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Winston-Salem
Regional Office

Luck's® Inc.

Heat and Eat Country Style™ Foods

P.O. Box 510, Hwy. 705 E.
Seagrove, NC 27341-0510
(919) 873-7211

September 3, 1992

Ms. Sherri V. Knight
Regional Groundwater Supervisor
State of North Carolina
Division of Environmental Management
8025 North Point Blvd.
Winston-Salem, NC 27106

RE: Soil Sampling - 1990 UST Closure - McLaren Hart Report

Dear Ms. Knight:

The attached report is the soil samplings results from sampling activities conducted on July 7th and 8th, 1992 by McLaren/Hart at the Luck's, Inc. facility in Seagrove, NC. The report contains background information, site investigations, information on the current investigation and conclusions/recommendations.

Data from this sampling continues to support Luck's knowledge that the tanks had not leaked and any oil and grease detected by sampling methods 9071 and 9073 are not from the tanks or their removal since no detectable levels of petroleum hydrocarbons have been found.

I am enclosing six pictures of the three #5 fuel oil tanks and the ground under each of these tanks that clearly show no evidence of oil leakage.

I trust this will be helpful.

Sincerely,

Darius Luck

Darius Luck
Vice-President/
Plant Manager

DL/jk

Attachment

cc: C. Elmendorf - McLaren Hart
G. Belardo - NYO
T. Anderson
B. McKenzie

UNDERGROUND STORAGE TANK CLOSURE REPORT
LUCKS, INC., SEAGROVE, N.C.

I. Background:

Three underground storage tanks (USTs), including two 10,000 gallon USTs and one 8,000 gallon UST, formerly used to store No. 5 fuel oil were removed by Luck's Incorporated during April 25 and 26, 1990. The USTs were excavated from an area measuring approximately 30 feet by 30 feet, to a depth of approximately 20 feet. The tanks were excavated and disposed of in accordance with applicable state and federal regulations. The tank excavation area is shown on the site map in Figure 1.

There was no visual evidence of free product observed in the tank excavation. Additionally, there was no evidence of fuel oil odor from the soils in the excavation. Based on verbal accounts from facility personnel, the USTs were in good condition and did not appear to be leaking when they were removed. Post-excavation samples were collected from the base of the excavation in August 1990 and analyzed. The excavation was then backfilled with clean soil.

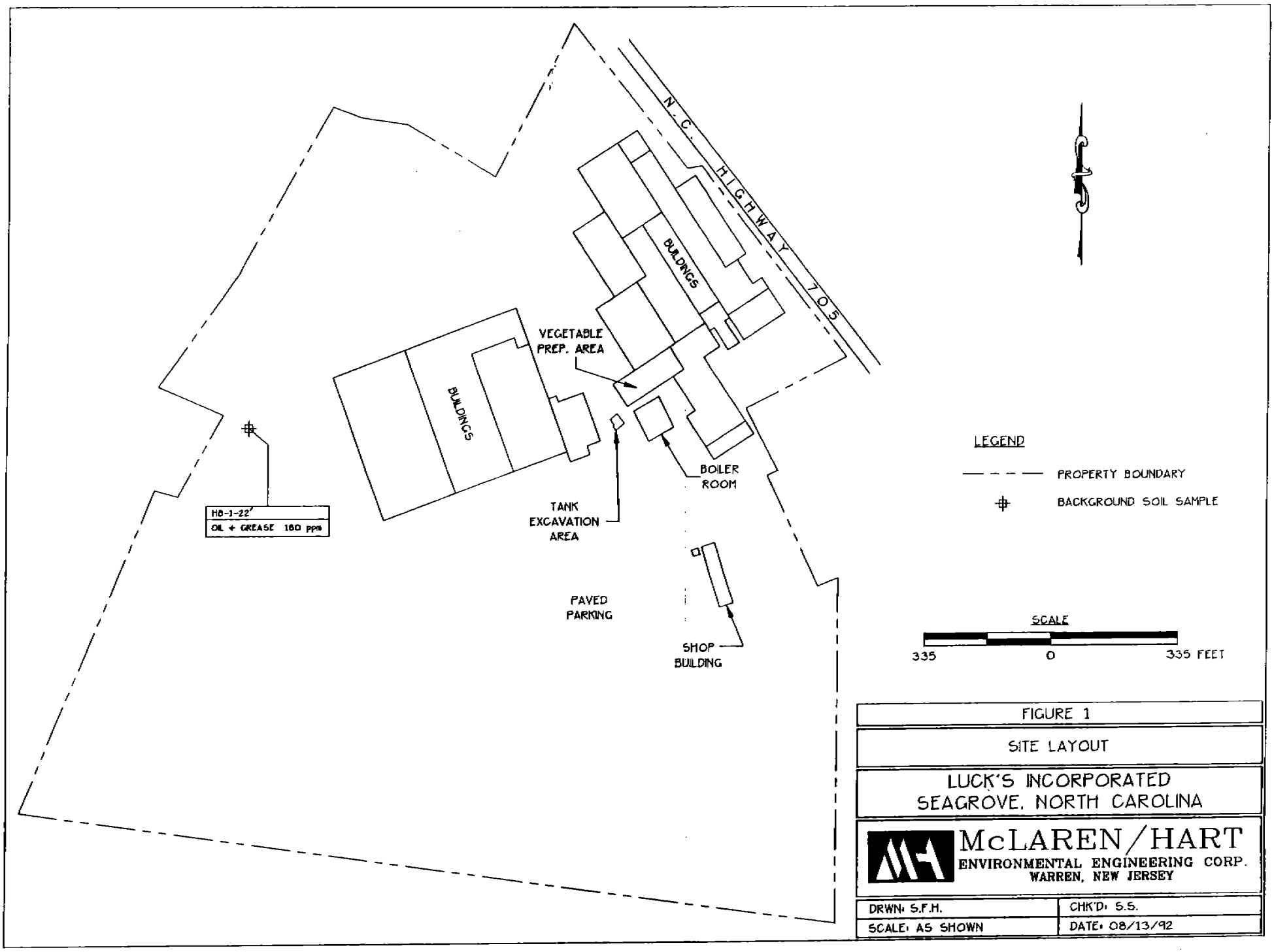
This report summarizes earlier investigations of the UST area, and describes the results of recent resampling of soils in this area, which was conducted in July 1992 by McLaren/Hart in order to satisfy the requirements of the North Carolina Department of Environmental Health and Natural Resources (NCDEHNR) for tank closure.

II. Site geology

According to the North Carolina State Geological Survey, the site lies within the east-west trending Carolina Slate Belt physiographic province. Bedrock underlying the site is predominantly made up of Precambrian age metamorphosed volcanics of the Uwharrie Formation. The Uwharrie Formation consists of light-gray to greenish-gray felsic metavolcanics (meta dacitic-rhyolitic flows and tuffs) interbedded with mafic and intermediate volcanic rock and meta-argillites and meta-mudstones.

A review of available boring logs indicate that the depth to bedrock is approximately 50 feet beneath the site. The bedrock surface is highly weathered (saprolites), with a high silt and clay content. The bedrock is overlain by sand which ranges in thickness from 15 to 45 feet. Overlying the sand deposits are approximately 5 to 10 feet of red-brown clay.

Groundwater was measured beneath the property at a depth of approximately 43.5 feet below grade in a background monitoring well located approximately 500 feet west of the tank excavation. A review of the monitoring well record indicated that this well was completed within the overburden sand deposits to a depth of 50 feet below grade (Attachment 1).


N. C. HIGHWAY 705
BUILDINGS
BUILDINGS
VEGETABLE PREP. AREA
BOILER ROOM
TANK EXCAVATION AREA
PAVED PARKING
SHOP BUILDING
HB-1-22'
OIL + GREASE 160 ppm
LEGEND
PROPERTY BOUNDARY
BACKGROUND SOIL SAMPLE
SCALE
335
0
335 FEET
FIGURE 1
SITE LAYOUT
LUCK'S INCORPORATED
SEAGROVE, NORTH CAROLINA
McLAREN/HART
ENVIRONMENTAL ENGINEERING CORP.
WARREN, NEW JERSEY
DRWN: S.F.H.
CHK'D: S.S.
SCALE: AS SHOWN
DATE: 08/13/92

III. Summary of previous investigations

In August of 1990, a total of seven post-excavation soil samples (HA-3A, HA-3B, HA-4A, HA-4B, HA-5A, HA-5B and HA-5C) were collected from the base of the excavation at locations shown in Figure 2. The soil samples were collected from a depth of approximately 21.5 feet below grade by Spatco Environmental of Kernersville, North Carolina. The soil samples were analyzed by Law and Co. Laboratories for oil and grease using method 9071, as required by NCDEHNR for No 5 fuel oil USTs investigations. The oil and grease analytical results are summarized in Figure 2. Four samples (HB-3B, HB-4B, HB-5A and HB-5C) yielded concentrations of oil and grease ranging from 80 ppm (HA-5A) to 643 ppm (HB-3B).

At the request of NCDEHNR (letter dated September 11, 1991; provided in Attachment 2), and to establish the baseline concentration of naturally occurring oil and grease beneath the site, a background sample (HB-1-22') was collected by Spatco on October 23, 1991 and analyzed using method 9071 for oil and grease. Sample HB-1-22' was collected from a depth of 22 feet at a location away from existing underground storage tanks and manufacturing/processing areas, as shown on the site map. The background sample was analyzed by EnviroTech Mid-Atlantic of Blacksburg, Virginia. HB-1-22' yielded a background oil and grease concentration of 180 ppm.

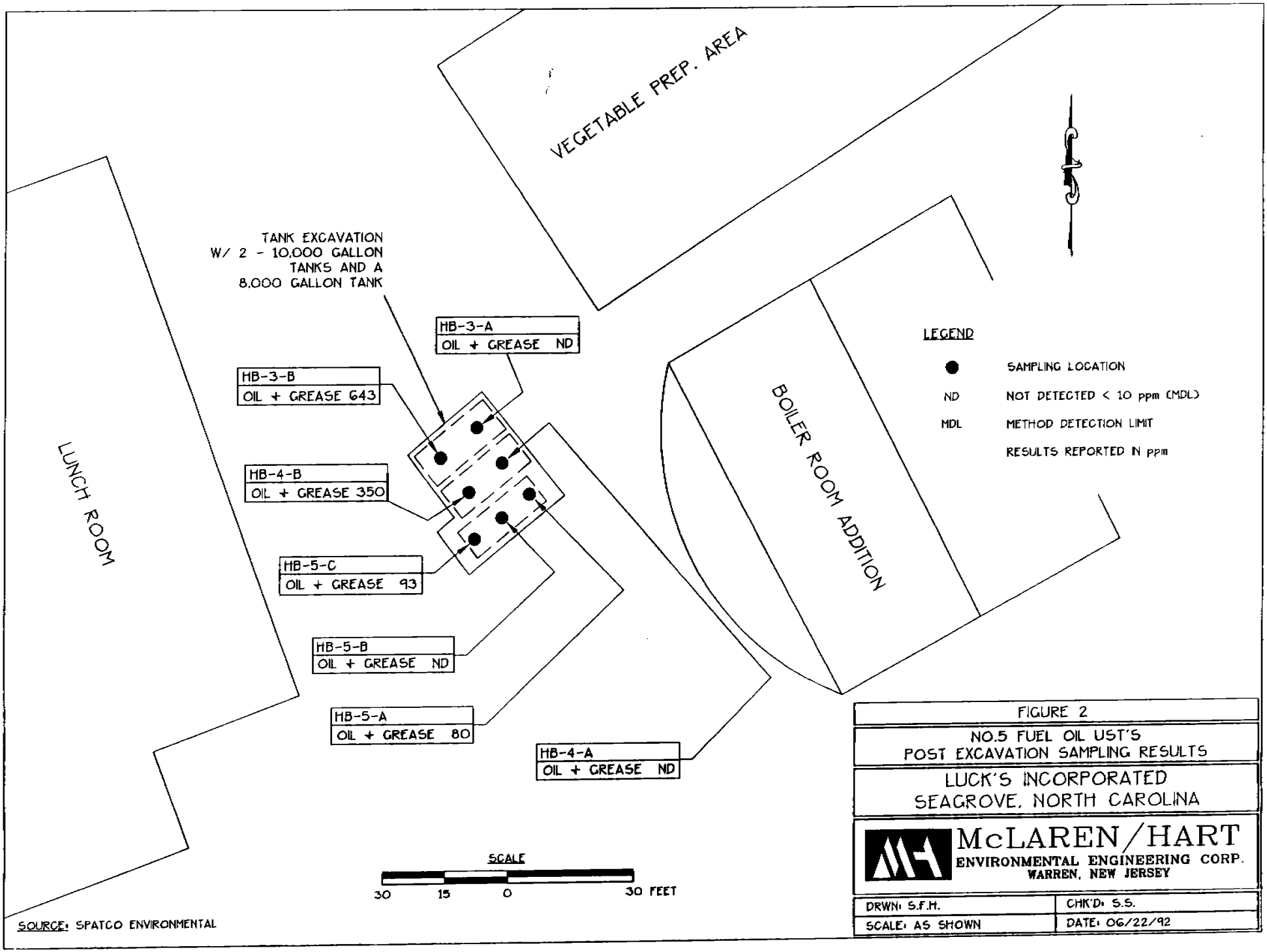
Based on the results of the background sample , NCDEHNR has assigned a remedial level of 190 ppm oil and grease for soils beneath the site (letter dated November 14, 1991; provided in Attachment 3). Two post-excavation soil samples, HB-3B (643 ppm oil and grease) and HB-4B (350 ppm oil and grease) exceeded the background concentration of 190 ppm oil and grease. As such, Luck's Inc was directed by NCDEHNR to conduct one of the following actions:

- excavate all soils with oil and grease concentrations greater than 190 ppm and conduct post-excavation sampling, or

- install a groundwater monitoring well downgradient of the tank excavation and sample annually for base neutral compounds using method 625, including a ten-peak library search.

IV. Current investigation

Based on analytical results from previous soil samples, it appeared that the detected concentrations of oil and grease may not be related to the USTs, but due to naturally occurring organic oil and grease. At the request of NCDEHNR, Luck's Inc., resampled soils in the area of the four samples (HB-3B, HB-4B, HB-5A and HB-5C) which indicated detectable levels of oil and grease. The four samples were analyzed for total petroleum hydrocarbons (TPH) using method 9073 and base neutrals using method 8270, as directed by NCDEHNR. Method 9073 is a non EPA approved method which involves a "natural oil and grease" clean-up process. According to NCDEHNR, method 9073 will indicate whether the detected levels of oil and grease are naturally occurring or from the No.5 fuel oil.


VEGETABLE PREP. AREA
LUNCH ROOM
BOILER ROOM ADDITION
TANK EXCAVATION
W/ 2 - 10,000 GALLON
TANKS AND A
8,000 GALLON TANK
HB-3-A
OIL + GREASE ND
HB-3-B
OIL + GREASE 643
HB-4-B
OIL + GREASE 350
HB-5-C
OIL + GREASE 93
HB-5-B
OIL + GREASE ND
HB-5-A
OIL + GREASE 80
HB-4-A
OIL + GREASE ND
LEGEND
SAMPLING LOCATION
ND NOT DETECTED < 10 ppm (MDL)
MDL METHOD DETECTION LIMIT
RESULTS REPORTED IN ppm
SCALE
30 15 0 30 FEET
SOURCE: SPATCO ENVIRONMENTAL
FIGURE 2
NO.5 FUEL OIL UST'S
POST EXCAVATION SAMPLING RESULTS
LUCK'S INCORPORATED
SEAGROVE, NORTH CAROLINA
McLAREN/HART
ENVIRONMENTAL ENGINEERING CORP.
WARREN, NEW JERSEY
DRWN: S.F.H.
CHK'D: S.S.
SCALE: AS SHOWN
DATE: 06/22/92

Four soil borings were completed using a hollow stem auger drill rig at locations shown in Figure 3. Soil samples **HB-3B-1, HB-4B-1, HB-5A-1** and **HB-5C-1** were collected from a depth of approximately 21.5 feet, at the former sample locations HB-3B, HB-4B, HB-5A and HB-5C, respectively. The soil samples were collected in accordance with the Standard Penetration Test Procedure (ASTM D 1586/67) using a properly decontaminated two-inch outer diameter split-spoon sampler. All split spoon samples were collected by McLaren/Hart Environmental Engineering Corp., of Warren, New Jersey.

In addition, one field blank (FB070792) was also collected and analyzed for the same parameters for QA/QC purposes. All analyses were performed by General Engineering Laboratories (GEL) of Charleston, South Carolina. GEL is certified by the state of North Carolina (certification No. 233).

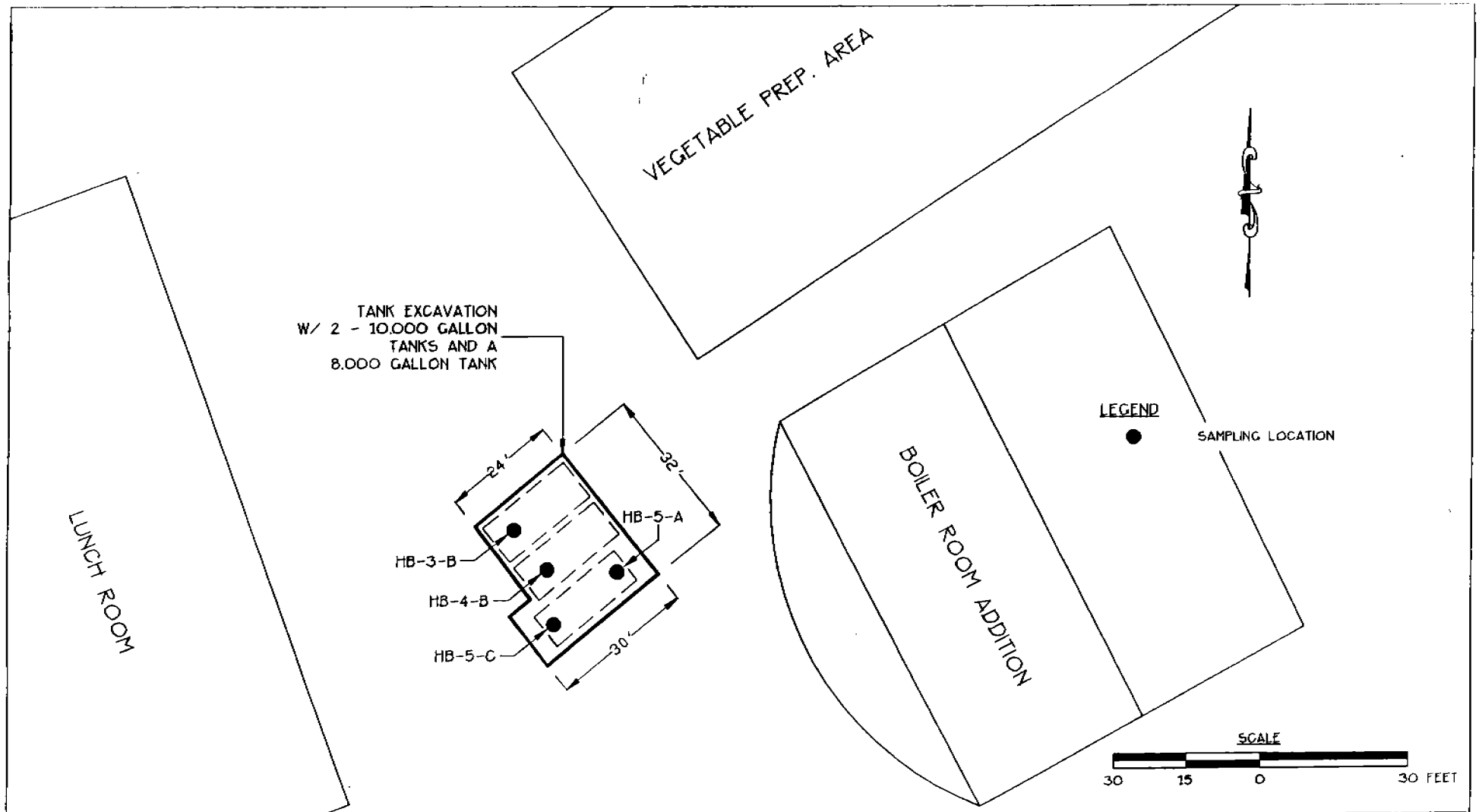
Soil samples HB-3B-1, HB-4B-1, HB-5A-1 and HB-5C-1 yielded detectable concentrations of TPH at 36 ppm, 204 ppm, 31 ppm, and 22 ppm, respectively. The field blank yielded non-detectable concentrations of TPH. TPH analytical results are presented in Table 1 and summarized in Figure 3. Complete laboratory analytical results are provided in Attachment 4.

Soil samples HB-3B-1, HB-4B-1, HB-5A-1 and HB-5C-1 and the field blank yielded non-detectable concentrations of targeted base neutral compounds. Further review of the chromatograms indicate that all five samples did not contain any non-targeted hydrocarbon compounds. Base neutrals analytical results are presented in Table 1 and summarized in Figure 3. Complete laboratory analytical results, including chromatograms for HB-3B-1 and HB-4B-1 are provided in Attachment 4.

V. Conclusions

A thorough investigation has been conducted to address the potential impact on soils from the operation of three No.5 fuel oil USTs. The tanks and surrounding soils were excavated and disposed of in accordance with applicable state and federal regulations. Verbal accounts from facility personnel indicated that the USTs were in good condition and did not appear to be leaking. Recent sample results indicate concentrations of TPH at 204 ppm (HB-4B) or less in the soils in this area. Analysis of these soil samples for base neutral compounds show non-detectable levels of these compounds. Based on these data, it is apparent that the soils in the area of the tanks are not contaminated with fuel oil. During the excavation there was no evidence of fuel product in the soils. In addition, there was no evidence of fuel oil odor from the soils in the excavation.

Considering the type of material stored in the USTs (No. 5 fuel oil) and its physical properties, any discharge which may have occurred from the USTs, would have remained in close proximity to the tank system and subsequently been removed during excavation. In addition, any discharge of the No. 5 fuel oil would be apparent both visually and by smell in the surrounding soils or in the samples collected; neither were observed.



Sample Number	HB-3B	HB-4B	HB-5A	HB-5C
Spatco, August 1990 Oil and Grease: method 9071	643	350	80	93
McLaren/Hart, July 1992 TPH: method 9073 Base neutrals: method 8270 (625)	36 ND	204 ND	31 ND	22 ND

Note:
Results reported in ppm, above MDL
ND:Not detected above method detection limits
Background concentration of oil and grease = 180 ppm yielded by sample HB-1-22 analyzed via method 9071, during October 1991.

FIGURE 3

NO.5 FUEL OIL UST'S EXCAVATION
SUMMARY OF SOIL ANALYTICAL RESULTS

LUCK'S INCORPORATED
SEAGROVE, NORTH CAROLINA

McLAREN/HART
ENVIRONMENTAL ENGINEERING CORP.
WARREN, NEW JERSEY

DRWN: S.F.H.	CHK'D: S.S.
SCALE: AS SHOWN	DATE: 08/13/92

TABLE 1
No.5 Fuel Oil USTs Excavation
Summary of Post Excavation Soil Sampling Results
Luck's Inc., Seagrove, NC

Sample Number	HB−3B	HB−4B	HB−5A	HB−5C
Sample Depth (ft)	21.5	21.5	21.5	21.5

Parameters (Results reported in parts per million)

Spatco, August 1990				
Oil and Grease: method 9071 *	643	350	80	93

Background concentration of oil and grease = 180 ppm ; yielded by sample HB−1−22 analyzed via method 9071, during October 1991.

McLaren/Hart, July 1992				
TPH: method 9073 **	36	204	31	22
Base neutrals: method 8270 (625)	ND	ND	ND	ND

Note:
* : Samples analyzed for oil and grease
** : TPH, less natural oil and grease
ND: Not detected above method detection limits

Because of the low levels of TPH detected, the low mobility of TPH in soils, the absence of base neutral compounds, and the depth to groundwater (approximately 43 feet below grade in the vicinity of the tank excavation), the concentrations of TPH detected do not represent an apparent contamination problem. Therefore no further investigation or excavation is warranted and tank closure activities are considered complete at this time.

VI. Recommendations

Based on a review of available hydrogeologic information, the physical properties of No.5 fuel oil, depth to groundwater in excess of 40 feet, and the analytical results of two rounds of post-excavation soil sampling, there is no indication of a contamination problem resulting from the operation of the underground tank farm. McLaren/Hart therefore recommends that this investigation be considered complete.

Attachment 1
Background monitoring well record

NORTH CAROLINA DEPARTMENT OF NATURAL RESOURCES & COMMUNITY DEVELOPMENT
DIVISION OF ENVIRONMENTAL MANAGEMENT
P. O. Box 27687 — RALEIGH, N.C. 27611 919-733-2020

WELL RECORD

DRILLING CONTRACTOR McCall Bros. REG. NO. 3 WELL CONSTRUCTION PERMIT NO. _____

1. WELL LOCATION: (Show sketch of the location below)
   Nearest Town: Seagrove N.C. County: _____
   Quadrangle No. _____
   (Road, Community or Subdivision and Lot No.)
2. OWNER: Lucks Foods
3. ADDRESS: Seagrove N.C.
4. TOPOGRAPHY: draw, valley, slope, hilltop, flat (circle one)
5. USE OF WELL: Monitoring DATE: 1-24-86
6. DOES THIS WELL REPLACE AN EXISTING WELL? No
7. TOTAL DEPTH: 50' RIG TYPE OR METHOD: Air
8. FORMATION SAMPLES COLLECTED: YES ____ NO X
9. CASING:

	Depth	Inside Dia.	Wall thick. or weight/ft.	type
From 0 to 45 ft		4"		PVC

10. GROUT:

	Depth	Material	Method
From 0 to 5 ft		Cement	Poured

11. SCREEN:

	Depth	Dia.	Type & Opening
From 45 to 50 ft		4"	020

12. GRAVEL:

	Depth	Size	Material
From 40 to 50 ft			Stone

13. WATER ZONES (depth): _____
14. STATIC WATER LEVEL: 40 ft. above/below top of casing
    Casing is ____ ft. above land surface ELEV: ____
15. YIELD (gpm): _____ METHOD OF TESTING: _____
16. PUMPING WATER LEVEL: _____ ft.
    after _____ hours at _____ gpm.
17. CHLORINATION: Type _____ Amount _____
18. WATER QUALITY: _____ TEMPERATURE (°F) ____
19. PERMANENT PUMP: Date Installed _____
    Type _____ Capacity _____ (gpm) HP _____
    Make _____ Intake Depth _____
    Airline Depth _____
20. HAS THE OWNER BEEN PROVIDED A COPY OF THIS RECORD AND INFORMED OF THE DEPARTMENTS REQUIREMENTS AND RECOMMENDATIONS? _____
21. REMARKS _____

DRILLING LOG

DEPTH FROM	TO	FORMATION DESCRIPTION
0	5'	Red Clay
5	50	Sand

If additional space is needed, use back of form

LOCATION SKETCH
(Show distance to numbered roads, or other map reference points)

I do hereby certify that this well was constructed in accordance with N.C. Well Construction Regulations and Standards and that this well record is true and exact.

McCall Brothers 3/25/86
SIGNATURE OF CONTRACTOR OR AGENT DATE

Submit original to Division of Environmental Management and copy to well owner

GW-1 Revised 10/1/80

Attachment 2
NCDEHNR correspondence
September 11, 1991

9-13-91 ok

State of North Carolina
Department of Environment, Health, and Natural Resources
Winston-Salem Regional Office

James G. Martin, Governor
William W. Cobey, Jr., Secretary

Margaret Plemmons Foster
Regional Manager

DIVISION OF ENVIRONMENTAL MANAGEMENT
GROUNDWATER SECTION

September 11, 1991

Darian Luck
Luck's Inc
P.O. Box 510
Highway 705 East
Seagrove, NC 27341-0510

SUBJECT: UST Closure at Luck's Plant, Highway 705 East, Seagrove, North Carolina

Dear Mr. Luck:

The Groundwater Section of the Division of Environmental Management at this office is presently reviewing the tank closure assessment for the subject location which was received in this office on June 8, 1990. We apologize for the delay in our response to your assessment. Due to reservation about the accuracy of EPA Test Method 9070 expressed to Mr. Steve Kay of the Groundwater Section by yourself and your consultant, we have been researching alternative soil sample test methods. The research of other methods, combined the overwhelming number of UST closures in the Winston-Salem Region, is responsible for our delay in responding.

If 9071 is to be used it will be necessary to take a back ground sample. The acceptable cleanup will be 100 PPM T O & G or background which ever is lower. If you wish to look into alternative test methods, there maybe one acceptable test. Please contact Connie Crosley at (919) 733-7015 for more information on acceptable test methods.

If you have any questions, please feel free to contact me at the letterhead address or telephone number.

Sincerely,

Thomas A. Salley

Thomas A. Salley
Hydrogeological Technician

EPA 625 Test 10 most Peaks

TAS/ahl
cc: WSRO

Attachment 3
NCDEHNR Correspondence
November 14, 1991

Attachment 4
Laboratory Analytical Results

GENERAL ENGINEERING LABORATORIES

Environmental Engineering and Analytical Services

Molly F. Greene
President

George C. Greene, P.E., Ph.D.
Vice President
SC Registration No. 9103

Laboratory	Certifications:
FL	E87156/87294
NC	233
SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali

Date: 08/03/92

Released by: QA/QC Officer

cc: MLRN00191 **Project Manager:** Winter Seibert **Page No.:** 1

Sample ID	: HB-3B-1
Lab ID	: 9207192-04
Matrix	: Soil
Date Collected	: 07/08/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Organic Prep

Evaporative Loss @ 105 C	21.0 wt%

Extractable Organics

Method 8270 Base/Neutral Compounds

1,2,4,5-Tetrachlorobenzene	< 330 ppb
1,2,4-Trichlorobenzene	< 330 ppb
1,2-Dichlorobenzene	< 330 ppb
1,3-Dichlorobenzene	< 330 ppb
1,4-Dichlorobenzene	< 330 ppb
1,4-Dioxane	< 422 ppb
1,4-Naphthoquinone	< 422 ppb
1-Naphthylamine	< 330 ppb
2,4-Dinitrotoluene	< 330 ppb
2,6-Dinitrotoluene	< 330 ppb
2-Acetylaminofluorene	< 747 ppb
2-Chloronaphthalene	< 330 ppb
2-Methylnaphthalene	< 330 ppb
2-Naphthylamine	< 330 ppb
2-Picoline	< 330 ppb
3,3'-Dichlorobenzidine	< 660 ppb
3,3'-Dimethylbenzidine	< 422 ppb
3-Methylcholanthrene	< 422 ppb
4,4'-DDD	< 422 ppb
4,4'-DDE	< 422 ppb
4,4'-DDT	< 422 ppb
4-Bromophenyl phenyl ether	< 422 ppb

P.O. Box 30712 • Charleston, SC 29417
Phone (803) 556-8171 • FAX (803) 766-1178

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Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

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Lab ID	: 9207192-04
Matrix	: Soil
Date Collected	: 07/08/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

4-Chloroaniline	< 330 ppb
4-Chlorophenyl phenyl ether	< 330 ppb
4-Nitroquinoline	< 1340 ppb
4-aminobiphenyl	< 422 ppb
5-Nitro-o-toluidine	< 330 ppb
7,12-Dimethylbenz(a)anthracene	< 464 ppb
Acenaphthene	< 330 ppb
Acenaphthylene	< 330 ppb
Acetophenone	< 330 ppb
Aldrin	< 709 ppb
Aniline	< 330 ppb
Anthracene	< 330 ppb
Aramite	< 330 ppb
Benzidine	< 422 ppb
Benzo(a)anthracene	< 330 ppb
Benzo(a)pyrene	< 330 ppb
Benzo(b)fluoranthene	< 395 ppb
Benzo(ghi)perylene	< 330 ppb
Benzo(k)fluoranthene	< 527 ppb
Benzyl Alcohol	< 330 ppb
Butyl benzyl phthalate	< 330 ppb
Chlordane	< 422 ppb
Chlorobenzilate	< 330 ppb
Chrysene	< 422 ppb
Di-n-butyl phthalate	< 422 ppb
Di-n-octyl phthalate	< 422 ppb
Diallate	< 330 ppb
Dibenzo(a,h)anthracene	< 330 ppb
Dibenzofuran	< 330 ppb
Dieldrin	< 772 ppb

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Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-3B-1
Lab ID	: 9207192-04
Matrix	: Soil
Date Collected	: 07/08/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Diethyl phthalate	< 330 ppb
Dimethoate	< 330 ppb
Dimethyl phthalate	< 330 ppb
Diphenylamine	< 330 ppb
Disulfoton	< 422 ppb
Endosulfan I	< 422 ppb
Endosulfan II	< 422 ppb
Endosulfan sulfate	< 751 ppb
Endrin	< 793 ppb
Endrin aldehyde	< 1440 ppb
Ethyl Methanesulfonate	< 422 ppb
Famphur	< 330 ppb
Fluoranthene	< 330 ppb
Fluorene	< 330 ppb
Heptachlor	< 941 ppb
Heptachlor epoxide	< 421 ppb
Hexachlorobenzene	< 330 ppb
Hexachlorobutadiene	< 330 ppb
Hexachlorocyclopentadiene	< 460 ppb
Hexachloroethane	< 330 ppb
Hexachloropropene	< 330 ppb
Indeno(1,2,3-c,d)pyrene	< 422 ppb
Isodrin	< 330 ppb
Isophorone	< 330 ppb
Isosafrole	< 330 ppb
Kepone	< 330 ppb
Methapyrilene	< 330 ppb
Methyl Methanesulfonate	< 422 ppb
N-Nitrosodi-n-butylamine	< 422 ppb
N-Nitrosodiethylamine	< 422 ppb

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Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

N-Nitrosodimethylamine	< 330 ppb
N-Nitrosodiphenylamine	< 422 ppb
N-Nitrosodipropylamine	< 330 ppb
N-Nitrosomethylethylamine	< 2950 ppb
N-Nitrosomorpholine	< 330 ppb
N-Nitrosopiperidine	< 330 ppb
N-Nitrosopyrolidine	< 330 ppb
Naphthalene	< 330 ppb
Nitrobenzene	< 330 ppb
O,O,O-Triethylphosphorothioate	< 422 ppb
Parathion, methyl	< 330 ppb
Pentachlorobenzene	< 330 ppb
Pentachloroethane	< 330 ppb
Pentachloronitrobenzene	< 330 ppb
Phenacetin	< 330 ppb
Phenanthrene	< 330 ppb
Phorate	< 330 ppb
Pronamide	< 330 ppb
Pyrene	< 330 ppb
Pyridine	< 789 ppb
Safrole	< 330 ppb
Sulfotepp	< 422 ppb
Thionazin	< 422 ppb
Toxaphene	< 422 ppb
Tributylphosphate	< 422 ppb
a-,a-Dimethylphenethylamine	< 422 ppb
alpha-BHC	< 422 ppb
beta-BHC	< 422 ppb
bis(2-Chloroethoxy)methane	< 330 ppb
bis(2-Chloroethyl) ether	< 330 ppb

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Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-3B-1
Lab ID	: 9207192-04
Matrix	: Soil
Date Collected	: 07/08/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

bis(2-Chloroisopropyl)ether	< 330 ppb
bis(2-Ethylhexyl)phthalate	< 422 ppb
delta-BHC	< 422 ppb
gamma-BHC	< 1220 ppb
m-Dinitrobenzene	< 330 ppb
m-Nitroaniline	< 2110 ppb
o-Nitroaniline	< 330 ppb
o-Toluidine	< 330 ppb
p-Dimethylaminoazobenzene	< 422 ppb
p-Nitroaniline	< 1810 ppb
p-Phenylenediamine	< 330 ppb
sym-Trinitrobenzene	< 422 ppb

General Chemistry

Total Petroleum Hydrocarbons	35.6 ppm

The following preparation procedures were performed:

Ext. & Conc. - B/N Compounds

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Sample: 9207192-04 Client ID: Matrix: Soil Description: HB-3B-1
Sampled: 07/08/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
TC: B/N EXT Method: EPA 8270					Batch:	Run Analyst:		DOA:	TOA:	NYD
TC: EVAP LOSS Method: EPA 160.3					Batch: 16959	Run Analyst: BDK		DOA: 07/11/92	TOA: 1200	OK
Evaporative Loss @ 105 C	21.0	wt%		0.00	wt%	1:1		0.00	0.00%	N/C
TC: 8270 B/N Method: EPA 8270					Batch: 17692	Run Analyst: AGW		DOA: 07/16/92	TOA: 1848	OK
						Prep Analyst: BDK		DOP: 07/15/92	TOP: 1600	OK
1,2,4,5-Tetrachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,2,4-Trichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,2-Dichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,3-Dichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,4-Dichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,4-Dioxane	0.00	ppm	<	422.	ppb	1:1		< 422.		
1,4-Naphthoquinone	0.00	ppm	<	422.	ppb	1:1		< 422.		
1-Naphthylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
2,4-Dinitrotoluene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	48.0%
2,6-Dinitrotoluene	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Acetylaminofluorene	0.00	ppm	<	747.	ppb	1:1		< 747.		
2-Chloronaphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Fluorobiphenyl	1.05	ppb		422.	ppb	1:1	49.8%	< 422.		
2-Methylnaphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Naphthylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Picoline	0.00	ppm	<	330.	ppb	1:1		< 330.		
3,3'-Dichlorobenzidine	0.00	ppm	<	660.	ppb	1:1		< 660.		
3,3'-Dimethylbenzidine	0.00	ppm	<	422.	ppb	1:1		< 422.		
3-Methylcholanthrene	0.00	ppm	<	422.	ppb	1:1		< 422.		
4,4'-DDD	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
4,4'-DDE	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
4,4'-DDT	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
4-Bromophenyl phenyl ether	0.00	ppm	<	422.	ppb	1:1		< 422.		
4-Chloroaniline	0.00	ppm	<	330.	ppb	1:1		< 330.		
4-Chlorophenyl phenyl ether	0.00	ppm	<	330.	ppb	1:1		< 330.		
4-Nitroquinoline	0.00	ppm	<	1340	ppb	1:1		< 1340		
4-aminobiphenyl	0.00	ppm	<	422.	ppb	1:1		< 422.		
5-Nitro-o-toluidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
7,12-Dimethylbenz(a)anthrace	0.00	ppm	<	464.	ppb	1:1		N/C		
Acenaphthene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	64.0%
Acenaphthylene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Acetophenone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Aldrin	0.00	ppm	<	709.	ppb	1:1		< 709.	0.00%	N/C
Aniline	0.00	ppm	<	330.	ppb	1:1		< 330.		
Anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Aramite	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzidine	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
Benzo(a)anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzo(a)pyrene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzo(b)fluoranthene	0.00	ppm	<	395.	ppb	1:1		< 395.		
Benzo(ghi)perylene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzo(k)fluoranthene	0.00	ppm	<	527.	ppb	1:1		< 527.		
Benzyl Alcohol	0.00	ppm	<	330.	ppb	1:1		< 330.		
Butyl benzyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Chlordane	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
Chlorobenzilate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Chrysene	0.00	ppm	<	422.	ppb	1:1		< 422.		
Di-n-butyl phthalate	0.00	ppm	<	422.	ppb	1:1		< 422.		
Di-n-octyl phthalate	0.00	ppm	<	422.	ppb	1:1		< 422.		

Sample: 9207192-04 Client ID: Matrix: Soil Description: HB-3B-1
Sampled: 07/08/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Diallate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzo(a,h)anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzofuran	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dieldrin	0.00	ppm	<	772.	ppb	1:1		< 772.	0.00%	N/C
Diethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethoate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Diphenylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Disulfoton	0.00	ppm	<	422.	ppb	1:1		< 422.		
Endosulfan I	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
Endosulfan II	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
Endosulfan sulfate	0.00	ppm	<	751.	ppb	1:1		< 751.	0.00%	N/C
Endrin	0.00	ppm	<	793.	ppb	1:1		< 793.	0.00%	N/C
Endrin aldehyde	0.00	ppm	<	1440	ppb	1:1		< 1440	0.00%	N/C
Ethyl Methanesulfonate	0.00	ppm	<	422.	ppb	1:1		< 422.		
Famphur	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluoranthene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluorene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Heptachlor	0.00	ppm	<	941.	ppb	1:1		< 941.	0.00%	N/C
Heptachlor epoxide	0.00	ppm	<	421.	ppb	1:1		< 421.	0.00%	N/C
Hexachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorobutadiene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorocyclopentadiene	0.00	ppm	<	460.	ppb	1:1		< 460.		
Hexachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachloropropene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Indeno(1,2,3-c,d)pyrene	0.00	ppm	<	422.	ppb	1:1		< 422.		
Isodrin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isophorone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isosafrole	0.00	ppm	<	330.	ppb	1:1		< 330.		
Kepone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methapyrilene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methyl Methanesulfonate	0.00	ppm	<	422.	ppb	1:1		< 422.		
N-Nitrosodi-n-butylamine	0.00	ppm	<	422.	ppb	1:1		< 422.		
N-Nitrosodiethylamine	0.00	ppm	<	422.	ppb	1:1		< 422.		
N-Nitrosodimethylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosodiphenylamine	0.00	ppm	<	422.	ppb	1:1		< 422.		
N-Nitrosodipropylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosomethylethylamine	0.00	ppm	<	2950	ppb	1:1		< 2950		
N-Nitrosomorpholine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopiperidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopyrolidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Naphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene-d5	0.759	ppb		422.	ppb	1:1	36.0%	< 422.		
O,O,O-Triethylphosphorothioa	0.00	ppm	<	422.	ppb	1:1		N/C		
Parathion, methyl	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloronitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenacetin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenanthrene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phorate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pronamide	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pyrene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	68.0%

Sample: 9207192-04 Client ID: Matrix: Soil Description: HB-3B-1
Sampled: 07/08/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Pyridine	0.00	ppm	<	789.	ppb	1:1		< 789.		
Safrole	0.00	ppm	<	330.	ppb	1:1		< 330.		
Sulfotepp	0.00	ppm	<	422.	ppb	1:1		< 422.		
Thionazin	0.00	ppm	<	422.	ppb	1:1		< 422.		
Toxaphene	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
Tributylphosphate	0.00	ppm	<	422.	ppb	1:1		< 422.		
a-,a-Dimethylphenethylamine	0.00	ppm	<	422.	ppb	1:1		< 422.		
alpha-BHC	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
beta-BHC	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
bis(2-Chloroethoxy)methane	0.00	ppm	<	330.	ppb	1:1		< 330.		
bis(2-Chloroethyl) ether	0.00	ppm	<	330.	ppb	1:1		< 330.		
bis(2-Chloroisopropyl)ether	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	N/C
bis(2-Ethylhexyl)phthalate	0.00	ppm	<	422.	ppb	1:1		< 422.		
delta-BHC	0.00	ppm	<	422.	ppb	1:1		< 422.	0.00%	N/C
gamma-BHC	0.00	ppm	<	1220	ppb	1:1		< 1220	0.00%	N/C
m-Dinitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
m-Nitroaniline	0.00	ppm	<	2110	ppb	1:1		< 2110		
o-Nitroaniline	0.00	ppm	<	330.	ppb	1:1		< 330.		
o-Toluidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
p-Dimethylaminoazobenzene	0.00	ppm	<	422.	ppb	1:1		< 422.		
p-Nitroaniline	0.00	ppm	<	1810	ppb	1:1		< 1810		
p-Phenylenediamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
p-Terphenyl-d14	1.69	ppb		422.	ppb	1:1	80.1%	< 422.		
sym-Trinitrobenzene	0.00	ppm	<	422.	ppb	1:1		< 422.		
TC: TPH-9073 Method: EPA 9073					Batch: 17388 Run Analyst: SJ DOA: 07/23/92 TOA: 1200 OK					
Total Petroleum Hydrocarbons	35.6	ppm		10.0	ppm	1:1		< 10.0	2.16%	103%

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SC Registration No. 9103

Laboratory	Certifications:
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NC	233
SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali

Date: 08/03/92

Released by: QA/QC Officer

cc: MLRN00191 **Project Manager:** Winter Seibert **Page No.:** 1

Sample ID	: HB-4B-1
Lab ID	: 9207192-03
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Organic Prep

Evaporative Loss @ 105 C	15.0 wt%

Extractable Organics

Method 8270 Base/Neutral Compounds

1,2,4,5-Tetrachlorobenzene	< 330 ppb
1,2,4-Trichlorobenzene	< 330 ppb
1,2-Dichlorobenzene	< 330 ppb
1,3-Dichlorobenzene	< 330 ppb
1,4-Dichlorobenzene	< 330 ppb
1,4-Dioxane	< 392 ppb
1,4-Naphthoquinone	< 392 ppb
1-Naphthylamine	< 330 ppb
2,4-Dinitrotoluene	< 330 ppb
2,6-Dinitrotoluene	< 330 ppb
2-Acetylaminofluorene	< 694 ppb
2-Chloronaphthalene	< 330 ppb
2-Methylnaphthalene	< 330 ppb
2-Naphthylamine	< 330 ppb
2-Picoline	< 330 ppb
3,3'-Dichlorobenzidine	< 660 ppb
3,3'-Dimethylbenzidine	< 392 ppb
3-Methylcholanthrene	< 392 ppb
4,4'-DDD	< 392 ppb
4,4'-DDE	< 392 ppb
4,4'-DDT	< 392 ppb
4-Bromophenyl phenyl ether	< 392 ppb

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TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali
cc: MLRN00191 **Project Manager:** Winter Seibert

Date: 08/03/92

Sample ID	: HB-4B-1
Lab ID	: 9207192-03
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

4-Chloroaniline	< 330 ppb
4-Chlorophenyl phenyl ether	< 330 ppb
4-Nitroquinoline	< 1250 ppb
4-aminobiphenyl	< 392 ppb
5-Nitro-o-toluidine	< 330 ppb
7,12-Dimethylbenz(a)anthracene	< 431 ppb
Acenaphthene	< 330 ppb
Acenaphthylene	< 330 ppb
Acetophenone	< 330 ppb
Aldrin	< 659 ppb
Aniline	< 330 ppb
Anthracene	< 330 ppb
Aramite	< 330 ppb
Benzidine	< 392 ppb
Benzo(a)anthracene	< 330 ppb
Benzo(a)pyrene	< 330 ppb
Benzo(b)fluoranthene	< 367 ppb
Benzo(ghi)perylene	< 330 ppb
Benzo(k)fluoranthene	< 490 ppb
Benzyl Alcohol	< 330 ppb
Butyl benzyl phthalate	< 330 ppb
Chlordane	< 392 ppb
Chlorobenzilate	< 330 ppb
Chrysene	< 392 ppb
Di-n-butyl phthalate	< 392 ppb
Di-n-octyl phthalate	< 392 ppb
Diallate	< 330 ppb
Dibenzo(a,h)anthracene	< 330 ppb
Dibenzofuran	< 330 ppb
Dieldrin	< 718 ppb

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WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-4B-1
Lab ID	: 9207192-03
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Diethyl phthalate	< 330 ppb
Dimethoate	< 330 ppb
Dimethyl phthalate	< 330 ppb
Diphenylamine	< 330 ppb
Disulfoton	< 392 ppb
Endosulfan I	< 392 ppb
Endosulfan II	< 392 ppb
Endosulfan sulfate	< 698 ppb
Endrin	< 737 ppb
Endrin aldehyde	< 1340 ppb
Ethyl Methanesulfonate	< 392 ppb
Famphur	< 330 ppb
Fluoranthene	< 330 ppb
Fluorene	< 330 ppb
Heptachlor	< 875 ppb
Heptachlor epoxide	< 391 ppb
Hexachlorobenzene	< 330 ppb
Hexachlorobutadiene	< 330 ppb
Hexachlorocyclopentadiene	< 427 ppb
Hexachloroethane	< 330 ppb
Hexachloropropene	< 330 ppb
Indeno(1,2,3-c,d)pyrene	< 392 ppb
Isodrin	< 330 ppb
Isophorone	< 330 ppb
Isosafrole	< 330 ppb
Kepone	< 330 ppb
Methapyrilene	< 330 ppb
Methyl Methanesulfonate	< 392 ppb
N-Nitrosodi-n-butylamine	< 392 ppb
N-Nitrosodiethylamine	< 392 ppb

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TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali
cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-4B-1
Lab ID	: 9207192-03
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

N-Nitrosodimethylamine	< 330 ppb
N-Nitrosodiphenylamine	< 392 ppb
N-Nitrosodipropylamine	< 330 ppb
N-Nitrosomethylethylamine	< 2750 ppb
N-Nitrosomorpholine	< 330 ppb
N-Nitrosopiperidine	< 330 ppb
N-Nitrosopyrolidine	< 330 ppb
Naphthalene	< 330 ppb
Nitrobenzene	< 330 ppb
O,O,O-Triethylphosphorothioate	< 392 ppb
Parathion, methyl	< 330 ppb
Pentachlorobenzene	< 330 ppb
Pentachloroethane	< 330 ppb
Pentachloronitrobenzene	< 330 ppb
Phenacetin	< 330 ppb
Phenanthrene	< 330 ppb
Phorate	< 330 ppb
Pronamide	< 330 ppb
Pyrene	< 330 ppb
Pyridine	< 733 ppb
Safrole	< 330 ppb
Sulfotepp	< 392 ppb
Thionazin	< 392 ppb
Toxaphene	< 392 ppb
Tributylphosphate	< 392 ppb
a-,a-Dimethylphenethylamine	< 392 ppb
alpha-BHC	< 392 ppb
beta-BHC	< 392 ppb
bis(2-Chloroethoxy)methane	< 330 ppb
bis(2-Chloroethyl) ether	< 330 ppb

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FL	E87156/87294
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SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-4B-1
Lab ID	: 9207192-03
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

bis(2-Chloroisopropyl)ether	< 330 ppb
bis(2-Ethylhexyl)phthalate	745 ppb
delta-BHC	< 392 ppb
gamma-BHC	< 1130 ppb
m-Dinitrobenzene	< 330 ppb
m-Nitroaniline	< 1960 ppb
o-Nitroaniline	< 330 ppb
o-Toluidine	< 330 ppb
p-Dimethylaminoazobenzene	< 392 ppb
p-Nitroaniline	< 1690 ppb
p-Phenylenediamine	< 330 ppb
sym-Trinitrobenzene	< 392 ppb

General Chemistry

Total Petroleum Hydrocarbons	204 ppm

The following preparation procedures were performed:

Ext. & Conc. - B/N Compounds

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Sample: 9207192-03 Client ID: Matrix: Soil Description: M8-48-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units	CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
TC: B/N EXT Method: EPA 8270				Batch:	Run	Analyst:	DOA:	TOA:	NYD
TC: EVAP LOSS Method: EPA 160.3				Batch: 16959	Run	Analyst: BDK	DOA: 07/11/92	TOA: 1200	OK
Evaporative Loss @ 105 C	15.0	wt%	0.00	wt%	1:1		0.00	0.00%	N/C
TC: 8270 B/N Method: EPA 8270				Batch: 17692	Run	Analyst: AGW	DOA: 07/16/92	TOA: 1756	OK
					Prep	Analyst: BDK	DOP: 07/15/92	TOP: 1600	NOT OK
1,2,4,5-Tetrachlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,2,4-Trichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,2-Dichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,3-Dichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,4-Dichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,4-Dioxane	0.00	ppm	< 392.	ppb	1:1		< 392.		
1,4-Naphthoquinone	0.00	ppm	< 392.	ppb	1:1		< 392.		
1-Naphthylamine	0.00	ppm	< 330.	ppb	1:1		< 330.		
2,4-Dinitrotoluene	0.00	ppm	< 330.	ppb	1:1		< 330.	0.00%	48.0%
2,6-Dinitrotoluene	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Acetylaminofluorene	0.00	ppm	< 694.	ppb	1:1		< 694.		
2-Chloronaphthalene	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Fluorobiphenyl	1.33	ppb	392.	ppb	1:1	67.8%	< 392.		
2-Methylnaphthalene	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Naphthylamine	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Picoline	0.00	ppm	< 330.	ppb	1:1		< 330.		
3,3'-Dichlorobenzidine	0.00	ppm	< 660.	ppb	1:1		< 660.		
3,3'-Dimethylbenzidine	0.00	ppm	< 392.	ppb	1:1		< 392.		
3-Methylcholanthrene	0.00	ppm	< 392.	ppb	1:1		< 392.		
4,4'-DDD	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
4,4'-DDE	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
4,4'-DDT	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
4-Bromophenyl phenyl ether	0.00	ppm	< 392.	ppb	1:1		< 392.		
4-Chloroaniline	0.00	ppm	< 330.	ppb	1:1		< 330.		
4-Chlorophenyl phenyl ether	0.00	ppm	< 330.	ppb	1:1		< 330.		
4-Nitroquinoline	0.00	ppm	< 1250	ppb	1:1		< 1250		
4-aminobiphenyl	0.00	ppm	< 392.	ppb	1:1		< 392.		
5-Nitro-o-toluidine	0.00	ppm	< 330.	ppb	1:1		< 330.		
7,12-Dimethylbenz(a)anthrace	0.00	ppm	< 431.	ppb	1:1		N/C		
Acenaphthene	0.00	ppm	< 330.	ppb	1:1		< 330.	0.00%	64.0%
Acenaphthylene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Acetophenone	0.00	ppm	< 330.	ppb	1:1		< 330.		
Aldrin	0.00	ppm	< 659.	ppb	1:1		< 659.	0.00%	N/C
Aniline	0.00	ppm	< 330.	ppb	1:1		< 330.		
Anthracene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Aramite	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzidine	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
Benzo(a)anthracene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzo(a)pyrene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzo(b)fluoranthene	0.00	ppm	< 367.	ppb	1:1		< 367.		
Benzo(ghi)perylene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzo(k)fluoranthene	0.00	ppm	< 490.	ppb	1:1		< 490.		
Benzyl Alcohol	0.00	ppm	< 330.	ppb	1:1		< 330.		
Butyl benzyl phthalate	0.00	ppm	< 330.	ppb	1:1		< 330.		
Chlordane	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
Chlorobenzilate	0.00	ppm	< 330.	ppb	1:1		< 330.		
Chrysene	0.00	ppm	< 392.	ppb	1:1		< 392.		
Di-n-butyl phthalate	0.00	ppm	< 392.	ppb	1:1		< 392.		
Di-n-octyl phthalate	0.00	ppm	< 392.	ppb	1:1		< 392.		

Sample: 9207192-03 Client ID: Matrix: Soil Description: HB-4B-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Diallate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzo(a,h)anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzofuran	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dieldrin	0.00	ppm	<	718.	ppb	1:1		< 718.	0.00%	N/C
Diethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethoate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Diphenylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Disulfoton	0.00	ppm	<	392.	ppb	1:1		< 392.		
Endosulfan I	0.00	ppm	<	392.	ppb	1:1		< 392.	0.00%	N/C
Endosulfan II	0.00	ppm	<	392.	ppb	1:1		< 392.	0.00%	N/C
Endosulfan sulfate	0.00	ppm	<	698.	ppb	1:1		< 698.	0.00%	N/C
Endrin	0.00	ppm	<	737.	ppb	1:1		< 737.	0.00%	N/C
Endrin aldehyde	0.00	ppm	<	1340	ppb	1:1		< 1340	0.00%	N/C
Ethyl Methanesulfonate	0.00	ppm	<	392.	ppb	1:1		< 392.		
Famphur	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluoranthene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluorene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Heptachlor	0.00	ppm	<	875.	ppb	1:1		< 875.	0.00%	N/C
Heptachlor epoxide	0.00	ppm	<	391.	ppb	1:1		< 391.	0.00%	N/C
Hexachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorobutadiene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorocyclopentadiene	0.00	ppm	<	427.	ppb	1:1		< 427.		
Hexachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachloropropene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Indeno(1,2,3-c,d)pyrene	0.00	ppm	<	392.	ppb	1:1		< 392.		
Isodrin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isophorone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isosafrole	0.00	ppm	<	330.	ppb	1:1		< 330.		
Kepone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methapyrilene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methyl Methanesulfonate	0.00	ppm	<	392.	ppb	1:1		< 392.		
N-Nitrosodi-n-butylamine	0.00	ppm	<	392.	ppb	1:1		< 392.		
N-Nitrosodiethylamine	0.00	ppm	<	392.	ppb	1:1		< 392.		
N-Nitrosodimethylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosodiphenylamine	0.00	ppm	<	392.	ppb	1:1		< 392.		
N-Nitrosodipropylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosomethylethylamine	0.00	ppm	<	2750	ppb	1:1		< 2750		
N-Nitrosomorpholine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopiperidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopyrolidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Naphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene-d5	1.18	ppb		392.	ppb	1:1	60.2%	< 392.		
O,O,O-Triethylphosphorothioa	0.00	ppm	<	392.	ppb	1:1		N/C		
Parathion, methyl	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloronitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenacetin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenanthrene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phorate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pronamide	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pyrene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	68.0%

Sample: 9207192-03 Client ID: Matrix: Soil Description: HB-4B-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units	CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Pyridine	0.00	ppm	< 733.	ppb	1:1		< 733.		
Safrole	0.00	ppm	< 330.	ppb	1:1		< 330.		
Sulfotepp	0.00	ppm	< 392.	ppb	1:1		< 392.		
Thionazin	0.00	ppm	< 392.	ppb	1:1		< 392.		
Toxaphene	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
Tributylphosphate	0.00	ppm	< 392.	ppb	1:1		< 392.		
a-,a-Dimethylphenethylamine	0.00	ppm	< 392.	ppb	1:1		< 392.		
alpha-BHC	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
beta-BHC	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
bis(2-Chloroethoxy)methane	0.00	ppm	< 330.	ppb	1:1		< 330.		
bis(2-Chloroethyl) ether	0.00	ppm	< 330.	ppb	1:1		< 330.		
bis(2-Chloroisopropyl)ether	0.00	ppm	< 330.	ppb	1:1		< 330.	0.00%	N/C
bis(2-Ethylhexyl)phthalate	0.745	ppm	< 392.	ppb	1:1		< 392.		
delta-BHC	0.00	ppm	< 392.	ppb	1:1		< 392.	0.00%	N/C
gamma-BHC	0.00	ppm	< 1130	ppb	1:1		< 1130	0.00%	N/C
m-Dinitrobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
m-Nitroaniline	0.00	ppm	< 1960	ppb	1:1		< 1960		
o-Nitroaniline	0.00	ppm	< 330.	ppb	1:1		< 330.		
o-Toluidine	0.00	ppm	< 330.	ppb	1:1		< 330.		
p-Dimethylaminoazobenzene	0.00	ppm	< 392.	ppb	1:1		< 392.		
p-Nitroaniline	0.00	ppm	< 1690	ppb	1:1		< 1690		
p-Phenylenediamine	0.00	ppm	< 330.	ppb	1:1		< 330.		
p-Terphenyl-d14	1.69	ppb	392.	ppb	1:1	86.2%	< 392.		
sym-Trinitrobenzene	0.00	ppm	< 392.	ppb	1:1		< 392.		
TC: TPH-9073 Method: EPA 9073				Batch: 17388	Run Analyst:	SJ	DOA: 07/23/92	TOA: 1200	OK
Total Petroleum Hydrocarbons	204.	ppm	10.0	ppm	1:1		< 10.0	2.16%	103%

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NC	233
SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali

Date: 08/03/92

Released by: QA/QC Officer

cc: MLRN00191 **Project Manager:** Winter Seibert **Page No.:** 1

Sample ID	: HB-5C-1
Lab ID	: 9207192-02
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Organic Prep

Evaporative Loss @ 105 C	19.0 wt%

Extractable Organics

Method 8270 Base/Neutral Compounds

1,2,4,5-Tetrachlorobenzene	< 330 ppb
1,2,4-Trichlorobenzene	< 330 ppb
1,2-Dichlorobenzene	< 330 ppb
1,3-Dichlorobenzene	< 330 ppb
1,4-Dichlorobenzene	< 330 ppb
1,4-Dioxane	< 412 ppb
1,4-Naphthoquinone	< 412 ppb
1-Naphthylamine	< 330 ppb
2,4-Dinitrotoluene	< 330 ppb
2,6-Dinitrotoluene	< 330 ppb
2-Acetylaminofluorene	< 728 ppb
2-Chloronaphthalene	< 330 ppb
2-Methylnaphthalene	< 330 ppb
2-Naphthylamine	< 330 ppb
2-Picoline	< 330 ppb
3,3'-Dichlorobenzidine	< 660 ppb
3,3'-Dimethylbenzidine	< 412 ppb
3-Methylcholanthrene	< 412 ppb
4,4'-DDD	< 412 ppb
4,4'-DDE	< 412 ppb
4,4'-DDT	< 412 ppb
4-Bromophenyl phenyl ether	< 412 ppb

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WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali
cc: MLRN00191 **Project Manager:** Winter Seibert

Date: 08/03/92

Sample ID	: HB-5C-1
Lab ID	: 9207192-02
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

4-Chloroaniline	< 330 ppb
4-Chlorophenyl phenyl ether	< 330 ppb
4-Nitroquinoline	< 1310 ppb
4-aminobiphenyl	< 412 ppb
5-Nitro-o-toluidine	< 330 ppb
7,12-Dimethylbenz(a)anthracene	< 453 ppb
Acenaphthene	< 330 ppb
Acenaphthylene	< 330 ppb
Acetophenone	< 330 ppb
Aldrin	< 691 ppb
Aniline	< 330 ppb
Anthracene	< 330 ppb
Aramite	< 330 ppb
Benzidine	< 412 ppb
Benzo(a)anthracene	< 330 ppb
Benzo(a)pyrene	< 330 ppb
Benzo(b)fluoranthene	< 385 ppb
Benzo(ghi)perylene	< 330 ppb
Benzo(k)fluoranthene	< 514 ppb
Benzyl Alcohol	< 330 ppb
Butyl benzyl phthalate	< 330 ppb
Chlordane	< 412 ppb
Chlorobenzilate	< 330 ppb
Chrysene	< 412 ppb
Di-n-butyl phthalate	< 412 ppb
Di-n-octyl phthalate	< 412 ppb
Diallate	< 330 ppb
Dibenzo(a,h)anthracene	< 330 ppb
Dibenzofuran	< 330 ppb
Dieldrin	< 753 ppb

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TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert **Page No.:** 3

Sample ID	: HB-5C-1
Lab ID	: 9207192-02
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Diethyl phthalate	< 330 ppb
Dimethoate	< 330 ppb
Dimethyl phthalate	< 330 ppb
Diphenylamine	< 330 ppb
Disulfoton	< 412 ppb
Endosulfan I	< 412 ppb
Endosulfan II	< 412 ppb
Endosulfan sulfate	< 733 ppb
Endrin	< 774 ppb
Endrin aldehyde	< 1410 ppb
Ethyl Methanesulfonate	< 412 ppb
Famphur	< 330 ppb
Fluoranthene	< 330 ppb
Fluorene	< 330 ppb
Heptachlor	< 918 ppb
Heptachlor epoxide	< 411 ppb
Hexachlorobenzene	< 330 ppb
Hexachlorobutadiene	< 330 ppb
Hexachlorocyclopentadiene	< 449 ppb
Hexachloroethane	< 330 ppb
Hexachloropropene	< 330 ppb
Indeno(1,2,3-c,d)pyrene	< 412 ppb
Isodrin	< 330 ppb
Isophorone	< 330 ppb
Isosafrole	< 330 ppb
Kepone	< 330 ppb
Methapyrilene	< 330 ppb
Methyl Methanesulfonate	< 412 ppb
N-Nitrosodi-n-butylamine	< 412 ppb
N-Nitrosodiethylamine	< 412 ppb

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NC	233
SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali
cc: MLRN00191 Project Manager: Winter Seibert

Date: 08/03/92

Page No.: 4

Sample ID	: HB-5C-1
Lab ID	: 9207192-02
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

N-Nitrosodimethylamine	< 330 ppb
N-Nitrosodiphenylamine	< 412 ppb
N-Nitrosodipropylamine	< 330 ppb
N-Nitrosomethylethylamine	< 2880 ppb
N-Nitrosomorpholine	< 330 ppb
N-Nitrosopiperidine	< 330 ppb
N-Nitrosopyrolidine	< 330 ppb
Naphthalene	< 330 ppb
Nitrobenzene	< 330 ppb
O,O,O-Triethylphosphorothioate	< 412 ppb
Parathion, methyl	< 330 ppb
Pentachlorobenzene	< 330 ppb
Pentachloroethane	< 330 ppb
Pentachloronitrobenzene	< 330 ppb
Phenacetin	< 330 ppb
Phenanthrene	< 330 ppb
Phorate	< 330 ppb
Pronamide	< 330 ppb
Pyrene	< 330 ppb
Pyridine	< 770 ppb
Safrole	< 330 ppb
Sulfotepp	< 412 ppb
Thionazin	< 412 ppb
Toxaphene	< 412 ppb
Tributylphosphate	< 412 ppb
a-,a-Dimethylphenethylamine	< 412 ppb
alpha-BHC	< 412 ppb
beta-BHC	< 412 ppb
bis(2-Chloroethoxy)methane	< 330 ppb
bis(2-Chloroethyl) ether	< 330 ppb

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GENERAL ENGINEERING LABORATORIES

Environmental Engineering and Analytical Services

Molly F. Greene
President

George C. Greene, P.E., Ph.D.
Vice President
SC Registration No. 9103

Laboratory	Certifications:
FL	E87156/87294
NC	233
SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-5C-1
Lab ID	: 9207192-02
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

bis(2-Chloroisopropyl)ether	< 330 ppb
bis(2-Ethylhexyl)phthalate	< 412 ppb
delta-BHC	< 412 ppb
gamma-BHC	< 1190 ppb
m-Dinitrobenzene	< 330 ppb
m-Nitroaniline	< 2060 ppb
o-Nitroaniline	< 330 ppb
o-Toluidine	< 330 ppb
p-Dimethylaminoazobenzene	< 412 ppb
p-Nitroaniline	< 1770 ppb
p-Phenylenediamine	< 330 ppb
sym-Trinitrobenzene	< 412 ppb

General Chemistry

Total Petroleum Hydrocarbons	21.7 ppm

The following preparation procedures were performed:

Ext. & Conc. - B/N Compounds

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Sample: 9207192-02 Client ID: Matrix: Soil Description: MB-5C-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units	CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
TC: B/N EXT Method: EPA 8270				Batch:	Run	Analyst:	DOA:	TOA:	NYD
TC: EVAP LOSS Method: EPA 160.3				Batch: 16959	Run	Analyst: BDK	DOA: 07/11/92	TOA: 1200	OK
Evaporative Loss @ 105 C	19.0	wt%	0.00	wt%	1:1		0.00	0.00%	N/C
TC: 8270 B/N Method: EPA 8270				Batch: 17692	Run	Analyst: AGW	DOA: 07/16/92	TOA: 1704	OK
					Prep	Analyst: BDK	DOP: 07/15/92	TOP: 1600	NOT OK
1,2,4,5-Tetrachlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,2,4-Trichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,2-Dichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,3-Dichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,4-Dichlorobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
1,4-Dioxane	0.00	ppm	< 412.	ppb	1:1		< 412.		
1,4-Naphthoquinone	0.00	ppm	< 412.	ppb	1:1		< 412.		
1-Naphthylamine	0.00	ppm	< 330.	ppb	1:1		< 330.		
2,4-Dinitrotoluene	0.00	ppm	< 330.	ppb	1:1		< 330.	0.00%	48.0%
2,6-Dinitrotoluene	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Acetylaminofluorene	0.00	ppm	< 728.	ppb	1:1		< 728.		
2-Chloronaphthalene	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Fluorobiphenyl	1.48	ppb	412.	ppb	1:1	71.9%	< 412.		
2-Methylnaphthalene	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Naphthylamine	0.00	ppm	< 330.	ppb	1:1		< 330.		
2-Picoline	0.00	ppm	< 330.	ppb	1:1		< 330.		
3,3'-Dichlorobenzidine	0.00	ppm	< 660.	ppb	1:1		< 660.		
3,3'-Dimethylbenzidine	0.00	ppm	< 412.	ppb	1:1		< 412.		
3-Methylcholanthrene	0.00	ppm	< 412.	ppb	1:1		< 412.		
4,4'-DDD	0.00	ppm	< 412.	ppb	1:1		< 412.	0.00%	N/C
4,4'-DDE	0.00	ppm	< 412.	ppb	1:1		< 412.	0.00%	N/C
4,4'-DDT	0.00	ppm	< 412.	ppb	1:1		< 412.	0.00%	N/C
4-Bromophenyl phenyl ether	0.00	ppm	< 412.	ppb	1:1		< 412.		
4-Chloroaniline	0.00	ppm	< 330.	ppb	1:1		< 330.		
4-Chlorophenyl phenyl ether	0.00	ppm	< 330.	ppb	1:1		< 330.		
4-Nitroquinoline	0.00	ppm	< 1310	ppb	1:1		< 1310		
4-aminobiphenyl	0.00	ppm	< 412.	ppb	1:1		< 412.		
5-Nitro-o-toluidine	0.00	ppm	< 330.	ppb	1:1		< 330.		
7,12-Dimethylbenz(a)anthrace	0.00	ppm	< 453.	ppb	1:1		N/C		
Acenaphthene	0.00	ppm	< 330.	ppb	1:1		< 330.	0.00%	64.0%
Acenaphthylene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Acetophenone	0.00	ppm	< 330.	ppb	1:1		< 330.		
Aldrin	0.00	ppm	< 691.	ppb	1:1		< 691.	0.00%	N/C
Aniline	0.00	ppm	< 330.	ppb	1:1		< 330.		
Anthracene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Aramite	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzidine	0.00	ppm	< 412.	ppb	1:1		< 412.	0.00%	N/C
Benzo(a)anthracene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzo(a)pyrene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzo(b)fluoranthene	0.00	ppm	< 385.	ppb	1:1		< 385.		
Benzo(ghi)perylene	0.00	ppm	< 330.	ppb	1:1		< 330.		
Benzo(k)fluoranthene	0.00	ppm	< 514.	ppb	1:1		< 514.		
Benzyl Alcohol	0.00	ppm	< 330.	ppb	1:1		< 330.		
Butyl benzyl phthalate	0.00	ppm	< 330.	ppb	1:1		< 330.		
Chlordane	0.00	ppm	< 412.	ppb	1:1		< 412.	0.00%	N/C
Chlorobenzilate	0.00	ppm	< 330.	ppb	1:1		< 330.		
Chrysene	0.00	ppm	< 412.	ppb	1:1		< 412.		
Di-n-butyl phthalate	0.00	ppm	< 412.	ppb	1:1		< 412.		
Di-n-octyl phthalate	0.00	ppm	< 412.	ppb	1:1		< 412.		

Sample: 9207192-02 Client ID: Matrix: Soil Description: HB-5C-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Diallate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzo(a,h)anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzofuran	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dieldrin	0.00	ppm	<	753.	ppb	1:1		< 753.	0.00%	N/C
Diethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethoate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Diphenylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Disulfoton	0.00	ppm	<	412.	ppb	1:1		< 412.		
Endosulfan I	0.00	ppm	<	412.	ppb	1:1		< 412.	0.00%	N/C
Endosulfan II	0.00	ppm	<	412.	ppb	1:1		< 412.	0.00%	N/C
Endosulfan sulfate	0.00	ppm	<	733.	ppb	1:1		< 733.	0.00%	N/C
Endrin	0.00	ppm	<	774.	ppb	1:1		< 774.	0.00%	N/C
Endrin aldehyde	0.00	ppm	<	1410	ppb	1:1		< 1410	0.00%	N/C
Ethyl Methanesulfonate	0.00	ppm	<	412.	ppb	1:1		< 412.		
Famphur	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluoranthene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluorene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Heptachlor	0.00	ppm	<	918.	ppb	1:1		< 918.	0.00%	N/C
Heptachlor epoxide	0.00	ppm	<	411.	ppb	1:1		< 411.	0.00%	N/C
Hexachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorobutadiene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorocyclopentadiene	0.00	ppm	<	449.	ppb	1:1		< 449.		
Hexachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachloropropene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Indeno(1,2,3-c,d)pyrene	0.00	ppm	<	412.	ppb	1:1		< 412.		
Isodrin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isophorone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isosafrole	0.00	ppm	<	330.	ppb	1:1		< 330.		
Kepone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methapyrilene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methyl Methanesulfonate	0.00	ppm	<	412.	ppb	1:1		< 412.		
N-Nitrosodi-n-butylamine	0.00	ppm	<	412.	ppb	1:1		< 412.		
N-Nitrosodiethylamine	0.00	ppm	<	412.	ppb	1:1		< 412.		
N-Nitrosodimethylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosodiphenylamine	0.00	ppm	<	412.	ppb	1:1		< 412.		
N-Nitrosodipropylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosomethylethylamine	0.00	ppm	<	2880	ppb	1:1		< 2880		
N-Nitrosomorpholine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopiperidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopyrolidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Naphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene-d5	1.32	ppb		412.	ppb	1:1	64.2%	< 412.		
O,O,O-Triethylphosphorothioa	0.00	ppm	<	412.	ppb	1:1		N/C		
Parathion, methyl	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloronitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenacetin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenanthrene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phorate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pronamide	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pyrene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	68.0%

Sample: 9207192-02 Client ID: Matrix: Soil Description: MB-5C-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Pyridine	0.00	ppm	<	770.	ppb	1:1		< 770.		
Safrole	0.00	ppm	<	330.	ppb	1:1		< 330.		
Sulfotepp	0.00	ppm	<	412.	ppb	1:1		< 412.		
Thionazin	0.00	ppm	<	412.	ppb	1:1		< 412.		
Toxaphene	0.00	ppm	<	412.	ppb	1:1		< 412.	0.00%	N/C
Tributylphosphate	0.00	ppm	<	412.	ppb	1:1		< 412.		
a-,a-Dimethylphenethylamine	0.00	ppm	<	412.	ppb	1:1		< 412.		
alpha-BHC	0.00	ppm	<	412.	ppb	1:1		< 412.	0.00%	N/C
beta-BHC	0.00	ppm	<	412.	ppb	1:1		< 412.	0.00%	N/C
bis(2-Chloroethoxy)methane	0.00	ppm	<	330.	ppb	1:1		< 330.		
bis(2-Chloroethyl) ether	0.00	ppm	<	330.	ppb	1:1		< 330.		
bis(2-Chloroisopropyl)ether	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	N/C
bis(2-Ethylhexyl)phthalate	0.00	ppm	<	412.	ppb	1:1		< 412.		
delta-BHC	0.00	ppm	<	412.	ppb	1:1		< 412.	0.00%	N/C
gamma-BHC	0.00	ppm	<	1190	ppb	1:1		< 1190	0.00%	N/C
m-Dinitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
m-Nitroaniline	0.00	ppm	<	2060	ppb	1:1		< 2060		
o-Nitroaniline	0.00	ppm	<	330.	ppb	1:1		< 330.		
o-Toluidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
p-Dimethylaminoazobenzene	0.00	ppm	<	412.	ppb	1:1		< 412.		
p-Nitroaniline	0.00	ppm	<	1770	ppb	1:1		< 1770		
p-Phenylenediamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
p-Terphenyl-d14	1.69	ppb		412.	ppb	1:1	82.1%	< 412.		
sym-Trinitrobenzene	0.00	ppm	<	412.	ppb	1:1		< 412.		
TC: TPH-9073 Method: EPA 9073					Batch: 17388	Run	Analyst:	SJ DOA: 07/23/92	TOA: 1200	OK
Total Petroleum Hydrocarbons	21.7	ppm		10.0	ppm	1:1		< 10.0	2.16%	103%

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NC	233
SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali

Date: 08/03/92

Released by: QA/QC Officer

cc: MLRN00191 **Project Manager:** Winter Seibert **Page No.:** 1

Sample ID	: HB-5A-1
Lab ID	: 9207192-01
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Organic Prep

Evaporative Loss @ 105 C	17.0 wt%

Extractable Organics

Method 8270 Base/Neutral Compounds

1,2,4,5-Tetrachlorobenzene	< 330 ppb
1,2,4-Trichlorobenzene	< 330 ppb
1,2-Dichlorobenzene	< 330 ppb
1,3-Dichlorobenzene	< 330 ppb
1,4-Dichlorobenzene	< 330 ppb
1,4-Dioxane	< 402 ppb
1,4-Naphthoquinone	< 402 ppb
1-Naphthylamine	< 330 ppb
2,4-Dinitrotoluene	< 330 ppb
2,6-Dinitrotoluene	< 330 ppb
2-Acetylaminofluorene	< 711 ppb
2-Chloronaphthalene	< 330 ppb
2-Methylnaphthalene	< 330 ppb
2-Naphthylamine	< 330 ppb
2-Picoline	< 330 ppb
3,3'-Dichlorobenzidine	< 660 ppb
3,3'-Dimethylbenzidine	< 402 ppb
3-Methylcholanthrene	< 402 ppb
4,4'-DDD	< 402 ppb
4,4'-DDE	< 402 ppb
4,4'-DDT	< 402 ppb
4-Bromophenyl phenyl ether	< 402 ppb

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CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali
cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-5A-1
Lab ID	: 9207192-01
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

4-Chloroaniline	< 330 ppb
4-Chlorophenyl phenyl ether	< 330 ppb
4-Nitroquinoline	< 1280 ppb
4-aminobiphenyl	< 402 ppb
5-Nitro-o-toluidine	< 330 ppb
7,12-Dimethylbenz(a)anthracene	< 442 ppb
Acenaphthene	< 330 ppb
Acenaphthylene	< 330 ppb
Acetophenone	< 330 ppb
Aldrin	< 675 ppb
Aniline	< 330 ppb
Anthracene	< 330 ppb
Aramite	< 330 ppb
Benzidine	< 402 ppb
Benzo(a)anthracene	< 330 ppb
Benzo(a)pyrene	< 330 ppb
Benzo(b)fluoranthene	< 376 ppb
Benzo(ghi)perylene	< 330 ppb
Benzo(k)fluoranthene	< 502 ppb
Benzyl Alcohol	< 330 ppb
Butyl benzyl phthalate	< 330 ppb
Chlordane	< 402 ppb
Chlorobenzilate	< 330 ppb
Chrysene	< 402 ppb
Di-n-butyl phthalate	< 402 ppb
Di-n-octyl phthalate	< 402 ppb
Diallate	< 330 ppb
Dibenzo(a,h)anthracene	< 330 ppb
Dibenzofuran	< 330 ppb
Dieldrin	< 735 ppb

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CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-5A-1
Lab ID	: 9207192-01
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Diethyl phthalate	< 330 ppb
Dimethoate	< 330 ppb
Dimethyl phthalate	< 330 ppb
Diphenylamine	< 330 ppb
Disulfoton	< 402 ppb
Endosulfan I	< 402 ppb
Endosulfan II	< 402 ppb
Endosulfan sulfate	< 715 ppb
Endrin	< 755 ppb
Endrin aldehyde	< 1370 ppb
Ethyl Methanesulfonate	< 402 ppb
Famphur	< 330 ppb
Fluoranthene	< 330 ppb
Fluorene	< 330 ppb
Heptachlor	< 896 ppb
Heptachlor epoxide	< 401 ppb
Hexachlorobenzene	< 330 ppb
Hexachlorobutadiene	< 330 ppb
Hexachlorocyclopentadiene	< 438 ppb
Hexachloroethane	< 330 ppb
Hexachloropropene	< 330 ppb
Indeno(1,2,3-c,d)pyrene	< 402 ppb
Isodrin	< 330 ppb
Isophorone	< 330 ppb
Isosafrole	< 330 ppb
Kepone	< 330 ppb
Methapyrilene	< 330 ppb
Methyl Methanesulfonate	< 402 ppb
N-Nitrosodi-n-butylamine	< 402 ppb
N-Nitrosodiethylamine	< 402 ppb

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CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-5A-1
Lab ID	: 9207192-01
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

N-Nitrosodimethylamine	< 330 ppb
N-Nitrosodiphenylamine	< 402 ppb
N-Nitrosodipropylamine	< 330 ppb
N-Nitrosomethylethylamine	< 2810 ppb
N-Nitrosomorpholine	< 330 ppb
N-Nitrosopiperidine	< 330 ppb
N-Nitrosopyrolidine	< 330 ppb
Naphthalene	< 330 ppb
Nitrobenzene	< 330 ppb
O,O,O-Triethylphosphorothioate	< 402 ppb
Parathion, methyl	< 330 ppb
Pentachlorobenzene	< 330 ppb
Pentachloroethane	< 330 ppb
Pentachloronitrobenzene	< 330 ppb
Phenacetin	< 330 ppb
Phenanthrene	< 330 ppb
Phorate	< 330 ppb
Pronamide	< 330 ppb
Pyrene	< 330 ppb
Pyridine	< 751 ppb
Safrole	< 330 ppb
Sulfotepp	< 402 ppb
Thionazin	< 402 ppb
Toxaphene	< 402 ppb
Tributylphosphate	< 402 ppb
a-,a-Dimethylphenethylamine	< 402 ppb
alpha-BHC	< 402 ppb
beta-BHC	< 402 ppb
bis(2-Chloroethoxy)methane	< 330 ppb
bis(2-Chloroethyl) ether	< 330 ppb

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Laboratory	Certifications:
FL	E87156/87294
NC	233
SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: HB-5A-1
Lab ID	: 9207192-01
Matrix	: Soil
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

bis(2-Chloroisopropyl)ether	< 330 ppb
bis(2-Ethylhexyl)phthalate	1490 ppb
delta-BHC	< 402 ppb
gamma-BHC	< 1160 ppb
m-Dinitrobenzene	< 330 ppb
m-Nitroaniline	< 2010 ppb
o-Nitroaniline	< 330 ppb
o-Toluidine	< 330 ppb
p-Dimethylaminoazobenzene	< 402 ppb
p-Nitroaniline	< 1730 ppb
p-Phenylenediamine	< 330 ppb
sym-Trinitrobenzene	< 402 ppb

General Chemistry

Total Petroleum Hydrocarbons	30.6 ppm

The following preparation procedures were performed:

Ext. & Conc. - B/N Compounds

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Sample: 9207192-01 Client ID: Matrix: Soil Description: H8-5A-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
TC: B/N EXT	Method: EPA 8270				Batch:	Run	Analyst:	DOA:	TOA:	NYD
TC: EVAP LOSS	Method: EPA 160.3				Batch: 16959	Run	Analyst: BDK	DOA: 07/11/92	TOA: 1200	OK
Evaporative Loss @ 105 C	17.0	wt%		0.00	wt%	1:1		0.00	0.00%	N/C
TC: 8270 B/N	Method: EPA 8270				Batch: 17692	Run	Analyst: AGW	DOA: 07/16/92	TOA: 1612	OK
						Prep	Analyst: BDK	DOP: 07/15/92	TOP: 1600	NOT OK
1,2,4,5-Tetrachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,2,4-Trichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,2-Dichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,3-Dichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,4-Dichlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
1,4-Dioxane	0.00	ppm	<	402.	ppb	1:1		< 402.		
1,4-Naphthoquinone	0.00	ppm	<	402.	ppb	1:1		< 402.		
1-Naphthylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
2,4-Dinitrotoluene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	48.0%
2,6-Dinitrotoluene	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Acetylaminofluorene	0.00	ppm	<	711.	ppb	1:1		< 711.		
2-Chloronaphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Fluorobiphenyl	1.33	ppb		402.	ppb	1:1	66.2%	< 402.		
2-Methylnaphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Naphthylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
2-Picoline	0.00	ppm	<	330.	ppb	1:1		< 330.		
3,3'-Dichlorobenzidine	0.00	ppm	<	660.	ppb	1:1		< 660.		
3,3'-Dimethylbenzidine	0.00	ppm	<	402.	ppb	1:1		< 402.		
3-Methylcholanthrene	0.00	ppm	<	402.	ppb	1:1		< 402.		
4,4'-DDD	0.00	ppm	<	402.	ppb	1:1		< 402.	0.00%	N/C
4,4'-DDE	0.00	ppm	<	402.	ppb	1:1		< 402.	0.00%	N/C
4,4'-DDT	0.00	ppm	<	402.	ppb	1:1		< 402.	0.00%	N/C
4-Bromophenyl phenyl ether	0.00	ppm	<	402.	ppb	1:1		< 402.		
4-Chloroaniline	0.00	ppm	<	330.	ppb	1:1		< 330.		
4-Chlorophenyl phenyl ether	0.00	ppm	<	330.	ppb	1:1		< 330.		
4-Nitroquinoline	0.00	ppm	<	1280	ppb	1:1		< 1280		
4-aminobiphenyl	0.00	ppm	<	402.	ppb	1:1		< 402.		
5-Nitro-o-toluidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
7,12-Dimethylbenz(a)anthrace	0.00	ppm	<	442.	ppb	1:1		N/C		
Acenaphthene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	64.0%
Acenaphthylene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Acetophenone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Aldrin	0.00	ppm	<	675.	ppb	1:1		< 675.	0.00%	N/C
Aniline	0.00	ppm	<	330.	ppb	1:1		< 330.		
Anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Aramite	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzidine	0.00	ppm	<	402.	ppb	1:1		< 402.	0.00%	N/C
Benzo(a)anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzo(a)pyrene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzo(b)fluoranthene	0.00	ppm	<	376.	ppb	1:1		< 376.		
Benzo(ghi)perylene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Benzo(k)fluoranthene	0.00	ppm	<	502.	ppb	1:1		< 502.		
Benzyl Alcohol	0.00	ppm	<	330.	ppb	1:1		< 330.		
Butyl benzyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Chlordane	0.00	ppm	<	402.	ppb	1:1		< 402.	0.00%	N/C
Chlorobenzilate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Chrysene	0.00	ppm	<	402.	ppb	1:1		< 402.		
Di-n-butyl phthalate	0.00	ppm	<	402.	ppb	1:1		< 402.		
Di-n-octyl phthalate	0.00	ppm	<	402.	ppb	1:1		< 402.		

Sample: 9207192-01 Client ID: Matrix: Soil Description: MB-5A-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Diallate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzo(a,h)anthracene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dibenzofuran	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dieldrin	0.00	ppm	<	735.	ppb	1:1		< 735.	0.00%	N/C
Diethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethoate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Dimethyl phthalate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Diphenylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Disulfoton	0.00	ppm	<	402.	ppb	1:1		< 402.		
Endosulfan I	0.00	ppm	<	402.	ppb	1:1		< 402.	0.00%	N/C
Endosulfan II	0.00	ppm	<	402.	ppb	1:1		< 402.	0.00%	N/C
Endosulfan sulfate	0.00	ppm	<	715.	ppb	1:1		< 715.	0.00%	N/C
Endrin	0.00	ppm	<	755.	ppb	1:1		< 755.	0.00%	N/C
Endrin aldehyde	0.00	ppm	<	1370	ppb	1:1		< 1370	0.00%	N/C
Ethyl Methanesulfonate	0.00	ppm	<	402.	ppb	1:1		< 402.		
Famphur	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluoranthene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Fluorene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Heptachlor	0.00	ppm	<	896.	ppb	1:1		< 896.	0.00%	N/C
Heptachlor epoxide	0.00	ppm	<	401.	ppb	1:1		< 401.	0.00%	N/C
Hexachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorobutadiene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachlorocyclopentadiene	0.00	ppm	<	438.	ppb	1:1		< 438.		
Hexachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Hexachloropropene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Indeno(1,2,3-c,d)pyrene	0.00	ppm	<	402.	ppb	1:1		< 402.		
Isodrin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isophorone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Isosafrole	0.00	ppm	<	330.	ppb	1:1		< 330.		
Kepone	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methapyrilene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Methyl Methanesulfonate	0.00	ppm	<	402.	ppb	1:1		< 402.		
N-Nitrosodi-n-butylamine	0.00	ppm	<	402.	ppb	1:1		< 402.		
N-Nitrosodiethylamine	0.00	ppm	<	402.	ppb	1:1		< 402.		
N-Nitrosodimethylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosodiphenylamine	0.00	ppm	<	402.	ppb	1:1		< 402.		
N-Nitrosodipropylamine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosomethylethylamine	0.00	ppm	<	2810	ppb	1:1		< 2810		
N-Nitrosomorpholine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopiperidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
N-Nitrosopyrolidine	0.00	ppm	<	330.	ppb	1:1		< 330.		
Naphthalene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Nitrobenzene-d5	1.12	ppb		402.	ppb	1:1	55.8%	< 402.		
O,O,O-Triethylphosphorothioa	0.00	ppm	<	402.	ppb	1:1		N/C		
Parathion, methyl	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachlorobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloroethane	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pentachloronitrobenzene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenacetin	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phenanthrene	0.00	ppm	<	330.	ppb	1:1		< 330.		
Phorate	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pronamide	0.00	ppm	<	330.	ppb	1:1		< 330.		
Pyrene	0.00	ppm	<	330.	ppb	1:1		< 330.	0.00%	68.0%

Sample: 9207192-01 Client ID: Matrix: Soil Description: MB-5A-1
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRM00191 Manager: WAS

Parameter	Result	Units	CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Pyridine	0.00	ppm	< 751.	ppb	1:1		< 751.		
Safrole	0.00	ppm	< 330.	ppb	1:1		< 330.		
Sulfotepp	0.00	ppm	< 402.	ppb	1:1		< 402.		
Thionazin	0.00	ppm	< 402.	ppb	1:1		< 402.		
Toxaphene	0.00	ppm	< 402.	ppb	1:1		< 402.	0.00%	N/C
Tributylphosphate	0.00	ppm	< 402.	ppb	1:1		< 402.		
a-,a-Dimethylphenethylamine	0.00	ppm	< 402.	ppb	1:1		< 402.		
alpha-BHC	0.00	ppm	< 402.	ppb	1:1		< 402.	0.00%	N/C
beta-BHC	0.00	ppm	< 402.	ppb	1:1		< 402.	0.00%	N/C
bis(2-Chloroethoxy)methane	0.00	ppm	< 330.	ppb	1:1		< 330.		
bis(2-Chloroethyl) ether	0.00	ppm	< 330.	ppb	1:1		< 330.		
bis(2-Chloroisopropyl)ether	0.00	ppm	< 330.	ppb	1:1		< 330.	0.00%	N/C
bis(2-Ethylhexyl)phthalate	1.49	ppm	< 402.	ppb	1:1		< 402.		
delta-BHC	0.00	ppm	< 402.	ppb	1:1		< 402.	0.00%	N/C
gamma-BHC	0.00	ppm	< 1160	ppb	1:1		< 1160	0.00%	N/C
m-Dinitrobenzene	0.00	ppm	< 330.	ppb	1:1		< 330.		
m-Nitroaniline	0.00	ppm	< 2010	ppb	1:1		< 2010		
o-Nitroaniline	0.00	ppm	< 330.	ppb	1:1		< 330.		
o-Toluidine	0.00	ppm	< 330.	ppb	1:1		< 330.		
p-Dimethylaminoazobenzene	0.00	ppm	< 402.	ppb	1:1		< 402.		
p-Nitroaniline	0.00	ppm	< 1730	ppb	1:1		< 1730		
p-Phenylenediamine	0.00	ppm	< 330.	ppb	1:1		< 330.		
p-Terphenyl-d14	1.49	ppb	402.	ppb	1:1	74.2%	< 402.		
sym-Trinitrobenzene	0.00	ppm	< 402.	ppb	1:1		< 402.		
TC: TPH-9073 Method: EPA 9073				Batch: 17388 Run Analyst: SJ DOA: 07/23/92 TOA: 1200 OK					
Total Petroleum Hydrocarbons	30.6	ppm	10.0	ppm	1:1		< 10.0	2.16%	103%

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SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali

Date: 08/03/92

Released by: QA/QC Officer

cc: MLRN00191 **Project Manager:** Winter Seibert **Page No.:** 1

Sample ID	: FB070792
Lab ID	: 9207192-05
Matrix	: GroundH2O
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Extractable Organics

Method 8270 Base/Neutral Compounds

1,2,4,5-Tetrachlorobenzene	< 10.0 ppb
1,2,4-Trichlorobenzene	< 10.0 ppb
1,2-Dichlorobenzene	< 10.0 ppb
1,3-Dichlorobenzene	< 10.0 ppb
1,4-Dichlorobenzene	< 10.0 ppb
1,4-Dioxane	< 10.0 ppb
1,4-Naphthoquinone	< 10.0 ppb
1-Naphthylamine	< 10.0 ppb
2,4-Dinitrotoluene	< 10.0 ppb
2,6-Dinitrotoluene	< 10.0 ppb
2-Acetylaminofluorene	< 10.8 ppb
2-Chloronaphthalene	< 10.0 ppb
2-Methylnaphthalene	< 10.0 ppb
2-Naphthylamine	< 10.0 ppb
2-Picoline	< 10.0 ppb
3,3'-Dichlorobenzidine	< 20.0 ppb
3,3'-Dimethylbenzidine	< 10.0 ppb
3-Methylcholanthrene	< 10.0 ppb
4,4'-DDD	< 10.0 ppb
4,4'-DDE	< 10.0 ppb
4,4'-DDT	< 10.0 ppb
4-Bromophenyl phenyl ether	< 10.0 ppb
4-Chloroaniline	< 10.0 ppb
4-Chlorophenyl phenyl ether	< 10.0 ppb
4-Nitroquinoline	< 19.5 ppb

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Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert **Page No.:** 2

Sample ID	: FB070792
Lab ID	: 9207192-05
Matrix	: GroundH2O
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

4-aminobiphenyl	< 10.0 ppb
5-Nitro-o-toluidine	< 10.0 ppb
7,12-Dimethylbenz(a)anthracene	< 10.0 ppb
Acenaphthene	< 10.0 ppb
Acenaphthylene	< 10.0 ppb
Acetophenone	< 10.0 ppb
Aldrin	< 10.3 ppb
Aniline	< 10.0 ppb
Anthracene	< 10.0 ppb
Aramite	< 10.2 ppb
Benzidine	< 10.0 ppb
Benzo(a)anthracene	< 10.0 ppb
Benzo(a)pyrene	< 10.0 ppb
Benzo(b)fluoranthene	< 10.0 ppb
Benzo(ghi)perylene	< 10.0 ppb
Benzo(k)fluoranthene	< 10.0 ppb
Benzyl Alcohol	< 10.0 ppb
Butyl benzyl phthalate	< 10.0 ppb
Chlordane	< 10.0 ppb
Chlorobenzilate	< 10.0 ppb
Chrysene	< 10.0 ppb
Di-n-butyl phthalate	< 10.0 ppb
Di-n-octyl phthalate	< 10.0 ppb
Diallate	< 10.0 ppb
Dibenzo(a,h)anthracene	< 10.0 ppb
Dibenzofuran	< 10.0 ppb
Dieldrin	< 11.2 ppb
Diethyl phthalate	< 10.0 ppb
Dimethoate	< 10.0 ppb
Dimethyl phthalate	< 10.0 ppb

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CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059
Contact: Mr. Sid Syedali
cc: MLRN00191 **Project Manager:** Winter Seibert

Date: 08/03/92
Page No.: 3

Sample ID	: FB070792
Lab ID	: 9207192-05
Matrix	: GroundH2O
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

Diphenylamine	< 10.0 ppb
Disulfoton	< 10.0 ppb
Endosulfan I	< 10.0 ppb
Endosulfan II	< 10.0 ppb
Endosulfan sulfate	< 10.9 ppb
Endrin	< 11.5 ppb
Endrin aldehyde	< 20.9 ppb
Ethyl Methanesulfonate	< 10.0 ppb
Famphur	< 10.0 ppb
Fluoranthene	< 10.0 ppb
Fluorene	< 10.0 ppb
Heptachlor	< 13.7 ppb
Heptachlor epoxide	< 10.0 ppb
Hexachlorobenzene	< 10.0 ppb
Hexachlorobutadiene	< 10.0 ppb
Hexachlorocyclopentadiene	< 10.0 ppb
Hexachloroethane	< 10.0 ppb
Hexachloropropene	< 10.0 ppb
Indeno(1,2,3-c,d)pyrene	< 10.0 ppb
Isodrin	< 10.0 ppb
Isophorone	< 10.0 ppb
Isosafrole	< 10.0 ppb
Kepone	< 10.0 ppb
Methapyrilene	< 10.0 ppb
Methyl Methanesulfonate	< 10.0 ppb
N-Nitrosodi-n-butylamine	< 10.0 ppb
N-Nitrosodiethylamine	< 10.0 ppb
N-Nitrosodimethylamine	< 10.0 ppb
N-Nitrosodiphenylamine	< 10.0 ppb
N-Nitrosodipropylamine	< 10.0 ppb

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Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali

cc: MLRN00191 **Project Manager:** Winter Seibert

Sample ID	: FB070792
Lab ID	: 9207192-05
Matrix	: GroundH2O
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

N-Nitrosomethylethylamine	< 42.9 ppb
N-Nitrosomorpholine	< 10.0 ppb
N-Nitrosopiperidine	< 10.0 ppb
N-Nitrosopyrolidine	< 10.0 ppb
Naphthalene	< 10.0 ppb
Nitrobenzene	< 10.0 ppb
O,O,O-Triethylphosphorothioate	< 10.0 ppb
Parathion, methyl	< 10.0 ppb
Pentachlorobenzene	< 10.0 ppb
Pentachloroethane	< 10.0 ppb
Pentachloronitrobenzene	< 10.0 ppb
Phenacetin	< 10.0 ppb
Phenanthrene	< 10.0 ppb
Phorate	< 10.2 ppb
Pronamide	< 10.0 ppb
Pyrene	< 10.0 ppb
Pyridine	< 11.4 ppb
Safrole	< 10.0 ppb
Sulfotepp	< 10.0 ppb
Thionazin	< 10.0 ppb
Toxaphene	< 10.0 ppb
Tributylphosphate	< 10.0 ppb
a-,a-Dimethylphenethylamine	< 10.0 ppb
alpha-BHC	< 10.0 ppb
beta-BHC	< 10.0 ppb
bis(2-Chloroethoxy)methane	< 10.0 ppb
bis(2-Chloroethyl) ether	< 10.0 ppb
bis(2-Chloroisopropyl)ether	< 10.0 ppb
bis(2-Ethylhexyl)phthalate	11.2 ppb
delta-BHC	< 10.0 ppb

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SC	10120
VA	00151
TN	02934
WI	99988779

CERTIFICATE OF ANALYSIS

Client: Mclaren/Hart
25 Independence Boulevard
Warren, New Jersey 07059

Date: 08/03/92

Contact: Mr. Sid Syedali
cc: MLRN00191 **Project Manager:** Winter Seibert

Page No.: 5

Sample ID	: FB070792
Lab ID	: 9207192-05
Matrix	: GroundH2O
Date Collected	: 07/07/92
Date Received	: 07/09/92
Priority	: Routine
Collector	: Client

gamma-BHC	< 17.7 ppb
m-Dinitrobenzene	< 10.0 ppb
m-Nitroaniline	< 30.6 ppb
o-Nitroaniline	< 10.0 ppb
o-Toluidine	< 10.0 ppb
p-Dimethylaminoazobenzene	< 10.0 ppb
p-Nitroaniline	< 26.3 ppb
p-Phenylenediamine	< 10.0 ppb
sym-Trinitrobenzene	< 10.0 ppb

General Chemistry

Total Petroleum Hydrocarbons	< 1.00 ppm

The following preparation procedures were performed:

Ext. & Conc. - B/N Compounds

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Sample: 9207192-05 Client ID: Matrix: GroundH2O Description: FB070792
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
TC: B/N EXT Method:					Batch:	Run	Analyst:	DOA:	TOA:	MYD
TC: 8270 B/N Method:					Batch: 17694	Run	Analyst: AGW	DOA: 07/16/92	TOA: 2129	OK
						Prep	Analyst: CBL	DOP: 07/13/92	TOP: 1600	OK
1,2,4,5-Tetrachlorobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
1,2,4-Trichlorobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
1,2-Dichlorobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
1,3-Dichlorobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
1,4-Dichlorobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
1,4-Dioxane	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
1,4-Naphthoquinone	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
1-Naphthylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
2,4-Dinitrotoluene	0.00	ppm	<	10.0	ppb	1:1		< 10.0	1.75%	56.0%
2,6-Dinitrotoluene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
2-Acetylaminofluorene	0.00	ppm	<	10.8	ppb	1:1		< 10.8		
2-Chloronaphthalene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
2-Fluorobiphenyl	0.0520	ppm		0.00	ppm	1:1	102%	0.0460		
2-Methylnaphthalene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
2-Naphthylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
2-Picoline	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
3,3'-Dichlorobenzidine	0.00	ppm	<	20.0	ppb	1:1		< 20.0		
3,3'-Dimethylbenzidine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
3-Methylcholanthrene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
4,4'-DDD	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
4,4'-DDE	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
4,4'-DDT	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
4-Bromophenyl phenyl ether	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
4-Chloroaniline	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
4-Chlorophenyl phenyl ether	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
4-Nitroquinoline	0.00	ppm	<	19.5	ppb	1:1		< 19.5		
4-aminobiphenyl	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
5-Nitro-o-toluidine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
7,12-Dimethylbenz(a)anthrace	0.00	ppm	<	10.0	ppb	1:1		N/C		
Acenaphthene	0.00	ppm	<	10.0	ppb	1:1		< 10.0	1.12%	88.0%
Acenaphthylene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Acetophenone	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Aldrin	0.00	ppm	<	10.3	ppb	1:1		< 10.3	0.00%	N/C
Aniline	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Anthracene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Aramite	0.00	ppm	<	10.2	ppb	1:1		< 10.2		
Benzidine	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
Benzo(a)anthracene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Benzo(a)pyrene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Benzo(b)fluoranthene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Benzo(ghi)perylene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Benzo(k)fluoranthene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Benzyl Alcohol	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Butyl benzyl phthalate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Chlordane	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
Chlorobenzilate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Chrysene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Di-n-butyl phthalate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Di-n-octyl phthalate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Diallate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Dibenzo(a,h)anthracene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		

Sample: 9207192-05 Client ID: Matrix: GroundH20 Description: FB070792
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRN00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Dibenzofuran	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Dieldrin	0.00	ppm	<	11.2	ppb	1:1		< 11.2	0.00%	N/C
Diethyl phthalate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Dimethoate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Dimethyl phthalate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Diphenylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Disulfoton	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Endosulfan I	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
Endosulfan II	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
Endosulfan sulfate	0.00	ppm	<	10.9	ppb	1:1		< 10.9	0.00%	N/C
Endrin	0.00	ppm	<	11.5	ppb	1:1		< 11.5	0.00%	N/C
Endrin aldehyde	0.00	ppm	<	20.9	ppb	1:1		< 20.9	0.00%	N/C
Ethyl Methanesulfonate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Famphur	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Fluoranthene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Fluorene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Heptachlor	0.00	ppm	<	13.7	ppb	1:1		< 13.7	0.00%	N/C
Heptachlor epoxide	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
Hexachlorobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Hexachlorobutadiene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Hexachlorocyclopentadiene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Hexachloroethane	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Hexachloropropene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Indeno(1,2,3-c,d)pyrene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Isodrin	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Isophorone	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Isosafrole	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Kepone	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Methapyrilene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Methyl Methanesulfonate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosodi-n-butylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosodiethylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosodimethylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosodiphenylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosodipropylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosomethylethylamine	0.00	ppm	<	42.9	ppb	1:1		< 42.9		
N-Nitrosomorpholine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosopiperidine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
N-Nitrosopyrolidine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Naphthalene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Nitrobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Nitrobenzene-d5	0.0388	ppm		0.00	ppm	1:1	76.0%	0.0350		
O,O,O-Triethylphosphorothioa	0.00	ppm	<	10.0	ppb	1:1		N/C		
Parathion, methyl	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Pentachlorobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Pentachloroethane	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Pentachloronitrobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Phenacetin	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Phenanthrene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Phorate	0.00	ppm	<	10.2	ppb	1:1		< 10.2		
Pronamide	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Pyrene	0.00	ppm	<	10.0	ppb	1:1		< 10.0	12.1%	58.0%
Pyridine	0.00	ppm	<	11.4	ppb	1:1		< 11.4		
Safrole	0.00	ppm	<	10.0	ppb	1:1		< 10.0		

Sample: 9207192-05 Client ID: Matrix: GroundH2O Description: FB070792
Sampled: 07/07/92 Received: 07/09/92 Due: 07/23/92 Priority: 3
Project: MLRW00191 Manager: WAS

Parameter	Result	Units		CRDL	Units	DF	Surr. %	Blank	RPD	%Rec
Sulfotepp	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Thionazin	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
Toxaphene	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
Tributylphosphate	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
a-,a-Dimethylphenethylamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
alpha-BHC	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
beta-BHC	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
bis(2-Chloroethoxy)methane	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
bis(2-Chloroethyl) ether	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
bis(2-Chloroisopropyl)ether	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
bis(2-Ethylhexyl)phthalate	0.0112	ppm	<	10.0	ppb	1:1		< 10.0		
delta-BHC	0.00	ppm	<	10.0	ppb	1:1		< 10.0	0.00%	N/C
gamma-BHC	0.00	ppm	<	17.7	ppb	1:1		< 17.7	0.00%	N/C
m-Dinitrobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
m-Nitroaniline	0.00	ppm	<	30.6	ppb	1:1		< 30.6		
o-Nitroaniline	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
o-Toluidine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
p-Dimethylaminoazobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
p-Nitroaniline	0.00	ppm	<	26.3	ppb	1:1		< 26.3		
p-Phenylenediamine	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
p-Terphenyl-d14	0.0347	ppm		0.00	ppm	1:1	68.0%	0.0540		
sym-Trinitrobenzene	0.00	ppm	<	10.0	ppb	1:1		< 10.0		
TC: TPH Method: EPA 418.1					Batch: 17161 Run Analyst: SJ DOA: 07/16/92 TOA: 1500 OK					
Total Petroleum Hydrocarbons	0.200	ppm	<	1.00	ppm	1:1		< 1.00	-0.198%	97.6%

McLaren/Hart Environmental Engineering Corp. Inc. D.C. No.: № 458

Name: Tony Maria
Affiliation: McLaren/Hart
Phone: 908-647-8111
Address: 25 Independence Blvd., Warren, NJ 07059
Client/Job No: American Home Products
Job Name: AHP/Luck's Location: Seagrove, NJ

CHAIN OF CUSTODY RECORD

	Sample No.	Lab I.D. No.	Date	Time	Matrix	No.of Cont.	Preservatives	Analysis Requested/Remarks
1207 192 -01	HB-5A-1		7/7/92	9:30	soil	1	cold	B.N. + oil and grease
-02	HB-5C-1		7/7/92	10:25	soil	1	cold	BN + oil and grease
-03	HB-4B-1		7/7/92	5:30	soil	1	cold	BN + oil and grease
-04	HB-3B-1		7/8/92	8:00	soil	1	cold	BN + oil and grease
-05	FB070792		7/7/92	11:40	water	3	cold	2 - BN 1 - oil + grease

Comments:

Relinquished by: T. Maria Date: 7/8/92 Shipment Method: Fed.ex
Time: 11:00 Airbill No.:

Received by: [illegible] Date: 7-9-92 Relinquished by: Date:
Time: 930 Time:

Received by: Date: Relinquished by: Date:
Time: Time:

Final Disposition of Samples:

Received by: Date: Time:

Page ____ of ____ Pink/Field Copy - Yellow/Lab Copy - White/Original - To be returned with results.

HB-3B-1

QUANT REPORT

Operator ID: BDK
Output File: ^JJ411::OU
Data File: >JJ411::D1
Name: 9207192-04 MLRN A/BN
Misc: BATCH 17152

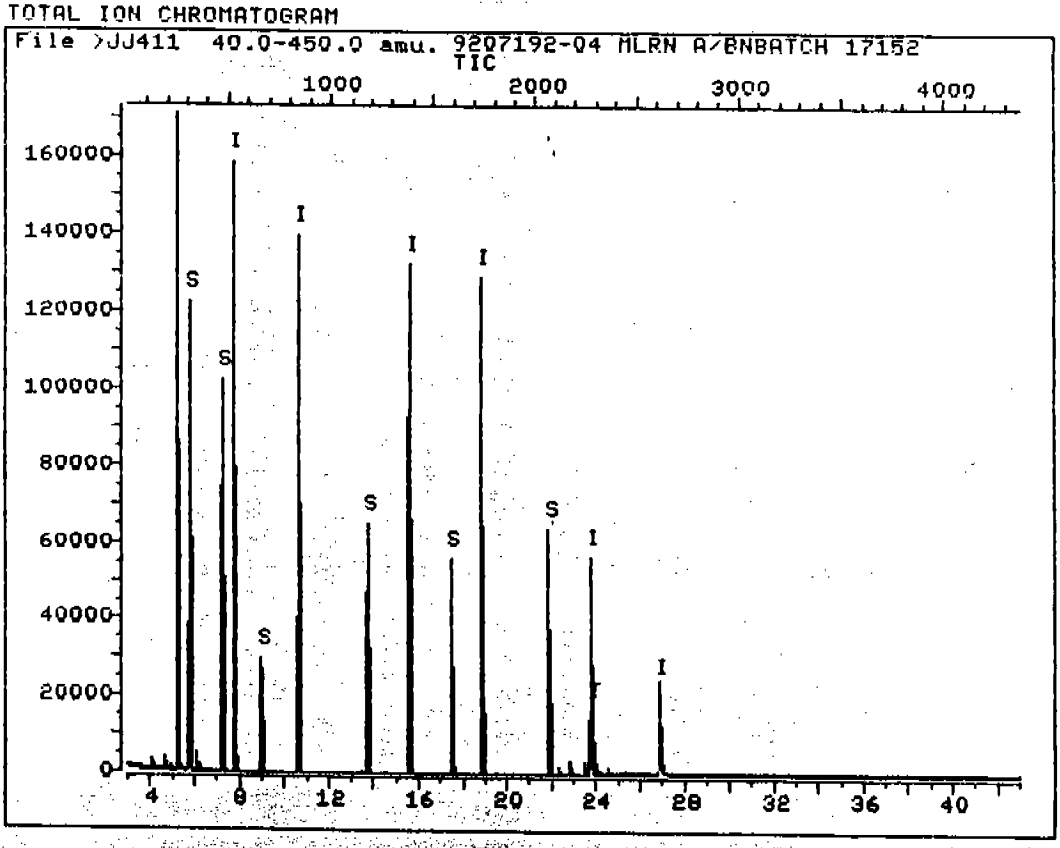
Quant Rev: 7
Quant Time: 920716 19:32
Injected at: 920716 18:48
Dilution Factor: 1.00000
Instrument ID: MSD1
BTL#11

ID File: ID_BN1::DB
Title: General Engineering BNA Standards
Last Calibration: 920714 17:16
Last Qcal Time: 920716 08:29

	Compound	R.T.	Q ion	Area	Conc	Units	q
1)	*1,4-Dichlorobenzene-d4	7.70	152.0	46272	40.00	ng/uL	88
3)	2-Fluorophenol	5.71	112.0	52167	43.72	ng/uL	86
4)	Phenol-d5	7.18	99.0	64086	48.08	ng/uL	88
17)	*Naphthalene-d8	10.58	136.0	133899	40.00	ng/uL	100
18)	Nitrobenzene-d5	8.91	82.0	16817	18.55	ng/uL	88
32)	*Acenaphthene-d10	15.52	164.0	63013	40.00	ng/uL	86
37)	2-Fluorobiphenyl	13.69	172.0	45389	25.23	ng/uL	96
52)	*Phenanthrene-d10	18.81	188.0	93947	40.00	ng/uL	99
55)	2,4,6-Tribromophenol	17.41	330.0	11340	62.89	ng/uL	86
63)	*Chrysene-d12	23.69	240.0	44308	40.00	ng/uL	99
66)	Terphenyl-d14	21.84	244.0	35631	39.66	ng/uL	90
71)	Bis(2-Ethylhexyl)phthalate	23.79	149.0	7424	6.85	ng/uL	97
72)	*Perylene-d12	26.84	264.0	32619	40.00	ng/uL	92

* Compound is ISTD

TOTAL ION CHROMATOGRAM

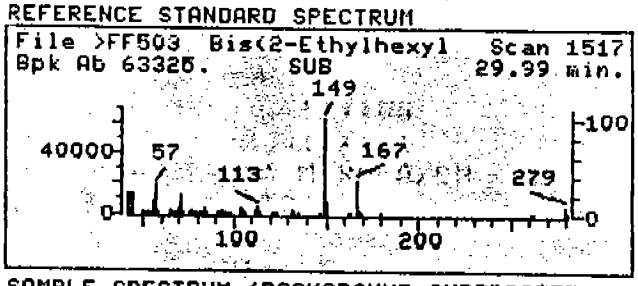


Data File: >JJ411::D1 Quant Output File: ^JJ411::OU
Name: 9207192-04 MLRN A/BN Instrument ID: MSD1
Misc: BATCH 17152 BTL#11

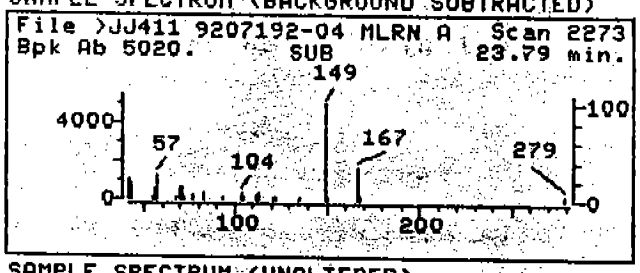
Id File: ID_BN1::DB
Title: General Engineering BNA Standards
Last Calibration: 920714 17:16 Last Qcal Time: 920716 08:29

Operator ID: BDK
Quant Time : 920716 19:32
Injected at: 920716 18:48

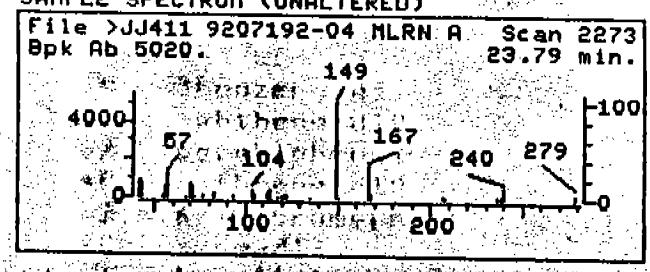
REFERENCE STANDARD SPECTRUM

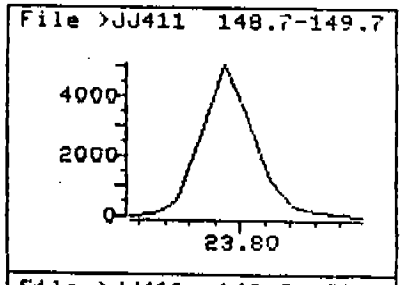


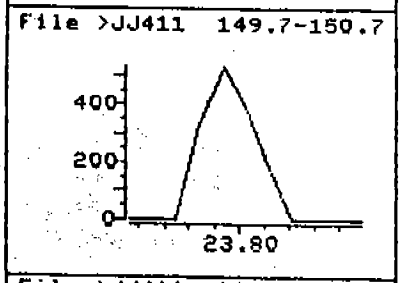
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)

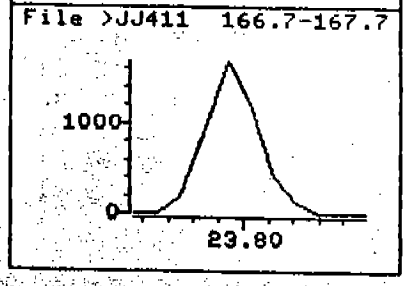


SAMPLE SPECTRUM (UNALTERED)









Data File: >JJ411::D1
Name: 9207192-04 MLRN A/BN
Misc: BATCH 17152
Quant Time: 920716 19:32
Injected at: 920716 18:48
Last Qcal Time: 920716 08:29

Quant Output File: ^JJ411::OU
Instrument ID: MSD1
BTL#11
Quant ID File: ID_BN1::DB
Last Calibration: 920714 17:16

Compound No : 71
Compound Name : Bis(2-Ethylhexyl)phthalate
Scan Number : 2273
Retention Time: 23.79 min.
Quant Ion : 149.0
Area : 7424
Concentration : 6.85 ng/uL
q-value : 97

HB-4B-1

QUANT REPORT

Operator ID: BDK
Output File: ^JJ410::OU
Data File: >JJ410::D1
Name: 9207192-03 MLRN A/BN
Misc: BATCH 17152

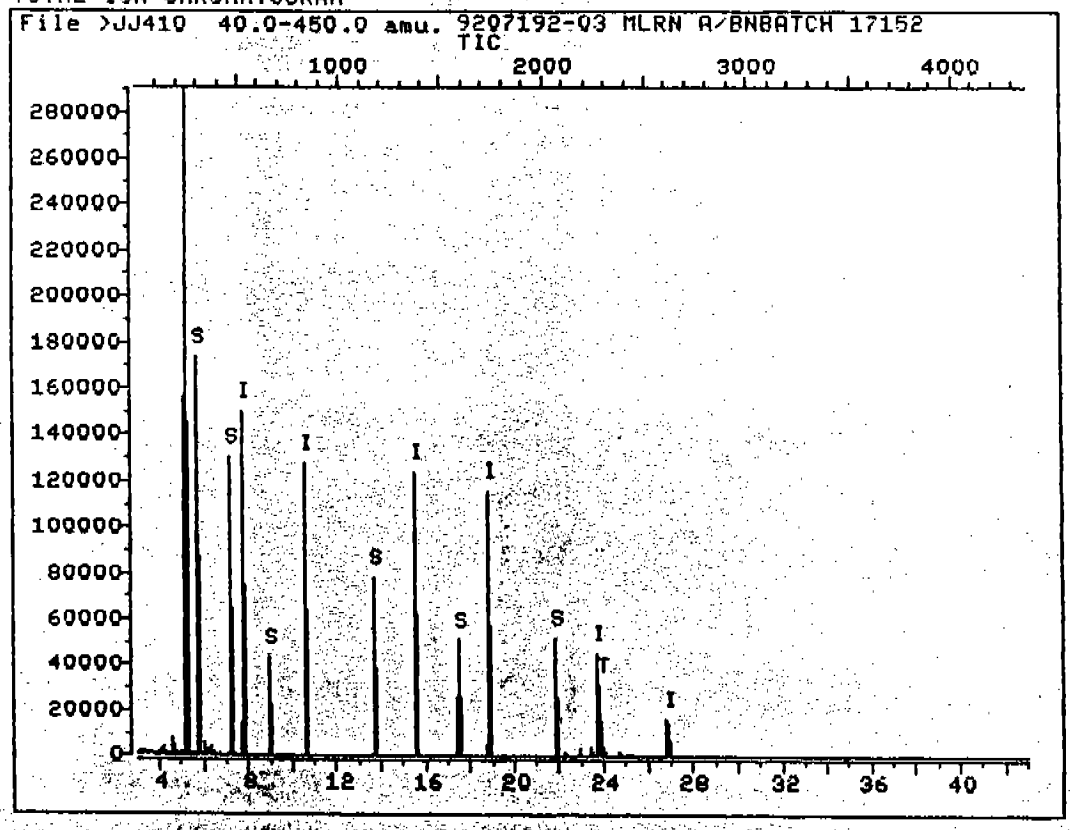
Quant Rev: 7
Quant Time: 920716 18:40
Injected at: 920716 17:56
Dilution Factor: 1.00000
Instrument ID: MSD1
BTL#10

ID File: ID_BN1::DB
Title: General Engineering BNA Standards
Last Calibration: 920714 17:16
Last Qcal Time: 920716 08:29

	Compound	R.T.	Q ion	Area	Conc	Units	q
1)	*1,4-Dichlorobenzene-d4	7.70	152.0	45043	40.00	ng/uL	88
3)	2-Fluorophenol	5.72	112.0	70521	60.71	ng/uL	84
4)	Phenol-d5	7.18	99.0	79808	61.51	ng/uL	90
17)	*Naphthalene-d8	10.58	136.0	127362	40.00	ng/uL	100
18)	Nitrobenzene-d5	8.91	82.0	26166	30.35	ng/uL	88
32)	*Acenaphthene-d10	15.52	164.0	61836	40.00	ng/uL	86
37)	2-Fluorobiphenyl	13.69	172.0	60961	34.53	ng/uL	94
52)	*Phenanthrene-d10	18.80	188.0	87100	40.00	ng/uL	99
55)	2,4,6-Tribromophenol	17.41	330.0	10596	63.38	ng/uL	89
63)	*Chrysene-d12	23.69	240.0	33504	40.00	ng/uL	99
66)	Terphenyl-d14	21.84	244.0	29511	43.44	ng/uL	92
71)	Bis(2-Ethylhexyl)phthalate	23.80	149.0	15585	19.02	ng/uL	95
72)	*Perylene-d12	26.85	264.0	23080	40.00	ng/uL	96

* Compound is ISTD

TOTAL ION CHROMATOGRAM

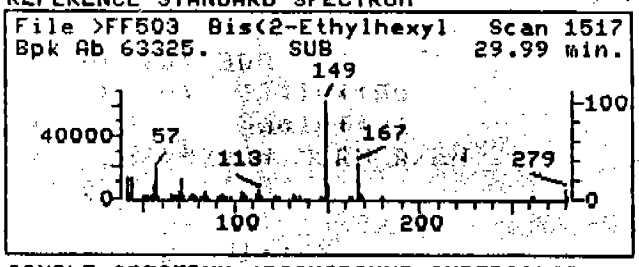


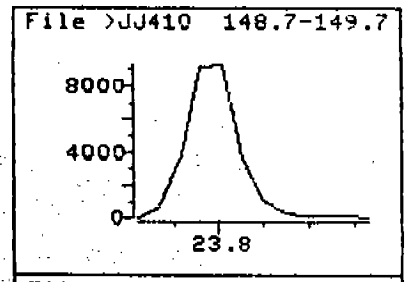
Data File: >JJ410::D1 Quant Output File: ^JJ410::OU
Name: 9207192-03 MLRN A/BN Instrument ID: MSD1
Misc: BATCH 17152 BTL#10

Id File: ID_BN1::DB
Title: General Engineering BNA Standards
Last Calibration: 920714 17:16 Last Qcal Time: 920716 08:29

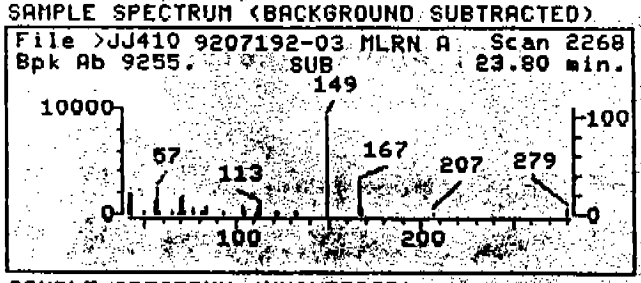
Operator ID: BDK
Quant Time : 920716 18:40
Injected at: 920716 17:56

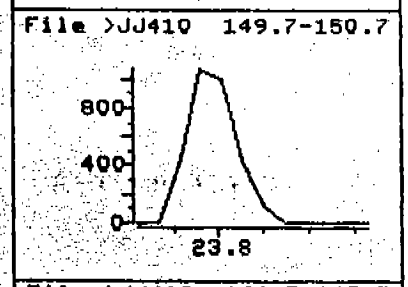
REFERENCE STANDARD SPECTRUM



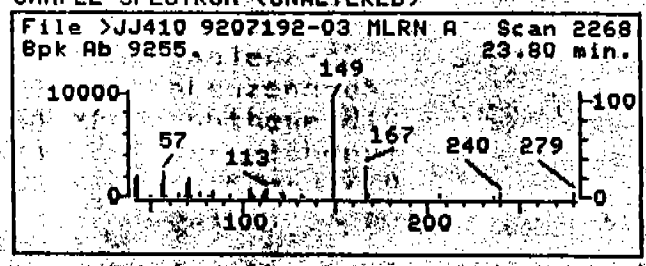


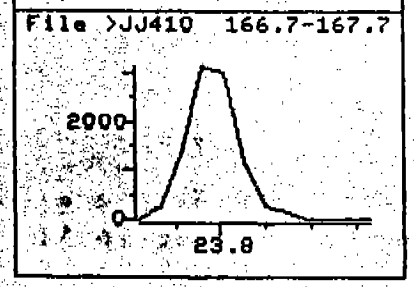
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)





SAMPLE SPECTRUM (UNALTERED)





Data File: >JJ410::D1
Name: 9207192-03 MLRN A/BN
Misc: BATCH 17152
Quant Time: 920716 18:40
Injected at: 920716 17:56
Last Qcal Time: 920716 08:29

Quant Output File: ^JJ410::OU
Instrument ID: MSD1
BTL#10
Quant ID File: ID_BN1::DB
Last Calibration: 920714 17:16

Compound No : 71
Compound Name : Bis(2-Ethylhexyl)phthalate
Scan Number : 2268
Retention Time: 23.80 min.
Quant Ion : 149.0
Area : 15585
Concentration : 19.02 ng/uL
q-value : 95