Harding Lawson Associates



June 16, 1992

I.P. No. 16419

Alameda County Department of Environmental Health 80 Swan Way, Room 210 Oakland, California 94621

Attention:

Ms. Susan Hugo

California Regional Water Quality Control Board San Francisco Bay Region 2101 Webster Street, Suite 500 Oakland, California 94612

Attention:

Mr. Rich Hiett

Dear Ms. Hugo and Mr. Hiett:

Response to Request for Work Plan Hydrocarbon Release from Underground Storage Tanks Chinatown Redevelopment Project Area Oakland, California

This letter report has been prepared by Harding Lawson Associates (HLA) on behalf of the Redevelopment Agency of the City of Oakland (Agency). It responds to a letter dated January 13, 1992, from the Alameda County Department of Environmental Health (ACDEH) to Pacific Renaissance Associates II (PRA II) and the City of Oakland (City). The letter requested, among other items, submittal of a work plan to characterize the extent of soil contamination associated with the release of petroleum hydrocarbons from underground storage tanks (USTs) near the corner of Franklin Street and the former location of 10th Street, Oakland (Plate 1) in the Chinatown Redevelopment Project Area (Project Area), and to investigate whether groundwater has been affected by releases from the USTs. a contents? undersom

BACKGROUND

During excavation for installation of service ducts as part of construction at the Pacific Renaissance Plaza (PRP) site, two USTs were discovered in the sidewalk area adjacent to the site, on the east side of Franklin Street and north of the former location of 10th Street. The tanks were approximately 10 feet long by 6 feet in diameter and buried approximately 6 to 12 feet below ground surface (bgs).

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After removal of the tanks, three samples were collected, one of stockpiled soil from above the tanks, and one from below the former location of each tank. Samples were analyzed by EPA Test Methods 3550/5030, 8020, 8240, and 8270, for hazardous waste characteristics, and for cyanide, cadmium, chromium, lead, nickel, and zinc. Analyses 400 ppm 0+6 were performed by NET Pacific, Inc., a California-certified laboratory in Santa Rosa, California. Sample results indicated the presence of petroleum hydrocarbons in the gasoline and diesel ranges and trace levels of the chlorinated hydrocarbons trichloroethene (TCE) and 1,1,1-trichloroethane (1,1,1-TCA). Results, presented in NET Pacific's laboratory report dated December 27, 1991, are attached to this letter (Attachment A). 🗁

At the request of ACDEH staff and as discussed at a December 24, 1991, meeting among representatives of ACDEH, the California Regional Water Quality Control Board, San Francisco Bay Region (RWQCB), PRA II, and HLA, additional samples were collected at the base of the duct excavation, approximately 20 feet bgs. Four soil samples were collected in each of four areas and were composited by NET into four samples which were analyzed by EPA Test Methods 3550/5030, 8010, 8020, and 8270. Results indicated nondetectable or low (<2 milligrams/kilogram [mg/kg]) concentrations of petroleum hydrocarbons in three of the four composites. The 4. A tomposite from the northernmost area indicated the presence of petroleum hydrocarbons as gasoline and diesel at concentrations of 420 and 52 mg/kg.

respectively. No chlorinated hydrocarbons were detected in the samples. Laboratory analytical results are presented in NET's laboratory report dated January 22, 1992, and and S. 330 ppm benzene 900 ppm naphalene (.910 ppm 2-nitroaniline are included in this letter as Attachment B.

OBJECTIVES AND SCOPE OF THIS REPORT

The overall objective of this report is to respond to the request for a work plan presented in Item 3) of a letter dated January 13, 1992, from ACDEH to PRA II and the City. In discussions with Paul Smith of ACDEH, it was agreed that an acceptable response to the request could consist of a plan that does not propose additional soil borings or monitoring well installations, with the rationale to support such a plan. This approach has been followed in this report. The scope of this report includes the following items:

- Description of the known extent of petroleum hydrocarbons in soil and groundwater near the former location of the USTs;
- Discussion of groundwater flow directions in the Project Area in general and 0 near the former tank locations in particular; and
- Presentation of recommendations for additional monitoring in the Project Area 0 and a rationale indicating that the existing characterization of the extent of hydrocarbons is sufficient.

bottom.

19 000 ppm gas 4, 100 ppm chisel 53 ppm benzene 6.3 ppm rapthalone 2.1 ppm 2-methyl "

.028 ppm 111 fcA .050ppm TCE

sampled 12-30-91

It was agreed that if HLA's contention was that mws were not needed they Should present then in a formal document supporting Aon. The rationale as to why .

Discussions of the hydrocarbon extent and groundwater flow direction draw on:

- Observations and analytical results from excavation and shoring stages of construction of the East Bay Municipal Utility District (EBMUD) administration building and the adjacent PRP building:
- Observations and analytical results from excavation for removal of the USTs at the PRP site; and
- Results of periodic groundwater monitoring in the Project Area from March 1988 to the present, as reported to the RWQCB. 0

NATURE AND EXTENT OF PETROLEUM HYDROCARBONS IN SOIL

Results of laboratory analyses of soil samples from near the USTs indicate the presence of petroleum hydrocarbons in the gasoline and diesel ranges (Attachment A). The area containing soils showing the presence of hydrocarbons based on laboratory analysis is closely correlated with soils showing sensory evidence of hydrocarbons such as odor or staining. The area containing these soils is referred to here as the release area. With the exception of one sample collected immediately below one UST, laboratory results do not indicate chlorinated hydrocarbons or other chemicals of concern were present in the release area.

The current and historical areal extent of elevated levels of petroleum hydrocarbons in soil in the vicinity of the USTs is presented on Plate 1. The historical extent of hydrocarbons within EBMUD and PRP property boundaries is based on observations made during excavation activities when soils containing petroleum hydrocarbons were encountered between 20 to 30 feet bgs. No soil containing petroleum hydrocarbons remains within the property boundaries, because the excavation depth for each building exceeded 30 feet (approximately 40 feet bgs for EBMUD and approximately 38 feet bgs for PRP).

The lateral extent of hydrocarbons in soil beyond the property boundaries has been estimated based on observations made during shoring construction activities for the two buildings and known migration patterns of hydrocarbons released at other locations in the Project Area. The lateral extent beneath Franklin Street, west of the EBMUD building, has been estimated based on observations of soil cutting returns from borings for soil nails, which are approximately 25 fact land and it. for soil nails, which are approximately 35 feet long and were drilled at an angle of approximately 30° below the horizontal. The lateral extent west of the PRP site has been estimated based on observations from 60-foot-long tieback borings, also drilled at angles of approximately 30° below the horizontal. The lateral extent is constrained to the southwest by the location of Monitoring Well MW-7. Observations made during installation of this well did not indicate that elevated levels of hydrocarbons were present in the soil (HLA, 1989).

Hydrocarbon releases in the Project Area tend to migrate vertically downward to the water table and then laterally in the general direction of groundwater flow in the uppermost shallow aquifer. In this area, groundwater flow in the shallow aquifer prior to the construction of the two buildings is estimated to have been westerly (HLA, 1989); migration of gasoline released from USTs at the southeast corner of the PRP site (at 9th and Webster Streets) was also generally westerly (HLA, 1988). The estimation of the extent of hydrocarbons in soil (shown on Plate 1) draws on these previous observations.

The ambient westerly direction of groundwater flow was altered by EBMUD and PRP construction dewatering activities from March 1988 to July 1991. Recent measurements of groundwater elevations in the Project Area indicate that foundation and shoring features of the buildings have also had permanent effects on local groundwater flow directions, as shown on Plate 1. The primary contributor to these effects is the PRP building. The shoring design for the building includes a low-permeability soil-and-concrete wall that extends along the entire 9th Street property boundary and approximately 150 feet north along both Franklin and Webster Streets. The excavation for the building removed soil to a depth of 38 feet bgs to the top of a clay layer underlying the silty sands that comprise the uppermost, shallow aquifer in the Project Area. The building foundation was constructed directly on native soil at 38 feet bgs. For these reasons, the building acts as a barrier to groundwater flow, with shallow aquifer groundwater flowing around the structure.

The EBMUD building has a key difference in its foundation as compared to the PRP building: a 1-foot-thick drain blanket of coarse gravel was placed between native soil and the concrete mat foundation of the building. This gravel layer appears to provide good hydraulic connection beneath the EBMUD building; a groundwater flow pathway is believed to exist beneath the building.

The potentiometric contours shown on Plate 1 were constructed using water-level elevations measured in December 1991 and incorporate the estimated effect of the PRP building on groundwater flow. Measurements made in September 1991 and March 1992 show a generally similar pattern. Groundwater flow lines were constructed using the potentiometric contours, standard techniques of flow net construction, and Darcy's Law for groundwater flow. Key assumptions reflected in the flow net presented on Plate 1 include:

- The PRP building acts as an impermeable barrier to regional groundwater flow. Groundwater flow theory states that flow lines will divide and flow around such an obstacle.
- Flow around an impermeable barrier must be parallel to the boundary of the obstacle; otherwise, there would be a component of flow through the obstacle, which contradicts the fact that the building foundation is an impermeable

barrier. This results in two flow lines, diverging around the building and running parallel to the building footprint.

Flow lines are constructed perpendicular to potentiometric contours. This assumes that the aquifer of interest is homogeneous in its properties. Based on numerous soil borings, well installations, and foundation excavations in the Project Area, the shallow aquifer has been observed to comprise silty sands in this area and can be considered homogeneous.

Several key points can be drawn from examination of the resultant flow net.

- o Groundwater flow is generally westerly in the Project Area but is diverted around the PRP building.
- Well MW-7 is positioned approximately 60 feet downgradient of the estimated lateral extent of the soil release area.

The observation that MW-7 is downgradient of the soil release area is further supported by groundwater chemistry data from samples collected in December 1991 and March 1992. The data indicate slightly elevated levels of total petroleum hydrocarbons (TPH) and gasoline constituents including benzene, toluene, ethylbenzene, and xylenes (BTEX) in groundwater at this location (HLA, 1992).

RESPONSE TO REQUEST FOR WORK PLAN

The following activities are proposed to monitor hydrocarbons in groundwater in the release area.

- Continue to monitor groundwater chemistry conditions at Well MW-7.
- o Continue to monitor groundwater flow directions in the Project Area.

Additional soil borings or monitoring wells are not proposed in the release area. Further characterization of the extent of hydrocarbons in soil below Franklin Street is not proposed for the following reasons.

- The nature and extent of hydrocarbons in soil are adequately characterized for purposes of assessing the need for and possible types of remedial action.
- The objectives and benefits of further characterization have not been established. Discussion of possible responses to the existence of hydrocarbons beneath the public right-of-way is a prerequisite to assessing the need for and benefits of further characterization. In particular, it does not appear that further characterization will provide regulatory decision-makers with any

information that will materially change the understanding of the problem or the options available for addressing the problem.

- o Groundwater elevation and chemistry data indicate that Well MW-7 is properly positioned to monitor groundwater conditions downgradient of the soil release area.
- Subsurface exploration in the street is dangerous with respect to both health and safety and because of the presence of underground utilities. Franklin Street is an urban, high-volume traffic area; work in the street area exposes drilling crews and the public to unnecessary hazards. With respect to underground utilities, the Pacific Bell fiber optic cable beneath Franklin Street is of particular concern; this cable is a main telecommunications service link in the East Bay.

The Agency would be pleased to discuss this letter, its contents and any related issues at a time convenient for ACDEH and the RWQCB, and would be particularly interested in discussing ACDEH and RWQCB objectives with respect to further characterization of hydrocarbons in the release area.

If you have questions regarding this letter please call David Leland at (415) 899-7352 or Peter Chen of the Agency at (510) 238-3692.

Yours very truly,

HARDING LAWSON ASSOCIATES

David F. Leland, P.E. Associate Engineer

R. Bruce Scheibach
Principal Hydrogeologist

DFL/RBS/mm/A22505-H

Attachments: A Laboratory Analytical Results of Soil Samples Below USTs B Laboratory Analytical Results of Soil Samples Collected at Bottom of

Duct Excavation

Plate 1 - Plan of Sites and Vicinity and Groundwater Flow Net -December 1991

cc: Ariu Levi, Alameda County Department of Environmental Health Lester Feldman, California Regional Water Quality Control Board Peter Chen, Redevelopment Agency of the City of Oakland Donnell Choy, City of Oakland Attorney's Office Julie Carver, City of Oakland Office of Public Works, Real Estate

References:

Harding Lawson Associates, 1988. Site Characterization, Pacific Renaissance Plaza, Chinatown Redevelopment Project Area, Oakland, California. December 22.

Harding Lawson Associates, 1989. A-Aquifer Monitoring Report, Chinatown Redevelopment Project Area, Oakland, California. January 31.

Harding Lawson Associates, 1992. Report of Groundwater Monitoring, December 1991, Chinatown Redevelopment Project Area, Oakland, California. May 6.



NATIONAL ENVIRONMENTAL TESTING, INC.

NET Pacific, Inc. 435 Tesconi Circle Santa Rosa, CA 95401 Tel: (707) 526-7200

Fax: (707) 526-7200

Dave Leland Harding Lawson Associates 200 Rush Landing Novato, CA 94947 Date: 12/27/1991

NET Client Acct. No: 28100 NET Pacific Log No: 91.1274

Received: 12/16/1991

Client Reference Information

Franklin St., Oakland, Job: 91005,928.02

Sample analysis in support of the project referenced above has been completed and results are presented on following pages. Please refer to the enclosed "Key to Abbreviations" for definition of terms. Should you have questions regarding procedures or results, please feel welcome to contact Client Services.

Approved by:

Jules Skamarack Laboratory Manager

Enclosure(s)



Client Name: Harding Lawson Associates

NET Log No: 91.1274

Date: 12/27/1991

Page: 2

Ref: Franklin St., Oakland, Job: 91005,928.02

umposite

SAMPLE DESCRIPTION: 91121601

Date Taken: 12/16/1991 Time Taken: 12:05

LAB Job No: (-108002)

100 000 /			_		
Parameter		Method	Reporting Limit	Results	Units
pH - Corro		150.1	N/A	7.5	pH units
	/Ignitability	1010	80	>140	F
	se (IR,Total)	5520C	50	6,300	mg/Kg
	se (IR,Non-Polar)5520CEF	50	6,300	mg/Kg
Cyanide (To	otal)	9010	0.2	ND	mg/Kg
Cadmium	(ICP)	EPA 6010	2.0	ND	mg/Kg
Chromium	(ICP)	EPA 6010	2.0	40	mg/Kg
Lead	(GFAA)	EPA 7421	0.2	7.8	mg/Kg
Nickel	(ICP)	EPA 6010	5.0	29	mg/Kg
Zinc	(ICP)	EPA 6010	2.0	29	mg/Kg
DATE ANAI DILUTION as Gasol: METHOD 802 DATE ANAI DILUTION Benzene Ethlyben: Toluene Xylenes	30 (GC,FID) LYZED FACTOR* ine 20 (GC,Solid) LYZED FACTOR* zene	5030 8020 8020 8020 8020	1 2.5 2.5 2.5 2.5 2.5	 12-18-91 1.5 12-18-91 1 ND ND 3.2 2.5	mg/Kg ug/Kg ug/Kg ug/Kg ug/Kg
METHOD 3550 DILUTION FA DATE EXTRAC DATE ANALYZ as Creosot as Diesel as Motor C	ACTOR* CTED KED :e	3550 3550 3550	10 10 10	1 12-17-91 12-18-91 ND 16 83	mg/Kg mg/Kg mg/Kg



Client Name: Harding Lawson Associates

NET Log No: 91.1274

Date: 12/27/1991

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121602

Date Taken: 12/16/1991 Time Taken: 14:40

LAB Job No: (-108003)Reporting Parameter Method Limit Results Units pH - Corrosivity 150.1 N/A 8.0 pH units Flashpoint/Ignitability 80 1010 88 Oil & Grease (IR, Total) 5520C 50 400 mg/Kg Oil & Grease (IR, Non-Polar) 5520CEF 50 310 mg/Kg Cyanide (Total) 9010 0.02 ND mg/Kg Cadmium (ICP) **EPA** 6010 2.0 2.8 mg/Kg Chromium (ICP) **EPA 6010** 2.0 45 mg/Kg Lead (GFAA) EPA 7421 0.2 4.2 mg/Kg Nickel (ICP) **EPA** 6010 5.0 36 mg/Kg Zinc (ICP) EPA 6010 2.0 23 mg/Kg TPH (Gas/BTXE, Solid) METHOD 5030 (GC,FID) DATE ANALYZED 12-18-91 DILUTION FACTOR* 2,000 as Gasoline 5030 1 19,000 mg/Kg METHOD 8020 (GC, Solid) DATE ANALYZED 12-18-91 DILUTION FACTOR* 2,000 Benzene 8020 2.5 53,000 ug/Kg Ethlybenzene 8020 2.5 150,000 ug/Kg Toluene 8020 2.5 340,000 ug/Kg Xylenes (Total) 8020 2.5 580,000 ug/Kg METHOD 3550 (GC, FID) DILUTION FACTOR* 100 DATE EXTRACTED 12-17-91 DATE ANALYZED 12-18-91 as Creosote 3550 10 ND mg/Kg as Diesel 3550 10 4,100 mg/Kg as Motor Oil 3550 10 ND mg/Kg



Client Name: Harding Lawson Associates

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121602

Date Taken: 12/16/1991

Time Taken: 14:40

LAB Job No: (-108003)

TAR 100 NO: (-108	003)				
Parameter	Method	Reporting Limit	Results	Units	
rarameter	mechod	LIMIC	KeBuits	Units	
METHOD 8240 (GCMS, Solid)		•			
DATE ANALYZED			12-19-91		
DILUTION FACTOR*			1250		
Benzene	8240	5	ND	ug/Kg	
Acetone	8240	10	ND	ug/Kg	
Bromodichloromethane	8240	5	ND	ug/Kg	
Bromoform	8240	5	ND	ug/Kg	
Bromomethane	8240	5	ND	ug/Kg	
2-Butanone	8240	10	ND	ug/Kg	
Carbon disulfide	8240	5	ND	ug/Kg	
Carbon tetrachloride	8240	5	ND	ug/Kg	
Chlorobenzene	8240	5	ND	ug/Kg	
Chloroethane	8240	5	ND	ug/Kg	
2-Chloroethyl vinyl ether		10	ND	ug/Kg	
Chloroform	8240	5	ND	ug/Kg ug/Kg	
Chloromethane	8240	5	ND	ug/Kg	
Dibromochloromethane	8240	5	ND	ug/Kg	
1,2-Dichlorobenzene	8240	5	ND	ug/Kg ug/Kg	
1,3-Dichlorobenzene	8240	5	ND	ug/Kg ug/Kg	
1,4-Dichlorobenzene	8240	5	ND	ug/Kg	
1,1-Dichloroethane	8240	5	ND	ug/Kg	
1,2-Dichloroethane	8240	5	ND	ug/Kg	
1,1-Dichloroethene	8240	5	ND	ug/Kg	
trans-1,2-Dichloroethene	8240	5	ND	ug/Kg	
1,2-Dichloropropane	8240	5	ND	ug/Kg	
cis-1,3-Dichloropropene	8240	5	ND	ug/Kg	
trans-1,3-Dichloropropene		5	ND	ug/Kg	
Ethyl benzene	8240	5	33,000	ug/Kg	
2-Hexanone	8240	10	ND		
Methylene chloride	8240	5	ND	ug/Kg	
4-Methyl-2-pentanone	8240	10	ND	ug/Kg ug/Kg	
Styrene	8240	5	ND	ug/Kg	
1,1,2,2-Tetrachloroethane		5	ND	ug/Kg	
Tetrachloroethene	8240	5	ND	• •	
Toluene	8240	5	96,000	ug/Kg ug/Kg	
1,1,1-Trichloroethane	8240	5	ND	ug/Kg	
1,1,2-Trichloroethane	8240	5	ND	ug/Kg	
Trichloroethene	8240	5	ND	ug/Kg	
Trichlorofluoromethane	8240	5	ND	ug/Kg	
Vinyl acetate	8240	10	ND	ug/Kg	
Vinyl chloride	8240	5	ND	ug/Kg ug/Kg	
Xylenes (total)	8240	5	460,000	ug/Kg	
SURROGATE RESULTS	-210	J		ag/vg	
Toluene-d8	8240		109	8	
Bromofluorobenzene	8240		117	3 %	
1,2-Dichloroethane-d4	8240		74	*5 *8	
-,	0470		7 T	•	



Client Name: Harding Lawson Associates

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121602
Date Taken: 12/16/1991
Time Taken: 14:40

LAB Job No: (-108003)

	Reporting			
Parameter	Method	Limit	Results	Units
METHOD 8270(GCMS, Solid)				
DATE EXTRACTED			12-17-91	
DATE ANALYZED			12-18-91	
DILUTION FACTOR*			1	
Acenaphthene	8270	330	ND	ug/Kg
Acenaphthylene	8270	330	ND	ug/Kg
Aldrin	8270	1600	ND	ug/Kg
Anthracene	8270	330	ND	ug/Kg
Benzidine	8270	1600	ND	ug/Kg
Benzo(a)anthracene	8270	330	ND	ug/Kg
Benzo(b) fluoranthene	8270	330	ND	ug/Kg
Benzo(k)fluoranthene	8270	330	ND	ug/Kg
Benzo(a)pyrene	8270	330	ND	ug/Kg
Benzo(g,h,i)perylene	8270	330	ND	ug/Kg
Benzoic acid	8270	1600	ND	ug/Kg
Benyzl alcohol	8270	330	ND	ug/Kg
Butyl benzyl phthalate	8270	330	ND	ug/Kg
delta-BHC	8270	1600	ND	ug/Kg
gamma-BHC	8270	1600	ND	ug/Kg
bis(2-Chloroethyl)ether	8270	330	ND	ug/Kg
bis(2-Chloroethoxy)methan		330	ND	ug/Kg
bis(2-Chloroisopropyl)eth		330	ND	ug/Kg
bis(2-Ethylhexyl)phthalat		330	ND	ug/Kg
4-Bromophenyl phenyl ethe		330	ND	ug/Kg
4-Chloroanaline	8270	330	ND	ug/Kg
2-Chloronaphthalene	8270	330	ND	ug/Kg
4-Chlorophenyl phenyl eth	8270	330	ND	ug/Kg
Chrysene	8270	330	ND	ug/Kg
4,4'-DDD	8270	1600	ND	ug/Kg
4,4'-DDE	8270	1600	ND	ug/Kg
4,4'-DDT	8270	1600	ND	ug/Kg
Dibenzo(a,h)anthracene	8270	330	ND	ug/Kg
Dibenzofuran	8270	330	ND	ug/Kg
Di-n-butylphthalate	8270	330	ND	ug/Kg
1,2-Dichlorobenzene	8270	330	ND	ug/Kg
1,3-Dichlorobenzene	8270	330	ND	ug/Kg ug/Kg
1,4-Dichlorobenzene	8270	330	ND	ug/Kg ug/Kg
3,3'-Dichlorobenzidine	8270	660	ND	ug/Kg
Dieldrin	8270	1600	ND	ug/Kg
Diethylphthalate	8270	330	ND	ug/Kg
Dimethyl phthalate	8270	330	ND	ug/Kg
2,4-Dinitrotoluene	8270	330	ND	ug/Kg
2,6-Dinitrotoluene	8270	330	ND	ug/Kg
	-			-9/ Mg



Client Name: Harding Lawson Associates

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Date: 12/27/1991

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121602

Date Taken: 12/16/1991 Time Taken: 14:40 LAB Job No: (-108003)

THE DOD NO: (-108003)		Danashisa		
Parameter	Method	Reporting Limit	Results	Units
	Mechod	Fillif	Results	Onica
Di-n-octyl phthalate	8270	330	ND	ug/Kg
Endrin aldehyde	8270	1600	ND	ug/Kg
Fluoranthene	8270	330	ND	ug/Kg
Fluorene	8270	330	ND	ug/Kg
Heptachlor	8270	1600	ND	ug/Kg
Heptachlor epoxide	8270	1600	ND	ug/Kg
Hexachlorobenzene	8270	330	ND	ug/Kg
Hexachlorobutadiene	8270	330	ND	ug/Kg
Hexachlorocyclopentadiene	8270	330	ND	ug/Kg
Hexachloroethane	8270	330	ND	ug/Kg
Indeno(1,2,3-cd)pyrene	8270	330	ND	ug/Kg
Isophorone	8270	330	ND	ug/Kg
2-Methylnaphthalene	8270	330	2,100	ug/Kg
Naphthalene	8270	330	2,300	ug/Kg
2-Nitroaniline	8270	1600	ND	ug/Kg
3-Nitroaniline	8270	1600	ND	ug/Kg
4-Nitroaniline	8270	1600	ND	ug/Kg
Nitrobenzene	8270	330	ND	ug/Kg
N-Nitroso-Di-N-propylamin	8270	330	ND	ug/Kg
N-Nitrosodiphenylamine	8270	330	ND	ug/Kg
Phenanthrene	8270	330	ND	ug/Kg
Pyrene	8270	330	ND	ug/Kg
1,2,4-Trichlorobenzene	8270	330	ND	ug/Kg
ACID EXTRACTABLES				37 3
4-Chloro-3-methylphenol	8270	330	ND	ug/Kg
2-Chlorophenol	8270	330	ND	ug/Kg
2,4-Dichlorophenol	8270	330	ND	ug/Kg
2,4-Dimethylphenol	8270	330	ND	ug/Kg
2,4-Dinitrophenol	8270	1600	ND	ug/Kg
4,6-Dinitro-2-methylpheno	8270	1600	ND	ug/Kg
2-Nitrophenol	8270	330	ND	ug/Kg
4-Nitrophenol	8270	1600	ND	ug/Kg
Pentachlorophenol	8270	1600	ND	ug/Kg
Phenol	8270	330	ND	ug/Kg
2,4,6-Trichlorophenol	8270	330	ND	ug/Kg
2-Methylphenol	8270	330	ND	ug/Kg
4-Methylphenol	8270	330	ND	ug/Kg
2,4,5-Trichlorophenol	8270	1600	ND	ug/Kg
• -				~3\ tra



Client Name: Harding Lawson Associates NET Log No: 91.1274

Date: 12/27/1991

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121602

Date Taken: 12/16/1991 Time Taken: 14:40 LAB Job No: (-108003)

Parameter	Method	Reporting Limit	Results	Units
SURROGATE RESULTS				
Nitrobenzene-d5	8270		46	*
2-Fluorobiphenyl	8270		46	*
p-Terphenyl-d14	8270		58	8
Phenol-d5	8270		37	%
2-Fluorophenol	8270		39	9,
2,4,6-Tribromophenol	8270		56	8



Client Name: Harding Lawson Associates

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121602

Date Taken: 12/16/1991 Time Taken: 14:40 LAB Job No: (-108003)

		Reporting		
Parameter	Method	Limit	Results	Units
METHOD 8270(GCMS, Soli	d, PCB			
DATE EXTRACTED			12-17-91	
DATE ANALYZED			12-18-91	
DILUTION FACTOR*			1	
POLYCHLORINATED BIPH	ENYLS			
Aroclor 1016	8270	660	ND	ug/Kg
Aroclor 1221	8270	660	ND	ug/Kg
Aroclor 1232	8270	660	ND	ug/Kg
Aroclor 1242	8270	660	ND	ug/Kg
Aroclor 1248	8270	660	ND	ug/Kg
Aroclor 1254	8270	660	ND	ug/Kg
Aroclor 1260	8270	660	ND	ug/Kg



Client Name: Harding Lawson Associates

NET Log No: 91.1274

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121603

Date Taken: 12/16/1991 Time Taken: 14:30 LAB Job No: (-108004)

Parameter	Method	Reporting Limit	Results	Units	
Oil & Grease (IR, Total) Oil & Grease (IR, Non-Pola	r)	50 50	ND ND	mg/Kg mg/Kg	
Cadmium (ICP) Chromium (ICP) Lead (GFAA) Nickel (ICP) Zinc (ICP)	EPA 6010 EPA 6010 EPA 7421 EPA 6010 EPA 6010	2.0 2.0 0.2 5.0 2.0	2.6 46 4.7 36	mg/Kg mg/Kg mg/Kg mg/Kg mg/Kg	
TPH (Gas/BTXE,Solid) METHOD 5030 (GC,FID) DATE ANALYZED DILUTION FACTOR* as Gasoline METHOD 8020 (GC,Solid)	5030	1	 12-18-91 1 3.1	mg/Kg	
DATE ANALYZED DILUTION FACTOR* Benzene Ethlybenzene Toluene	8020 8020 8020	2.5 2.5 2.5	12-18-91 1 46 77 130	ug/Kg ug/ Kg ug/Kg	= .046 ppm
Xylenes (Total) METHOD 3550 (GC,FID) DILUTION FACTOR* DATE EXTRACTED DATE ANALYZED as Creosote as Diesel as Motor Oil	8020 3550 3550 3550	10 10 10	370 1 12-17-91 12-18-91 ND 5.9	ug/Kg mg/Kg mg/Kg mg/Kg	



Client Name: Harding Lawson Associates

NET Log No: 91.1274

Date: 12/27/1991

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121603
Date Taken: 12/16/1991
Time Taken: 14:30

LAB Job No: (-108004)

Reporting				
Parameter	Method	Limit	Results	Units
METHOD 8240 (GCMS, Solid)				
DATE ANALYZED			10 10 01	
DILUTION FACTOR*			12-19-91	
Benzene	8240	_	5	
Acetone	8240	5	35	ug/Kg
Bromodichloromethane	8240	10	ND	ug/Kg
Bromoform	8240	5 5	ND	ug/Kg
Bromomethane	8240	5	ND	ug/Kg
2-Butanone	8240	10	ND ND	ug/Kg
Carbon disulfide	8240	5	ND	ug/Kg
Carbon tetrachloride	8240	5	ND	ug/Kg
Chlorobenzene	8240	5	ND	ug/Kg
Chloroethane	8240	5	ND	ug/Kg
2-Chloroethyl vinyl ether		10	ND	ug/Kg
Chloroform	8240	5	ND	ug/Kg
Chloromethane	8240	5	ND	ug/Kg
Dibromochloromethane	8240	5	ND	ug/Kg
1,2-Dichlorobenzene	8240	5		ug/Kg
1,3-Dichlorobenzene	8240	5	ND ND	ug/Kg
1,4-Dichlorobenzene	8240	5		ug/Kg
1,1-Dichloroethane	8240	5	ND ND	ug/Kg
1,2-Dichloroethane	8240	5		ug/Kg
1,1-Dichloroethene	8240	5	ND ND	ug/Kg
trans-1,2-Dichloroethene	8240	5	ND	ug/Kg
1,2-Dichloropropane	8240	5	ND	ug/Kg
cis-1,3-Dichloropropene	8240	5		ug/Kg
trans-1,3-Dichloropropene		5	ND ND	ug/Kg
Ethyl benzene	8240	5	70	ug/Kg
2-Hexanone	8240	10	ND	ug/Kg
Methylene chloride	8240	5	ND	ug/Kg
4-Methyl-2-pentanone	8240	10	ND	ug/Kg
Styrene	8240	5	ND	ug/Kg
1,1,2,2-Tetrachloroethane		5	ND	ug/Kg
Tetrachloroethene	8240	5	ND	ug/Kg
Toluene	8240	5	340	ug/Kg
1,1,1-Trichloroethane	8240	5	28	ug/Kg
1,1,2-Trichloroethane	8240	5	ND	ug/Kg
Trichloroethene	8240	5	50	ug/Kg
Trichlorofluoromethane	8240	5	ND	ug/Kg ug/Kg
Vinyl acetate	8240	10	ND	ug/Kg ug/Kg
Vinyl chloride	8240	5	ND	ug/Kg
Xylenes (total)	8240	5	800	
SURROGATE RESULTS		-		ug/Kg
Toluene-d8	8240		98	*
Bromofluorobenzene	8240		107	8
1,2-Dichloroethane-d4	8240		70	8
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Client Name: Harding Lawson Associates

NET Log No: 91.1274

Date: 12/27/1991

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121603
Date Taken: 12/16/1991

Time Taken: 14:30 LAB Job No: (-108004)

Parameter	Method	Reporting Limit	Results	**- * *-
ratanic cer	Mechod	TIWIE	Kesnics	Units
METHOD 8270 (GCMS, Solid)				
DATE EXTRACTED			12-07-91	
DATE ANALYZED			12-18-91	
DILUTION FACTOR*			1	
Acenaphthene	8270	330	ND	ug/Kg
Acenaphthylene	8270	330	ND	ug/Kg
Aldrin	8270	1600	ND	ug/Kg
Anthracene	8270	330	ND	ug/Kg
Benzidine	8270	1600	ND	ug/Kg
Benzo(a)anthracene	8270	330	ND	ug/Kg
Benzo(b) fluoranthene	8270	330	ND	ug/Kg
Benzo(k)fluoranthene	8270	330	ND	ug/Kg
Benzo(a)pyrene	8270	330	ND	ug/Kg
Benzo(g,h,i)perylene	8270	330	ND	ug/Kg
Benzoic acid	8270	1600	ND	ug/Kg
Benyzl alcohol	8270	330	ND	ug/Kg
Butyl benzyl phthalate	8270	330	ND	ug/Kg
delta-BHC	8270	1600	ND	
gamma-BHC	8270	1600	ND	ug/Kg
bis(2-Chloroethyl)ether	8270	330	ND	ug/Kg
bis(2-Chloroethoxy)methan		330	ND ND	ug/Kg
bis(2-Chloroisopropyl)eth		330	ND ND	ug/Kg
bis(2-Ethylhexyl)phthalat		330		ug/Kg
4-Bromophenyl phenyl ethe		330	ND	ug/Kg
4-Chloroanaline	8270 8270		ND	ug/Kg
2-Chloronaphthalene		330	ND	ug/Kg
4-Chlorophenyl phenyl eth	8270	330	ND	ug/Kg
		330	ND	ug/Kg
Chrysene	8270	330	ND	ug/Kg
4,4'-DDD	8270	1600	ND	ug/Kg
4,4'-DDE	8270	1600	ND	ug/Kg
4,4'-DDT	8270	1600	ND	ug/Kg
Dibenzo(a,h)anthracene	8270	330	ND	ug/Kg
Dibenzofuran	8270	330	ND	ug/Kg
Di-n-butylphthalate	8270	330	ND	ug/Kg
1,2-Dichlorobenzene	8270	330	ND	ug/Kg
1,3-Dichlorobenzene	8270	330	ND	ug/Kg
1,4-Dichlorobenzene	8270	330	ND	ug/Kg
3,3'-Dichlorobenzidine	8270	660	ND	ug/Kg
Dieldrin	8270	1600	ND	ug/Kg
Diethylphthalate	8270	330	ND	ug/Kg
Dimethyl phthalate	8270	330	ND	ug/Kg
2,4-Dinitrotoluene	8270	330	ND	ug/Kg
2,6-Dinitrotoluene	8270	330	ND	ug/Kg



Client Name: Harding Lawson Associates NET Log No: 91.1274

Date: 12/27/1991

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121603

Date Taken: 12/16/1991 Time Taken: 14:30 LAB Job No: (-108004)

•				
_	Reporting			
Parameter	Method	Limit	Results	Units
Di-n-octyl phthalate	8270	330	ND	ug/Kg
Endrin aldehyde	8270	1600	ND	ug/Kg
Fluoranthene	8270	330	ND	ug/Kg
Fluorene	8270	330	ND	ug/Kg
Heptachlor	8270	1600	ND	ug/Kg
Heptachlor epoxide	8270	1600	ND	ug/Kg
Hexachlorobenzene	8270	330	ND	ug/Kg
Hexachlorobutadiene	8270	330	ND	ug/Kg
Hexachlorocyclopentadiene	8270	330	ND	ug/Kg
Hexachloroethane	8270	330	ND	ug/Kg
Indeno(1,2,3-cd)pyrene	8270	330	ND	ug/Kg
Isophorone	8270	330	ND	ug/Kg
2-Methylnaphthalene	8270	330	6,400	ug/Kg
Naphthalene	8270	330	6,300	ug/Kg
2-Nitroaniline	8270	1600	ND	ug/Kg
3-Nitroaniline	8270	1600	ND	ug/Kg
4-Nitroaniline	8270	1600	ND	ug/Kg
Nitrobenzene	8270	330	ND	ug/Kg
N-Nitroso-Di-N-propylamin	8270	330	ND	ug/Kg
N-Nitrosodiphenylamine	8270	330	ND	ug/Kg
Phenanthrene	8270	330	ND	ug/Kg
Pyrene	8270	330	ND	ug/Kg
1,2,4-Trichlorobenzene	8270	330	ND	ug/Kg
ACID EXTRACTABLES				<i>3,</i> 3
4-Chloro-3-methylphenol	8270	330	ND	ug/Kg
2-Chlorophenol	8270	330	ND	ug/Kg
2,4-Dichlorophenol	8270	330	ND	ug/Kg
2,4-Dimethylphenol	8270	330	ND	ug/Kg
2,4-Dinitrophenol	8270	1600	ND	ug/Kg
4,6-Dinitro-2-methylpheno	8270	1600	ND	ug/Kg
2-Nitrophenol	8270	330	ND	ug/Kg
4-Nitrophenol	8270	1600	ND	ug/Kg
Pentachlorophenol	8270	1600	ND	ug/Kg
Phenol	8270	330	ND	ug/Kg
2,4,6-Trichlorophenol	8270	330	ND	ug/Kg
2-Methylphenol	8270	330	ИD	ug/Kg
4-Methylphenol	8270	330	ИD	ug/Kg



Client Name: Harding Lawson Associates

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121603

Date Taken: 12/16/1991 Time Taken: 14:30 LAB Job No: (-108004)

Parameter	Method	Reporting Limit	Results	Units	
SURROGATE RESULTS					
Nitrobenzene-d5	8270		44	*	
2-Fluorobiphenyl	8270		59	*	
p-Terphenyl-d14	8270		69	8	
Phenol-d5	8270		36	%	
2-Fluorophenol	8270		49	8	
2,4,6-Tribromophenol	8270		93	*	



Client Name: Harding Lawson Associates

NET Log No: 91.1274

Date: 12/27/1991

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Ref: Franklin St., Oakland, Job: 91005,928.02

SAMPLE DESCRIPTION: 91121603

Date Taken: 12/16/1991 Time Taken: 14:30 LAB Job No: (-108004)

Parameter	Method	Reporting Limit	Results	Units
METHOD 8270(GCMS, Solid,	РСВ		·	
DATE EXTRACTED			12-17-91	
DATE ANALYZED			12-18-91	
DILUTION FACTOR*			1	
POLYCHLORINATED BIPHEN	YLS			
Aroclor 1016	8270	660	ND	ug/Kg
Aroclor 1221	8270	660	ND	ug/Kg
Aroclor 1232	8270	660	ND	ug/Kg
Aroclor 1242	8270	660	ND	ug/Kg
Aroclor 1248	8270	660	ND	ug/Kg
Aroclor 1254	8270	660	ND	ug/Kg
Aroclor 1260	8270	660	ND	ug/Kg



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NET Log No: 91.1274

Date: 12/27/1991 Page: 15

Ref: Franklin St., Oakland, Job: 91005,928.02

QUALITY CONTROL DATA

Parameter	Reporting Limits	:	Cal Verf Stand % Recovery	Blank Data	Spike %	Duplicate Spike % Recovery	RPD
Diesel	1	mg/Kg	91	ND	N/A	N/A	9.3
Motor Oil	10	mg/Kg	116	ND	N/A	N/A	N/A
O&G (Total)	50	mg/Kg	102	ND	92	98	6.1
O&G(Non-Pola	ar)50	mg/Kg	111	ND	N/A	N/A	N/A
Gasoline	1	mg/Kg	89	ND	91	86	4.8
pH	N/A	pH unito	100	N/A	N/A	N/A	< 1
Flashpoint	80	F	100	ND	N/A	N/A	4.5
Cyanide	0.2	mg/Kg	98	ND	82	91	11
Cadmium	2.0	mg/Kg	107	ND	88	87	1.1
Chromium	2.0	mg/Kg	107	ND	87	87	< 1
Lead	0.2	mg/Kg	109	ND	92	90	2.9
Nickel	5.0	mg/Kg	110	ND	90	90	< 1
Zinc	2.0	mg/Kg	106	ND	84	88	3.7

QUALITY CONTROL DATA

Parameter	Reporting Limits		Cal Verf Stand % Recovery		Spike %	Duplicate Spike % Recovery	RPD
Chlorobenzene	5	ug/Kg	94	ND	111	111	< 1
1,1-Dichloroethene	5	ug/Kg	76	ND	117	119	1.0
Trichloroethene	5	ug/Kg	94	ND	113	116	3.0
Toluene	5	ug/Kg	99	ND	102	99	3.0
Benzene	5	ug/Kg	96	ND	96	97	< 1
COMMENT:	Blank Res	ults w	ere ND on	other	analytes	tested.	
Phenol	330	ug/Kg	112	ND	43	42	3.0
2-Chlorophenol	330	ug/Kg	108	ND	43	42	3.0
1,4-Dichlorobenzene	≥ 330	ug/Kg	107	ND	42	40	4.0
1,2,4-Trichlorobens	zene 330	ug/Kg	77	ND	76	61	22
Pyrene	330	ug/Kg	99	ND	84	74	13

COMMENT: Blank Results were ND on other analytes tested.



KEY TO ABBREVIATIONS and METHOD REFERENCES

<	:	Less than; When appearing in results column indicates analyte
		not detected at the value following. This datum supercedes
		the listed Reporting Limit.

Reporting Limits are a function of the dilution factor for any given sample. To obtain the actual reporting limits for this sample, multiply the stated Reporting Limits by the dilution factor (but do not multiply reported values).

ICVS : Initial Calibration Verification Standard (External Standard).

mean : Average; sum of measurements divided by number of measurements.

mg/Kg (ppm): Concentration in units of milligrams of analyte per kilogram of sample, wet-weight basis (parts per million).

mg/L : Concentration in units of milligrams of analyte per liter of sample.

mL/L/hr : Milliliters per liter per hour.

MPN/100 mL : Most probable number of bacteria per one hundred milliliters

of sample.

N/A : Not applicable.

NA : Not analyzed.

ND : Not detected; the analyte concentration is less than applicable

listed reporting limit.

NTU : Nephelometric turbidity units.

RPD : Relative percent difference, 100 [Value 1 - Value 2]/mean value.

SNA : Standard not available.

ug/Kg (ppb) : Concentration in units of micrograms of analyte per kilogram

of sample, wet-weight basis (parts per billion).

ug/L : Concentration in units of micrograms of analyte per liter of

sample.

umhos/cm : Micromhos per centimeter.

Method References

Methods 100 through 493: see "Methods for Chemical Analysis of Water & Wastes", U.S. EPA, 600/4-79-020, rev. 1983.

Methods 601 through 625: see "Guidelines Establishing Test Procedures for the Analysis of Pollutants" U.S. EPA, 40 CFR, Part 136, rev. 1988.

Methods 1000 through 9999: see "Test Methods for Evaluating Solid Waste", U.S. EPA SW-846, 3rd edition, 1986.

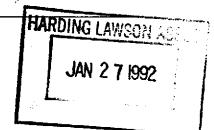
 $\underline{\text{SM}}\colon$ see "Standard Methods for the Examination of Water & Wastewater, 16th Edition, APHA, 1985.

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NATIONAL ENVIRONMENTAL TESTING, INC. NET Pacific, Inc. 435 Tesconi Circle Santa Rosa, CA 95401

Tel: (707) 526-7200 Fax: (707) 526-9623



David Leland Harding Lawson Associates 200 Rush Landing Novato, CA 94947

Date: 01/22/1992

NET Client Acct No: 28100 NET Pacific Log No: 91.1676

Received: 01/03/1992

Client Reference Information

PRP (PRA II), 91005,928.02

Sample analysis in support of the project referenced above has been completed and results are presented on following pages. Please refer to the enclosed "Key to Abbreviations" for definition of terms. Should you have questions regarding procedures or results, please feel welcome to contact Client Services.

Approved by:

Jules Skamarack Laboratory Manager

JS:rct Enclosure(s)



Client Name: Harding Lawson Associates

NET Log No: 91.1676

Date: 01/22/1992

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Page: 2

Ref: PRP (PRA II), 91005,928.02

Parameter	Method	Reporting Limit	91123001-911 12/30/1991 14:13 110105 1821TE 1 (No	91123005-91: 12/30/1991 14:27 ひかかがアモ 110106	_
TPH (Gas/BTXE, Solid) METHOD 5030 (GC, FID)					
DATE ANALYZED			01-10-92	01-10-92	
DILUTION FACTOR*			100	1	
as Gasoline		1	420	1.4	mg/Kg
METHOD 8020 (GC, Solid)					
DATE ANALYZED			01-10-91	01-10-91	
DILUTION FACTOR*			100	1	~ _ £00
Benzene		2.5	330	ND	ug/Kg
Ethylbenzene		2.5	3,700	3.5	ug/Kg
Toluene		2.5	7,100	ND	ug/Kg
Xylenes (Total)		2.5	19,000	12	ug/Kg
METHOD 3550 (GC,FID)					
DILUTION FACTOR*			1	1	
DATE EXTRACTED			01-13-92	01-13-92	
DATE ANALYZED			01-14-92	01-14-92	
as Diesel		1	52 **	1.5	mg/Kg

^{**} Petroleum hydrocarbons quantified as diesel appear to be a lighter hydrocarbon than diesel.



Client Name: Harding Lawson Associates

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Ref: PRP (PRA II), 91005,928.02

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			91123001-911 12/30/1991 14:13	91123005-911 12/30/1991 14:27	
Parameter	Method	Reporting Limit	110105	110106	Units
METHOD 8010 (GC, Solid)		· · · · · · · · · · · · · · · · · · ·		· · · · · · · · · · · · · · · · · · ·	
DATE ANALYZED			01-14-92	01-14-92	
DILUTION FACTOR*			1	1	
Bromodichloromethane		2.0	ND	ND	ug/Kg
Bromoform		2.0	ND	ND	ug/Kg
Bromomethane		2.0	ND	ND	ug/Kg
Carbon tetrachloride		2.0	ND	ND	ug/Kg
Chlorobenzene		2.0	ND .	ND	ug/Kg
Chloroethane		2.0	ND	ND	ug/Kg
2-Chloroethylvinyl ether		5.0	ND	ND	ug/Kg
Chloroform		2.0	ND	ND	ug/Kg
Chloromethane		2.0	ND	ND	ug/Kg
Dibromochloromethane		2.0	ND	ND	ug/Kg
1,2-Dichlorobenzene		2.0	ND	ND	ug/Kg
1,3-Dichlorobenzene		2.0	ND	ND	ug/Kg
1,4-Dichlorobenzene		2.0	ND	ND	ug/Kg
Dichlorodifluoromethane		2.0	ND	ND	ug/Kg
1,1-Dichloroethane		2.0	ND	ND	ug/Kg
1,2-Dichloroethane		2.0	ND	ND	ug/Kg
1,1-Dichloroethene		2.0	ND	ND	ug/Kg
trans-1,2-Dichloroethene		2.0	ND	ND	ug/Kg
1,2-Dichloropropane		2.0	ND	ND	ug/Kg
cis-1,3-Dichloropropene		2.0	ND	ND	ug/Kg
trans-1,3-Dichloropropene	:	2.0	ND	ND	ug/Kg
Methylene chloride		50	ND	ND	ug/Kg
1,1,2,2-Tetrachloroethane	ł	2.0	ND	ND	ug/Kg
Tetrachloroethene		2.0	ND	ND	ug/Kg
1,1,1-Trichloroethane		2.0	ND	ND	ug/Kg
1,1,2-Trichloroethane		2.0	ND	ND	ug/Kg
Trichloroethene		2.0	ИĎ	ND	ug/Kg
Trichlorofluoromethane		2.0	ND	ND	ug/Kg
Vinyl chloride		2.0	ND	ND	ug/Kg



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Parameter	Method	Reporting Limit	91123001-911 12/30/1991 14:13	91123005-911 12/30/1991 14:27 110106	Units
rarameter	WACHOO	TIMIC	110103	110100	UIILE
METHOD 8270 (GCMS, Solid)				·-··	
DATE EXTRACTED			1/08/92	1/08/92	
DATE ANALYZED			1/13/92	1/13/92	
DILUTION FACTOR*			1	1	
Acenaphthene		330	ND	ND	ug/Kg
Acenaphthylene		330	ND	ND	ug/Kg
Aldrin		1600	ND	ПD	ug/Kg
Anthracene		330	ND	ND	ug/Kg
Benzidine		1600	ND .	ND	ug/Kg
Benzo(a)anthracene		330	ND	ND	ug/Kg
Benzo(b)fluoranthene		330	ND	ND	ug/Kg
Benzo(k)fluoranthene		330	ND	ND	ug/Kg
Benzo(a)pyrene		330	ND	ND	ug/Kg
Benzo(g,h,i)perylene		330	ND	ND	ug/Kg
Benzoic acid		1600	ND	ND	ug/Kg
Benyzl alcohol		330	מא	ND	ug/Kg
Butyl benzyl phthalate		330	ND	ND	ug/Kg
delta-BHC		1600	ND	ND	ug/Kg
gamma-BHC		1600	ND	ND	ug/Kg
bis(2-Chloroethyl)ether		330	ND	ND	ug/Kg
bis(2-Chloroethoxy)methan		330	ND	ND	ug/Kg
bis(2-Chloroisopropyl)eth		330	ND	ND	ug/Kg
bis(2-Ethylhexyl)phthalat		330	ND	ND	ug/Kg
4-Bromophenyl phenyl ethe		330	ND	ND	ug/Kg
4-Chloroanaline		330	ND	ND	ug/Kg
2-Chloronaphthalene		330	ND	ND	ug/Kg
4-Chlorophenyl phenyl eth		330	ND	ND	ug/Kg
Chrysene		330	ND	ND	ug/Kg
4,4'~DDD		1600	ND	ND	ug/Kg
4,4'-DDE		1600	ND	ND	ug/Kg
4,4'-DDT		1600	ND	ND	ug/Kg
Dibenzo(a,h)anthracene		330	ND	ND	ug/Kg
Dibenzofuran		330	ND	ND	ug/Kg
Di-n-butylphthalate		330	ND	ND	ug/Kg
1,2-Dichlorobenzene		330	ND	ND	ug/Kg
1,3-Dichlorobenzene		330	ND	ND	ug/Kg
1,4-Dichlorobenzene		330	ND	ND	ug/Kg
3,3'-Dichlorobenzidine		660	ND	ND	ug/Kg
Dieldrin		1600	ND	ND	ug/Kg
Diethylphthalate		330	ND	ND	ug/Kg
Dimethyl phthalate		330	ND	ND	ug/Kg
2,4-Dinitrotoluene		330	ND	ND	ug/Kg
2,6-Dinitrotoluene		330	ND	ND	ug/Kg
Di-n-octyl phthalate		330	ND	ND	ug/Kg
Endrin aldehyde		1600	ND	ND	ug/Kg



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		Reporting	91123001-911 12/30/1991 14:13	91123005-911 12/30/1991 14:27	
Parameter	Method	Limit	110105	110106	Units
Fluoranthene		330	ND	ND	ug/Kg
Fluorene		330	ND	ND	ug/Kg
Heptachlor		1600	ND	ND	ug/Kg
Heptachlor epoxide		1600	ND	ND	ug/Kg
Hexachlorobenzene		330	ND	ND	ug/Kg
Hexachlorobutadiene		330	ND	ND	ug/Kg
Hexachlorocyclopentadiene		330	ND	ND	ug/Kg
Hexachloroethane		330	ND	ND	ug/Kg
Indeno(1,2,3-cd)pyrene		330	ND	ND	ug/Kg
Isophorone		330	ND	ND	ug/Kg
2-Methylnaphthalene		330	ND	ND	ug/Kg
Naphthalene		330	900	ND	ug/Kg
2-Nitroaniline		1600	910	ND	ug/Kg
3-Nitroaniline		1600	ND	ND	ug/Kg
4-Nitroaniline		1600	ND	ND	ug/Kg
Nitrobenzene		330	ND	ND	ug/Kg
N-Nitroso-Di-N-propylamin		330	ND	ND	ug/Kg
N-Nitrosodiphenylamine		330	ND	ND	ug/Kg
Phenanthrene		330	ND	ND	ug/Kg
Pyrene		330	ND	ND	ug/Kg
1,2,4-Trichlorobenzene		330	ND	ND	ug/Kg
ACID EXTRACTABLES		550			~9/ .·9
4-Chloro-3-methylphenol		330	ND	ND	ug/Kg
2-Chlorophenol		330	ND	ND	ug/Kg
2,4-Dichlorophenol		330	ND	ND	
2,4-Dimethylphenol		330	ND ND		ug/Kg
2,4-Dimethylphenol				ND	ug/Kg
		1600	ND	ND	ug/Kg
4,6-Dinitro-2-methylpheno		1600	ND	ND	ug/Kg
2-Nitrophenol		330	ND	ND	ug/Kg
4-Nitrophenol		1600	ND	ND	ug/Kg
Pentachlorophenol		1600	ND	ND	ug/Kg
Phenol		330	ND	ND	ug/Kg
2,4,6-Trichlorophenol		330	ND	ND	ug/Kg
2-Methylphenol		330	ND	ND	ug/Kg
4-Methylphenol		330	ND	ND	ug/Kg
2,4,5-Trichlorophenol SURROGATE RESULTS		1600	ND 	ND 	ug/Kg
Nitrobenzene-d5			46	49	\$ rec
2-Fluorobiphenyl			43	50	% rec
p-Terphenyl-d14			79	79	% rec
Phenol-d5			44	54	% rec
2-Fluorophenol			41	50	% rec
2,4,6-Tribromophenol			58	59	% rec



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Ref: PRP (PRA II), 91005,928.02

Parameter	Method	Reporting	12/30/1991 14:42 110107	91123013-91 12/30/1991 15:14 110108 Composiff 4	units (よつい下計)
		20,110			(300111)
TPH (Gas/BTXE, Solid)	•				
METHOD 5030 (GC,FID) DATE ANALYZED			01-10-92	01-10-92	
DILUTION FACTOR*			1	1	
as Gasoline		1	ND	ND	mg/Kg
METHOD 8020 (GC, Solid)		•			mg/ 1.g
DATE ANALYZED			01-10-91	01-10-91	
DILUTION FACTOR*			1	1	
Benzene		2.5	ND .	ND	ug/Kg
Ethylbenzene		2.5	ND	ND	ug/Kg
Toluene		2.5	ND	ND	ug/Kg
Xylenes (Total)		2.5	ND	ND	ug/Kg
METHOD 3550 (GC,FID)					
DILUTION FACTOR*			1	1	
DATE EXTRACTED			01-13-92	01-13-92	
DATE ANALYZED			01-14-92	01-14-92	
as Diesel		1	ND	ND	mg/Kg



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Ref: PRP (PRA II), 91005,928.02

			91123009-911	91123013-911				
			12/30/1991	12/30/1991				
			14:42	15:14				
		Reporting						
Parameter	Method	Limit	110107	110108	Unit			
METHOD 8010 (GC, Solid)				<u>, , , , , , , , , , , , , , , , , , , </u>				
DATE ANALYZED			01-14-92	01-14-92				
DILUTION FACTOR*			1	1				
Bromodichloromethane		2.0	ND	ND	ug/Kg			
Bromoform		2.0	ND	ND	ug/Kg			
Bromomethane		2.0	ND	ND	ug/Kg			
Carbon tetrachloride		2.0	ND	ND	ug/Kg			
Chlorobenzene		2.0	ND	ND	ug/Kg			
Chloroethane		2.0	ND .	ND	ug/Kg			
2-Chloroethylvinyl ether		5.0	ND	ND	ug/Kg			
Chloroform		2.0	ND	ND	ug/Kg			
Chloromethane		2.0	ND	ND	ug/Kg			
Dibromochloromethane		2.0	ND	ND	ug/Kg			
1,2-Dichlorobenzene		2.0	ND	ND	ug/Kg			
1,3-Dichlorobenzene		2.0	ND	ND	ug/Kg			
1,4-Dichlorobenzene		2.0	ND	ND	ug/Kg			
Dichlorodifluoromethane		2.0	ND	ND	ug/Kg			
1,1-Dichloroethane		2.0	ND	ND	ug/Kg			
1,2-Dichloroethane		2.0	ND	ND	ug/Kg			
1,1-Dichloroethene		2.0	ND	ND	ug/Kg			
trans-1,2-Dichloroethene		2.0	ND	ND	ug/Kg			
1,2-Dichloropropane		2.0	ND	ND	ug/Kg			
cis-1,3-Dichloropropene		2.0	ND	ND	ug/Kg			
trans-1,3-Dichloropropene	ł	2.0	ND	ND	ug/Kg			
Methylene chloride		50	ND	ND	ug/Kg			
1,1,2,2-Tetrachloroethane	!	2.0	ND	ND	ug/Kg			
Tetrachloroethene		2.0	ND	ND	ug/Kg			
1,1,1-Trichloroethane		2.0	ND	ND	ug/Kg			
1,1,2-Trichloroethane		2.0	ND	ND	ug/Kg			
Trichloroethene		2.0	ND	ND	ug/Kg			
Trichlorofluoromethane		2.0	ND	מא	ug/Kg			
Vinyl chloride		2.0	ND	ND	ug/Kg			



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Date: 01/22/1992

Client No: 28100 Client Name: Harding Lawson Associates

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Descriptor, Lab No. and Results

91123009-911 91123013-911

			12/30/1991	12/30/1991	
			14:42	15:14	
		Reporting			
Parameter	Method	Limit	110107	110108	Units
METHOD 8270 (GCMS, Solid)				· · · · · · · · · · · · · · · · · · ·	
DATE EXTRACTED			1/08/92	1/08/92	
DATE ANALYZED			1/13/92	1/13/92	
DILUTION FACTOR*			1	1	
Acenaphthene		330	ND	ND	ug/Kg
Acenaphthylene		330	ND	ND	ug/Kg
Aldrin		1600	ND	ND	ug/Kg
Anthracene		330	ND	ND	ug/Kg
Benzidine		1600	ND	ND	ug/Kg
Benzo(a)anthracene		330	ND	ND	ug/Kg
Benzo(b)fluoranthene		330	ND	ND	ug/Kg
Benzo(k)fluoranthene		330	ND	ND	ug/Kg
Benzo(a)pyrene		330	ND	ND	ug/Kg
Benzo(g,h,i)perylene		330	ND	ND	ug/Kg
Benzoic acid		1600	ND	ND	ug/Kg
Benyzl alcohol		330	ND	ND	ug/Kg
Butyl benzyl phthalate		330	ND	ND	ug/Kg
delta-BHC		1600	ND	ND	ug/Kg
gamma-BHC		1600	ND	ND	ug/Kg
bis(2-Chloroethyl)ether		330	ND	ND	ug/Kg
bis(2-Chloroethoxy)methan		330	ND	ND	ug/Kg
bis(2-Chloroisopropyl)eth	ļ	330	ND	ND	ug/Kg
bis(2-Ethylhexyl)phthalat	;	330	ND	ND	ug/Kg
4-Bromophenyl phenyl ethe	t	330	ND	ND	ug/Kg
4-Chloroanaline		330	ND	ND	ug/Kg
2-Chloronaphthalene		330	ND	ND	ug/Kg
4-Chlorophenyl phenyl eth	ı	330	ND	ND	ug/Kg
Chrysene		330	ND	ND	ug/Kg
4,4'-DDD		1600	ИD	ND	ug/Kg
4,4'-DDE		1600	ND	ND	ug/Kg
4,4'-DDT		1600	ND	ND	ug/Kg
Dibenzo(a,h)anthracene		330	ND	ND	ug/Kg
Dibenzofuran		330	ND	ND	ug/Kg
Di-n-butylphthalate		330	ND	ND	ug/Kg
1,2-Dichlorobenzene		330	ND	ND	ug/Kg
1,3-Dichlorobenzene		330	ND	ND	ug/Kg
1,4-Dichlorobenzene		330	ND	ND	ug/Kg
3,3'-Dichlorobenzidine		660	ND	ND	ug/Kg
Dieldrin		1600	ND	ND	ug/Kg
Diethylphthalate		330	ND	ND	ug/Kg
Dimethyl phthalate		330	ND	ND	ug/Kg
2,4-Dinitrotoluene		330	ND	ND	ug/Kg
2,6-Dinitrotoluene		330	ND	ND	ug/Kg
Di-n-octyl phthalate		330	ND	ND	ug/Kg
Endrin aldehyde		1600	ND	ND	ug/Kg



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Descriptor, Lab No. and Results

91123009-911 91123013-911 12/30/1991 12/30/1991

			14:42	15:14	
		Reporting	•		
Parameter	Method	Limit	110107	110108	Units
Fluoranthene	- .	330	ND	ND	ug/Kg
Fluorene		330	ND	ND	ug/Kg
Heptachlor		1600	ND	ND	ug/Kg
Heptachlor epoxide		1600	ND	ND	ug/Kg
Hexachlorobenzene		330	ND	ND	ug/Kg
Hexachlorobutadiene		330	ND	ND	ug/Kg
Hexachlorocyclopentadiene	•	330	ND	ND	ug/Kg
Hexachloroethane		330	ND	ND	ug/Kg
Indeno(1,2,3-cd)pyrene		330	ND	ND	ug/Kg
Isophorone		330	ND	ND	ug/Kg
2-Methylnaphthalene		330	ND	ND	ug/Kg
Naphthalene		330	ND	ND	ug/Kg
2-Nitroaniline		1600	ND	ND	ug/Kg
3-Nitroaniline		1600	ND	ND	ug/Kg
4-Nitroaniline		1600	ND	ND	ug/Kg
Nitrobenzene		330	ND	ND	ug/Kg
N-Nitroso-Di-N-propylamin		330	ND	ND	ug/Kg
N-Nitrosodiphenylamine		330	ND	ND	ug/Kg
Phenanthrene		330	ND	ND	ug/Kg
Pyrene		330	ND	ND	ug/Kg
1,2,4-Trichlorobenzene		330	ND	ND	ug/Kg
ACID EXTRACTABLES					-919
4-Chloro-3-methylphenol		330	ND	ND	ug/Kg
2-Chlorophenol		330	ND	ND	ug/Kg
2,4-Dichlorophenol		330	ND	ND	ug/Kg
2,4-Dimethylphenol		330	ND	ND	ug/Kg
2,4-Dinitrophenol		1600	ND	ND	ug/Kg
4,6-Dinitro-2-methylpheno		1600	ND	ND	ug/Kg
2-Nitrophenol		330	ND	ND	ug/Kg
4-Nitrophenol		1600	ND	ND	ug/Kg
Pentachlorophenol		1600	ND	ND	ug/Kg
Phenol		330	ND	ND	ug/Kg
2,4,6-Trichlorophenol		330	ND	ND	ug/Kg
2-Methylphenol		330	ND	ND	ug/Kg
4-Methylphenol		330	ND	ND	ug/Kg
2,4,5-Trichlorophenol		1600	ND	ND	
SURROGATE RESULTS		1900	ND	ND 	ug/Kg
Nitrobenzene-d5			37	47	% rec
2-Fluorobiphenyl			46	4 / 50	% rec
p-Terphenyl-d14				- -	
Phenol-d5			81 49	84 53	% rec
2-Fluorophenol					% rec
			48	46	% rec
2,4,6-Tribromophenol			58	62	% rec



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QUALITY CONTROL DATA

<u>Parameter</u>	Reporting Limits	<u>Units</u>	Cal Verf Stand % Recovery	Blank Data	Spike %	Duplicate Spike % Recovery	RPD
Gasoline	1.0	mg/Kg	102	ND	95	94	< 1
Benzene	2.5	ug/Kg	98	ND	80	81	< 1
Toluene	2.5	ug/Kg ug/Kg	104	מא	92	93	1.3
TOIGENE	2.5	ug/kg	104	RU	72	73	1.3
Diesel	1	mg/Kg	110	ND	84	95	12
Motor Oil	10	mg/Kg	121	ND	N/A	N/A	N/A
					-	-	-
Benzene	2.0	mg/Kg	116	ND	242	266	9.3
Toluene	2.0	mg/Kg	111	ND	390	394	1.2
1,1-DCE	2.0	mg/Kg	100	ND	95	96	1.3
TCE	2.0	mg/Kg	100	ND	92	99	7.2
Chlorobenzene	∍2.0	mg/Kg	110	ND	104	104	< 1
Phenol	330	ug/Kg	112	ND	46	48	3.0
2-Chlorophen	330	ug/Kg	106	ND	43	45	3.0
1,4-DCB	330	ug/Kg	114	ND	36	36	1.0
N-Nitroso	330	ug/Kg	127	ND	47	46	1.0
1,2,4-TCB	330	ug/Kg	109	ND	43	49	12
4-Chloro-3	330	ug/Kg	100	ND	84	75	11
Acenaphthene	330	ug/Kg	115	ND	67	68	2.0
4-Nitrophen	1600	ug/Kg	43	ND	114	98	15
2,4-Dinitro	1600	ug/Kg	96	ND	78	70	10
Pentachloro	1600	ug/Kg	105	ND	93	89	4.0
Pyrene	330	ug/Kg	98	ND	86	80	7.0

COMMENT: Blank Results were ND on other analytes tested.



KEY TO ABBREVIATIONS and METHOD REFERENCES

<	:	Less than; When appearing in results column indicates analyte
		not detected at the value following. This datum supercedes
		the listed Reporting Limit.

: Reporting Limits are a function of the dilution factor for any given sample. To obtain the actual reporting limits for this sample, multiply the stated Reporting Limits by the dilution factor (but do not multiply reported values).

ICVS : Initial Calibration Verification Standard (External Standard).

mean : Average; sum of measurements divided by number of measurements.

mg/Kg (ppm): Concentration in units of milligrams of analyte per kilogram of sample, wet-weight basis (parts per million).

mg/L : Concentration in units of milligrams of analyte per liter of sample.

mL/L/hr : Milliliters per liter per hour.

MPN/100 mL : Most probable number of bacteria per one hundred milliliters

of sample.

N/A : Not applicable.

NA : Not analyzed.

ND : Not detected; the analyte concentration is less than applicable

listed reporting limit.

NTU : Nephelometric turbidity units.

RPD : Relative percent difference, 100 [Value 1 - Value 2]/mean value.

SNA : Standard not available.

ug/Kg (ppb) : Concentration in units of micrograms of analyte per kilogram

of sample, wet-weight basis (parts per billion).

ug/L : Concentration in units of micrograms of analyte per liter of

sample.

umhos/cm : Micromhos per centimeter.

Method References

Methods 100 through 493: see "Methods for Chemical Analysis of Water & Wastes", U.S. EPA, 600/4-79-020, rev. 1983.

Methods 601 through 625: see "Guidelines Establishing Test Procedures for the Analysis of Pollutants" U.S. EPA, 40 CFR, Part 136, rev. 1988.

Methods 1000 through 9999: see "Test Methods for Evaluating Solid Waste", U.S. EPA SW-846, 3rd edition, 1986.

 \underline{SM} : see "Standard Methods for the Examination of Water & Wastewater, $\overline{16}$ th Edition, APHA, 1985.

HL	P.O. Box 5 Novato, CA (415) 892-	CA 94948	Samplers:	Robert L. Nober	2 CONTRACTOR STEED
loh		91005,928.07	oampiei o. <u></u>		
Nom	number:_ ie/Locatio	on: PRP (PRA II)	<u> </u>	S. 184	—
Proj	ect Mana	ager: David Leland	Recorder:	Robert L. Nelman Signature Required)	Metals Metals drocarb
GE	MATRIX	#CONTAINERS SAMPLE & PRESERV. NUMBER OR LAB	DATE	STATION DESCRIPTION/ NOTES	EPA 601/8010) EPA 602/8020 EPA 624/8240 EPA 625/8270) X Priority Pilmt. Metals Benzene/Toluene/Xylene Total Petrol. Hydrocarb. ### Con & Dank ####################################
SOURCE CODE	Water Sediment Soil Oil	NUMBER Yr Wk Se	R Seq Yr Mo Dy Time		EPA 60 EPA 62 EPA 62 EPA 62 EPA 62 Priority TOTA Benzence
46	X			3	
48	X	 -	0029//230/4/	<u> </u>	X X X X Composer
48	<u> </u>	▎▘▊▕▎▘▎▐▕▋▕▍▕▍▕▍▕▍	003911230142		
48	X	 	049/1230142	5	
48	X			ŽĮ ————————————————————————————————————	
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