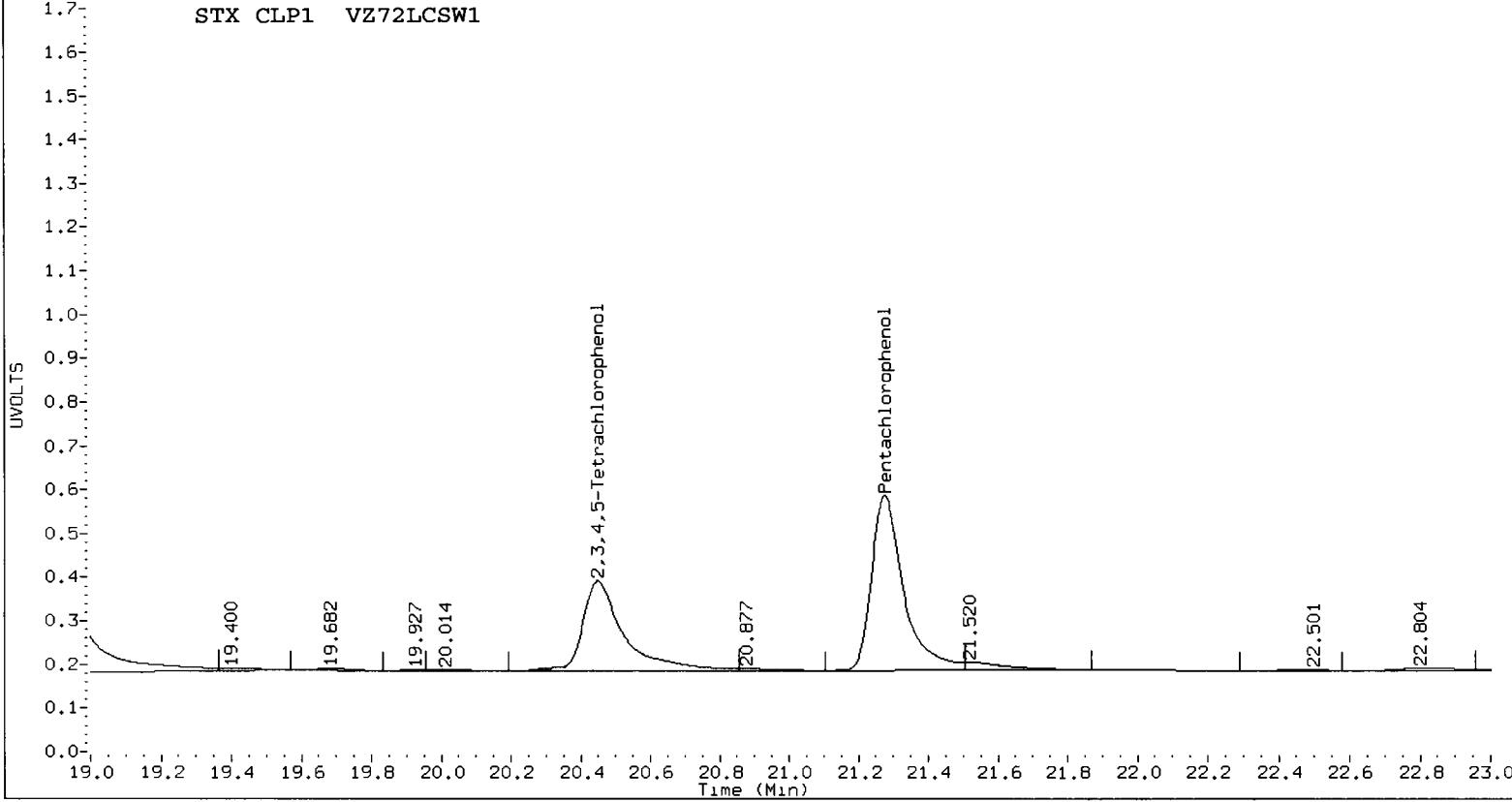
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STX CLP2 VZ72LCSW1



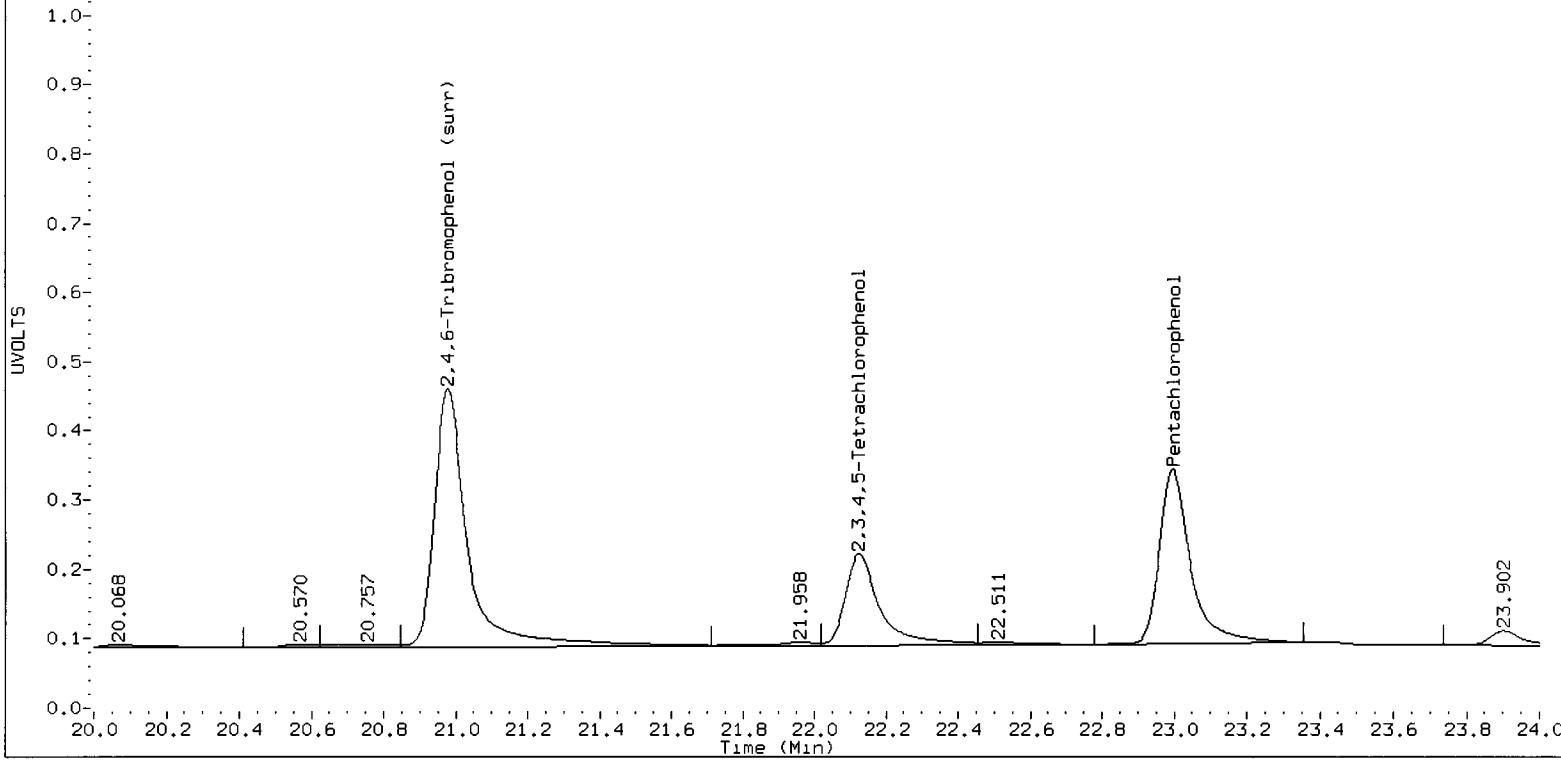
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STX CLP1 VZ72LCSW1



AIA 0122A005.cdf

/chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A005.d
STX CLP2 VZ72LCSW1



VZ72LCSW1

Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Data file 1: /chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A006.d ARI ID: VZ72LCSDW1
 Data file 2: /chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A006.d Client ID: VZ72LCSDW1
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 22-JAN-2013 14:32
 Compound Sublist: pcpca1 Report Date: 01/23/2013 10:56
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

VZ 01/23/13

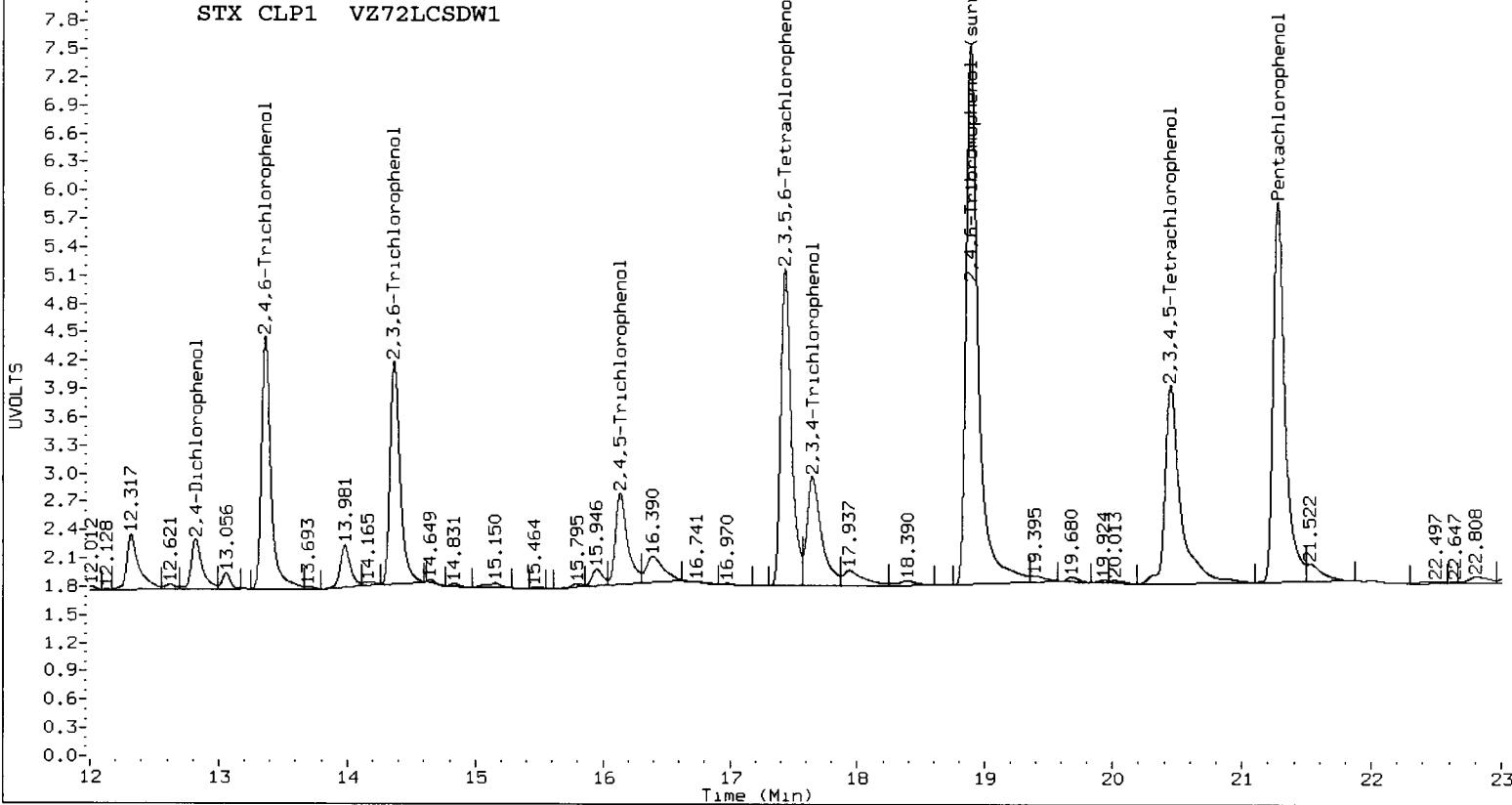
RT	STX CLP1 Col		STX CLP2 Col		on col	on col	RPD	STX CLP2 Compound
	Shift	Response	RT	Shift	Response			
21.274	0.038	1279983	22.995	0.031	763113	21.9010	21.3824	2.4 Pentachlorophenol
13.361	0.031	698629	14.367	0.029	400410	21.9333	16.3404	29.2 2,4,6-Trichlorophenol
14.365	0.035	633906	15.612	0.030	380715	18.3077	17.4518	4.8 2,3,6-Trichlorophenol
16.133	0.045	322197	17.533	0.035	181787	14.6330	11.9706	20.0 2,4,5-Trichlorophenol
17.648	0.046	455423	19.082	0.035	303302	17.3084	16.7907	3.0 2,3,4-Trichlorophenol
17.431	0.039	982215	18.859	0.032	578179	20.9013	20.0162	4.3 2,3,5,6-Tetrachlorophene
20.448	0.041	882537	22.123	0.032	439793	24.1865	21.2229	13.1 2,3,4,5-Tetrachlorophene
12.819	0.035	162051	13.882	0.030	86355	94.4002	71.5843	27.5 2,4-Dichlorophenol
18.885	0.043	1930592	20.977	0.033	1177442	45.0	39.0	14.2 2,4,6-Tribromophenol

PERCENT RECOVERY

COMPOUND	Col1	Col2
Pentachlorophenol	87.6	85.5
2,4,6-Trichlorophenol	87.7	65.4
2,3,6-Trichlorophenol	73.2	69.8
2,4,5-Trichlorophenol	58.5	47.9
2,3,4-Trichlorophenol	69.2	67.2
2,3,5,6-Tetrachlorophenol	83.6	80.1
2,3,4,5-Tetrachlorophenol	96.7	84.9
2,4-Dichlorophenol	37.8	28.6
2,4,6-TBP (surr)	90.1	78.1

/chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A006.cdf

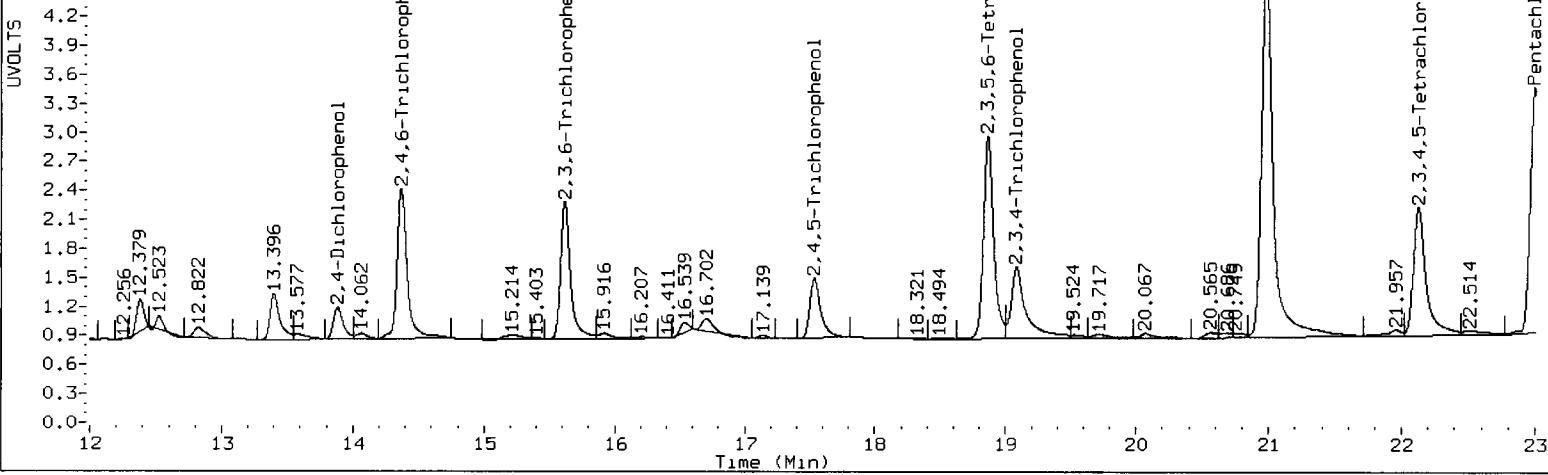
STX CLP1 VZ72LCSDW1



AIA 0122A006.cdf

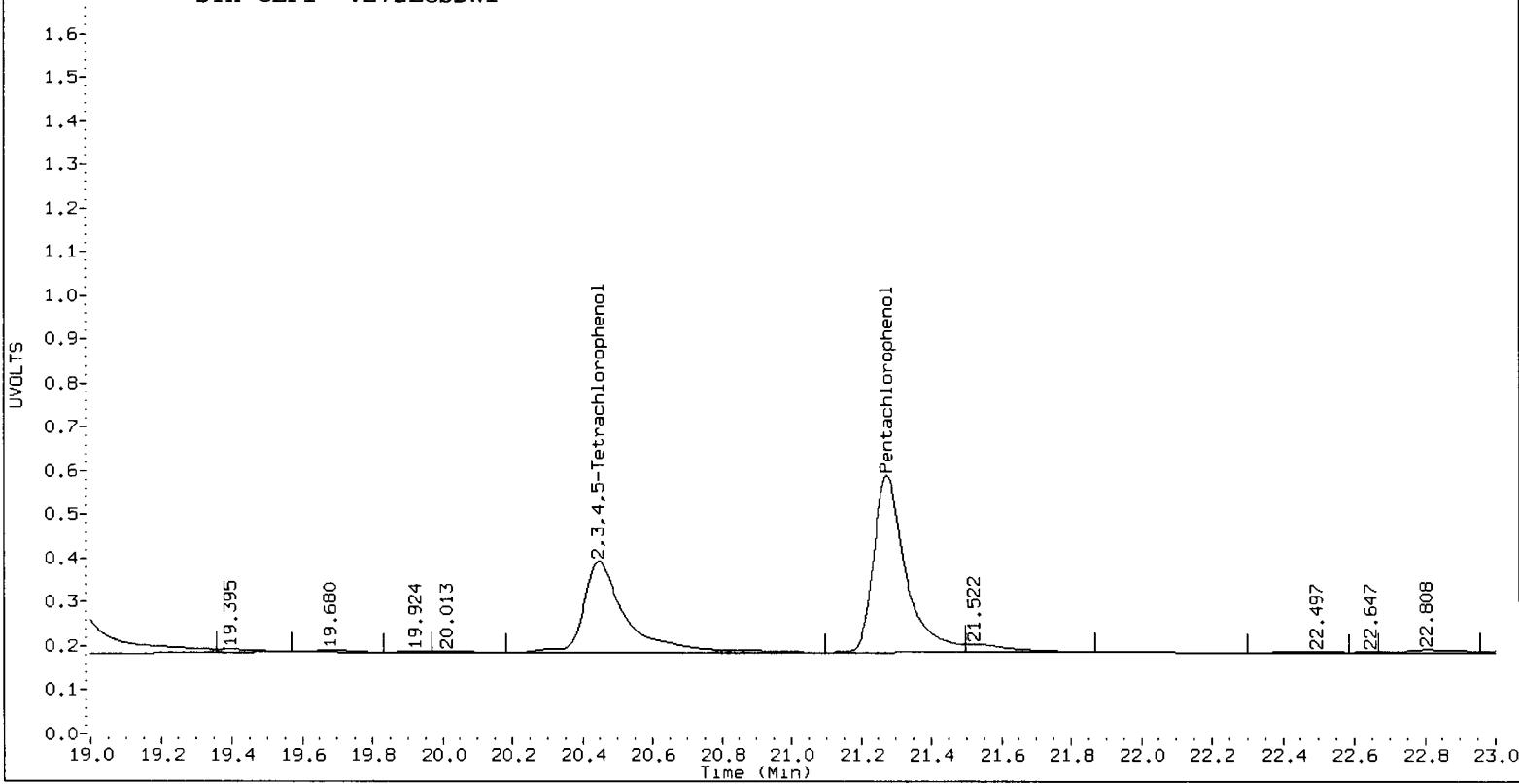
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STX CLP2 VZ72LCSDW1



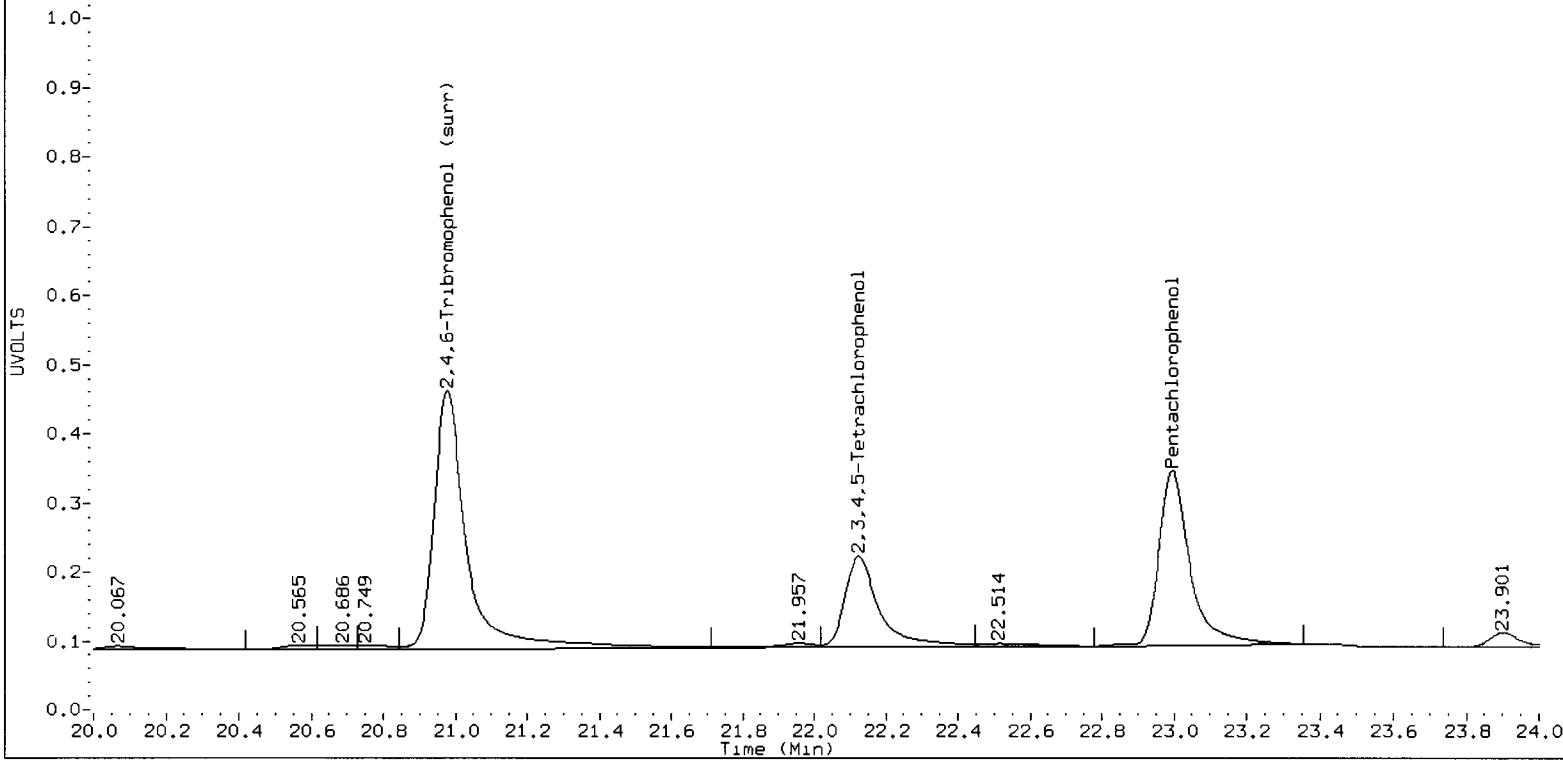
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STX CLP1 VZ72LCSDW1



AIA 0122A006.cdf

/chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A006.d
STX CLP2 VZ72LCSDW1



Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Data file 1: /chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A008.d ARI ID: VZ72A
 Data file 2: /chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A008.d Client ID: LLMW175-130111-W
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 22-JAN-2013 15:45
 Compound Sublist: pcpca1 Report Date: 01/23/2013 11:01
 Instrument: ecd1.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

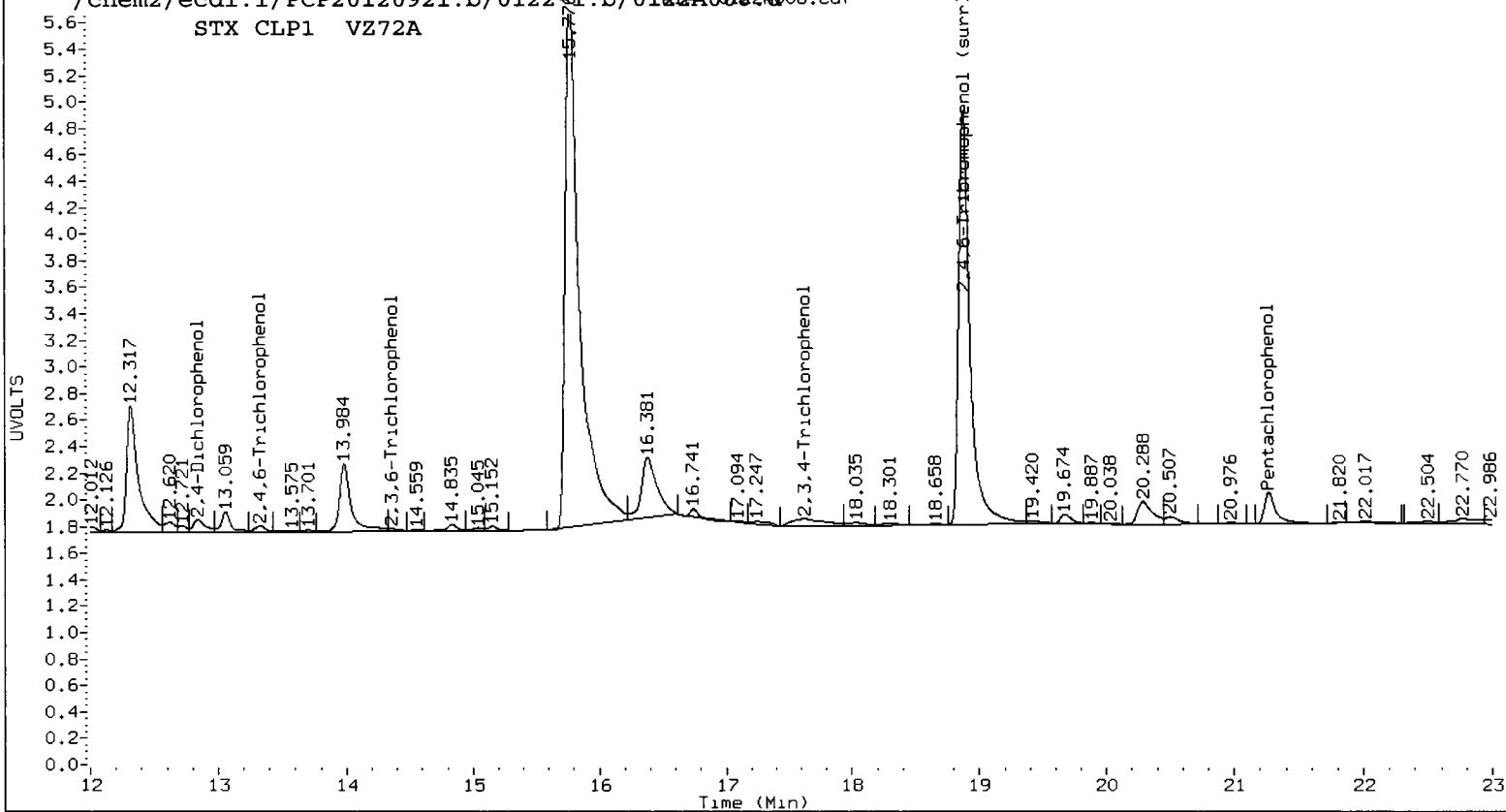
STX CLP1 Col			STX CLP2 Col			STX CLP1	STX CLP2		
RT	Shift	Response	RT	Shift	Response	on col	on col	RPD	Compound
21.271	0.035	74869	22.994	0.030	46446	1.1289	1.3014	14.2	Pentachlorophenol
13.327	-0.003	11486	14.377	0.039	5202	0.3606	0.1887	62.6*	2,4,6-Trichlorophenol
14.359	0.029	3564	15.539	-0.043	8836	0.0826	0.3467	123.0*	2,3,6-Trichlorophenol
----	----					0.0000	0.0000	---	2,4,5-Trichlorophenol
17.629	0.028	50285	----			1.5906	0.0000	---	2,3,4-Trichlorophenol
----			18.869	0.042	6713	0.0000	0.2324	---	2,3,5,6-Tetrachlorophene
----			22.108	0.017	819	0.0000	0.0395	---	2,3,4,5-Tetrachlorophene
12.843	0.058	29653	13.912	0.061	2774	16.3394	2.0863	154.7*	2,4-Dichlorophenol
18.883	0.040	987956	20.976	0.032	594075	20.5	19.7	4.1	2,4,6-Tribromophenol (s)

PERCENT RECOVERY

COMPOUND	Col1	Col2
2,4,6-TBP (surr)	82.1	78.8

/chem2/ecdl.i/PCP20120921.b/0122-1.b/0122A008.cdf

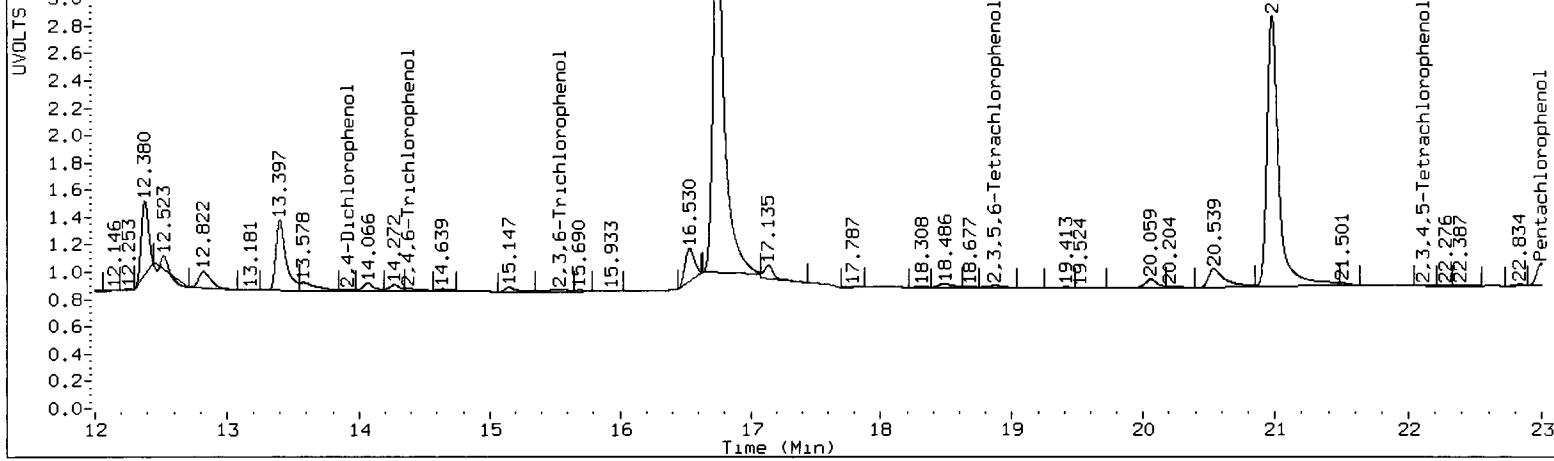
STX CLP1 VZ72A



AIA 0122A008.cdf

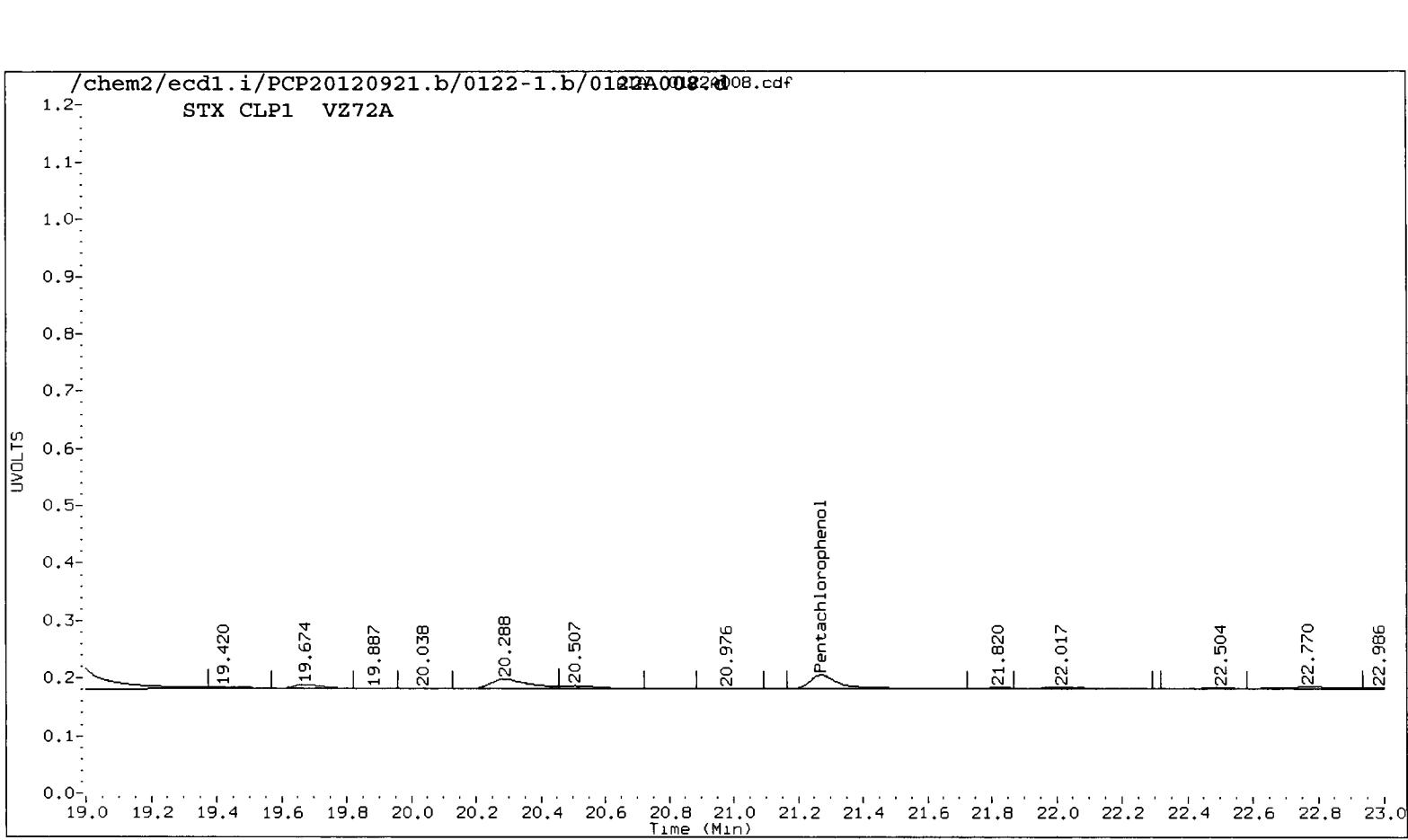
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STX CLP2 VZ72A



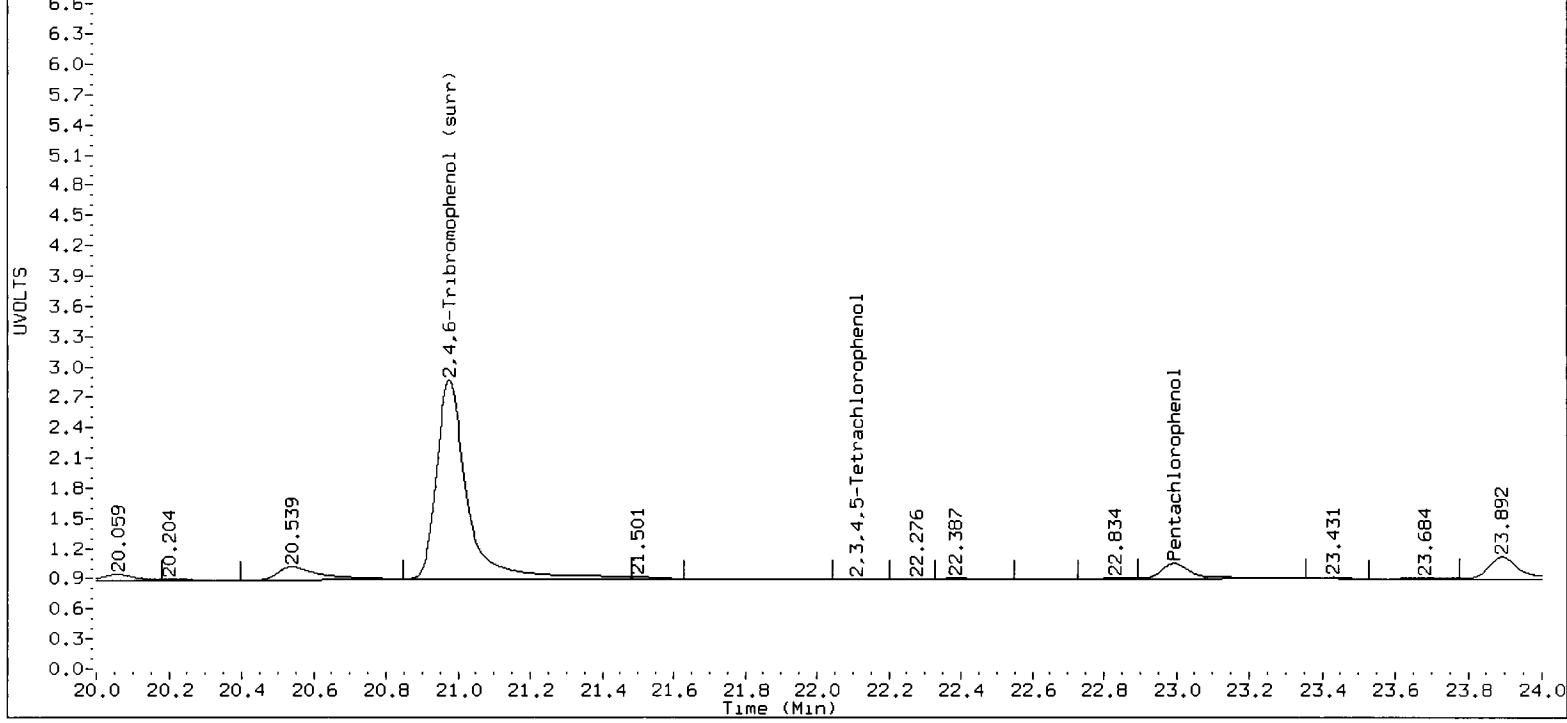
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STX CLP1 VZ72A



AIA 0122A008.cdf

/chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A008.d
STX CLP2 VZ72A



Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

12 02/23

Data file 1: /chem2/ecdl.i/PCP20120921.b/0122-1.b/0122A009.d ARI ID: VZ72B
 Data file 2: /chem2/ecdl.i/PCP20120921.b/0122-2.b/0122A009.d Client ID: LLMW17D-130111-W
 Method: /chem2/ecdl.i/PCP20120921.b/PCP.m Injection Date: 22-JAN-2013 16:21
 Compound Sublist: pcpcal Report Date: 01/23/2013 11:01
 Instrument: ecdl.i Matrix: WATER
 Operator: ar Dilution Factor: 1.000

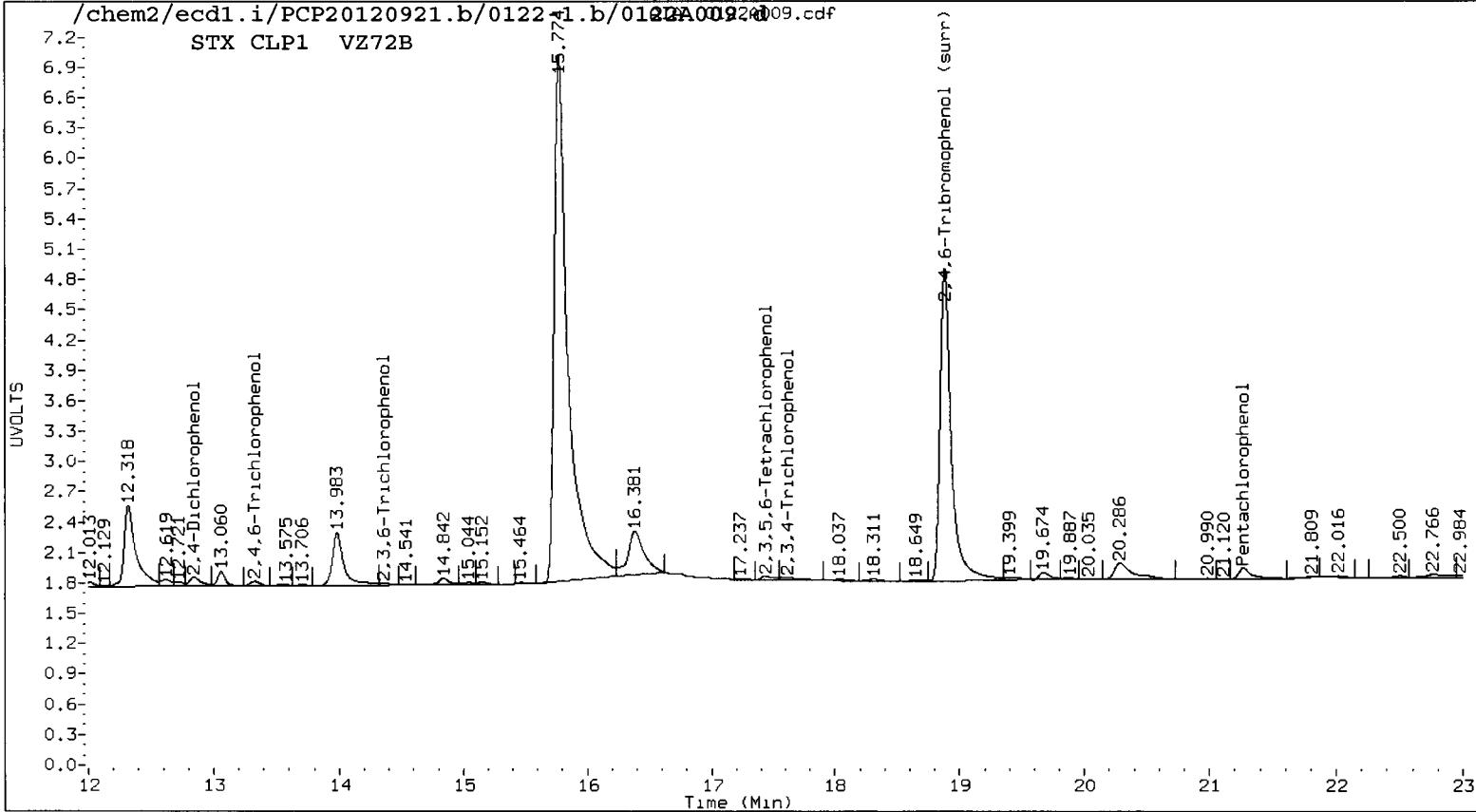
RT	STX CLP1 Col			STX CLP2 Col			on col	on col	RPD	STX CLP2 Compound
	Shift	Response	RT	Shift	Response					
21.271	0.034	36146	22.993	0.029	18711	0.5427	0.5243	3.4	Pentachlorophenol	
13.331	0.000	12882	14.277	-0.061	27794	0.4044	1.0153	86.1*	2,4,6-Trichlorophenol	
14.357	0.027	4648	15.610	0.028	5374	0.1078	0.2106	64.6*	2,3,6-Trichlorophenol	
----	----		----	----		0.0080	0.0000	---	2,4,5-Trichlorophenol	
17.615	0.013	13891	----	----		0.4314	0.0000	---	2,3,4-Trichlorophenol	
17.437	0.045	12513	18.865	0.038	9980	0.2663	0.3455	25.9	2,3,5,6-Tetrachlorophene	
----	----		22.109	0.017	1243	0.0000	0.0600	---	2,3,4,5-Tetrachlorophene	
12.844	0.059	27869	13.915	0.064	2191	15.3446	1.6466	161.2*	2,4-Dichlorophenol	
18.882	0.040	974023	20.975	0.031	596347	20.2	19.8	2.1	2,4,6-Tribromophenol (s)	

PERCENT RECOVERY

COMPOUND	Col1	Col2
2,4,6-TBP (surr)	80.8	79.1

/chem2/ecdl.i/PCP20120921.b/0122-1.b/0122A009.cdf

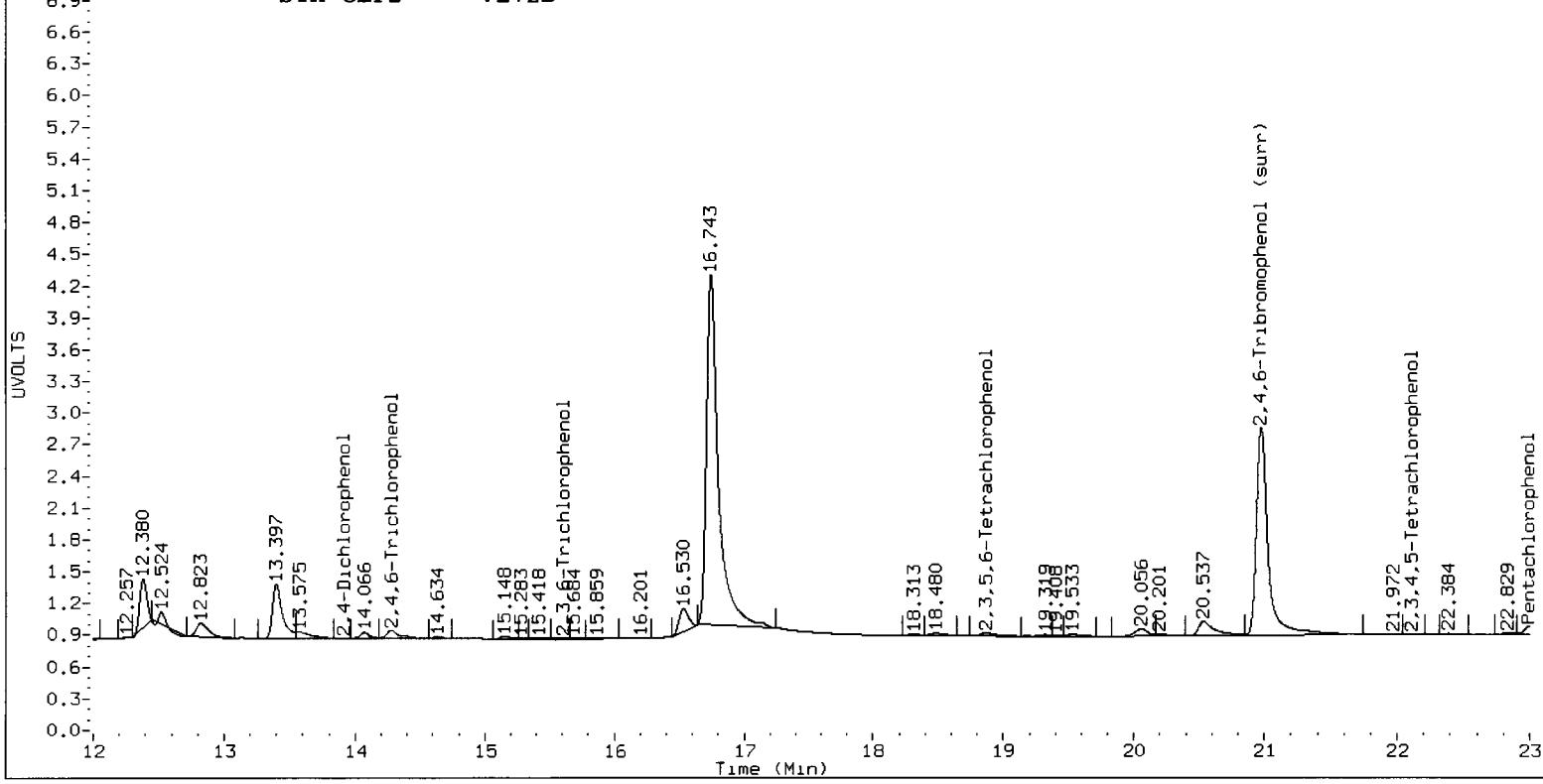
STX CLP1 VZ72B



AIA 0122A009.cdf

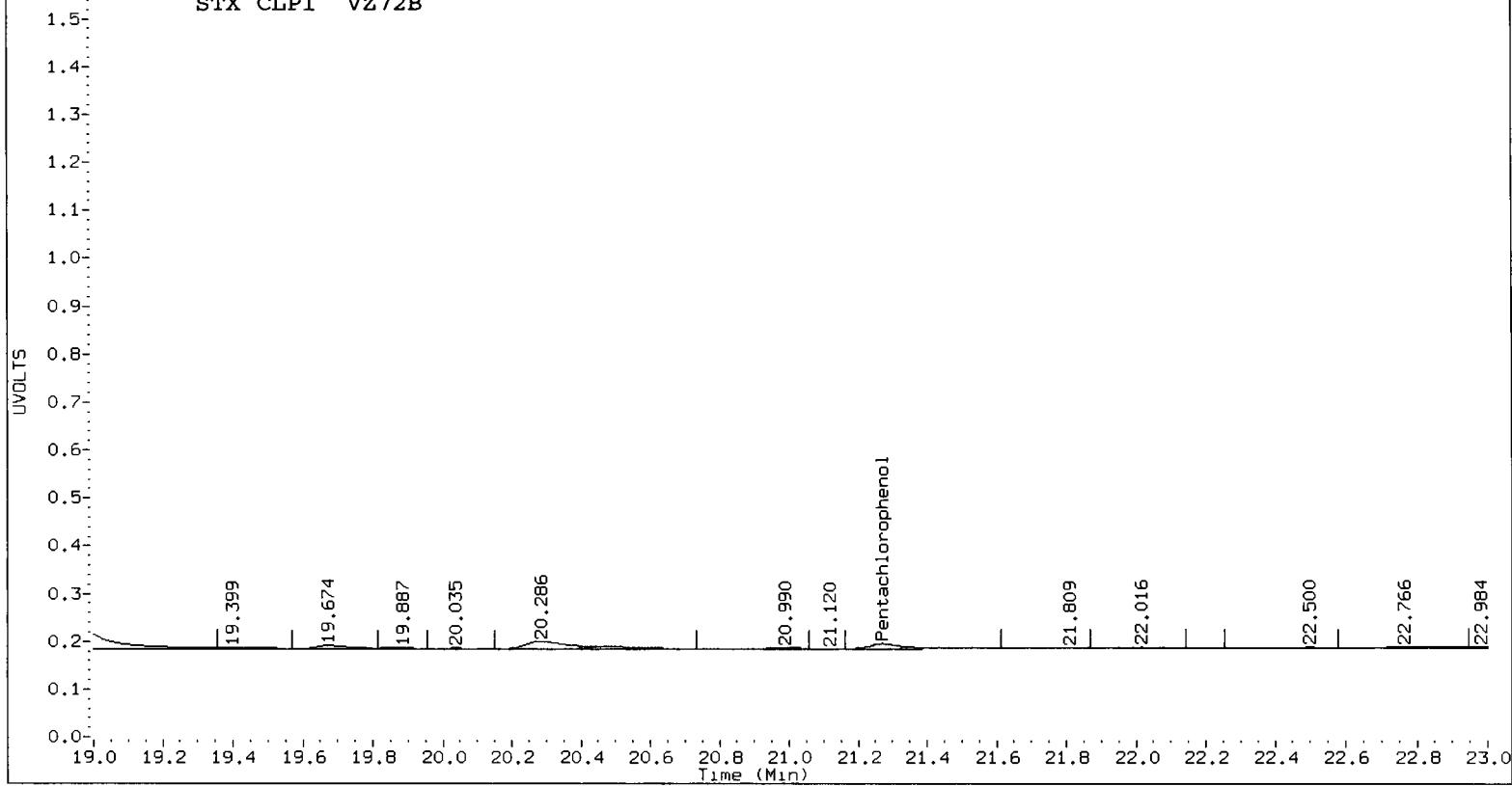
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STX CLP2 VZ72B



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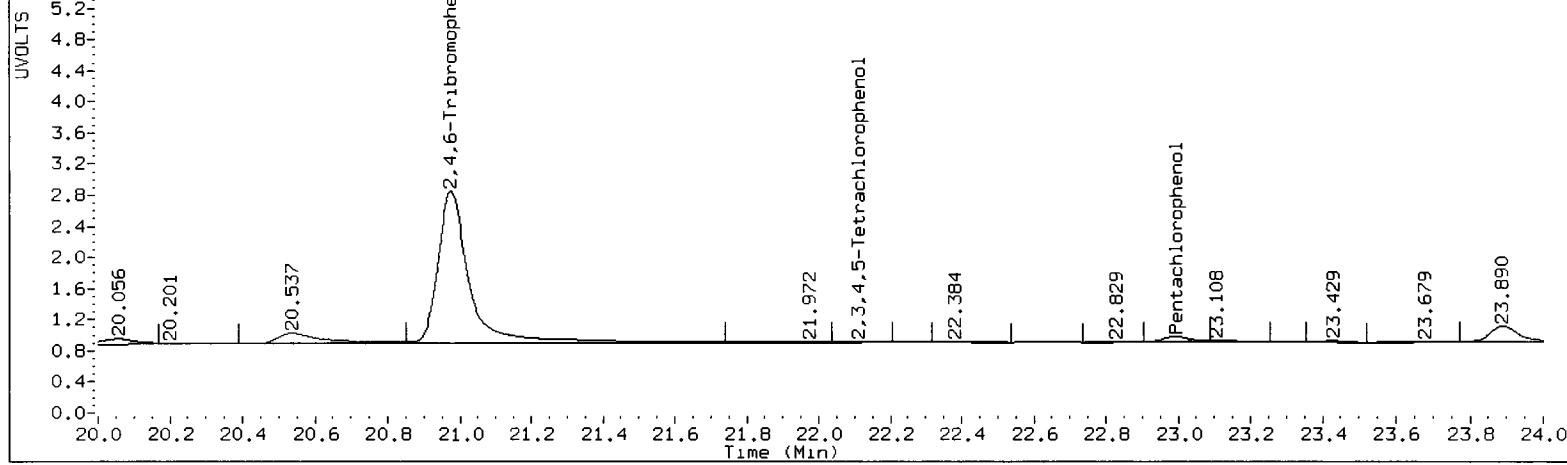
STX CLP1 VZ72B



AIA 0122A009.cdf

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STX CLP2 VZ72B



Analytical Resources Inc.
Dual Column 8041 Chlorinated Phenols Quantitation Report

Data file 1: /chem2/ecd1.i/PCP20120921.b/0122-1.b/0122A011.d ARI ID: PCPCCAL *Y2 01/23/13*
 Data file 2: /chem2/ecd1.i/PCP20120921.b/0122-2.b/0122A011.d Client ID:
 Method: /chem2/ecd1.i/PCP20120921.b/PCP.m Injection Date: 22-JAN-2013 17:33
 Compound Sublist: pcpca1 Report Date: 01/23/2013 11:16
 Instrument: ecd1.i Matrix: NONE
 Operator: ar Dilution Factor: 1.000

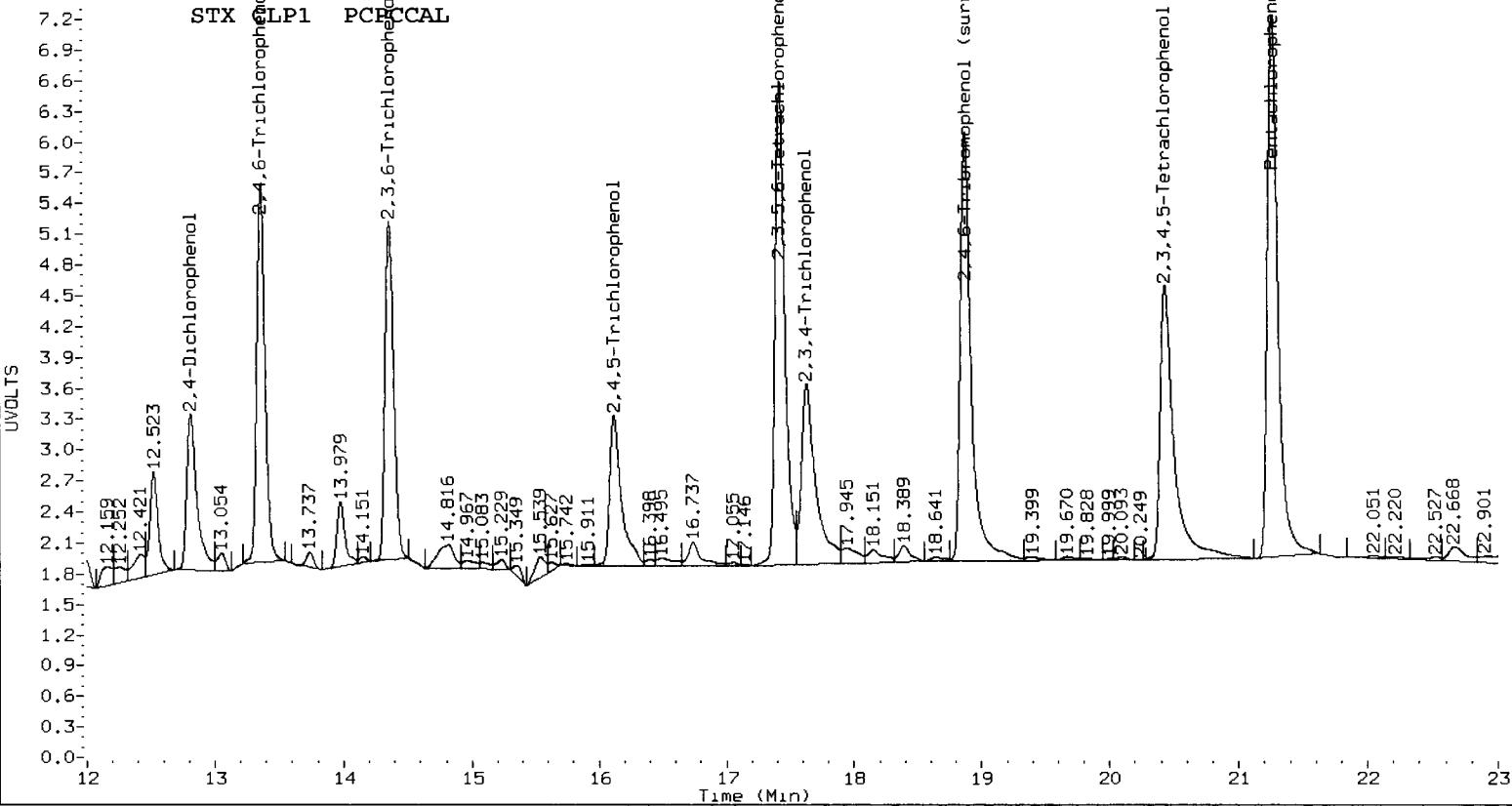
RT	STX CLP1 Col		STX CLP2 Col		on col	on col	RPD	STX CLP2 Compound
	Shift	Response	RT	Shift	Response			
21.267	0.031	1523920	22.990	0.026	992127	26.7017	27.7994	4.0 Pentachlorophenol
13.360	0.030	820568	14.365	0.027	559990	25.7616	23.8791	7.6 2,4,6-Trichlorophenol
14.361	0.031	754536	15.610	0.028	565397	22.6150	27.7701	20.5 2,3,6-Trichlorophenol
16.121	0.033	472096	17.526	0.028	287339	23.5454	21.9865	6.8 2,4,5-Trichlorophenol
17.635	0.033	676059	19.073	0.027	378150	28.0400	22.9037	20.2 2,3,4-Trichlorophenol
17.424	0.032	1265458	18.855	0.028	741998	26.9287	25.6875	4.7 2,3,5,6-Tetrachloroph
20.438	0.031	997483	22.118	0.026	556221	27.9409	26.8413	4.0 2,3,4,5-Tetrachlorophe
12.815	0.031	424066	13.879	0.028	239863	273.4789	232.6967	16.1 2,4-Dichlorophenol
18.875	0.032	1271994	20.971	0.027	784100	27.4	26.0	5.3 2,4,6-Tribromophenol (

PERCENT RECOVERY

COMPOUND	Col1	Col2
Pentachlorophenol	106.8	111.2
2,4,6-Trichlorophenol	103.0	95.5
2,3,6-Trichlorophenol	90.5	111.1
2,4,5-Trichlorophenol	94.2	87.9
2,3,4-Trichlorophenol	112.2	91.6
2,3,5,6-Tetrachlorophenol	107.7	102.8
2,3,4,5-Tetrachlorophenol	111.8	107.4
2,4-Dichlorophenol	109.4	93.1
2,4,6-TBP (surr)	109.6	104.0

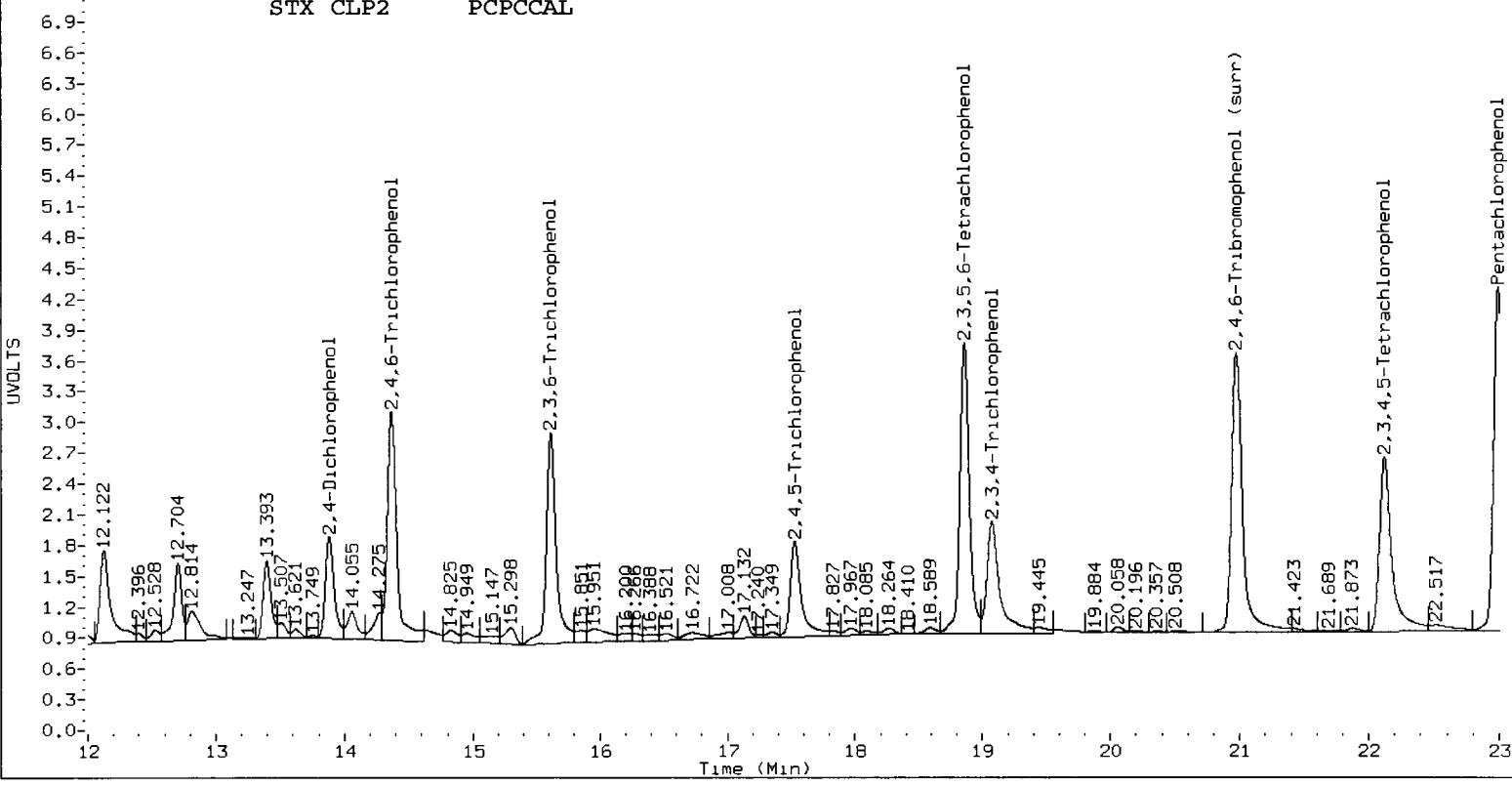
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STX CLP1 PCPCCAL



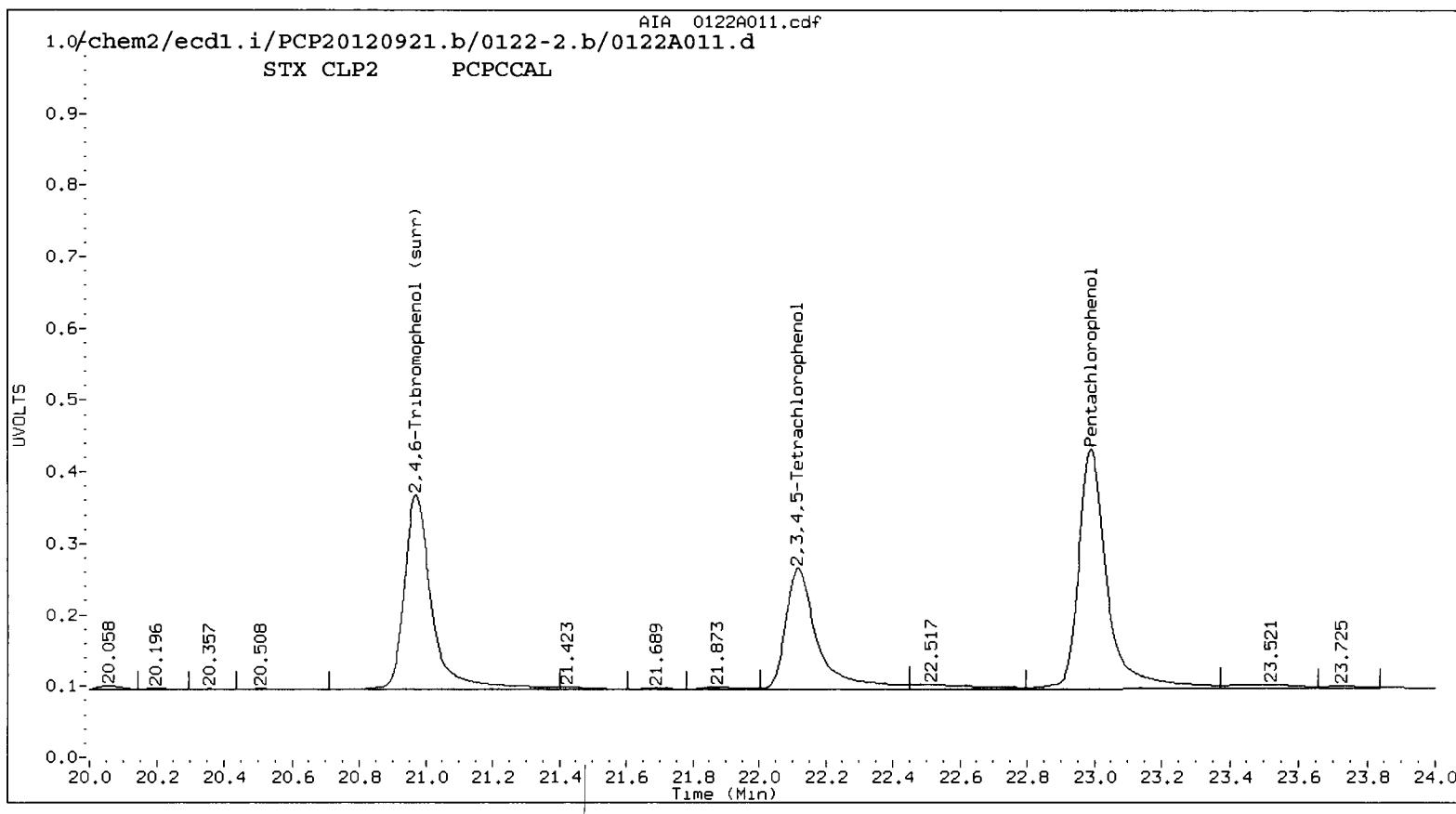
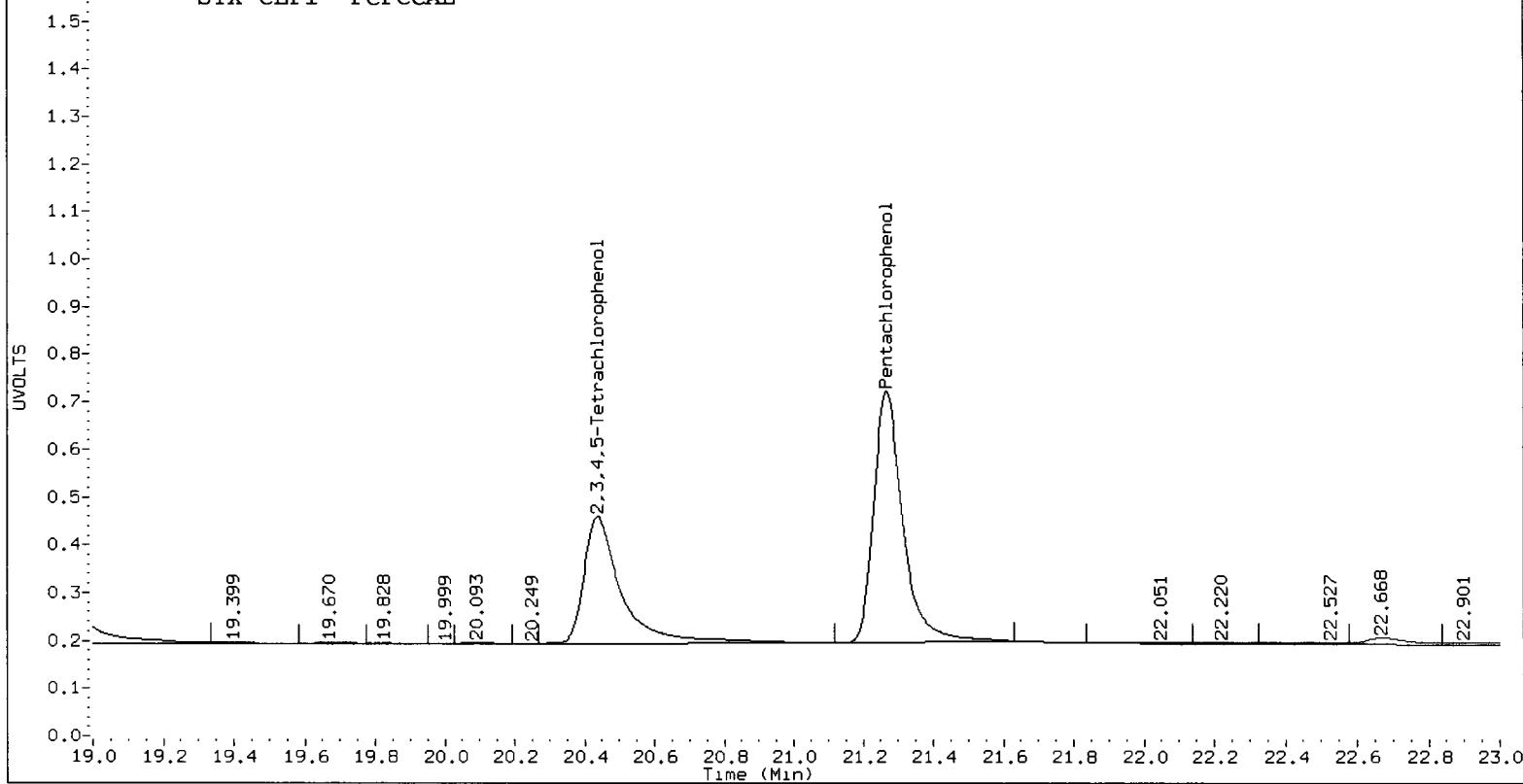
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STX CLP2 PCPCCAL



/chem2/ecdl.i/PCP20120921.b/0122-1.b/0122A011.cdf

STX CLP1 PCPCCAL



Attachment 2: Weyerhaeuser Results

Weyerhaeuser Analytical & Testing Services
 32901 Weyerhaeuser Way South
 Federal Way, WA 98003

Service Request 13-
 0088
 WA Cert.# C1219

Report
Everett East (Mill B) Sample Splits with ARI
Unit in ug/L
Method: 8151M

Client ID	Sample		Lab	Pentachlorophenol 87-86-5	Total Tetrachlorophenols	2,4,5 Trichlorophenol 88-06-2	DCAA	% Recovery		Date
	Date	Time						Surrogates	Extracted	
LLMW17-S	01/11/13	10:20	001	<0.50	<0.50	<0.50	103%	102%	01/16/13	01/18/13
LLMW17-D	01/11/13	10:50	002	<0.50	<0.50	<0.50	100%	92%	01/16/13	01/18/13
Method Blank				<0.50	<0.50	<0.50	98%	96%	01/16/13	01/18/13
Lab Control Spike			LCS	82%	158%	81%	95%	96%	01/16/13	01/18/13
LLMW17-S	01/11/13	10:20	001MS	89%	121%	91%	101%	93%	01/16/13	01/18/13
LLMW17-S	01/11/13	10:20	001MSD	77%	125%	84%	98%	91%	01/16/13	01/18/13