



Earth & Environmental

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AGRA Earth &
Environmental, Inc.
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Fairbanks, Alaska
U.S.A. 99701
Tel (907) 479-7586
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July 18, 1996

6-024-01173-1 Task 05

Seekins Ford-Lincoln-Mercury
1625 Old Steese Highway
Fairbanks, Alaska 99712

Attention: Al Haynes

RE: GROUNDWATER SAMPLING RESULTS, MAY 1996
SEEKINS FORD-LINCOLN-MERCURY, NRO File No. 100.26.131

Dear Al:

In May 1996, AGRA Earth & Environmental, Inc. (AEE) completed the first quarterly sampling event under the *Interim Corrective Action Plan (ICAP)* approved for Seekins Ford-Lincoln-Mercury in Fairbanks, Alaska. The plan was approved by the Alaska Department of Environmental Conservation (ADEC) and the Environmental Protection Agency (EPA) in April 1996. Groundwater sampling of the monitoring well network under the ICAP will again be conducted in August and November 1996 and in February 1997. This letter summarizes the results of the first sampling event. The locations of the monitoring wells are shown on the attached Figure 1.

FIELD METHODS

Prior to sampling, AEE field personnel measured the depth to water in each well using an Associated Remedial Technologies, Inc. Model 15101-E Hydrocarbon Interface Probe. This measurement was then subtracted from the surveyed elevation of the top of the well casing to yield the elevation of the water surface within the well. The survey data and measurements obtained during sample collection efforts are included in the attached Table 1.

During well sampling, AEE personnel used the decontamination, purging, and sampling procedures outlined in the ADEC *Underground Storage Tanks Procedures Manual*. Disposable polyethylene bailers were used to retrieve the water samples. One sample was collected as a Quality Assurance (QA) duplicate sample. The water samples and QA duplicate were placed in a chilled cooler and shipped to the AEE Environmental Chemistry Laboratory in Portland, Oregon.

RESULTS

The original samples and the QA duplicate were analyzed for benzene, toluene, ethylbenzene, and xylenes (BTEX) by Alaska Test Method AK101, gasoline range petroleum hydrocarbons (GRPH) by Alaska Test Method AK101, diesel range petroleum hydrocarbons (DRPH) by Alaska Test Method AK102, and volatile organic compounds (VOCs) by EPA Method 624. In addition, select samples were analyzed for polynuclear aromatic hydrocarbons (PAHs) based on the proximity of the monitoring wells to the former injection well area. Copies of the laboratory analytical reports for these samples are included with this letter. In addition to the above samples, AEE field personnel submitted a trip blank sample, which accompanied the samples throughout the sampling event. This sample exhibited non-detectable concentrations of all tested compounds. Sampling data and analytical results are summarized in Tables 1 and 2.

The QA duplicate (DUP-1) from monitoring well MW-1 was submitted as a quality control indicator. The analytical results for the two samples showed close agreement for all tested compounds. The reported results indicated a relative percent difference (RPD) for the two samples that is within the allowable RPD limit of 30 percent for all tested compounds.

CLOSURE

AEE is pleased to be of continued service to Seekins. We will collect water level measurements at each of the well locations on a monthly basis. These data will be included in future quarterly letter reports. Our next sampling event is scheduled to occur in August 1996. We will forward the results of testing as soon as we have compiled that information. If you have any questions or comments regarding this letter, please contact our office.

Sincerely,

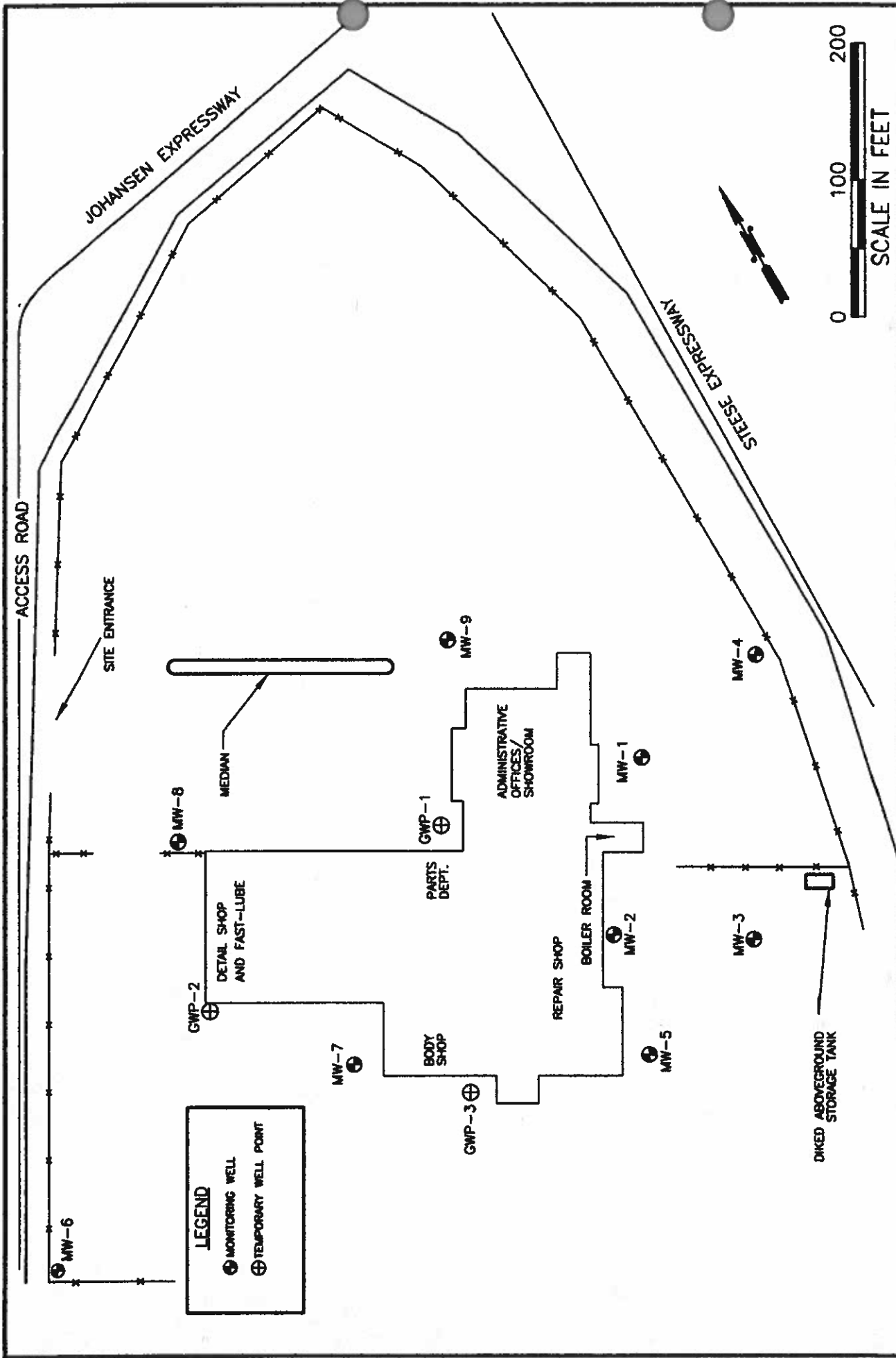
AGRA Earth & Environmental, Inc.



James A. Spontak
Project Manager

Atch: Figure 1, Site Plan
 Tables 1 and 2, Summary of Groundwater Analytical Data
 Laboratory Report

c: Chet Halcomb, ADEC - Fairbanks
 Jonathan Williams, EPA - Seattle



LEGEND
 ⊕ MONITORING WELL
 ⊕ TEMPORARY WELL POINT

**SEEKINS FORD-LINCOLN-MERCURY
 1626 OLD STEESE HIGHWAY
 FAIRBANKS, ALASKA**

W.O. 6-024-01173-1
 DESIGN MPP
 DRAWN MPP / LOCWELL
 DATE 06/06/96
 SCALE 1"=100'

AGRA
Earth & Environmental
 3504 Industrial Ave. Suite 5
 Fairbanks, Alaska 99701

**SITE PLAN
 FIGURE 1**

TABLE 1
Summary of Groundwater Analytical Data for Seekins Ford-Lincoln-Mercury
 BTEX, GRPH, and DRPH Results

Well Number	Date	MP Elevation	DTW	Ground Water Elevation	EPA 8020 (ug/L)	EPA 8100 (mg/L)		
					AK101	Total BTEX	GRPH	DRPH
Maximum Contaminant Level (ug/L)					5			
MW-1	7/21/95	97.13	14.20	82.93	12,000	71,300	180	5.4
	5/1/96	97.11	15.70	81.41	7,500	97,300	240	6.2
MW-2	7/21/95	97.57	14.64	82.93	ND(0.5)	6.8	0.15	0.35
	5/1/96	97.57	16.13	81.44	1.69	26.3	0.26	0.74
MW-3	7/21/95	96.62	13.70	82.92	8	NT	2.8	13
	5/1/96	96.62	15.18	81.44	ND(1)	151	0.99	5.4
MW-4	7/21/95	95.88	12.93	82.95	ND(0.5)	ND	ND(0.05)	ND(0.1)
	5/1/96	95.88	14.43	81.45	ND(1)	ND	ND(0.05)	0.24
MW-5	7/21/95	---	---	---	MONITORING WELL INSTALLED IN APRIL 1996.			
	5/1/96	97.08	15.68	81.40	2.49	33.9	0.17	1.1
MW-6	7/21/95	---	---	---	MONITORING WELL INSTALLED IN APRIL 1996.			
	5/1/96	97.14	15.78	81.36	ND(1)	23.3	0.12	0.94
MW-7	7/21/95	---	---	---	MONITORING WELL INSTALLED IN APRIL 1996.			
	5/1/96	97.70	16.29	81.41	ND(1)	48.3	0.26	0.47
MW-8	7/21/95	---	---	---	MONITORING WELL INSTALLED IN APRIL 1996.			
	5/1/96	97.85	16.49	81.36	8.39	110.2	0.35	0.69
MW-9	7/21/95	---	---	---	MONITORING WELL INSTALLED IN APRIL 1996.			
	5/1/96	97.37	15.95	81.42	ND(1)	8.8	0.06	0.84
GWP-1	7/21/95	97.47	14.62	82.85	1,500	1,722	4	0.19
	5/1/96	97.53	16.11	81.42	117	134.3	0.34	0.48
GWP-2	7/21/95	97.75	15.02	82.73	ND(0.5)	ND	ND(0.05)	ND(0.1)
	5/1/96	97.91	16.54	81.37	ND(1)	ND	ND(0.05)	0.35
GWP-3	7/21/95	97.02	14.18	82.84	ND(0.5)	NT	ND(0.05)	ND(0.1)
	5/1/96	97.14	15.71	81.43	ND(1)	ND	ND(0.05)	0.17
DUP-1	5/1/96	---	---	---	10,000	124,600	220	5.6

NT INDICATES THAT THESE MONITORING WELLS WERE NOT TESTED FOR THIS PARAMETER.
 ND INDICATES THAT THE RESULT WAS NON-DETECTABLE ABOVE THE LIMIT SHOWN.

TABLE 2
Summary of Groundwater Analytical Data for Seekins Ford-Lincoln-Mercury
VOC and PAH Compounds

Well Number	Date	EPA 824 (ug/L)																	EPA 8310 (ug/L)									
		Trichloro-fluoro-methane	Tetra-chloro-ethene	1,2-Dichloro-benzene	Tri-chloro-ethene	1,1,1-Trichloro-ethane	1,1-Dichloro-methane	Chloroform	Carbon Tetrachloride	Acetone	Methylene Chloride	2-Butanone (MEK)	Isopropyl Benzene	1,2,3-Trichloro propane	n-Propyl benzene	1,3,5-Trimethyl benzene	1,2,4-Trimethyl benzene	sec-Butyl benzene	4-Isopropyl toluene	n-Butyl benzene	Naphthalene	Fluorene	Phenanthrene	Fluoranthene	Pyrene	Naphthalene		
MM-1	7/21/95	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	
MM-2	7/21/95	6.3	20	1.71	ND(3)	57	9.39	ND(3)	1.88	ND(10)	ND(1)	8.2	ND(40)	NT	ND(20)	NT	ND(10)	3.53	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
	5/1/96	ND(1)	20	1.71	ND(3)	57	9.39	ND(3)	1.88	ND(10)	ND(1)	8.2	ND(40)	NT	ND(20)	NT	ND(10)	3.53	NT	NT	NT	NT	NT	NT	NT	NT	NT	
MM-3	7/21/95	8.8	29	90	12	6.5	2.40	ND(3)	2.40	ND(3)	5.5	ND(3)	ND(40)	NT	ND(20)	NT	ND(10)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
	5/1/96	ND(1)	23	86	22	6.5	2.40	ND(3)	2.40	ND(3)	5.5	ND(3)	ND(40)	NT	ND(20)	NT	ND(10)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
MM-4	7/21/95	ND(3)	ND(3)	ND(3)	ND(3)	ND(3)	ND(10)	5.5	ND(10)	5.5	ND(1)	ND(3)	ND(40)	NT	ND(20)	NT	ND(10)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(10)	5.5	ND(10)	5.5	ND(1)	ND(3)	ND(40)	NT	ND(20)	NT	ND(10)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
MM-5	7/21/95	ND(1)	ND(1)	2.71	ND(1)	1.45	NT	ND(1)	1.01	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	ND(1)	ND(1)	2.71	ND(1)	1.45	NT	ND(1)	1.01	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
MM-6	7/21/95	10.1	5.23	ND(1)	ND(1)	2.28	2.93	NT	1.01	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	3.11	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	
	5/1/96	10.1	5.23	ND(1)	ND(1)	2.28	2.93	NT	1.01	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	3.11	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	
MM-7	7/21/95	13.4	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	13.4	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
MM-8	7/21/95	16.4	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	16.4	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
MM-9	7/21/95	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	1.85	NT	NT	NT	NT	NT	NT	NT	NT	NT	
GWP-1	7/21/95	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	3.70	NT	NT	NT	NT	NT	NT	NT	NT	NT	
GWP-2	7/21/95	33.4	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	33.4	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	3.70	NT	NT	NT	NT	NT	NT	NT	NT	NT	
GWP-3	7/21/95	2.22	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	2.22	ND(1)	ND(1)	ND(1)	ND(1)	NT	ND(1)	ND(1)	ND(1)	ND(1)	ND(1)	ND(20)	NT	ND(10)	NT	ND(1)	4.91	NT	NT	NT	NT	NT	NT	NT	NT	NT	
DUP-1	7/21/95	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT	
	5/1/96	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	NT	ND(300)	ND(300)	ND(300)	ND(300)	ND(300)	ND(6000)	NT	ND(3000)	NT	ND(300)	NT	ND(300)	NT	NT	NT	NT	NT	NT	NT	NT	

NT INDICATES THAT THE MONITORING WELL WAS NOT TESTED FOR THIS PARAMETER.
ND INDICATES THAT THE PARAMETER WAS NON-DETECTABLE ABOVE THE LIMIT SHOWN.
ONLY THOSE ANALYTES REPORTED IN DETECTABLE CONCENTRATIONS ARE LISTED IN THIS TABLE. ALL OTHER VOC AND PAH COMPOUNDS WERE NON-DETECTABLE.



AGRA Earth & Environmental, Inc.
7477 SW Tech Center Drive
Portland, Oregon
U.S.A. 97223-8025
Tel (503) 639-3400
Fax (503) 620-7892

May 14, 1996

AGRA Earth & Environmental
3504 Industrial Way, Suite 5
Fairbanks, AK 99701

Attention: Mr. James Spontak

Dear Mr. Spontak:

RE: Analytical Results For Project 6-024-01173-1T04

Attached are the results for the samples submitted on May 3, 1996 from the above referenced project. For your reference, our project number associated with these samples is AK960277.

The samples were analyzed for volatile organic compounds by GC/MSD, diesel range organics, and gasoline range organics at the AGRA Earth & Environmental Portland Chemistry Laboratory. Samples MW-2, MW-3, and GWP-3 were subcontracted to American Environmental Network (NET) for PAH analysis. The NET data will be reported under separate cover at a later date.

All analyses were conducted in accordance with applicable QA/QC guidelines. The results apply only to the samples submitted.

Please feel free to contact me if you have any questions regarding this report, or if I can be of any assistance in any other matter.

Respectfully submitted,

AGRA Earth & Environmental

A handwritten signature in cursive script that reads 'Sean Gormley'.

Sean Gormley
Laboratory Manager
Laboratory ID # UST-008

Project: Seekins Ford - Lincoln Mercury
 Project No.: 6-024-01173-0T04
 Project Manager: James Spontak
 Sample Matrix: Water

Service Request No.: AK960277
 Report Date: 5/9/96
 Report No.: 96027703
 C.O.C. No.: 02920

Volatile Organic Compounds by GC/MSD
EPA Method 624
 ug/L(ppb)

Sample Name: Lab Code:	(a)	(b)							Reporting Limit
	MW-1 0277-1	MW-2 0277-2	MW-3 0277-3	MW-4 0277-4	MW-5 0277-5	MW-6 0277-6	MW-7 0277-7	MW-8 0277-8	
Dichlorodifluoromethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Chloromethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Vinyl Chloride	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Bromomethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Chloroethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Trichlorofluoromethane	<300	ND	ND	ND	ND	10.1	13.4	16.4	1.0
1,1-Dichloroethene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Acetone	<6000	ND	ND	ND	ND	27.2	ND	36.2	20
Carbon Disulfide	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Methylene Chloride	ND (c)	ND	230	ND	ND	ND	ND	4.01	1.0
trans-1,2-Dichloroethene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,1-Dichloroethane	<300	1.88	ND	ND	1.45	2.93	ND	ND	1.0
2,2-Dichloropropane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
cis-1,2-Dichloroethene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
2-Butanone(MEK)	<3000	ND	ND	ND	ND	ND	ND	18.0	10
Bromochloromethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Chloroform	<300	ND	ND	ND	ND	1.01	ND	ND	1.0
1,1,1-Trichloroethane	<300	9.39	ND	ND	ND	2.28	ND	ND	1.0
Carbon Tetrachloride	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,1-Dichloropropene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Benzene	7500	1.69	ND	ND	2.49	ND	ND	8.39	1.0
1,2-Dichloroethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Trichloroethene	<300	ND	22	ND	ND	ND	ND	ND	1.0
1,2-Dichloropropane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Dibromomethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Bromodichloromethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
cis-1,3-Dichloropropene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
4-Methyl-2-Pentanone(MIBK)	<3000	ND	ND	ND	ND	ND	ND	ND	10
Toluene	54,000	1.00	48	ND	7.82	5.37	16.9	10.2	1.0
trans-1,3-Dichloropropene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,1,2-Trichloroethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Tetrachloroethene	<300	1.71	23	ND	ND	5.23	ND	ND	1.0
2-Hexanone	<3000	ND	ND	ND	ND	ND	ND	ND	10
1,3-Dichloropropane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Dibromochloromethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,2-Dibromoethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Chlorobenzene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,1,1,2-Tetrachloroethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Ethylbenzene	5600	10.2	14	ND	2.49	2.71	4.77	6.35	1.0
m,p-Xylene	21,000	13.4	56	ND	13.6	10.4	18.3	24.5	1.0
o-Xylene	9200	ND	33	ND	7.52	4.84	8.34	12.5	1.0
Styrene	<300	ND	ND	ND	ND	ND	ND	ND	1.0

ND Not Detected

a) Results are from a 1:300 dilution.

b) Results are from a 1:5 dilution.

c) Analyzed on 5/8/96 with a ND result.

Project: Seekins Ford - Lincoln Mercury
 Project No.: 6-024-01173-0T04
 Project Manager: James Spontak
 Sample Matrix: Water

Service Request No.: AK960277
 Report Date: 5/9/96
 Report No.: 96027703
 C.O.C. No.: 02920

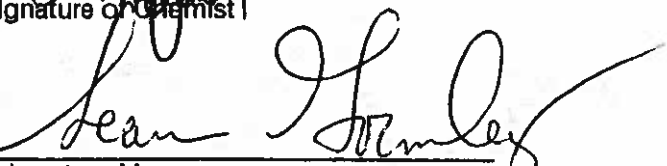
Volatile Organic Compounds by GC/MSD
EPA Method 624
 ug/L(ppb)

Sample Name:	MW-1	MW-2	MW-3	MW-4	MW-5	MW-6	MW-7	MW-8	Reporting Limit
Lab Code:	0277-1	0277-2	0277-3	0277-4	0277-5	0277-6	0277-7	0277-8	
Bromoform	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Isopropylbenzene	<300	3.53	ND	ND	1.85	3.11	2.50	3.84	1.0
Bromobenzene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,1,2,2-Tetrachloroethane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,2,3-Trichloropropane	<300	ND	ND	ND	ND	ND	ND	1.14	1.0
n-Propylbenzene	420	9.59	6.4	ND	4.43	8.96	5.78	9.67	1.0
2-Chlorotoluene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
4-Chlorotoluene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,3,5-Trimethylbenzene	740	16.2	18	ND	15.0	11.7	7.45	13.0	1.0
tert-Butylbenzene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,2,4-Trimethylbenzene	3000	22.2	49	ND	36.2	37.7	25.8	41.8	1.0
sec-Butylbenzene	<300	ND	ND	ND	2.68	4.73	1.33	2.64	1.0
1,3-Dichlorobenzene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
4-Isopropyltoluene	<300	1.30	7.6	ND	3.50	8.89	2.47	4.06	1.0
1,4-Dichlorobenzene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,2-Dichlorobenzene	<300	ND	86	ND	2.71	ND	ND	ND	1.0
n-Butylbenzene	<300	ND	ND	ND	ND	11.0	2.42	5.06	1.0
1,2-Dibromo-3-Chloropropane	<300	ND	ND	ND	ND	ND	ND	ND	1.0
1,2,4-Trichlorobenzene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Hexachlorobutadiene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Naphthalene	<300	ND	ND	ND	18.0	35.2	12.0	20.4	5.0
1,2,3-Trichlorobenzene	<300	ND	ND	ND	ND	ND	ND	ND	1.0
Sample Date:	5/1/96	5/1/96	5/1/96	5/1/96	5/1/96	5/1/96	5/1/96	5/1/96	
Analysis Date:	5/7/96	5/7/96	5/7/96	5/7/96	5/7/96	5/7/96	5/7/96	5/7/96	
Surrogate Recoveries:									EPA %Recovery Acceptance
Dibromofluoromethane:	101%	98.1%	101%	98.7%	102%	104%	102%	99.1%	86%-118%
Toluene-d8:	100%	98.5%	97.9%	98.6%	97.8%	98.8%	98.8%	99.0%	88%-110%
4-Bromofluorobenzene:	108%	108%	107%	110%	106%	104%	104%	105%	86%-115%

ND Not Detected



 Signature of Chemist



 Laboratory Manager

Project: Seekins Ford - Lincoln Mercury
 Project No.: 6-024-01173-0T04
 Project Manager: James Spontak
 Sample Matrix: Water

Service Request No.: AK960277
 Report Date: 5/9/96
 Report No.: 96027704
 C.O.C. No.: 02920

Volatile Organic Compounds by GC/MSD
EPA Method 624
 ug/L(ppb)

Sample Name: Lab Code:	MW-9 0277-9	GWP-1 0277-10	GWP-2 0277-11	GWP-3 0277-12	(a) DUP-1 0277-13	Trip Blank 0277-14	Lab Blank 0277-MB	Reporting Limit
Dichlorodifluoromethane	ND	ND	ND	ND	<300	ND	ND	1.0
Chloromethane	ND	ND	ND	ND	<300	ND	ND	1.0
Vinyl Chloride	ND	ND	ND	ND	<300	ND	ND	1.0
Bromomethane	ND	ND	ND	ND	<300	ND	ND	1.0
Chloroethane	ND	ND	ND	ND	<300	ND	ND	1.0
Trichlorofluoromethane	ND	ND	33.4	2.22	<300	ND	ND	1.0
1,1-Dichloroethene	ND	ND	ND	ND	<300	ND	ND	1.0
Acetone	ND	ND	ND	ND	<6000	ND	ND	20
Carbon Disulfide	ND	ND	ND	ND	<300	ND	ND	1.0
Methylene Chloride	3.56	3.63	3.87	4.91	ND (c)	ND	ND	1.0
trans-1,2-Dichloroethene	ND	ND	ND	ND	<300	ND	ND	1.0
1,1-Dichloroethane	ND	ND	ND	ND	<300	ND	ND	1.0
2,2-Dichloropropane	ND	ND	ND	ND	<300	ND	ND	1.0
cis-1,2-Dichloroethene	ND	ND	ND	ND	<300	ND	ND	1.0
2-Butanone(MEK)	ND	ND	ND	ND	<3000	ND	ND	10
Bromochloromethane	ND	ND	ND	ND	<300	ND	ND	1.0
Chloroform	ND	ND	ND	ND	<300	ND	ND	1.0
1,1,1-Trichloroethane	ND	ND	ND	ND	<300	ND	ND	1.0
Carbon Tetrachloride	ND	ND	ND	ND	<300	ND	ND	1.0
1,1-Dichloropropene	ND	ND	ND	ND	<300	ND	ND	1.0
Benzene	ND	117	ND	ND	10,000	ND	ND	1.0
1,2-Dichloroethane	ND	ND	ND	ND	<300	ND	ND	1.0
Trichloroethene	ND	ND	ND	ND	<300	ND	ND	1.0
1,2-Dichloropropane	ND	ND	ND	ND	<300	ND	ND	1.0
Dibromomethane	ND	ND	ND	ND	<300	ND	ND	1.0
Bromodichloromethane	ND	ND	ND	ND	<300	ND	ND	1.0
cis-1,3-Dichloropropene	ND	ND	ND	ND	<300	ND	ND	1.0
4-Methyl-2-Pentanone(MIBK)	ND	ND	ND	ND	<3000	ND	ND	10
Toluene	2.53	ND	ND	ND	72,000	ND	ND	1.0
trans-1,3-Dichloropropene	ND	ND	ND	ND	<300	ND	ND	1.0
1,1,2-Trichloroethane	ND	ND	ND	ND	<300	ND	ND	1.0
Tetrachloroethene	ND	ND	ND	ND	<300	ND	ND	1.0
2-Hexanone	ND	ND	ND	ND	<3000	ND	ND	10
1,3-Dichloropropane	ND	ND	ND	ND	<300	ND	ND	1.0
Dibromochloromethane	ND	ND	ND	ND	<300	ND	ND	1.0
1,2-Dibromoethane	ND	ND	ND	ND	<300	ND	ND	1.0
Chlorobenzene	ND	ND	ND	ND	<300	ND	ND	1.0
1,1,1,2-Tetrachloroethane	ND	ND	ND	ND	<300	ND	ND	1.0
Ethylbenzene	1.00	1.36	ND	ND	6600	ND	ND	1.0
m,p-Xylene	3.62	15.9	ND	ND	25,000	ND	ND	1.0
o-Xylene	1.66	ND	ND	ND	11,000	ND	ND	1.0
Styrene	ND	ND	ND	ND	<300	ND	ND	1.0

ND Not Detected

(a) Results are from a 1:300 dilution.

(b) Analyzed on 5/8/96 with a ND result.

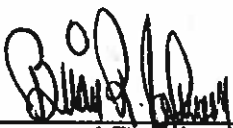
Project: Seekins Ford - Lincoln Mercury
 Project No.: 6-024-01173-0T04
 Project Manager: James Spontak
 Sample Matrix: Water

Service Request No.: AK960277
 Report Date: 5/9/96
 Report No.: 96027704
 C.O.C. No.: 02920

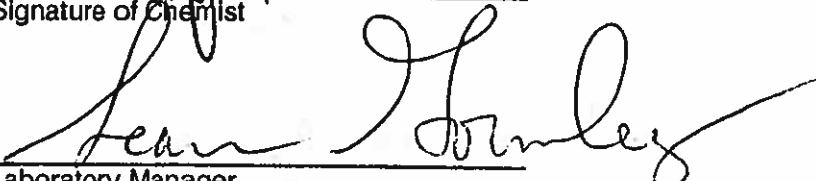
Volatile Organic Compounds by GC/MSD
EPA Method 824
 ug/L(ppb)

Sample Name: Lab Code:	MW-9 0277-9	GWP-1 0277-10	GWP-2 0277-11	GWP-3 0277-12	(a) DUP-1 0277-13	Trip Blank 0277-14	Lab Blank 0277-MB	Reporting Limit
Bromoform	ND	ND	ND	ND	<300	ND	ND	1.0
Isopropylbenzene	1.08	3.70	ND	ND	<300	ND	ND	1.0
Bromobenzene	ND	ND	ND	ND	<300	ND	ND	1.0
1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	<300	ND	ND	1.0
1,2,3-Trichloropropane	ND	ND	ND	ND	<300	ND	ND	1.0
n-Propylbenzene	2.95	1.53	ND	ND	410	ND	ND	1.0
2-Chlorotoluene	ND	ND	ND	ND	<300	ND	ND	1.0
4-Chlorotoluene	ND	ND	ND	ND	<300	ND	ND	1.0
1,3,5-Trimethylbenzene	3.43	ND	ND	ND	780	ND	ND	1.0
tert-Butylbenzene	ND	ND	ND	ND	<300	ND	ND	1.0
1,2,4-Trimethylbenzene	11.32	ND	ND	ND	3100	ND	ND	1.0
sec-Butylbenzene	1.67	ND	ND	ND	<300	ND	ND	1.0
1,3-Dichlorobenzene	ND	ND	ND	ND	<300	ND	ND	1.0
4-Isopropyltoluene	2.11	ND	ND	ND	<300	ND	ND	1.0
1,4-Dichlorobenzene	ND	ND	ND	ND	<300	ND	ND	1.0
1,2-Dichlorobenzene	ND	ND	ND	ND	<300	ND	ND	1.0
n-Butylbenzene	3.72	ND	ND	ND	<300	ND	ND	1.0
1,2-Dibromo-3-Chloropropane	ND	ND	ND	ND	<300	ND	ND	1.0
1,2,4-Trichlorobenzene	ND	ND	ND	ND	<300	ND	ND	1.0
Hexachlorobutadiene	ND	ND	ND	ND	<300	ND	ND	1.0
Naphthalene	10.9	ND	ND	ND	<300	ND	ND	5.0
1,2,3-Trichlorobenzene	ND	ND	ND	ND	<300	ND	ND	1.0
Sample Date:	5/1/96	5/1/96	5/1/96	5/1/96	5/1/96	5/1/96	5/7/96	
Analysis Date:	5/7/96	5/7/96	5/7/96	5/7/96	5/7/96	5/7/96	5/7/96	
Surrogate Recoveries:								EPA %Recovery Acceptance
Dibromofluoromethane:	100%	101%	102%	100%	102%	98.7%	102%	86%-118%
Toluene-d8:	98.3%	99.3%	99.4%	99.2%	99.5%	98.4%	98.0%	88%-110%
4-Bromofluorobenzene:	105%	106%	105%	108%	107%	108%	108%	86%-115%

ND Not Detected



 Signature of Chemist



 Laboratory Manager

Project: Seekins Ford - Lincoln Mercury
 Project No.: 6-024-01173-0T04
 Project Manager: James Spontak
 Sample Matrix: Water

Service Request No.: AK960277
 Report Date: 5/9/96
 Report No.: 96027705
 C.O.C.: 02920

**QC Data Report
 MS/MSD Summary
 Volatile Organic Compounds by GC/MS
 EPA Method 624
 ug/L(ppb)**

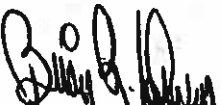
Sample Name:	MW4	Spike Level	Matrix Spike	Percent Recovery	Matrix Spike Duplicate	Percent Recovery	EPA % Recovery Acceptance Criteria (a)	Relative Percent Difference (RPD)
Lab Code:	0277-4	(ug/L)		(MS)		(DMS)		
1,1 - Dichloroethene	ND	50	55.6	111	54.7	109	75% - 113%	1.8
Benzene	ND	50	54.2	108	54.3	109	77% - 117%	<1.0
Trichloroethene	ND	50	51.5	103	50.6	101	70% - 110%	2.0
Toluene	ND	50	54.9	110	55.0	110	77% - 126%	<1.0
Chlorobenzene	ND	50	52.6	105	54.8	110	80% - 116%	4.6

Sample Date: 5/1/96 ~ 5/1/96 ~ 5/1/96 ~ ~
 Analysis Date: 5/7/96 ~ 5/7/96 ~ 5/7/96 ~ ~

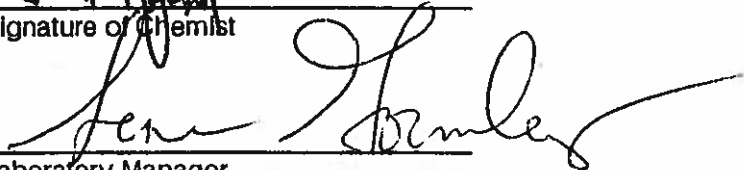
Surrogate Recovery:						Control Limits
Dibromofluoromethane:	98.7%	~	102%	~	101%	~ 86%-118%
Toluene-d ₈ :	98.6%	~	99.0%	~	100%	~ 88%-110%
4-Bromofluorobenzene:	110%	~	106%	~	107%	~ 86%-115%

ND Not Detected

(a) Criteria from EPA Method 8260A, Table 6.



 Signature of Chemist



 Laboratory Manager

Project: Seekins Ford - Lincoln Mercury
Project No.: 6-024-01173-0T04
Project Manager: James Spontak
Sample Matrix: Water

Service Request No.: AK960277
Report Date: 5/9/96
Report No.: 96027701
C.O.C. No.: 02920

Gasoline Range Organics
ADEC Method AK 101
ug/L(ppb)

Sample Name	Lab Code	Sample Date	Analysis Date	Method Reporting Limit	Result	Surrogate Recovery a,a,a - TFT
MW-1	0277-1	5/1/96	5/7/96	10,000(a)	240,000(b)	93.1
MW-2	0277-2	5/1/96	5/6/96	50.0	259	103
MW-3	0277-3	5/1/96	5/7/96	50.0	996(c)	93.2
MW-4	0277-4	5/1/96	5/7/96	50.0	ND	93.0
MW-5	0277-5	5/1/96	5/7/96	50.0	169(c)	95.3
MW-6	0277-6	5/1/96	5/7/96	50.0	118(c)	92.9
MW-7	0277-7	5/1/96	5/7/96	50.0	258(c)	94.6
MW-8	0277-8	5/1/96	5/7/96	50.0	354(c)	94.0
MW-9	0277-9	5/1/96	5/7/96	50.0	55.8(c)	93.9
GWP-1	0277-10	5/1/96	5/7/96	50.0	340(d)	93.5
GWP-2	0277-11	5/1/96	5/7/96	50.0	ND	92.0
GWP-3	0277-12	5/1/96	5/7/96	50.0	ND	91.8
DUP-1	0277-13	5/1/96	5/7/96	10,000(a)	220,000(b)	92.6
Trip Blank	0277-14	5/1/96	5/7/96	50.0	ND	91.5
Lab Blank	0277-MB	5/6/96	5/6/96	50.0	ND	92.1
Lab Blank	0277-MB	5/7/96	5/7/96	50.0	ND	92.4

ND Not Detected

Acceptance Criteria 50%-150%

- (a) Elevated method reporting limit due to a 1:200 dilution.
- (b) Result is from a 1:200 dilution.
- (c) Chromatographic peaks were quantified as gasoline; however, the chromatographic peak pattern suggests a heavier organic material such as diesel or mineral spirits.
- (d) Chromatographic peaks were quantified as gasoline; however, the chromatographic peak pattern indicates that benzene produced the majority of the quantifiable area for gasoline.

Signature of Chemist

Laboratory Manager

Project: Seekins Ford - Lincoln Mercury
 Project No.: 6-024-01173-0T04
 Project Manager: James Spontak
 Sample Matrix: Water

Service Request No.: AK960277
 Report Date: 5/9/96
 Report No.: 96027702
 C.O.C. No.: 02920

QC Data Report
Matrix Spike Recoveries
Gasoline Range Organics
ADEC Method AK101
ug/L(ppb)

Sample Name:	MW-2	Spike Level (ug/L)	Matrix Spike (MS)	Percent Recovery (MS)	Matrix Spike Duplicate (MSD)	Percent Recovery (MSD)	Relative Percent Difference
Gasoline:	281	638	859	90.6	891	95.6	3.66
Acceptance Limits:	~	~	~	75%-125%	~	75%-125%	<25
Sample Date:	5/1/96	~	5/1/96	~	5/1/96	~	~
Analysis Date:	5/6/96	~	5/6/96	~	5/6/96	~	~
Surrogate Recovery:							Control Limits
a,a,a-Trifluorotoluene:	103%	~	111%	~	110%	~	38%-150%

ND Not Detected

Spike Source: ERA Gasoline Standard, Lot # 50008

Signature of Chemist

Sean Hanley

Laboratory Manager

Project: Seekins Ford - Lincoln Mercury
Project No.: 6-024-01173-0T04
Project Manager: James Spontak
Sample Matrix: Water

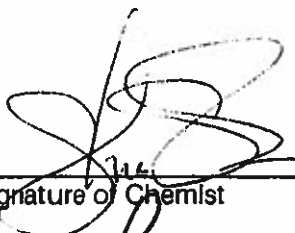
Service Request No.: AK960277
Report Date: 5/10/96
Report No.: 96027706
C.O.C. No.: 02920

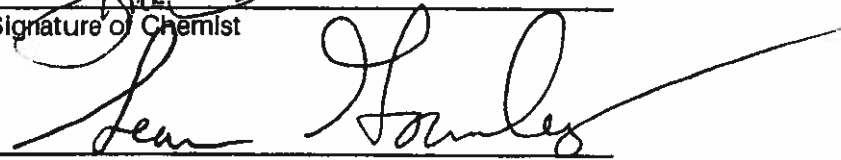
Diesel Range Organics
ADEC Method AK 102
mg/L(ppm)

Sample Name	Lab Code	Sample Date	Extraction Date	Analysis Date	Diesel Result	Surrogate Recovery 2-Fluorobiphenyl
MW-1	0277-1	5/1/96	5/7/96	5/8/96	6.2(a)(b)	(c)
MW-2	0277-2	5/1/96	5/7/96	5/8/96	0.74	(c)
MW-3	0277-3	5/1/96	5/7/96	5/8/96	5.4(a)	(c)
MW-4	0277-4	5/1/96	5/7/96	5/8/96	0.24	(c)
MW-5	0277-5	5/1/96	5/7/96	5/8/96	1.1(b)	(c)
MW-6	0277-6	5/1/96	5/7/96	5/9/96	0.94(b)	(c)
MW-7	0277-7	5/1/96	5/7/96	5/9/96	0.47(b)	(c)
MW-8	0277-8	5/1/96	5/7/96	5/9/96	0.69(b)	(c)
MW-9	0277-9	5/1/96	5/7/96	5/9/96	0.84(b)	(c)
GWP-1	0277-10	5/1/96	5/7/96	5/9/96	0.48(a)(b)	(c)
GWP-2	0277-11	5/1/96	5/7/96	5/9/96	0.35(b)	(c)
GWP-3	0277-12	5/1/96	5/7/96	5/9/96	0.17(b)	(c)
DUP-1	0277-13	5/1/96	5/7/96	5/9/96	5.6(a)(b)	(c)
Lab Blank	0277-MB	5/7/96	5/7/96	5/8/96	<0.05	120

Acceptance Criteria: 50%-150%

- (a) Result is from a 1:5 dilution.
- (b) Chromatographic peaks were quantitated as diesel; however, the chromatographic peak pattern does not resemble the diesel standard.
- (c) Surrogate recovery cannot be determined due to the presence of target and non-target analytes within the retention time window of 2-Fluorobiphenyl.


Signature of Chemist


Laboratory Manager

Project: Seekins Ford - Lincoln Mercury
 Project No.: 6-024-01173-0T04
 Project Manager: James Spontak
 Sample Matrix: Water

Service Request No.: AK960277
 Report Date: 5/10/96
 Report No.: 96027707
 C.O.C. No.: 02920


**Quality Assurance Data Report
 Laboratory Control Sample Summary
 Diesel Range Organics
 ADEC Method AK 102
 mg/L(ppm)**

Standard Source	Lab Code	True Value	LCS Result	LCS Recovery	LCS Dup Result	LCS Dup Recovery	% Recovery Acceptance Limits	RPD LCS/LCS Dup
Restek *	0277-MB	0.50	0.48	96	0.50	100	75-125	4.1

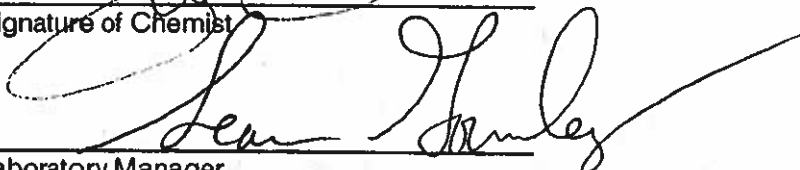
Date Extracted: ~ ~ 5/7/96 ~ 5/7/96 ~
Date Analyzed: ~ ~ 5/8/96 ~ 5/8/96 ~

Surrogate Recovery: 2-Fluorobiphenyl: ~ ~ 88% ~ 90% ~ **ADEC Acceptance Criteria**
 50%-150%

* Restek Diesel Fuel (Lot # A005769)



 Signature of Chemist



 Laboratory Manager

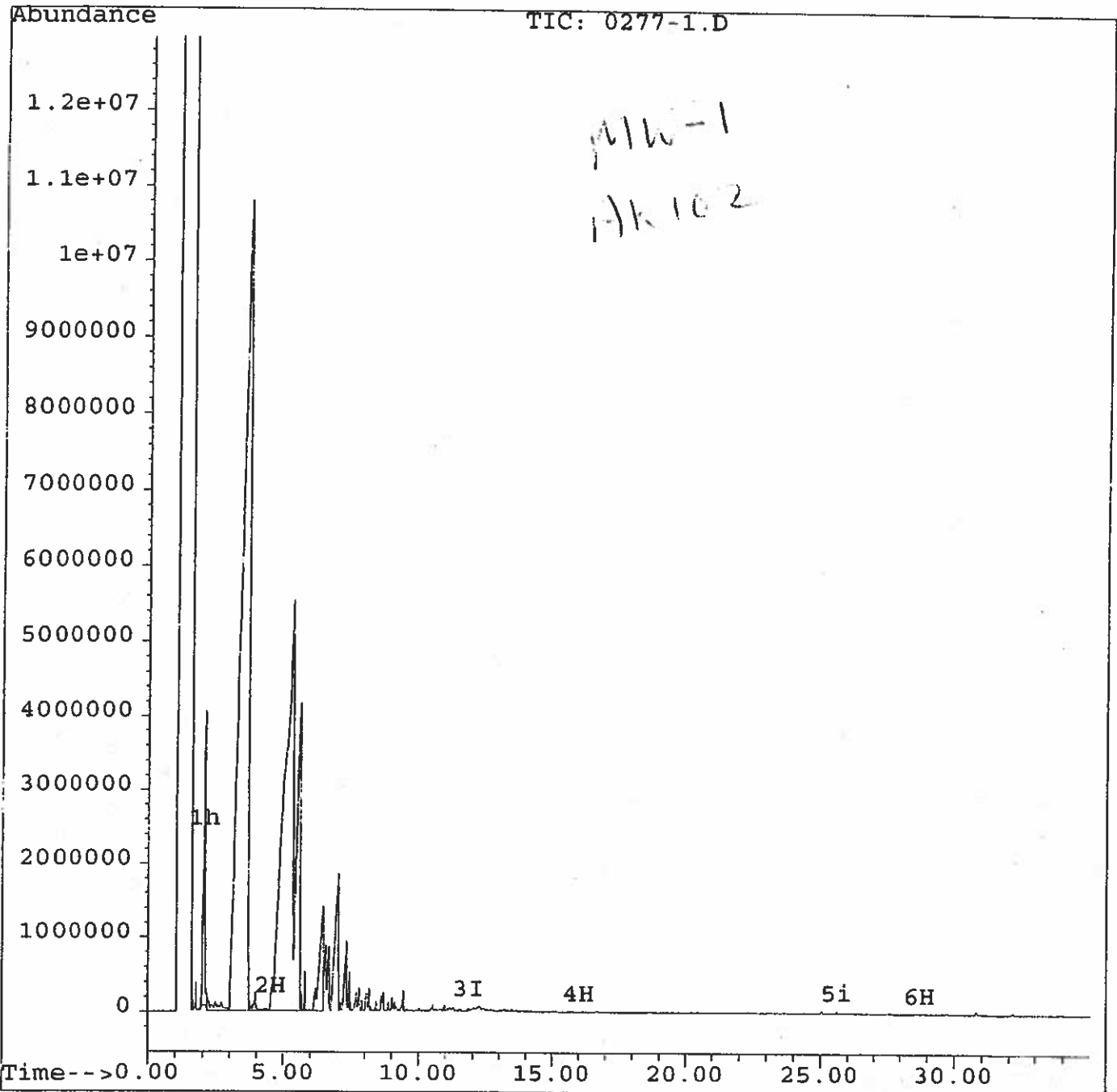
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-1.D
Acq On : 08 May 96 08:11 AM
Sample : 0277-1 1L TO 1ML
Misc :
Quant Time: May 8 9:11 1996

Vial: 8
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Tue Apr 30 13:02:45 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



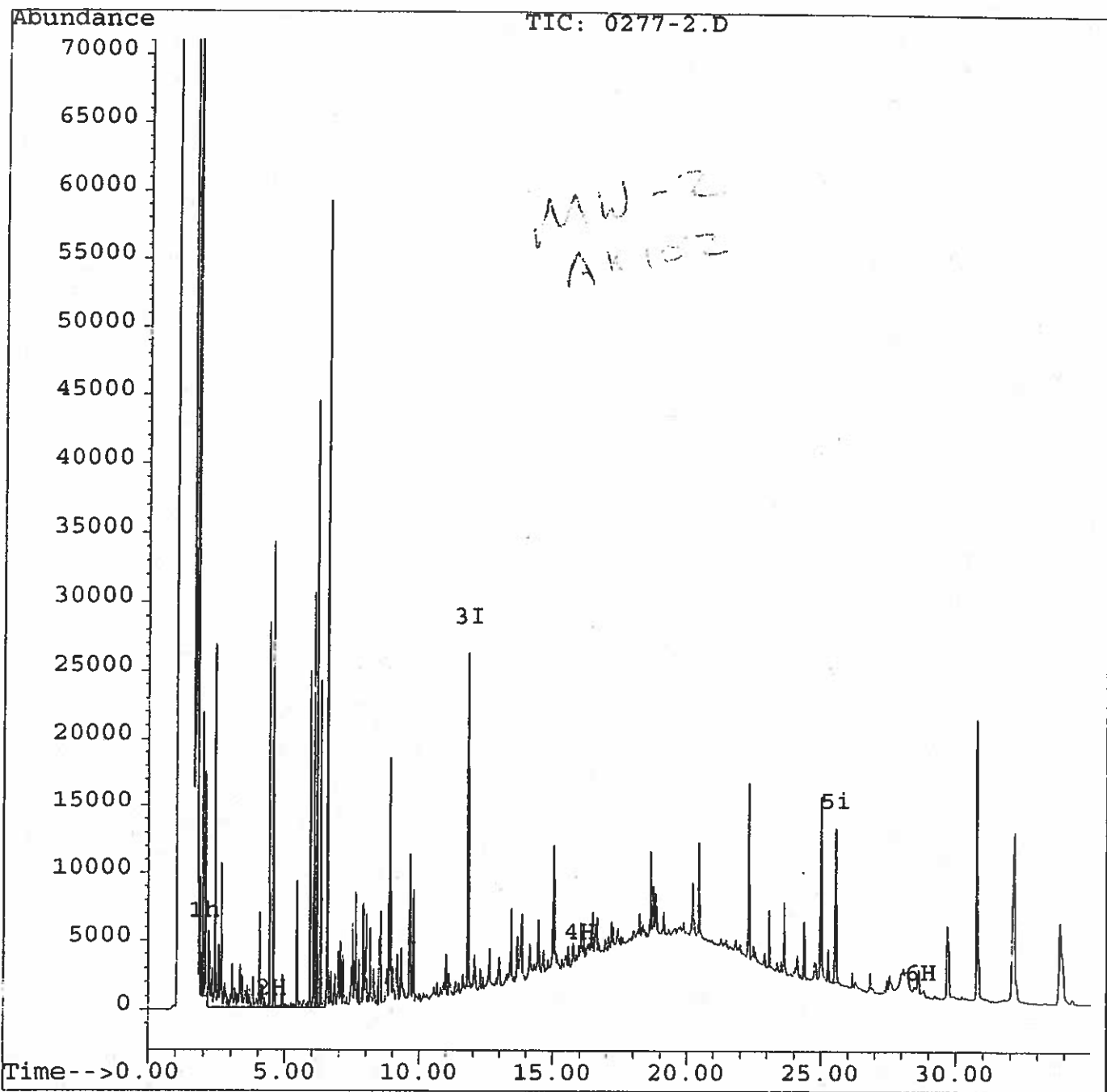
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-2.D
Acq On : 08 May 96 08:56 AM
Sample : 0277-2 1L TO 1ML
Misc :
Quant Time: May 8 9:42 1996

Vial: 9
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Tue Apr 30 13:02:45 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



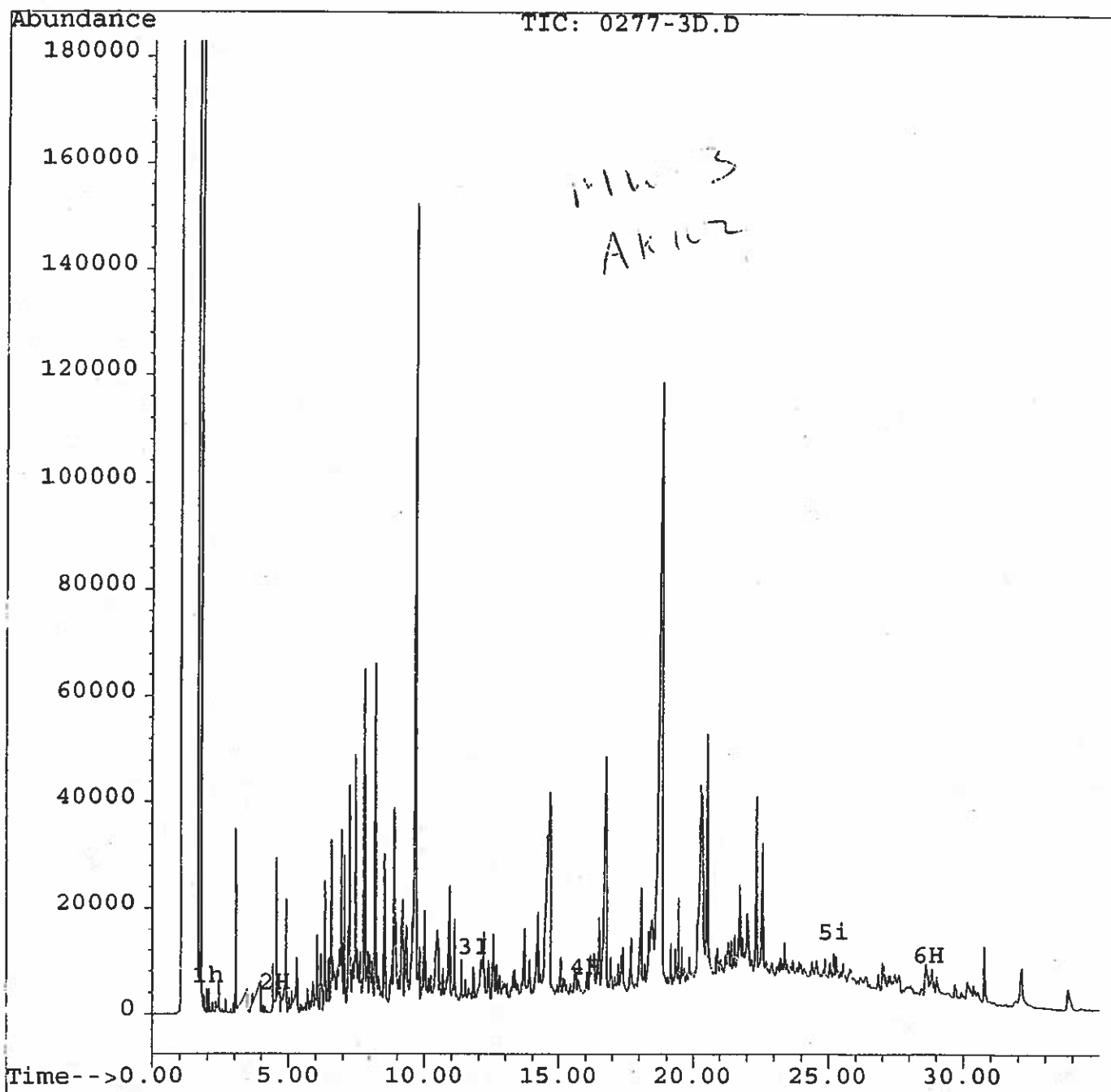
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-3D.D
Acq On : 08 May 96 10:25 AM
Sample : 0277-3 1L TO 1ML X5
Misc :
Quant Time: May 8 11:09 1996

Vial: 11
Operator:
Inst : FID-ECD
Multiplr: 0.01

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Tue Apr 30 13:02:45 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



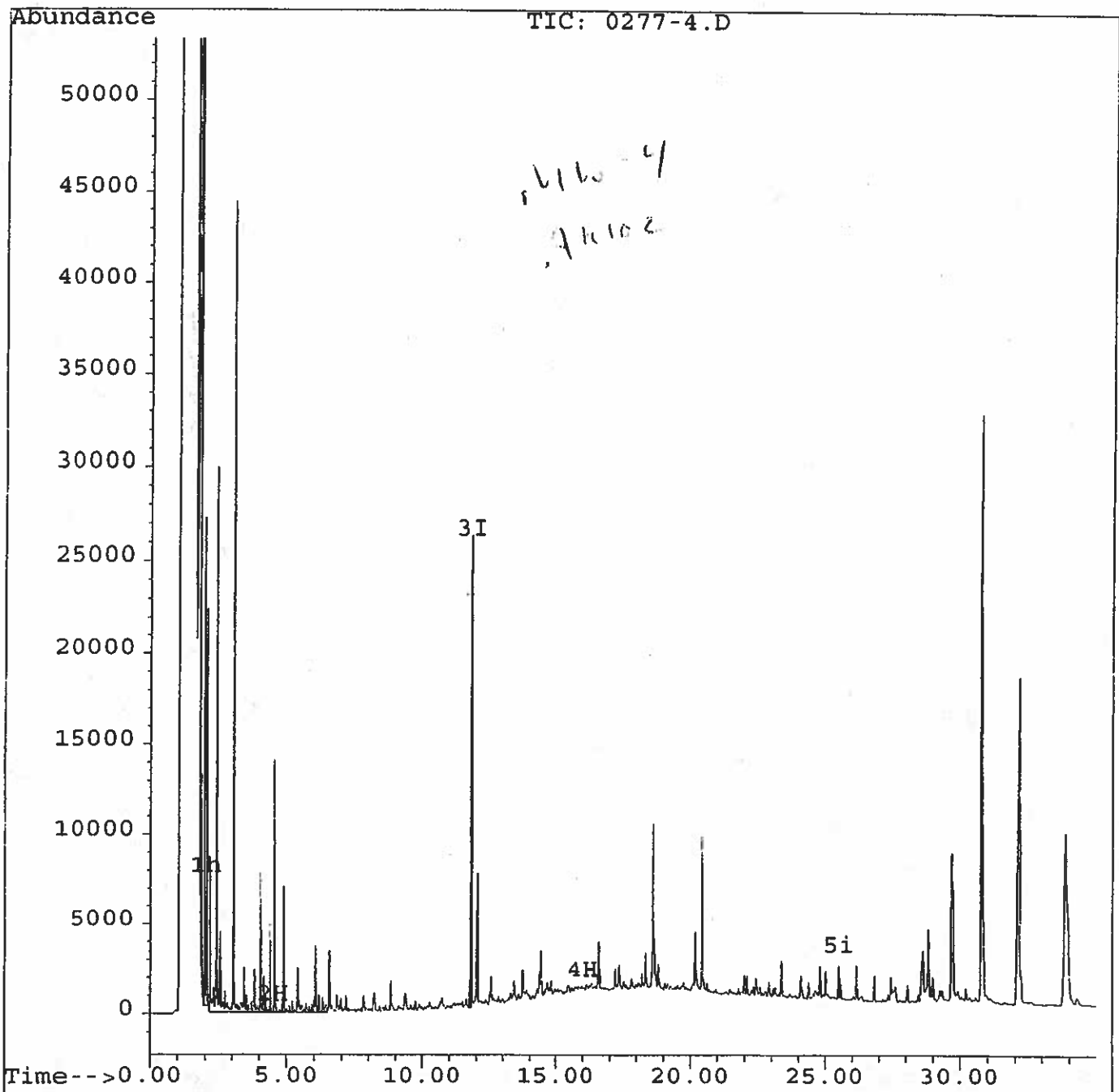
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-4.D
Acq On : 08 May 96 11:54 AM
Sample : 0277-4 1L TO 1ML
Misc :
Quant Time: May 8 13:13 1996

Vial: 12
Operator:
Inst : FID-ECD
Multiplr: 0.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Tue Apr 30 13:02:45 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



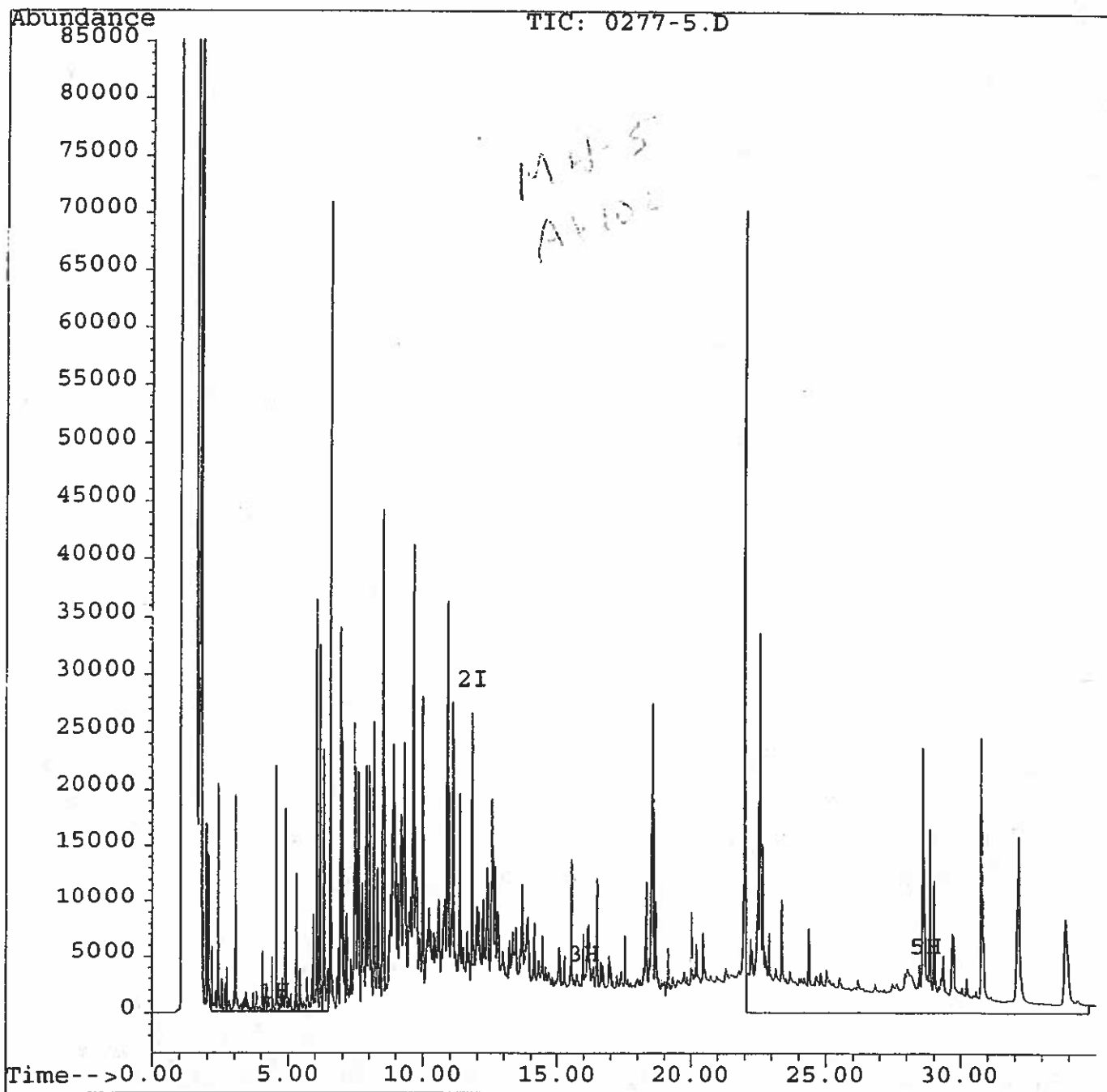
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-5.D
Acq On : 08 May 96 12:40 PM
Sample : 0277-5 1L TO 1ML
Misc :
Quant Time: May 9 12:06 1996

Vial: 13
Operator:
Inst : FID-ECD
Multiplr: 0.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



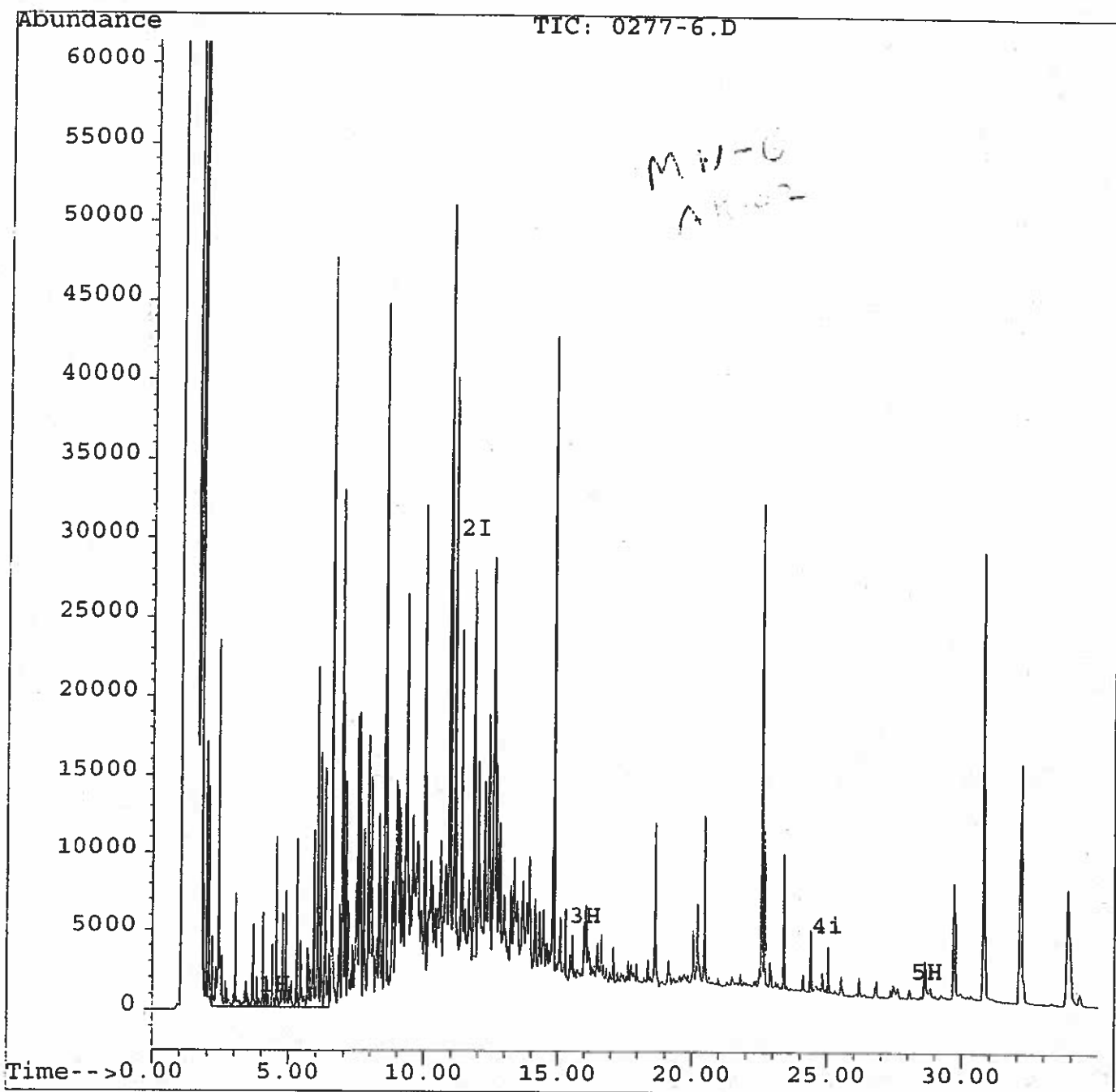
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-6.D
Acq On : 09 May 96 00:03 AM
Sample : 0277-6 1L TO 1ML
Misc :
Quant Time: May 9 12:20 1996

Vial: 14
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



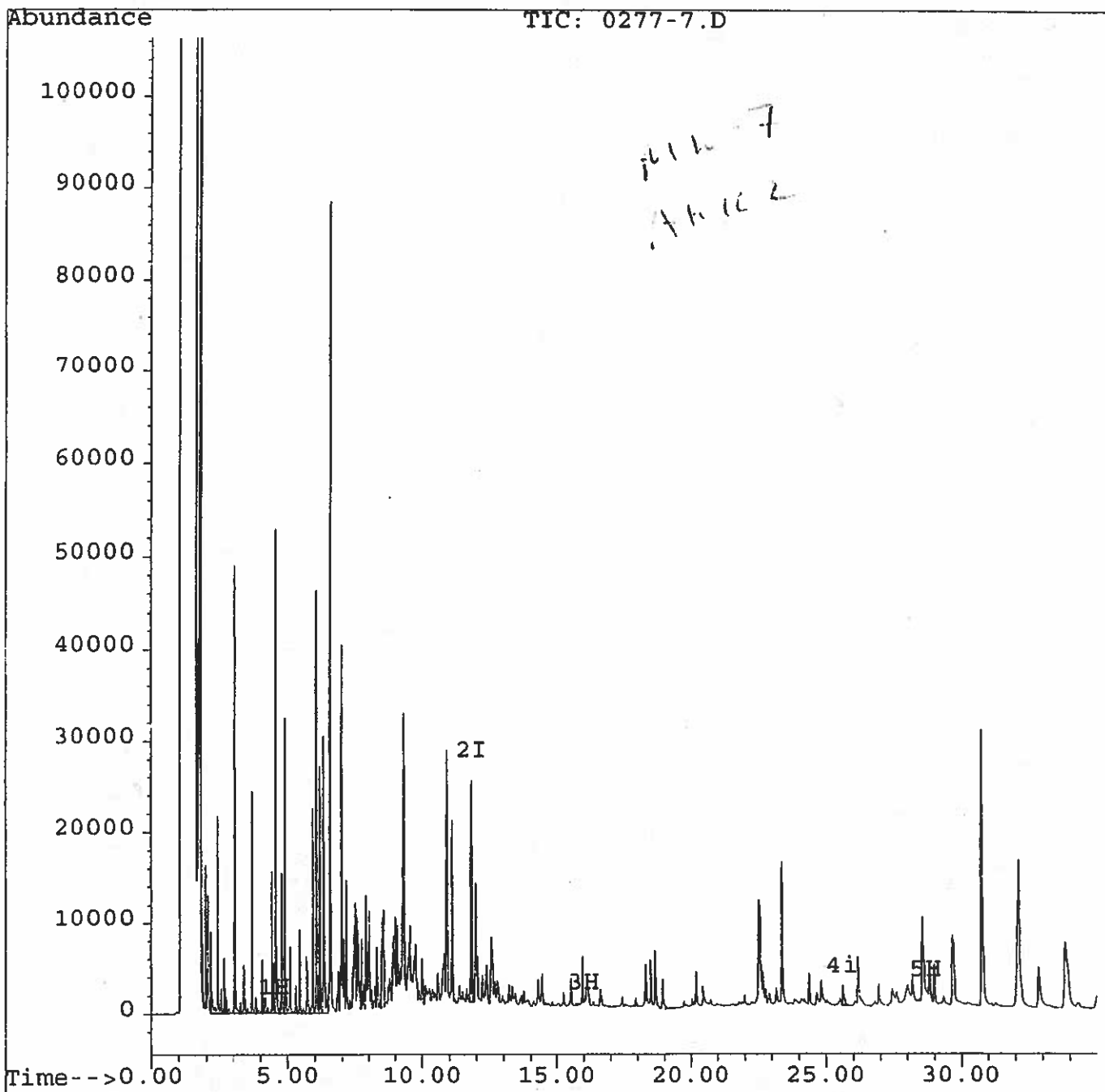
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-7.D
Acq On : 09 May 96 00:48 AM
Sample : 0277-7 1L TO 1ML
Misc :
Quant Time: May 9 12:21 1996

Vial: 15
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



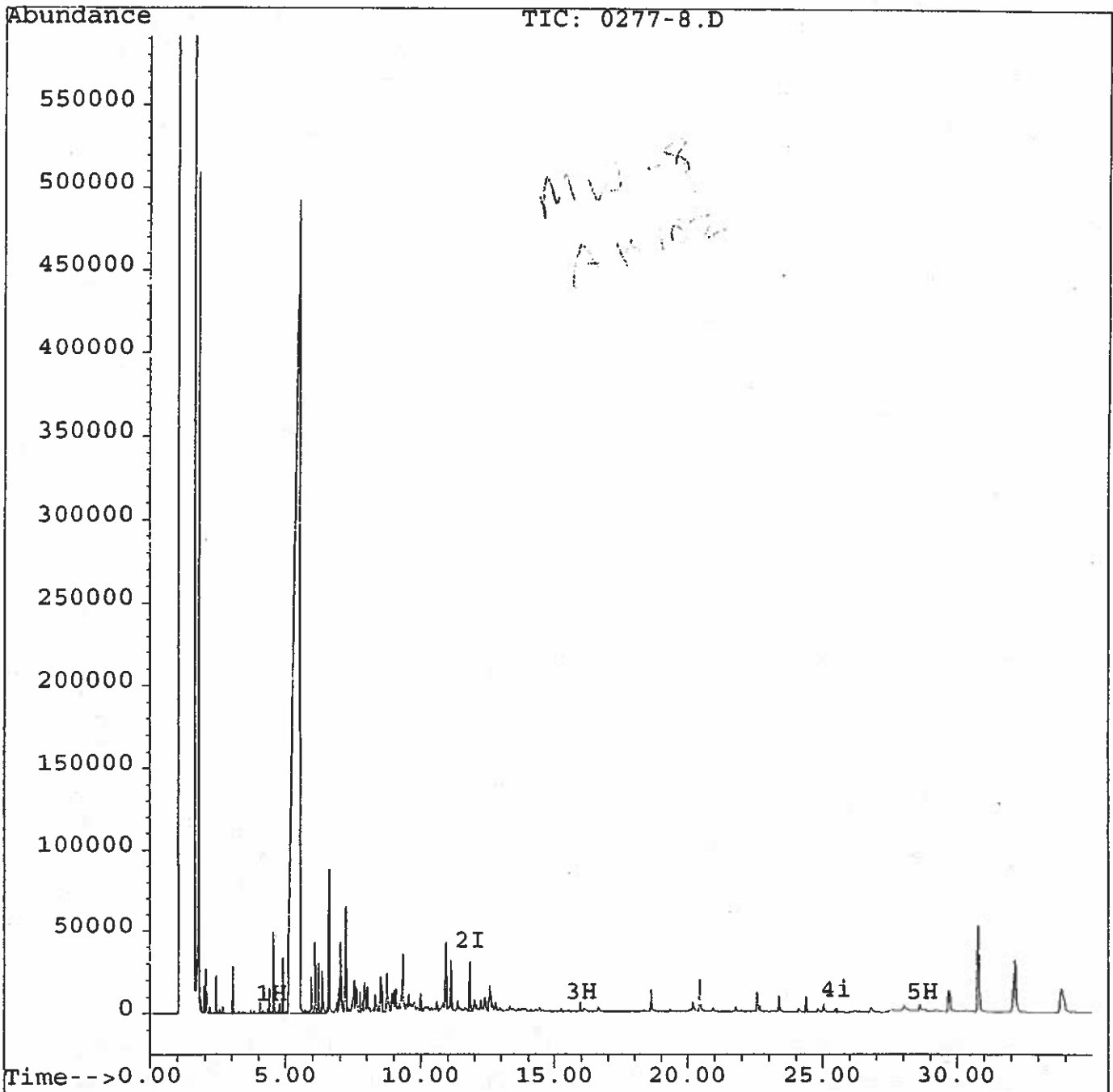
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-8.D
Acq On : 09 May 96 01:33 AM
Sample : 0277-8 1L TO 1ML
Misc :
Quant Time: May 9 12:22 1996

Vial: 16
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



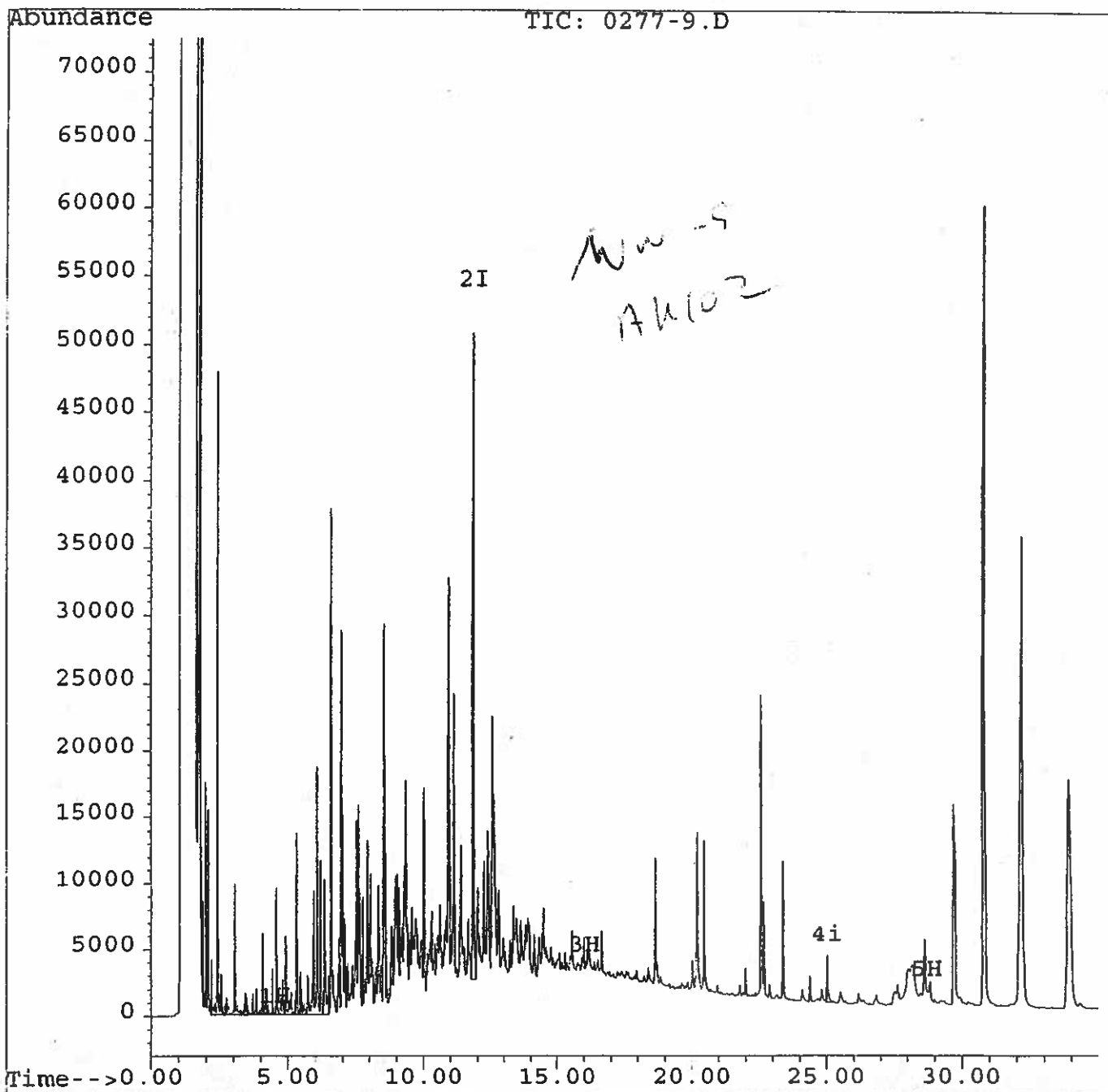
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-9.D
Acq On : 09 May 96 02:19 AM
Sample : 0277-9 1L TO 1ML
Misc :
Quant Time: May 9 12:23 1996

Vial: 17
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



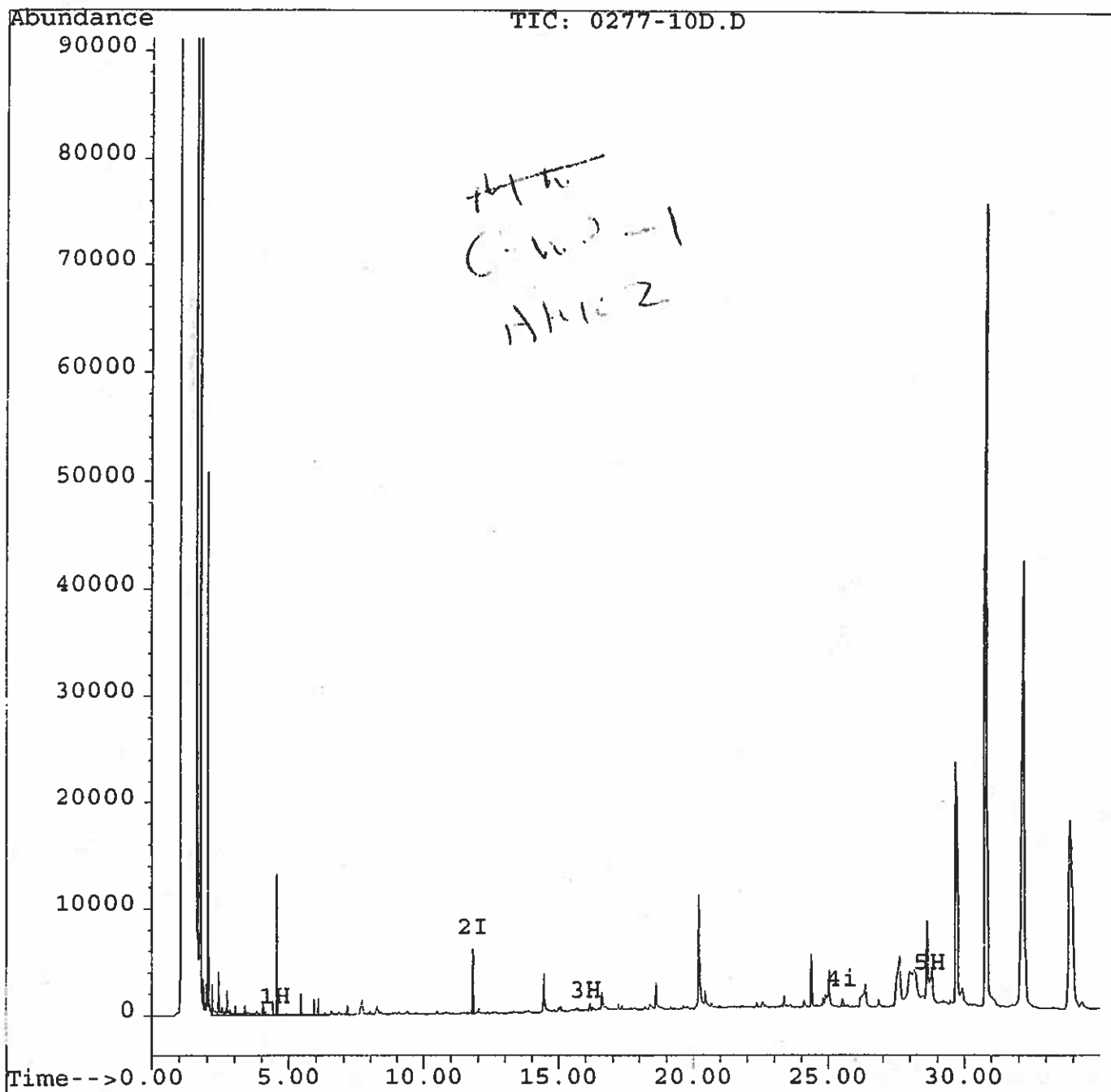
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-10D.D
Acq On : 09 May 96 03:04 AM
Sample : 0277-10 1L TO 1ML X5
Misc :
Quant Time: May 9 12:23 1996

Vial: 18
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



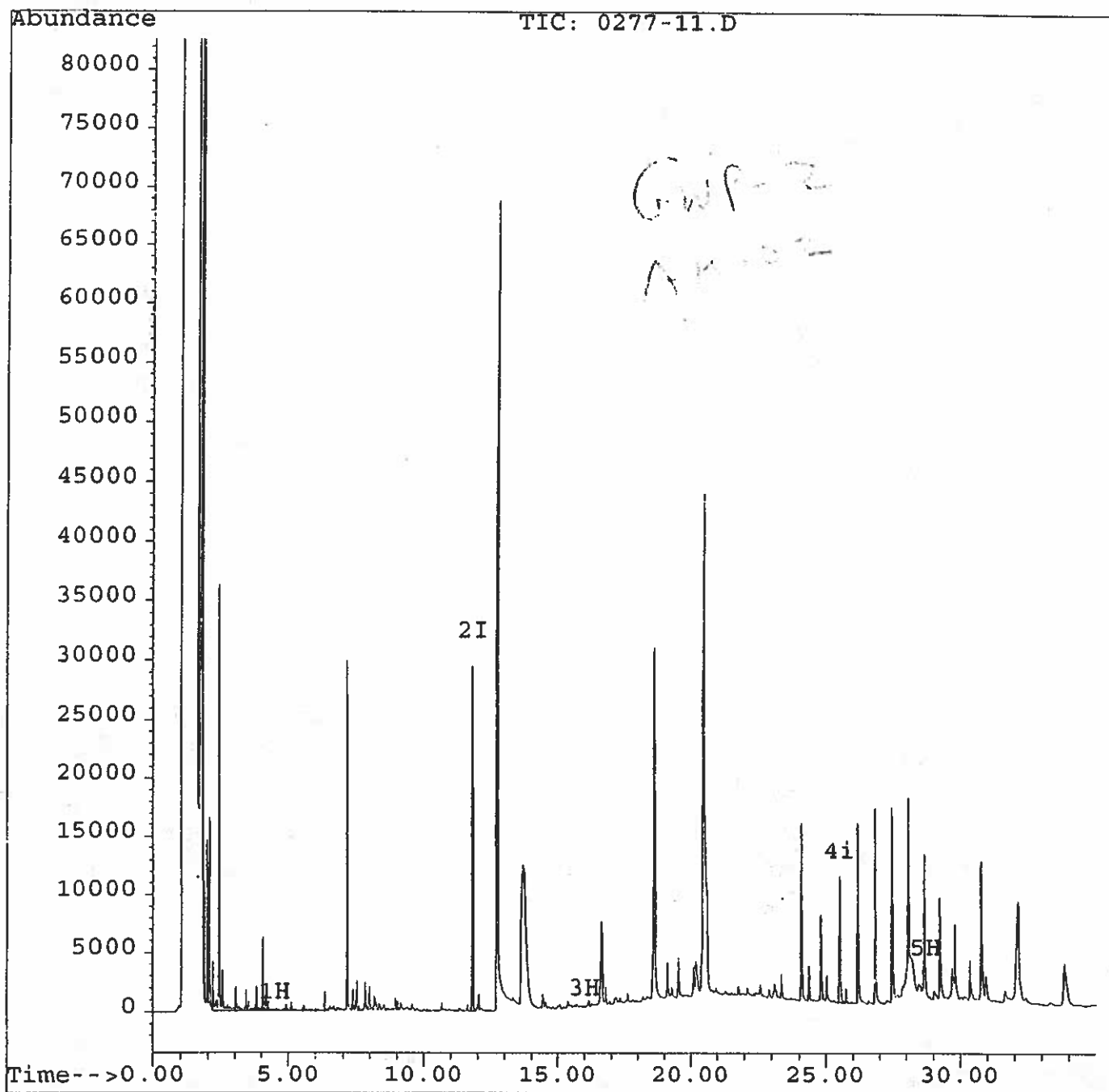
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-11.D
Acq On : 09 May 96 03:49 AM
Sample : 0277-11 1L TO 1ML
Misc :
Quant Time: May 9 12:07 1996

Vial: 19
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



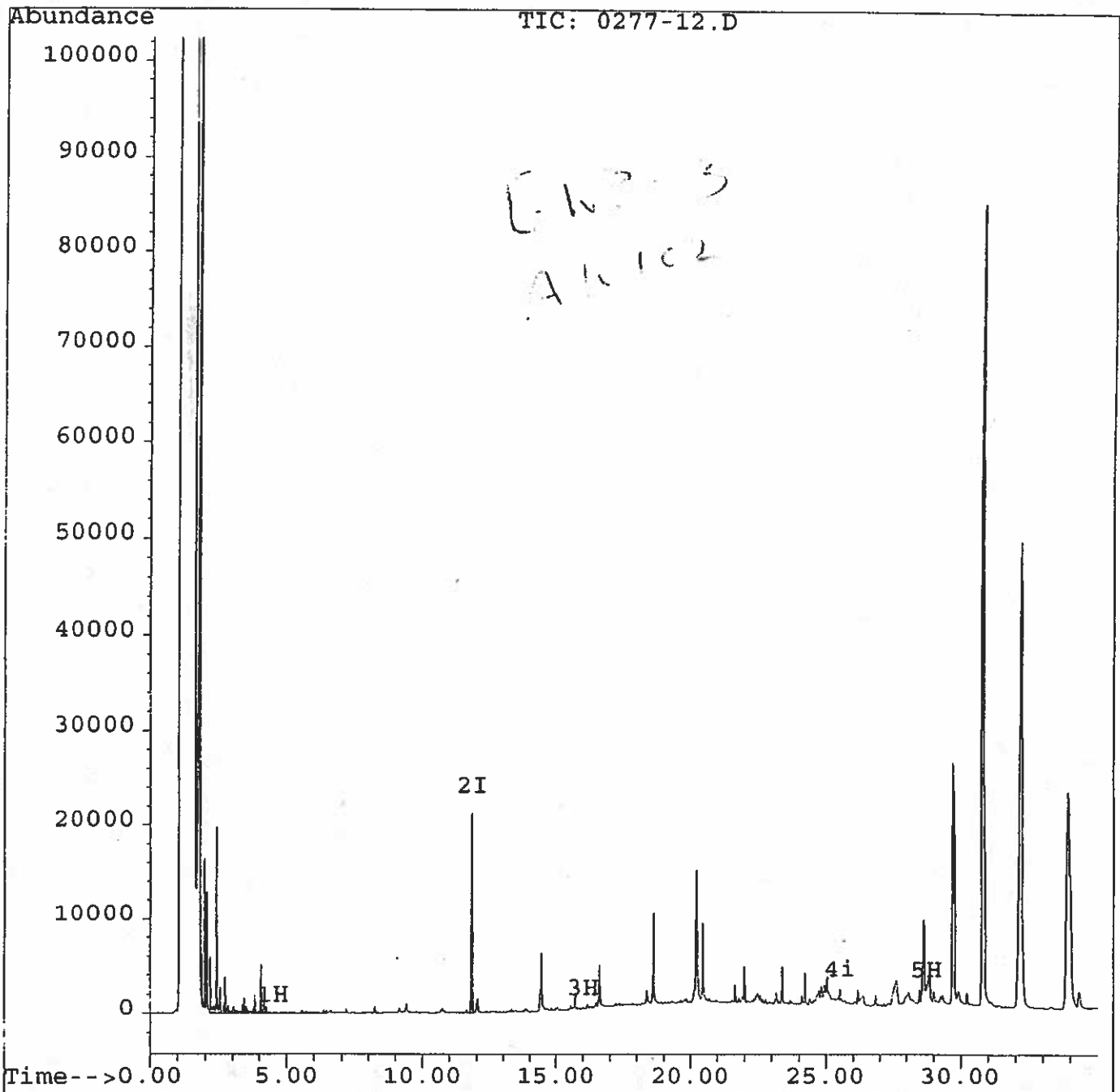
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-12.D
Acq On : 09 May 96 04:34 AM
Sample : 0277-12 1L TO 1ML
Misc :
Quant Time: May 9 12:24 1996

Vial: 20
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :



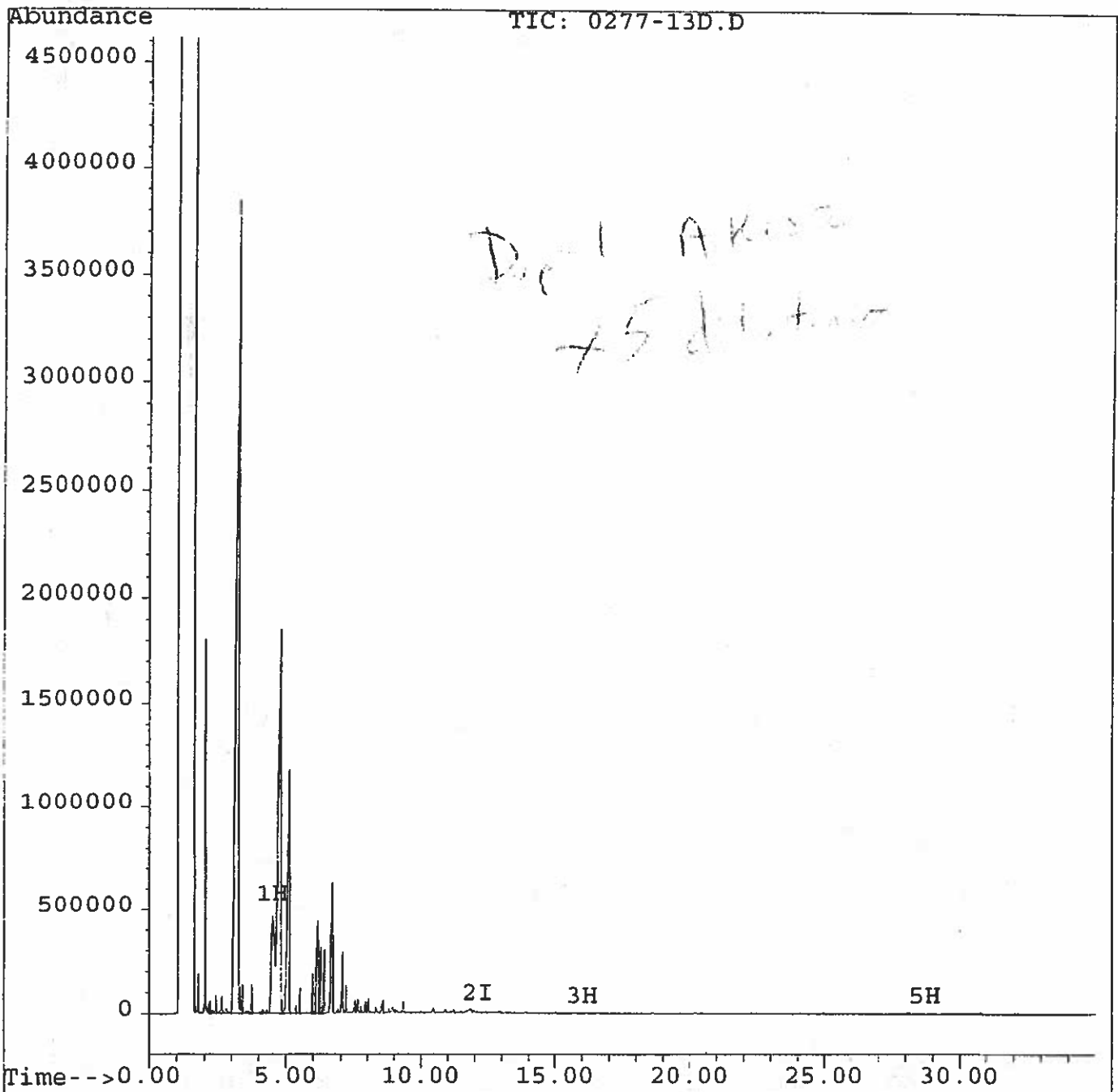
Quantitation Report

Data File : C:\HPCHEM\5\DATA\050796\0277-13D.D
Acq On : 09 May 96 05:19 AM
Sample : 0277-13 1L TO 1ML X5
Misc :
Quant Time: May 9 12:25 1996

Vial: 21
Operator:
Inst : FID-ECD
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\ADECTPHD.M
Title : 8015MOD
Last Update : Wed May 08 15:51:07 1996
Response via : Multiple Level Calibration

Volume Inj. :
Signal Phase :
Signal Info :





AGRA
Earth & Environmental
 3504 Industrial Avenue, Suite 5
 Fairbanks, Alaska, U.S.A. 99701
 Tel (907) 479-7586 Fax (907) 479-0193

02923

CHAIN OF CUSTODY

PROJECT	PROJECT NO.		CLIENT		PRESERVATIVE		MATRIX		DATE		TIME		CONTAINERS		ANALYSIS REQUESTED (circle, check box or write preferred method in box)															
	6-024-01173-1 Tox		SISKINS FORD - LINCOLN MERRURY		AL HAYNES		HCL		W		5/1/96		1425		1 L		BTEX by 5030 / 8020	GRPH by 5030 / 8015	DRPH by 3550 / 8100	BTEX/GRPH Combo by 5030 / 8020-8015	TPH by 3550 / 418.1	Halogenated Volatiles by 5030 / 8010	WTPH-418.1 MODIFIED	Aromatics by 602	Polynuclear Aromatics by 610 or 8310	Total Halogens (TOX) by 9078	Total Metals by ICP AA	Purgeable Organics GC/MS by 8240 or 824	Base/NeuVacid/Organics GC/MS by 825 or 8270	PCB by 8080
CLIENT	PHONE NO.		PHONE NO.		PHONE NO.		PHONE NO.		PHONE NO.		PHONE NO.		PHONE NO.		PHONE NO.															
PROJECT MANAGER	407) 479-7586		SAME		SAME		SAME		SAME		SAME		SAME		SAME															
SAMPLERS NAME (please print)	Keri DePalma		STEPHANIE FORESTER		STEPHANIE FORESTER		STEPHANIE FORESTER		STEPHANIE FORESTER		STEPHANIE FORESTER		STEPHANIE FORESTER		STEPHANIE FORESTER															
SAMPLERS SIGNATURE	<i>[Signature]</i>		<i>[Signature]</i>		<i>[Signature]</i>		<i>[Signature]</i>		<i>[Signature]</i>		<i>[Signature]</i>		<i>[Signature]</i>		<i>[Signature]</i>															
SAMPLE I.D.	1. GWP-2		2. MW-8		3.		4.		5.		6.		7.		8.		9.		10.											

SAMPLE RECEIPT	LABORATORY		TURNAROUND TIME		SPECIAL INSTRUCTIONS / ADDITIONAL COMMENTS	
	AEE PORTLAND		<input type="checkbox"/> 6 HOUR <input type="checkbox"/> 24 HOUR <input checked="" type="checkbox"/> 1 WEEK <input type="checkbox"/> 2 WEEK (standard) <input type="checkbox"/> OTHER		Please extract upon arrival to lab	
TOTAL # CONTAINERS	SHIPPING I.D. / AIRBILL #		ACCEPTED BY / AFFILIATION		DATE	
CONDITION OF CONTAINERS	CARRIER		TIME		DATE	
CONDITION OF SEALS	DOT DESIGNATION		TIME		DATE	
RELINQUISHED BY / AFFILIATION	DATE		TIME		DATE	
1. J. DePalma / AEE, FRKS	5/6/96		0800		5/7/96	
2.					930am	
3.					PAGE 1 OF 1	

AGRA Earth & Environmental, Inc. (7/84) DISTRIBUTION: White, Yellow - Laboratory, Pink - Originator AKG100077



AGRA Earth & Environmental, Inc.
7477 SW Tech Center Drive
Portland, Oregon
U.S.A. 97223-8025
Tel (503) 639-3400
Fax (503) 620-7892

May 28, 1996

AGRA Earth & Environmental
3504 Industrial Way, Suite 5
Fairbanks, AK 99701

Attention: Mr. James Spontak

Dear Mr. Spontak:

RE: Analytical Results For Project 6-024-01173-1T04

Attached are the PAH results for the samples submitted on May 3, 1996 from the above referenced project. For your reference, our project number associated with these samples is AK960277.

Volatile organic compounds by GC/MSD, diesel range organics, and gasoline range organics results were reported under separate cover in a package dated May 14, 1996.

All analyses were conducted in accordance with applicable QA/QC guidelines. The results apply only to the samples submitted.

Please feel free to contact me if you have any questions regarding this report, or if I can be of any assistance in any other matter.

Respectfully submitted,

AGRA Earth & Environmental

A handwritten signature in black ink, appearing to read "Sean Gormley", written over a horizontal line.

Sean Gormley
Laboratory Manager
Laboratory ID # UST-008

American Environmental Network, Inc.

17400 SW Upper Boones Ferry Rd. • Suite 270 • Durham, OR 97224 • (503) 684-0447
(Formerly Analytical Technologies, Inc.)

AEN I.D. 605531

May 20, 1996

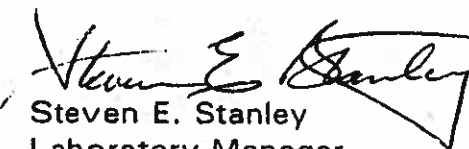
Sean Gormley
AGRA Earth & Environmental
7477 SW Tech Center Dr. Suite 135
Portland, OR 97223

Project Name / Number: Seekins Ford - Lincoln / 6-024-01173-1 T04

Attention: Sean Gormley

On May 6, 1996, American Environmental Network (OR), Inc. (ADEC UST-028) received three water samples for analysis for the above listed project. The samples were analyzed with EPA methodology or equivalent methods. The results of these analyses and the quality control data, which follow each set of analyses, are enclosed. The results from these samples relate only to the items tested. This report shall not be reproduced, except in full, without the written approval of the laboratory.

If you have any questions or comments, please do not hesitate to contact us at (503)684-0447.


Steven E. Stanley
Laboratory Manager

SES:alm
Enclosure

SAMPLE CROSS REFERENCE SHEET

CLIENT: AGRA Earth & Environmental **AEN I.D.:** 605531
PROJECT #: 6-024-01173-1T04
PROJECT NAME: Seekins Ford Lincoln Mercury **MATRIX:** WATER

<u>AEN #</u>	<u>CLIENT DESCRIPTION</u>	<u>DATE SAMPLED</u>
605531-1	MW-2	5/01/96
605531-2	MW-3	5/01/96
605531-3	GWP-3	5/01/96

-----TOTALS-----

<u>MATRIX</u>	<u># SAMPLES</u>
WATER	3

AEN STANDARD DISPOSAL PRACTICE

The samples from this project will be disposed of in thirty (30) days from the date of the report. If an extended storage period is required, please contact our sample control department before the scheduled disposal date.

ANALYTICAL SCHEDULE

CLIENT: AGRA Earth & Environmental AEN I.D.: 605531
PROJECT #: 6-024-01173-1T04
PROJECT NAME: Seekins Ford Lincoln Mercury

ANALYSIS	TECHNIQUE	REFERENCE	LAB
PAHs	HPLC/UV/FLUOR	EPA 8310	PLD

PLD = AEN - Portland
PHX = AEN - Phoenix
PNR = AEN - Pensacola
CA = AEN - Pleasant Hill
MD = AEN - Columbia
SUB = Subcontract

CASE NARRATIVE FOR EPA METHOD 8310

Polynuclear Aromatic Hydrocarbons

AEN ACCESSION NUMBER 605531

A 1.0 liter aliquot of each aqueous sample was extracted with methylene chloride by liquid:liquid method (EPA method 3510). The methylene chloride was reduced and exchanged to acetonitrile. Final volume was 1.0 ml.

The extracts were analyzed by HPLC with UV (254 nm) and fluorescence detectors in series. A gradient mobile phase of water and acetonitrile was utilized for the analysis. Five and six point calibration curves were generated and utilized for the analysis.

The sample GW-3 (AEN ID: 605531-3) yielded a surrogate recovery that was below the accepted criteria. This low recovery was most likely caused by interference from an emulsion that formed during the extraction.

Mark Ritzheimer

LIQUID CHROMATOGRAPHY RESULTS

METHOD:	8310	AEN I.D.:	605531-0
CLIENT I.D.:	Method Blank	DATE SAMPLED:	NA
CLIENT:	AGRA Earth & Environmental	DATE RECEIVED:	NA
PROJECT #:	6-024-01173-1T04	DATE EXTRACTED:	05/08/96
PROJECT NAME:	Seekins Ford Lincoln Mercury	DATE ANALYZED:	05/15/96
SAMPLE MATRIX:	WATER	DILUTION FACTOR:	1
		UNITS:	ug/L

PARAMETER	RESULTS
NAPHTHALENE	< 0.5
ACENAPHTHYLENE	< 1.0
ACENAPHTHENE	< 1.0
FLUORENE	< 0.1
PHENANTHRENE	< 0.05
ANTHRACENE	< 0.05
FLUORANTHENE	< 0.1
PYRENE	< 0.1
BENZO(a)ANTHRACENE	< 0.1
CHRYSENE	< 0.1
BENZO(b)FLUORANTHENE	< 0.1
BENZO(k)FLUORANTHENE	< 0.1
BENZO(a)PYRENE	< 0.1
DIBENZO(a,h)ANTHRACENE	< 0.1
BENZO(g,h,i)PERYLENE	< 0.1
INDENO(1,2,3-cd)PYRENE	< 0.1

SURROGATE:
 BIPHENYL (28%-125%) 45%

Analyst: GTC 5/20/96
 Reviewer: AH 5-20-96

LIQUID CHROMATOGRAPHY RESULTS

METHOD:	8310	AEN I.D.:	605531-1
CLIENT I.D.:	MW-2	DATE SAMPLED:	05/01/96
CLIENT:	AGRA Earth & Environmental	DATE RECEIVED:	05/06/96
PROJECT #:	6-024-01173-1T04	DATE EXTRACTED:	05/08/96
PROJECT NAME:	Seekins Ford Lincoln Mercury	DATE ANALYZED:	05/16/96
SAMPLE MATRIX:	WATER	DILUTION FACTOR:	1
		UNITS:	ug/L

PARAMETER	RESULTS
NAPHTHALENE	< 0.5
ACENAPHTHYLENE	< 1.0
ACENAPHTHENE	< 1.0
FLUORENE	< 0.1
PHENANTHRENE	< 0.05
ANTHRACENE	< 0.05
FLUORANTHENE	< 0.1
PYRENE	< 0.1
BENZO(a)ANTHRACENE	< 0.1
CHRYSENE	< 0.1
BENZO(b)FLUORANTHENE	< 0.1
BENZO(k)FLUORANTHENE	< 0.1
BENZO(a)PYRENE	< 0.1
DIBENZO(a,h)ANTHRACENE	< 0.1
BENZO(g,h,i)PERYLENE	< 0.1
INDENO(1,2,3-cd)PYRENE	< 0.1

SURROGATE:
 BIPHENYL (28%-125%) 53%

Analyst: CE 5/29/96

Reviewer: BIT 5-20-96

LIQUID CHROMATOGRAPHY RESULTS

METHOD:	8310	AEN I.D.:	605531-2
CLIENT I.D.:	MW-3	DATE SAMPLED:	05/01/96
CLIENT:	AGRA Earth & Environmental	DATE RECEIVED:	05/06/96
PROJECT #:	6-024-01173-1T04	DATE EXTRACTED:	05/08/96
PROJECT NAME:	Seakins Ford Lincoln Mercury	DATE ANALYZED:	05/16/96
SAMPLE MATRIX:	WATER	DILUTION FACTOR:	1
		UNITS:	ug/L

PARAMETER	RESULTS
NAPHTHALENE	9.0
ACENAPHTHYLENE	< 1.0
ACENAPHTHENE	< 1.0
FLUORENE	1.4
PHENANTHRENE	1.7
ANTHRACENE	< 0.05
FLUORANTHENE	1.1
PYRENE	0.3
BENZO(a)ANTHRACENE	< 0.1
CHRYSENE	< 0.1
BENZO(b)FLUORANTHENE	< 0.1
BENZO(k)FLUORANTHENE	< 0.1
BENZO(a)PYRENE	< 0.1
DIBENZO(a,h)ANTHRACENE	< 0.1
BENZO(g,h,i)PERYLENE	< 0.1
INDENO(1,2,3-cd)PYRENE	< 0.1

SURROGATE:
 BIPHENYL (28%-125%) 54%

Analyst: GTE 5/22/96

Reviewer: RAH 5-20-96

LIQUID CHROMATOGRAPHY RESULTS

METHOD:	8310	AEN I.D.:	605531-3
CLIENT I.D.:	GWP-3	DATE SAMPLED:	05/01/96
CLIENT:	AGRA Earth & Environmental	DATE RECEIVED:	05/06/96
PROJECT #:	6-024-01173-1T04	DATE EXTRACTED:	05/08/96
PROJECT NAME:	Seekins Ford Lincoln Mercury	DATE ANALYZED:	05/16/96
SAMPLE MATRIX:	WATER	DILUTION FACTOR:	1
		UNITS:	ug/L

PARAMETER	RESULTS
NAPHTHALENE	< 0.5
ACENAPHTHYLENE	< 1.0
ACENAPHTHENE	< 1.0
FLUORENE	< 0.1
PHENANTHRENE	< 0.05
ANTHRACENE	< 0.05
FLUORANTHENE	< 0.1
PYRENE	< 0.1
BENZO(a)ANTHRACENE	< 0.1
CHRYSENE	< 0.1
BENZO(b)FLUORANTHENE	< 0.1
BENZO(k)FLUORANTHENE	< 0.1
BENZO(a)PYRENE	< 0.1
DIBENZO(a,h)ANTHRACENE	< 0.1
BENZO(g,h,i)PERYLENE	< 0.1
INDENO(1,2,3-cd)PYRENE	< 0.1

SURROGATE:
 BIPHENYL (28 %-125%) 11% H

H = Out of limits

Analyst: GE 5/20/96
 Reviewer: BL 5-20-96

LIQUID CHROMATOGRAPHY SPIKE RESULTS

METHOD:	8310	ATI I.D.:	605531
CLIENT:	AGRA Earth & Environmental	QC SAMPLE:	605518-7
PROJECT #:	6-024-01173-1T04	DATE EXTRACTED:	05/08/96
PROJECT NAME:	Seekins Ford Lincoln Mercury	DATE ANALYZED:	05/16/96
SAMPLE MATRIX:	WATER	DILUTION FACTOR:	1
		UNITS:	ug/L

PARAMETER		SAMPLE RESULT	SPIKE CONC.	SPIKED RESULT	% REC.	DUP. SPIKED RESULT	DUP. % REC.	RPD
ACENAPHTHYLENE	<	1.0	10.0	3.1	31%	3.3	33%	6
PHENANTHRENE	<	0.05	1.00	0.54	54%	0.52	52%	4
PYRENE	<	0.1	1.0	0.5	50%	0.5	50%	0
BENZO(k)FLUORANTHENE	<	0.1	1.0	0.4	40%	0.3	30%	29
DIBENZO(a,h)ANTHRACENE	<	0.1	1.0	0.4	40%	0.3	30%	29
SURROGATE:								
BIPHENYL (28% - 125%)				42%		45%		

CONTROL LIMITS

	% REC	RPD
ACENAPHTHYLENE	27 - 117	42
PHENANTHRENE	42 - 127	31
PYRENE	50 - 125	35
BENZO(k)FLUORANTHENE	30 - 125	33
DIBENZO(a,h)ANTHRACENE	30 - 146	37

Analyst: MR 5/20/96

Reviewer: BA 5/20/96