

GW - 32

REPORTS

YEAR(S):

1991 ANALYTICAL
DATA 4/91 → 5/91



Route 3, Box 7
Gallup, New Mexico
87301

505
722-3833

June 12, 1991

Roger Anderson
New Mexico Oil Conservation Division
P.O. Box 2088
State Land Office Building
Santa Fe, New Mexico 87504

RE: Analytical Data for GW-32 Renewal Request

Dear Roger:

The attached data includes the analytical associated with all samples collected by Giant Refining Company during the April 30, 1991 to May 3, 1991 audit and inspection conducted by your office. The attachments also include the "Ground Water Monitoring Sample Record" or field log. This log includes field data such as water levels, pH, conductance, temperature, etc.. The ground water samples consist of well samples, two equipment washes and a duplicate sample of OW-1, labeled as OW-30.

This data should aid your office in the preparation of comments for the renewal request of Ground Water Discharge Permit GW-32.

If you have any questions, contact my office at (505) 722-0217.

Sincerely,

Claud Rosendale
Environmental Manager
Ciniza Refinery

List of Attachments: ATI Analytical and Field Sample Logs for MW-4, OW-1, OW-2, OW-5, OW-7, OW-9, OW-10, OW-12, OW-13, OW-14, OW-20, OW-16, OW-25, OW-26, Duplicate of OW-1, Equipment Wash (2) and Trip Blanks (2). ATI Analytical for Aeration Lagoon Effluents

cc w/attachments: Kim Bullerdick - Corporate Counsel
Giant Industries Arizona, Inc.

91 JUN 18 10 11 AM '91
OIL CONSERVATION DIVISION
RECEIVED

Ground Water Monitoring
Sample Record

Plant Refinery
(allip), New Mexico

Job Number AEI-SM#2
Location Grat-Liniz
Date 4-30-91
Weather _____

Purpose AEI-SM#2 * PCD

NOTE: 2-in Sch 40=0.163 gal/ft or 0
4-in Sch 40=0.653 gal/ft or 0
5-in Sch 40=1.020 gal/ft or 1

Sample on 4-29-91 Sampled on 4-30-91

Well Number	Water Level	Casing Storage	Purge Time	Total Purged (gal)	Sample Time	PH	Temp	Cord	Sample collection Method
<u>Egyp. Blank</u>	Total Depth, TOC, ft <u>X</u> Depth to Water, TOC, ft <u>X</u> Length of Water Column _____	Casing Diam _____ 1 Casing Vol (gal) _____	Begin Purge <u>X</u> End Purge _____	<u>X</u> Dry Y N	<u>0700</u>	<u>5.04</u> <u>X</u>	<u>39</u> <u>X</u>	<u>60</u> <u>X</u>	<u>Baker</u>
<u>Form DW-1</u>	Total Depth, TOC, ft _____ Depth to Water, TOC, ft _____ Length of Water Column _____	Casing Diam _____ 1 Casing Vol (gal) _____	Begin Purge _____ End Purge _____	_____ Dry Y N	_____	_____	_____	_____	_____
	Total Depth, TOC, ft _____ Depth to Water, TOC, ft _____ Length of Water Column _____	Casing Diam _____ 1 Casing Vol (gal) _____	Begin Purge _____ End Purge _____	_____ Dry Y N	_____	_____	_____	_____	_____
	Total Depth, TOC, ft _____ Depth to Water, TOC, ft _____ Length of Water Column _____	Casing Diam _____ 1 Casing Vol (gal) _____	Begin Purge _____ End Purge _____	_____ Dry Y N	_____	_____	_____	_____	_____
	Total Depth, TOC, ft _____ Depth to Water, TOC, ft _____ Length of Water Column _____	Casing Diam _____ 1 Casing Vol (gal) _____	Begin Purge _____ End Purge _____	_____ Dry Y N	_____	_____	_____	_____	_____

Ground Water Monitoring
Sample Record

Job Number RI-I SWM#2
 Location Grapt-Ciaia
 Date 4-30-91
 Weather _____

Grant Refinery
 Gulfport, New Mexico

NOTE: 2-in Sch 40=0.163 gal/ft or
 4-in Sch 40=0.653 gal/ft or
 5-in Sch 40=1.020 gal/ft or

Purpose RI-I-SWM#2 & OGD

Sampled on 4-29-91 Sampled on 4-30-91

Well Number	Water Level	Casing Storage	Purge Time	Total Purged (gal)	Sample Time	pH	Temp	Cord	Sample Collection Method
MW-4	Total Depth, TOC, ft <u>129.99</u> Depth to Water, TOC, ft <u>5.80</u> Length of Water Column <u>124.19</u>	Casing Diam <u>4"</u> 1 Casing Vol (gal) <u>5.88</u> <u>0.74</u>	Begin Purge <u>2:25</u> End Purge <u>2:53</u>	<u>280</u> Dry Vol <u>(M)</u>	<u>0850</u>	<u>8.68</u>	<u>49</u>	<u>1220</u>	Ba. 1
OW-5	Total Depth, TOC, ft <u>23.67</u> Depth to Water, TOC, ft <u>14.02</u> Length of Water Column <u>9.65</u>	Casing Diam <u>4"</u> 1 Casing Vol (gal) <u>15.68</u> <u>0.74</u>	Begin Purge <u>2:09</u> End Purge <u>2:11</u>	<u>18</u> Dry Vol <u>(M)</u>	<u>0755</u>	<u>6.47</u> <u>6.68</u>	<u>48</u> <u>49</u>	<u>30200</u> <u>39700</u>	Ba. 1
OW-7	Total Depth, TOC, ft <u>25.35</u> Depth to Water, TOC, ft <u>3.79</u> Length of Water Column <u>21.56</u>	Casing Diam <u>4"</u> 1 Casing Vol (gal) <u>4.12</u> <u>0.74</u>	Begin Purge <u>1:42</u> End Purge <u>1:59</u>	<u>170</u> Dry Vol <u>(M)</u>	<u>0915</u>	<u>8.75</u> <u>8.82</u>	<u>47</u> <u>48</u>	<u>1230</u> <u>1280</u>	Ba. 1
OW-1	Total Depth, TOC, ft <u>99.73</u> Depth to Water, TOC, ft <u>0</u> Length of Water Column <u>99.73</u>	Casing Diam <u>4"</u> 1 Casing Vol (gal) <u>58.11</u> <u>0.74</u>	Begin Purge <u>1:17</u> End Purge <u>1:25</u>	<u>80</u> Dry Vol <u>(M)</u>	<u>7:08</u> <u>Am</u>	<u>8.68</u> <u>8.64</u>	<u>51</u> <u>51</u>	<u>1262</u> <u>1275</u>	Ba. 1
OW-30	Total Depth, TOC, ft _____ Depth to Water, TOC, ft _____ Length of Water Column _____	Casing Diam <u>4"</u> 1 Casing Vol (gal) <u>0.74</u>	Begin Purge _____ End Purge _____	<u>(M)</u> Dry Vol _____	<u>0730</u>	<u>8.63</u>	<u>51</u>	<u>1276</u>	Ba. 1

OW-1 Duplicate

FIGURE 1

Ground Water Monitoring
Sample Record
Glant Refinery
Gallup, New Mexico

Job Number _____
Location _____
Date 1/16/87
Weather SE Wind 5-15 mph

Purpose _____

NOTE: 2-in Sch 40=0.163 gal/ft or 0
4-in Sch 40=0.653 gal/ft or 0
5-in Sch 40=1.020 gal/ft or 1

Well Number	Water Level	Casing Storage	Purge Time	Total Purged (gal)	Sample Time	pH	Temp	Cord	Sample Collection Method
OU 2	Total Depth, TOC, ft <u>64.41</u> Depth to Water, TOC, ft <u>28.80</u> Length of Water Column <u>35.61</u>	Casing Diam <u>4</u> 1 Casing Vol (gal) <u>74</u> <u>28.85</u>	Begin Purge <u>4:27</u> End Purge <u>4:48</u>	<u>125</u> Dry Y/N <u>N</u>	<u>0915</u>	<u>7.80</u>	<u>50</u>	<u>1360</u>	
OU 9	Total Depth, TOC, ft <u>45.12</u> Depth to Water, TOC, ft <u>.9</u> Length of Water Column <u>44.22</u>	Casing Diam <u>4</u> 1 Casing Vol (gal) <u>74</u> <u>.45</u>	Begin Purge <u>8:06</u> End Purge <u>5:17</u>	<u>125</u> Dry Y/N <u>N</u>	<u>0835</u>	<u>7.74</u>	<u>45</u>	<u>1840</u>	
OU 10	Total Depth, TOC, ft <u>61.29</u> Depth to Water, TOC, ft <u>0</u> Length of Water Column <u>61.29</u>	Casing Diam <u>4</u> 1 Casing Vol (gal) <u>74</u> <u>0"</u>	Begin Purge <u>5:24</u> End Purge <u>5:35</u>	<u>100</u> Dry Y/N <u>N</u>	<u>0815</u>	<u>7.46</u>	<u>52</u>	<u>1010</u>	
KGUP V65h	Total Depth, TOC, ft <u>X</u> Depth to Water, TOC, ft <u>X</u> Length of Water Column <u>X</u>	Casing Diam <u>X</u> 1 Casing Vol (gal) <u>X</u>	Begin Purge <u>X</u> End Purge <u>From DW-9</u>	<u>X</u> Dry Y/N <u>N</u>	<u>0850</u>	<u>5.08</u>	<u>47</u>	<u>60</u>	
	Total Depth, TOC, ft _____ Depth to Water, TOC, ft _____ Length of Water Column _____	Casing Diam _____ 1 Casing Vol (gal) _____	Begin Purge _____ End Purge _____	_____					

Ground Water Monitoring
Sample Record

Glant Refinery

Callup, New Mexico

Job Number _____

Location _____

Date _____

Weather _____

NOTE: 2-in Sch 40=0.163 gal/ft or
4-in Sch 40=0.653 gal/ft or
5-in Sch 40=1.020 gal/ft or

Purpose _____

Well Number	Water Level	Casing Storage	Purge Time	Total Purged (gal)	Sample Time	pH	Temp	Cond	Sample Collection Method
OW12	Total Depth, TOC, ft 131.51 Depth to Water, TOC, ft 47.38 Length of Water Column 84.13 HNV-0	Casing Diam 4 1 Casing Vol (gal) 16.09 1.74	Begin Purge 9:50 End Purge 9:00 75 GAL	400 DRY Y N	1:45	11.33	57	1770	
OW13	Total Depth, TOC, ft 105.60 Depth to Water, TOC, ft 23.11 Length of Water Column 82.49 HNV-0	Casing Diam 4 1 Casing Vol (gal) 23.39 1.74	Begin Purge 8:57 End Purge 9:19 220 GAL	- DRY Y N	2:25	7.99	53	1194	
OW14	Total Depth, TOC, ft 46.71 Depth to Water, TOC, ft 25.93 Length of Water Column 20.78 HNV-0	Casing Diam 4 1 Casing Vol (gal) 20.78 1.74	Begin Purge 10:17 End Purge 10:24 60 GAL	- DRY Y N	2:55	7.55	54	1571	
OW20	Total Depth, TOC, ft 66.10 Depth to Water, TOC, ft 55.84 Length of Water Column 10.26 HNV-2	Casing Diam 4 1 Casing Vol (gal) 2.74	Begin Purge 2:35 End Purge 2:32 12 GAL	- DRY Y N	3:30	12.06	56	5410	
	Total Depth, TOC, ft _____ Depth to Water, TOC, ft _____ Length of Water Column _____	Casing Diam _____ 1 Casing Vol (gal) _____	Begin Purge _____ End Purge _____	_____ DRY Y N					

Ground Water Monitoring
Sample Record

Job Number DCD Discharge

Location Discharge

Date 4-29

Weather Sunny Wind West 10-30 mph
250°F

Plant Refinery

(Phillips), New Mexico

NOTE: 2-in Sch 40=0.163 gal/ft or 0
4-in Sch 40=0.653 gal/ft or 0
5-in Sch 40=1.020 gal/ft or 1

Purpose DCD

Well Number	Water Level	Casing Storage	Total		Sample Purged (gal)	Sample Time	pH	Temp	Cond	Sample Collection Method
			Purge Time	Purge Line						
DW-26	Total Depth, TOC, ft 53.12	Casing Diam 4" 1 Casing Vol (gal) 0.74	Begin Purge 0815	2.69 / 0830	7.12	54	1.910	Pump		
	Depth to Water, TOC, ft 31.93		End Purge 0821							
	Length of Water Column 21.19		contained visible oily hydrocarbons							
	TLV - 10.1 gpc									
DW-16	Total Depth, TOC, ft 56.81	Casing Diam 4" 1 Casing Vol (gal) 0.74	Begin Purge 0845	6.99 / 0815	8.41	52	1.210	Pump		
	Depth to Water, TOC, ft 27.39		End Purge 0856							
	Length of Water Column 29.42		clear water							
	TLV - 0									
DW-25	Total Depth, TOC, ft 51.98	Casing Diam 4" 1 Casing Vol (gal) 0.74	Begin Purge 0930	6.99 / 1000	7.92	46	1.280	Pump		
	Depth to Water, TOC, ft 45.27		End Purge 0933							
	Length of Water Column 6.71		clear water							
	TLV - 0									
	Total Depth, TOC, ft	Casing Diam	Begin Purge	Dry Y N						
	Depth to Water, TOC, ft		End Purge							
	Length of Water Column									
	Total Depth, TOC, ft	Casing Diam	Begin Purge	Dry Y N						
	Depth to Water, TOC, ft		End Purge							
	Length of Water Column									

FIGURE 1



Analytical**Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 105511

June 10, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: RFI

Attention: Claud Rosendale

On 05/01/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

DATE RECEIVED : 05/01/91

REPORT DATE : 05/29/91

ATI I.D. : 105511

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	OW-1	AQUEOUS	04/30/91
02	OW-30	AQUEOUS	04/30/91
03	EQUIP. WASH	AQUEOUS	04/30/91

----- TOTALS -----

MATRIX	# SAMPLES
AQUEOUS	3

ATI STANDARD DISPOSAL PRACTICE

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



Analytical Technologies, Inc.

GENERAL CHEMISTRY RESULTS

ATI I.D. : 105511

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 05/01/91

PROJECT # : (NONE)

PROJECT NAME : RFI

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02	03
PH	UNITS	8.86	8.85	5.79



Analytical **Technologies**, Inc.

GENERAL CHEMISTRY - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

ATI I.D. : 105511

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
PH	UNITS	10553702	8.36	8.37	0.1	NA	NA	NA

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

METALS RESULTS

ATI I.D. : 105511

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 05/01/91

PROJECT # : (NONE)

PROJECT NAME : RFI

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02	03
ARSENIC	MG/L	<0.005	<0.005	<0.005
BARIUM	MG/L	0.128	0.130	<0.010
BERYLLIUM	MG/L	<0.005	<0.005	<0.005
CADMIUM	MG/L	0.006	<0.005	<0.005
COBALT	MG/L	<0.010	<0.010	<0.010
CHROMIUM	MG/L	<0.010	<0.050	<0.010
COPPER	MG/L	0.017	0.019	<0.010
MERCURY	MG/L	<0.0002	<0.0002	<0.0002
POTASSIUM	MG/L	2.6	2.2	<1.0
NICKEL	MG/L	<0.020	<0.020	<0.020
LEAD	MG/L	0.008	0.010	<0.002
ANTIMONY	MG/L	<0.05	<0.05	<0.05
SELENIUM	MG/L	0.007	0.007	<0.005
VANADIUM	MG/L	0.050	0.051	<0.010
ZINC	MG/L	0.033	0.036	0.012



Analytical Technologies, Inc.

METALS - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

ATI I.D. : 105511

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
ARSENIC	MG/L	10551904	<0.005	<0.005	NA	0.040	0.050	80
ARSENIC	MG/L	10551103	<0.005	<0.005	NA	0.038	0.050	76
BARIUM	MG/L	10551003	0.067	0.068	1	1.01	1.00	94
BERYLLIUM	MG/L	10551103	<0.005	<0.005	NA	0.465	0.500	93
CADMIUM	MG/L	10551003	<0.005	<0.005	NA	0.490	0.500	98
COBALT	MG/L	10551003	<0.010	<0.010	NA	0.960	1.00	96
CHROMIUM	MG/L	10551003	<0.010	<0.010	NA	0.929	1.00	93
COPPER	MG/L	10551003	0.018	0.018	0	0.502	0.500	97
MERCURY	MG/L	10491402	<0.0002	<0.0002	NA	0.0049	0.0050	98
POTASSIUM	MG/L	10551101	2.6	2.6	0	56.4	50.0	108
NICKEL	MG/L	10551003	<0.020	<0.020	NA	0.949	1.00	95
LEAD	MG/L	10551102	0.010	0.008	22	0.061	0.050	102
LEAD	MG/L	10551103	<0.002	<0.002	NA	0.047	0.050	94
ANTIMONY	MG/L	10551003	<0.05	<0.05	NA	0.94	1.00	94
SELENIUM	MG/L	10488708	<0.005	<0.005	NS	0.037	0.050	74
SELENIUM	MG/L	10551103	<0.005	<0.005	NA	0.049	0.050	98
VANADIUM	MG/L	10551003	0.024	0.022	9	1.00	1.00	98
ZINC	MG/L	10551003	0.015	0.011	31	0.506	0.500	98

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



GCMS - RESULTS

ATI I.D. : 10551101

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-1	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	105
BROMOFLUOROBENZENE (%)	82
TOLUENE-D8 (%)	106



GCMS - RESULTS

ATI I.D. : 10551102

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-30	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
-----------	---------

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	110
BROMOFLUOROBENZENE (%)	106
TOLUENE-D8 (%)	108



GCMS - RESULTS

ATI I.D. : 10551103

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: EQUIP. WASH	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	105
BROMOFLUOROBENZENE (%)	90
TOLUENE-D8 (%)	109



Analytical Technologies, Inc.

GCMS - RESULTS

REAGENT BLANK

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 105511
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/13/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/13/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	106
BROMOFLUOROBENZENE (%)	101
TOLUENE-D8 (%)	105



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 105511

TEST : VOLATILE ORGANICS (EPA 8240)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599906

DATE ANALYZED : 05/13/91

SAMPLE MATRIX : SOIL

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	DUP. %		DUP. %	RPD
	RESULT	SPIKED		REC. SAMPLE	REC. SAMPLE		
1,1-DICHLOROETHENE	<1	50	59	118	53	106	11
TRICHLOROETHENE	<1	50	47	94	51	102	8
CHLOROBENZENE	<5	50	49	98	53	106	8
TOLUENE	<5	50	53	106	58	116	9
BENZENE	<5	50	51	102	51	102	0

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

TR - Compound detected at an unquantifiable trace level



GCMS - RESULTS

ATI I.D. : 10551101

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/01/91
CLIENT I.D.	: OW-1	DATE ANALYZED	: 05/11/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO(a)ANTHRACENE	<5
BIS(2-ETHYLHEXYL)PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO(b)FLUORANTHENE	<5
BENZO(k)FLUORANTHENE	<5
BENZO(a)PYRENE	<5
DIBENZO(a,h)ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO(A,J)ACRIDINE	<5 *
7,12-DIMETHYLBENZ(a)ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

ADDITIONAL COMPOUNDS

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10551101

COMPOUNDS

RESULTS

ADDITIONAL COMPOUND

HYDROCARBON C10

40

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)

54

2-FLUOROBIPHENYL (%)

58

TERPHENYL (%)

83

PHENOL-D6 (%)

33

2-FLUOROPHENOL (%)

47

2,4,6-TRIBROMOPHENOL (%)

42



GCMS - RESULTS

ATI I.D. : 10551102

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI
CLIENT I.D. : OW-30
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/30/91
DATE RECEIVED : 05/01/91
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/11/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS

RESULTS

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10551102

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	63
2-FLUOROBIPHENYL (%)	55
TERPHENYL (%)	92
PHENOL-D6 (%)	35
2-FLUOROPHENOL (%)	49
2,4,6-TRIBROMOPHENOL (%)	46



GCMS - RESULTS

ATI I.D. : 10551103

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/01/91
CLIENT I.D.	: EQUIP. WASH	DATE ANALYZED	: 05/11/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10551103

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	50
2-FLUOROBIPHENYL (%)	48
TERPHENYL (%)	111
PHENOL-D6 (%)	34
2-FLUOROPHENOL (%)	52
2,4,6-TRIBROMOPHENOL (%)	35



GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 105511
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/01/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/10/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<25
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<5
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10551101

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	41
2-FLUOROBIPHENYL (%)	49
TERPHENYL (%)	96
PHENOL-D6 (%)	57
2-FLUOROPHENOL (%)	71
2,4,6-TRIBROMOPHENOL (%)	40



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 105511

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599911

DATE ANALYZED : 05/09/91

SAMPLE MATRIX :

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED RESULT	SPIKED SAMPLE	% REC.	DUP.	DUP.	RPD
	RESULT	SPIKED				SPIKED SAMPLE	% REC.	
1,2,4-TRICHLOROBENZENE	<10	50	37	74	34	68		8
ACENAPHTHENE	<10	50	36	72	41	82		13
2,4-DINITROTOLUENE	<25	50	31	62	34	68		9
PYRENE	<5	50	63	126	57	114		10
N-NITROSO-DI-N-PROPYL AMINE	<10	50	38	76	43	86		12
1,4-DICHLOROBENZENE	<5	50	42	84	41	82		2
PENTACHLOROPHENOL	<50	100	88	88	89	89		1
PHENOL	<5	100	77	77	79	79		3
2-CHLOROPHENOL	<10	100	84	84	82	82		2
4-CHLORO-3-METHYLPHENOL	<10	100	80	80	75	75		6
4-NITROPHENOL	<50	100	40	40	47	47		16

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$



DATE 4-30-91 PAGE 1 OF 1

ATI Labs: San Diego (619)458-9141 • Phoenix (602)438-1530 • Seattle (206)228-8335 • Pensacola (904)474-1001

DISTRIBUTION: White, Canary - ANALYTICAL TECHNOLOGIES, INC. • Pink - ORIGINATOR

TABLE -1
BACKGROUND METALS

Total Metals

<u>Parameter</u>	<u>Analytical Method</u>	<u>Reporting Limit mg/kg</u>
✓Antimony	6010	6.0
✓Arsenic	7060	0.5
✓Barium	6010	1.0
✓Beryllium	6010	0.2
✓Cadmium	6010	0.5
✓Chromium	6010	1.0
✓Cobalt	6010	1.0
✓Copper	6010	2.0
✓Lead	6010	5.0
✓Mercury	7471	0.2
✓Nickel	6010	4.0
✓Potassium	6010	500
✓Selenium	7740	0.5
✓Vanadium	6010	1.0
✓Zinc	6010	2.0

TABLE-4
SKINNER LIST

METHOD 8240

<u>Parameter</u>	<u>Reporting Limit (ug/kg)</u>
Benzene	500
Carbon disulfide	500
Chlorobenzene	500
2-Chloroethylvinyl ether	1,000
1,2-Dibromomethane	1,000
1,2-Dichloroethane	500
1,4-Dioxane	50,000
Ethyl Benzene	500
Methyl ethyl ketone (2-butanone)	1,000
Styrene	500
Toluene	500
Xylenes	500

METHOD 8270

Anthracene	5,000
Benzenethiol	-
Benzo(a)anthracene	5,000
Benzo(b)fluoranthene	5,000
Benzo(k)fluoranthene	5,000
Benzo(a)pyrene	5,000
Bis(2-ethylhexyl)phthalate	5,000
Butyl benzyl phthalate	5,000
Chrysene	5,000
Dibenzo(a,X)acridine	-
Dibenzo(a,h)anthracene	5,000
Di-n-butylphthalate	5,000
1,2-Dichlorobenzene	5,000
1,3-Dichlorobenzene	5,000
1,4-Dichlorobenzene	5,000
Diethyl phthalate	5,000
7,12-Dimethylbenz(a)anthracene	5,000
2,4-Dimethylphenol	5,000
Dimethyl phthalate	5,000
2,4-Dinitrophenol	25,000
Di-n-octyl phthalate	5,000
Fluoranthene	5,000
Indene	5,000
Methylchrysene	-
1-Methylnaphthalene	5,000
2-Methylphenol	5,000
3-Methylphenol	5,000
4-Methylphenol	5,000
Naphthalene	5,000
4-Nitrophenol	25,000
Phenanthrene	5,000

TABLE-4 Continued

Phenol.	5,000
Pyrene	5,000
Pyridine	10,000
Quinoline	25,000



Analytical **Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 105537

June 10, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: RFI

Attention: Claud Rosendale

On 05/02/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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 14 day volatile holding time for samples OW-9, OW-10, and the trip blank has been surpassed by one day. The samples were maintained at 4°C with hydrochloric acid so the data is thought to be reliable. The results for the 8240 analysis are provided at no charge.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

DATE RECEIVED : 05/02/91

REPORT DATE : 05/29/91

ATI I.D. : 105537

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	OW-9	AQUEOUS	05/01/91
02	OW-10	AQUEOUS	05/01/91
03	TRIP BLANK	AQUEOUS	05/01/91

----- TOTALS -----

MATRIX	# SAMPLES
AQUEOUS	3

ATI STANDARD DISPOSAL PRACTICE

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



Analytical **Technologies**, Inc.

GENERAL CHEMISTRY RESULTS

ATI I.D. : 105537

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 05/02/91

PROJECT # : (NONE)

PROJECT NAME : RFI

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02
PH	UNITS	8.48	8.36



Analytical Technologies, Inc.

GENERAL CHEMISTRY - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

ATI I.D. : 105537

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE CONC	SPIKE	% REC
PH	UNITS	10553702	8.36	8.37	0.1	NA	NA	NA

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

METALS RESULTS

ATI I.D. : 105537

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 05/02/91

PROJECT # : (NONE)

PROJECT NAME : RFI

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02
ARSENIC	MG/L	<0.005	<0.005
BARIUM	MG/L	0.044	0.081
BERYLLIUM	MG/L	<0.005	<0.005
CADMIUM	MG/L	<0.005	0.013
COBALT	MG/L	<0.010	<0.010
CHROMIUM	MG/L	<0.010	<0.010
COPPER	MG/L	0.014	<0.010
MERCURY	MG/L	<0.0002	<0.0002
POTASSIUM	MG/L	1.0	1.8
NICKEL	MG/L	<0.020	<0.020
LEAD	MG/L	0.002	<0.002
ANTIMONY	MG/L	<0.05	<0.05
SELENIUM	MG/L	0.013	0.024
VANADIUM	MG/L	0.012	<0.010
ZINC	MG/L	0.023	0.012



Analytical Technologies, Inc.

METALS - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

ATI I.D. : 105537

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
ARSENIC	MG/L	10554901	<0.005	<0.005	NA	0.052	0.050	104
BARIUM	MG/L	10553701	0.044	0.042	5	0.987	1.00	94
BARIUM	MG/L	10553702	0.081	0.080	1	0.178	0.100	97
BERYLLIUM	MG/L	10553401	0.007	0.007	0	0.105	0.100	98
BERYLLIUM	MG/L	10599912	<0.005	<0.005	NA	0.112	0.100	112
CADMIUM	MG/L	10553701	<0.005	<0.005	NA	0.500	0.500	100
CADMIUM	MG/L	10553702	0.013	0.014	7	0.109	0.100	96
COBALT	MG/L	10553701	<0.010	<0.010	NA	1.00	1.00	100
COBALT	MG/L	10553702	<0.010	<0.010	NA	0.110	0.100	110
CHROMIUM	MG/L	10553701	<0.010	<0.010	NA	0.944	1.00	94
CHROMIUM	MG/L	10553702	<0.010	<0.010	NA	0.089	0.100	89
COPPER	MG/L	10553701	0.014	0.014	0	0.504	0.500	98
COPPER	MG/L	10553702	<0.010	<0.010	NA	0.097	0.100	97
MERCURY	MG/L	10557001	<0.0002	<0.0002	NA	0.0048	0.0050	96
POTASSIUM	MG/L	10551101	2.6	2.6	0	56.4	50.0	108
POTASSIUM	MG/L	10557001	3.9	3.8	3	59.5	50.0	111
NICKEL	MG/L	10553701	<0.020	<0.020	NA	0.960	1.00	96
NICKEL	MG/L	10553702	<0.020	<0.020	NA	0.100	0.100	100
LEAD	MG/L	10551103	<0.002	<0.002	NA	0.047	0.050	94
ANTIMONY	MG/L	10553701	<0.05	<0.05	NA	0.93	1.00	93
ANTIMONY	MG/L	10599906	<0.05	<0.05	NA	0.09	0.10	90
SELENIUM	MG/L	10554901	<0.005	<0.005	NA	0.031	0.050	62
VANADIUM	MG/L	10553701	0.012	0.011	9	0.990	1.00	98
VANADIUM	MG/L	10553702	<0.010	<0.010	NA	0.104	0.100	104
ZINC	MG/L	10553701	0.023	0.023	0	0.527	0.500	101
ZINC	MG/L	10553702	0.012	0.012	0	0.122	0.100	110

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$

GCMS - RESULTS

ATI I.D. : 10553701

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-9	DATE ANALYZED	: 05/16/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	102
BROMOFLUOROBENZENE (%)	90
TOLUENE-D8 (%)	102



Analytical Technologies, Inc.

GCMS - RESULTS

ATI I.D. : 10553702

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-10	DATE ANALYZED	: 05/16/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	103
BROMOFLUOROBENZENE (%)	91
TOLUENE-D8 (%)	101



GCMS - RESULTS

ATI I.D. : 10553703

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: TRIP BLANK	DATE ANALYZED	: 05/16/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	103
BROMOFLUOROBENZENE (%)	87
TOLUENE-D8 (%)	99



GCMS - RESULTS

REAGENT BLANK

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 105537
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/16/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/16/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	99
BROMOFLUOROBENZENE (%)	88
TOLUENE-D8 (%)	102



Analytical Technologies, Inc.

QUALITY CONTROL DATA

TEST : VOLATILE ORGANICS (EPA 8240)

ATI I.D. : 105537

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599907

DATE ANALYZED : 05/16/91

SAMPLE MATRIX : SOIL

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	DUP.		DUP.	RPD
	RESULT	SPIKED		% SPIKED REC.	% SPIKED REC.		
1,1-DICHLOROETHENE	<1	50	55	110	54	108	2
TRICHLOROETHENE	<1	50	49	98	49	98	0
CHLOROBENZENE	<5	50	51	102	51	102	0
TOLUENE	<5	50	51	102	51	102	0
BENZENE	<5	50	41	82	41	82	0

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

TR - Compound detected at an unquantifiable trace level



GCMS - RESULTS

ATI I.D. : 10553701

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/06/91
CLIENT I.D.	: OW-9	DATE ANALYZED	: 05/15/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS

RESULTS

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO(a) ANTHRACENE	<5
BIS(2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO(b) FLUORANTHENE	<5
BENZO(k) FLUORANTHENE	<5
BENZO(a) PYRENE	<5
DIBENZO(a,h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO(A,J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ(a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10553701

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	45
2-FLUOROBIPHENYL (%)	46
TERPHENYL (%)	87
PHENOL-D6 (%)	37
2-FLUOROPHENOL (%)	57
2,4,6-TRIBROMOPHENOL (%)	42

GCMS - RESULTS

ATI I.D. : 10553702

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/06/91
CLIENT I.D.	: OW-10	DATE ANALYZED	: 05/15/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10553702

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	52
2-FLUOROBIPHENYL (%)	57
TERPHENYL (%)	107
PHENOL-D6 (%)	37
2-FLUOROPHENOL (%)	51
2,4,6-TRIBROMOPHENOL (%)	39



GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 105537
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/06/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/15/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
-----------	---------

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROTOLUENE	<5
2,6-DINITROTOLUENE	<5
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 105537

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	67
2-FLUOROBIPHENYL (%)	67
TERPHENYL (%)	107
PHENOL-D6 (%)	45
2-FLUOROPHENOL (%)	56
2,4,6-TRIBROMOPHENOL (%)	38



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 105537

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599915

DATE ANALYZED : 05/15/91

SAMPLE MATRIX :

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	% REC.	DUP. SPIKED %		RPD
	RESULT	SPIKED			SAMPLE	REC.	
1,2,4-TRICHLOROBENZENE	<10	50	31	62	33	66	6
ACENAPHTHENE	<10	50	32	64	37	74	14
2,4-DINITROTOLUENE	<25	50	26	52	31	62	18
PYRENE	<5	50	52	104	55	110	6
N-NITROSO-DI-N-PROPYL AMINE	<10	50	26	52	29	58	11
1,4-DICHLOROBENZENE	<5	50	37	74	38	76	3
PENTACHLOROPHENOL	<50	100	61	61	65	65	6
PHENOL	<5	100	50	50	52	104	4
2-CHLOROPHENOL	<10	100	52	52	58	58	11
4-CHLORO-3-METHYLPHENOL	<10	100	47	47	47	47	0
4-NITROPHENOL	<50	100	40	40	41	41	2

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$



Chain of Custody

DATE 5-1-91 PAGE 1 OF 1

[illegible]

ATI Labs: San Diego (619) 458-9141 • Phoenix (602) 438-1530 • Seattle (206) 228-8335 • Pensacola (904) 474-1001
DISTRIBUTION: White, Canary - ANALYTICAL TECHNOLOGIES, INC. • Pink - ORIGINATOR

TABLE -1
BACKGROUND METALS

Total Metals

<u>Parameter</u>	<u>Analytical Method</u>	<u>Reporting Limit mg/kg</u>
✓Antimony	6010	6.0
✓Arsenic	7060	0.5
✓Barium	6010	1.0
✓Beryllium	6010	0.2
✓Cadmium	6010	0.5
✓Chromium	6010	1.0
✓Cobalt	6010	1.0
✓Copper	6010	2.0
✓Lead	6010	5.0
✓Mercury	7471	0.2
✓Nickel	6010	4.0
✓Potassium	6010	500
✓Selenium	7740	0.5
✓Vanadium	6010	1.0
✓Zinc	6010	2.0

TABLE-4
SKINNER LIST

METHOD 8240

<u>Parameter</u>	<u>Reporting Limit (ug/kg)</u>
Benzene	500
Carbon disulfide	500
Chlorobenzene	500
2-Chloroethylvinyl ether	1,000
1,2-Dibromomethane	1,000
1,2-Dichloroethane	500
1,4-Dioxane	50,000
Ethyl Benzene	500
Methyl ethyl ketone (2-butanone)	1,000
Styrene	500
Toluene	500
Xylenes	500

METHOD 8270

Anthracene	5,000
Benzenethiol	-
Benzo(a)anthracene	5,000
Benzo(b)fluoranthene	5,000
Benzo(k)fluoranthene	5,000
Benzo(a)pyrene	5,000
Bis(2-ethylhexyl)phthalate	5,000
Butyl benzyl phthalate	5,000
Chrysene	5,000
Dibenzo(a,X)acridine	-
Dibenzo(a,h)anthracene	5,000
Di-n-butylphthalate	5,000
1,2-Dichlorobenzene	5,000
1,3-Dichlorobenzene	5,000
1,4-Dichlorobenzene	5,000
Diethyl phthalate	5,000
7,12-Dimethylbenz(a)anthracene	5,000
2,4-Dimethylphenol	5,000
Dimethyl phthalate	5,000
2,4-Dinitrophenol	25,000
Di-n-octyl phthalate	5,000
Fluoranthene	5,000
Indene	5,000
Methylchrysene	-
1-Methylnaphthalene	5,000
2-Methylphenol	5,000
3-Methylphenol	5,000
4-Methylphenol	5,000
Naphthalene	5,000
4-Nitrophenol	25,000
Phenanthrene	5,000

TABLE-4 Continued

Phenol.	5,000
Pyrene	5,000
Pyridine	10,000
Quinoline	25,000



Analytical**Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 105510

June 10, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: RFI

Attention: Claud Rosendale

On 05/01/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

Sample OW-5 had matrix interferences for several metals (excluding As, Pb, K, and Mercury) analyses. These samples had to be diluted to be analyzed. Detection limits have been raised accordingly.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

Lorraine Davis
QA Coordinator

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

DATE RECEIVED : 05/01/91

REPORT DATE : 05/29/91

ATI I.D. : 105510

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	MW-4	AQUEOUS	04/30/91
02	OW-5	AQUEOUS	04/30/91
03	OW-7	AQUEOUS	04/30/91
04	TRIP BLANK	AQUEOUS	04/30/91

----- TOTALS -----

MATRIX	# SAMPLES
AQUEOUS	4

ATI STANDARD DISPOSAL PRACTICE

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



Analytical Technologies, Inc.

GENERAL CHEMISTRY RESULTS

ATI I.D. : 105510

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 05/01/91

PROJECT # : (NONE)

PROJECT NAME : RFI

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02	03
PH	UNITS	8.77	6.98	8.87



Analytical **Technologies**, Inc.

GENERAL CHEMISTRY - QUALITY CONTROL

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

ATI I.D. : 105510

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
PH	UNITS	10551003	8.87	8.87	0	NA	NA	NA

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

METALS RESULTS

ATI I.D. : 105510

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 05/01/91

PROJECT # : (NONE)

PROJECT NAME : RFI

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02	03
ARSENIC	MG/L	<0.005	<0.005	<0.005
BARIUM	MG/L	0.024	0.085	0.067
BERYLLIUM	MG/L	<0.005	<0.010	<0.005
CADMIUM	MG/L	<0.005	<0.025	<0.005
COBALT	MG/L	<0.010	<0.050	<0.010
CHROMIUM	MG/L	<0.010	<0.010	<0.010
COPPER	MG/L	0.379	<0.050	0.018
MERCURY	MG/L	<0.0002	<0.0002	<0.0002
POTASSIUM	MG/L	1.4	7.9	2.0
NICKEL	MG/L	<0.020	<0.10	<0.020
LEAD	MG/L	0.002	0.021	0.003
ANTIMONY	MG/L	<0.05	<0.25	<0.05
SELENIUM	MG/L	<0.005	<0.05	<0.005
VANADIUM	MG/L	<0.010	<0.50	0.024
ZINC	MG/L	0.038	<0.050	0.015



Analytical Technologies, Inc.

METALS - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

ATI I.D. : 105510

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
ARSENIC	MG/L	10551904	<0.005	<0.005	NA	0.040	0.050	80
BARIUM	MG/L	10551003	0.067	0.068	1	1.01	1.00	94
BERYLLIUM	MG/L	10551103	<0.005	<0.005	NA	0.465	0.500	93
CADMIUM	MG/L	10551003	<0.005	<0.005	NA	0.490	0.500	98
COBALT	MG/L	10551003	<0.010	<0.010	NA	0.960	1.00	96
CHROMIUM	MG/L	10551003	<0.010	<0.010	NA	0.929	1.00	93
COPPER	MG/L	10551003	0.018	0.018	0	0.502	0.500	97
MERCURY	MG/L	10491402	<0.0002	<0.0002	NA	0.0049	0.0050	98
POTASSIUM	MG/L	10551101	2.6	2.6	0	56.4	50.0	108
NICKEL	MG/L	10551003	<0.020	<0.020	NA	0.949	1.00	95
LEAD	MG/L	10490101	0.009	0.009	0	0.057	0.050	96
ANTIMONY	MG/L	10551003	<0.05	<0.05	NA	0.94	1.00	94
SELENIUM	MG/L	10488708	<0.005	<0.005	NA	0.037	0.050	74
VANADIUM	MG/L	10551003	0.024	0.022	9	1.00	1.00	98
ZINC	MG/L	10551003	0.015	0.011	31	0.506	0.500	98

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

GCMS - RESULTS

ATI I.D. : 10551001

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: MW-4	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
-----------	---------

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLEETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	96
BROMOFLUOROBENZENE (%)	94
TOLUENE-D8 (%)	105



GCMS - RESULTS

ATI I.D. : 10551002

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-5	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	103
BROMOFLUOROBENZENE (%)	104
TOLUENE-D8 (%)	109



GCMS - RESULTS

ATI I.D. : 10551003

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-7	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	103
BROMOFLUOROBENZENE (%)	90
TOLUENE-D8 (%)	110



GCMS - RESULTS

ATI I.D. : 10551004

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: TRIP BLANK	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	106
BROMOFLUOROBENZENE (%)	92
TOLUENE-D8 (%)	107



Analytical Technologies, Inc.

GCMS - RESULTS

REAGENT BLANK

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	ATI I.D. :	105510
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/13/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/13/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLEETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	106
BROMOFLUOROBENZENE (%)	101
TOLUENE-D8 (%)	105



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 105510

TEST : VOLATILE ORGANICS (EPA 8240)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599906

DATE ANALYZED : 05/13/91

SAMPLE MATRIX : SOIL

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	%	DUP.		RPD
	RESULT	SPIKED			SPIKED SAMPLE	% REC.	
1,1-DICHLOROETHENE	<1	50	59	118	53	106	11
TRICHLOROETHENE	<1	50	47	94	51	102	8
CHLOROBENZENE	<5	50	49	98	53	106	8
TOLUENE	<5	50	53	106	58	116	9
BENZENE	<5	50	51	102	51	102	0

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

TR - Compound detected at an unquantifiable trace level

GCMS - RESULTS

ATI I.D. : 10551001

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/01/91
CLIENT I.D.	: MW-4	DATE ANALYZED	: 05/11/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO(a)ANTHRACENE	<5
BIS(2-ETHYLHEXYL)PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO(b)FLUORANTHENE	<5
BENZO(k)FLUORANTHENE	<5
BENZO(a)PYRENE	<5
DIBENZO(a,h)ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO(A,J)ACRIDINE	<5 *
7,12-DIMETHYLBENZ(a)ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10551001

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	42
2-FLUOROBIPHENYL (%)	46
TERPHENYL (%)	92
PHENOL-D6 (%)	38
2-FLUOROPHENOL (%)	51
2,4,6-TRIBROMOPHENOL (%)	44



Analytical Technologies, Inc.

GCMS - RESULTS

ATI I.D. : 10551002

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/01/91
CLIENT I.D.	: OW-5	DATE ANALYZED	: 05/11/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10551002

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	60
2-FLUOROBIPHENYL (%)	56
TERPHENYL (%)	97
PHENOL-D6 (%)	32
2-FLUOROPHENOL (%)	44
2,4,6-TRIBROMOPHENOL (%)	42



GCMS - RESULTS

ATI I.D. : 10551003

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/30/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/01/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/01/91
CLIENT I.D.	: OW-7	DATE ANALYZED	: 05/11/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO(a)ANTHRACENE	<5
BIS(2-ETHYLHEXYL)PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO(b)FLUORANTHENE	<5
BENZO(k)FLUORANTHENE	<5
BENZO(a)PYRENE	<5
DIBENZO(a,h)ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO(A,J)ACRIDINE	<5 *
7,12-DIMETHYLBENZ(a)ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10551003

COMPOUNDS

RESULTS

PYRIDINE

<5

QUINOLINE

<25

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)

59

2-FLUOROBIPHENYL (%)

58

TERPHENYL (%)

106

PHENOL-D6 (%)

10

2-FLUOROPHENOL (%)

21

2,4,6-TRIBROMOPHENOL (%)

14



GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 105510
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/01/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/10/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
-----------	---------

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO(a)ANTHRACENE	<5
BIS(2-ETHYLHEXYL)PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO(b)FLUORANTHENE	<5
BENZO(k)FLUORANTHENE	<5
BENZO(a)PYRENE	<5
DIBENZO(a,h)ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO(A,J)ACRIDINE	<5 *
7,12-DIMETHYLBENZ(a)ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 105510

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	41
2-FLUOROBIPHENYL (%)	49
TERPHENYL (%)	96
PHENOL-D6 (%)	57
2-FLUOROPHENOL (%)	71
2,4,6-TRIBROMOPHENOL (%)	40



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 105510

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599911

DATE ANALYZED : 05/09/91

SAMPLE MATRIX :

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	% REC.	DUP. SPIKED %		RPD
	RESULT	SPIKED			SAMPLE	REC.	
1,2,4-TRICHLOROBENZENE	<10	50	37	74	34	68	8
ACENAPHTHENE	<10	50	36	72	41	82	13
2,4-DINITROTOLUENE	<25	50	31	62	34	68	9
PYRENE	<5	50	63	126	57	114	10
N-NITROSO-DI-N-PROPYL AMINE	<10	50	38	76	43	86	12
1,4-DICHLOROBENZENE	<5	50	42	84	41	82	2
PENTACHLOROPHENOL	<50	100	88	88	89	89	1
PHENOL	<5	100	77	77	79	79	3
2-CHLOROPHENOL	<10	100	84	84	82	82	2
4-CHLORO-3-METHYLPHENOL	<10	100	80	80	75	75	6
4-NITROPHENOL	<50	100	40	40	47	47	16

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$



DATE 4-30-99 PAGE 1 OF 1

ATI Labs: San Diego (619)458-9141 • Phoenix (602)438-1530 • Seattle (206)228-8335 • Pensacola (904)474-1001
DISTRIBUTION: White, Canary - ANALYTICAL TECHNOLOGIES, INC. • Pink - ORIGINATOR

TABLE -1
BACKGROUND METALS

Total Metals

<u>Parameter</u>	<u>Analytical Method</u>	<u>Reporting Limit mg/kg</u>
✓Antimony	6010	6.0
✓Arsenic	7060	0.5
✓Barium	6010	1.0
✓Beryllium	6010	0.2
✓Cadmium	6010	0.5
✓Chromium	6010	1.0
✓Cobalt	6010	1.0
✓Copper	6010	2.0
✓Lead	6010	5.0
✓Mercury	7471	0.2
✓Nickel	6010	4.0
✓Potassium	6010	500
✓Selenium	7740	0.5
✓Vanadium	6010	1.0
✓Zinc	6010	2.0

TABLE-4
SKINNER LIST

METHOD 8240

<u>Parameter</u>	<u>Reporting Limit (ug/kg)</u>
Benzene	500
Carbon disulfide	500
Chlorobenzene	500
2-Chloroethylvinyl ether	1,000
1,2-Dibromomethane	1,000
1,2-Dichloroethane	500
1,4-Dioxane	50,000
Ethyl Benzene	500
Methyl ethyl ketone (2-butanone)	1,000
Styrene	500
Toluene	500
Xylenes	500

METHOD 8270

Anthracene	5,000
Benzenethiol	-
Benzo(a)anthracene	5,000
Benzo(b)fluoranthene	5,000
Benzo(k)fluoranthene	5,000
Benzo(a)pyrene	5,000
Bis(2-ethylhexyl)phthalate	5,000
Butyl benzyl phthalate	5,000
Chrysene	5,000
Dibenzo(a,X)acridine	-
Dibenzo(a,h)anthracene	5,000
Di-n-butylphthalate	5,000
1,2-Dichlorobenzene	5,000
1,3-Dichlorobenzene	5,000
1,4-Dichlorobenzene	5,000
Diethyl phthalate	5,000
7,12-Dimethylbenz(a)anthracene	5,000
2,4-Dimethylphenol	5,000
Dimethyl phthalate	5,000
2,4-Dinitrophenol	25,000
Di-n-octyl phthalate	5,000
Fluoranthene	5,000
Indene	5,000
Methylchrysene	-
1-Methylnaphthalene	5,000
2-Methylphenol	5,000
3-Methylphenol	5,000
4-Methylphenol	5,000
Naphthalene	5,000
4-Nitrophenol	25,000
Phenanthrene	5,000

TABLE-4 Continued

Phenol.	5,000
Pyrene	5,000
Pyridine	10,000
Quinoline	25,000



Analytical **Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 105534

June 10, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: RFI

Attention: Claud Rosendale

On 05/02/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

Due to matrix interferences, sample OW-1 had to be diluted for Selenium analysis, causing the detection limit to be raised.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

DATE RECEIVED : 05/02/91

REPORT DATE : 05/29/91

ATI I.D. : 105534

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	OW-2	AQUEOUS	05/01/91
02	EQUIP WASH #2	AQUEOUS	05/01/91

----- TOTALS -----

MATRIX	# SAMPLES
-----	-----
AQUEOUS	2

ATI STANDARD DISPOSAL PRACTICE

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



Analytical **Technologies**, Inc.

GENERAL CHEMISTRY RESULTS

ATI I.D. : 105534

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 05/02/91

PROJECT # : (NONE)

PROJECT NAME : RFI

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02
PH	UNITS	8.50	5.79



Analytical Technologies, Inc.

GENERAL CHEMISTRY - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

ATI I.D. : 105534

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
PH	UNITS	10553702	8.36	8.37	0.1	NA	NA	NA

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

METALS RESULTS

ATI I.D. : 105534

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

DATE RECEIVED : 05/02/91

REPORT DATE : 05/29/91

PARAMETER	UNITS	01	02
ARSENIC	MG/L	0.006	<0.005
BARIUM	MG/L	0.497	0.057
BERYLLIUM	MG/L	0.007	<0.005
CADMIUM	MG/L	<0.005	<0.005
COBALT	MG/L	<0.010	<0.010
CHROMIUM	*MG/L	0.023	<0.010
COPPER	MG/L	0.044	0.017
MERCURY	MG/L	<0.0002	<0.0002
POTASSIUM	MG/L	2.7	<1.0
NICKEL	MG/L	<0.020	<0.020
LEAD	*MG/L	0.026	<0.002
ANTIMONY	MG/L	<0.05	<0.05
SELENIUM	MG/L	<0.010	<0.005
VANADIUM	MG/L	0.051	<0.010
ZINC	MG/L	0.063	0.014



Analytical Technologies, Inc.

METALS - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : RFI

ATI I.D. : 105534

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
ARSENIC	MG/L	10554901	<0.005	<0.005	NA	0.052	0.050	104
BARIUM	MG/L	10553106	<0.010	<0.010	NA	0.098	0.100	98
BERYLLIUM	MG/L	10553401	0.007	0.007	0	0.105	0.100	98
CADMIUM	MG/L	10553701	<0.005	<0.005	NA	0.500	0.500	100
COBALT	MG/L	10553701	<0.010	<0.010	NA	1.00	1.00	100
CHROMIUM	MG/L	10553106	<0.010	<0.010	NA	0.088	0.100	88
COPPER	MG/L	10553106	<0.010	<0.010	NA	0.090	0.100	90
COPPER	MG/L	10553701	0.014	0.014	0	0.504	0.500	98
MERCURY	MG/L	10557001	<0.0002	<0.0002	NA	0.0048	0.0050	96
POTASSIUM	MG/L	10551101	2.6	2.6	0	56.4	50.0	108
NICKEL	MG/L	10553701	<0.020	<0.020	NA	0.960	1.00	96
LEAD	MG/L	10551103	<0.002	<0.002	NA	0.047	0.050	94
ANTIMONY	MG/L	10553701	<0.05	<0.05	NA	0.93	1.00	93
SELENIUM	MG/L	10490101	<0.005	<0.005	NA	0.030	0.050	60
VANADIUM	MG/L	10553701	0.012	0.011	9	0.990	1.00	98
ZINC	MG/L	10553106	0.015	0.016	6	0.123	0.100	108

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$

GCMS - RESULTS

ATI I.D. : 10553401

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-2	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	108
BROMOFLUOROBENZENE (%)	81
TOLUENE-D8 (%)	108

GCMS - RESULTS

ATI I.D. : 10553402

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: N/A
CLIENT I.D.	: EQUIP WASH #2	DATE ANALYZED	: 05/13/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	108
BROMOFLUOROBENZENE (%)	86
TOLUENE-D8 (%)	102



GCMS - RESULTS

REAGENT BLANK

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 105534
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/13/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/13/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<2.5

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	106
BROMOFLUOROBENZENE (%)	101
TOLUENE-D8 (%)	105



Analytical Technologies, Inc.

QUALITY CONTROL DATA

TEST : VOLATILE ORGANICS (EPA 8240)

ATI I.D. : 105534

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599906

DATE ANALYZED : 05/13/91

SAMPLE MATRIX : SOIL

UNITS : UG/L

COMPOUNDS	SAMPLE RESULT	CONC. SPIKED	SPIKED SAMPLE	% REC.	DUP.	DUP.	RPD
					SPIKED	% REC.	
1,1-DICHLOROETHENE	<1	50	59	118	53	106	11
TRICHLOROETHENE	<1	50	47	94	51	102	8
CHLOROBENZENE	<5	50	49	98	53	106	8
TOLUENE	<5	50	53	106	58	116	9
BENZENE	<5	50	51	102	51	102	0

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

TR - Compound detected at an unquantifiable trace level



GCMS - RESULTS

ATI I.D. : 10553401

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/06/91
CLIENT I.D.	: OW-2	DATE ANALYZED	: 05/15/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
-----------	---------

PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10553401

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	53
2-FLUOROBIPHENYL (%)	58
TERPHENYL (%)	114
PHENOL-D6 (%)	36
2-FLUOROPHENOL (%)	51
2,4,6-TRIBROMOPHENOL (%)	35

GCMS - RESULTS

ATI I.D. : 10553402

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/01/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/02/91
PROJECT NAME	: RFI	DATE EXTRACTED	: 05/06/91
CLIENT I.D.	: EQUIP WASH #2	DATE ANALYZED	: 05/15/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 10553402

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	56
2-FLUOROBIPHENYL (%)	59
TERPHENYL (%)	128
PHENOL-D6 (%)	38
2-FLUOROPHENOL (%)	54
2,4,6-TRIBROMOPHENOL (%)	34



GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 105534
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/06/91
PROJECT NAME	: RFI	DATE ANALYZED	: 05/15/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
PHENOL	<5
1,3-DICHLOROBENZENE	<5
1,4-DICHLOROBENZENE	<5
1,2-DICHLOROBENZENE	<5
2-METHYLPHENOL	<5
4-METHYLPHENOL	<5
2,4-DIMETHYLPHENOL	<5
NAPHTHALENE	<5
DIMETHYLPHTHALATE	<5
2,4-DINITROPHENOL	<25
4-NITROPHENOL	<25
DIETHYLPHTHALATE	<5
PHENANTHRENE	<5
ANTHRACENE	<5
DI-N-BUTYLPHTHALATE	<5
FLUORANTHENE	<5
PYRENE	<5
BUTYLBENZYLPHTHALATE	<5
BENZO (a) ANTHRACENE	<5
BIS (2-ETHYLHEXYL) PHTHALATE	<5
CHRYSENE	<5
DI-N-OCTYLPHTHALATE	<5
BENZO (b) FLUORANTHENE	<5
BENZO (k) FLUORANTHENE	<5
BENZO (a) PYRENE	<5
DIBENZO (a, h) ANTHRACENE	<5
BENZENETHIOL	<5
DIBENZO (A, J) ACRIDINE	<5 *
7,12-DIMETHYLBENZ (a) ANTHRACENE	<5
INDENE	<5
METHYLCHRYSENE	<5
1-METHYLNAPHTHALENE	<5
3-METHYLPHENOL	<5
PYRIDINE	<5
QUINOLINE	<25

* ESTIMATED PRACTICAL QUANTITATION LIMIT BASED ON A REVIEW OF THE MASS SPECTRA DATA.



Analytical Technologies, Inc.

TEST : SEMI-VOLATILE ORGANICS (SKINNER LIST) EPA 8270

ATI I.D. : 105534

COMPOUNDS

RESULTS

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	67
2-FLUOROBIPHENYL (%)	67
TERPHENYL (%)	107
PHENOL-D6 (%)	45
2-FLUOROPHENOL (%)	56
2,4,6-TRIBROMOPHENOL (%)	38



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D.

: 105534

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : RFI

REF I.D. : 10599915

DATE ANALYZED : 05/15/91

SAMPLE MATRIX :

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED RESULT	SPIKED SAMPLE	DUP. %		RPD
	RESULT	SPIKED			SAMPLE	REC.	
1,2,4-TRICHLOROBENZENE	<10	50	31	62	33	66	6
ACENAPHTHENE	<10	50	32	64	37	74	14
2,4-DINITROTOLUENE	<25	50	26	52	31	62	18
PYRENE	<5	50	52	104	55	110	6
N-NITROSO-DI-N-PROPYL AMINE	<10	50	26	52	29	58	11
1,4-DICHLOROBENZENE	<5	50	37	74	38	76	3
PENTACHLOROPHENOL	<50	100	61	61	65	65	6
PHENOL	<5	100	50	50	52	104	4
2-CHLOROPHENOL	<10	100	52	52	58	58	11
4-CHLORO-3-METHYLPHENOL	<10	100	47	47	47	47	0
4-NITROPHENOL	<50	100	40	40	41	41	2

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$



DATE 5-1-91 PAGE 1 OF 1

ATI Labs: San Diego (619) 458-9141 • Phoenix (602) 438-1530 • Seattle (206) 228-8335 • Pensacola (904) 474-1001
DISTRIBUTION: White, Canary - ANALYTICAL TECHNOLOGIES, INC.; Pink - ORIGINATOR

TABLE -1
BACKGROUND METALS

Total Metals

<u>Parameter</u>	<u>Analytical Method</u>	<u>Reporting Limit mg/kg</u>
✓Antimony	6010	6.0
✓Arsenic	7060	0.5
✓Barium	6010	1.0
✓Beryllium	6010	0.2
✓Cadmium	6010	0.5
✓Chromium	6010	1.0
✓Cobalt	6010	1.0
✓Copper	6010	2.0
✓Lead	6010	5.0
✓Mercury	7471	0.2
✓Nickel	6010	4.0
✓Potassium	6010	500
✓Selenium	7740	0.5
✓Vanadium	6010	1.0
✓Zinc	6010	2.0

TABLE-4
SKINNER LIST

METHOD 8240

<u>Parameter</u>	<u>Reporting Limit (ug/kg)</u>
Benzene	500
Carbon disulfide	500
Chlorobenzene	500
2-Chloroethylvinyl ether	1,000
1,2-Dibromomethane	1,000
1,2-Dichloroethane	500
1,4-Dioxane	50,000
Ethyl Benzene	500
Methyl ethyl ketone (2-butanone)	1,000
Styrene	500
Toluene	500
Xylenes	500

METHOD 8270

Anthracene	5,000
Benzenethiol	-
Benzo(a)anthracene	5,000
Benzo(b)fluoranthene	5,000
Benzo(k)fluoranthene	5,000
Benzo(a)pyrene	5,000
Bis(2-ethylhexyl)phthalate	5,000
Butyl benzyl phthalate	5,000
Chrysene	5,000
Dibenzo(a,X)acridine	-
Dibenzo(a,h)anthracene	5,000
Di-n-butylphthalate	5,000
1,2-Dichlorobenzene	5,000
1,3-Dichlorobenzene	5,000
1,4-Dichlorobenzene	5,000
Diethyl phthalate	5,000
7,12-Dimethylbenz(a)anthracene	5,000
2,4-Dimethylphenol	5,000
Dimethyl phthalate	5,000
2,4-Dinitrophenol	25,000
Di-n-octyl phthalate	5,000
Fluoranthene	5,000
Indene	5,000
Methylchrysene	-
1-Methylnaphthalene	5,000
2-Methylphenol	5,000
3-Methylphenol	5,000
4-Methylphenol	5,000
Naphthalene	5,000
4-Nitrophenol	25,000
Phenanthrene	5,000

TABLE-4 Continued

Phenol.	5,000
Pyrene	5,000
Pyridine	10,000
Quinoline	25,000



Analytical**Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 105576

June 10, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: OCD

Attention: Claud Rosendale

On 05/04/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

8240 analyses were performed by ATI, Fort Collins. The detection limit they are able to report for EDB is 10 $\mu\text{g/l}$, four times higher than that reportable by ATI, Phoenix

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : OCD

DATE RECEIVED : 05/04/91

REPORT DATE : 05/29/91

ATI I.D. : 105576

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	OW-12	AQUEOUS	05/02/91
02	OW-13	AQUEOUS	05/02/91
03	OW-14	AQUEOUS	05/02/91
04	OW-20	AQUEOUS	05/02/91
05	TRIP BLANK	AQUEOUS	03/25/91

----- TOTALS -----

MATRIX	# SAMPLES
AQUEOUS	5

ATI STANDARD DISPOSAL PRACTICE

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



GCMS - RESULTS

ATI I.D. : 10557601

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/02/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/04/91
PROJECT NAME	: OCD	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-12	DATE ANALYZED	: 05/17/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS	RESULTS
-----------	---------

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<10

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	110
BROMOFLUOROBENZENE (%)	101
TOLUENE-D8 (%)	94



Analytical Technologies, Inc.

GCMS - RESULTS

ATI I.D. : 10557602

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/02/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/04/91
PROJECT NAME	: OCD	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-13	DATE ANALYZED	: 05/17/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<10

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	108
BROMOFLUOROBENZENE (%)	100
TOLUENE-D8 (%)	103



Analytical Technologies, Inc.

GCMS - RESULTS

ATI I.D. : 10557603

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/02/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/04/91
PROJECT NAME	: OCD	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-14	DATE ANALYZED	: 05/17/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<10

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	114
BROMOFLUOROBENZENE (%)	99
TOLUENE-D8 (%)	101



Analytical Technologies, Inc.

GCMS - RESULTS

ATI I.D. : 10557604

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 05/02/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/04/91
PROJECT NAME	: OCD	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-20	DATE ANALYZED	: 05/16/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	490
BENZENE	130
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	170
CHLOROBENZENE	<5
ETHYLBENZENE	5
STYRENE	<5
TOTAL XYLENES	29
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<10

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	106
BROMOFLUOROBENZENE (%)	101
TOLUENE-D8 (%)	103



GCMS - RESULTS

ATI I.D. : 10557605

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 03/25/91
PROJECT #	: (NONE)	DATE RECEIVED	: 05/04/91
PROJECT NAME	: OCD	DATE EXTRACTED	: N/A
CLIENT I.D.	: TRIP BLANK	DATE ANALYZED	: 05/16/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYL VINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<10

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	107
BROMOFLUOROBENZENE (%)	101
TOLUENE-D8 (%)	96



GCMS - RESULTS

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	ATI I.D. :	105576
PROJECT #	: (NONE)	DATE EXTRACTED	: 05/16/91
PROJECT NAME	: OCD	DATE ANALYZED	: 05/16/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYL VINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<10

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	100
BROMOFLUOROBENZENE (%)	96
TOLUENE-D8 (%)	110



Analytical Technologies, Inc.

GCMS - RESULTS

TEST : VOLATILE ORGANICS (SKINNER LIST) EPA 8240

CLIENT	: GIANT REFINING CO.	ATI I.D. : 105576
PROJECT #	: (NONE)	DATE EXTRACTED : 05/17/91
PROJECT NAME	: OCD	DATE ANALYZED : 05/17/91
CLIENT I.D.	: REAGENT BLANK	UNITS : UG/L
		DILUTION FACTOR : N/A

COMPOUNDS	RESULTS
-----------	---------

CARBON DISULFIDE	<5
1,2-DICHLOROETHANE	<5
2-BUTANONE (MEK)	<5
BENZENE	<5
2-CHLOROETHYLVINYLETHER	<5
TOLUENE	<5
CHLOROBENZENE	<5
ETHYLBENZENE	<5
STYRENE	<5
TOTAL XYLENES	<5
1,4-DIOXANE	<10
1,2-DIBROMOETHANE (EDB)	<10

SURROGATE PERCENT RECOVERIES

1,2-DICHLOROETHANE-D4 (%)	108
BROMOFLUOROBENZENE (%)	106
TOLUENE-D8 (%)	102



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 105576

TEST : VOLATILE ORGANICS (EPA 8240)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : OCD

REF I.D. : 10557601

DATE ANALYZED : 05/17/91

SAMPLE MATRIX : AQUEOUS

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	DUP.		DUP.	RPD
	RESULT	SPIKED		% SPIKED SAMPLE REC.	% SAMPLE REC.		
1,1-DICHLOROETHENE	18	50	70	104	75	114	7
TRICHLOROETHENE	<5	50	57	114	51	102	11
CHLOROBENZENE	<5	50	51	102	48	96	6
TOLUENE	<5	50	53	106	49	98	8
BENZENE	<5	50	49	98	47	94	4

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

Chain of Custody

DATE 5-3-91 PAGE 1 OF 1

PROJECT MANAGER: <u>Bill Proffitt</u>				ANALYSIS REQUEST			
COMPANY: <u>Great Air Fining Co</u>							
ADDRESS: <u>873 Box 7</u>							
<u>Gallego, NM 87301</u>							
BILL TO:							
COMPANY: <u>Samuel</u>							
ADDRESS:							
SAMPLES: (Signature) <u>Michael K. Kunkel</u> PHONE NUMBER <u>(505) 222-0217</u>							
SAMPLE ID	DATE	TIME	MATRIX	LAB ID			
<u>OW-12</u>	<u>5-2</u>	<u>1:35p</u>	<u>Water</u>	<u>1</u>			
<u>OW-13</u>	<u>5-2</u>	<u>2:25p</u>	<u>Water</u>	<u>2</u>			
<u>OW-14</u>	<u>5-2</u>	<u>3:55p</u>	<u>Water</u>	<u>3</u>			
<u>OW-20</u>	<u>5-2</u>	<u>3:30</u>	<u>Water</u>	<u>4</u>			
<u>Y-04101</u>	<u>3-25</u>		<u>Water</u>	<u>5</u>			
PROJECT INFORMATION					SAMPLE RECEIPT		
PROJECT NO: <u>—</u>	TOTAL NO. OF CONTAINERS			9			
PROJECT NAME: <u>OCRD</u>	CHAIN OF CUSTODY SEALS			N			
P.O. NO. <u>01413</u>	INACT? <u>Y</u>						
SHIPPED VIA: <u>Fed Ex</u>	RECEIVED GOOD COND./COLD			Y			
SAMPLE DISPOSAL INSTRUCTIONS	LAB NUMBER			<u>105576</u>			
☒ ATT ☐ RETURN							
PRIOR AUTHORIZATION IS REQUIRED FOR RUSH PROJECTS							
TAT: (NORMAL) ☒ (RUSH) ☐ 24 ☐ 48 ☐ 72 ☐ 1 WEEK							
Comments: <u>see attached for analytical</u>							
<u>Regulation 15</u>							
RELINQUISHED BY: 1.				RELINQUISHED BY: 2.			
Signature: <u>Michael K. Kunkel</u> Time: <u>0900</u>				Signature: _____ Time: _____			
Printed Name: <u>Michael K. Kunkel</u> Date: <u>5-3-91</u>				Printed Name: _____ Date: _____			
Company: <u>Great Air Fining Co</u>				Company: _____			
RECEIVED BY: 1. Signature: _____ Time: _____				RECEIVED BY: 2. Signature: _____ Time: _____			
Printed Name: _____ Date: _____				Printed Name: _____ Date: _____			
Company: _____				Company: _____			
RECEIVED BY: 3. Signature: _____ Time: _____				RECEIVED BY: 4. Signature: _____ Time: _____			
Printed Name: <u>Wanda Eschelmann</u> Date: <u>5/4/91</u>				Printed Name: _____ Date: _____			
Company: <u>Analytical Technologies, Inc.</u>				Company: _____			
Petroleum Hydrocarbons (418.1)							
(MOD 8015) Gas/Diesel							
Diesel/Gasoline/BTXE (MOD 8015/8020)							
BTXE (8020)							
Chlorinated Hydrocarbons (601/8010)							
Aromatic Hydrocarbons (602/8020)							
MTBE							
Pesticides/PCB (608/8080)							
Herbicides (615/8150)							
Base/Neutral/Acid Compounds GC/MS (625/8270)							
Volatile Organics GC/MS (625/8240)					<u>see attached</u>		
SDWA Primary Standards							
SDWA Secondary Standards							
SDWA Volatiles (502.1/503.1)							
The 13 Priority Pollutant Metals							
The 8 EP Tox Metals by EP Tox Prep. (1310)							
The 8 EP Tox Metals by Total Digestion							
The 8 EP Tox Metals by TCLP (1311)							
NUMBER OF CONTAINERS							

TABLE-4
SKINNER LIST

METHOD 8240

<u>Parameter</u>	<u>Reporting Limit (ug/kg)</u>
Benzene	500
Carbon disulfide	500
Chlorobenzene	500
2-Chloroethylvinyl ether	1,000
1,2-Dibromomethane	1,000
1,2-Dichloroethane	500
1,4-Dioxane	50,000
Ethyl Benzene	500
Methyl ethyl ketone (2-butanone)	1,000
Styrene	500
Toluene	500
Xylenes	500

METHOD 8270

Anthracene	5,000
Benzenethiol	-
Benzo(a)anthracene	5,000
Benzo(b)fluoranthene	5,000
Benzo(k)fluoranthene	5,000
Benzo(a)pyrene	5,000
Bis(2-ethylhexyl)phthalate	5,000
Butyl benzyl phthalate	5,000
Chrysene	5,000
Dibenzo(a,h)acridine	-
Dibenzo(a,h)anthracene	5,000
Di-n-butylphthalate	5,000
1,2-Dichlorobenzene	5,000
1,3-Dichlorobenzene	5,000
1,4-Dichlorobenzene	5,000
Diethyl phthalate	5,000
7,12-Dimethylbenz(a)anthracene	5,000
2,4-Dimethylphenol	5,000
Dimethyl phthalate	5,000
2,4-Dinitrophenol	25,000
Di-n-octyl phthalate	5,000
Fluoranthene	5,000
Indene	5,000
Methylchrysene	-
1-Methylnaphthalene	5,000
2-Methylphenol	5,000
3-Methylphenol	5,000
4-Methylphenol	5,000
Naphthalene	5,000
4-Nitrophenol	25,000
Phenanthrene	5,000



Analytical**Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 104901

May 17, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: OCD Plan 997-9008-28

Attention: Claud Rosendale

On 04/30/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

DATE RECEIVED : 04/30/91

REPORT DATE : 05/16/91

ATI I.D. : 104901

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	OW-26	AQUEOUS	04/29/91

----- TOTALS -----

MATRIX	# SAMPLES
-----	-----
AQUEOUS	1

ATI STANDARD DISPOSAL PRACTICE

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



Analytical Technologies, Inc.

GENERAL CHEMISTRY - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

ATI I.D. : 104901

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
CARBONATE	MG/L	10491101	<1	<1	NA	NA	NA	NA
BICARBONATE	MG/L		126	125	1	NA	NA	NA
HYDROXIDE	MG/L		<1	<1	NA	NA	NA	NA
TOTAL ALKALINITY	MG/L		126	125	1	NA	NA	NA
CHLORIDE	MG/L	10555503	390	390	0	1150	750	101
CONDUCTIVITY (UMHOS/CM)		10490201	1190	1240	4	NA	NA	NA
PH	UNITS	10563801	8.2	8.3	1	NA	NA	NA
PHENOLICS, TOTAL	MG/L	10550401	<0.02	<0.02	NA	0.25	0.25	100
SULFATE	MG/L	10488717	130	140	7	280	140	100
TOTAL DISSOLVED SOLIDS	MG/L	10490101	1100	1100	0	NA	NA	NA

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

METALS RESULTS

ATI I.D. : 104901

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 04/30/91

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

REPORT DATE : 05/16/91

PARAMETER	UNITS	01
SILVER	MG/L	<0.010
ARSENIC	MG/L	0.008
BARIUM	MG/L	4.16
CALCIUM	MG/L	89.4
CADMIUM	MG/L	0.008
CHROMIUM	MG/L	<0.010
MANGANESE	MG/L	2.26
SODIUM	MG/L	289
LEAD	MG/L	0.009
SELENIUM	MG/L	<0.005



Analytical Technologies, Inc.

METALS - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

ATI I.D. : 104901

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
SILVER	MG/L	10489603	<0.010	<0.010	NA	0.440	0.500	88
ARSENIC	MG/L	10490101	0.008	0.008	0	0.042	0.050	68
BARIUM	MG/L	10489603	0.041	0.040	2	0.962	1.00	92
CALCIUM	MG/L	10491401	30.6	30.7	0.3	83.9	50.0	107
CADMIUM	MG/L	10489603	<0.005	0.007	NA	0.482	0.500	96
CHROMIUM	MG/L	10489603	<0.010	<0.010	NA	0.876	1.00	88
MANGANESE	MG/L	10489603	2.32	2.33	0.4	3.26	1.00	94
SODIUM	MG/L	10491401	34.1	35.5	4	85.4	50.0	103
LEAD	MG/L	10490101	0.009	0.009	0	0.057	0.050	96
SELENIUM	MG/L	10490101	<0.005	<0.005	NA	0.030	0.050	60

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

GAS CHROMATOGRAPHY - RESULTS

ATI I.D. : 10490101

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : OW-26
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : N/A
DATE ANALYZED : 05/09/91
UNITS : UG/L
DILUTION FACTOR : 500

COMPOUNDS

RESULTS

BENZENE	5500
BROMODICHLOROMETHANE	<100
BROMOFORM	<100
BROMOMETHANE	<100
CARBON TETRACHLORIDE	<100
CHLOROBENZENE	<250
CHLOROETHANE	<100
CHLOROFORM	<100
CHLOROMETHANE	<100
DIBROMOCHLOROMETHANE	<100
1,3-DICHLOROBENZENE	<250
2-CHLOROETHYL VINYL ETHER	<250
1,2 & 1,4-DICHLOROBENZENE	<250
DICHLORODIFLUOROMETHANE	<100
1,1-DICHLOROETHANE	<100
1,2-DICHLOROETHANE	<100
1,1-DICHLOROETHENE	<100
1,2-DICHLOROETHENE (TOTAL)	<100
1,2-DICHLOROPROPANE	<100
CIS-1,3-DICHLOROPROPENE	<100
TRANS-1,3-DICHLOROPROPENE	<100
ETHYLBENZENE	<250
METHYLENE CHLORIDE	<1000
1,1,2,2-TETRACHLOROETHANE	<100
TETRACHLOROETHENE	<100
TOLUENE	3000
1,1,1-TRICHLOROETHANE	<100
1,1,2-TRICHLOROETHANE	<100
TRICHLOROETHENE	<100
TRICHLOROFLUOROMETHANE	<250
VINYL CHLORIDE	<100
TOTAL XYLENES	2050
TRICHLOROFLUOROMETHANE	<100

SURROGATE PERCENT RECOVERIES

BROMOCHLOROMETHANE (%)	110
BROMOFLUOROBENZENE (%)	88



Analytical Technologies, Inc.

GAS CHROMATOGRAPHY - RESULTS

REAGENT BLANK

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 104901
PROJECT #	: 997-9008-28	DATE EXTRACTED	: 05/09/91
PROJECT NAME	: OCD PLAN	DATE ANALYZED	: 05/09/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

BENZENE	<0.5
BROMODICHLOROMETHANE	<0.2
BROMOFORM	<0.2
BROMOMETHANE	<0.2
CARBON TETRACHLORIDE	<0.2
CHLOROBENZENE	<0.5
CHLOROETHANE	<0.2
CHLOROFORM	<0.2
CHLOROMETHANE	<0.2
DIBROMOCHLOROMETHANE	<0.2
1,3-DICHLOROBENZENE	<0.5
2-CHLOROETHYL VINYL ETHER	<0.5
1,2 & 1,4-DICHLOROBENZENE	<0.5
DICHLORODIFLUOROMETHANE	<0.2
1,1-DICHLOROETHANE	<0.2
1,2-DICHLOROETHANE	<0.2
1,1-DICHLOROETHENE	<0.2
1,2-DICHLOROETHENE (TOTAL)	<0.2
1,2-DICHLOROPROPANE	<0.2
CIS-1,3-DICHLOROPROPENE	<0.2
TRANS-1,3-DICHLOROPROPENE	<0.2
ETHYLBENZENE	<0.5
METHYLENE CHLORIDE	<2.0
1,1,2,2-TETRACHLOROETHANE	<0.2
TETRACHLOROETHENE	<0.2
TOLUENE	<0.5
1,1,1-TRICHLOROETHANE	<0.2
1,1,2-TRICHLOROETHANE	<0.2
TRICHLOROETHENE	<0.2
TRICHLOROFLUOROMETHANE	<0.5
VINYL CHLORIDE	<0.2
TOTAL XYLENES	<0.5
TRICHLOROFLUOROMETHANE	<0.2

SURROGATE PERCENT RECOVERIES

BROMOCHLOROMETHANE (%)	106
BROMOFLUOROBENZENE (%)	98



Analytical Technologies, Inc.

QUALITY CONTROL DATA

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020) ATI I.D. : 104901

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
REF I.D. : 10489901

DATE ANALYZED : 05/10/91
SAMPLE MATRIX : AQUEOUS
UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	% REC.	DUP. SPIKED		RPD
	RESULT	SPIKED			SAMPLE	% REC.	
1,1-DICHLOROETHENE	<0.2	20	20	100	21	105	5
TRICHLOROETHENE	<0.2	20	19	95	20	100	5
TETRACHLOROETHENE	<0.2	20	18	90	19	95	5
BENZENE	<0.5	20	18	90	19	95	5
BROMODICHLOROMETHANE	<0.2	20	19	95	19	95	0
CHLOROFORM	<0.2	20	19	95	19	95	0
1,1,1-TRICHLOROETHANE	<0.2	20	19	95	19	95	0
TOLUENE	<0.5	20	19	95	18	90	5
CHLOROBENZENE	<0.5	20	18	90	18	90	0
M-XYLENE	<0.5	20	18	90	18	90	0

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

GCMS - RESULTS

ATI I.D. : 10490101

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : OW-26
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/11/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS	RESULTS
-----------	---------

ACETOPHENONE	<10
2-ACETYLAMINOFLUORENE	<10
a,a-DIMETHYLPHENETHYLAMINE	<10
4-AMINOBIIPHENYL	<10
ARAMITE	<50
CHLOROBENZILATE	<10
DIALATE	<10
2,6-DICHLOROPHENOL	<10
DIMETHOATE	<10
p-(DIMETHYLAMINO)AZOBENZENE	<10
7,12-DIMETHYLBENZO(a)ANTHRACENE	<10
3,3'-DIMETHYLBENZIDINE	<10
m-DINITROBENZENE	<10
DIPHENYLAMINE	<10
ETHYL METHANESULFONATE	<10
HEXACHLOROPHENE	<10
HEXACHLOROPROPENE	<10
ISOSAFROLE	<10
METHAPYRILENE	<10
3-METHYLCHOLANTHRENE	<10
METHYL METHANESULFONATE	<10
3-METHYLPHENOL (m-CRESOL)	<10
1,4-NAPHTHOQUINONE	<10
1-NAPHTHYLAMINE	<10
2-NAPHTHYLAMINE	<10
5-NITRO-O-TOLUIDINE	<10
4-NITROQUINOLINE-1-OXIDE	<10
N-NITROSODIETHYLAMINE	<10
N-NITROSODI-BUTYLAMINE	<10
N-NITROSOMETHYLETHYLAMINE	<200
N-NITROSOMORPHOLINE	<10
N-NITROSOPIPERIDINE	<10
N-NITROSOPYRROLIDINE	<10
PENTACHLOROBENZENE	<10
PENTACHLORONITROBENZENE	<10
PENTACHLOROETHANE	<10
PHENACETIN	<10
p-PHENYLENEDIAMINE	<10
2-PICOLINE	<10
PRONAMIDE	<10
SAFROLE	<10

(CONTINUED NEXT PAGE)



GCMS - RESULTS

ATI I.D. : 10490101

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
DIETHYLPHTHALATE	<10
4-CHLOROPHENYL-PHENYLETHER	<10
FLUORENE	<10
4-NITROANILINE	<50
4,6-DINITRO-2-METHYLPHENOL	<50
N-NITROSODIPHENYLAMINE	<10
4-BROMOPHENYL-PHENYLETHER	<10
HEXACHLOROBENZENE	<10
PENTACHLOROPHENOL	<50
PHENANTHRENE	<10
ANTHRACENE	<10
DI-N-BUTYLPHTHALATE	<10
FLUORANTHENE	<10
BENZIDINE	<100
PYRENE	<10
BUTYLBENZYLPHTHALATE	<10
3,3'-DICHLOROBENZIDINE	<20
BENZO(a)ANTHRACENE	<10
BIS(2-ETHYLHEXYL)PHTHALATE	<10
CHRYSENE	<10
DI-N-OCTYLPHTHALATE	<10
BENZO(b)FLUORANTHENE	<10
BENZO(k)FLUORANTHENE	<10
BENZO(a)PYRENE	<10
INDENO(1,2,3-cd)PYRENE	<10
DIBENZO(a,h)ANTHRACENE	<10
BENZO(g,h,i)PERYLENE	<10

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	110
2-FLUOROBIPHENYL (%)	59
TERPHENYL (%)	90
PHENOL-D6 (%)	12
2-FLUOROPHENOL (%)	24
2,4,6-TRIBROMOPHENOL (%)	39



Analytical **Technologies**, Inc.

GCMS - RESULTS

ATI I.D. : 10490101

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
1,2,4,5-TETRACHLOROBENZENE	<10
2,3,4,6-TETRACHLOROPHENOL	<10
o-TOLUIDINE	<10
1,3,5-TRINITROBENZENE	<10
PYRIDINE	<10



GCMS - RESULTS

ATI I.D. : 10490101

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : OW-26
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/11/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS	RESULTS
-----------	---------

N-NITROSODIMETHYLAMINE	<10
PHENOL	<10
ANILINE	<10
BIS(2-CHLOROETHYL) ETHER	<10
2-CHLOROPHENOL	<10
1,3-DICHLOROBENZENE	<10
1,4-DICHLOROBENZENE	<10
BENZYL ALCOHOL	<10
1,2-DICHLOROBENZENE	<10
2-METHYLPHENOL	18
BIS(2-CHLOROISOPROPYL) ETHER	<10
4-METHYLPHENOL	<10
N-NITROSO-DI-N-PROPYLAMINE	<10
HEXACHLOROETHANE	<10
NITROBENZENE	<10
ISOPHORONE	<10
2-NITROPHENOL	<10
2,4-DIMETHYLPHENOL	<10
BENZOIC ACID	54
BIS(2-CHLOROETHOXY) METHANE	<10
2,4-DICHLOROPHENOL	<10
1,2,4-TRICHLOROBENZENE	<10
NAPHTHALENE	TR
4-CHLOROANILINE	<10
HEXACHLOROBUTADIENE	<10
4-CHLORO-3-METHYLPHENOL	<10
2-METHYLNAPHTHALENE	208
HEXACHLOROCYCLOPENTADIENE	<10
2,4,6-TRICHLOROPHENOL	<10
2,4,5-TRICHLOROPHENOL	<50
2-CHLORONAPHTHALENE	<10
2-NITROANILINE	<50
DIMETHYLPHTHALATE	<10
ACENAPHTHYLENE	<10
3-NITROANINLINE	<50
ACENAPHTHENE	<10
2,4-DINITROPHENOL	<50
4-NITROPHENOL	<50
DIBENZOFURAN	<10
2,4-DINITROTOLUENE	<10
2,6-DINITROTOLUENE	<10

TR - Compound detected at an unquantifiable trace level

(CONTINUED NEXT PAGE)



Analytical**Technologies**, Inc.

ADDITIONAL COMPOUNDS (SEMI-QUANTITATED)

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

ATI I.D. : 10490101

COMPOUNDS	RESULTS
SUBSTITUTED BENZENES	5000
SUBSTITUTED NAPHTHALENES	300
TOTAL ORGANIC ACIDS	10000



GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : REAGENT BLANK

ATI I.D. : 104901
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/10/91
UNITS : UG/L
DILUTION FACTOR : N/A

COMPOUNDS

RESULTS

ACETOPHENONE	<10
2-ACETYLAMINOFLUORENE	<10
a,a-DIMETHYLPHENETHYLAMINE	<10
4-AMINOBIIPHENYL	<10
ARAMITE	<50
CHLOROBENZILATE	<10
DIALATE	<10
2,6-DICHLOROPHENOL	<10
DIMETHOATE	<10
p-(DIMETHYLAMINO)AZOBENZENE	<10
7,12-DIMETHYLBENZO(a)ANTHRACENE	<10
3,3'-DIMETHYLBENZIDINE	<10
m-DINITROBENZENE	<10
DIPHENYLAMINE	<10
ETHYL METHANESULFONATE	<10
HEXACHLOROPHENE	<10
HEXACHLOROPROPENE	<10
ISOSAFROLE	<10
METHAPYRILENE	<10
3-METHYLCHOLANTHRENE	<10
METHYL METHANESULFONATE	<10
3-METHYLPHENOL (m-CRESOL)	<10
1,4-NAPHTHOQUINONE	<10
1-NAPHTHYLAMINE	<10
2-NAPHTHYLAMINE	<10
5-NITRO-O-TOLUIDINE	<10
4-NITROQUINOLINE-1-OXIDE	<10
N-NITROSODIETHYLAMINE	<10
N-NITROSODI-BUTYLAMINE	<10
N-NITROSOMETHYLETHYLAMINE	<200
N-NITROSOMORPHOLINE	<10
N-NITROSOPIPERIDINE	<10
N-NITROSOPYRROLIDINE	<10
PENTACHLOROBENZENE	<10
PENTACHLORONITROBENZENE	<10
PENTACHLOROETHANE	<10
PHENACETIN	<10
p-PHENYLENEDIAMINE	<10
2-PICOLINE	<10
PRONAMIDE	<10
SAFROLE	<10
1,2,4,5-TETRACHLOROBENZENE	<10
2,3,4,6-TETRACHLOROPHENOL	<10

(CONTINUED NEXT PAGE)



Analytical Technologies, Inc.

GCMS - RESULTS

REAGENT BLANK

ATI I.D. : 104901

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
FLUORENE	<10
4-NITROANILINE	<50
4,6-DINITRO-2-METHYLPHENOL	<50
N-NITROSODIPHENYLAMINE	<10
4-BROMOPHENYL-PHENYLEETHER	<10
HEXACHLOROBENZENE	<10
PENTACHLOROPHENOL	<50
PHENANTHRENE	<10
ANTHRACENE	<10
DI-N-BUTYLPHTHALATE	<10
FLUORANTHENE	<10
BENZIDINE	<100
PYRENE	<10
BUTYLBENZYLPHTHALATE	<10
3,3'-DICHLOROBENZIDINE	<20
BENZO(a)ANTHRACENE	<10
BIS(2-ETHYLHEXYL)PHTHALATE	<10
CHRYSENE	<10
DI-N-OCTYLPHTHALATE	<10
BENZO(b)FLUORANTHENE	<10
BENZO(k)FLUORANTHENE	<10
BENZO(a)PYRENE	<10
INDENO(1,2,3-cd)PYRENE	<10
DIBENZO(a,h)ANTHRACENE	<10
BENZO(g,h,i)PERYLENE	<10

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	41
2-FLUOROBIPHENYL (%)	49
TERPHENYL (%)	96
PHENOL-D6 (%)	57
2-FLUOROPHENOL (%)	71
2,4,6-TRIBROMOPHENOL (%)	40

GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 104901
PROJECT #	: 997-9008-28	DATE EXTRACTED	: 05/01/91
PROJECT NAME	: OCD PLAN	DATE ANALYZED	: 05/10/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

N-NITROSODIMETHYLAMINE	<10
PHENOL	<10
ANILINE	<10
BIS(2-CHLOROETHYL) ETHER	<10
2-CHLOROPHENOL	<10
1,3-DICHLOROBENZENE	<10
1,4-DICHLOROBENZENE	<10
BENZYL ALCOHOL	<10
1,2-DICHLOROBENZENE	<10
2-METHYLPHENOL	<10
BIS(2-CHLOROISOPROPYL) ETHER	<10
4-METHYLPHENOL	<10
N-NITROSO-DI-N-PROPYLAMINE	<10
HEXACHLOROETHANE	<10
NITROBENZENE	<10
ISOPHORONE	<10
2-NITROPHENOL	<10
2,4-DIMETHYLPHENOL	<10
BENZOIC ACID	<50
BIS(2-CHLOROETHOXY) METHANE	<10
2,4-DICHLOROPHENOL	<10
1,2,4-TRICHLOROBENZENE	<10
NAPHTHALENE	<10
4-CHLOROANILINE	<10
HEXACHLOROBUTADIENE	<10
4-CHLORO-3-METHYLPHENOL	<10
2-METHYLNAPHTHALENE	<10
HEXACHLOROCYCLOPENTADIENE	<10
2,4,6-TRICHLOROPHENOL	<10
2,4,5-TRICHLOROPHENOL	<50
2-CHLORONAPHTHALENE	<10
2-NITROANILINE	<50
DIMETHYLPHTHALATE	<10
ACENAPHTHYLENE	<10
3-NITROANINLINE	<50
ACENAPHTHENE	<10
2,4-DINITROPHENOL	<50
4-NITROPHENOL	<50
DIBENZOFURAN	<10
2,4-DINITROTOLUENE	<10
2,6-DINITROTOLUENE	<10
DIETHYLPHTHALATE	<10
4-CHLOROPHENYL-PHENYLETHER	<10

(CONTINUED NEXT PAGE)



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 104901

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
REF I.D. : 10599911

DATE ANALYZED : 05/09/91
SAMPLE MATRIX : AQUEOUS
UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	% REC.	DUP.	DUP.	RPD
	RESULT	SPIKED			SPIKED SAMPLE	% REC.	
1,2,4-TRICHLOROBENZENE	<10	50	37	74	34	68	8
ACENAPHTHENE	<10	50	36	72	41	82	13
2,4-DINITROTOLUENE	<10	50	31	62	34	68	9
PYRENE	<10	56	63	126	57	114	10
N-NITROSO-DI-N-PROPYL AMINE	<10	50	38	76	43	86	12
1,4-DICHLOROBENZENE	<10	50	42	84	41	82	2
PENTACHLOROPHENOL	<50	100	88	88	89	89	1
PHENOL	<10	100	77	77	79	79	3
2-CHLOROPHENOL	<10	100	84	84	82	82	2
4-CHLORO-3-METHYLPHENOL	<10	100	80	80	75	75	6
4-NITROPHENOL	<50	100	40	40	47	47	16

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$



Analytical Technologies, Inc.
Phoenix, Arizona

Chain of Custody

DATE 4-29-91 / PAGE 1 OF 1

PROJECT INFORMATION				SAMPLE RECEIPT				RELINQUISHED BY:		RELINQUISHED BY:		RELINQUISHED BY:	
PROJECT NO:	TOTAL NO. OF CONTAINERS	CHAIN OF CUSTODY SEALS	INTACT?	RECEIVED GOOD COND./COLD	LAB NUMBER	SIGNATURE	TIME	SIGNATURE	TIME	SIGNATURE	TIME	SIGNATURE	TIME
PROJECT NAME: C.D. Plan	4	4	4	4	104901	Signature	Time	Signature	Time	Signature	Time	Signature	Time
P.O. NO. 997-9005-25						Printed Name: C.D. Plan	Date: 4/29/91	Printed Name: C.D. Plan	Date: 4/29/91	Printed Name: C.D. Plan	Date: 4/29/91	Printed Name: C.D. Plan	Date: 4/29/91
SHIPPED VIA: Fed Ex						Company: C.D. Plan		Company: C.D. Plan		Company: C.D. Plan		Company: C.D. Plan	
SAMPLE DISPOSAL INSTRUCTIONS						RECEIVED BY: (LAB)	3	RECEIVED BY: (LAB)	3	RECEIVED BY: (LAB)	3	RECEIVED BY: (LAB)	3
☒ ATI ☐ RETURN						Signature: [Signature]	Time: [Time]	Signature: [Signature]	Time: [Time]	Signature: [Signature]	Time: [Time]	Signature: [Signature]	Time: [Time]
PRIOR AUTHORIZATION IS REQUIRED FOR RUSH PROJECTS													
TAT: (NORMAL) ☒ (RUSH) ☐ 24 ☐ 48 ☐ 72 ☐ 1 WEEK													
Comments: Subcontractors for special analysis													

ANALYSIS REQUEST														
PROJECT MANAGER: Bob Hartlett														
COMPANY: Great Refining Co.														
ADDRESS: Box 7301, Gallup, NM 87301														
BILL TO: [Signature]														
COMPANY: [Signature]														
ADDRESS: [Signature]														
SAMPLERS: (Signature) [Signature] (505) 722-0217														
PHONE NUMBER														
SAMPLE ID	DATE	TIME	MATRIX	LAB ID										
001-26	4/29/91	0830	Water	1										
					Petroleum Hydrocarbons (418.1)	X								
					(MOD 8015) Gas/Liquid	X								
					Diesel/Gasoline/BTXE (MOD 8015/8020)	X								
					BTXE (8020)	X								
					Chlorinated Hydrocarbons (801/8010)	X								
					Aromatic Hydrocarbons (802/8020)	X								
					MTBE	X								
					Pesticides/PCB (608/8080)	X								
					Herbicides (615/8150)	X								
					Base/Neutral/Acid Compounds GC/MS (625/8270)	X								
					Volatile Organics GC/MS (624/8240)	X								
					SDWA Primary Standards	X								
					SDWA Secondary Standards	X								
					SDWA Volatiles (502.1/503.1)	X								
					The 13 Priority Pollutant Metals	X								
					The 8 EP Tox Metals by EP Tox Prep. (1310)	X								
					The 8 EP Tox Metals by Total Digestion	X								
					The 8 EP Tox Metals by TCLP (1311)	X								
					NUMBER OF CONTAINERS	4								

8
12 miles
for pens
Baker

Table 5

GIANT REFINING GALLUP, NEW MEXICO

General Inorganics

Parameter	Units	Reporting Limit
Alkalinity, Total as CaCO ₃ at pH 4.5	mg/L	5.0
Alkalinity, Bicarb. as CaCO ₃ at pH 4.5	mg/L	5.0
Alkalinity, Carb. as CaCO ₃ at pH 8.3	mg/L	5.0
Alkalinity, Hydrox. as CaCO ₃	mg/L	5.0
✓ Chloride	mg/L	3.0
✓ pH	units	--
✓ Phenolics	mg/L	0.010
✓ Sulfate	mg/L	5.0
✓ Specific Conductance at 25 deg.C	uans/c	1.0
✓ Total Dissolved Solids	mg/L	10.0

Table 6

GIANT REFINING GALLUP, NEW MEXICO

METALS

~~BASELINE~~ METALS

10771

Parameter

- ✓ Arsenic
- ✓ Barium
- ✓ Cadmium
- ✓ Calcium
- ✓ Chromium
- ✓ Lead
- ✓ Manganese
- ✓ Selenium
- ✓ Silver
- ✓ Sodium

Units Reporting
Limit

mg/L	0.0050
mg/L	0.010
mg/L	0.0050
mg/L	0.20 (?)
mg/L	0.010
mg/L	0.010
mg/L	0.010
mg/L	0.0050
mg/L	0.010
mg/L	5.0

Table 7

GIANT REFINING GALLUP, NEW MEXICO

AROMATIC VOLATILE ORGANICS

Parameter	Units	Reporting
		Limit
Benzene	ug/L	0.50
Toluene	ug/L	0.50
Chlorobenzene	ug/L	0.50
Ethyl benzene	ug/L	0.50
Total xylenes	ug/L	1.0
1,3-Dichlorobenzene	ug/L	0.50
1,4-Dichlorobenzene	ug/L	0.50
1,2-Dichlorobenzene	ug/L	0.50

Table 8

GIANT REFINING GALLUP, NEW MEXICO

Halogenated Volatile Organics

Parameter	Units	Reporting Limit
Chloromethane	ug/L	5.0
Bromomethane	ug/L	5.0
Vinyl chloride	ug/L	1.0
Chloroethane	ug/L	5.0
Methylene chloride	ug/L	0.50
1,1-Dichloroethene	ug/L	0.50
1,1-Dichloroethane	ug/L	1.0
1,2-Dichloroethane	ug/L	0.50
trans-1,2-Dichloroethene	ug/L	0.50
Chloroform		
1,1,2-Trichloro-1,2,2-trifluoroethane	ug/L	0.50
1,1,1-Trichloroethane	ug/L	0.50
Carbon tetrachloride	ug/L	1.0
Bromodichloroethane	ug/L	1.0
1,2-Dichloropropane	ug/L	5.0
Bromoform	ug/L	1.0
1,1,2,2-Tetrachloroethane	ug/L	0.50
Tetrachloroethene	ug/L	2.0
Chlorobenzene		

Table 9

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	nits	Reporting Limit
Acenaphthene	ug/L	10
Acenaphthylene	ug/L	10
Acetophenone	ug/L	10
2-Acetylaminofluorene	ug/L	10
4-Aminobiphenyl	ug/L	10
Aniline	ug/L	10
Anthracene	ug/L	10
Aramite	ug/L	10
Benzo(a)anthracene	ug/L	10
Benzo(b)fluoranthene	ug/L	10
Benzo(k)fluoranthene	ug/L	10
Benzo(g,h,i)perylene	ug/L	10
Benzo(a)pyrene	ug/L	10
Benzyl alcohol		
bis(2-Chloroethoxy)- methane	ug/L	10
bis(2-Chloroethyl)ether	ug/L	10
bis(2-Chloroisopropyl) ether	ug/L	10
bis(2-Ethylhexyl) phthalate	ug/L	10
4-Bromophenyl phenyl ether	ug/L	10
Butyl benzyl phthalate		
2sec-Butyl-4,6-dinitro- phenol (Dinoseb)	ug/L	10
4-Chloroaniline	ug/L	10
4-Chloro-3-methylphenol	ug/L	10
2-Chloronaphthalene	ug/L	10
2-Chlorophenol		
4-Chlorophenyl phenyl ether	ug/L	10
o-Cresol	ug/L	10
m & p-Cresol(s)	ug/L	10
Chrysene	ug/L	10
Dibenz(a,h)anthracene	ug/L	10
Dibenzofuran	ug/L	10
Di-n-butyl phthalate	ug/L	10
1,2-Dichlorobenzene	ug/L	10
1,3-Dichlorobenzene	ug/L	10
1,4-Dichlorobenzene	ug/L	10

Table 9 cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting Limit
3,3'-Dichlorobenzidine	ug/L	20
2,4-Dichlorophenol	ug/L	10
2,6-Dichlorophenol	ug/L	10
Diethyl phthalate	ug/L	10
Dimethoate	ug/L	10
p-Dimethylaminoazobenzene	ug/L	10
7,12-Dimethylbenzanthracene	ug/L	10
3,3'-Dimethylbenzidine	ug/L	10
a,a-Dimethylphenethylamine	ug/L	10
2,4-Dimethylphenol	ug/L	10
Dimethyl phthalate	ug/L	10
1,3-Dinitrobenzene	ug/L	10
4,6-Dinitro-2-methylphenol	ug/L	10
4,6-Dinitro-o-cresol	ug/L	50
2,4-Dinitrophenol	ug/L	50
2,4-Dinitrotoluene	ug/L	10
2,6-Dinitrotoluene	ug/L	10
Di-n-octyl phthalate	ug/L	10
Diphenylamine	ug/L	10
Disulfoton	ug/L	50
bis(2-Ethylhexyl) phthalate	ug/L	10
Ethyl methanesulfonate	ug/L	10
Famphur	ug/L	--
Flouranthene	ug/L	10
Flourene	ug/L	10
Hexachlorobenzene	ug/L	10
Hexachlorobutadiene	ug/L	10
Hexachlorocyclopentadiene	ug/L	10
Hexachloroethane		
Hexachlorophene	ug/L	--
Hexachloropropene	ug/L	10
Indeno(1,2,3-c,d)pyrene	ug/L	10
Isophorone	ug/L	10
Isosafrole	ug/L	20
Methapyrilene	ug/L	10
3-Methylcholanthrene	ug/L	10
Methyl methanesulfonate	ug/L	10
2-Methylnaphthalen	ug/L	10
Methyl parathion	ug/L	50
2-Methylphenol	ug/L	10
3/4-Methylphenol	ug/L	10
Methyl methacrylate	ug/L	10
Napthalene	ug/L	10

Table 9 Cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting
		Limit
1,4-Naphthaquinone	ug/L	10
1-Naphthylamine	ug/L	10
2-Naphthylamine	ug/L	10
2-Nitroaniline	ug/L	50
3-Nitroaniline	ug/L	50
4-Nitroaniline	ug/L	50
Nitrobenzene	ug/L	10
2-Nitrophenol	ug/L	10
4-Nitrophenol	ug/L	50
4-Nitroquinoline-1-oxide	ug/L	--
N-Nitroso-di-n-butylamine	ug/L	10
N-Nitrosodiethylamine	ug/L	10
N-Nitrosodimethylamine	ug/L	10
N-Nitrosodiphenylamine	ug/L	10
N-Nitroso-di-n-propylamine	ug/L	10
N-Nitrosomethylethylamine	ug/L	10
N-Nitrosomorpholine	ug/L	10
N-Nitrosopiperidine	ug/L	10
5-Nitro-o-toluidine	ug/L	10
N-Nitrosopyrrolidine	ug/L	10
Parathion	ug/L	50
Pentachlorobenzene	ug/L	10
Pentachlorethane	ug/L	10
Pentachloronitrobenzene	ug/L	50
Pentachlorophenol	ug/L	50
Phenacetin	ug/L	10
Phenanthrene	ug/L	10
Phenol	ug/L	10
4-Phenylenediamine	ug/L	--
Phorate	ug/L	100
2-Picoline	ug/L	10
Pronamide	ug/L	10
Pyrene	ug/L	10
Pyridine	ug/L	20
Safrole	ug/L	10
Sulfotapp	ug/L	50
1,2,4,5-Tetrachloro- benzene	ug/L	10
2,3,4,6-Tetrachlorophenol	ug/L	50
Thionazin	ug/L	50

2/12/96 cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting Limit
sym-Trinitrobenzene	ug/L	10
2-Toluidine	ug/L	10
1,2,4-Trichlorobenzene	ug/L	10
2,4,5-Trichlorophenol	ug/L	50
0,0,0-Triethylphosphoro- thioate	ug/L	10
2,4,6-Trichlorophenol	ug/L	10
1,3,5-Trinitrobenzene	ug/L	10
Ethyl methacrylate	ug/L	10



Analytical **Technologies, Inc.**

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 104899

May 17, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: OCD Plan 997-9008-28

Attention: Claud Rosendale

On 04/30/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

DATE RECEIVED : 04/30/91

REPORT DATE : 05/16/91

ATI I.D. : 104899

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	OW-16	AQUEOUS	04/29/91
02	TRIP BLANK	AQUEOUS	04/23/91

----- TOTALS -----

MATRIX	# SAMPLES
-----	-----
AQUEOUS	2

ATI STANDARD DISPOSAL PRACTICE

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



Analytical **Technologies**, Inc.

GENERAL CHEMISTRY RESULTS

ATI I.D. : 104899

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 04/30/91

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

REPORT DATE : 05/16/91

PARAMETER	UNITS	01
CARBONATE (CACO3)	MG/L	14
BICARBONATE (CACO3)	MG/L	282
HYDROXIDE (CACO3)	MG/L	<1
TOTAL ALKALINITY (AS CACO3)	MG/L	296
CHLORIDE	MG/L	170
CONDUCTIVITY, (UMHOS/CM)		1130
PH	UNITS	8.63
PHENOLICS, TOTAL	MG/L	<0.02
SULFATE	MG/L	32
TOTAL DISSOLVED SOLIDS	MG/L	710



Analytical Technologies, Inc.

GENERAL CHEMISTRY - QUALITY CONTROL

CLIENT : GIANT REFINING CO.

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

ATI I.D. : 104899

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
CARBONATE	MG/L	10491101	<1	<1	NA	NA	NA	NA
BICARBONATE	MG/L		126	125	1	NA	NA	NA
HYDROXIDE	MG/L		<1	<1	NA	NA	NA	NA
TOTAL ALKALINITY	MG/L		126	125	1	NA	NA	NA
CHLORIDE	MG/L	10555503	390	390	0	1150	750	101
CONDUCTIVITY(UMHOS/CM)		10488708	993	997	0.4	NA	NA	NA
PH	UNITS	10563801	8.2	8.3	1	NA	NA	NA
PHENOLICS, TOTAL	MG/L	10550401	<0.02	<0.02	NA	0.25	0.25	100
SULFATE	MG/L	10488717	130	140	7	280	140	100
TOTAL DISSOLVED SOLIDS	MG/L	10490101	1100	1100	0	NA	NA	NA

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical**Technologies**, Inc.

METALS RESULTS

ATI I.D. : 104899

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 04/30/91

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

REPORT DATE : 05/16/91

PARAMETER	UNITS	01
SILVER	MG/L	<0.010
ARSENIC	MG/L	<0.005
BARIUM	MG/L	0.043
CALCIUM	MG/L	5.4
CADMIUM	MG/L	0.008
CHROMIUM	MG/L	<0.010
MANGANESE	MG/L	0.022
SODIUM	MG/L	226
LEAD	MG/L	<0.002
SELENIUM	MG/L	0.030



Analytical Technologies, Inc.

METALS - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

ATI I.D. : 104899

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
SILVER	MG/L	10489603	<0.010	<0.010	NA	0.440	0.500	88
ARSENIC	MG/L	10551103	<0.005	<0.005	NA	0.038	0.050	76
BARIUM	MG/L	10488708	0.036	0.035	3	0.135	0.100	99
CALCIUM	MG/L	10491401	30.6	30.7	0.3	83.9	50.0	107
CADMIUM	MG/L	10489603	<0.005	0.007	NA	0.482	0.500	96
CHROMIUM	MG/L	10489603	<0.010	<0.010	NA	0.876	1.00	88
MANGANESE	MG/L	10488708	<0.010	<0.010	NA	0.097	0.100	97
SODIUM	MG/L	10491401	34.1	35.5	4	85.4	50.0	103
LEAD	MG/L	10489601	<0.002	<0.002	NA	0.046	0.050	92
SELENIUM	MG/L	10488707	<0.005	<0.005	NA	0.027	0.050	54

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical Technologies, Inc.

GAS CHROMATOGRAPHY - RESULTS

ATI I.D. : 10489901

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : OW-16
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : N/A
DATE ANALYZED : 05/09/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS RESULTS

BENZENE	<0.5
BROMODICHLOROMETHANE	<0.2
BROMOFORM	<0.2
BROMOMETHANE	<0.2
CARBON TETRACHLORIDE	<0.2
CHLOROBENZENE	<0.5
CHLOROETHANE	<0.2
CHLOROFORM	<0.2
CHLOROMETHANE	<0.2
DIBROMOCHLOROMETHANE	<0.2
1,3-DICHLOROBENZENE	<0.5
2-CHLOROETHYL VINYL ETHER	<0.5
1,2 & 1,4-DICHLOROBENZENE	<0.5
DICHLORODIFLUOROMETHANE	<0.2
1,1-DICHLOROETHANE	0.8
1,2-DICHLOROETHANE	5.7
1,1-DICHLOROETHENE	<0.2
1,2-DICHLOROETHENE (TOTAL)	<0.2
1,2-DICHLOROPROPANE	<0.2
CIS-1,3-DICHLOROPROPENE	<0.2
TRANS-1,3-DICHLOROPROPENE	<0.2
ETHYLBENZENE	<0.5
METHYLENE CHLORIDE	<2.0
1,1,2,2-TETRACHLOROETHANE	<0.2
TETRACHLOROETHENE	<0.2
TOLUENE	<0.5
1,1,1-TRICHLOROETHANE	<0.2
1,1,2-TRICHLOROETHANE	<0.2
TRICHLOROETHENE	<0.2
TRICHLOROFLUOROMETHANE	<0.5
VINYL CHLORIDE	<0.2
TOTAL XYLENES	<0.5
TRICHLOROFLUOROMETHANE	<0.2

SURROGATE PERCENT RECOVERIES

BROMOCHLOROMETHANE (%)	98
BROMOFLUOROBENZENE (%)	99

GAS CHROMATOGRAPHY - RESULTS

ATI I.D. : 10489902

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : TRIP BLANK
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/23/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : N/A
DATE ANALYZED : 05/09/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS RESULTS

BENZENE	<0.5
BROMODICHLOROMETHANE	<0.2
BROMOFORM	<0.2
BROMOMETHANE	<0.2
CARBON TETRACHLORIDE	<0.2
CHLOROBENZENE	<0.5
CHLOROETHANE	<0.2
CHLOROFORM	<0.2
CHLOROMETHANE	<0.2
DIBROMOCHLOROMETHANE	<0.2
1,3-DICHLOROBENZENE	<0.5
2-CHLOROETHYL VINYL ETHER	<0.5
1,2 & 1,4-DICHLOROBENZENE	<0.5
DICHLORODIFLUOROMETHANE	<0.2
1,1-DICHLOROETHANE	<0.2
1,2-DICHLOROETHANE	<0.2
1,1-DICHLOROETHENE	<0.2
1,2-DICHLOROETHENE (TOTAL)	<0.2
1,2-DICHLOROPROPANE	<0.2
CIS-1,3-DICHLOROPROPENE	<0.2
TRANS-1,3-DICHLOROPROPENE	<0.2
ETHYLBENZENE	<0.5
METHYLENE CHLORIDE	<2.0
1,1,2,2-TETRACHLOROETHANE	<0.2
TETRACHLOROETHENE	<0.2
TOLUENE	<0.5
1,1,1-TRICHLOROETHANE	<0.2
1,1,2-TRICHLOROETHANE	<0.2
TRICHLOROETHENE	<0.2
TRICHLOROFLUOROMETHANE	<0.5
VINYL CHLORIDE	<0.2
TOTAL XYLENES	<0.5
TRICHLOROFLUOROMETHANE	<0.2

SURROGATE PERCENT RECOVERIES

BROMOCHLOROMETHANE (%)	118
BROMOFLUOROBENZENE (%)	99



Analytical Technologies, Inc.

GAS CHROMATOGRAPHY - RESULTS

REAGENT BLANK

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 104899
PROJECT #	: 997-9008-28	DATE EXTRACTED	: 05/09/91
PROJECT NAME	: OCD PLAN	DATE ANALYZED	: 05/09/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

BENZENE	<0.5
BROMODICHLOROMETHANE	<0.2
BROMOFORM	<0.2
BROMOMETHANE	<0.2
CARBON TETRACHLORIDE	<0.2
CHLOROBENZENE	<0.5
CHLOROETHANE	<0.2
CHLOROFORM	<0.2
CHLOROMETHANE	<0.2
DIBROMOCHLOROMETHANE	<0.2
1,3-DICHLOROBENZENE	<0.5
2-CHLOROETHYL VINYL ETHER	<0.5
1,2 & 1,4-DICHLOROBENZENE	<0.5
DICHLORODIFLUOROMETHANE	<0.2
1,1-DICHLOROETHANE	<0.2
1,2-DICHLOROETHANE	<0.2
1,1-DICHLOROETHENE	<0.2
1,2-DICHLOROETHENE (TOTAL)	<0.2
1,2-DICHLOROPROPANE	<0.2
CIS-1,3-DICHLOROPROPENE	<0.2
TRANS-1,3-DICHLOROPROPENE	<0.2
ETHYLBENZENE	<0.5
METHYLENE CHLORIDE	<2.0
1,1,2,2-TETRACHLOROETHANE	<0.2
TETRACHLOROETHENE	<0.2
TOLUENE	<0.5
1,1,1-TRICHLOROETHANE	<0.2
1,1,2-TRICHLOROETHANE	<0.2
TRICHLOROETHENE	<0.2
TRICHLOROFLUOROMETHANE	<0.5
VINYL CHLORIDE	<0.2
TOTAL XYLENES	<0.5
TRICHLOROFLUOROMETHANE	<0.2

SURROGATE PERCENT RECOVERIES

BROMOCHLOROMETHANE (%)	106
BROMOFLUOROBENZENE (%)	98



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 104899

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT : GIANT REFINING CO.

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

REF I.D. : 10489901

DATE ANALYZED : 05/10/91

SAMPLE MATRIX : AQUEOUS

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	% REC.	DUP. SPIKED %		RPD
	RESULT	SPIKED			SAMPLE	REC.	
1,1-DICHLOROETHENE	<0.2	20	20	100	21	105	5
TRICHLOROETHENE	<0.2	20	19	95	20	100	5
TETRACHLOROETHENE	<0.2	20	18	90	19	95	5
BENZENE	<0.5	20	18	90	19	95	5
BROMODICHLOROMETHANE	<0.2	20	19	95	19	95	0
CHLOROFORM	<0.2	20	19	95	19	95	0
1,1,1-TRICHLOROETHANE	<0.2	20	19	95	19	95	0
TOLUENE	<0.5	20	19	95	18	90	5
CHLOROBENZENE	<0.5	20	18	90	18	90	0
M-XYLENE	<0.5	20	18	90	18	90	0

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

GCMS - RESULTS

ATI I.D. : 10489901

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : OW-16
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/11/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS RESULTS

ACETOPHENONE	<10
2-ACETYLAMINOFLUORENE	<10
a,a-DIMETHYLPHENETHYLAMINE	<10
4-AMINOBIIPHENYL	<10
ARAMITE	<50
CHLOROBENZILATE	<10
DIALATE	<10
2,6-DICHLOROPHENOL	<10
DIMETHOATE	<10
p-(DIMETHYLAMINO)AZOBENZENE	<10
7,12-DIMETHYLBENZO(a)ANTHRACENE	<10
3,3'-DIMETHYLBENZIDINE	<10
m-DINITROBENZENE	<10
DIPHENYLAMINE	<10
ETHYL METHANESULFONATE	<10
HEXACHLOROPHENE	<10
HEXACHLOROPROPENE	<10
ISOSAFROLE	<10
METHAPYRILENE	<10
3-METHYLCHOLANTHRENE	<10
METHYL METHANESULFONATE	<10
3-METHYLPHENOL (m-CRESOL)	<10
1,4-NAPHTHOQUINONE	<10
1-NAPHTHYLAMINE	<10
2-NAPHTHYLAMINE	<10
5-NITRO-O-TOLUIDINE	<10
4-NITROQUINOLINE-1-OXIDE	<10
N-NITROSODIETHYLAMINE	<10
N-NITROSODI-BUTYLAMINE	<10
N-NITROSOMETHYLETHYLAMINE	<200
N-NITROSOMORPHOLINE	<10
N-NITROSOPIPERIDINE	<10
N-NITROSOPIRROLIDINE	<10
PENTACHLOROBENZENE	<10
PENTACHLORONITROBENZENE	<10
PENTACHLOROETHANE	<10
PHENACETIN	<10
p-PHENYLENEDIAMINE	<10
2-PICOLINE	<10
PRONAMIDE	<10
SAFROLE	<10

(CONTINUED NEXT PAGE)



GCMS - RESULTS

ATI I.D. : 10489901

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
DIETHYLPHTHALATE	<10
4-CHLOROPHENYL-PHENYLETHER	<10
FLUORENE	<10
4-NITROANILINE	<50
4,6-DINITRO-2-METHYLPHENOL	<50
N-NITROSODIPHENYLAMINE	<10
4-BROMOPHENYL-PHENYLETHER	<10
HEXACHLOROBENZENE	<10
PENTACHLOROPHENOL	<50
PHENANTHRENE	<10
ANTHRACENE	<10
DI-N-BUTYLPHTHALATE	<10
FLUORANTHENE	<10
BENZIDINE	<100
PYRENE	<10
BUTYLBENZYLPHTHALATE	<10
3,3'-DICHLOROBENZIDINE	<20
BENZO(a)ANTHRACENE	<10
BIS(2-ETHYLHEXYL)PHTHALATE	10
CHRYSENE	<10
DI-N-OCTYLPHTHALATE	<10
BENZO(b)FLUORANTHENE	<10
BENZO(k)FLUORANTHENE	<10
BENZO(a)PYRENE	<10
INDENO(1,2,3-cd)PYRENE	<10
DIBENZO(a,h)ANTHRACENE	<10
BENZO(g,h,i)PERYLENE	<10

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	47
2-FLUOROBIPHENYL (%)	45
TERPHENYL (%)	88
PHENOL-D6 (%)	39
2-FLUOROPHENOL (%)	53
2,4,6-TRIBROMOPHENOL (%)	44



Analytical **Technologies**, Inc.

GCMS - RESULTS

ATI I.D. : 10489901

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
1,2,4,5-TETRACHLOROBENZENE	<10
2,3,4,6-TETRACHLOROPHENOL	<10
o-TOLUIDINE	<10
1,3,5-TRINITROBENZENE	<10
PYRIDINE	<10

GCMS - RESULTS

ATI I.D. : 10489901

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : OW-16
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/11/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS

RESULTS

N-NITROSODIMETHYLAMINE	<10
PHENOL	<10
ANILINE	<10
BIS(2-CHLOROETHYL) ETHER	<10
2-CHLOROPHENOL	<10
1,3-DICHLOROBENZENE	<10
1,4-DICHLOROBENZENE	<10
BENZYL ALCOHOL	<10
1,2-DICHLOROBENZENE	<10
2-METHYLPHENOL	<10
BIS(2-CHLOROISOPROPYL) ETHER	<10
4-METHYLPHENOL	<10
N-NITROSO-DI-N-PROPYLAMINE	<10
HEXACHLOROETHANE	<10
NITROBENZENE	<10
ISOPHORONE	<10
2-NITROPHENOL	<10
2,4-DIMETHYLPHENOL	<10
BENZOIC ACID	<50
BIS(2-CHLOROETHOXY) METHANE	<10
2,4-DICHLOROPHENOL	<10
1,2,4-TRICHLOROBENZENE	<10
NAPHTHALENE	<10
4-CHLOROANILINE	<10
HEXACHLOROBUTADIENE	<10
4-CHLORO-3-METHYLPHENOL	<10
2-METHYLNAPHTHALENE	<10
HEXACHLOROCYCLOPENTADIENE	<10
2,4,6-TRICHLOROPHENOL	<10
2,4,5-TRICHLOROPHENOL	<50
2-CHLORONAPHTHALENE	<10
2-NITROANILINE	<50
DIMETHYLPHTHALATE	<10
ACENAPHTHYLENE	<10
3-NITROANINLINE	<50
ACENAPHTHENE	<10
2,4-DINITROPHENOL	<50
4-NITROPHENOL	<50
DIBENZOFURAN	<10
2,4-DINITROTOLUENE	<10
2,6-DINITROTOLUENE	<10

(CONTINUED NEXT PAGE)



GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : REAGENT BLANK

ATI I.D. : 104899
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/10/91
UNITS : UG/L
DILUTION FACTOR : N/A

COMPOUNDS	RESULTS
ACETOPHENONE	<10
2-ACETYLAMINOFLUORENE	<10
a,a-DIMETHYLPHENETHYLAMINE	<10
4-AMINOBIIPHENYL	<10
ARAMITE	<50
CHLOROBENZILATE	<10
DIALATE	<10
2,6-DICHLOROPHENOL	<10
DIMETHOATE	<10
p-(DIMETHYLAMINO)AZOBENZENE	<10
7,12-DIMETHYLBENZO(a)ANTHRACENE	<10
3,3'-DIMETHYLBENZIDINE	<10
m-DINITROBENZENE	<10
DIPHENYLAMINE	<10
ETHYL METHANESULFONATE	<10
HEXACHLOROPHENE	<10
HEXACHLOROPROPENE	<10
ISOSAFROLE	<10
METHAPYRILENE	<10
3-METHYLCHOLANTHRENE	<10
METHYL METHANESULFONATE	<10
3-METHYLPHENOL (m-CRESOL)	<10
1,4-NAPHTHOQUINONE	<10
1-NAPHTHYLAMINE	<10
2-NAPHTHYLAMINE	<10
5-NITRO-O-TOLUIDINE	<10
4-NITROQUINOLINE-1-OXIDE	<10
N-NITROSODIETHYLAMINE	<10
N-NITROSODI-BUTYLAMINE	<10
N-NITROSOMETHYLETHYLAMINE	<200
N-NITROSOMORPHOLINE	<10
N-NITROSOPIPERIDINE	<10
N-NITROSOPYRROLIDINE	<10
PENTACHLOROBENZENE	<10
PENTACHLORONITROBENZENE	<10
PENTACHLOROETHANE	<10
PHENACETIN	<10
p-PHENYLENEDIAMINE	<10
2-PICOLINE	<10
PRONAMIDE	<10
SAFROLE	<10
1,2,4,5-TETRACHLOROBENZENE	<10
2,3,4,6-TETRACHLOROPHENOL	<10

(CONTINUED NEXT PAGE)



Analytical Technologies, Inc.

GCMS - RESULTS

REAGENT BLANK

ATI I.D. : 104899

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
FLUORENE	<10
4-NITROANILINE	<50
4,6-DINITRO-2-METHYLPHENOL	<50
N-NITROSODIPHENYLAMINE	<10
4-BROMOPHENYL-PHENYLEETHER	<10
HEXACHLOROBENZENE	<10
PENTACHLOROPHENOL	<50
PHENANTHRENE	<10
ANTHRACENE	<10
DI-N-BUTYLPHTHALATE	<10
FLUORANTHENE	<10
BENZIDINE	<100
PYRENE	<10
BUTYLBENZYLPHTHALATE	<10
3,3'-DICHLOOROBENZIDINE	<20
BENZO(a)ANTHRACENE	<10
BIS(2-ETHYLHEXYL)PHTHALATE	<10
CHRYSENE	<10
DI-N-OCTYLPHTHALATE	<10
BENZO(b)FLUORANTHENE	<10
BENZO(k)FLUORANTHENE	<10
BENZO(a)PYRENE	<10
INDENO(1,2,3-cd)PYRENE	<10
DIBENZO(a,h)ANTHRACENE	<10
BENZO(g,h,i)PERYLENE	<10

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	41
2-FLUOROBIPHENYL (%)	49
TERPHENYL (%)	96
PHENOL-D6 (%)	57
2-FLUOROPHENOL (%)	71
2,4,6-TRIBROMOPHENOL (%)	40



GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : REAGENT BLANK

ATI I.D. : 104899
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/10/91
UNITS : UG/L
DILUTION FACTOR : N/A

COMPOUNDS

RESULTS

N-NITROSODIMETHYLAMINE	<10
PHENOL	<10
ANILINE	<10
BIS(2-CHLOROETHYL) ETHER	<10
2-CHLOROPHENOL	<10
1,3-DICHLOROBENZENE	<10
1,4-DICHLOROBENZENE	<10
BENZYL ALCOHOL	<10
1,2-DICHLOROBENZENE	<10
2-METHYLPHENOL	<10
BIS(2-CHLOROISOPROPYL) ETHER	<10
4-METHYLPHENOL	<10
N-NITROSO-DI-N-PROPYLAMINE	<10
HEXACHLOROETHANE	<10
NITROBENZENE	<10
ISOPHORONE	<10
2-NITROPHENOL	<10
2,4-DIMETHYLPHENOL	<10
BENZOIC ACID	<50
BIS(2-CHLOROETHOXY) METHANE	<10
2,4-DICHLOROPHENOL	<10
1,2,4-TRICHLOROBENZENE	<10
NAPHTHALENE	<10
4-CHLOROANILINE	<10
HEXACHLOROBUTADIENE	<10
4-CHLORO-3-METHYLPHENOL	<10
2-METHYLNAPHTHALENE	<10
HEXACHLOROCYCLOPENTADIENE	<10
2,4,6-TRICHLOROPHENOL	<10
2,4,5-TRICHLOROPHENOL	<50
2-CHLORONAPHTHALENE	<10
2-NITROANILINE	<50
DIMETHYLPHTHALATE	<10
ACENAPHTHYLENE	<10
3-NITROANINLINE	<50
ACENAPHTHENE	<10
2,4-DINITROPHENOL	<50
4-NITROPHENOL	<50
DIBENZOFURAN	<10
2,4-DINITROTOLUENE	<10
2,6-DINITROTOLUENE	<10
DIETHYLPHTHALATE	<10
4-CHLOROPHENYL-PHENYLETHER	<10

(CONTINUED NEXT PAGE)



Analytical **Technologies**, Inc.

GCMS - RESULTS

REAGENT BLANK

ATI I.D. : 104899

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS

RESULTS

o-TOLUIDINE

<10

1,3,5-TRINITROBENZENE

<10

PYRIDINE

<10



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 104899

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
REF I.D. : 10599911

DATE ANALYZED : 05/09/91
SAMPLE MATRIX : AQUEOUS
UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	% REC.	DUP.	DUP.	RPD
	RESULT	SPIKED			SPIKED SAMPLE	% REC.	
1,2,4-TRICHLOROBENZENE	<10	50	37	74	34	68	8
ACENAPHTHENE	<10	50	36	72	41	82	13
2,4-DINITROTOLUENE	<10	50	31	62	34	68	9
PYRENE	<10	56	63	126	57	114	10
N-NITROSO-DI-N-PROPYL AMINE	<10	50	38	76	43	86	12
1,4-DICHLOROBENZENE	<10	50	42	84	41	82	2
PENTACHLOROPHENOL	<50	100	88	88	89	89	1
PHENOL	<10	100	77	77	79	79	3
2-CHLOROPHENOL	<10	100	84	84	82	82	2
4-CHLORO-3-METHYLPHENOL	<10	100	80	80	75	75	6
4-NITROPHENOL	<50	100	40	40	47	47	16

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

Chain of Custody

DATE 4-29-91 PAGE 1 OF 1

PROJECT INFORMATION				SAMPLE RECEIPT				ANALYSIS REQUEST												RELINQUISHED BY: 1.				RELINQUISHED BY: 2.				RELINQUISHED BY: 3.							
PROJECT NO:	TOTAL NO. OF CONTAINERS	CHAIN OF CUSTODY SEALS	INTACT?	RECEIVED GOOD COND./COLD	LAB NUMBER	SAMPLE ID	DATE	TIME	MATRIX	LAB ID	Petroleum Hydrocarbons (418.1)	(MOD 8015) Gas/Diesel	Diesel/Gasoline/BTXE (MOD 8015/8020)	BTXE (8020)	Chlorinated Hydrocarbons (808010)	Aromatic Hydrocarbons (808020)	MTBE	Pesticides/PCB (608/8080)	Herbicides (615/8150)	Base/Neutral/Acid Compounds GC/MS (625/8270)	Volatile Organics GC/MS (624/8240)	SDWA Primary Standards	SDWA Secondary Standards	SDWA Volatiles (502.1/503.1)	The 13 Priority Pollutant Metals	The 8 EP Tox Metals by EP Tox Prep. (1310)	The 8 EP Tox Metals by Total Digestion	The 8 EP Tox Metals by TCLP (1311)	NUMBER OF CONTAINERS						
PROJECT MANAGER: Ruth Proffert COMPANY: Great Refining Co. ADDRESS: Rt 3 Box 7 Gallup NM 87301 BILL TO: [Signature] COMPANY: [Signature] ADDRESS: [Signature]											SAMPLERS (Signature): [Signature] (505) 722-0217 PHONE NUMBER											RELINQUISHED BY: 1. Signature: [Signature] Time: 11:00 Printed Name: [Signature] Date: 4-29-91 Company: Great Refining				RELINQUISHED BY: 2. Signature: [Signature] Time: [Signature] Printed Name: [Signature] Date: 4-29-91 Company: [Signature]				RELINQUISHED BY: 3. Signature: [Signature] Time: 10:55 Printed Name: Linda Estelmen Date: 4/30/91 Company: Analytical Technologies, Inc.					
PROJECT NO: — PROJECT NAME: DCD Plan P.O. NO.: 997-9008-28 SHIPPED VIA: Fed Ex SAMPLE DISPOSAL INSTRUCTIONS: [Signature] ☒ ATI ☐ RETURN						LAB NUMBER: 104899 PRIOR AUTHORIZATION IS REQUIRED FOR RUSH PROJECTS TAT: (NORMAL) ☒ (RUSH) ☐ 24 ☐ 48 ☐ 72 ☐ 1 WEEK Comments: See Attachments for Specific Analytical Parameters 1 bottle broke for 8270						[Grid for Analysis Request with handwritten 'X' marks in various columns]												RELINQUISHED BY: 1. Signature: [Signature] Time: [Signature] Printed Name: [Signature] Date: [Signature] Company: [Signature]				RELINQUISHED BY: 2. Signature: [Signature] Time: [Signature] Printed Name: [Signature] Date: [Signature] Company: [Signature]				RELINQUISHED BY: 3. Signature: [Signature] Time: [Signature] Printed Name: [Signature] Date: [Signature] Company: [Signature]			

Table 5

GIANT REFINING GALLUP, NEW MEXICO

General Inorganics

Parameter	Units	Reporting Limit
Alkalinity, Total as CaCO ₃ at pH 4.5	mg/L	5.0
Alkalinity, Bicarb. as CaCO ₃ at pH 4.5	mg/L	5.0
Alkalinity, Carb. as CaCO ₃ at pH 8.3	mg/L	5.0
Alkalinity, Hydrox. as CaCO ₃	mg/L	5.0
✓ Chloride	mg/L	3.0
✓ pH	units	--
✓ Phenolics	mg/L	0.010
✓ Sulfate	mg/L	5.0
✓ Specific Conductance at 25 deg.C	uamhos/c	1.0
✓ Total Dissolved Solids	mg/L	10.0

Table 6

GIANT REFINING GALLUP, NEW MEXICO

METALS

~~PERMITTED~~ METALS

Tv+71

Parameter	Units	Reporting Limit
✓ Arsenic	mg/L	0.0050
✓ Barium	mg/L	0.010
✓ Cadmium	mg/L	0.0050
✓ Calcium	mg/L	0.20 (?)
✓ Chromium	mg/L	0.010
✓ Lead	mg/L	0.010
✓ Manganese	mg/L	0.010
✓ Selenium	mg/L	0.0050
✓ Silver	mg/L	0.010
✓ Sodium	mg/L	5.0

Table 7

GIANT REFINING GALLUP, NEW MEXICO

AROMATIC VOLATILE ORGANICS

Parameter	Units	Reporting Limit
Benzene	ug/L	0.50
Toluene	ug/L	0.50
Chlorobenzene	ug/L	0.50
Ethyl benzene	ug/L	0.50
Total xylenes	ug/L	1.0
1,3-Dichlorobenzene	ug/L	0.50
1,4-Dichlorobenzene	ug/L	0.50
1,2-Dichlorobenzene	ug/L	0.50

Table 8

GIANT REFINING GALLUP, NEW MEXICO

Halogenated Volatile Organics

Parameter	Units	Reporting Limit
Chloromethane	ug/L	5.0
Bromomethane	ug/L	5.0
Vinyl chloride	ug/L	1.0
Chloroethane	ug/L	5.0
Methylene chloride	ug/L	5.0
1,1-Dichloroethene	ug/L	0.50
1,1-Dichloroethane	ug/L	0.50
1,2-Dichloroethane	ug/L	1.0
trans-1,2-Dichloroethene	ug/L	0.50
Chloroform	ug/L	0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ug/L	0.50
1,1,1-Trichloroethane	ug/L	0.50
Carbon tetrachloride	ug/L	1.0
Bromodichloromethane	ug/L	1.0
1,2-Dichloropropane	ug/L	5.0
Bromoform	ug/L	1.0
1,1,2,2-Tetrachloroethane	ug/L	0.50
Tetrachloroethene	ug/L	2.0
Chlorobenzene		

Table 9

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	nits	Reporting Limit
Acenaphthene	ug/L	10
Acenaphthylene	ug/L	10
Acetophenone	ug/L	10
2-Acetylaminofluorene	ug/L	10
4-Aminobiphenyl	ug/L	10
Aniline	ug/L	10
Anthracene	ug/L	10
Aramite	ug/L	10
Benzo(a)anthracene	ug/L	10
Benzo(b)fluoranthene	ug/L	10
Benzo(k)fluoranthene	ug/L	10
Benzo(g,h,i)perylene	ug/L	10
Benzo(a)pyrene	ug/L	10
Benzyl alcohol	ug/L	10
bis(2-Chloroethoxy)- methane	ug/L	10
bis(2-Chloroethyl)ether	ug/L	10
bis(2-Chloroisopropyl) ether	ug/L	10
bis(2-Ethylhexyl) phthalate	ug/L	10
4-Bromophenyl phenyl ether	ug/L	10
Butyl benzyl phthalate	ug/L	10
2sec-Butyl-4,6-dinitro- phenol (Dinoseb)	ug/L	10
4-Chloroaniline	ug/L	10
4-Chloro-3-methylphenol	ug/L	10
2-Chloronaphthalene	ug/L	10
2-Chlorophenol	ug/L	10
4-Chlorophenyl phenyl ether	ug/L	10
o-Cresol	ug/L	10
m & p-Cresol(s)	ug/L	10
Chrysene	ug/L	10
Dibenz(a,h)anthracene	ug/L	10
Dibenzofuran	ug/L	10
Di-n-butyl phthalate	ug/L	10
1,2-Dichlorobenzene	ug/L	10
1,3-Dichlorobenzene	ug/L	10
1,4-Dichlorobenzene	ug/L	10

Table 9 cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting Limit
3,3'-Dichlorobenzidine	ug/L	20
2,4-Dichlorophenol	ug/L	10
2,6-Dichlorophenol	ug/L	10
Diethyl phthalate	ug/L	10
Dimethoate	ug/L	10
p-Dimethylaminoazobenzene	ug/L	10
7,12-Dimethylbenz-anthracene	ug/L	10
3,3'-Dimethylbenzidine	ug/L	10
a,a-Dimethylphen-ethylamine	ug/L	10
2,4-Dimethylphenol	ug/L	10
Dimethyl phthalate	ug/L	10
1,3-Dinitrobenzene	ug/L	10
4,6-Dinitro-2-methylphenol	ug/L	10
4,6-Dinitro-o-cresol	ug/L	50
2,4-Dinitrophenol	ug/L	50
2,4-Dinitrotoluene	ug/L	10
2,6-Dinitrotoluene	ug/L	10
Di-n-octyl phthalate	ug/L	10
Diphenylamine	ug/L	10
Disulfoton	ug/L	50
bis(2-Ethylhexyl) phthalate	ug/L	10
Ethyl methanesulfonate	ug/L	10
Famphur	ug/L	--
Flouranthene	ug/L	10
Flourene	ug/L	10
Hexachlorobenzene	ug/L	10
Hexachlorobutadiene	ug/L	10
Hexachlororocyclopentadiene	ug/L	10
Hexachloroethane	ug/L	--
Hexachlorophene	ug/L	10
Hexachloropropene	ug/L	10
Indeno(1,2,3-c,d)pyrene	ug/L	10
Isophorone	ug/L	10
Isosafrole	ug/L	20
Methapyrilene	ug/L	10
3-Methylcholanthrene	ug/L	10
Methyl methanesulfonate	ug/L	10
2-Methylnaphthalen	ug/L	10
Methyl parathion	ug/L	50
2-Methylphenol	ug/L	10
3/4-Methylphenol	ug/L	10
Methyl methacrylate	ug/L	10
Napthalene	ug/L	10

Table 9 Conti

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Reporting	
	Units	Limit
1,4-Naphthaquinone	ug/L	10
1-Naphthylamine	ug/L	10
2-Naphthylamine	ug/L	10
2-Nitroaniline	ug/L	50
3-Nitroaniline	ug/L	50
4-Nitroaniline	ug/L	50
Nitrobenzene	ug/L	10
2-Nitrophenol	ug/L	10
4-Nitrophenol	ug/L	50
4-Nitroquinoline-1-oxide	ug/L	--
N-Nitroso-di-n-butylamine	ug/L	10
N-Nitrosodiethylamine	ug/L	10
N-Nitrosodimethylamine	ug/L	10
N-Nitrosodiphenylamine	ug/L	10
N-Nitroso-di-n-propylamine	ug/L	10
N-Nitrosomethylethylamine	ug/L	10
N-Nitrosomorpholine	ug/L	10
N-Nitrosopiperidine	ug/L	10
5-Nitro-o-toluidine	ug/L	10
N-Nitrosopyrrolidine	ug/L	10
Parathion	ug/L	50
Pentachlorobenzene	ug/L	10
Pentachlorethane	ug/L	10
Pentachloronitrobenzene	ug/L	50
Pentachlorophenol	ug/L	50
Phenacetin	ug/L	10
Phenanthrene	ug/L	10
Phenol	ug/L	10
4-Phenylenediamine	ug/L	--
Phorate	ug/L	100
2-Picoline	ug/L	10
Pronamide	ug/L	10
Pyrene	ug/L	10
Pyridine	ug/L	20
Safrole	ug/L	10
Sulfotapp	ug/L	50
1,2,4,5-Tetrachloro- benzene	ug/L	10
2,3,4,6-Tetrachlorophenol	ug/L	50
Thionazin	ug/L	50

Table 9 Cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting Limit
sym-Trinitrobenzene	ug/L	10
2-Toluidine	ug/L	10
1,2,4-Trichlorobenzene	ug/L	10
2,4,5-Trichlorophenol	ug/L	50
0,0,0-Triethylphosphoro- thioate	ug/L	10
2,4,6-Trichlorophenol	ug/L	10
1,3,5-Trinitrobenzene	ug/L	10
Ethyl methacrylate	ug/L	10



Analytical **Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 104902

May 17, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: OCD Plan 997-9008-28

Attention: Claud Rosendale

On 04/30/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

DATE RECEIVED : 04/30/91

REPORT DATE : 05/16/91

ATI I.D. : 104902

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	OW-25	AQUEOUS	04/29/91

----- TOTALS -----

MATRIX	# SAMPLES
AQUEOUS	1

ATI STANDARD DISPOSAL PRACTICE

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



Analytical **Technologies**, Inc.

GENERAL CHEMISTRY RESULTS

ATI I.D. : 104902

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

DATE RECEIVED : 04/30/91

REPORT DATE : 05/16/91

PARAMETER	UNITS	01
CARBONATE (CACO3)	MG/L	<1
BICARBONATE (CACO3)	MG/L	525
HYDROXIDE (CACO3)	MG/L	<1
TOTAL ALKALINITY (AS CACO3)	MG/L	525
CHLORIDE	MG/L	91
CONDUCTIVITY, (UMHOS/CM)		1190
PH	UNITS	8.51
PHENOLICS, TOTAL	MG/L	<0.02
SULFATE	MG/L	27
TOTAL DISSOLVED SOLIDS	MG/L	780



Analytical **Technologies**, Inc.

GENERAL CHEMISTRY - QUALITY CONTROL

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN

ATI I.D. : 104902

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
CARBONATE	MG/L	10491101	<1	<1	NA	NA	NA	NA
BICARBONATE	MG/L		126	125	1	NA	NA	NA
HYDROXIDE	MG/L		<1	<1	NA	NA	NA	NA
TOTAL ALKALINITY	MG/L		126	125	1	NA	NA	NA
CHLORIDE	MG/L	10555503	390	390	0	1150	750	101
CONDUCTIVITY(UMHOS/CM)		10490201	1190	1240	4	NA	NA	NA
PH	UNITS	10563801	8.2	8.3	1	NA	NA	NA
PHENOLICS, TOTAL	MG/L	10550401	<0.02	<0.02	NA	0.25	0.25	100
SULFATE	MG/L	10488717	130	140	7	280	140	100
TOTAL DISSOLVED SOLIDS	MG/L	10490101	1100	1100	0	NA	NA	NA

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$



Analytical **Technologies**, Inc.

METALS RESULTS

ATI I.D. : 104902

CLIENT : GIANT REFINING CO.

DATE RECEIVED : 04/30/91

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

REPORT DATE : 05/16/91

PARAMETER	UNITS	01
SILVER	MG/L	<0.010
ARSENIC	MG/L	<0.005
BARIUM	MG/L	0.148
CALCIUM	MG/L	15.6
CADMIUM	MG/L	0.008
CHROMIUM	MG/L	<0.010
MANGANESE	MG/L	0.032
SODIUM	MG/L	266
LEAD	MG/L	0.002
SELENIUM	MG/L	<0.005



Analytical Technologies, Inc.

METALS - QUALITY CONTROL

CLIENT : GIANT REFINING CO.

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

ATI I.D. : 104902

PARAMETER	UNITS	ATI I.D.	SAMPLE RESULT	DUP. RESULT	RPD	SPIKED SAMPLE	SPIKE CONC	% REC
SILVER	MG/L	10489603	<0.010	<0.010	NA	0.440	0.500	88
ARSENIC	MG/L	10551103	<0.005	<0.005	NA	0.038	0.050	76
BARIUM	MG/L	10489603	0.041	0.040	2	0.962	1.00	92
CALCIUM	MG/L	10491401	30.6	30.7	0.3	83.9	50.0	107
CADMIUM	MG/L	10489603	<0.005	0.007	NA	0.482	0.500	96
CHROMIUM	MG/L	10489603	<0.010	<0.010	NA	0.876	1.00	88
MANGANESE	MG/L	10489603	2.32	2.33	0.4	3.26	1.00	94
SODIUM	MG/L	10491401	34.1	35.5	4	85.4	50.0	103
LEAD	MG/L	10490101	0.009	0.009	0	0.057	0.050	96
SELENIUM	MG/L	10490101	<0.005	<0.005	NA	0.030	0.050	60

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative Percent Difference)} = \frac{(\text{Sample Result} - \text{Duplicate Result})}{\text{Average Result}} \times 100$$

GAS CHROMATOGRAPHY - RESULTS

ATI I.D. : 10490201

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT	: GIANT REFINING CO.	DATE SAMPLED	: 04/29/91
PROJECT #	: 997-9008-28	DATE RECEIVED	: 04/30/91
PROJECT NAME	: OCD PLAN	DATE EXTRACTED	: N/A
CLIENT I.D.	: OW-25	DATE ANALYZED	: 05/09/91
SAMPLE MATRIX	: AQUEOUS	UNITS	: UG/L
		DILUTION FACTOR	: 1

COMPOUNDS

RESULTS

BENZENE	<0.5
BROMODICHLOROMETHANE	<0.2
BROMOFORM	<0.2
BROMOMETHANE	<0.2
CARBON TETRACHLORIDE	<0.2
CHLOROBENZENE	<0.5
CHLOROETHANE	<0.2
CHLOROFORM	<0.2
CHLOROMETHANE	<0.2
DIBROMOCHLOROMETHANE	<0.2
1,3-DICHLOROBENZENE	<0.5
2-CHLOROETHYL VINYL ETHER	<0.5
1,2 & 1,4-DICHLOROBENZENE	<0.5
DICHLORODIFLUOROMETHANE	<0.2
1,1-DICHLOROETHANE	1.4
1,2-DICHLOROETHANE	21.5
1,1-DICHLOROETHENE	<0.2
1,2-DICHLOROETHENE (TOTAL)	<0.2
1,2-DICHLOROPROPANE	<0.2
CIS-1,3-DICHLOROPROPENE	<0.2
TRANS-1,3-DICHLOROPROPENE	<0.2
ETHYLBENZENE	<0.5
METHYLENE CHLORIDE	<2.0
1,1,2,2-TETRACHLOROETHANE	<0.2
TETRACHLOROETHENE	<0.2
TOLUENE	35.8
1,1,1-TRICHLOROETHANE	<0.2
1,1,2-TRICHLOROETHANE	<0.2
TRICHLOROETHENE	0.2
TRICHLOROFLUOROMETHANE	<0.5
VINYL CHLORIDE	<0.2
TOTAL XYLENES	<0.5
TRICHLOROFLUOROMETHANE	<0.2

SURROGATE PERCENT RECOVERIES

BROMOCHLOROMETHANE (%)	106
BROMOFLUOROBENZENE (%)	100

Analytical**Technologies**, Inc.

GAS CHROMATOGRAPHY - RESULTS

REAGENT BLANK

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 104902
PROJECT #	: 997-9008-28	DATE EXTRACTED	: 05/09/91
PROJECT NAME	: OCD PLAN	DATE ANALYZED	: 05/09/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

BENZENE	<0.5
BROMODICHLOROMETHANE	<0.2
BROMOFORM	<0.2
BROMOMETHANE	<0.2
CARBON TETRACHLORIDE	<0.2
CHLOROBENZENE	<0.5
CHLOROETHANE	<0.2
CHLOROFORM	<0.2
CHLOROMETHANE	<0.2
DIBROMOCHLOROMETHANE	<0.2
1,3-DICHLOROBENZENE	<0.5
2-CHLOROETHYL VINYL ETHER	<0.5
1,2 & 1,4-DICHLOROBENZENE	<0.5
DICHLORODIFLUOROMETHANE	<0.2
1,1-DICHLOROETHANE	<0.2
1,2-DICHLOROETHANE	<0.2
1,1-DICHLOROETHENE	<0.2
1,2-DICHLOROETHENE (TOTAL)	<0.2
1,2-DICHLOROPROPANE	<0.2
CIS-1,3-DICHLOROPROPENE	<0.2
TRANS-1,3-DICHLOROPROPENE	<0.2
ETHYLBENZENE	<0.5
METHYLENE CHLORIDE	<2.0
1,1,2,2-TETRACHLOROETHANE	<0.2
TETRACHLOROETHENE	<0.2
TOLUENE	<0.5
1,1,1-TRICHLOROETHANE	<0.2
1,1,2-TRICHLOROETHANE	<0.2
TRICHLOROETHENE	<0.2
TRICHLOROFLUOROMETHANE	<0.5
VINYL CHLORIDE	<0.2
TOTAL XYLENES	<0.5
TRICHLOROFLUOROMETHANE	<0.2

SURROGATE PERCENT RECOVERIES

BROMOCHLOROMETHANE (%)	106
BROMOFLUOROBENZENE (%)	98



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 104902

TEST : VOLATILE HALOCARBONS/AROMATICS (EPA 8010/8020)

CLIENT : GIANT REFINING CO.

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

REF I.D. : 10489901

DATE ANALYZED : 05/10/91

SAMPLE MATRIX : AQUEOUS

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	% REC.	DUP. SPIKED %		RPD
	RESULT	SPIKED			SAMPLE	REC.	
1,1-DICHLOROETHENE	<0.2	20	20	100	21	105	5
TRICHLOROETHENE	<0.2	20	19	95	20	100	5
TETRACHLOROETHENE	<0.2	20	18	90	19	95	5
BENZENE	<0.5	20	18	90	19	95	5
BROMODICHLOROMETHANE	<0.2	20	19	95	19	95	0
CHLOROFORM	<0.2	20	19	95	19	95	0
1,1,1-TRICHLOROETHANE	<0.2	20	19	95	19	95	0
TOLUENE	<0.5	20	19	95	18	90	5
CHLOROBENZENE	<0.5	20	18	90	18	90	0
M-XYLENE	<0.5	20	18	90	18	90	0

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$

GCMS - RESULTS

ATI I.D. : 10490201

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
 PROJECT # : 997-9008-28
 PROJECT NAME : OCD PLAN
 CLIENT I.D. : OW-25
 SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
 DATE RECEIVED : 04/30/91
 DATE EXTRACTED : 05/01/91
 DATE ANALYZED : 05/11/91
 UNITS : UG/L
 DILUTION FACTOR : 1

COMPOUNDS

RESULTS

COMPOUNDS	RESULTS
ACETOPHENONE	<10
2-ACETYLAMINOFLUORENE	<10
a,a-DIMETHYLPHENETHYLAMINE	<10
4-AMINOBIIPHENYL	<10
ARAMITE	<50
CHLOROBENZILATE	<10
DIALATE	<10
2,6-DICHLOROPHENOL	<10
DIMETHOATE	<10
p-(DIMETHYLAMINO)AZOBENZENE	<10
7,12-DIMETHYLBENZO(a)ANTHRACENE	<10
3,3'-DIMETHYLBENZIDINE	<10
m-DINITROBENZENE	<10
DIPHENYLAMINE	<10
ETHYL METHANESULFONATE	<10
HEXACHLOROPHENE	<10
HEXACHLOROPROPENE	<10
ISOSAFROLE	<10
METHAPYRILENE	<10
3-METHYLCHOLANTHRENE	<10
METHYL METHANESULFONATE	<10
3-METHYLPHENOL (m-CRESOL)	<10
1,4-NAPHTHOQUINONE	<10
1-NAPHTHYLAMINE	<10
2-NAPHTHYLAMINE	<10
5-NITRO-O-TOLUIDINE	<10
4-NITROQUINOLINE-1-OXIDE	<10
N-NITROSODIETHYLAMINE	<10
N-NITROSODI-BUTYLAMINE	<10
N-NITROSOMETHYLETHYLAMINE	<200
N-NITROSOMORPHOLINE	<10
N-NITROSOPIPERIDINE	<10
N-NITROSOPYRROLIDINE	<10
PENTACHLOROBENZENE	<10
PENTACHLORONITROBENZENE	<10
PENTACHLOROETHANE	<10
PHENACETIN	<10
p-PHENYLENEDIAMINE	<10
2-PICOLINE	<10
PRONAMIDE	<10
SAFROLE	<10

(CONTINUED NEXT PAGE)



GCMS - RESULTS

ATI I.D. : 10490201

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
DIETHYLPHTHALATE	<10
4-CHLOROPHENYL-PHENYLETHER	<10
FLUORENE	<10
4-NITROANILINE	<50
4,6-DINITRO-2-METHYLPHENOL	<50
N-NITROSODIPHENYLAMINE	<10
4-BROMOPHENYL-PHENYLETHER	<10
HEXACHLOROBENZENE	<10
PENTACHLOROPHENOL	<50
PHENANTHRENE	<10
ANTHRACENE	<10
DI-N-BUTYLPHTHALATE	<10
FLUORANTHENE	<10
BENZIDINE	<100
PYRENE	<10
BUTYLBENZYLPHTHALATE	<10
3,3'-DICHLOOROBENZIDINE	<20
BENZO(a)ANTHRACENE	<10
BIS(2-ETHYLHEXYL)PHTHALATE	<10
CHRYSENE	<10
DI-N-OCTYLPHTHALATE	<10
BENZO(b)FLUORANTHENE	<10
BENZO(k)FLUORANTHENE	<10
BENZO(a)PYRENE	<10
INDENO(1,2,3-cd)PYRENE	<10
DIBENZO(a,h)ANTHRACENE	<10
BENZO(g,h,i)PERYLENE	<10

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	66
2-FLUOROBIPHENYL (%)	60
TERPHENYL (%)	114
PHENOL-D6 (%)	52
2-FLUOROPHENOL (%)	60
2,4,6-TRIBROMOPHENOL (%)	61



Analytical **Technologies**, Inc.

GCMS - RESULTS

ATI I.D. : 10490201

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
1,2,4,5-TETRACHLOROBENZENE	<10
2,3,4,6-TETRACHLOROPHENOL	<10
o-TOLUIDINE	<10
1,3,5-TRINITROBENZENE	<10
PYRIDINE	<10

GCMS - RESULTS

ATI I.D. : 10490201

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : OW-25
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 04/29/91
DATE RECEIVED : 04/30/91
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/11/91
UNITS : UG/L
DILUTION FACTOR : 1

COMPOUNDS	RESULTS
-----------	---------

N-NITROSODIMETHYLAMINE	<10
PHENOL	<10
ANILINE	<10
BIS(2-CHLOROETHYL) ETHER	<10
2-CHLOROPHENOL	<10
1,3-DICHLOROBENZENE	<10
1,4-DICHLOROBENZENE	<10
BENZYL ALCOHOL	<10
1,2-DICHLOROBENZENE	<10
2-METHYLPHENOL	<10
BIS(2-CHLOROISOPROPYL) ETHER	<10
4-METHYLPHENOL	<10
N-NITROSO-DI-N-PROPYLAMINE	<10
HEXACHLOROETHANE	<10
NITROBENZENE	<10
ISOPHORONE	<10
2-NITROPHENOL	<10
2,4-DIMETHYLPHENOL	<10
BENZOIC ACID	<50
BIS(2-CHLOROETHOXY) METHANE	<10
2,4-DICHLOROPHENOL	<10
1,2,4-TRICHLOROBENZENE	<10
NAPHTHALENE	<10
4-CHLOROANILINE	<10
HEXACHLOROBUTADIENE	<10
4-CHLORO-3-METHYLPHENOL	<10
2-METHYLNAPHTHALENE	<10
HEXACHLOROCYCLOPENTADIENE	<10
2,4,6-TRICHLOROPHENOL	<10
2,4,5-TRICHLOROPHENOL	<50
2-CHLORONAPHTHALENE	<10
2-NITROANILINE	<50
DIMETHYLPHTHALATE	<10
ACENAPHTHYLENE	<10
3-NITROANINLINE	<50
ACENAPHTHENE	<10
2,4-DINITROPHENOL	<50
4-NITROPHENOL	<50
DIBENZOFURAN	<10
2,4-DINITROTOLUENE	<10
2,6-DINITROTOLUENE	<10

(CONTINUED NEXT PAGE)

GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.
PROJECT # : 997-9008-28
PROJECT NAME : OCD PLAN
CLIENT I.D. : REAGENT BLANK

ATI I.D. : 104902
DATE EXTRACTED : 05/01/91
DATE ANALYZED : 05/10/91
UNITS : UG/L
DILUTION FACTOR : N/A

COMPOUNDS	RESULTS
ACETOPHENONE	<10
2-ACETYLAMINOFLUORENE	<10
a,a-DIMETHYLPHENETHYLAMINE	<10
4-AMINOBIIPHENYL	<10
ARAMITE	<50
CHLOROBENZILATE	<10
DIALATE	<10
2,6-DICHLOROPHENOL	<10
DIMETHOATE	<10
p-(DIMETHYLAMINO)AZOBENZENE	<10
7,12-DIMETHYLBENZO(a)ANTHRACENE	<10
3,3'-DIMETHYLBENZIDINE	<10
m-DINITROBENZENE	<10
DIPHENYLAMINE	<10
ETHYL METHANESULFONATE	<10
HEXACHLOROPHENE	<10
HEXACHLOROPROPENE	<10
ISOSAFROLE	<10
METHAPYRILENE	<10
3-METHYLCHOLANTHRENE	<10
METHYL METHANESULFONATE	<10
3-METHYLPHENOL (m-CRESOL)	<10
1,4-NAPHTHOQUINONE	<10
1-NAPHTHYLAMINE	<10
2-NAPHTHYLAMINE	<10
5-NITRO-O-TOLUIDINE	<10
4-NITROQUINOLINE-1-OXIDE	<10
N-NITROSODIETHYLAMINE	<10
N-NITROSODI-BUTYLAMINE	<10
N-NITROSOMETHYLETHYLAMINE	<200
N-NITROSOMORPHOLINE	<10
N-NITROSOPIPERIDINE	<10
N-NITROSOPYRROLIDINE	<10
PENTACHLOROBENZENE	<10
PENTACHLORONITROBENZENE	<10
PENTACHLOROETHANE	<10
PHENACETIN	<10
p-PHENYLENEDIAMINE	<10
2-PICOLINE	<10
PRONAMIDE	<10
SAFROLE	<10
1,2,4,5-TETRACHLOROBENZENE	<10
2,3,4,6-TETRACHLOROPHENOL	<10

(CONTINUED NEXT PAGE)



Analytical Technologies, Inc.

GCMS - RESULTS

REAGENT BLANK

ATI I.D. : 104902

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS	RESULTS
FLUORENE	<10
4-NITROANILINE	<50
4,6-DINITRO-2-METHYLPHENOL	<50
N-NITROSODIPHENYLAMINE	<10
4-BROMOPHENYL-PHENYLETHER	<10
HEXACHLOROBENZENE	<10
PENTACHLOROPHENOL	<50
PHENANTHRENE	<10
ANTHRACENE	<10
DI-N-BUTYLPHTHALATE	<10
FLUORANTHENE	<10
BENZIDINE	<100
PYRENE	<10
BUTYLBENZYLPHTHALATE	<10
3,3'-DICHLOOROBENZIDINE	<20
BENZO(a)ANTHRACENE	<10
BIS(2-ETHYLHEXYL)PHTHALATE	<10
CHRYSENE	<10
DI-N-OCTYLPHTHALATE	<10
BENZO(b)FLUORANTHENE	<10
BENZO(k)FLUORANTHENE	<10
BENZO(a)PYRENE	<10
INDENO(1,2,3-cd)PYRENE	<10
DIBENZO(a,h)ANTHRACENE	<10
BENZO(g,h,i)PERYLENE	<10

SURROGATE PERCENT RECOVERIES

NITROBENZENE-D5 (%)	41
2-FLUOROBIPHENYL (%)	49
TERPHENYL (%)	96
PHENOL-D6 (%)	57
2-FLUOROPHENOL (%)	71
2,4,6-TRIBROMOPHENOL (%)	40



Analytical **Technologies**, Inc.

GCMS - RESULTS

REAGENT BLANK

ATI I.D. : 104902

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

COMPOUNDS

RESULTS

o-TOLUIDINE

<10

1,3,5-TRINITROBENZENE

<10

PYRIDINE

<10

GCMS - RESULTS

REAGENT BLANK

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT	: GIANT REFINING CO.	ATI I.D.	: 104902
PROJECT #	: 997-9008-28	DATE EXTRACTED	: 05/01/91
PROJECT NAME	: OCD PLAN	DATE ANALYZED	: 05/10/91
CLIENT I.D.	: REAGENT BLANK	UNITS	: UG/L
		DILUTION FACTOR	: N/A

COMPOUNDS	RESULTS
-----------	---------

N-NITROSODIMETHYLAMINE	<10
PHENOL	<10
ANILINE	<10
BIS(2-CHLOROETHYL) ETHER	<10
2-CHLOROPHENOL	<10
1,3-DICHLOROBENZENE	<10
1,4-DICHLOROBENZENE	<10
BENZYL ALCOHOL	<10
1,2-DICHLOROBENZENE	<10
2-METHYLPHENOL	<10
BIS(2-CHLOROISOPROPYL) ETHER	<10
4-METHYLPHENOL	<10
N-NITROSO-DI-N-PROPYLAMINE	<10
HEXACHLOROETHANE	<10
NITROBENZENE	<10
ISOPHORONE	<10
2-NITROPHENOL	<10
2,4-DIMETHYLPHENOL	<10
BENZOIC ACID	<50
BIS(2-CHLOROETHOXY) METHANE	<10
2,4-DICHLOROPHENOL	<10
1,2,4-TRICHLOROBENZENE	<10
NAPHTHALENE	<10
4-CHLOROANILINE	<10
HEXACHLOROBUTADIENE	<10
4-CHLORO-3-METHYLPHENOL	<10
2-METHYLNAPHTHALENE	<10
HEXACHLOROCYCLOPENTADIENE	<10
2,4,6-TRICHLOROPHENOL	<10
2,4,5-TRICHLOROPHENOL	<50
2-CHLORONAPHTHALENE	<10
2-NITROANILINE	<50
DIMETHYLPHTHALATE	<10
ACENAPHTHYLENE	<10
3-NITROANINLINE	<50
ACENAPHTHENE	<10
2,4-DINITROPHENOL	<50
4-NITROPHENOL	<50
DIBENZOFURAN	<10
2,4-DINITROTOLUENE	<10
2,6-DINITROTOLUENE	<10
DIETHYLPHTHALATE	<10
4-CHLOROPHENYL-PHENYLETHER	<10

(CONTINUED NEXT PAGE)



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 104902

TEST : SEMI-VOLATILE ORGANICS (EPA 8270)

CLIENT : GIANT REFINING CO.

PROJECT # : 997-9008-28

PROJECT NAME : OCD PLAN

REF I.D. : 10599911

DATE ANALYZED : 05/09/91

SAMPLE MATRIX : AQUEOUS

UNITS : UG/L

COMPOUNDS	SAMPLE RESULT	CONC. SPIKED	SPIKED SAMPLE	DUP.		RPD
				% SPIKED REC.	% SPIKED REC.	
1,2,4-TRICHLOROBENZENE	<10	50	37	74	34	8
ACENAPHTHENE	<10	50	36	72	41	13
2,4-DINITROTOLUENE	<10	50	31	62	34	9
PYRENE	<10	56	63	126	57	10
N-NITROSO-DI-N-PROPYL AMINE	<10	50	38	76	43	12
1,4-DICHLOROBENZENE	<10	50	42	84	41	2
PENTACHLOROPHENOL	<50	100	88	88	89	1
PHENOL	<10	100	77	77	79	3
2-CHLOROPHENOL	<10	100	84	84	82	2
4-CHLORO-3-METHYLPHENOL	<10	100	80	80	75	6
4-NITROPHENOL	<50	100	40	40	47	16

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$



DATE 4-29-99 / PAGE 1 OF 1

ATTI Labs: ~~San~~ **San Diego** (619)458-9141 • **Phoenix** (602)438-1530 • **Seattle** (206)228-8335 • **Pensacola** (904)474-1001

Table 9

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	nits	Reporting Limit
✓Acenaphthene	ug/L	10
✓Acenaphthylene	ug/L	10
✓Acetophenone	ug/L	10
✓2-Acetylaminofluorene	ug/L	10
✓4-Aminobiphenyl	ug/L	10
✓Aniline	ug/L	10
✓Anthracene	ug/L	10
✓Aranite	ug/L	10
✓Benzo(a)anthracene	ug/L	10
✓Benzo(b)fluoranthene	ug/L	10
✓Benzo(k)fluoranthene	ug/L	10
✓Benzo(g,h,i)perylene	ug/L	10
✓Benzo(a)pyrene	ug/L	10
✓Benzyl alcohol	ug/L	10
✓bis(2-Chloroethoxy)- methane	ug/L	10
✓bis(2-Chloroethyl)ether	ug/L	10
✓bis(2-Chloroisopropyl) ether	ug/L	10
✓bis(2-Ethylhexyl) phthalate	ug/L	10
✓4-Bromophenyl phenyl ether	ug/L	10
✓Butyl benzyl phthalate	ug/L	10
✓2sec-Butyl-4,6-dinitro- phenol (Dinoseb)	ug/L	10
✓4-Chloroaniline	ug/L	10
✓4-Chloro-3-methylphenol	ug/L	10
✓2-Chloronaphthalene	ug/L	10
✓2-Chlorophenol	ug/L	10
✓4-Chlorophenyl phenyl ether	ug/L	10
o-Cresol <i>ok 2-methyl</i>	ug/L	10
✓m & p-Cresol(s)	ug/L	10
✓Chrysene	ug/L	10
✓Dibenz(a,h)anthracene	ug/L	10
✓Dibenzofuran	ug/L	10
✓Di-n-butyl phthalate	ug/L	10
✓1,2-Dichlorobenzene	ug/L	10
✓1,3-Dichlorobenzene	ug/L	10
✓1,4-Dichlorobenzene	ug/L	10

Table 9 cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting Limit
✓3,3'-Dichlorobenzidine	ug/L	20
✓2,4-Dichlorophenol	ug/L	10
✓2,6-Dichlorophenol	ug/L	10
✓Diethyl phthalate	ug/L	10
✓Dimethoate	ug/L	10
✓p-Dimethylaminoazobenzene	ug/L	10
✓7,12-Dimethylbenzanthracene	ug/L	10
✓3,3'-Dimethylbenzidine	ug/L	10
✓4,4-Dimethylphenethylamine	ug/L	10
✓2,4-Dimethylphenol	ug/L	10
✓Diethyl phthalate	ug/L	10
✓1,3-Dinitrobenzene	ug/L	10
✓4,6-Dinitro-2-ethylphenol	ug/L	10
4,6-Dinitro-o-cresol <i>ok</i>	ug/L	50
✓2,4-Dinitrophenol	ug/L	50
✓2,4-Dinitrotoluene	ug/L	10
✓2,6-Dinitrotoluene	ug/L	10
✓Di-n-octyl phthalate	ug/L	10
✓Diphenylamine	ug/L	10
<i>de</i> Disulfoton <i>X</i>	ug/L	50
✓Bis(2-Ethylhexyl) phthalate	ug/L	10
✓Ethyl methanesulfonate	ug/L	10
<i>de</i> Faaphur <i>X</i>	ug/L	--
✓Flouranthene	ug/L	10
✓Flourene	ug/L	10
✓Hexachlorobenzene	ug/L	10
✓Hexachlorobutadiene	ug/L	10
✓Hexachlorocyclopentadiene	ug/L	10
✓Hexachloroethane	ug/L	--
✓Hexachlorophene	ug/L	10
✓Hexachloropropene	ug/L	10
✓Indeno(1,2,3-c,d)pyrene	ug/L	10
✓Isophorone	ug/L	10
✓Isosafrole	ug/L	20
✓Methapyrilene	ug/L	10
✓3-Methylcholanthrene	ug/L	10
✓Methyl methanesulfonate	ug/L	10
✓2-Methylnaphthalen	ug/L	10
<i>de</i> Methyl parathion <i>X</i>	ug/L	50
✓2-Methylphenol	ug/L	10
✓3/4-Methylphenol	ug/L	10
Methyl methacrylate. <i>voA</i>	ug/L	10
✓Naphthalene	ug/L	10

Table 9 Cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting Limit
✓ 1,4-Naphthaquinone	ug/L	10
✓ 1-Naphthylamine	ug/L	10
✓ 2-Naphthylamine	ug/L	10
✓ 2-Nitroaniline	ug/L	50
✓ 3-Nitroaniline	ug/L	50
✓ 4-Nitroaniline	ug/L	50
✓ Nitrobenzene	ug/L	10
✓ 2-Nitrophenol	ug/L	10
✓ 4-Nitrophenol OK	ug/L	50
✓ 4-Nitroquinoline-1-oxide	ug/L	--
✓ N-Nitroso-di-n-butylamine	ug/L	10
✓ N-Nitrosodiethylamine	ug/L	10
✓ N-Nitrosodisethylamine	ug/L	10
✓ N-Nitrosodiphenylamine	ug/L	10
✓ N-Nitroso-di-n-propylamine	ug/L	10
✓ N-Nitrosomethylethylamine	ug/L	10
✓ N-Nitrosomorpholine	ug/L	10
✓ N-Nitrosopiperidine	ug/L	10
✓ 5-Nitro-o-toluidine	ug/L	10
✓ N-Nitrosopyrrolidine	ug/L	10
✓ Parathion	ug/L	50
✓ Pentachlorobenzene	ug/L	10
✓ Pentachlorethane	ug/L	10
✓ Pentachloronitrobenzene	ug/L	50
✓ Pentachlorophenol	ug/L	50
✓ Phenacetin	ug/L	10
✓ Phenanthrene	ug/L	10
✓ Phenol	ug/L	10
✓ 4-Phenylenediamine	ug/L	--
✓ Phorate	ug/L	100
✓ 2-Picoline	ug/L	10
✓ Pronamide	ug/L	10
✓ Pyrene	ug/L	10
✓ Pyridine	ug/L	20
✓ Saffrole	ug/L	10
✓ Sulfotepp	ug/L	50
1,2,4,5-Tetrachloro- benzene	ug/L	10
✓ 2,3,4,6-Tetrachlorophenol	ug/L	50
✓ Thionazin	ug/L	50

WA ✓ - 5 ppb

Thiophosphoric acid tetraethyl ester

O-2-pyrazinyl phosphorothioate

These 2 compounds are part of App IX 8140 list in the library

Table 9 Cont.

GIANT REFINING GALLUP, NEW MEXICO

APPENDIX IX SEMIVOLATILE ORGANICS

Parameter	Units	Reporting Limit
<i>h</i> sym-Trinitrobenzene	ug/L	10
✓ 2-Toluidine	ug/L	10
✓ 1,2,4-Trichlorobenzene	ug/L	10
<i>c</i> 2,4,5-Trichlorophenol	ug/L	50
<i>h</i> 0,0,0-Triethylphosphorochioate	ug/L	10
<i>c</i> 2,4,6-Trichlorophenol	ug/L	10
✓ 1,3,5-Trinitrobenzene	ug/L	10
Ethyl methacrylate VOA	ug/L	10



Analytical **Technologies**, Inc.

9830 S. 51st Street Suite B-113 Phoenix, AZ 85044 (602) 496-4400

ATI I.D. 105575

May 14, 1991

Giant Refining Company
Route 3, P.O. Box 7
Gallup, NM 87301

Project Name/Number: OCD

Attention: Claud Rosendale

On 05/04/91, Analytical Technologies, Inc. received a request to analyze aqueous sample(s). The sample(s) 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.

If you have any questions or comments, please do not hesitate to contact us at (602) 496-4400.

Elizabeth Proffitt
Senior Project Manager

Robert V. Woods
Laboratory Manager

RVW:clf
Enclosure



Analytical Technologies, Inc.

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : OCD

DATE RECEIVED : 05/04/91

REPORT DATE : 05/13/91

ATI I.D. : 105575

ATI #	CLIENT DESCRIPTION	MATRIX	DATE COLLECTED
01	AER. LAGOON #1 EFF	AQUEOUS	05/03/91
02	AER. LAGOON #2 EFF	AQUEOUS	05/03/91

----- TOTALS -----

MATRIX	# SAMPLES
AQUEOUS	2

ATI STANDARD DISPOSAL PRACTICE

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



Analytical Technologies, Inc.

GAS CHROMATOGRAPHY - RESULTS

ATI I.D. : 10557501

TEST : BTEX & MTBE (EPA METHOD 602)

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : OCD
CLIENT I.D. : AER. LAGOON #1 EFF
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 05/03/91
DATE RECEIVED : 05/04/91
DATE EXTRACTED : N/A
DATE ANALYZED : 05/09/91
UNITS : UG/L
DILUTION FACTOR : 5

COMPOUNDS

RESULTS

BENZENE	200
TOLUENE	320
ETHYLBENZENE	31
TOTAL XYLENES	200
METHYL-t-BUTYL ETHER	<12.5

SURROGATE PERCENT RECOVERIES

BROMOFLUOROBENZENE (%)	93
------------------------	----



Analytical Technologies, Inc.

GAS CHROMATOGRAPHY - RESULTS

ATI I.D. : 10557502

TEST : BTEX & MTBE (EPA METHOD 602)

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : OCD
CLIENT I.D. : AER. LAGOON #2 EFF
SAMPLE MATRIX : AQUEOUS

DATE SAMPLED : 05/03/91
DATE RECEIVED : 05/04/91
DATE EXTRACTED : N/A
DATE ANALYZED : 05/09/91
UNITS : UG/L
DILUTION FACTOR : 5

COMPOUNDS

RESULTS

BENZENE	26
TOLUENE	40
ETHYLBENZENE	5.2
TOTAL XYLENES	40
METHYL-t-BUTYL ETHER	<12.5

SURROGATE PERCENT RECOVERIES

BROMOFLUOROBENZENE (%)	86
------------------------	----



Analytical Technologies, Inc.

GAS CHROMATOGRAPHY - RESULTS

REAGENT BLANK

TEST : BTEX & MTBE (EPA METHOD 602)

CLIENT : GIANT REFINING CO.
PROJECT # : (NONE)
PROJECT NAME : OCD
CLIENT I.D. : REAGENT BLANK

ATI I.D. : 105575
DATE EXTRACTED : 05/09/91
DATE ANALYZED : 05/09/91
UNITS : UG/L
DILUTION FACTOR : N/A

COMPOUNDS	RESULTS
BENZENE	<0.5
TOLUENE	<0.5
ETHYLBENZENE	<0.5
TOTAL XYLENES	<0.5
METHYL-t-BUTYL ETHER	<2.5

SURROGATE PERCENT RECOVERIES

BROMOFLUOROBENZENE (%)	100
------------------------	-----



Analytical Technologies, Inc.

QUALITY CONTROL DATA

ATI I.D. : 105575

TEST : BTEX & MTBE (EPA METHOD 602)

CLIENT : GIANT REFINING CO.

PROJECT # : (NONE)

PROJECT NAME : OCD

REF I.D. : 10599905

DATE ANALYZED : 05/06/91

SAMPLE MATRIX : AQUEOUS

UNITS : UG/L

COMPOUNDS	SAMPLE CONC.		SPIKED SAMPLE	DUP. % SPIKED		DUP. %	RPD
	RESULT	SPIKED		REC.	SAMPLE REC.		
BENZENE	<0.5	20	20	100	19	95	5
TOLUENE	<0.5	20	20	100	19	95	5
ETHYLBENZENE	<0.5	20	20	100	18	90	11
TOTAL XYLENES	<0.5	60	59	98	54	92	9
METHYL-t-BUTYL ETHER	<2.5	40	41	103	38	95	8

$$\% \text{ Recovery} = \frac{(\text{Spike Sample Result} - \text{Sample Result})}{\text{Spike Concentration}} \times 100$$

$$\text{RPD (Relative \% Difference)} = \frac{(\text{Spiked Sample Result} - \text{Duplicate Spike Sample Result})}{\text{Average of Spiked Sample}} \times 100$$



DATE 5-3-91 PAGE 1 OF 1

ATI Labs: San Diego (619)458-9141 • Phoenix (602)438-1530 • Seattle (206)228-8335 • Pensacola (904)474-1001
DISTRIBUTION: White, Canary - ANALYTICAL TECHNOLOGIES, INC.; Pink - ORIGINATOR