

December 13, 1996

Alameda County Health Care Services 1131 Harbor Bay Parkway Alameda, California 94502

RE:

Unocal Service Station #1871 96 MacArthur Boulevard Oakland, California

Per the request of the Unocal Corporation Project Manager, Ms. Tina R. Berry, enclosed please find our report (MPDS-UN1871-13) dated December 2, 1996 for the above referenced site.

Should you have any questions regarding the reporting of data, please feel free to call our office at (510) 602-5120. Any other questions may be directed to the Project Manager at (510) 277-2321.

Sincerely,

MPDS Services, Inc.

Jarrel F. Crider

/jfc

Enclosure

cc: Ms. Tina R. Berry



MPDS-UN1871-13 December 2, 1996

Unocal Corporation 2000 Crow Canyon Place, Suite 400 P.O. Box 5155 San Ramon, California 94583

Attention: Mr. Robert A. Boust

RE: Quarterly Data Report

Unocal Service Station #1871 96 MacArthur Boulevard Oakland, California

Dear Mr. Boust:

This data report presents the results of the most recent quarter of monitoring and sampling of the monitoring wells at the referenced site by MPDS Services, Inc.

RECENT FIELD ACTIVITIES

The monitoring wells that were monitored and sampled during this quarter are indicated in Table 1. Prior to sampling, the wells were checked for depth to water and the presence of free product or sheen. The monitoring data and the ground water elevations are summarized in Table 1. The ground water flow direction during the most recent quarter is shown on the attached Figure 1.

Ground water samples were collected on October 24, 1996. Prior to sampling, the wells were each purged of between 4.5 and 27.5 gallons of water. During purging operations, the field parameters pH, temperature, and electrical conductivity were recorded on the purging/sampling data sheets which are attached to this report. Once the field parameters were observed to stabilize, and where possible, a minimum of approximately three casing volumes had been removed from each well, samples were then collected using a clean Teflon bailer. The samples were decanted into clean VOA vials and/or one-liter amber bottles, as appropriate, which were then sealed with Teflon-lined screw caps, labeled, and stored in a cooler, on ice, until delivery to a state-certified laboratory. Field blank and Trip blank samples (denoted as ES1 and ES2, respectively) were also collected for quality assurance and control. MPDS Services, Inc. transported the purged ground water to the Unocal Refinery located in Rodeo, California, for treatment and discharge to San Pablo Bay under NPDES permit.

ANALYTICAL RESULTS

The ground water samples were analyzed at Sequoia Analytical Laboratory and were accompanied by properly executed Chain of Custody documentation. The analytical results of the ground water samples collected to date are summarized in Table 2. The concentrations of Total Petroleum Hydrocarbons (TPH) as gasoline and benzene detected in the ground water samples collected this

MPDS-UN1871-13 December 2, 1996 Page 2

quarter are shown on the attached Figure 2. Copies of the laboratory analytical results and the Chain of Custody documentation are attached to this report.

LIMITATIONS

Environmental changes, either naturally-occurring or artificially-induced, may cause changes in ground water levels and flow paths, thereby changing the extent and concentration of any contaminants.

DISTRIBUTION

A copy of this report should be sent to the Alameda County Health Care Services Agency.

If you have any questions regarding this report, please do not hesitate to call Mr. Joel G. Greger at (510) 602-5120.

JOEL G. GREGER
No. EG 1633
CERTIFIED
ENGINEERING
GEOLOGIST

Sincerely,

MPDS Services, Inc.

Haig (Gary) Tejirian Senior Staff Geologist

Joel G. Greger, C.E.G. Senior Engineering Geologist

License No. EG 1633 Exp. Date 8/31/98

Attachments: Tables 1 & 2

Location Map Figures 1 & 2

Laboratory Analyses

Chain of Custody documentation Purging/Sampling Data Sheets

cc: Mr. Thomas J. Berkins, Kaprealian Engineering, Inc.

Table 1
Summary of Monitoring Data

	Ground Water	Depth to	Total Well	Product		Water
Well#	Elevation (feet)	Water (feet)+	Depth (feet)*	Thickness (feet)	Sheen	Purged (galions)
AACH W	(teet)	(ICCI)*	(1001)*	TEST	. PHI VII	(Secontral)
		(Monitored an	d Sampled on O	tober 24, 1996)		
MW-1	71.39	14.85	24.18	0	Yes	18.5
MW-2	70.88	10.78	24.74	0	No	27.5
MW-3	69.90	12.65	23.70	0	No	22
MW-4	70.90	11.14	19.57	0	No	6
MW-5	70.40	11.40	20.00	0	No	4.5
		(Monitored	and Sampled on .	July 24, 1996)		
MW-1	72.09	14.15	24,10	0	No	26
MW-2	71.64	10.02	24.70	0	No	25
MW-3	70.38	12.17	23.68	0	No	30
MW-4	71.57	10.47	19.56	0	No	6.5
MW-5	71.00	10.80	20.00	0	No	6.5
		(Monitored a	and Sampled on A	April 18, 1996)		
MW-1	72.84	13.40	24.20	0	No	23
MW-2	72.39	9.27	24.80	0	No	41
MW-3	71.25	11,30	23,77	0	No	33
MW-4	72.21	9.83	19.61	0	No	7.5
MW-5	72.15	9.65	20.05	0	No	7.5
		(Monitored ar	id Sampled on Ja	muary 18, 1996)		
MW-1	66,97	14.21	24.13	0	No	17
MW-2	66.50	10.11	24.74	0	No	38
MW-3	65.69	11.79	23.71	0	No	31

	Well Casing Elevation	Well Casing Elevation
Well#	(feei)**	(feet)*
MW-1	81.18	86.24
MW-2	76.61	81.66
MW-3	77.48	82.55
MW-4	N/A	82.04
MW-5	N/A	81.80

Table 1 Summary of Monitoring Data

- The depth to water level and total well depth measurements were taken from the top of the well casings.
- * The top of casing elevations were re-surveyed by Kier & Wright in May, 1996, per City of Oakland Benchmark No. 2310, a cut square in concrete curb at mid point of return at the northeast corner of El Dorado and Fairmont Streets (elevation = 77.53 feet MSL). These well casing elevations are used beginning with the April 18, 1996 monitoring event.
- ** The elevations of the top of the well casings, used prior to April 18, 1996, were surveyed by Roux Associates relative to Mean Sea Level (benchmark unknown).

Table 2
Summary of Laboratory Analyses
Water

Well#	Date	TPH as Diesel	TPH as Gasoline	Benzene	Toluene	Ethyl- Benzene	Xylenes	MTBE
17. 14. II		30,70						
MW-1	11/3/92		260,000	2,300	4,600	3,700	17,000	
	1/25/93		120,000	2,100	4,600	4,900	22,000	
	4/29/93		100,000	850	2,000	4,300	19,000	
	7/16/93		29,000	590	560	980	4,200	
	10/19/93		67,000	1,400	2,600	2,900	5,000	
	1/20/94		92,000	1,200	3,000	3,400	17,000	
	4/13/94		51,000	1,000	2,600	3,200	15,000	
	7/13/94		35,000	550	150	1,400	5,700	
	10/10/94		52,000	1,000	810	3,300	12,000	
	1/10/95		810	16	18	59	250	
	4/17/95		48,000	880	530	2,500	11,000	
	7/24/95		48,000	1,500	420	2,700	9,700	
	10/23/95		47,000	780	210	2,100	11,000	270
	1/18/96		30,000	1,500	500	3,500	13,000	2,400
	4/18/96		66,000	2,700	2,200	3,100	13,000	57,000
	7/24/96		5,600	2,100	ND	160	160	24,000
	10/24/96		110,000	7,500	8,000	3,300	14,000	58,000
MW-2	11/3/92		140	2.2	ND	ND	2.0	
	1/25/93		2,100	56	1.1	90	140	
	4/29/93		1,500	290	ND	33	11	
	7/16/93		510*	17	0.60	3.2	2.5	
	10/19/93		670	24	1.1	7.7	23	
	1/20/94		820	97	ND	12	ND	
	4/13/94		550	71	ND	5.1	1.3	
	7/13/94		2,000	490	ND	17	13	
	10/10/94		2,300	340	ND	25	ND	
	1/10/95		850	3.8	ND	8.5	1.3	
	4/17/95		1,300	4.7	ND	8.3	1.2	
	7/24/95		960	20	ND	4.2	6.2	
	10/23/95		ND	ND	ND	ND	ND	19
•	1/18/96		900	300	86	7.6	18	4,300
	4/18/96		18,000	3,600	680	890	4,100	19,000
	7/24/96		100,000	13,000	21,000	2,700	16,000	120,000
	10/24/96		800	110	17	11	20	20,000
MW-3	11/3/92		2,100	120	15	38	200	
	1/25/93		2,300	80	1	55	52	-
	4/29/93		4,500	1,700	ND	200	140	
	7/16/93		4,000*	1,100	28	52	70	
	10/19/93		3,800	42	ND	50	56	
	1/20/94		4,200	11	ND	21	15	
	4/13/94		4,200	210	ND	36	53	

Table 2
Summary of Laboratory Analyses
Water

		TPH as	TPH as			Ethyl-		
Well#	Date	Diesel	Gasoline	Benzene	Tolpene	Benzene	Xylenes	МТВЕ
MW-3	7/13/94		1,800**	16	16	ND	21	
(Cont.)	10/10/94		4,300	11	ND	12	ND	
(соца,)	1/10/95		310	4.6	ND	3.5	2.1	
	4/17/95		7,800	ND	4.6	300	450	
	7/24/95		3,200	170	ND	22	16	
	10/23/95		3,900	55	ND	19	11	4,500
	1/18/96		2,200	270	33	26	18	5,500
	4/18/96		6,000	1,800	ND	100	230	48,000
	7/24/96		ND	2,500	ND	ND	ND	71,000
	10/24/96		3,800	660	ND	15	ND	65,000
MW-4	4/18/96◆	110†	ND	630	ND	ND	ND	18,000
747 44	7/24/96♦♦	ND	ND	ND	ND	ND	5.2	3,900
	10/24/96★	ND	ND	ND	ND	ND	ND	6,300
MW-5	4/18/96		31,000	5,500	1,400	1,700	8,100	66,000
1.1 1. 0	7/24/96		32,000	6,400	ND	1,600	6,100	120,000
	10/24/96		17,000	6,900	ND	970	130	84,000

- * Primarily due to the presence of discrete peaks not indicative of gasoline.
- ** Sequoia Analytical Laboratory reported that the hydrocarbons detected appeared to be a gasoline and non-gasoline mixture.
- † Sequoia Analytical Laboratory reported that the hydrocarbons detected did not appear to contain diesel.
- ♦ Total Oil & Grease and all EPA Method 8010 constituents were non-detectable.
- ♦♦ Total Oil & Grease and all EPA Method 8010 and 8270 constituents were non-detectable.
- ★ Total Oil and Grease and all EPA Method 8010 constituents were non-detectable. All EPA Method 8270 constituents were non-detectable, except for Bis (2-ethylhexyl) phthalate which was detected at a concentration of 14 µg/L.
- -- Indicates analysis was not performed.

ND = Non-detectable.

Results are in micrograms per liter (µg/L), unless otherwise indicated.

Table 2 Summary of Laboratory Analyses Water

Note:

The detection limit for results reported as ND by Sequoia Analytical Laboratory is equal to the stated detection limit times the dilution factor indicated on the laboratory analytical sheets.

Prior to August 1, 1995, the total purgeable petroleum hydrocarbon (TPH as gasoline) quantification range used by Sequoia Analytical Laboratory was C4 - C12. Since August 1, 1995, the quantification range used by Sequoia Analytical Laboratory is C6 - C12.

Laboratory analyses data prior to October 19, 1993, were provided by GeoStrategies, Inc.

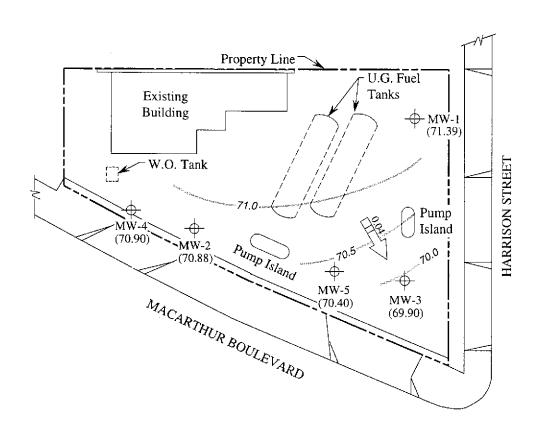


Base modified from 7.5 minute U.S.G.S. Oakland East and West Quadrangles (both photorevised 1980)

0 2000 4000 Approx. scale feet



UNOCAL SERVICE STATION # 1871 96 MACARTHUR BOULEVARD OAKLAND, CALIFORNIA LOCATION MAP



LEGEND

→ Monitoring well

() Ground water elevation in feet above Mean Sea Level

Direction of ground water flow with approximate hydraulic gradient

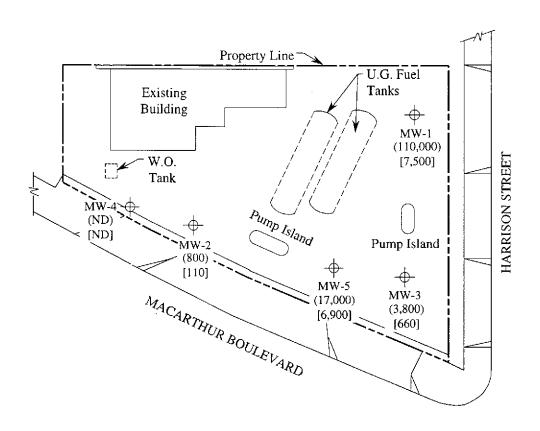
Contours of ground water elevation



GROUND WATER FLOW DIRECTION MAP FOR THE OCTOBER 24, 1996 MONITORING EVENT

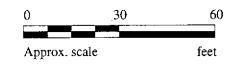


UNOCAL SERVICE STATION # 1871 96 MACARTHUR BOULEVARD OAKLAND, CALIFORNIA FIGURE



LEGEND

- Monitoring well
- () Concentration of TPH as gasoline in μ g/L
- [] Concentration of benzene in μ g/L
- ND Non-detectable



PETROLEUM HYDROCARBON CONCENTRATIONS IN GROUND WATER ON OCTOBER 24, 1996



UNOCAL SERVICE STATION # 1871 96 MACARTHUR BOULEVARD OAKLAND, CALIFORNIA **FIGURE**

2



680 Chesapeake Drive 404 N. Wiget Lane 819 Striker Avenue, Suite 8

Redwood City, CA 94063 Walnut Creek, CA 94598 Sacramento, CA 95834

(415) 364-9600 (510) 988-9600 (916) 921-9600 FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services

2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider

Client Project ID: Matrix Descript:

Unocal #1871, 96 Mac Arthur Blvd. Oakland

Water

Analysis Method: EPA 5030/8015 Mod./8020 First Sample #: 610-1411

Sampled:

Oct 24, 1996

Received: Oct 24, 1996 Reported:

Nov 4, 1996

TOTAL PURGEABLE PETROLEUM HYDROCARBONS with BTEX DISTINCTION

Sample Number	Sample Description	Purgeable Hydrocarbons μg/L	Benzene μg/L	Toluene μg/L	Ethyl Benzene μg/L	Total Xylenes μg/L
610-1411	MW-1	110,000	7,500	8,000	3,300	14,000
610-1412	MW-2	800	110	17	11	20
610-1413	MW-3	3,800	660	ND	15	ND .
610-1414	MW-4	ND	ND	ND	ND	ND
610-1415	MW-5	17,000	6,900	ND	970	130
610-1416	ES-1	ND	ND	ND	ND	ND
610-1417	ES-2	ND	ND	ND	ND	ND

Detection Limits:	50	0.50	0.50	0.50	0.50	
1111						

Total Purgeable Petroleum Hydrocarbons are quantitated against a fresh gasoline standard. Analytes reported as ND were not present above the stated limit of detection.

SEQUOIA ANALYTICAL, #1271

Signature on File







Redwood City, CA 94063 Walnut Creek, CA 94598 Sacramento, CA 95834

(415) 364-9600 (510) 988-9600 (916) 921-9600 FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services 2401 Stanwell Dr., Ste. 300

Concord, CA 94520 Attention: Jarrel Crider Matrix Descript:

First Sample #:

Services Client Project ID: Unocal #1871, 96 Mac Arthur Blvd. Oakland Sampled:

Water

EPA 5030/8015 Mod./8020

Analysis Method: 610-1411

Received:

Reported:

Oct 24, 1996 Oct 24, 1996

Nov 4, 1996

TOTAL PURGEABLE PETROLEUM HYDROCARBONS with BTEX DISTINCTION

Sample Number	Sample Description	Chromatogram Pattern	DL Mult. Factor	Date Analyzed	Instrument ID	Surrogate Recovery, % QC Limits: 70-130
610-1411	MW-1	Gasoline	400	10/31/96	HP-4	94
610-1412	MW-2	Gasoline	10	10/31/96	HP-4	92
610-1413	MW-3	Gasoline	10	10/31/96	HP-4	78
610-1414	MW-4		1.0	10/31/96	HP-4	96
610-1415	MW-5	Gasoline	100	11/2/96	HP-4	91
610-1416	ES-1		1.0	10/31/96	HP-4	97
610-1417	ES-2		1.0	10/31/96	HP-4	94

SEQUOIA ANALYTICAL, #1271

Signature on File



Redwood City, CA 94063 Walnut Creek, CA 94598 Sacramento, CA 95834 (415) 364-9600 (510) 988-9600 (916) 921-9600 FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services 2401 Stanwell Dr., Ste. 300

Concord, CA 94520 Attention: Jarrel Crider Client Project ID: Sample Descript:

Unocal #1871, 96 Mac Arthur Blvd. Oakland Water

MTBE (Modified EPA 8020)

Analysis for: MTBE (M First Sample #: 610-1411

Sampled: Oct 24, 1996 Received: Oct 24, 1996

Analyzed: Oct 31-Nov 2, 96 Reported: Nov 4, 1996

LABORATORY ANALYSIS FOR: MTBE (Modified EPA 8020)

Sample Number	Sample Description	Detection Limit μg/L	Sample Result $\mu \mathrm{g}/\mathrm{L}$
610-1411	MW-1	2,000	58,000
610-1412	MW-2	50	20,000
610-1413	MW-3	50	65,000
610-1414	MW-4	5.0	6,300
610-1415	MW-5	500	84,000

Analytes reported as N.D. were not present above the stated limit of detection.

SEQUOIA ANALYTICAL, #1271

Signature on File





Redwood City, CA 94063 Walnut Creek, CA 94598 Sacramento, CA 95834 (415) 364-9600 (510) 988-9600 (916) 921-9600 FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services

2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider

Chromatogram Pattern:

Client Project ID: Sample Matrix:

: Unocal #1871, 96 Mac Arthur Blvd. Oakland

Water

Analysis Method: EPA 3510/8015 Mod.

First Sample #: 610-1414

Sampled: Oc

Received:

Reported:

Oct 24, 1996 Oct 24, 1996

Nov 4, 1996

TOTAL EXTRACTABLE PETROLEUM HYDROCARBONS

Analyte	Reporting Limit μg/L	Sample I.D. 610-1414 MW-4	
Extractable Hydrocarbons	50	N.D.	

Quality Control Data

Report Limit Multiplication Factor:

1.0

Date Extracted:

10/29/96

Date Analyzed:

10/30/96

Instrument Identification:

HP-3B

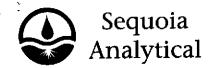
Extractable Hydrocarbons are quantitated against a fresh diesel standard.

Analytes reported as N.D. were not detected above the stated reporting limit.

SEQUOIA ANALYTICAL, #1271

Signature on File





Redwood City, CA 94063 Walnut Creek, CA 94598 Sacramento, CA 95834

(415) 364-9600 (510) 988-9600 (916) 921-9600 FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services 2401 Stanwell Dr., Ste. 300 Concord, CA 94520

Client Project ID: Matrix Descript:

Unocal #1871, 96 Mac Arthur Blvd. Oakland Water

Sampled: Received:

Oct 24, 1996 Oct 24, 1996

Attention: Jarrel Crider

Analysis Method: First Sample #:

SM 5520 B&F (Gravimetric) 610-1414

Extracted: Analyzed: Oct 24, 1996 Oct 25, 1996

Reported:

Nov 4, 1996

TOTAL RECOVERABLE PETROLEUM OIL

Sample Number	Sample Description	Oil & Grease mg/L (ppm)	Detection Limit Multiplication Factor
610-1414	MW-4	N.D.	1.0

Detection Limits:

5.0

Analytes reported as N.D. were not present above the stated limit of detection.

SEQUOIA ANALYTICAL, #1271

Signature on File



680 Chesapeake Drive 404 N. Wiget Lane 819 Striker Avenue, Suite 8

Redwood City, CA 94063 Walnut Creek, CA 94598 Sacramento, CA 95834

(415) 364-9600 (510) 988-9600 (916) 921-9600

FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services 2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider

Client Project ID: Unocal #1871, 96 Mac Arthur Blvd. Oakland Sample Descript: Water, MW-4

Analysis Method: EPA 5030/8010 Lab Number: 610-1414

Sampled: Oct 24, 1996 Oct 24, 1996 Received: Analyzed: Oct 29, 1996 Nov 4, 1996 Reported:

HALOGENATED VOLATILE ORGANICS (EPA 8010)

Analyte	Detection Limit		Sample Results
•	μg/L		µg/L
Bromodichloromethane	1.0		N.D.
Bromoform	1.0		N.D.
Bromomethane	2.0	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	N.D.
Carbon tetrachloride	1.0		N.D.
Chlorobenzene	1.0		N.D.
Chloroethane	2.0		N.D.
2-Chloroethylvinyl ether	2.0		N.D.
Chloroform	1.0		N.D.
Chloromethane	2.0		N.D.
Dibromochloromethane	1.0		N.D.
1,3-Dichlorobenzene	1.0		N.D.
1,4-Dichlorobenzene	1.0		N.D.
1,2-Dichlorobenzene	1.0		N.D.
1,1-Dichloroethane	1.0		N.D.
1,2-Dichloroethane	1.0	***************************************	N.D.
1,1-Dichloroethene	1.0	***************************************	N.D.
cis-1,2-Dichloroethene	1.0	***************************************	N.D.
trans-1,2-Dichloroethene	1.0	***************************************	N.D.
1,2-Dichloropropane	1.0		N.D.
cis-1,3-Dichloropropene	1.0	***************************************	N.D.
trans-1,3-Dichloropropene	1.0		N.D.
Methylene chloride	10	,	N.D.
1,1,2,2-Tetrachloroethane	1.0	***************************************	N.D.
Tetrachloroethene	1.0		N.D.
1,1,1-Trichloroethane	1.0	***************************************	N.D.
1,1,2-Trichloroethane	1.0	***************************************	N.D.
Trichloroethene	1.0		N.D.
Trichlorofluoromethane	1.0		N.D.
Vinyl chloride	2.0	***************************************	N.D.

Analytes reported as N.D. were not present above the stated limit of detection. Because matrix effects and/or other factors required additional sample dilution, detection limits for this sample have been raised.

SEQUOIA ANALYTICAL, #1271

Signature on File





Redwood City, CA 94063 Walnut Creek, CA 94598 Sacramento, CA 95834

(415) 364-9600 (510) 988-9600 (916) 921-9600 FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services 2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider

Client Project ID: Sample Descript: Analysis Method: Lab Number:

Unocal #1871, 96 Mac Arthur Blvd. Oakland Water, MW-4

EPA 8270 610-1414

Oct 24, 1996 Sampled: Received: Oct 24, 1996 Extracted: Oct 29, 1996 Analyzed: Nov 11, 1996

Nov 18, 1996

Reported:

SEMI-VOLATILE ORGANICS by GC/MS (EPA 8270)

Acenaphthylene 2.0 N.D. Acenaphthylene 2.0 N.D. Anlline 2.0 N.D. Anlline 2.0 N.D. Anthracene 2.0 N.D. Benzidine 50 N.D. Benzo(a) Acid 10 N.D. Benzo(a) Acid 10 N.D. Benzo(a) Acid 10 N.D. Benzo(a) Acid 10 N.D. Benzo(b) Liporanthene 2.0 N.D. Benzo(gh) Iporylene 2.0 N.D. Bis(2-chlorostoylene) 2.0 N.D. Bis(2-chlorostoylene) 2.0 N.D. Bis(2-chlorostoylene	Analyte	Detection Limit µg/L		Sample Results μg/L
Aniline	Acenaphthene	2.0	******	
Anthracene. 2.0 N.D. Benzidine. 50 N.D. Benzo(e) Acid 10 N.D. Benzo(b)fluoranthene. 2.0 N.D. Benzo(b)fluoranthene. 2.0 N.D. Benzo(g, h,i)perylene. 2.0 N.D. Benzo(a) pyrene. 2.0 N.D. Benzo(a) pyrene. 2.0 N.D. Benzyl alcohol. 2.0 N.D. Bis(2-chloroethoxy)methane. 2.0 N.D. Bis(2-chloroethyy)methane. 2.0 N.D. Bis(2-chloroethyl)pether. 2.0 N.D. Bis(2-chlorosty)methane. 2.0 N.D. Bis(2-chloropethylpheryl	Acenaphthylene	2.0		
Benzici Acid	Aniline	2.0		
Benzoic Acid	Anthracene	2.0		N.D.
Benzoic Acid. 10 N.D. Benzo(a)anthracene. 2.0 N.D. Benzo(b)fluoranthene. 2.0 N.D. Benzo(g),hi)perylene. 2.0 N.D. Benzo(g),hi)perylene. 2.0 N.D. Benzo(a) pyrene. 2.0 N.D. Benzy alcohol. 2.0 N.D. Bis(2-chloroethoxy)methane. 2.0 N.D. Bis(2-chloroethoxy)methane. 2.0 N.D. Bis(2-chlorostopropyl)ether. 2.0 N.D. Bis(2-chlorostopropyl)ether. 2.0 N.D. Bis(2-chlorostopropyl)ether. 2.0 N.D. 4-Bromophenyl phenyl ether. 2.0 N.D. 4-Bromophenyl phenyl ether. 2.0 N.D. 4-Chloro-3-methylphenol. 2.0 N.D. 4-Chloro-3-methylphenol. 2.0 N.D. 4-Chlorophenol. 2.0 N.D. 4-Chlorophenol. 2.0 N.D. 4-Chlorophenol. 2.0 N.D. Dibenzofuran. 2.0 N.D. D	Benzidine	50		N.D.
Benzo(b)fluoranthene. 2.0 N.D.		10	***************************************	N.D.
Benzo(b)fluoranthene. 2.0 N.D.	Benzo(a)anthracene	2.0	***************************************	N.D.
Benzo(k)fluoranthene. 2.0 N.D. Benzo(g,h,l)perylene. 2.0 N.D. Benzo(a)pyrene. 2.0 N.D. Benzyl alcohol. 2.0 N.D. Bis(2-chloroethoxy)methane. 2.0 N.D. Bis(2-chloroethyl)ether 2.0 N.D. Bis(2-chloroisopropyl)ether. 2.0 N.D. Bis(2-chloroisopropyl)ether. 2.0 N.D. 4-Bromophenyl phenyl ether. 2.0 N.D. 4-Bromophenyl phenyl ether. 2.0 N.D. 4-Chloroanlline. 2.0 N.D. 4-Chloroaphthalene. 2.0 N.D. 4-Chloroaphthalene. 2.0 N.D. 4-Chloro-3-methylphenol. 2.0 N.D. 2-Chloronaphthalene. 2.0 N.D. 4-Chloro-3-methylphenol. 2.0 N.D. 2-Chlorophenol. 2.0 N.D. 4-Chlorophenyl phenyl ether. 2.0 N.D. 10-rophenyl phenyl ether. 2.0 N.D. 2-Chloronaphthalene. 2.0 N.D.	Benzo(b)fluoranthene	2.0		N.D.
Benzo(g,h,i)perylene. 2.0 N.D. Benzo(a)pyrene. 2.0 N.D. Berzy alcohol. 2.0 N.D. Bis(2-chloroethoxy)methane. 2.0 N.D. Bis(2-chloroethyl)ether. 2.0 N.D. Bis(2-chloroethyl)ether. 2.0 N.D. Bis(2-chlorospoyl)ether. 2.0 N.D. Bis(2-chlyflexyl)phthalate. 10 14 4-Bromophenyl phenyl ether. 2.0 N.D. Butyl benzyl phthalate. 2.0 N.D. 4-Chloroanlline. 2.0 N.D. 2-Chloropaphthalene. 2.0 N.D. 4-Chloroa-methylphenol. 2.0 N.D. 2-Chlorophenol. 2.0 N.D. 4-Chlorophenyl phenyl ether. 2.0 N.D. 4-Chlorophenyl phenyl ether. 2.0 N.D. 2-Chlorophenyl phenyl ether. 2.0 N.D. Dibenz(a,h)anthracene. 2.0 N.D. Dibenz(a)ranthracene. 2.0 N.D. Dibenz(a)ranthracene. 2.0 N.D.	• •	2.0		N.D.
Benzy (a) pyrene. 2.0 N.D. Benzy (a) (a) (b) (c) (c) (c) (c) (c) (c) (c) (c) (c) (c	• /	2.0	*********	N.D.
Benzyl alcohol 2.0 N.D.	-	2.0	***************************************	N.D.
Bis(2-chloroethoxy)methane. 2.0 N.D.			***************************************	N.D.
Bis(2-chloroethyl)ether. 2.0 N.D.				N.D.
Bis(2-chloroisopropyl)ether		2.0		N.D.
Bis(2-ethylnexyl)phthalate. 10				N.D.
A-Bromophenyl phenyl ether. 2.0 N.D.				. 14
Butyl benzyl phthalate 2.0 N.D. 4-Chloroanliine 2.0 N.D. 2-Chloronaphthalene 2.0 N.D. 4-Chloro-3-methylphenol 2.0 N.D. 2-Chlorophenol 2.0 N.D. 4-Chlorophenyl phenyl ether 2.0 N.D. 4-Chlorophenyl phenyl ether 2.0 N.D. Chrysene 2.0 N.D. Dibenz(a, h)anthracene 2.0 N.D. Dibenzofuran 2.0 N.D. Di-N-butyl phthalate 10 N.D. 1,3-Dichlorobenzene 2.0 N.D. 1,3-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzene 2.0 N.D. 2,4-Dichlorobenzidine 10 N.D. 2,4-Dichlorobenzidine 2.0 N.D. 2,4-Dinitroluene 2.0 N.D. 2,4-Dimethylphenol 2.0 N.D. 2,4-Dimitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0		2.0	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	N.D.
4-Chloroanilline 2.0 N.D. 2-Chloronaphthalene 2.0 N.D. 4-Chloro-3-methylphenol 2.0 N.D. 2-Chlorophenol 2.0 N.D. 4-Chlorophenyl phenyl ether 2.0 N.D. Chrysene 2.0 N.D. Dibenz(a, h)anthracene 2.0 N.D. Dibenzofuran 2.0 N.D. Di-N-butyl phthalate 10 N.D. 1,3-Dichlorobenzene 2.0 N.D. 1,4-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzene 2.0 N.D. 3,3-Dichlorobenzidine 10 N.D. 2,4-Dichlorophenol 2.0 N.D. 2,4-Dichlorophenol 2.0 N.D. 2,4-Dimethyl phthalate 2.0 N.D. 2,4-Dimethyl phthalate 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.				N.D.
2-Chloronaphthalene. 2.0 N.D. 4-Chloro-3-methylphenol. 2.0 N.D. 2-Chlorophenol. 2.0 N.D. 4-Chlorophenyl phenyl ether. 2.0 N.D. Chrysene. 2.0 N.D. Dibenz(a,h)anthracene. 2.0 N.D. Dibenzofuran. 2.0 N.D. Di-N-butyl phthalate. 10 N.D. 1,3-Dichlorobenzene. 2.0 N.D. 1,4-Dichlorobenzene. 2.0 N.D. 1,2-Dichlorobenzene. 2.0 N.D. 3,3-Dichlorobenzidine. 10 N.D. 2,4-Dichlorophenol. 2.0 N.D. Diethyl phthalate. 2.0 N.D. 2,4-Dimethyl phthalate. 2.0 N.D. 2,4-Dimitro-2-methylphenol. 10 N.D. 2,4-Dinitro-2-methylphenol. 10 N.D. 2,4-Dinitro-toluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D.				N.D.
4-Chloro-3-methylphenol 2.0 N.D. 2-Chlorophenol 2.0 N.D. 4-Chlorophenyl phenyl ether 2.0 N.D. Chrysene 2.0 N.D. Dibenz(a,h)anthracene 2.0 N.D. Dibenzofuran 2.0 N.D. Di-N-butyl phthalate 10 N.D. 1,3-Dichlorobenzene 2.0 N.D. 1,4-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzidine 10 N.D. 2,4-Dichlorophenol 2.0 N.D. 2,4-Dimethyl phthalate 2.0 N.D. 2,4-Dimethyl phthalate 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.				N.D.
2-Chlorophenol 2.0 N.D. 4-Chlorophenyl phenyl ether 2.0 N.D. Chrysene 2.0 N.D. Dibenz(a,h)anthracene 2.0 N.D. Dibenzofuran 2.0 N.D. Di-N-butyl phthalate 10 N.D. 1,3-Dichlorobenzene 2.0 N.D. 1,4-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzene 2.0 N.D. 3,3-Dichlorobenzidine 10 N.D. 2,4-Dichlorophenol 2.0 N.D. Diethyl phthalate 2.0 N.D. 2,4-Dimethylphenol 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.				N.D.
4-Chlorophenyl phenyl ether. 2.0 N.D. Chrysene. 2.0 N.D. Dibenzofuran. 2.0 N.D. Di-N-butyl phthalate. 10 N.D. 1,3-Dichlorobenzene. 2.0 N.D. 1,4-Dichlorobenzene. 2.0 N.D. 1,2-Dichlorobenzene. 2.0 N.D. 3,3-Dichlorobenzidine. 10 N.D. 2,4-Dichlorophenol. 2.0 N.D. 2,4-Dimethylphenol. 2.0 N.D. 2,4-Dimethylphenol. 2.0 N.D. 4,6-Dinitro-2-methylphenol. 10 N.D. 2,4-Dinitrophenol. 10 N.D. 2,4-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D.				N.D.
Chrysene. 2.0 N.D. Dibenz(a,h)anthracene. 2.0 N.D. Dibenzofuran. 2.0 N.D. Di-N-butyl phthalate. 10 N.D. 1,3-Dichlorobenzene. 2.0 N.D. 1,4-Dichlorobenzene. 2.0 N.D. 1,2-Dichlorobenzidine. 10 N.D. 3,3-Dichlorobenzidine. 10 N.D. 2,4-Dichlorophenol. 2.0 N.D. 2,4-Dimethylphthalate. 2.0 N.D. 2,4-Dimethylphenol. 2.0 N.D. 2,4-Dinitro-2-methylphenol. 10 N.D. 2,4-Dinitro-2-methylphenol. 10 N.D. 2,4-Dinitrophenol. 10 N.D. 2,4-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D.				
Dibenz(a,h)anthracene. 2.0 N.D. Dibenzofuran. 2.0 N.D. Di-N-butyl phthalate. 10 N.D. 1,3-Dichlorobenzene. 2.0 N.D. 1,4-Dichlorobenzene. 2.0 N.D. 1,2-Dichlorobenzene. 2.0 N.D. 3,3-Dichlorobenzidine. 10 N.D. 2,4-Dichlorophenol. 2.0 N.D. Diethyl phthalate. 2.0 N.D. 2,4-Dimethylphenol. 2.0 N.D. Dimethyl phthalate. 2.0 N.D. 4,6-Dinitro-2-methylphenol. 10 N.D. 2,4-Dinitrophenol. 10 N.D. 2,4-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D.				N.D.
Dibenzofuran 2.0 N.D. Di-N-butyl phthalate 10 N.D. 1,3-Dichlorobenzene 2.0 N.D. 1,4-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzene 2.0 N.D. 3,3-Dichlorobenzidine 10 N.D. 2,4-Dichlorophenol 2.0 N.D. 2,4-Dinthyl phthalate 2.0 N.D. 2,4-Dimethylphenol 2.0 N.D. Dimethyl phthalate 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.	•			N.D.
Di-N-butyl phthalate 10 N.D. 1,3-Dichlorobenzene 2.0 N.D. 1,4-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzidine 2.0 N.D. 3,3-Dichlorobenzidine 10 N.D. 2,4-Dichlorophenol 2.0 N.D. Diethyl phthalate 2.0 N.D. 2,4-Dimethylphenol 2.0 N.D. Dimethyl phthalate 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.				N.D.
1,3-Dichlorobenzene 2.0 N.D. 1,4-Dichlorobenzene 2.0 N.D. 1,2-Dichlorobenzene 2.0 N.D. 3,3-Dichlorobenzidine 10 N.D. 2,4-Dichlorophenol 2.0 N.D. Diethyl phthalate 2.0 N.D. 2,4-Dimethylphenol 2.0 N.D. Dimethyl phthalate 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.				
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1,2-Dichlorobenzene	·			
3,3-Dichlorobenzidine 10 N.D. 2,4-Dichlorophenol 2.0 N.D. Diethyl phthalate 2.0 N.D. 2,4-Dimethylphenol 2.0 N.D. Dimethyl phthalate 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.	· ·		•	
2,4-Dichlorophenol. 2.0 N.D. Diethyl phthalate. 2.0 N.D. 2,4-Dimethylphenol. 2.0 N.D. Dimethyl phthalate. 2.0 N.D. 4,6-Dinitro-2-methylphenol. 10 N.D. 2,4-Dinitrophenol. 10 N.D. 2,4-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D. 2,6-Dinitrotoluene. 2.0 N.D.	·			
Diethyl phthalate	·		• • •	
2,4-Dimethylphenol 2.0 N.D. Dimethyl phthalate 2.0 N.D. 4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.	·		***************************************	
Dimethyl phthalate				
4,6-Dinitro-2-methylphenol 10 N.D. 2,4-Dinitrophenol 10 N.D. 2,4-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D. 2,6-Dinitrotoluene 2.0 N.D.			·	
2,4-Dinitrophenol				
2,4-Dinitrotoluene				
2,6-Dinitrotoluene 2.0 N.D.				
	,			
Di-N-octyl phthalate		2.0		N.D.



680 Chesapeake Drive 404 N. Wiget Lane

Redwood City, CA 94063 Walnut Creek, CA 94598 819 Striker Avenue, Suite 8 Sacramento, CA 95834

(415) 364-9600 (510) 988-9600 (916) 921-9600 FAX (415) 364-9233 FAX (510) 988-9673 FAX (916) 921-0100

MPDS Services 2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider

Client Project ID: Sample Descript: Analysis Method:

Lab Number:

Unocal #1871, 96 Mac Arthur Blvd. Oakland Water, MW-4

EPA 8270 610-1414

Sampled: Oct 24, 1996 Oct 24, 1996 Received: Extracted: Oct 29, 1996 Nov 11, 1996 Analyzed: Nov 18, 1996 Reported:

SEMI-VOLATILE ORGANICS by GC/MS (EPA 8270)

Analyte	Detection Limit		Sample Results
	μg/L		μg/L
Fluoranthene	2.0		N.D.
Fluorene	2.0		N.D.
Hexachlorobenzene	2.0		Ň.D.
Hexachlorobutadiene.	2.0		N.D.
Hexachlorocyclopentadiene	2.0		N.D.
Hexachloroethane	2.0		N.D.
Indeno(1,2,3-cd)pyrene	2.0		N.D.
Isophorone	2.0		N.D.
2-Methylnaphthalene	2.0		N.D.
2-Methylphenol	2.0		N.D.
4-Methylphenol	2.0		N.D.
Naphthalene	2.0		N.D.
2-Nitroaniline	10		N.D.
3-Nitroaniline	10		N.D.
4-Nitroaniline	10	:	N.D.
Nitrobenzene	2.0		N.D.
2-Nitrophenol	2.0		N.D.
4-Nitrophenol	10		N.D.
N-Nitrosodimethylamine	2.0		N.D.
N-Nitrosodiphenylamine	2.0		N.D.
N-Nitroso-di-N-propylamine	2.0		N.D.
Pentachlorophenol	10		N.D.
Phenanthrene	2.0		N.D.
Phenol	2.0		N.D.
Pyrene	2.0		N.D.
1,2,4-Trichlorobenzene	2.0		N.D.
2,4,5-Trichlorophenol	10		N.D.
2,4,6-Trichlorophenol	2.0		N.D.

Analytes reported as N.D. were not present above the stated limit of detection.

SEQUOIA ANALYTICAL, #1271

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MPDS Services

2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider Client Project ID:

Matrix:

Unocal #1871, 96 Mac Arthur Blvd. Oakland

Liquid

QC Sample Group: 6101411-417

Reported:

Nov 18, 1996

QUALITY CONTROL DATA REPORT

ANALYTE	Benzene	Toluene	Ethyl	Xylenes	Diesel	Oil &
			Benzene		•	Grease
Method:	EPA 8020	EPA 8020	EPA 8020	EPA 8020	EPA 8015	SM 5520
Analyst:	D. Newcomb	D. Newcomb	D. Newcomb	D. Newcomb	D. Sharma	D. Newcomb
MS/MSD		•				•
Batch#:	6101133	6101133	6101133	6101133	BLK102996	BLK102496
Date Prepared:	10/31/96	10/31/96	10/31/96	10/31/96	10/29/96	10/24/96
Date Analyzed:	10/31/96	10/31/96	10/31/96	10/31/96	10/30/96	10/25/96
nstrument l.D.#:	HP-4	HP-4	HP-4	HP-4	HP-3B	Manual
Conc. Spiked:	20 μg/L	20 μg/L	$20\mu\mathrm{g/L}$	60 μg/L	300 μg/L	100 mg/L
Matrix Spike % Recovery:	100	80	80	83	93	84
Matrix Spike Duplicate % Recovery:	100	. 85	85	85	93	94
Relative % Difference:	0.0	6.1	6.1	2.0	0.0	11
LCS Batch#:	4LCS103196	4LCS103196	4LCS103196	4LCS103196	LCS102996	BLK102496
Date Prepared:	10/31/96	10/31/96	10/31/96	10/31/96	10/29/96	10/24/96
Date Analyzed:	10/31/96	10/31/96	10/31/96	10/31/96	10/30/96	10/25/96
nstrument l.D.#:	HP-4	HP-4	HP-4	HP-4	HP-3B	Manual
LCS %						
Recovery:	90	95	100	103	97	84
% Recovery		.		· · ·		
Control Limits:	60-140	60-140	60-140	60-140	60-140	60-140

The

SEQUOIA ANALYTICAL, #1271

Signature on File

Alan B. Kemp Project Manager Please Note:

The LCS is a control sample of known, interferent free matrix that is analyzed using the same reagents, preparation, and analytical methods employed for the samples. The matrix spike is an aliquot of sample fortified with known quantities of specific compounds and subjected to the entire analytical procedure. If the recovery of analytes from the matrix spike does not fall within specified control limits due to matrix interference, the LCS recovery is to be used to validate the batch.





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MPDS Services

2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider Client Project ID:

Unocal #1871, 96 Mac Arthur Blvd. Oakland

Matrix: Liquid

QC Sample Group: 6101411-417

Reported: Nov 18, 1996

QUALITY CONTROL DATA REPORT

ANALYTE	Benzene	Toluene	Ethyl	Xylenes	
			Benzene		
Method:	EPA 8020	EPA 8020	EPA 8020	EPA 8020	
Analyst:	D. Newcomb	D. Newcomb	D. Newcomb	D. Newcomb	
MS/MSD					
Batch#:	6110196	6110196	6110196	6110196	
_					
Date Prepared:	11/1/96	11/1/96	11/1/96	11/1/96	
Date Analyzed:	11/1/96	11/1/96	11/1/96	11/1/96	
nstrument I.D.#:	HP-4	HP-4	HP-4	HP-4	
Conc. Spiked:	20 μg/L	20 μg/L	$20\mu \mathrm{g/L}$	60 μg/L	
Matrix Spike					
% Recovery:	125	95	90	93	
Matrix Spike					
Duplicate %					
Recovery:	125	95	90	92	
Difference:	0.0	0.0	0.0	1.8	
Relative % Difference:	0.0	0.0	0.0	1.8	

LCS Batch#:	4LCS110196	4LCS110196	4LCS110196	4LCS110196
Date Prepared:	11/1/96	11/1/96	11/1/96	11/1/96
Date Analyzed:	11/1/96	11/1/96	11/1/96	11/1/96
Instrument I.D.#:	HP-4	HP-4	HP-4	HP-4
LCS %				
Recovery:	125	95	90	93
% Recovery				
Control Limits:	60-140	60-140	60-140	60-140

SEQUOIA ANALYTICAL, #1271

Signature on File

Alan B. Kemp Project Manager Please Note:

The LCS is a control sample of known, interferent free matrix that is analyzed using the same reagents, preparation, and analytical methods employed for the samples. The matrix spike is an aliquot of sample fortified with known quantities of specific compounds and subjected to the entire analytical procedure. If the recovery of analytes from the matrix spike does not fall within specified control limits due to matrix interference, the LCS recovery is to be used to validate the batch.





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MPDS Services

2401 Stanwell Dr., Ste. 300 Concord, CA 94520 Attention: Jarrel Crider Client Project ID:

Unocal #1871, 96 Mac Arthur Blvd. Oakland

Matrix:

QC Sample Group: 6101411-417

Reported:

Nov 18, 1996

QUALITY CONTROL DATA REPORT

Liquid

ANALYTE	1,1-Dichloro-	Trichloro-	Chlaro-	
	ethene	ethene	benzene	
Method:	EPA 8010	EPA 8010	EPA 8010	
Analyst:	P. Horton	P. Horton	P. Horton	
MS/MSD				
Batch#:	6101463	6101463	6101463	
Date Prepared:	10/28/96	10/28/96	10/28/96	
Date Analyzed:	10/28/96	10/28/96	10/28/96	
Instrument I.D.#:	HP-7	HP-7	HP-7	
Conc. Spiked:	10 μg/L	10 μg/L	10 μg/L	
Matrix Spike				
% Recovery:	65	97	85	
Matrix Spike				
Duplicate % Recovery:	74	94	91	
1130010191	, ,	31	5 1	
Relative %				
Difference:	13	3.1	6.8	

LCS Batch#:	LCS102896	LCS102896	LCS102896
Date Prepared:	10/28/96	10/28/96	10/28/96
Date Analyzed:	10/28/96	10/28/96	10/28/96
Instrument I.D.#:	HP-7	HP-7	HP-7
LCS % Recovery:	83	122	90
% Recovery Control Limits:	60-140	60-140	60-140

SEQUOIA ANALYTICAL, #1271

Signature on File

Alan B. Kemp Project Manager Please Note:

The LCS is a control sample of known, interferent free matrix that is analyzed using the same reagents, preparation, and analytical methods employed for the samples. The matrix spike is an aliquot of sample fortified with known quantities of specific compounds and subjected to the entire analytical procedure. If the recovery of analytes from the matrix spike does not fall within specified control limits due to matrix interference, the LCS recovery is to be used to validate the batch.





Tel: (510) 602-5100, Fax: (510) 689-1918

CHAIN OF CUSTODY

ANALYSES REQUESTED UNOCAL SAMPLER TURN AROUND TIME: S/S # 187/ CITY: OAKLAND TPH-GAS BTEX STEVE BALIAN 0 5_DAYS 仪 TPH-DIESEL ADDRESS: 96 MAC ARTHUR RIVY ATBI WITNESSING AGENCY d TOG 8010 REMARKS SAMPLING LOCATION COMP NO. OF CONT. DATE TIME WATER GRAB SAMPLE ID NO. 61014116 10-24-96 13:40 WELL REGULAR 11 13:22 11 8010 \searrow 6101413 TIME 13:25 6101411 AG X Χ 10:20 6101415 X 14:10 // THE FOLLOWING MUST BE COMPLETED BY THE LABORATORY ACCEPTING SAMPLES FOR ANALYSES: DATE/TIME RECEIVED BY: DATE/TIME RELINQUISHED BY: 15:50 1. HAVE ALL SAMPLES RECEIVED FOR ANALYSIS BEEN STORED ON ICE? STEVE BALIAN 2. WILL SAMPLES REMAIN REFRIGERATED UNTIL ANALYZED? (SIGNATUBE) (SIGNATURE) (SIGNATURE) 4. WERE SAMPLES IN APPROPRIATE CONTAINERS AND PROPERLY PACKAGED? (SIGNATURE) (SIGNATURE) SIGNATURE: 10/2× (SIGNATURE) (SIGNATURE) 1550

Note: All water containers to be sampled for TPHG/BTEX, 8010 & 8240 are preserved with HCL. All water containers to be sampled for Lead or Metals are preserved with HN03. All other containers are unpreserved.



CHAIN OF CUSTODY

SAMPLER			UNOC	CAL # <u>/8</u>	7/	_ CITY: <u>0AK(A</u>	ם מ			ĀN	ALYSES	REQUES	l'ED	1		TURN AROUND TIME:
STEVE]						MAC ARTHUR	AIUJ.	甲田田	TPH- DIESEL	<u>ت</u>	0,					REGULAR
SAMPLE ID NO.	DATE	TIME	WATER	GRAB	сомр	NO. OF CONT.	SAMPLING	HH	TP	TOG	8010					REMARKS
ESI	10-24-96		X	X		1		X					<u> </u>	-		-
ES2	"		X	X		1		X						-		-
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RELINQUIS	SHED BY:	DATE/1	TIME] F	RECEIVED BY:	D	ATE/TIME	THE FO	DLLOWING	MUST BE	COMPLETE	D BY THE I	LABORATO	RY ACCEP	TING SAMPLES FOR ANALYSES:
	. ,	/5:°							1. HAVE	ALL SAMI	PLES RECEIV	VED FOR A	NALYSIS E	BEEN STOR	ED ON ICE	·
STEVE B	AUAN	10-20	1-16	ISIGN	ATURE				2. WILL	SAMPLES	remain re	FRIGERATI	ED UNTIL A	NALYZED	·	
(SIGNATURE)				ISIGN	ATURE	:)		·····	3. DID A	NY SAMPI	£\$ RECEIVI	ED FOR AN	IALYSIS H	AVE HEAD	SPACE? _	
(SIGNATURE)				ISIGN	BRUTAI	1					IN APPROF	PRIATE CO	NTAINERS			
(SIGNATURE)			-	ISIGN	ĀTURI	M		16/1×	SIGNAT	TURE:	184	1		A.	TLE: MYST	DATE:

Note: All water containers to be sampled for TPHG/BTEX, 8010 & 8240 are preserved with HCL. All water containers to be sampled for Lead or Metals are preserved with HN03. All other containers are unpreserved.

MRDS Services Inc.

2401 Stanwell Drive Concord, California 94520 Tel: (510) 602-5120 Fax: (510) 689-1918

SAMPLING LOCATION: # 187/ OAKIAND	DATE & A.M. TIME SAMPLED 10-24-96 13:40 P.M.
96 MAC ARTHUR BILLS	FIELD TECHNICIAN STEUR BALIAM
PURGE METHOD PUMP	DATE(S) PURGED _/0 _ 24-96
WELL NUMBERMw_/	
WATER LEVEL-INITIAL /4.85	SAMPLING METHOD BALL
WATER LEVEL-FINAL	CONTAINERS 4
WELL DEPTH 24.18	PRESERVATIVES Hol
WELL CASING VOLUME 6.06	tCASING DIAMETER 4

TIME	GALLONS PURGED	TEMPERATURE (°F) (± 1°F)	ELECTRICAL CONDUCTIVITY ([\mu mhos/cm]x100) (± 10% of TOTAL	pH (± 0.2)
//:00	Ð	76.6	565 US	6.64
	6	76.1	533 US	6.55
	12	76.0	537 01	6.53
	18.5 DE	WATERED		
12:05	18.5	75.8	544 01	6.57
		·		

† Correction Factors:	Well Diameter 2" 3" 4" 4.5"	Factor 0.17 0.37 0.65 0.82
	4"	0.65
	•	
	6"	1.46
	8"	2.6
	12"	5.87

◆240'1 Stanwell Drive Concord, California 94520 Tel: (510) 602-5120 Fax: (510) 689-1918

SAMPLING LOCATION: # 187/ OAK(AND	DATE & TIME SAMPLED 10-24-96 13:55 P.M.
96 MAR ARTHUR Blue	FIELD TECHNICIAN STEVE BALIAN
PURGE METHOD PUMP	DATE(S) PURGED _/0 - 24 - 96
WELL NUMBER MW-2	
WATER LEVEL-INITIAL 10.78	SAMPLING METHOD BAIL
WATER LEVEL-FINAL 11.69	CONTAINERS 4
WELL DEPTH	PRESERVATIVES Hel
WELL CASING VOLUME9.07	tCASING DIAMETER 4"

TIME	GALLONS PURGED	TEMPERATURE (°F) (± 1°F)	ELECTRICAL CONDUCTIVITY ([µmhos/cm]x100) (± 10% of TOTAL	pH (± 0.2)
12:15	0	73.4	437 US	6.72
	9	74./	407 US	6.79
	18	73.6	407 US	6.77
	DE	UATERED		
12:45	27.5	73.9	40/ US	6.73

Correction Factors:	Well Diameter 2" 3" 4" 4.5"	Factor 0.17 0.37 0.65 0.82
	4.5 " 6"	0.82 1.46
	8"	2.6
	12"	5.87

• 2401 Stanwell Drive Concord, California 94520

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SAMPLING LOCATION:#1871 DAKLAND	DATE & A.M. TIME SAMPLED 10-24-96 13:25 P.M.
96 MC ARTHUR BIVOL	FIELD TECHNICIAN STEVE BALIAN
PURGE METHOD PUMP	DATE(S) PURGED 10-24-96
WELL NUMBER MW-3	
WATER LEVEL-INITIAL 12.65	SAMPLING METHOD BAIL
WATER LEVEL-FINAL 14.71	CONTAINERS 4
WELL DEPTH 23.70	PRESERVATIVES Hel
	tCASING DIAMETER 4"

TIME	GALLONS PURGED	TEMPERATURE (°F) (± 1°F)	ELECTRICAL CONDUCTIVITY ([µmhos/cm]x100) (± 10% of TOTAL	pH (± 0.2)
10:30	О	72.9	525 US	6.56
	7.5	75.2	563 US	6.56
	15	75.6	538 US	6.58
10:50	22	75.8	531 05	6.55

† Correction Factors:	Well Diameter	<u>Factor</u>
	2"	0.17
	3"	0.37
	4"	0.65
	4.5"	0.82
	6"	1.46
	8"	2.6
	12"	5.87

WPDS Services Inc. 2401 Stanwell Drive Concord, California 94520

Tel: (510) 602-5120 Fax: (510) 689-1918

SAMPLING LOCATION: # /87/ OAKLAND	DATE & A.M. TIME SAMPLED (0-24-96 10:20 P.M.
96 MAC ARTHUR BIUG	FIELD TECHNICIAN STEVE BALIAGE
PURGE METHOD PUMP	DATE(S) PURGED _ 10 - 24 - 96
WELL NUMBER MW-4	
WATER LEVEL-INITIAL //./4	SAMPLING METHOD <u>BAIL</u>
WATER LEVEL-FINAL //. 23	containers 9
WELL DEPTH	PRESERVATIVES Hel
WELL CASING VOLUME /. 43	tCASING DIAMETER _2"

TIME	GALLONS PURGED	TEMPERATURE (°F) (± 1°F)	ELECTRICAL CONDUCTIVITY ([µmhos/cm]x100) (± 10% of TOTAL		pH (± 0.2)
9:55	D	66.3	588	٥.>	7.88
	1.5	68.0	452	υς	7.50
	3	69.8	420	رں	7.23
	4.5	70.2	399	US	7.00
10:05	6	70.3	408	(L	6.95
<u></u>					

t	Correction Factors:	Well Diameter	<u>Factor</u>
		2"	0.17
		3"	0.37
		4"	0.65
		4.5"	0.82
		6"	1.46
		8"	2.6
		12"	5.87

MPÓS Services Inc. 2401 Stanwell Drive Concord, California 94520

Tel: (510) 602-5120 Fax: (510) 689-1918

SAMPLING LOCATION: #1871 OAK(AND	DATE & A.M. TIME SAMPLED 10-24-96 14:10 (P.M.)
96 MAC ARTHUR BILLY.	FIELD TECHNICIAN STEVE BALLAN
PURGE METHOD PUMP	DATE(S) PURGED 10-24-96
WELL NUMBER <u>Mw_5</u>	
WATER LEVEL-INITIAL //. 40	SAMPLING METHOD BAIL
WATER LEVEL-FINAL 43.07	CONTAINERS 4
WELL DEPTH	PRESERVATIVES Hel
WELL CASING VOLUME/. 46	tCASING DIAMETER

TIME	GALLONS PURGED	TEMPERATURE (°F) (± 1°F)	ELECTRICAL CONDUCTIVITY ([µmhos/cm]x100) (± 10% of TOTAL	pH (± 0.2)
/3:00	٥	75.2	484 US	6.79
	1.5	75.8	465 US	6.79
	3	78.3	493 US	6.75
13:05	4.5	75.1	499 US	6.74

t	Correction Factors:	Well Diameter	<u>Factor</u>
		2"	0.17
		3"	0.37
		4"	0.65
		4.5"	0.82
		6"	1.46
		8"	2.6
		12"	5.87