

JAN 09 1979

→ Rosen - 12  
Reference 7

3717

Southern District Headquarters  
Route 3, Wakanda Drive  
Waunakee, WI 53597

January 2, 1979

4410

Mr. Ray Washkovick, Sr.  
Route 2  
Ripon, WI 54971

Dear Mr. Washkovick:

This letter is to address your complaint regarding toxic industrial waste from the Speed Queen plant located in the City of Ripon being deposited at the City of Ripon solid waste disposal site.

Our Southern District Office has recently received an application for a license to operate an incinerator at the Speed Queen plant. The proposed unit is designed to dispose of oils, paint sludge, solvents and other flammable wastes which are presently being deposited at the City of Ripon solid waste disposal site. If plans proceed on schedule, the unit should be on-line by June 1, 1979.

Your concern for the environment is appreciated. If I can be of any further assistance, please feel free to contact me. My telephone number is 266-0889.

Sincerely,

James Bakken  
Solid Waste Specialist

JB:cb  
Bureau of Solid Waste Mgt. - Pyare  
Horicon Area Office

*... sparks or flame*  
*... municipal refuse at this*  
*... this is the*  
*start of a new production process*

GENERATOR OF WASTE <i>Speed Queen Corp.</i>		TELEPHONE NUMBER <i>414-787-3717</i>
STREET OR ROUTE <i>205 E. Main</i>		TYPE OF PROCESS WHICH PRODUCES WASTE <i>Gear production - balancing</i>
CITY, STATE, ZIP CODE <i>Wautoma, WI.</i>		DNR LICENSE NUMBER <i>1228</i>
DISPOSAL FACILITY <i>City Wautoma Landfill</i>		COUNTY <i>Waushara</i>
SITE LOCATION <i>SW 1/4 SW 1/4 SECTION 22 T. 19N R. 10E</i>		CONTAINER SIZE <input checked="" type="checkbox"/> DRUMS <input type="checkbox"/> BAGS <input type="checkbox"/> CARTONS <input type="checkbox"/> OTHERS
AMOUNT AND FREQUENCY OF DISPOSAL <i>55 gal. once per month</i>		
NUMBER OF CONTAINERS <i>one</i>		

WASTE COMPONENTS	% BY WEIGHT OR VOLUME	TOXICITY DATA		
		LD <sub>50</sub> - ORAL	LD <sub>50</sub> - DERMAL	LD <sub>50</sub> - INHALATION
<i>magnesium filings</i>				

FLASH POINT	NFPA - HAZARD CLASSIFICATIONS:		
SP. GRAVITY	<input type="checkbox"/> HEALTH	<input type="checkbox"/> REACTIVITY	<input checked="" type="checkbox"/> FLAMMABILITY
pH	OTHERS		

SPECIAL PROPERTIES AND HANDLING INSTRUCTIONS (I.E., LABELING OR OSHA DATA SHEETS):

- protect from sparks or flame
- filings will be mixed in with municipal refuse at disposal site and covered daily

1. IS FUTURE USE OF MATERIAL A POSSIBILITY?	YES <input checked="" type="checkbox"/>	NO <input type="checkbox"/>
2. CAN THE MATERIAL STILL BE USED AS ORIGINALLY INTENDED?	<input type="checkbox"/>	<input checked="" type="checkbox"/>
3. CAN IT BE RETURNED TO MANUFACTURER FOR DISPOSAL OR RECYCLING OR CAN IT BE RECYCLED ON A LOCAL BASIS?	<input type="checkbox"/>	<input type="checkbox"/>
4. HAS ANY DISPOSAL ACTION BEEN TAKEN TO DATE?	<input type="checkbox"/>	<input checked="" type="checkbox"/>
5. WHAT ARE PREVIOUSLY EMPLOYED METHODS OF DISPOSAL? EXPLAIN:	<i>none - this is the start of a new production process</i>	

# Donohue

REFERENCE 8  
SITE NAME Ripon Landfill  
SITE ID WTD 980616190

DC

November 17, 1981

State of Wisconsin  
Department of Natural Resources  
P. O. Box 7921  
Madison, WI 53707

Attn: Ms. Barbara Bickford  
Bureau of Solid Waste Management

Re: City of Ripon Landfill  
License No. 467  
Donohue Project No. 11829

NOV 19 1981  
response due 12/19/81

Dear Ms. Bickford:

In accordance with the authorization of the City of Ripon, we have completed installation of new monitoring wells at the referenced site. The purpose of the monitoring wells is to determine the position of the phreatic surface beneath the site and to provide groundwater sampling points for long-term water quality analyses.

The monitoring well array consists of three wells completed at varying depths so that ten feet of screen occurs below the water table and five feet above. Our original intention was to place four wells at the corners of the site to account for all spatial conditions. During field efforts, however, extreme difficulty was encountered in the northwest corner, precluding completion of that well. Subsequent review of the water table and geologic data indicates pronounced uniformity of stratigraphic conditions and low gradients slightly falling to the southwest. We feel that the three well array will provide acceptable monitoring control within the operative budget constraints. Complete soil boring logs and well installation diagrams are attached.

Water analytic parameters will be performed in accordance with NR 180.13, (11), (a), 7 and will include water elevation, field pH, field conductivity, COD, dissolved iron, hardness, chloride, sulfate, and alkalinity. We anticipate acquisition of the first series of samples during the third week of November and will forward the data when complete.

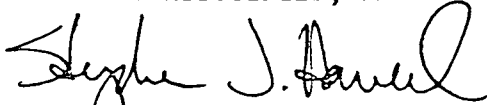
Donohue & Associates, Inc.  
4738 North 40th Street  
Sheboygan, Wisconsin 53081  
Engineers & Architects  
414 458 8711

# Donohue

I trust you will find this discussion in order. If you have any questions, please feel free to call me.

Sincerely,

DONOHUE & ASSOCIATES, INC.



Stephen J. Haverl, P.G.  
Senior Engineering Geologist



Robert Klink, P.E.  
Project Engineer

SJH/slg

enc: Soil Boring Logs  
Well Installation Diagrams

cc: Ken Hein, DNR Solid Waste Specialist  
Claude Lee, City of Ripon

Refer to

Sept 25, 1981 letter from  
Bob Klink to Claude Lee (Ripon)  
for map of well locations  
(1974)



# ecology and environment, inc.

111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-663-9415  
International Specialists in the Environment

REFERENCE 9  
SITE NAME Ripon Landfill  
SITE ID WID 980610190

## M E M O R A N D U M

DATE: July 22, 1985  
TO: File  
FROM: Robert Hingtgen *RH*  
SUBJECT: Wisconsin/R05-8310-01A-37  
Ripon/City of Ripon Landfill  
WID980610190

On June 26, 1985, the author, Ron Bock, and Arlene Pratl met with Dave Edwards, Wisconsin Department of Natural Resources, and Phillip Hoopman, Ripon Wastewater Treatment Plant Superintendent, in order to resample wells at the Ripon Landfill.

Mr. Hoopman had an assistant, John Wendler, take us to the landfill and open the wells for us. The following table contains the wells which were sampled and the water levels in each well.

<u>Well</u>	<u>Depth to Water (Ft.)</u>	<u>Total Depth (Ft.)</u>
5A	50.85	69.00
5	51.33	56.00
6	38.69	47.00
8	23.61	34.00

Well 5 produced a reading of 1000 ppm methane in the well, and 4 ppm in the breathing zone, on the OVA.

Samples were collected June 27, 1985. At the request of Mr. Hoopman, samples were also collected for the wastewater treatment plant. This involved filling one plastic milk jug to approximately one-half volume for each well.

Ripon Landfill has been covered with a clay, but a large leachate seep was observed on the east border of the landfill between wells 6 and 8. The landfill is easily accessible as it borders county road NN. No barrier exists between the landfill and road.

28A:4M

CORRESPONDENCE/MEMORANDUM


DC

Date: August 26, 1981

File Ref. 4440

To: Ken Hein, Horicon Area

REFERENCE 11  
SITE NAME Ripon Landfill  
SITE ID WED 980610130

From:  Ellen Smith, SW/3, GEF II

Subject: City and Town of Ripon Landfill, Lic. #467

This memo is in response to your request that the Residuals Section provide technical assistance on the proper abandonment of existing wells at the Ripon landfill and on the installation of new monitoring points.

Apparently the four observation wells installed through waste in 1974 have since been buried or partially buried by refuse; only well #3 has been located. It may be possible to locate the remaining wells in order to abandon them. Bob Klink at Donohue has a surveyor's report dated 1974 showing boring locations (he's not sure if the survey was done before or after drilling), land surface contours, and boring elevations (don't know if they are pipe tops or ground surface). The Bureau doesn't have a copy of that report. If it proves difficult to locate the wells using this map, I don't feel an extensive search is justified to find them and abandon them. Soils beneath the waste are primarily sands (probably SP and SM) and groundwater was 1 to 8 feet below the waste in 1974, so groundwater contamination is likely with or without these wells.

Any well(s) which are located should be abandoned by grouting the hole(s) with bentonite-Portland cement slurry while removing the galvanized steel well casing. (A slurry pipe should be run through the well casing to just below the base of the casing, and the well casing should be pulled out gradually while slurry is pumped into the hole below the pipe bottom.)

It is extremely important that a groundwater monitoring system be established to define flow patterns, evaluate the extent of any existing contamination, and estimate directions of potential contaminant migration. This information is needed to plan for the orderly completion and proper abandonment of the landfill site. Regionally, the topography suggests that flow is toward the south and west, though local conditions may differ. If the 1974 boring elevations are for the points from which water levels were measured in 1974, these measurements also suggest a gentle southwestward gradient. Given the tentative nature of the water level data, I would suggest that 4 or 5 observation wells be installed just beyond the landfill perimeter on all sides of the landfill site. Well screens should be at least 15 feet in length and

Ken Hein, Horicon Area  
August 26, 1981

2.

positioned to intersect the groundwater table, which is estimated to be at around 820 to 830 feet elevation (USGS). Wells should be constructed of 2-inch diameter PVC pipe. After initial water elevations are known, it should be possible to define prevailing groundwater flow directions and determine additional monitoring needs. At least one down gradient well should be converted to a nest by adding a piezometer with a maximum 5-foot screen at least 15 feet below the bottom of the observation well's screen. It is likely that additional wells will also be needed at greater distances down gradient of the site to define the extent of contamination. Complete soil boring logs and well construction information (in accordance with s. NR 180.13(6)(b)3 c. and g.) and surveyed locations should be compiled for all newly constructed wells and provided to the Bureau as soon as they become available. When wells are in place, initial monitoring should include, at a minimum, the parameters specified in s. NR 180.13(11)(a)7.

Your July 27, 1981 letter to the City of Ripon encourages the City to develop an abandonment plan for the site. Given the site's poor hydrogeologic setting, I would add that abandonment should be planned so as to achieve positive drainage away from the fill area as soon as possible, and not delayed merely in order to prolong the site life.

Topic 3 of your July 27 letter needs clarification: 3:1 slopes are not the minimum requirement for final grades, but the maximum allowable landfill side slope under NR 180. Also, the City should be aware of NR 180 submittal requirements and fees for an abandonment plan.

Please keep me informed on the City's progress in this matter. I understand that you will be meeting with the City again on August 31.

ES:cm/0687Q

cc: Southern District Headquarters

*Handwritten note:*  
All of the official copies, on account  
of the...  
10/1/81




## CORRESPONDENCE/MEMORANDUM

file - 10

Date: September 22, 1981

File Ref: 4400

To: Ken Hein - Horicon Area

From: Ellen Smith - SW/3 

REFERENCE 13


SITE NAME Ripon Landfill

SITE ID WFD 980616190

Subject: Ripon Landfill, License #467

This memo is in response to your request for Bureau guidance on municipal sludge codisposal at the Ripon landfill.

In a few cases involving the codisposal of papermill sludge, the Department has required that sludge acceptance be limited initially to 1 cubic yard sludge for each 10 gate yards of refuse. Where this level of sludge acceptance is found to cause operational problems, sludge volumes are reduced. If no problems occur, increased sludge acceptance has been allowed on an experimental basis. Experience statewide has found that sludge codisposal is successful only with relatively dry sludges and only at landfills with provision for leachate containment and collection.

The Ripon sludge is at 9-10% solids, which usually corresponds to a liquid consistency. 

The Ripon landfill does not provide for leachate control. Because of these factors, continued disposal of any municipal sludge at the landfill is inadvisable.

If the city doesn't have capacity to dispose of all its sludge on farmland at present, some sludge may have to continue to go to the landfill while alternatives are developed. If this occurs, the 10:1 ratio may be a good guideline on a temporary basis. The refuse loading rate you estimated on the basis of population is 8,402,000 lbs/year. At 500 lbs/gate yard (a conservatively high assumption), this amounts to about 16,000 gate yards per year. With 10 yards refuse to 1 yard sludge, 1600 yards (or 323,000 gallons) of sludge could be taken to the landfill. At 1100 gallons per load, one load per day should be the limit on sludge acceptance. To minimize operating problems, sludge should only be disposed of at or near the close of the operating day, but before daily cover is applied. Shallow trenches may be dug into the day's refuse to help maximum absorption and prevent runoff of sludge.

Sludge acceptance at Ripon should be treated strictly as a temporary measure, since this material causes operating problems at the landfill, contributes to leachate production, and can be disposed of in an environmentally sound manner through land application. If the sludge is suitable, it may be desirable to develop sludge storage facilities to handle sludge during periods when land is unavailable for spreading. An NR 180 storage approval could be needed for storage facilities not on treatment plant property.

ES:ok

cc: Solid Waste Section - Southern District Headquarters  
Sludge Coordinator - Southern District Headquarters

DATE 10/8/86

TDD # B58310-015

TIME 12:45 pm

Page 1 of     

SITE Ripon Landfill

CONTACT Bob Lukoski - Fire Chief PHONE (414) 748-2888

City of Ripon

REFERENCE 16

SITE NAME Ripon Landfill

SITE ID WID980610190

SUBJECT: Fire or Explosion Threat

Mr. Lukoski stated that he did not consider The Ripon Landfill to be a Fire or Explosion hazard.

Robert Harty

Date: 10/8/86

DATE 10/8/86

TDD # RS8310-015

TIME 4:00 pm

Page of

SITE Ripon Landfill

CONTACT Dave Edwards

PHONE (414) 485-4435

WDNR

REFERENCE 17

SITE NAME Ripon Landfill

SITE ID WFD 980610190

SUBJECT: Ripon Landfill

Mr. Edwards stated that the Falkenberg well is located about 500' south of the landfill and the well log is as follows: 0-15' clay, 15'-20' sand and gravel, 20'-100' limestone, and 100'-150' sandstone. The well is cased to 44 feet, and the static water level is about 20 feet. Mr. Edwards also thought that the limestone and overlying drift are hydraulically connected as the limestone is quite fractured in that area. Mr. Edwards was not aware of any federally endangered species in the wetlands located northeast of the site.

Robert Hinton

Date: 10/8/86

NGV 23 1992

JUN 17 1977  
 WELL CONSTRUCTOR'S REPORT  
 Form 3300-15 Rev. 12-76

State of Wisconsin  
 Department of Natural Resources  
 Box 9021  
 Madison, Wisconsin 53707

COUNTY OF  
FOND DU LAC

White Copy  
 Green Copy  
 Yellow Copy

Division's Copy  
 Driller's Copy  
 Owner's Copy

1. COUNTY Green Lake CHECK (X) ONE:  Town  Village  City Name Ripon

2. LOCATION Section SE Section 7 Township 16 Range 14 NAME  OWNER  AGENT AT TIME OF DRILLING CHECK (X) ALLIANT ALLIANT BLDG

OR Grid or Street No. Street Name ADDRESS CT7 NW

AND Available subdivision name, lot & block No. POST OFFICE Green Lake, Wis

4. Distance in feet from well to nearest: (Record answer in appropriate place)

Sanitary Bldg. Drain	Sanitary Bldg. Sewer	Floor Drain Connected To:	Storm Bldg. Drain	Storm Bldg. Sewer
C.I.	Other	C.I. Other	C.I.	Other

Street Sewer Other Sewer Foundation Drain Connected to Sewage Sump Clearwater Septic Holding Sewage Absorption Unit  
 San. Storm C.I. Other Sewer Sump Clearwater Clearwater Sump Tank Tank Seepage Pit Seepage Bed Seepage Trench

Pit Noncontaminating Existing Subsurface Pump Room Barn Animal Animal Silo Glass Lined S/W Earthen Silage  
 Waste Well Noncontaminating Existing Surface Pump Room Gutters Barn Pen Yard With Pit Storage Facility w/o Earthen Storage  
 Tank Pump Tank Existing Existing Disposal Unit (Specify Type) Other (Give Description)

5. Well is intended to supply water for Future Home

6. DRILLHOLE

Dia. (in.)	From (ft.)	To (ft.)	Dia. (in.)	From (ft.)	To (ft.)
8	Surface	40			
6	120	102			

9. FORMATIONS

Kind	From (ft.)	To (ft.)
Clay	Surface	15
Sand rock	15	18

7. CASING LINER, CURBING AND SCREEN Material, Weight, Specification & Method of Assembly

Dia. (in.)	From (ft.)	To (ft.)
6	Surface	40

6" HDPE pipe  
1280 wall  
weld jts.  
P-53

10. TYPE OF DRILLING MACHINE USED

<input type="checkbox"/> Cable Tool	<input checked="" type="checkbox"/> Rotary hammer with mud & air	<input type="checkbox"/> Jetting with
<input type="checkbox"/> Rotary air w/drilling mud	<input type="checkbox"/> Rotary hammer & air	<input type="checkbox"/> Air
<input type="checkbox"/> Rotary w/drilling mud	<input type="checkbox"/> Reverse Rotary	<input type="checkbox"/> Water

8. GROUT OR OTHER SEALING MATERIAL

Kind	From (ft.)	To (ft.)
Mud	Surface	8
Cement	8	40

Well construction completed on 5-11 1977

Well is terminated 12 inches  above final grade  below

Well disinfected upon completion  Yes  No

Well sealed water tight upon completion  Yes  No

11. MISCELLANEOUS DATA

Yield Test: 2 Hrs. at 20 GPH

Depth from surface to normal water level 47 Ft.

Depth of water level when pumping 65 Ft. Stabilized  Yes  No

Water sample sent to Madison laboratory on 5-11 1977

Your opinion concerning other pollution hazards, information concerning difficulties encountered, and data relating to nearby wells, screens, seals, method finishing the well, amount of cement used in grouting, blasting, etc., should be given on reverse side.

Signature: Lawrence Halvick Complete Mail Address: \_\_\_\_\_  
 Registered Well Driller

NOV 23 1982

WELL CONSTRUCTOR'S REPORT TO WISCONSIN STATE BOARD OF HEALTH  
See Instructions on Reverse Side

1. County Fond du Lac (Town  Ripon, Village , City  Check on
2. Location Township 15 North Range 14 East Section 7  
Name of street and number of premises or Section, Town and Range numbers
3. Owner  or Agent  L. F. Lehman  
Name of individual, partnership or firm
4. Mail Address Route # 2 Ripon, Wisconsin Box 273  
Complete address required
5. From well to nearest: Building 5 ft; sewer None ft; drain None ft; septic tank None ft;  
dry well or filter bed None ft; abandoned well None ft
6. Well is intended to supply water for: Private Home

RECEIVED  
SEP 29 1958  
ENVIRONMENTAL  
SANITATION

7. DRILLHOLE:

Dia. (in.)	From (ft.)	To (ft.)	Dia. (in.)	From (ft.)	To (ft.)
10	0	23	6	23	153

10. FORMATIONS:

Kind	From (ft.)	To (ft.)
Sand	0	23
Gravel & hardpan	23	62
Limerock	62	82
Shale & sandrock	82	90
Limerock & shale	90	120
Sandrock	120	153

8. CASING AND LINER PIPE OR CURSING:

Dia. (in.)	Kind and Weight	From (ft.)	To (ft.)
6	Steel casing	0	63

9. GROUT:

Kind	From (ft.)	To (ft.)
Clay puddle & sandrock outtings	0	23

11. MISCELLANEOUS DATA:

Yield test: 6 Hrs. at 9 GPM.  
Depth from surface to water-level: 25 ft.  
Water-level when pumping: 30 ft.  
Water sample was sent to the state laboratory at:  
Laborsk on 9/22 1958  
City

Construction of the well was completed on:  
September 15 1958

The well is terminated 10 ft.  above, below  the permanent ground surface

Was the well disinfected upon completion?  
Yes  No

Was the well sealed watertight upon completion?  
Yes  No

Signature J. J. SCHAFFER & SONS Fremont, Wisconsin  
Registered Well Driller Complete Mail Address

Rec'd SI-22 1958 No. 4111

Anal'd Safe

Interpretation Safe

10 ml 10 ml 10 ml 10 ml

Gas 34 hrs. 00000

48 hrs. 00000

Confirm \_\_\_\_\_

B. Coli \_\_\_\_\_

82

NOTE:

White Copy - Division's Copy  
Green Copy - Driller's Copy  
Yellow Copy - Owner's Copy

1. COUNTY Fond du Lac CHECK (1) ONE:  Town  Village  City Argon Name Argon

2. LOCATION Section SE 7 Township 16N Range 14E 3. NAME  OWNER  AGENT AT TIME OF DRILLING CHECK (A OR B)  
OR Grid of Street No. Street Name 6 FF ADDRESS Argon Oldfield Rd  
AND If available subdivision name, lot & block No. POST OFFICE Argon

4. Distance in feet from well to nearest: (Record answer in appropriate block)

Building	Sanitary Bldg. Drain	Sanitary Bldg. Sewer	Floor Drain Connected To	Storm Bldg. Drain	Storm Bldg. Sewer
<u>15</u>	C.I. Other	C.I. Other	C.I. Sewer Other Sewer	C.I. Other	C.I. Other

Street Sewer:  San.  Storm  C.I.  Other

Other Sewers:  Foundation Drain Connected To  Sewage Sump  Clearwater Sump

Clearwater Sump: 100

Septic Tank:  Holding Tank:  Sewage Absorption Unit: 125

Seepage Pit:  Seepage Bed:  Seepage Trench:

Privy:  Pet. Waste Pit:  Pit:  Nonconforming Existing:  Subsurface Pumproom:  Barn:  Animal Barn:  Animal Yard:  Silo With Pit:  Glass Lined Storage Facility:  Silo w/o Pit:  Earthen Silage Storage Trench (PH):

Temporary Manure Stack:  Watertight Liquid Manure Tank:  Solid Manure Storage Structure:  Subsurface Gasoline or Oil Tank:  Waste Pond or Land Disposal Unit (Specify Type):  Other (Give Description):

5. Well is intended to supply water for: Home

6. DRILLHOLE

Dia. (in.)	From (ft.)	To (ft.)	Dia. (in.)	From (ft.)	To (ft.)
<u>10</u>	<u>Surface</u>	<u>54</u>			
<u>6</u>	<u>54</u>	<u>180</u>			

7. CASING, LINER, CURBING AND SCREEN  
Material, Weight, Specifications & Method of Assembly

Dia. (in.)	From (ft.)	To (ft.)
<u>6</u>	<u>Surface</u>	<u>54</u>

NEW BLACK STEEL FE-1857 ASTM A53

9. FORMATIONS

Kind	From (ft.)	To (ft.)
<u>Clay</u>	<u>Surface</u>	<u>15</u>
<u>Traveler Sand</u>	<u>15</u>	<u>54</u>
<u>Sandstone</u>	<u>54</u>	<u>140</u>
<u>Sandstone</u>	<u>140</u>	<u>180</u>
<u>Water Bearing</u>		

8. GROUT OR OTHER SEALING MATERIAL

Kind	From (ft.)	To (ft.)
<u>Shury Alg Drilling</u>	<u>Surface</u>	<u>54</u>

10. TYPE OF DRILLING MACHINE USED

Cable Tool  Rotary-hammer w/drilling mud & air  Jetting with

Rotary-air w/drilling mud  Rotary-hammer & air  Air

Rotary-w/drilling mud  Reverse Rotary  Water

11. MISCELLANEOUS DATA

Field Test: 8 Hrs. at 10 GPM

Depth from surface to normal water level: 30 Ft.

Depth of water level when pumping: 30 Ft. Stabilized  Yes  No

Well construction completed on July 10 1972

Well is terminated 10 inches  above  below final grade

Well disinfected upon completion  Yes  No

Well sealed watertight upon completion  Yes  No

Water sample sent to Doherty laboratory on July 30 1972

Your opinion concerning other pollution hazards, information concerning difficulties encountered, and data relating to nearby wells, streams, etc., should be given on reverse side.

Signature: William Clark Complete Mail Address: 5411 Argon Rd. Argon

Registered Well Driller

NOV 23 1962

WELL CONSTRUCTOR'S REPORT TO WISCONSIN STATE BOARD OF HEALTH

151 6

See Instructions on Reverse Side

NOV 23 1962

1. County Fond du Lac Town  Village  City  Ripon (check one and give name)

2. Location Rte 7 - SE 1/4 - 116 - 1145  
Name of street and number of premise of Section, Town and Range numbers

3. Owner  or Agent  W. Muscarish  
Name of individual, partnership or firm

4. Mail Address Ripon, Wis.  
Complete address required

RECEIVED

5. From well to nearest: Building 6 ft; sewer      ft; drain      ft; septic tank 50 ft;  
dry well or filter bed      ft; abandoned well 50 ft

SEP 2 1962

6. Well is intended to supply water for: Home

7. DRILLHOLE:

Dia. (in.)	From (ft.)	To (ft.)	Dia. (in.)	From (ft.)	To (ft.)
10	0	76			
6	76	116			

8. CASING AND LINER PIPE OR CURBING:

Dia. (in.)	Kind and Weight	From (ft.)	To (ft.)
6	Steel	0	76

9. GROUT:

Kind	From (ft.)	To (ft.)
Cement	0	76

11. MISCELLANEOUS DATA:

Yield test: 3 Hrs at 25 GPM  
 Depth from surface to water-level: 30 ft  
 Water-level when pumping: 17 ft  
 Water sample was sent to the state laboratory at:  
Waukegan City on 3, 75 1962

10. FORMATIONS:

Kind	From (ft.)	To (ft.)
clay	0	15
loose sand	15	28
hard sand	28	65
soft fine sand	65	70
fine rock from	70	115
and gravel	115	116
etc., bearing		

Construction of the well was completed on:

September 25 1962

The well is terminated 10 inches  above, below  the permanent ground surface.

Was the well disinfected upon completion?

Yes  No

Was the well sealed watertight upon completion?

Yes  No

Signature Walter Hub  
Registered Well Driller

Clark W. Dilling  
Complete Mail Address

Please do not write in space below

Rec'd No

Ans'd     

Interpretation     

10 ml 10 ml 10 ml 10 ml 10 ml

Gas 34 hrs

45 hrs

Confirm

B. Coli

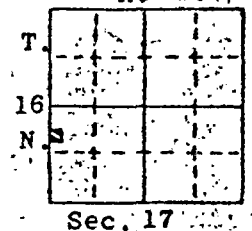
Examiner

Well name Wisconsin Power & Light Co. Well #9  
 Owner.... Wisconsin Power & Light Co.  
 Address.. 112 Watson St., P.O. Box 203  
 Ripon, WI 54971  
 Driller.. Layne-Northwest Co.  
 Engineer. Wisconsin Power & Light Co.  
 Madison, Wisconsin

County: Fond du Lac

R 14E

Completed... 6/72  
 Field check.  
 Altitude.... 843' ETM  
 Use..... Municipal  
 Static w.l.. 19' 5"  
 Spec. cap... 3.5 GPM/ft.



Quad. Ripon 7 1/2'

Drill Hole						Casing & Liner Pipe or Curbing							
Dia.	from	to	Dia.	from	to	Dia.	Wgt. & Kind	from	to	Dia.	Wgt. & Kind	from	to
24"	0	114'	23"	114'	320'	24"	3/8 wall steel bk A53 Grade B welded +6"	114'		16"	3/8 wall bk A53 Grade B welded	+12"	135'

Drilling method:  
 Samples from 0 to 320' Rec'd: 1/5/73  
 Studied by: Kathleen Massie

Grout	from	to
Neat Cement	0	135'

Formations: Surface, Drift, St. Peter Sandstone.

Issued: 5/6/83

Remarks: Well tested for 12 hours at 395 GPM with 110' 7" of drawdown.

LOG OF WELL:

Depth	Graphic Section	Rock Type	Color	Grain Size		Miscellaneous Characteristics
				Mode	Range	
0-5		Soil	Dk brown	—	—	Little gravel. Trace organic material.
5-10		Gravel	Mxd brown	L pnb	Gran/M pnb	Fos dolomite, dol, quartz, cht, dol cem ss, trap, granite. Ltl sand.
10-15		"	Mxd grey	M pnb	Gran/L pnb	Same but much sand.
15-20		Sand	"	C	Vfn/VC	Many dolomite frags. Much gravel. Ltl silt. Trace clay.
20-25		Gravel	"	M pnb	Gran/L pnb	Fos dol, dol, qtz, cht, grnt, trap. Much sand (most dol). Tr silt.
25-30		Sand	"	C	Vfn/VC	Many dolomite fragments. Much gravel. Trace silt.
30-35		Gravel	"	L pnb	Gran/M pnb	Fos dolomite, dolomite, quartz, chert, trap, granite. Much sand.
35-40		Sand	Brown	Fn	Vfn/VC	Dolomitic. Much silt. Little clay. Trace gravel.
40-45		"	"	C	"	Many dolomite fragments. Much gravel. Trace silt.
45-50		Clay	Gry brown	—	—	Calcareous. Much silt. Little sand.
50-55		Silt & cl	"	—	—	Calcareous. Little sand. Trace gravel.
55-60		Snd & silt	"	Fn	Vfn/VC	Calcareous. Much gravel, clay.
60-65		"	"	"	"	Same.
65-70		"	"	"	"	"
70-75		"	"	Fn/M	"	Same but little gravel.
75-80		Silt & cl	"	—	—	Calcareous. Much sand. Little gravel.
80-85		Snd & silt	"	M	Vfn/VC	Calcareous. Much clay. Little gravel.
85-90		Gravel	Bk & grey	Gran	Gran/M pnb	Gab, dior, fos dol, dol, qtz, grnt. Much sand.
90-95		Snd & silt	Gry brown	M	Vfn/VC	Calcareous. Much clay. Little gravel.
95-100		"	"	"	"	Same but much gravel.
100-105		"	Brown	"	"	Little dolomitic clay. Trace gravel.
105-110		Sandstone	Pl gry bn	M/C	"	Rounded. Much caved gravel.
110-115		"	V pl bn	M	"	Rounded. Tr G sil cem ss w/pnk chert. Trace caved ovl & sand.
115-120		"	Red	Fn	Vfn/VC	Srnd. Ltl V G sil cem, st. Mch hemic sh. Tr cvd ovl & end, wh
120-125		"	Yl red	M	"	See end of log. siliceous matx.
125-130		"	Red	"	"	Same but much hematitic shale.
130-135		"	Yl red	"	"	Srnd. Mch yl rd shale. Tr dk rd bn ss as above (cvd?), st, Fn
135-140		"	"	"	"	Same. glauconite.
140-145		"	Dull yl rd	"	"	Srnd. Ltl dull yl rd sh. Tr dk rd bn ss as above (cvd?), st, Fn/M
145-150		"	"	"	"	Same. glauconite.
150-155		"	Lt brown	"	"	Srnd to sang. Tr sil cem, Fn/M glauc, st, dk rd bn ss (as above, cvd)
155-160		"	Yl red	"	"	Srnd. Mch yl rd dolc sh, Ltl yl calcus sh (occ in layers?). st. Ltl pl yl shale. Trace Fn/M glauconite.



Well name: Wisconsin Power & Light Co. Well #9

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Depths	Graphic Section	Rock Type	Color	Grain Size		Miscellaneous Characteristics
				Mode	Range	
160-165		Sandstone	Yl red	Fn	Vfn/VC	Srnd. Mch yl rd dolc sh, st. Tr dol cem, Fn glauconite.
165-170		"	Rd yellow	M	"	Srnd. Mch red yellow dolc shale. Ltl silt. Tr Fn-glauconite.
170-175		"	"	Fn	"	Same plus trace good silica cement.
175-180		"	"	Fn&C	"	Srnd. Tr dol & lim cem, pnk cht, Fn glauc. Mch rd yl dolc sh, st.
180-185		"	"	C	"	Same plus trace pale yellow green dolomitic shale.
185-190		"	"	"	"	Same but no glauconite.
190-195		"	"	"	"	Srnd. Trace dolomite cem, pnk chert. Mch rd yl dolc sh, silt.
195-200		"	"	Fn&C	"	Same.
200-205		"	Strg bn	M	"	Same but strong brown shale.
205-210		"	Brown	C	"	Srnd. Tr G dol cem, Fn/M glauc, pl gn sh, pnk cht. Much dolc
210-215		"	"	"	"	Same. shale.
215-220		"	V pl brown	"	"	Srnd to sang. Tr P dol cem, Fn/M-glauc, mfc incl, pnk cht. Ltl
220-225		"	"	M/C	"	Srnd. Tr sil cem, Fn-glauc, mfc incl, pnk cht, silt. silt, shale.
225-230		"	"	"	"	Same.
230-235		"	"	"	"	Same but no cement.
235-240		"	"	C	"	Srnd. Tr pnk cht(w/tr fltg snd), Fn-glauc, mfc incl, pl gn micus
240-245		"	"	M/C	"	Same but much silt. shale. Ltl silt.
245-250		"	Lt rd brown	C	"	Srnd. Mch pnk dolc sh, pnk cht(w/tr rd bn hemic stng, bk hem, thin
250-255		"	"	"	"	Same plus tr pl gn shale. qtz layers, micus incl). Tr calc x. s.
255-260		Ss & chert	Rd yellow	M/C	"	See end of log.
260-265		"	"	"	"	Same plus tr mssv glauc, but tr dk rd bn hemic stng.
265-270		Chert	"	—	—	Ltl ss(as above), fltg snd, siliceous pnk sh. Few thin qtz layers
270-275		Sandstone	Pink	M	Vfn/C	Srnd. Tr G to F Tr st, mssv glauc, bk hem, dk rd bn hemic stng.
275-280		"	"	"	"	Same. sil cem, mssv glauc. Mch pnk sh, st. Ltl cht(as above)
280-285		"	Rd yellow	"	Vfn/VC	Srnd. Tr G sil cem, dol, rd yl cht. Mch siliceous rd yl sh. Ltl
285-290		"	Lt red	"	"	Same plus tr wh siliceous & rd bn hemic shale. silt.
290-295		"	"	"	"	Same but much silt.
295-300		"	Rd yellow	M/C	"	Srnd. Mch rd yl siliceous cl. Ltl st. Tr wh siliceous & rd bn
300-305		"	"	"	"	Same. hemic shale.
305-310		"	Pink	M	"	Srnd. Tr G sil cem, dol, pl gn shale. Ltl pnk siliceous shale.
310-315		"	Rd yellow	"	"	Srnd. Mch wh siliceous sh, pl gn micus sh, rd bn hemic sh. Tr st.
315-320		"	Pnk white	"	"	Subrounded. Little caved red yellow shale & sandstone. Trace silt, one quartz granule.

END OF LOG

120-125		Sandstone	Yl red	M	Vfn/VC	Subangular. Little good silica cement. Trace hematitic shale. Much hard dark red brown hematitic very glauconitic Fn sandstone with trace fossil fragments, also with trace silica cemented micaceous & glauconitic sandstone.
255-260		Ss & chert	Rd yellow	M/C	Vfn/VC	Subrounded. Trace calcite cement, black hematite & floating sand (chert), pale green shale. Few quartz layers (chert). Little dark red brown hematitic staining. Much dolomitic pink shale.

DATE 10/8/86

TDD # RS8310-015

TIME 11:35 AM

Page 1 of       

SITE Ripon Landfill

CONTACT Jim Hurlaud - County

PHONE (414) 929-3172

Extension Service.

REFERENCE 19

SITE NAME Ripon Landfill

SITE ID WFD 980610120

SUBJECT: Irrigation wells

Mr. Hurlaud stated he had no records concerning irrigation wells at his office, but to his knowledge did not think any irrigation wells existed in the area I described. He then forwarded my call to the county planner - sanitation office. The receptionist told the planner was on vacation, but to try to reach the sanitation at 1:00 pm or 4:00 pm at 414 929-3139.

<sup>10/9/86</sup> Was able to reach Wayne Rollin, County Planner at 929-3139. Mr. Rollin stated there were no irrigation wells near the Ripon Landfill.

Robert Hurlaud  
Date: 10/9/86



CITY OF RIPON  
RIPON, WISCONSIN 54971

OFFICE OF DIRECTOR OF PUBLIC WORKS

August 10, 1981

REFERENCE 20  
SITE NAME Ripon Landfill  
SITE ID WIP 980610190

RECEIVED  
AUG 11 1981  
HORICON AREA

REFERENCE \_\_\_\_\_  
SITE NAME \_\_\_\_\_

Mr. Kenneth E. Hein  
Dept. of Natural Resources  
Horicon Area Headquarters  
P.O. Box D  
Horicon, WI 53032

Dear Mr. Hein: Re: City of Ripon - Town of Ripon Landfill Lic.#467

This letter is to acknowledge your July 27, 1981 letter regarding 4410-1 and our meeting on July 7, 1981 held at the landfill site between you, Donohue's representatives and myself.

The information in your letter and the report from Donohue's was presented to the Board of Public Works at their August 3, 1981 meeting. There was a meeting date of August 31, 1981 set to meet with Donohue's representatives and you, if you can make that date. At that time, all facts will be presented and the necessary recommendations from the Board to the Council, to take whatever reasonable steps are required to improve present operations and also consideration for extending or expanding the existing landfill site.

Donohue's have completed the sampling and testing of the surface water at the site and there does not seem to be an over-abundance of contamination in what surface water is on site. Therefore, I will wait until I receive an answer as to whether or not the City should haul the water to the plant for processing or use the existing fill at site and fill area accordingly.

I would appreciate hearing from you regarding the August 31, 1981, 7 P.M. Meeting, and also the method you prefer for surface water handling.

If you need more information or have questions regarding the above mentioned, please contact my office.

Sincerely,

*Claude A. Bell*

Claude A. Bell  
Director of Public Works

CAL:sws  
Enc.

City of Ripon Wastewater Treatment Plant Sludge Application Record from Jan. 1, 1981 through June 30, 1981

Yr.	Farmland			Ripon Landfill Site		
	Loads	Gals.	% Solids	Loads	Gals.	% Sol.
1981	76	83,600	8.8	-		
"	45	49,500	9.1	24	26,400	9.1
"	3	3,300	9.5	67	73,700	9.5
"	-			80	88,000	9.7
y "	-			68	74,800	9.7
n. "	-			71	78,100	10.1

RECEIVED  
 AUG 11 1981  
 HORICON AREA



# ecology and environment, inc.

223 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60606, TEL. 312-663-9415

International Specialists in the Environmental Sciences

REFERENCE 21

SITE NAME Ripon Landfill

SITE ID WID 980610190

Date Received for Review: 8/9/85 Date Review Completed: 8/23/85

To: Bob Ningtgen

From: Cynthia Bachunas / ARIENE PRATE

Subject: Ripon L.F. R05-8510-CIR (UUC)

Sample Description: CASE # 4571 6 INCH MIXERS

Project Data Status: Still awaiting test water  
beginning

## FIT Data Review Findings:

All attached CR review.

As (by 5%), Pb, & Zn outside GC limits

Ca, Cu, Fe, Pb, Mg, Mn, Na+Zn detected in  
field blank

use these values with caution

## Additional Comments:

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

DATE: 8-6-85

SUBJECT: Review of Region V CLP Data  
Received for Review on 8/2/85

FROM: Curtis Ross, Director (SSCRL)  
Central Regional Laboratory Jay Thacker

TO: Data User: JH

We have reviewed the data for the following case(s).

SITE NAME: Ripon Landfill SMD Case No. 4571  
EPA Data Set No. SF 2477 No. of Samples: 6 D.U./Activity Numbers 1965/047500  
CRL No. 85FH10316 - 85FH10R06  
SMD Traffic No. MEC994 - MEC997 ; MEF209 - MEF210  
CLP Laboratory: RMAH Hrs. Required for Review: 1 hr.

Following are our findings.

This review covers 6 low concentration water samples. RECEIVED AUG 9 1985  
The spike recovery of T1 (42%) was biased low.

The EPA supplied 'true value' was not used to calculate the R on Form VII, the ICP Interference Check.

The following elements showed high (>10%) RPD's between the sample value and the serial dilution of that sample (MEF209), and the reported values for these elements are therefore estimates: Fe, Co, Mg, Mn and Na.

All other G.C. audits are O.K.

The field blank (MEF210) analysis shows high concentrations of Fe (1240 ug/L, J) and Cu (935 ug/L). The Cu and Fe data is questionable.

Jan F. Peis  
8/6/85

- ( ) Data are acceptable for use.
- (x) Data are acceptable for use with qualifications noted above.
- ( ) Data are preliminary - pending verification by Contractor Laboratory.
- ( ) Data are unacceptable.

cc: Dr. Alfred Haebeler/Joan Fisk/Gary Ward, EPA Support Services  
Ross K. Robeson, EMSL-Las Vegas  
Don Trees, CLP/Sample Management Office

U.S. EPA Contract Laboratory Program  
 Sample Management Office  
 P.O. Box 818 - Alexandria, VA 22313  
 703/557-2490 FTS: 8-557-2490

Date 7-31-85

COVER PAGE  
INORGANIC ANALYSIS DATA PACKAGE

Lab Name ROCKY MOUNTAIN ANALYTICAL  
 SOW No. 784

Case No. 4571  
 QC Report No. 55002

Sample Numbers

<u>EPA No.</u>	<u>Lab ID No.</u>	<u>EPA No.</u>	<u>Lab ID No.</u>
<u>MEC994</u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>MEC995</u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>MEC996</u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>MEF207</u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>MEF209</u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>MEF210</u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>[MEF999]</u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>                    </u>	<u>                    </u>	<u>                    </u>	<u>                    </u>
<u>                    </u>	<u>                    </u>	<u>                    </u>	<u>                    </u>

RECEIVED  
 JUL 31 1985  
 U.S. ENVIRONMENTAL PROTECTION AGENCY  
 550 S. CLARK STREET  
 CHICAGO, ILLINOIS 60605

Comments: 6 LOW WATERS FOR TASKS 1&2 ONLY.  
THE ICAP 10X SERIAL DILUTION FOR SAMPLE MEF209 IS IDENTIFIED AS  
IMFF9991.

ICP interelement and background corrections applied? Yes X No  
 if yes, corrections applied before X or after          generation of raw data.

Footnotes:

- NR - not required by contract at this time
- Form I: Value - If the result is a value greater than or equal to the instrument detection limit but less than the contract required detection limit, report the value in brackets (i.e., [10]). Indicate the method used with P (for ICP/Plasma AA) or F (for furnace).
- D - Indicates element was analyzed for but not detected. Report with the detection limit value (i.e., [ND]).
- I - Indicates a value estimated or not reported due to the presence of interference. Explanatory note included on cover page.
- S - Indicates value determined by Method of Standard Addition.
- R - Indicates spike sample recovery is not within control limits.
- X - Indicates duplicate analysis is not within control limits.
- + - Indicates the correlation coefficient for method of standard addition is less than 0.995.
- CV - Indicates Cold Vapor.
- AS - Indicates Automated Spectrophotometric.

J - Estimate due to high %D between sample & serial dilution of sample

DATA TRACKING - FORM I

CRL Data Set No. SF 2477 ERRIS No. \_\_\_\_\_

MO Case No. 4571 Site Name: Ripon Landfill

Name of Laboratory: RMAH Data User: Jit

No. of Samples: 6 Date Samples Received: 8/1/85

- 1. Have chain-of-custody records been received? YES  NO
- 2. Have Traffic reports been received? YES  NO
- 3. If no, are Traffic report numbers written on the chain-of-custody record? YES  NO
- 4. If no, which Traffic report numbers are missing?


5. Are basic data forms in? YES  NO

6. Number of samples claimed: 6 Number of samples received: 6

7. Checked by: Widia Feliciano Date: 8/2/85

8. Received by Contract Project Management Section: JD Date: 8.2.85

9. Review Started: 8/5/85 Reviewer Signature: Jan F. Pels

10. Total time spent on review: 1 hr Date review completed: 8/5/85

11. Copied (xeroxed) by: \_\_\_\_\_ Date: \_\_\_\_\_

12. Mailed to Data User by: \_\_\_\_\_ Date: \_\_\_\_\_

TO DATA USERS:

RECEIVED AUG 9 1985

Please fill in the blanks and return this form to:

Charles Elly, DPO, Region V, SSCRL

13. Data received by: Patricia Quinn Date: 8/21/85

14. Q.A. review received by: Patricia Quinn Date: 8/21/85

15. Received by CRL - CPM Section for file by: \_\_\_\_\_  
Date: \_\_\_\_\_



MW 5A

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

85FH10516

EPA Sample No.  
MBC994

Date 7-31-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME ROCKY MOUNTAIN ANALYTICAL  
SOW NO. 784  
LAB SAMPLE ID. NO. -

CASE NO. 4571

QC REPORT NO. 55062

Elements Identified and Measured

Concentration: Low X X Medium \_\_\_\_\_ Other \_\_\_\_\_  
Matrix: Water \_\_\_\_\_ Soil \_\_\_\_\_ Sludge \_\_\_\_\_

UG/L RECEIVED AUG 9 1985

1. ALUMINUM	25U	F	13. MAGNESIUM	55900	P J
2. ANTIMONY	31U	F	14. MANGANESE	67	P J
3. ARSENIC	3U	F	15. MERCURY	0.1U	CV
4. BARIUM	[60]	F	16. NICKEL	[7.6]	F
5. BERYLLIUM	0.3U	F	17. POTASSIUM	8300	F
6. CADMIUM	4U	F	18. SELENIUM	6U	F
7. CALCIUM	78500	P J	19. SILVER	3U	F
8. CHROMIUM	4U	F	20. SODIUM	84700	F J
9. COBALT	4U	F	21. THALLIUM	10U	F R
10. COPPER	[5.2]	F	22. TIN	17U	P R
11. IRON	2390	F J	23. VANADIUM	4U	F
12. LEAD	[3.6]	F	24. ZINC	[20]	P --

Cyanide NR Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
Lab Manager [Signature]

MWS

Form 1

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22318  
703/557-2490 FTS: 8-557-2490

85FH10517  
EPA Sample No.  
MEC995

Date 7-31-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME ROCKY MOUNTAIN ANALYTICAL  
SOW NO. 784  
LAB SAMPLE ID. NO. -

CASE NO. 4571  
QC REPORT NO. 55062

Elements Identified and Measured

Concentration: Low X X Medium \_\_\_\_\_ Other \_\_\_\_\_  
Matrix: Water \_\_\_\_\_ Soil \_\_\_\_\_ Sludge \_\_\_\_\_

UG/L RECEIVED AUG 9 1985

1. ALUMINUM	25U	P	13. MAGNESIUM	64700	P J
2. ANTIMONY	31U	P	14. MANGANESE	382	P J
3. ARSENIC	[6]	F	15. MERCURY	0.1U	CV
4. BARIUM	[89]	P	16. NICKEL	[15]	P
5. BERYLLIUM	0.3U	P	17. POTASSIUM	13300	P
6. CADMIUM	4U	P	18. SELENIUM	3U	F
7. CALCIUM	78300	P J	19. SILVER	3U	P
8. CHROMIUM	4U	P	20. SODIUM	82400	P J
9. COBALT	4U	P	21. THALLIUM	10U	F R
10. COPPER	[8]	P	22. TIN	17U	P R
11. IRON	4080	P J	23. VANADIUM	4U	F
12. LEAD	5.8	F	24. ZINC	78	P

Cyanide NR Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
Lab Manager D. H. Hatt

MWB

Form I

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

85FH10318  
EPA Sample No.  
MBC996

Date 7-31-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME ROCKY MOUNTAIN ANALYTICAL  
SOW NO. 784  
LAB SAMPLE ID. NO. -

CASE NO. 4571  
GC REPORT NO. 55052

Elements Identified and Measured

Concentration: Low X X Medium \_\_\_\_\_ Other \_\_\_\_\_  
Matrix: Water \_\_\_\_\_ Soil \_\_\_\_\_ Sludge \_\_\_\_\_

RECEIVED AUG 9 1985

UG/L

1. ALUMINUM	25U	P	13. MAGNESIUM	52000	P J
2. ANTIMONY	31U	P	14. MANGANESE	873	P J
3. ARSENIC	3U	F	15. MERCURY	0.1U	CV
4. BARIUM	(72)	P	16. NICKEL	(33)	P
5. BERYLLIUM	0.3U	P	17. POTASSIUM	10000	P
6. CADMIUM	4U	P	18. SELENIUM	3U	F
7. CALCIUM	121000	P J	19. SILVER	3U	P
8. CHROMIUM	4U	P	20. SODIUM	42800	P J
9. COBALT	(16)	P	21. THALLIUM	100U	F R
10. COPPER	(7.3)	P	22. TIN	17U	P R
11. IRON	182	P J	23. VANADIUM	4U	P
12. LEAD	<del>22</del> 21 SAM	F	24. ZINC	35	P

Cyanide NR Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: \_\_\_\_\_

Lab Manager ABhatt

MWB

U.S. EPA Contract Laboratory Program  
 Sample Management Office  
 P.O. Box 818 - Alexandria, VA 22313  
 703/557-2490 FTS: 8-557-2490

85FH10519  
 EPA Sample No.  
 M8C997

Date 7-31-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME ROCKY MOUNTAIN ANALYTICAL  
 SOW NO. 784  
 LAB SAMPLE ID. NO. -

CASE NO. 4571  
 QC REPORT NO. 55062

Elements Identified and Measured

Concentration: Low X X Medium \_\_\_\_\_ Other \_\_\_\_\_  
 Matrix: Water \_\_\_\_\_ Soil \_\_\_\_\_ Sludge \_\_\_\_\_

UG/L RECEIVED AUG 9 1985

1. ALUMINUM	25U	P	13. MAGNESIUM	110000	P J
2. ANTIMONY	31U	P	14. MANGANESE	457	P J
3. ARSENIC	[4.5]	F	15. MERCURY	0.1U	CV
4. BARIUM	[146]	P	16. NICKEL	74	P
5. BERYLLIUM	0.3U	P	17. POTASSIUM	5210	P
6. CADMIUM	4U	P	18. SELENIUM	5U	F
7. CALCIUM	113000	P J	19. SILVER	3U	P
8. CHROMIUM	[7.3]	P	20. SODIUM	435000	P J
9. COBALT	[29]	P	21. THALLIUM	30U	F R
10. COPPER	[21]	P	22. TIN	17U	P R
11. IRON	161	J P	23. VANADIUM	4U	P
12. LEAD	2U	F	24. ZINC	28	P

Cyanide NR Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: \_\_\_\_\_

Lab Manager [Signature]

U.S. EPA Contract Laboratory Program  
Sample Management Office  
P.O. Box 818 - Alexandria, VA 22313  
703/557-2490 FTS: 8-557-2490

85 FH10DA  
EPA Sample No.  
MEF209

Date 7-31-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME ROCKY MOUNTAIN ANALYTICAL  
SOW NO. 784  
LAB SAMPLE ID. NO. -

CASE NO. 4571  
QC REPORT NO. 55062

Elements Identified and Measured

Concentration: Low X X Medium \_\_\_\_\_ Other \_\_\_\_\_  
Matrix: Water \_\_\_\_\_ Soil \_\_\_\_\_ Sludge \_\_\_\_\_

UG/L **RECEIVED AUG 9 1985**

1. ALUMINUM	25U	P	13. MAGNESIUM	65500	P J
2. ANTIMONY	31U	P	14. MANGANESE	524	P J
3. ARSENIC	(7.5)	F	15. MERCURY	0.10	CV
4. BARIUM	(104)	P	16. NICKEL	(12)	P
5. BERYLLIUM	0.3U	P	17. POTASSIUM	13000	P
6. CADMIUM	(4.1)	P	18. SELENIUM	3U	F
7. CALCIUM	79400	P J	19. SILVER	3U	P
8. CHROMIUM	4U	P	20. SODIUM	82000	P J
9. COBALT	4U	P	21. THALLIUM	30U	F R
10. COPPER	3U	P	22. TIN	17U	P R
11. IRON	4360 J	F	23. VANADIUM	4U	P
12. LEAD	<del>1000</del> 2U <del>5PM</del>	P F	24. ZINC	36	P

Cyanide NR Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are encouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: \_\_\_\_\_

Lab Manager Shatt

U.S. EPA Contract Laboratory Program  
 Sample Management Office  
 P.O. Box 818 - Alexandria, VA 22315  
 703/557-2490 FTS: 8-557-2490

85FH10206  
 EPA Sample No.  
 MEF210

Date 7-31-85

INORGANIC ANALYSIS DATA SHEET

LAB NAME ROCKY MOUNTAIN ANALYTICAL  
 SOW NO. 784  
 LAB SAMPLE ID. NO. -

CASE NO. 4571  
 QC REPORT NO. 55062

Elements Identified and Measured

Concentration: Low X X Medium \_\_\_\_\_ Other \_\_\_\_\_  
 Matrix: Water \_\_\_\_\_ Soil \_\_\_\_\_ Sludge \_\_\_\_\_

UG/L

RECEIVED AUG 9 1985 3701

1. ALUMINUM	25U	P J	13. MAGNESIUM		P J
2. ANTIMONY	31U	P	14. MANGANESE	25	P J
3. ARSENIC	3U	P	15. MERCURY	0.1U	CV
4. BARIUM	12U	P	16. NICKEL	5U	P
5. BERYLLIUM	0.3U	P	17. POTASSIUM	990U	P
6. CADMIUM	4U	P	18. SELENIUM	3U	P
7. CALCIUM	[4150]	P J	19. SILVER	3U	P
8. CHROMIUM	4U	P	20. SODIUM	15100	P J
9. COBALT	4U	P	21. THALLIUM	3U	F R
10. COPPER	935	P	22. TIN	17U	P R
11. IRON	1240	P J	23. VANADIUM	4U	P
12. LEAD	[4]	F	24. ZINC	148	P

Cyanide NR Percent Solids (%) \_\_\_\_\_

Footnotes: For reporting results to EPA, standard result qualifiers are used as defined on Cover Page. Additional flags or footnotes explaining results are discouraged. Definition of such flags must be explicit and contained on Cover Page, however.

Comments: \_\_\_\_\_

Lab Manager J. Heath



# ecology and environment, inc.

111 WEST JACKSON BLVD., CHICAGO, ILLINOIS 60604, TEL. 312-853-9415

International Specialists in the Environment

Date Received for Review: 11/26/85 Date Review Completed: 12/9/85

To: Rob Hingtgen

From: Cynthia Pugh

Subject: Ripon Landfill (wisconsin)  
(RS-8310-1A)

Sample Description: Case # 4571: Low Water Organics

Project Data Status: COMPLETE

FIT Data Review Findings: Data Acceptable with Qualifications Below:

The following percent differences existed between Sample ED803 and Duplicate ED806: 25% vinylchloride; 24% Trans-1, 2-Dichloroethane; 25% toluene; 24% Total Xylenes; 40% Di-n-octylphthalate; and 20% bis (2-Ethylhexyl) phthalate. Benzene (9.0ug/l J) and Trichloroethene (22ug/l J) were detected in the sample (ED803) but were not detected in the Dup. Methylene chloride was detected in the Dup. but not in the sample (ED803.) Use data for above named compounds with caution.

Additional Comments:

The Attached Data has been received unreviewed by CRL. Although it has passed the CLP QA requirements, it will not be further reviewed to assure accuracy per regional requirements unless otherwise requested.

Book No. 4  
Page No. 278

UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION V

DATE: 9/9/85

RECEIVED NOV 26 1985

SUBJECT: Review of Region V CLP Data  
Received for Review on 7/29/85

FROM: Curtis Ross, Director (SSCRL) Francis Thomas  
Central Regional Laboratory

TO: Data User: Et

We have reviewed the data for the following case(s).

SITE NAME: Ripon Landfill SMO Case No. 4571  
EPA Data Set No. SF 2477 No. of D.U./Activity Samples: 6 Numbers 47051 048500  
CRL No. 85FH10316 - 85FH10319  
SMO Traffic No. ED802 - ED807  
CLP Laboratory: Wagleton Hrs. Required for Review: \_\_\_\_\_

Following are our findings.

Preliminary pending review by  
Peter Isaacson, Vian & Co.

- { } Data are acceptable for use.
- { } Data are acceptable for use with qualifications noted above.
- { } Data are preliminary - pending verification by Contractor Laboratory.
- { } Data are unacceptable.

cc: Dr. Alfred Haebeler/Joan Fisk/Gary Ward, EPA Support Services  
Ross K. Robeson, EMSL-Las Vegas  
Don Trees, CLP/Sample Management Office



July 23, 1985

Sample Management Office  
Viar and Company  
300 North Lee Street  
Alexandria, VA 22314

RECEIVED

U.S. ENVIRONMENTAL PROTECTION AGENCY  
LABORATORY

60605

Dear Gentlemen:

Enclosed please find the data packages for Case 4571 received on June 27, 1985 from Region V. All samples were analyzed according to the protocols provided under our Contract No. 68-01-6961.

Please note the following relating to the quality control items in this case:

- GC-MS Tuning - All tuning requirements for both BFB and DFTPP were within contract criteria.
- Instrumental Calibration - All initial and continuing calibration requirements for the volatile and semi-volatile fractions were within contract criteria. All initial and continuing calibration requirements for the pesticide analyses were met with the exception of the percent deviation of the individual mix B compound response factors run on 7-13-85 at 7:14:16. This case was run along with a second case immediately following it in sequence. The individual mix B was injected following three samples of the next case which contaminated (due to carry over) this standard. In this situation instead of rejecting the good run of all samples previously run under Case 4571 we reported the values of the contaminated standard but did not use it in the evaluation of the previously injected samples.
- Blanks - All reagent and instrumental blanks analyzed for all fractions were within contract criteria.
- Surrogate Recoveries - All surrogate recoveries for the volatile fraction were within contract criteria. Three surrogates were outside of contract criteria in the semi-volatile fraction d<sub>5</sub>-nitrobenzene (3% high)

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in both ED802MS and ED802MSD, as well as 2,4,6-Tribromophenol (7% high) in ED802MSD. One DBC recovery in the pesticide fraction was found outside the advisory limit, ED802MS (162%).

If you have any questions or if we may be of further service please feel free to call.

Sincerely,



David C. Hills  
Manager, Environmental Analysis


DCH/msw

cc: USEPA Region V  
USEPA EMSL-LV  
Central, File

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QC SUMMARY PACKAGE

Case No. 4571

 **HAZLETON** LABORATORIES AMERICA, INC  
Chemical & BioMedical Sciences Division  
3301 KINSMAN BLVD. P.O. BOX 7545  
MADISON, WISCONSIN 53707  
PHONE (608) 241-4471

# WATER SURROGATE PERCENT RECOVERY SUMMARY

Case No. 4571

Contract Laboratory HAZLETON

Contract No. 68-01-6961

SAS TRAFFIC NO.	VOLATILE						SEMI-VOLATILE				PESTICIDE
	TOLUENE-00 (00-110)	BFO (02-121)	1,2-DICHLORO-ETHANE-04 (17-120)	NITRO-BENZENE-08 (41-120)	2-FLUORO-BIPHENYL (44-119)	TERPHEYL-014 (33-120)	PHENOL-08 (10-103)	2-FLUORO-PHENOL (23-121)	2,4,6-TRIBROMO-PHENOL (10-130)	DIBUTYL-CHLOROPHATE (40-130)	
ED802	99.2	103	109	115	101	121	23.0	58.0	118	90	
ED803	97.6	103	110	112	98.2	119	33.5	78.4	129	132	
ED804	96.8	101	112	109	102	92.2	26.2	57.4	93.5	113	
ED805	98.2	102	117	115	99.0	72.8	39.1	23.0	114	87	
ED806	99.6	99.8	115	103	102	120	29.2	75.4	113	132	
ED807	99.2	99.2	111	91.5	86.6	124	39.4	111	111	107	
ED802HS	95.6	97.6	109	123*	100	117	28.3	49.5	128	162*	
ED802MSD	107	111	119	123*	99.7	121	26.6	50.9	137*	76	
MSL 28290	101	102	102	NR	NR	NR	NR	NR	NR	NR	
MS 28303	96.4	98.6	115	NR	NR	NR	NR	NR	NR	NR	
PAW 893	NR	NR	NR	92.4	81.5	121	40.4	118	111	NR	
PAW Post	NR	NR	NR	NR	NR	NR	NR	NR	NR	114	

RECEIVED NOV 26 1993

\* VALUES ARE OUTSIDE OF CONTRACT REQUIRED QC LIMITS  
 \*\* ADVISORY LIMITS ONLY

Volatiles: 0 out of 30 ; outside of QC limits  
 Semi-Volatiles: 3 out of 54 ; outside of QC limits  
 Pesticides: 1 out of 9 ; outside of QC limits

Comments: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

# WATER MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Case No. 4571

Contractor HAZLETON

Contract No. 68-01-6961

RECEIVED NOV 26 1985

FRACTION	COMPOUND	CONC. SPIKE ADDED (ug/L)	SAMPLE RESULT	CONC. MS	% REC	CONC. MSD	% REC	RPD	QC LIMITS*	
									RPD	RECOVERY
VOA SMO SAMPLE NO. <u>ED 802</u>	1,1-Dichloroethene	50.0	0.0	42.2	84.4	45.6	91.2	6.8	14	81-145
	Trichloroethene	50.0	0.0	47.2	94.4	53.9	108	13.6	14	71-120
	Chlorobenzene	50.0	0.0	48.8	97.6	53.6	107	9.4	13	75-130
	Toluene	50.0	0.0	49.6	99.2	66.0	112	12.8	13	78-125
	Benzene	50.0	0.0	49.8	99.6	64.2	108	8.4	11	78-127
B/N SMO SAMPLE NO. <u>ED 802</u>	1,2,4-Trichlorobenzene	100	0.0	92.9	92.9	93.3	93.3	0.42	28	39-98
	Acenaphthene	100	0.0	88.1	88.1	89.3	89.3	1.4	31	48-118
	2,4 Dinitrotoluene	100	0.0	36.8	36.8	45.0	45.0	20.0	38	24-96
	Di-n-Butylphthalate	100	2.7	89.4	89.4	89.0	89.0	0.22	40	11-117
	Pyrene	100	0.0	99.5	99.5	102	102	2.5	31	28-127
	N-Nitroso-Di-n-Propylamine	100	0.0	99.8	99.8	94.2	94.2	5.8	38	41-116
ACID SMO SAMPLE NO. <u>ED 802</u>	1,4-Dichlorobenzene	100	0.0	88.8	88.8	88.2	88.2	0.68	28	38-97
	Pentachlorophenol	200	0.0	168	78.8	156	78.1	0.89	50	9-103
	Phenol	200	0.0	48.1	24.1*	29.6	14.8*	48*	42	12-89
	2-Chlorophenol	200	0.0	106	83.0	110	65.0	3.7	40	27-123
	4-Chloro-3-Methylphenol	200	0.0	196	98.0	190	94.7	3.4	42	23-97
PEST SMO SAMPLE NO. <u>ED 802</u>	4-Nitrophenol	200	0.0	37	18.4	45.0	22.5	20.0	50	10-80
	Lindane	0.20	0.0	0.18	90.0	0.19	95.0	5.4	15	58-123
	Heptachlor	0.20	0.0	0.16	80.0	0.15	75.0	6.5	20	40-131
	Aldrin	0.20	0.0	0.13	65.0	0.11	55.0	17	22	40-120
	Dieldrin	0.50	0.0	0.42	84.0	0.36	72.0	15	18	52-126
	Endrin	0.50	0.0	0.51	102	0.45	90.0	13	21	58-121
	4,4'-DDT	0.50	0.0	0.37	74.0	0.32	64.0	15	27	38-127

\* ASTERISKED VALUES ARE OUTSIDE QC LIMITS.

RPD: VOAs 0 out of 5; outside QC limits  
 B/N 0 out of 7; outside QC limits  
 ACID 1 out of 5; outside QC limits  
 PEST 0 out of 6; outside QC limits

RECOVERY: VOAs 0 out of 10; outside QC limits  
 B/N 0 out of 14; outside QC limits  
 ACID 2 out of 10; outside QC limits  
 PEST 0 out of 12; outside QC limits

Comments: \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

# REAGENT BLANK SUMMARY

Case No. 4671

Contractor HAZLETON

Contract No. 68-01-6961

FILE ID	DATE OF ANALYSIS	FRACTION	MATRIX	CONC. LEVEL	INST. ID	CAS NUMBER	COMPOUND (HSL, TIC OR UNKNOWN)	CONC.	UNITS	CRED.
28290	7-3-85	VoA	H <sub>2</sub> O	LOW	HP5993	75-09-2	Methylene Chloride	4.0J	μg/L	5
	"	"	"	"	"	67-64-1	Acetone	18	"	10
BAN593	7-2-85	BAN	H <sub>2</sub> O	LOW	FINN 4500	PA 74-2	Di-n-butylphthalate	0.79	μg/L	10
	"	"	"	"	"	117-81-7	Bis-(2-ethyl hexyl) phthalate	26	"	10
RB1 Pest (7-D)	7-12-85	PEST.	H <sub>2</sub> O	LOW	5779	-	None Found	-	-	-

RECEIVED NOV 26 1985

Comments:

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**Bromofluorobenzene (BFB)** RECEIVED NOV 26 1985

Case No. 4571 Contractor HAZLETON Contract No. 68-01-6961  
 Instrument ID HP 5993 Date 5-10-85 Time 1130  
 Lab ID 27833 Data Release Authorized By: David C. Hill

m/e ION ABUNDANCE CRITERIA %RELATIVE ABUNDANCE

50	15.0 - 40.0% of the base peak	21.9	
75	30.0 - 60.0% of the base peak	53.7	
95	Base peak, 100% relative abundance	100.0	
96	5.0 - 9.0% of the base peak	8.80	
173	Less than 1.0% of the base peak	0.0	
174	Greater than 50.0% of the base peak	95.2	
175	5.0 - 9.0% of mass 174	7.50	(7.88) <sup>1</sup>
176	Greater than 85.0%, but less than 101.0% of mass 174	92.7	(97.4) <sup>1</sup>
177	5.0 - 9.0% of mass 176	7.50	(8.09) <sup>2</sup>

<sup>1</sup>Value in parenthesis is % mass 174.

<sup>2</sup>Value in parenthesis is % mass 176.

THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING SAMPLES, BLANKS AND STANDARDS.

SAMPLE ID	FRN	LAB ID	DATE OF ANALYSIS	TIME OF ANALYSIS
50 ppb Std	27836		5-10-85	1820
200 ppb Std	27837		"	1440
150 ppb Std	27838		"	1640
100 ppb Std	27839		"	1730
20 ppb Std	27840		"	1815

Bromofluorobenzene (BFB) RECEIVED NOV 26 1985

Case No. 4571 Contractor Hazleton Contract No. 68-01-6961  
 Instrument ID HP5993 Date 7-3-85 Time 1130  
 Lab ID 28289 Data Release Authorized By: Robert A. Thew

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE	
50	15.0 - 40.0% of the base peak	26.0	
75	30.0 - 60.0% of the base peak	58.1	
95	Base peak, 100% relative abundance	100	
96	5.0 - 9.0% of the base peak	8.8	
173	Less than 1.0% of the base peak	0	
174	Greater than 50.0% of the base peak	70.3	
175	5.0 - 9.0% of mass 174	5.5	(7.82) <sup>1</sup>
176	Greater than 95.0%, but less than 101.0% of mass 174	67.2	(95.6) <sup>1</sup>
177	5.0 - 9.0% of mass 176	5.6	(8.33) <sup>2</sup>

<sup>1</sup>Value in parenthesis is % mass 174.<sup>2</sup>Value in parenthesis is % mass 176.

THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING  
 SAMPLES, BLANKS AND STANDARDS.

SAMPLE ID	FRN	LAB ID	DATE OF ANALYSIS	TIME OF ANALYSIS
Inst. Blank	28290		7-3-85	1150
50ppb V STD	28291			1305
ED 803	28293			1440
ED 802	28294			1525
ED 802 MS	28296			1705
ED 802 MSD	28297			1745
ED 804	28298			1830
ED 805	28299			1920
ED 806	28301			2100
ED 807	28302			2145
Holding Blank	28303		↓	2230

6/84

FORM V



**GC/MS TUNING AND MASS CALIBRATION**  
**Decafluorotriphenylphosphine (DFTPP)**

Case No. 4571 Contractor HAZLETON Contract No. 68-01-6961  
 Instrument ID Finn 4500 Date 7/1/85 Time 1444  
 Lab ID DFTPP 70/850 Data Release Authorized By: [Signature]

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	54.3
68	less than 2.0% of mass 69	0.0 (0.0) <sup>1</sup>
69	mass 69 relative abundance	60.1
70	less than 2.0% of mass 69	0.0 (0.0) <sup>1</sup>
127	40.0 - 60.0% of mass 198	56.9
197	less than 1.0% of mass 198	0.0
198	base peak, 100% relative abundance	100
199	5.0 - 9.0% of mass 198	6.46
275	10.0 - 30.0% of mass 198	17.4
365	greater than 1.00% of mass 198	1.61
441	present, but less than mass 443	6.24
442	greater than 40.0% of mass 198	48.7
443	17.0 - 23.0% of mass 442	9.19 (8.9) <sup>2</sup>

<sup>1</sup> Value in parenthesis is % mass 69.  
<sup>2</sup> Value in parenthesis is % mass 442.

THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING  
 SAMPLES, BLANKS AND STANDARDS.

SAMPLE ID	LAB ID	DATE OF ANALYSIS	TIME OF ANALYSIS
BAN 585	10 PPM BAN STD	7/1/85	1615
BAN 587	25 PPM BAN STD	7/1/85	1810
BAN 588	40 PPM BAN STD		1907
BAN 589	60 PPM BAN STD		2005
BAN 590	80 PPM BAN STD	√	2102

**Decafluorotriphenylphosphine (DFTPP)**

Case No. 4571 Contractor HAZLETON Contract No. 68-01-6961  
 Instrument ID Finn 4500 Date 7/2/85 Time 1548  
 Lab ID DFTPP70285B Data Release Authorized By: John Mathews

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	48.1
68	less than 2.0% of mass 69	0 (0) <sup>1</sup>
69	mass 69 relative abundance	53.1
70	less than 2.0% of mass 69	0 (0) <sup>1</sup>
127	40.0 - 60.0% of mass 198	52.2
187	less than 1.0% of mass 198	0
198	base peak, 100% relative abundance	100
199	5.0 - 9.0% of mass 198	6.25
275	10.0 - 30.0% of mass 198	18.1
385	greater than 1.00% of mass 198	1.71
441	present, but less than mass 443	6.52
442	greater than 40.0% of mass 198	52.4
443	17.0 - 23.0% of mass 442	9.81 (8.7) <sup>2</sup>

<sup>1</sup> Value in parenthesis is % mass 69.  
<sup>2</sup> Value in parenthesis is % mass 442.

THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING SAMPLES, BLANKS AND STANDARDS.

SAMPLE ID	LAB ID	DATE OF ANALYSIS	TIME OF ANALYSIS
BAN 604	ONGOING CALIBRATION	7/2/85	1619
BAN 593	BLANK 7/1/85		1808
ED 807	BAN 594		1906
ED 806	BAN 595		2004
ED 804	BAN 597		2200
ED 803	BAN 598		2257

Decafluorotriphenylphosphine (DFTPP)

Case No. 4571 Contractor HAZLETON Contract No. 68-01-6961  
 Instrument ID  Finn 4500 Date 7/5/85 Time 8:14  
 Lab ID DFTPP 7085 Data Release Authorized By: John Mathew

m/e ION ABUNDANCE CRITERIA %RELATIVE ABUNDANCE

51	30.0 - 60.0% of mass 198	59.3	
68	less than 2.0% of mass 69	0.0	(0.0) <sup>1</sup>
69	mass 69 relative abundance	61.2	
70	less than 2.0% of mass 69	0.0	(0.0) <sup>1</sup>
127	40.0 - 60.0% of mass 198	59.3	
197	less than 1.0% of mass 198	0.0	
198	base peak, 100% relative abundance	100	
199	5.0 - 9.0% of mass 198	6.15	
275	10.0 - 30.0% of mass 198	17.5	
365	greater than 1.00% of mass 198	1.96	
441	present, but less than mass 443	6.07	
442	greater than 40.0% of mass 198	48.5	
443	17.0 - 23.0% of mass 442	8.95	(18.4) <sup>2</sup>

<sup>1</sup> Value in parenthesis is % mass 69.  
<sup>2</sup> Value in parenthesis is % mass 442.

THIS PERFORMANCE TUNE APPLIES TO THE FOLLOWING SAMPLES, BLANKS AND STANDARDS.

SAMPLE ID	LAB ID	DATE OF ANALYSIS	TIME OF ANALYSIS
BANG12	ONGOING CALIBRATION	7/5/85	10:41
ED802	BAN 608	7/5/85	12:21
ED805	BAN 609	7/5/85	13:19
ED802MS	BAN 610	7/5/85	14:17
ED802MSD	BAN 611	7/5/85	15:15

Case No. 4577 Laboratory ANALYTICAL  
 Contract No. 68-01-6961 GC Column 1.5% SP-2250/1.45% SP-2401 ON 100/120 SUPER COBALT  
 Date of Analysis 7-12-85 Instrument ID 5779

EVALUATION CHECK FOR LINEARITY

LABORATORY ID	397	398	399	
PESTICIDE	CALIBRATION FACTOR EVAL. MIX A	CALIBRATION FACTOR EVAL. MIX B	CALIBRATION FACTOR EVAL. MIX C	% RSD (≤10%)
ALDRIN	$8.4 \times 10^5$	$9.1 \times 10^5$	$8.6 \times 10^5$	3
ENDRIN	$5.3 \times 10^5$	$5.7 \times 10^5$	$5.6 \times 10^5$	3
4,4'-DDT <sup>(1)</sup>	$5.2 \times 10^5$	$6.4 \times 10^5$	$5.6 \times 10^5$	9
DIBUTYL CHLORENDATE	$7.1 \times 10^5$	*	$7.3 \times 10^5$	1

RECEIVED NOV 26 1985

EVALUATION CHECK FOR 4,4'-DDT/ENDRIN BREAKDOWN

LABORATORY ID	PERCENT BREAKDOWN EXPRESSED AS TOTAL DEGRADATION			
	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B
LABORATORY ID	398	422	437	
TIME OF ANALYSIS	10:18:40 7-12-85	3:03:08 7-13-85	22:24:42 7-14-85	
ENDRIN	3.9	CROSS CONTAMINATION FROM A PREVIOUS	1.0	
4,4'-DDT	7.3	SAMPLE PREVENTED 4,4'-DDT/ENDRIN BREAKDOWN CALCULATION	13	
COMBINED <sup>(2)</sup>	-	-	-	

EVALUATION OF RETENTION TIME SHIFT FOR DIBUTYLCHLORENDATE

SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF.	SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF.
BLANK (7-1-85)	411	19:22:42	0.2				
ED 802	412	20:04:33	0.1				
ED 802 MS	413	20:46:25	0.2				
ED 802 MSD	414	21:28:16	0.03				
ED 803	415	22:10:08	0.1				
ED 804	417	23:33:52	0.05				
ED 805	418	0:15:43	0.1				
ED 806	419	0:57:34	0.2				
ED 807	420	1:39:25	0.03				

(1) SEE EXHIBIT E, SECTION 7.5.4

(2) SEE EXHIBIT E, SECTION 7.3.1.2.2.1

\* DIBUTYL CHLORENDATE WAS LEFT OUT BECAUSE IT CO-ELUDES WITH ENDRIN KETONE (ON THIS COLUMN) AND PREVENTS ENDRIN BREAKDOWN CALCULATIONS.

Case No. 4571 Low water's Laboratory HAZLETON

Contract No. 68-01-6961 GC Column 330V-1

Date of Analysis 7-12-85 Instrument ID 278211

RECEIVED NOV 26 1985

EVALUATION CHECK FOR LINEARITY

LABORATORY ID	403	404	405	
PESTICIDE	CALIBRATION FACTOR EVAL. MIX A	CALIBRATION FACTOR EVAL. MIX B	CALIBRATION FACTOR EVAL. MIX C	% RSD (±10%)
ALDRIN	4.1x10 <sup>5</sup>	5.2x10 <sup>5</sup>	5.4x10 <sup>5</sup>	12
ENDRIN	3.1x10 <sup>5</sup>	3.8x10 <sup>5</sup>	3.9x10 <sup>5</sup>	10
4,4'-DDT <sup>(1)</sup>	2.1x10 <sup>5</sup>	2.4x10 <sup>5</sup>	2.6x10 <sup>5</sup>	8.7
DIBUTYL CHLORENDATE	4.7x10 <sup>5</sup>	*	4.7x10 <sup>5</sup>	0.0

EVALUATION CHECK FOR 4,4'-DDT/ENDRIN BREAKDOWN

LABORATORY ID	PERCENT BREAKDOWN EXPRESSED AS TOTAL DEGRADATION			
	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B
LABORATORY ID	404	428	440	
TIME OF ANALYSIS	11:30:48 7/12/85	4:12:26 7/13/85	12:33:54 7/13/85	
ENDRIN	—	—	—	
4,4'-DDT	—	—	—	
COMBINED <sup>(2)</sup>	5.8	10.4	9.6	

EVALUATION OF RETENTION TIME SHIFT FOR DIBUTYLCHLORENDATE

SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF.	SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF.
EVAL Mix A		10:49:12	0.0				
EVAL Mix C		12:12:24	0.5%				
EPA Mix A		12:54:00	0.5%				
Blank		20:33:51	0.6%				
ED 802		21:15:39	0.4%				
ED 802 MS		21:57:19	0.5%				
ED 802 MS D		22:38:59	0.7%				
ED 803		23:20:40	0.1%				
EPA Mix A		2:02:21	0.3%				
ED 804		0:44:01	0.2%				
ED 805		1:25:40	0.1%				
ED 806		2:07:21	0.1%				
ED 807		2:49:03	0.4%				

(1) SEE EXHIBIT E, SECTION 7.5.4

(2) SEE EXHIBIT E, SECTION 7.3.1.2.2.1

# Pesticide/PCB Standards Summary

Case No. 4571 Laboratory HAZLETON LABORATORIES  
 Contract No. 68-01-6961 GC Column 1.5% SP2250/1.95% SP2301 GC Instrument ID 5779

INDIVIDUAL MIXA

DATE OF ANALYSIS <u>① 7-12-85</u>	DATE OF ANALYSIS <u>② 7-13-85</u>
TIME OF ANALYSIS <u>11:42:20</u>	TIME OF ANALYSIS <u>1:54:04</u>
LABORATORY ID <u>400</u>	LABORATORY ID <u>428</u>

COMPOUND	RT	RETENTION TIME WINDOW	CALIBRATION FACTOR	CONF. OR QUANT.	RT	CALIBRATION FACTOR	CONF. OR QUANT.	PERCENT DIFF. **
alpha-BHC	2.30	±.03	1.1 x 10 <sup>6</sup>	QUANT.	2.29	1.1 x 10 <sup>6</sup>	QUANT.	0.0
beta-BHC	3.36	±.03	4.6 x 10 <sup>5</sup>		3.35	4.7 x 10 <sup>5</sup>		2.1
delta-BHC	3.92	±.03	8.6 x 10 <sup>5</sup>		3.91	8.7 x 10 <sup>5</sup>		1.1
gamma-BHC	2.92	±.04	8.7 x 10 <sup>5</sup>		2.90	8.9 x 10 <sup>5</sup>		2.2
Haptachlor	3.60	±.04	8.9 x 10 <sup>5</sup>		3.58	9.0 x 10 <sup>5</sup>		1.1
Aldrin	4.37	±.06	8.7 x 10 <sup>5</sup>		4.34	9.0 x 10 <sup>5</sup>		3.3
Haptachlor Epoxide	6.60	±.13	8.3 x 10 <sup>5</sup>		6.54	8.6 x 10 <sup>5</sup>		3.5
Endosulfan I	8.32	±.15	7.7 x 10 <sup>5</sup>		8.25	8.1 x 10 <sup>5</sup>		4.9
Dieldrin	10.22	±.17	8.2 x 10 <sup>5</sup>		10.14	8.6 x 10 <sup>5</sup>		5.0
4,4'-DDE	9.74	±.04	7.9 x 10 <sup>5</sup>		9.72	8.0 x 10 <sup>5</sup>		1.3
Endrin	12.46	±.04	5.7 x 10 <sup>5</sup>		12.44	9.1 x 10 <sup>5</sup>		37 *
Endosulfan II	15.15	±.25	8.2 x 10 <sup>5</sup>		15.03	1.0 x 10 <sup>6</sup>		18 *
4,4'-DDD	15.01	±.11	6.5 x 10 <sup>5</sup>		14.96	1.2 x 10 <sup>6</sup>		46 *
Endrin Aldehyde	19.85	±.34	6.9 x 10 <sup>5</sup>		19.68	1.1 x 10 <sup>6</sup>		37 *
Endosulfan Sulfate	23.92	±.13	5.1 x 10 <sup>5</sup>		23.86	6.4 x 10 <sup>5</sup>		20 *
4,4'-DDT	18.17	±.27	6.0 x 10 <sup>5</sup>		18.04	4.3 x 10 <sup>5</sup>		28 *
Methoxychlor	35.60	±.60	3.6 x 10 <sup>5</sup>		35.30	2.6 x 10 <sup>5</sup>		28 *
Endrin Ketone	32.91	±.11	8.5 x 10 <sup>5</sup>		32.86	8.5 x 10 <sup>5</sup>		0.0
Tech. Chlordane	nr	nr	nr					
alpha-Chlordane	nr	nr	nr					
gamma-Chlordane	nr	nr	nr					
Toxaphene	nr	nr	nr					
Aroclor - 1016	nr	nr	nr					
Aroclor - 1221	nr	nr	nr					
Aroclor - 1232	nr	nr	nr					
Aroclor - 1242	nr	nr	nr					
Aroclor - 1248	nr	nr	nr					
Aroclor - 1254	nr	nr	nr					
Aroclor - 1260	nr	nr	nr	↓			↓	

\* SEE EXHIBIT B, PART 7 \* DUE TO CROSS-CONTAMINATION OF HIGHLY CONCENTRATED SAMPLES - STANDARDS WERE NOT WITHIN CONTRACT LIMITS.

\*\* CONF. = CONFIRMATION (<20% DIFFERENCE)  
 QUANT. = QUANTITATION (<10% DIFFERENCE)

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26

**Pesticide/PCB Standards Summary**

Case No. 4571 Low waters Laboratory HAZLETON LABORATORIES  
 Contract No. 68-01-6961 GC Column 3%OV-1 GC Instrument ID 278211  
 Mix A Mix B Mix A Mix B

DATE OF ANALYSIS <u>July 12, 1985</u>	<u>July 12, 1985</u>	DATE OF ANALYSIS <u>7/13/85</u>	<u>7/13/85</u>
TIME OF ANALYSIS <u>12:54:00</u>	<u>13:35:39</u>	TIME OF ANALYSIS <u>0:02:21</u>	<u>8:22:38</u>
LABORATORY ID _____	_____	LABORATORY ID _____	_____

COMPOUND	RT	RETENTION TIME WINDOW	CALIBRATION FACTOR	CONF. OR QUANT.	RT	CALIBRATION FACTOR	CONF. OR QUANT.	PERCENT DIFF. **
alpha-BHC	2.10	±.02	5.5x10 <sup>5</sup>		2.10	5.2x10 <sup>5</sup>		5.4
beta-BHC	2.30	±.02	2.8x10 <sup>5</sup>		2.30	2.6x10 <sup>5</sup>		7.1
delta-BHC	2.63	±.02	4.4x10 <sup>5</sup>		2.63	4.3x10 <sup>5</sup>		2.3
gamma-BHC	2.49	±.02	5.2x10 <sup>5</sup>		2.50	4.5x10 <sup>5</sup>		13.5
Heptachlor	3.99	±.02	5.5x10 <sup>5</sup>		3.99	4.9x10 <sup>5</sup>		10.9
Aldrin	4.95	±.02	4.8x10 <sup>5</sup>		4.96	3.9x10 <sup>5</sup>		18.8
Heptachlor Epoxide	6.12	±.02	5.2x10 <sup>5</sup>		6.13	4.2x10 <sup>5</sup>		19.2
Endosulfan I	7.64	±.02	4.6x10 <sup>5</sup>		7.65	3.9x10 <sup>5</sup>		15.2
Dieldrin	8.96	±.03	4.6x10 <sup>5</sup>		8.98	3.9x10 <sup>5</sup>		15.2
4,4'-DDE	8.91	±.03	4.7x10 <sup>5</sup>		8.91	4.3x10 <sup>5</sup>		8.5
Endrin	10.08	±.03	3.6x10 <sup>5</sup>		10.08	3.8x10 <sup>5</sup>		5.6
Endosulfan II	10.42	±.03	4.9x10 <sup>5</sup>		10.44	4.2x10 <sup>5</sup>		14.3
4,4'-DDD	11.42	±.03	3.6x10 <sup>5</sup>		11.42	3.2x10 <sup>5</sup>		11.1
Endrin Aldehyde	11.68	±.03	4.9x10 <sup>5</sup>		11.70	4.2x10 <sup>5</sup>		14.3
Endosulfan Sulfate	13.56	±.02	3.2x10 <sup>5</sup>		13.55	3.4x10 <sup>5</sup>		6.2
4,4'-DDT	14.91	±.06	2.5x10 <sup>5</sup>		14.95	2.3x10 <sup>5</sup>		8.0
Methoxychlor	22.73	±.04	5.7x10 <sup>5</sup>		22.76	9.2x10 <sup>5</sup>		61.4*
Endrin Ketone	17.68	±.02	4.4x10 <sup>5</sup>		17.67	4.6x10 <sup>5</sup>		4.5
Tech. Chlordane	MR	MR	MR		MR	MR		
alpha-Chlordane	↓	↓	↓		↓	↓		
gamma-Chlordane	↓	↓	↓		↓	↓		
Toxaphene	↓	↓	↓		↓	↓		
Aroclor - 1016	↓	↓	↓		↓	↓		
Aroclor - 1221	↓	↓	↓		↓	↓		
Aroclor - 1232	↓	↓	↓		↓	↓		
Aroclor - 1242	↓	↓	↓		↓	↓		
Aroclor - 1248	↓	↓	↓		↓	↓		
Aroclor - 1254	↓	↓	↓		↓	↓		
Aroclor - 1260	↓	↓	↓		↓	↓		

\* SEE EXHIBIT B, PART 7 \* Chromatogram shows large interference  
 \*\* CONF. = CONFIRMATION (<20% DIFFERENCE)  
 QUANT. = QUANTITATION (<10% DIFFERENCE)

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Case No. 4571 Low waters

Laboratory

HAZLETON LABORATORIES

Contract No. 68-01-6961

GC Column

3% OV-1

GC Instrument ID 278211

Mix A

Mix B

Mix A

Mix B

DATE OF ANALYSIS July 12, 1985  
 TIME OF ANALYSIS 12:54:00  
 LABORATORY ID \_\_\_\_\_

DATE OF ANALYSIS 7/12/85  
 TIME OF ANALYSIS 0:02:21  
 LABORATORY ID \_\_\_\_\_

COMPOUND	RT	RETENTION TIME WINDOW	CALIBRATION FACTOR	(CONF. OR QUANT.)	RT	CALIBRATION FACTOR	(CONF. OR QUANT.)	PERCENT DIFF. **
alpha-BHC	2.10	±.02	5.5x10 <sup>5</sup>		2.10	5.2x10 <sup>5</sup>		5.4
beta-BHC	2.30	±.02	2.8x10 <sup>5</sup>		2.30	2.6x10 <sup>5</sup>		7.1
delta-BHC	2.63	±.02	4.4x10 <sup>5</sup>		2.63	4.3x10 <sup>5</sup>		2.3
gamma-BHC	2.79	±.02	5.2x10 <sup>5</sup>		2.50	4.5x10 <sup>5</sup>		13.5
Hepachlor	3.99	±.02	5.5x10 <sup>5</sup>		3.99	4.9x10 <sup>5</sup>		10.9
Aldrin	4.95	±.02	4.8x10 <sup>5</sup>		4.96	3.9x10 <sup>5</sup>		18.8
Hepachlor Epoxide	6.12	±.02	5.2x10 <sup>5</sup>		6.13	4.2x10 <sup>5</sup>		19.2
Endosulfan I	7.64	±.02	4.6x10 <sup>5</sup>		7.65	3.9x10 <sup>5</sup>		15.2
Dieldrin	8.96	±.03	4.6x10 <sup>5</sup>		8.98	3.9x10 <sup>5</sup>		15.2
4,4'-DDE	8.91	±.03	4.7x10 <sup>5</sup>		8.91	4.3x10 <sup>5</sup>		8.5
Endrin	10.08	±.03	3.6x10 <sup>5</sup>		10.08	3.9x10 <sup>5</sup>		5.6
Endosulfan II	10.42	±.03	4.9x10 <sup>5</sup>		10.44	4.2x10 <sup>5</sup>		14.3
4,4'-DDD	11.42	±.03	3.6x10 <sup>5</sup>		11.42	3.2x10 <sup>5</sup>		11.1
Endrin Aldehyde	11.68	±.03	4.9x10 <sup>5</sup>		11.70	4.2x10 <sup>5</sup>		14.3
Endosulfan Sulfate	13.56	±.02	3.2x10 <sup>5</sup>		13.55	3.4x10 <sup>5</sup>		6.2
4,4'-DDT	14.91	±.06	2.5x10 <sup>5</sup>		14.95	2.3x10 <sup>5</sup>		8.0
Methoxychlor	22.73	±.04	5.7x10 <sup>5</sup>		22.76	9.2x10 <sup>5</sup>		61.4*
Endrin Ketone	17.68	±.02	4.4x10 <sup>5</sup>		17.67	4.6x10 <sup>5</sup>		4.5
Tech. Chlordane	MR	MR	MR		MR	MR		
alpha-Chlordane*								
gamma-Chlordane*								
Toxaphene								
Aroclor - 1016								
Aroclor - 1221								
Aroclor - 1232								
Aroclor - 1242								
Aroclor - 1248								
Aroclor - 1254								
Aroclor - 1260								

\* SEE EXHIBIT B, PART 7 \* Chromatogram shows large interference  
 \*\* CONF. = CONFIRMATION (<20% DIFFERENCE)  
 QUANT. = QUANTITATION (<10% DIFFERENCE)



RECEIVED NOV 26 1985

Case Number  
 ED 802

Organics Analysis Data Sheet  
 (Page 1)

85FH10516

Laboratory Name: HAZLETON LABORATORIES  
 Lab Sample ID No: 50606606  
 Sample Matrix: WATER  
 Data Release Authorized By: David C. Kelly

Case No: 4571  
 QC Report No: \_\_\_\_\_  
 Contract No: 68-01-6261  
 Date Sample Received: 6-28-85

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 7-3-85 ERN 28294  
 Date Analyzed: 7-3-85 1525  
 Conc/Dil Factor: 1 pH \_\_\_\_\_  
 Percent Moisture: \_\_\_\_\_  
 Percent Moisture (Decanted): \_\_\_\_\_

CAS Number	Compound	ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	10 U
74-83-9	Bromomethane	10 U
75-01-4	Vinyl Chloride	10 U
75-00-3	Chloroethane	10 U
75-09-2	Methylene Chloride	5 U
67-64-1	Acetone	<u>(18 B)</u> 10 U
75-15-0	Carbon Disulfide	5 U
75-35-4	1, 1-Dichloroethane	5 U
75-34-3	1, 1-Dichloroethane	5 U
156-60-5	Trans-1, 2-Dichloroethane	5 U
67-66-3	Chloroform	5 U
107-06-2	1, 2-Dichloroethane	5 U
78-93-3	2-Butanone	<u>(V4)</u> 10 U
71-55-6	1, 1, 1-Trichloroethane	5 U
56-23-5	Carbon Tetrachloride	5 U
108-05-4	Vinyl Acetate	10 U
75-27-4	Bromodichloromethane	5 U

CAS Number	Compound	ug/l or ug/Kg (Circle One)
79-34-5	1, 1, 2, 2-Tetrachloroethane	5 U
78-87-5	1, 2-Dichloropropane	5 U
10061-02-6	Trans-1, 3-Dichloropropene	5 U
79-01-6	Trichloroethane	5 U
124-48-1	Dibromochloromethane	5 U
79-00-5	1, 1, 2-Trichloroethane	5 U
71-43-2	Benzene	5 U
10061-01-5	cis-1, 3-Dichloropropene	5 U
110-75-8	2-Chloroethylvinylether	10 U
75-25-2	Bromoform	5 U
591-78-6	2-Hexanone	10 U
108-10-1	4-Methyl-2-Pentanone	10 U
127-18-4	Tetrachloroethene	5 U
108-88-3	Toluene	5 U
108-90-7	Chlorobenzene	5 U
100-41-4	Ethylbenzene	5 U
100-42-5	Styrene	5 U
	Total Xylenes	5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- Value** If the result is a value greater than or equal to the detection limit, report the value.
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution actions. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10J)

- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/ul in the final extract should be confirmed by GC/MS.
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Laboratory Name: HAZLETONCase No: 4571

RECEIVED NOV 26 1985

Sample Number  
ED802Organics Analysis Data Sheet  
(Page 2)

## Semivolatile Compounds

Concentration: Low Medium (Circle One)Date Extracted/Prepared: 7/1/85 FEN BAN 608Date Analyzed: 7/5/85 1221Conc/Dil Factor: 1

CAS Number		ug/l or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20 U
108-95-2	Phenol	20 U
62-53-3	Aniline	20 U
111-64-4	bis(2-Chloroethyl)Ether	20 U
95-57-8	2-Chlorophenol	20 U
541-73-1	1, 3-Dichlorobenzene	20 U
106-46-7	1, 4-Dichlorobenzene	20 U
100-51-6	Benzyl Alcohol	20 U
95-50-1	1, 2-Dichlorobenzene	20 U
95-48-7	2-Methylphenol	20 U
39638-32-9	bis(2-chloroisopropyl)Ether	20 U
106-44-5	4-Methylphenol	20 U
621-64-7	N-Nitroso-Di-n-Propylamine	20 U
67-72-1	Hexachloroethane	20 U
98-95-3	Nitrobenzene	20 U
78-59-1	Isophorone	<u>0.83</u> 20 U
88-75-5	2-Nitrophenol	20 U
105-67-9	2, 4-Dimethylphenol	20 U
65-85-0	Benzoic Acid	100 U
111-91-1	bis(2-Chloroethoxy)Methane	20 U
120-83-2	2, 4-Dichlorophenol	20 U
120-82-1	1, 2, 4-Trichlorobenzene	20 U
91-20-3	Naphthalene	20 U
106-47-8	4-Chloroaniline	20 U
87-68-3	Hexachlorobutadiene	20 U
59-50-7	4-Chloro-3-Methylphenol	20 U
91-57-6	2-Methylnaphthalene	20 U
77-47-4	Hexachlorocyclopentadiene	20 U
88-06-2	2, 4, 6-Trichlorophenol	20 U
95-95-4	2, 4, 5-Trichlorophenol	100 U
91-58-7	2-Chloronaphthalene	20 U
88-74-4	2-Nitroaniline	100 U
131-11-3	Dimethyl Phthalate	20 U
208-96-8	Acenaphthylene	20 U
99-09-2	3-Nitroaniline	100 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	20 U
51-28-5	2, 4-Dinitrophenol	100 U
100-02-7	4-Nitrophenol	100 U
132-64-9	Dibenzofuran	20 U
121-14-2	2, 4-Dinitrotoluene	20 U
606-20-2	2, 6-Dinitrotoluene	20 U
84-66-2	Diethylphthalate	<u>0.78</u> 20 U
7005-72-3	4-Chlorophenyl-phenylether	20 U
86-73-7	Fluorene	20 U
100-01-6	4-Nitroaniline	100 U
534-52-1	4, 6-Dinitro-2-Methylphenol	100 U
86-30-6	N-Nitrosodiphenylamine (1)	20 U
101-55-3	4-Bromophenyl-phenylether	20 U
118-74-1	Hexachlorobenzene	20 U
87-86-5	Pentachlorophenol	100 U
85-01-8	Phenanthrene	20 U
120-12-7	Anthracene	20 U
84-74-2	Di-n-Butylphthalate	<u>2.9</u> 20 U
206-44-0	Fluoranthene	20 U
92-87-5	Benzdine	100 U
129-00-0	Pyrene	20 U
85-68-7	Butylbenzylphthalate	20 U
31-94-1	3, 3'-Dichlorobenzidine	20 U
56-55-3	Benzo(a)Anthracene	20 U
117-81-7	bis(2-Ethylhexyl)Phthalate	<u>11.8</u> 20 U
218-01-9	Chrysene	20 U
117-84-0	Di-n-Octyl Phthalate	20 U
205-99-2	Benzo(b)Fluoranthene	20 U
207-08-9	Benzo(k)Fluoranthene	20 U
50-32-8	Benzo(a)Pyrene	20 U
193-39-5	Indeno(1, 2, 3-cd)Pyrene	20 U
53-70-3	Dibenz(a, h)Anthracene	20 U
191-24-2	Benzo(g, h, i)Perylene	20 U

(1)-Cannot be separated from diphenylamine

Organics Analysis Data Sheet  
(Page 3)

Pesticide/PCEs

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 7-1-85  
 Date Analyzed: 7-12-85  
 Conc/Dil Factor: 1

CAS Number		<u>ug/L</u> or ug/Kg (Circle One)
319-84-6	Alpha-BHC	0.05 U
318-85-7	Beta-BHC	0.05 U
319-86-8	Delta-BHC	0.05 U
58-89-9	Gamma-BHC (Lindane)	0.05 U
76-44-8	Heptachlor	0.05 U
309-00-2	Aldrin	0.05 U
1024-57-3	Heptachlor Epoxide	0.05 U
358-98-8	Endosulfan I	0.05 U
50-57-1	Dieldrin	0.10 U
72-55-9	4, 4'-DDE	0.10 U
72-20-8	Endrin	0.10 U
33213-65-9	Endosulfan II	0.10 U
72-54-8	4, 4'-DDD	0.10 U
7421-93-4	Endrin Aldehyde	0.10 U
1031-07-8	Endosulfan Sulfate	0.10 U
50-29-3	4, 4'-DDT	0.10 U
72-43-5	Methoxychlor	0.50 U
53494-70-5	Endrin Ketone	0.10 U
57-74-9	Chlordane	0.50 U
8001-35-2	Toxaphene	1.0 U
12674-11-2	Aroclor-1016	0.50 U
11104-28-2	Aroclor-1221	0.50 U
11141-16-5	Aroclor-1232	0.50 U
53469-21-9	Aroclor-1242	0.50 U
12672-29-6	Aroclor-1248	0.50 U
11097-69-1	Aroclor-1254	1.0 U
11096-82-5	Aroclor-1260	1.0 U

$V_i$  = Volume of extract injected (ul)  
 $V_s$  = Volume of water extracted (ml)  
 $W_s$  = Weight of sample extracted (g)  
 $V_t$  = Volume of total extract (ul)

$V_s$  1000 or  $W_s$  —  $V_t$  10,000  $V_i$  4.0

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	Unknown	BNA	845	210
2.	Unknown	BNA	1570	43
3.	5353-25-3 Ethanol, 2-(9-octadecenyloxy)-, (Z)-	BNA	1745	46
4.	Unknown	BNA	1926	53
5.	Unknown	BNA	1934	46
6.	Unknown	BNA	1994	47
7.	Unknown	BNA	2069	31
8.	Unknown	BNA	2090	50
9.	Unknown	BNA	2097	58
10.	Unknown	BNA	2148	92
11.	Unknown Alkane C <sub>25</sub> -C <sub>26</sub>	BNA	2179	42
12.	Unknown	BNA	2250	43
13.	Unknown	BNA	2256	96
14.	Unknown	BNA	2261	30
15.	Unknown	BNA	2271	32
16.	Unknown	BNA	2321	32
17.	Unknown	BNA	2393	102
18.	Unknown	BNA	2424	50
19.	Unknown Alkane C <sub>14</sub> -C <sub>20</sub>	BNA	2451	63
20.	Unknown	BNA	2486	87
21.	Unknown	BNA	2592	20
22.				
23.	- None Found	VOA	-	-
24.				
25.				
26.				
27.				
28.				
29.				
30.				

Organics Analysis Data Sheet  
 (Page 1)

Laboratory Name: HAZLETON LABORATORIES  
 Lab Sample ID No: 50606607  
 Sample Matrix: Water  
 Data Release Authorized By: David C. Hill

Case No: 4571  
 QC Report No: \_\_\_\_\_  
 Contract No: 68-01-6961  
 Date Sample Received: 6-28-85

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 7-3-85 FRN28293  
 Date Analyzed: 7-3-85 1440  
 Conc/Dil Factor: 5 pH \_\_\_\_\_  
 Percent Moisture: \_\_\_\_\_  
 Percent Moisture (Decanted): \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	50u
74-83-9	Bromomethane	50u
75-01-4	Vinyl Chloride	(110)
75-00-3	Chloroethane	50u
75-09-2	Methylene Chloride	25u
67-64-1	Acetone	50u
75-15-0	Carbon Disulfide	25u
75-35-4	1, 1-Dichloroethene	25u
75-34-3	1, 1-Dichloroethane	25u
156-60-5	Trans-1, 2-Dichloroethene	(970)
67-66-3	Chloroform	25u
107-06-2	1, 2-Dichloroethane	25u
78-93-3	2-Butanone	50u
71-55-6	1, 1, 1-Trichloroethane	25u
56-23-5	Carbon Tetrachloride	25u
108-05-4	Vinyl Acetate	50u
75-27-4	Bromodichloromethane	25u

CAS Number		ug/l or ug/Kg (Circle One)
79-34-5	1, 1, 2, 2-Tetrachloroethane	25u
78-87-5	1, 2-Dichloropropane	25u
10061-02-6	Trans-1, 3-Dichloropropene	25u
79-01-6	Trichloroethene	(22 J)
124-48-1	Dibromochloromethane	25u
79-00-5	1, 1, 2-Trichloroethane	25u
71-43-2	Benzene	(9.0 J)
10061-01-5	cis-1, 3-Dichloropropene	25u
110-75-8	2-Chloroethylvinylether	50u
75-25-2	Bromoform	25u
591-78-6	2-Hexanone	50u
108-10-1	4-Methyl-2-Pentanone	50u
127-18-4	Tetrachloroethane	25u
108-88-3	Toluene	(10 J)
108-90-7	Chlorobenzene	25u
100-41-4	Ethylbenzene	25u
100-42-5	Styrene	25u
	Total Xylenes	(37)

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

**Value** If the result is a value greater than or equal to the detection limit, report the value.

**U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution actions. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

**J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10J)

**C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/ul in the final extract should be confirmed by GC/MS.

**B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

**Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Organics Analysis Data Sheet  
(Page 2)

## Semivolatile Compounds

Concentration: Low Medium (Circle One)Date Extracted/Prepared: 7/1/85 FRN BAN598Date Analyzed: 7/2/85 2257Conc/Dil Factor: 1

CAS Number		ug/l or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20 U
108-95-2	Phenol	20 U
62-53-3	Aniline	20 U
111-64-4	bis(2-Chloroethyl)Ether	20 U
95-57-8	2-Chlorophenol	20 U
541-73-1	1,3-Dichlorobenzene	20 U
106-46-7	1,4-Dichlorobenzene	20 U
100-51-6	Benzyl Alcohol	20 U
95-50-1	1,2-Dichlorobenzene	20 U
95-48-7	2-Methylphenol	20 U
39638-32-9	bis(2-chloroisopropyl)Ether	20 U
106-44-5	4-Methylphenol	20 U
621-64-7	N-Nitroso-Di-n-Propylamine	20 U
67-72-1	Hexachloroethane	20 U
98-95-3	Nitrobenzene	20 U
78-59-1	Isophorone	20 U
88-75-5	2-Nitrophenol	20 U
105-67-9	2,4-Dimethylphenol	20 U
65-85-0	Benzoic Acid	100 U
111-91-1	bis(2-Chloroethoxy)Methane	20 U
120-83-2	2,4-Dichlorophenol	20 U
120-82-1	1,2,4-Trichlorobenzene	20 U
91-20-3	Naphthalene	20 U
105-47-8	4-Chloroaniline	20 U
87-68-3	Hexachlorobutadiene	20 U
59-50-7	4-Chloro-3-Methylphenol	20 U
91-57-6	2-Methylnaphthalene	20 U
77-47-4	Hexachlorocyclopentadiene	20 U
88-06-2	2,4,6-Trichlorophenol	20 U
95-95-4	2,4,5-Trichlorophenol	100 U
91-58-7	2-Chloronaphthalene	20 U
88-74-4	2-Nitroaniline	100 U
131-11-3	Dimethyl Phthalate	20 U
208-96-8	Acenaphthylene	20 U
99-09-2	3-Nitroaniline	100 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	20 U
51-28-5	2,4-Dinitrophenol	100 U
100-02-7	4-Nitrophenol	100 U
132-64-9	Dibenzofuran	20 U
121-14-2	2,4-Dinitrotoluene	20 U
606-20-2	2,6-Dinitrotoluene	20 U
84-66-2	Diethylphthalate	20 U
7005-72-3	4-Chlorophenyl-phenylether	20 U
96-73-7	Fluorene	20 U
100-01-6	4-Nitroaniline	100 U
534-52-1	4,6-Dinitro-2-Methylphenol	100 U
96-30-6	N-Nitrosodiphenylamine (1)	20 U
101-55-3	4-Bromophenyl-phenylether	20 U
118-74-1	Hexachlorobenzene	20 U
87-86-5	Pentachlorophenol	100 U
85-01-8	Phenanthrene	20 U
120-12-7	Anthracene	20 U
94-74-2	Di-n-Butylphthalate	<u>1.45</u> 20 U
206-44-0	Fluoranthene	20 U
92-87-5	Benzidine	100 U
129-00-0	Pyrene	20 U
85-58-7	Butylbenzylphthalate	20 U
91-94-1	3,3'-Dichlorobenzidine	20 U
56-55-3	Benzo(a)Anthracene	20 U
117-81-7	bis(2-Ethylhexyl)Phthalate	<u>15.85</u> 20 U
218-01-9	Chrysene	20 U
117-84-0	Di-n-Octyl Phthalate	<u>13.55</u> 20 U
205-99-2	Benzo(b)Fluoranthene	20 U
207-08-9	Benzo(k)Fluoranthene	20 U
50-32-8	Benzo(a)Pyrene	20 U
193-39-5	Indeno(1,2,3-cd)Pyrene	20 U
53-70-3	Dibenz(a,h)Anthracene	20 U
191-24-2	Benzo(c,h,i)Perylene	20 U

(1)-Cannot be separated from diphenylamine

No: 4571

ED 803

Organics Analysis Data Sheet  
(Page 3)

## Pesticide/PCBs

Concentration: Low Medium (Circle One)Date Extracted/Prepared: 7-1-85Date Analyzed: 7-12-85Conc/Dil Factor: 1

CAS Number		<u>ug/g</u> or ug/Kg (Circle One)
319-64-6	Alpha-BHC	0.05 U
318-85-7	Beta-BHC	0.05 U
318-86-8	Delta-BHC	0.05 U
58-89-9	Gamma-BHC (Lindane)	0.05 U
78-44-8	Heptachlor	0.05 U
309-00-2	Aldrin	0.05 U
1024-57-3	Heptachlor Epoxide	0.05 U
259-98-8	Endosulfan I	0.05 U
50-57-1	Dieldrin	0.10 U
72-55-9	4, 4'-DDE	0.10 U
72-20-8	Endrin	0.10 U
33213-65-9	Endosulfan II	0.10 U
72-54-8	4, 4'-DDD	0.10 U
7421-93-4	Endrin Aldehyde	0.10 U
1031-07-8	Endosulfan Sulfate	0.10 U
50-29-3	4, 4'-DDT	0.10 U
72-43-5	Methoxychlor	0.50 U
53494-70-5	Endrin Ketone	0.10 U
57-74-9	Chlordane	0.50 U
8001-35-2	Toxaphene	1.0 U
12674-11-2	Aroclor-1016	0.50 U
11104-28-2	Aroclor-1221	0.50 U
11141-16-5	Aroclor-1232	0.50 U
53469-21-9	Aroclor-1242	0.50 U
12672-29-6	Aroclor-1248	0.50 U
11097-69-1	Aroclor-1254	1.0 U
11096-82-5	Aroclor-1260	1.0 U

 $V_i$  = Volume of extract injected (ul) $V_s$  = Volume of water extracted (ml) $W_s$  = Weight of sample extracted (g) $V_t$  = Volume of total extract (ul) $V_s$  1000 or  $W_s$  —  $V_i$  10,000  $V_t$  4.0

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/Tar ug/kg)
1. 95-47-6	Benzene, 1,2-dimethyl	BNA	275	8.3
2.	Unknown	BNA	821	190
3.	Unknown	BNA	1573	13
4. 5675-51-4	1,2 Dodecanediol	BNA	1749	12
5.	Unknown	BNA	1930	13
6.	Unknown	BNA	1937	8.3
7.	Unknown	BNA	1996	20
8.	Unknown	BNA	2072	11
9.	Unknown	BNA	2151	9.6
10.	Unknown	BNA	2270	12
11.	Unknown	BNA	2395	15
12.				
13. 60-29-7	Ethane, 1,1'-Oxybis	VOA	169	59
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				



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ED 804

Organics Analysis Data Sheet  
 (Page 1)

25F H 10518

Laboratory Name: HAZLETON LABORATORIES

Case No: 4571

Lab Sample ID No: 50606608

QC Report No: \_\_\_\_\_

Sample Matrix: WATER

Contract No: 68-01-6961

Data Release Authorized By: David C. Kelly

Date Sample Received: 6-28-85

Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7-3-85 FRN 2829

Date Analyzed: 7-3-85 1830

Conc/Dil Factor: 1 pH \_\_\_\_\_

Percent Moisture: \_\_\_\_\_

Percent Moisture (Decanted): \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	10 U
74-83-9	Bromomethane	10 U
75-01-4	Vinyl Chloride	10 U
75-00-3	Chloroethane	10 U
75-09-2	Methylene Chloride	5 U
67-64-1	Acetone	10 U
75-15-0	Carbon Disulfide	5 U
75-35-4	1, 1-Dichloroethane	5 U
75-34-3	1, 1-Dichloroethane	5 U
156-60-5	Trans-1, 2-Dichloroethane	5 U
67-66-3	Chloroform	5 U
107-06-2	1, 2-Dichloroethane	5 U
78-93-3	2-Butanone	<u>(20)</u> <del>10 U</del>
71-55-6	1, 1, 1-Trichloroethane	5 U
56-23-5	Carbon Tetrachloride	5 U
108-05-4	Vinyl Acetate	10 U
75-27-4	Bromodichloromethane	5 U

CAS Number		ug/l or ug/Kg (Circle One)
79-34-5	1, 1, 2, 2-Tetrachloroethane	5 U
78-87-5	1, 2-Dichloropropane	5 U
10061-02-6	Trans-1, 3-Dichloropropene	5 U
79-01-6	Trichloroethene	5 U
124-48-1	Dibromochloromethane	5 U
79-00-5	1, 1, 2-Trichloroethane	5 U
71-43-2	Benzene	5 U
10061-01-5	cis-1, 3-Dichloropropene	5 U
110-75-8	2-Chloroethylvinylether	10 U
75-25-2	Bromoform	5 U
591-78-6	2-Hexanone	10 U
108-10-1	4-Methyl-2-Pentanone	10 U
127-18-4	Tetrachloroethene	5 U
108-88-3	Toluene	<u>(1.05)</u> <del>5 U</del>
108-90-7	Chlorobenzene	5 U
100-41-4	Ethylbenzene	5 U
100-42-5	Styrene	5 U
	Total Xylenes	5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

**Value** If the result is a value greater than or equal to the detection limit, report the value.

**U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution actions. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum obtainable detection limit for the sample.

**J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10U)

**C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/ul in the final extract should be confirmed by GC/MS.

**B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

**Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Organics Analysis Data Sheet  
(Page 2)

## Semivolatile Compounds

Concentration: Low Medium (Circle One)Date Extracted/Prepared: 7/1/85 FRN BAN597Date Analyzed: 7/2/85 2200Conc/Dil Factor: 1

CAS Number		ug/l or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20 U
108-95-2	Phenol	20 U
62-53-3	Aniline	20 U
111-44-4	bis(2-Chloroethyl)Ether	20 U
95-57-8	2-Chlorophenol	20 U
541-73-1	1, 3-Dichlorobenzene	20 U
105-46-7	1, 4-Dichlorobenzene	20 U
100-51-6	Benzyl Alcohol	20 U
95-50-1	1, 2-Dichlorobenzene	20 U
95-48-7	2-Methylphenol	20 U
39638-32-9	bis(2-chloroisopropyl)Ether	20 U
106-44-5	4-Methylphenol	20 U
621-64-7	N-Nitroso-Di-n-Propylamine	20 U
67-72-1	Hexachloroethane	20 U
98-95-3	Nitrobenzene	20 U
78-59-1	Isophorone	20 U
88-75-5	2-Nitrophenol	20 U
105-67-9	2, 4-Dimethylphenol	20 U
65-85-0	Benzoic Acid	100 U
111-91-1	bis(2-Chloroethoxy)Methane	20 U
120-83-2	2, 4-Dichlorophenol	20 U
120-82-1	1, 2, 4-Trichlorobenzene	20 U
91-20-3	Naphthalene	20 U
106-47-8	4-Chloroaniline	20 U
87-68-3	Hexachlorobutadiene	20 U
59-50-7	4-Chloro-3-Methylphenol	20 U
91-57-6	2-Methylnaphthalene	20 U
77-47-4	Hexachlorocyclopentadiene	20 U
88-06-2	2, 4, 6-Trichlorophenol	20 U
95-95-4	2, 4, 5-Trichlorophenol	100 U
91-58-7	2-Chloronaphthalene	20 U
88-74-4	2-Nitroaniline	100 U
131-11-3	Dimethyl Phthalate	20 U
208-96-8	Acenaphthylene	20 U
99-09-2	3-Nitroaniline	100 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	20 U
51-28-5	2, 4-Dinitrophenol	100 U
100-02-7	4-Nitrophenol	100 U
132-64-9	Dibenzofuran	20 U
121-14-2	2, 4-Dinitrotoluene	20 U
606-20-2	2, 6-Dinitrotoluene	20 U
84-55-2	Diethylphthalate	20 U
7005-72-3	4-Chlorophenyl-phenylether	20 U
86-73-7	Fluorene	20 U
100-01-6	4-Nitroaniline	100 U
534-52-1	4, 6-Dinitro-2-Methylphenol	100 U
86-30-6	N-Nitrosodiphenylamine (1)	20 U
101-55-3	4-Bromophenyl-phenylether	20 U
118-74-1	Hexachlorobenzene	20 U
87-86-5	Pentachlorophenol	100 U
85-01-8	Phenanthrene	20 U
120-12-7	Anthracene	20 U
84-74-2	Di-n-Butylphthalate	1.25 <del>20 U</del>
206-44-0	Fluoranthene	20 U
92-87-5	Benzidine	100 U
129-00-0	Pyrene	20 U
85-68-7	Butylbenzylphthalate	20 U
91-94-1	3, 3'-Dichlorobenzidine	20 U
56-55-3	Benzo(a)Anthracene	20 U
117-81-7	bis(2-Ethylhexyl)Phthalate	6.4 <del>20 U</del>
218-01-9	Chrysene	20 U
117-84-0	Di-n-Octyl Phthalate	20 U
205-99-2	Benzo(b)Fluoranthene	20 U
207-08-9	Benzo(k)Fluoranthene	20 U
50-32-8	Benzo(e)Pyrene	20 U
193-39-5	Indeno(1, 2, 3-cd)Pyrene	20 U
53-70-3	Dibenz(a, h)Anthracene	20 U
191-24-2	Benzo(g, h, i)Perylene	20 U

(1)-Cannot be separated from diphenylamine

No: 4571

Sample Number  
ED 804

Organics Analysis Data Sheet  
(Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7-1-85

Date Analyzed: 7-12-85

Conc/Dil Factor: 1

CAS Number		<u>ug/L</u> or ug/kg (Circle One)
319-64-6	Alpha-BHC	0.05 U
319-85-7	Beta-BHC	0.05 U
319-86-8	Delta-BHC	0.05 U
58-89-9	Gamma-BHC (Lindane)	0.05 U
76-44-8	Heptachlor	0.05 U
309-00-2	Aldrin	0.05 U
1024-57-3	Heptachlor Epoxide	0.05 U
959-98-8	Endosulfan I	0.05 U
30-57-1	Dieldrin	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	Endrin	0.10 U
33213-85-9	Endosulfan II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	Endrin Aldehyde	0.10 U
1031-07-8	Endosulfan Sulfate	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	Methoxychlor	0.50 U
53494-70-5	Endrin Ketone	0.10 U
57-74-9	Chlordane	0.50 U
8001-35-2	Toxaphene	1.0 U
12674-11-2	Aroclor-1016	0.50 U
11104-28-2	Aroclor-1221	0.50 U
11141-18-5	Aroclor-1232	0.50 U
53469-21-9	Aroclor-1242	0.50 U
12672-29-6	Aroclor-1248	0.50 U
11097-69-1	Aroclor-1254	1.0 U
11096-82-5	Aroclor-1260	1.0 U

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  1000 or  $W_s$  —  $V_i$  10,000  $V_t$  4.0

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	Unknown	BNA	837	460
2.	Unknown Alkene	BNA	1568	19
3.	Unknown	BNA	1760	8.4
4.	Unknown	BNA	1937	14
5.	Unknown	BNA	2072	12
6.	Unknown Alkene C <sub>13</sub> → C <sub>43</sub>	BNA	2181	14
7.	Unknown Alkene C <sub>30</sub> → C <sub>40</sub>	BNA	2210	11
8.	Unknown	BNA	2259	24
9.				
10.	None Found	VOA	-	-
11.				
12.				
13.				
14.				
15.				
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ED 805  
 85FH105M

Organics Analysis Data Sheet  
 (Page 1)

Laboratory Name: HAZLETON LABORATORIES  
 Lab Sample ID No: 50606609  
 Sample Matrix: WATER  
 Data Release Authorized By: David C. Will

Case No: 4571  
 QC Report No: \_\_\_\_\_  
 Contract No: 68-01-6461  
 Date Sample Received: 6-28-85

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 7-3-85 FRN 28299  
 Date Analyzed: 7-3-85 1920  
 Conc/Dil Factor: 1 pH \_\_\_\_\_  
 Percent Moisture: \_\_\_\_\_  
 Percent Moisture (Decanted): \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	10 U
74-83-9	Bromomethane	10 U
75-01-4	Vinyl Chloride	10 U
75-00-3	Chloroethane	10 U
75-09-2	Methylene Chloride	5 U
67-64-1	Acetone	10 U
75-15-0	Carbon Disulfide	5 U
75-35-4	1, 1-Dichloroethene	5 U
75-34-3	1, 1-Dichloroethane	5 U
156-80-5	Trans-1, 2-Dichloroethene	5 U
67-66-3	Chloroform	5 U
107-06-2	1, 2-Dichloroethane	5 U
78-93-3	2-Butanone	10 U
71-55-6	1, 1, 1-Trichloroethane	5 U
56-23-5	Carbon Tetrachloride	5 U
103-05-4	Vinyl Acetate	10 U
75-27-4	Bromodichloromethane	5 U

CAS Number		ug/l or ug/Kg (Circle One)
79-34-5	1, 1, 2, 2-Tetrachloroethane	5 U
78-87-5	1, 2-Dichloropropane	5 U
10061-02-6	Trans-1, 3-Dichloropropene	5 U
79-01-6	Trichloroethene	5 U
124-48-1	Dibromochloromethane	5 U
79-00-5	1, 1, 2-Trichloroethane	5 U
71-43-2	Benzene	5 U
10061-01-5	cis-1, 3-Dichloropropene	5 U
110-75-8	2-Chloroethylvinylether	10 U
75-25-2	Bromoform	5 U
591-78-6	2-Hexanone	10 U
108-10-1	4-Methyl-2-Pentanone	10 U
127-18-4	Tetrachloroethene	5 U
108-88-3	Toluene	5 U
108-90-7	Chlorobenzene	5 U
100-41-4	Ethylbenzene	5 U
100-42-5	Styrene	5 U
	Total Xylenes	5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

- V** Value: If the result is a value greater than or equal to the detection limit, report the value.
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution factors. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10U)

- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng/ul in the final extract should be confirmed by GC/MS
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7/1/85 FRN BAN609

Date Analyzed: 7/5/85 1319

Conc/Dil Factor: 1

CAS Number		ug/l or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20 U
108-95-2	Phenol	20 U
62-53-3	Aniline	20 U
111-44-4	bis(2-Chloroethyl)Ether	20 U
95-57-8	2-Chlorophenol	20 U
541-73-1	1, 3-Dichlorobenzene	20 U
106-46-7	1, 4-Dichlorobenzene	20 U
100-51-6	Benzyl Alcohol	20 U
95-50-1	1, 2-Dichlorobenzene	20 U
95-48-7	2-Methylphenol	20 U
39638-32-9	bis(2-chloroisopropyl)Ether	20 U
106-44-5	4-Methylphenol	20 U
621-64-7	N-Nitroso-Di-n-Propylamine	20 U
67-72-1	Hexachloroethane	20 U
98-95-3	Nitrobenzene	20 U
78-59-1	Isophorone	<del>20 U</del>
88-75-5	2-Nitrophenol	20 U
105-67-9	2, 4-Dimethylphenol	20 U
65-85-0	Benzoic Acid	100 U
111-91-1	bis(2-Chloroethoxy)Methane	20 U
120-83-2	2, 4-Dichlorophenol	20 U
120-82-1	1, 2, 4-Trichlorobenzene	20 U
91-20-3	Naphthalene	20 U
106-47-8	4-Chloroaniline	20 U
87-68-3	Hexachlorobutadiene	20 U
59-50-7	4-Chloro-3-Methylphenol	20 U
91-57-6	2-Methylnaphthalene	20 U
77-47-4	Hexachlorocyclopentadiene	20 U
88-06-2	2, 4, 6-Trichlorophenol	20 U
95-95-4	2, 4, 5-Trichlorophenol	100 U
91-58-7	2-Chloronaphthalene	20 U
88-74-4	2-Nitroaniline	100 U
131-11-3	Dimethyl Phthalate	20 U
208-96-8	Acenaphthylene	20 U
99-09-2	3-Nitroaniline	100 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	20 U
51-28-5	2, 4-Dinitrophenol	100 U
100-02-7	4-Nitrophenol	100 U
132-64-9	Dibenzofuran	20 U
121-14-2	2, 4-Dinitrotoluene	20 U
606-20-2	2, 6-Dinitrotoluene	20 U
84-66-2	Diethylphthalate	<u>22 J</u>
7005-72-3	4-Chlorophenyl-phenylether	20 U
86-73-7	Fluorene	20 U
100-01-6	4-Nitroaniline	100 U
534-52-1	4, 6-Dinitro-2-Methylphenol	100 U
86-30-6	N-Nitrosodiphenylamine (1)	20 U
101-55-3	4-Bromophenyl-phenylether	20 U
118-74-1	Hexachlorobenzene	20 U
87-86-5	Pentachlorophenol	100 U
85-01-8	Phenanthrene	20 U
120-12-7	Anthracene	20 U
84-74-2	Di-n-Butylphthalate	20 U
206-44-0	Fluoranthene	20 U
92-87-5	Benzdine	100 U
129-00-0	Pyrene	20 U
85-68-7	Butylbenzylphthalate	20 U
91-94-1	3, 3'-Dichlorobenzidine	20 U
56-55-3	Benzo(a)Anthracene	20 U
117-81-7	bis(2-Ethylhexyl)Phthalate	<u>3.5 BJ</u>
218-01-9	Chrysene	20 U
117-84-0	Di-n-Octyl Phthalate	20 U
205-99-2	Benzo(b)Fluoranthene	20 U
207-08-9	Benzo(k)Fluoranthene	20 U
50-32-8	Benzo(a)Pyrene	20 U
193-39-5	Indeno(1, 2, 3-cd)Pyrene	20 U
53-70-3	Dibenzo(a, h)Anthracene	20 U
191-24-2	Benzo(g, h, i)Perylene	20 U

(1)-Cannot be separated from diphenylamine

Organics Analysis Data Sheet  
(Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7-1-85

Date Analyzed: 7-13-85

Conc/Dil Factor: 1

CAS Number		<u>ug/L or ug/Kg</u> (Circle One)
319-84-6	Alpha-BHC	0.05 U
319-85-7	Beta-BHC	0.05 U
319-86-8	Delta-BHC	0.05 U
58-89-9	Gamma-BHC (Lindane)	0.05 U
76-44-8	Heptachlor	0.05 U
309-00-2	Aldrin	0.05 U
1024-57-3	Heptachlor Epoxide	0.05 U
959-98-8	Endosulfan I	0.05 U
50-57-1	Dieldrin	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	Endrin	0.10 U
33213-65-9	Endosulfan II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	Endrin Aldehyde	0.10 U
1031-07-8	Endosulfan Sulfate	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	Methoxychlor	0.50 U
53494-70-5	Endrin Ketone	0.10 U
57-74-9	Chlordane	0.50 U
8001-35-2	Toxaphene	1.0 U
12674-11-2	Aroclor-1016	0.50 U
11104-28-2	Aroclor-1221	0.50 U
11141-16-5	Aroclor-1232	0.50 U
53469-21-9	Aroclor-1242	0.50 U
12672-29-6	Aroclor-1248	0.50 U
11097-69-1	Aroclor-1254	1.0 U
11096-82-5	Aroclor-1260	1.0 U

- $V_i$  = Volume of extract injected (ul)
- $V_s$  = Volume of water extracted (ml)
- $W_s$  = Weight of sample extracted (g)
- $V_t$  = Volume of total extract (ul)

$V_s$  1000 or  $W_s$  —  $V_i$  10,000  $V_t$  4.0

Organics Analysis Data Sheet  
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	Unknown	BNA	508	12
2. 617-94-7	Benzene methanol, .alpha., .alpha., -Dimethyl-	BNA	592	9.0
3.	Unknown	BNA	809	13
4.	Unknown	BNA	841	170
5. 55724-73-7	Butanoic Acid, 4-Butoxy -	BNA	965	45
6. 98-73-7	Benzoic Acid, 4(1,1)-dimethylethyl)-	BNA	1077	10
7. 80-59-7	Benzene sulfonamide, N-Ethyl-4-methyl	BNA	1230	16
8. 1421-49-4	Benzoic Acid, 3,5-bis(1,1-Dimethylethyl)-4-hydroxy	BNA	1443	8.8
9.	Unknown	BNA	2069	14
10.	Unknown	BNA	2193	12
11.	Unknown	BNA	2199	24
12.	Unknown	BNA	2235	109
13.	Unknown	BNA	2254	9.6
14.	Unknown	BNA	2302	42
15.	Unknown	BNA	2412	15
16.				
17. 60-29-7	Ethane, 1,1'-oxybis	VOA	169	6.0
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Sample Number  
**ED 806**

Organics Analysis Data Sheet  
 (Page 1)

85FH10017

Laboratory Name: HAZLETON LABORATORIES  
 Lab Sample ID No: 50006610  
 Sample Matrix: WATER  
 Data Release Authorized By: David C. [Signature]

Case No: 4571  
 QC Report No: \_\_\_\_\_  
 Contract No: 68-01-6961  
 Date Sample Received: 6-28-85

Volatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 7-3-85 FRN 28301  
 Date Analyzed: 7-8-85 2100  
 Gen Dil Factor: 5 pH \_\_\_\_\_  
 Percent Moisture: \_\_\_\_\_  
 Percent Moisture (Decanted): \_\_\_\_\_

Duplicate  
 of ED803

CAS Number	Compound	(ug/l) or ug/Kg (Circle One)
74-87-3	Chloromethane	50u
74-83-9	Bromomethane	50u
75-01-4	Vinyl Chloride	(82)
75-00-3	Chloroethane	50u
75-09-2	Methylene Chloride	(27 B)
67-64-1	Acetone	50u
75-15-0	Carbon Disulfide	25u
75-35-4	1, 1-Dichloroethane	25u
75-34-3	1, 1-Dichloroethane	25u
158-60-5	Trans-1, 2-Dichloroethane	(740)
67-86-3	Chloroform	25u
107-08-2	1, 2-Dichloroethane	25u
78-93-3	2-Butenone	50u
71-55-6	1, 1, 1-Trichloroethane	25u
56-23-5	Carbon Tetrachloride	25u
108-05-4	Vinyl Acetate	50u
75-27-4	Bromodichloromethane	25u

CAS Number	Compound	(ug/l) or ug/Kg (Circle One)
79-34-5	1, 1, 2, 2-Tetrachloroethane	25u
78-87-5	1, 2-Dichloropropane	25u
10061-02-6	Trans-1, 3-Dichloropropene	25u
79-01-6	Trichloroethene	25u
124-48-1	Dibromochloromethane	25u
79-00-5	1, 1, 2-Trichloroethane	25u
71-43-2	Benzene	25u
10061-01-5	cis-1, 3-Dichloropropene	25u
110-75-8	2-Chloroethylvinylether	50u
75-25-2	Bromoform	25u
591-78-6	2-Hexanone	50u
108-10-1	4-Methyl-2-Pentanone	50u
127-18-4	Tetrachloroethane	25u
108-88-3	Toluene	(7.5 J)
108-90-7	Chlorobenzene	25u
100-41-4	Ethylbenzene	25u
100-42-5	Styrene	25u
	Total Xylenes	(28)

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

**Value** If the result is a value greater than or equal to the detection limit, report the value.

**U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution actions. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

**J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero (e.g., 10U).

**C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/ul in the final extract should be confirmed by GC/MS.

**B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

**Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Organics Analysis Data Sheet  
 (Page 2)

RECEIVED NOV 26 1985

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 7/1/85 FRN BAN595  
 Date Analyzed: 7/2/85 2004  
Conc/Dil Factor 1

CAS Number		ug/l or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20 U
108-95-2	Phenol	20 U
62-53-3	Aniline	20 U
111-44-4	bis(2-Chloroethyl)Ether	20 U
95-57-8	2-Chlorophenol	20 U
541-73-1	1, 3-Dichlorobenzene	20 U
106-46-7	1, 4-Dichlorobenzene	20 U
100-51-6	Benzyl Alcohol	20 U
95-50-1	1, 2-Dichlorobenzene	20 U
95-48-7	2-Methylphenol	20 U
39638-32-9	bis(2-chloroisopropyl)Ether	20 U
106-44-5	4-Methylphenol	20 U
621-64-7	N-Nitroso-Di-n-Propylamine	20 U
67-72-1	Hexachloroethane	20 U
98-95-3	Nitrobenzene	20 U
78-59-1	Isophorone	20 U
88-75-5	2-Nitrophenol	20 U
105-67-9	2, 4-Dimethylphenol	20 U
65-85-0	Benzoic Acid	100 U
111-91-1	bis(2-Chloroethoxy)Methane	20 U
120-83-2	2, 4-Dichlorophenol	20 U
120-82-1	1, 2, 4-Trichlorobenzene	20 U
91-20-3	Naphthalene	20 U
105-47-8	4-Chloroaniline	20 U
87-68-3	Hexachlorobutadiene	20 U
59-50-7	4-Chloro-3-Methylphenol	20 U
91-57-6	2-Methylnaphthalene	20 U
77-47-4	Hexachlorocyclopentadiene	20 U
88-06-2	2, 4, 6-Trichlorophenol	20 U
95-95-4	2, 4, 5-Trichlorophenol	100 U
91-58-7	2-Chloronaphthalene	20 U
88-74-4	2-Nitroaniline	100 U
131-11-3	Dimethyl Phthalate	20 U
208-96-8	Acenaphthylene	20 U
99-09-2	3-Nitroaniline	100 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	20 U
51-28-5	2, 4-Dinitrophenol	100 U
100-02-7	4-Nitrophenol	100 U
132-64-9	Dibenzofuran	20 U
121-14-2	2, 4-Dinitrotoluene	20 U
606-20-2	2, 6-Dinitrotoluene	20 U
94-66-2	Diethylphthalate	20 U
7005-72-3	4-Chlorophenyl-phenylether	20 U
86-73-7	Fluorene	20 U
100-01-6	4-Nitroaniline	100 U
534-52-1	4, 6-Dinitro-2-Methylphenol	100 U
86-30-6	N-Nitrosodiphenylamine (1)	20 U
101-55-3	4-Bromophenyl-phenylether	20 U
118-74-1	Hexachlorobenzene	20 U
87-86-5	Pentachlorophenol	100 U
95-01-8	Phenanthrene	20 U
120-12-7	Anthracene	20 U
94-74-2	Di-n-Butylphthalate	1.3 BT
206-44-0	Fluoranthene	20 U
92-87-5	Benzidine	100 U
129-00-0	Pyrene	20 U
35-68-7	Butylbenzylphthalate	20 U
91-94-1	3, 3'-Dichlorobenzidine	20 U
56-55-3	Benzo(a)Anthracene	20 U
117-81-7	bis(2-Ethylhexyl)Phthalate	12 BT
218-01-9	Chrysene	20 U
117-84-0	Di-n-Octyl Phthalate	2.1 J
205-99-2	Benzo(b)Fluoranthene	20 U
207-08-9	Benzo(k)Fluoranthene	20 U
50-32-8	Benzo(a)Pyrene	20 U
193-39-5	Indeno(1, 2, 3-cd)Pyrene	20 U
53-70-3	Dibenz(a, h)Anthracene	20 U
191-24-2	Benzo(g, h, i)Perylene	20 U

(1)-Cannot be separated from diphenylamine

Laboratory Name: HAZLETON

Case No: 4571

Sample Number  
**ED 806**

Organics Analysis Data Sheet  
(Page 3)

RECEIVED NOV 26 1985

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7-1-85

Date Analyzed: 7-13-85

Conc/Dil Factor: 1

CAS Number (Circle One)  
ug/l or ug/kg

319-84-6	Alpha-BHC	0.05 U
319-85-7	Beta-BHC	0.05 U
319-86-8	Delta-BHC	0.05 U
58-89-9	Gamma-BHC (Lindane)	0.05 U
76-44-8	Heptachlor	0.05 U
309-00-2	Aldrin	0.05 U
1024-57-3	Heptachlor Epoxide	0.05 U
959-98-8	Endosulfan I	0.05 U
60-57-1	Dieldrin	0.10 U
72-55-9	4, 4'-DDE	0.10 U
72-20-8	Endrin	0.10 U
33213-65-9	Endosulfan II	0.10 U
72-54-8	4, 4'-DDD	0.10 U
7421-93-4	Endrin Aldehyde	0.10 U
1031-07-8	Endosulfan Sulfate	0.10 U
50-29-3	4, 4'-DDT	0.10 U
72-43-5	Methoxychlor	0.50 U
53494-70-5	Endrin Ketone	0.10 U
57-74-9	Chlordane	0.50 U
8001-35-2	Toxaphene	1.0 U
12674-11-2	Aroclor-1016	0.50 U
11104-28-2	Aroclor-1221	0.50 U
11141-16-5	Aroclor-1232	0.50 U
53469-21-9	Aroclor-1242	0.50 U
12672-29-6	Aroclor-1248	0.50 U
11097-69-1	Aroclor-1254	1.0 U
11096-82-5	Aroclor-1260	1.0 U

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$V_s$  1000 or  $W_s$  —  $V_t$  10,000  $V_i$  4.0

Organics Analysis Data Sheet  
(Page 4)

RECEIVED NOV 26 1985

## Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. 544-25-2	1,3,5 Cycloheptatriene	BNA	133	15
2. 95-47-6	1,2 Dimethyl benzene	BNA	238	8.3
3.	Unknown	BNA	766	35
4.	Unknown	BNA	822	260
5.	Unknown	BNA	1937	8.0
6.	Unknown	BNA	1996	14
7.	Unknown	BNA	2073	9.8
8.	Unknown	BNA	2397	9.7
9.				
10. 60-29-7	Ethane, 1,1'-oxybis-	VOA	169	46
11.	See sample EDBOS FRN 28299			
12.	for Library search information			
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14.				
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ED 807

Organics Analysis Data Sheet  
(Page 1)

RECEIVED NOV 26 1985  
SFH 10 R06

Laboratory Name: HAZLETON LABORATORIES

Case No: 4571

Lab Sample ID No: 50606611

QC Report No: \_\_\_\_\_

Sample Matrix: WATER

Contract No: 68-01-6961

Data Release Authorized By: David C. Gill

Date Sample Received: 6-28-85

Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7-3-85 FEN 28302

Date Analyzed: 7-3-85 2145

Conc/Dil Factor: 1 pH \_\_\_\_\_

Percent Moisture: \_\_\_\_\_

Percent Moisture (Decanted): \_\_\_\_\_

Blank

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	10 U
74-83-9	Bromomethane	10 U
75-01-4	Vinyl Chloride	10 U
75-00-3	Chloroethane	10 U
75-09-2	Methylene Chloride	5 U
67-64-1	Acetone	10 U
75-15-0	Carbon Disulfide	5 U
75-35-4	1, 1-Dichloroethene	5 U
75-34-3	1, 1-Dichloroethane	5 U
156-80-5	Trans-1, 2-Dichloroethene	5 U
67-66-3	Chloroform	5 U
107-05-2	1, 2-Dichloroethane	5 U
78-93-3	2-Butanone	10 U
71-55-6	1, 1, 1-Trichloroethane	5 U
56-23-5	Carbon Tetrachloride	5 U
108-05-4	Vinyl Acetate	10 U
75-27-4	Bromodichloromethane	5 U

CAS Number		ug/l or ug/Kg (Circle One)
79-34-5	1, 1, 2, 2-Tetrachloroethane	5 U
78-87-5	1, 2-Dichloropropane	5 U
10061-02-6	Trans-1, 3-Dichloropropene	5 U
79-01-6	Trichloroethene	5 U
124-48-1	Dibromochloromethane	5 U
79-00-5	1, 1, 2-Trichloroethane	5 U
71-43-2	Benzene	5 U
10061-01-5	cis-1, 3-Dichloropropene	5 U
110-75-8	2-Chloroethylvinylether	10 U
75-25-2	Bromoform	5 U
591-78-6	2-Hexanone	10 U
108-10-1	4-Methyl-2-Pentanone	10 U
127-18-4	Tetrachloroethene	5 U
108-88-3	Toluene	5 U
108-90-7	Chlorobenzene	5 U
100-41-4	Ethylbenzene	5 U
100-42-5	Styrene	5 U
	Total Xylenes	5 U

Data Reporting Qualifiers

For reporting results to EPA, the following results qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

**Value** If the result is a value greater than or equal to the detection limit, report the value.

**U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution factors. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.

**J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicates the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10U)

**C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/ul in the final extract should be confirmed by GC/MS.

**B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.

**Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Organics Analysis Data Sheet  
(Page 3)

RECEIVED NOV 26 1985

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7-1-85

Date Analyzed: 7-13-85

Conc/Dil Factor: 1

CAS Number (ug/l or ug/Kg)  
(Circle One)

319-84-6	Alpha-BHC	0.05 U
319-85-7	Beta-BHC	0.05 U
319-86-8	Delta-BHC	0.05 U
58-89-9	Gamma-BHC (Lindane)	0.05 U
76-44-8	Heptachlor	0.05 U
309-00-2	Aldrin	0.05 U
1024-57-3	Heptachlor Epoxide	0.05 U
959-98-8	Endosulfan I	0.05 U
50-57-1	Dieldrin	0.10 U
72-55-9	4, 4'-DDE	0.10 U
72-20-8	Endrin	0.10 U
33213-65-9	Endosulfan II	0.10 U
72-54-8	4, 4'-DDD	0.10 U
7421-93-4	Endrin Aldehyde	0.10 U
1031-07-8	Endosulfan Sulfate	0.10 U
50-29-3	4, 4'-DDT	0.10 U
72-43-5	Methoxychlor	0.50 U
53494-70-5	Endrin Ketone	0.10 U
57-74-9	Chlordane	0.50 U
8001-35-2	Toxaphene	1.0 U
12674-11-2	Aroclor-1016	0.50 U
11104-28-2	Aroclor-1221	0.50 U
11141-16-5	Aroclor-1232	0.50 U
53469-21-9	Aroclor-1242	0.50 U
12672-29-6	Aroclor-1248	0.50 U
11097-69-1	Aroclor-1254	1.0 U
11096-82-5	Aroclor-1260	1.0 U

$V_i$  = Volume of extract injected (ul)

$V_s$  = Volume of water extracted (ml)

$W_s$  = Weight of sample extracted (g)

$V_t$  = Volume of total extract (ul)

$v_s$  1000 or  $W_s$  —  $v_t$  10,000  $v_i$  4.0

Laboratory Name: HAZLETON

Case No: 4571

Sample Number  
**ED 807**

Organics Analysis Data Sheet  
(Page 2)

RECEIVED NOV 26 1985

Semivolatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 7/1/85 FRN BAN 594

Date Analyzed: 7/2/85 1906

Conc/Dil Factor: 1

CAS Number		ug/l or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	20 U
108-95-2	Phenol	20 U
62-53-3	Aniline	20 U
111-44-4	bis(2-Chloroethyl)Ether	20 U
95-57-8	2-Chlorophenol	20 U
541-73-1	1, 3-Dichlorobenzene	20 U
106-46-7	1, 4-Dichlorobenzene	20 U
100-51-6	Benzyl Alcohol	20 U
95-50-1	1, 2-Dichlorobenzene	20 U
95-48-7	2-Methylphenol	20 U
39638-32-9	bis(2-chloroisopropyl)Ether	20 U
106-44-5	4-Methylphenol	20 U
621-64-7	N-Nitroso-Di-n-Propylamine	20 U
67-72-1	Hexachloroethane	20 U
98-95-3	Nitrobenzene	20 U
78-59-1	Isophorone	20 U
88-75-5	2-Nitrophenol	20 U
105-67-9	2, 4-Dimethylphenol	20 U
65-85-0	Benzoic Acid	100 U
111-91-1	bis(2-Chloroethoxy)Methane	20 U
120-83-2	2, 4-Dichlorophenol	20 U
120-82-1	1, 2, 4-Trichlorobenzene	20 U
91-20-3	Naphthalene	20 U
106-47-8	4-Chloroaniline	20 U
87-68-3	Hexachlorobutadiene	20 U
59-50-7	4-Chloro-3-Methylphenol	20 U
91-57-6	2-Methylnaphthalene	20 U
77-47-4	Hexachlorocyclopentadiene	20 U
88-06-2	2, 4, 6-Trichlorophenol	20 U
95-95-4	2, 4, 5-Trichlorophenol	100 U
91-58-7	2-Chloronaphthalene	20 U
88-74-4	2-Nitroaniline	100 U
131-11-3	Dimethyl Phthalate	20 U
208-96-8	Acenaphthylene	20 U
99-09-2	3-Nitroaniline	100 U

CAS Number		ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthene	20 U
51-28-5	2, 4-Dinitrophenol	100 U
100-02-7	4-Nitrophenol	100 U
132-64-9	Dibenzofuran	20 U
121-14-2	2, 4-Dinitrotoluene	20 U
606-20-2	2, 6-Dinitrotoluene	20 U
84-66-2	Diethylphthalate	20 U
7005-72-3	4-Chlorophenyl-phenylether	20 U
86-73-7	Fluorene	20 U
100-01-6	4-Nitroaniline	100 U
534-52-1	4, 6-Dinitro-2-Methylphenol	100 U
86-30-6	N-Nitrosodiphenylamine (1)	20 U
101-55-3	4-Bromophenyl-phenylether	20 U
118-74-1	Hexachlorobenzene	20 U
87-86-5	Pentachlorophenol	100 U
85-01-8	Phenanthrene	20 U
120-12-7	Anthracene	20 U
94-74-2	Di-n-Butylphthalate	100 U
206-44-0	Fluoranthene	20 U
92-87-5	Benzidine	100 U
129-00-0	Pyrene	20 U
95-68-7	Butylbenzylphthalate	20 U
91-94-1	3, 3'-Dichlorobenzidine	20 U
56-55-3	Benzo(a)Anthracene	20 U
117-81-7	bis(2-Ethylhexyl)Phthalate	20 U
218-01-9	Chrysene	20 U
117-84-0	Di-n-Octyl Phthalate	20 U
205-99-2	Benzo(b)Fluoranthene	20 U
207-03-9	Benzo(k)Fluoranthene	20 U
50-32-8	Benzo(a)Pyrene	20 U
193-39-5	Indeno(1, 2, 3-cd)Pyrene	20 U
53-70-3	Dibenz(a, h)Anthracene	20 U
191-24-2	Benzo(g, h, i)Perylene	20 U

(1) - Cannot be separated from diphenylamine

Organics Analysis Data Sheet  
(Page 4)

RECEIVED NOV 26 1985

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1. 513360	Propane, 1-chloro-2-methyl-	VOA	63	18.
2.				
3. —	None Found	BAN	—	—
4.				
5.				
6.				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				



MEMORANDUM

RECEIVED NOV 26 1985

DATE: 9/9/85

TO: Data Review Team  
Sample Management Office

FROM: Chuck Eddy for  
USEPA Region V

SUBJECT: Data Review Request

COPIES: RENEE MAYS, E & E  
CURTIS ROSS, CRL

Please review the data from the following SMO Case:

SMO Case No.: 4571

Site Name: RIPON LANDFILL

Lab Name(s): HAZELTON

\_\_\_\_\_

\_\_\_\_\_

L Sample Information:

- A. Number of Samples in Case: 6
- B. Number of Samples to be Reviewed: 6

(List Numbers if Not All)


- C. Organics to be Reviewed? Yes X No
- D. Inorganics to be Reviewed? Yes    No X

Site Name / TDD#: RIPON LANDFILL / R5-8310-01A  
Case Number: 4571  
Sampling Date: 6-27-85  
Sampling Time: 08:25  
Sample/Station Location: Sample #1 / G-5A (Matrix Spike Duplicate)

Organic Traffic Number E 0802  
Inorganic Traffic Number MEC994  
High Hazard Traffic Number E

Physical Description

At time of collection: clear

Physical Changes (if any)

From time of collection until shipment: NONE

Instrument Readings (i.e. - pH, conductivity...):

pH = 7 cond = 600  $\mu$ mhos

Sampling Date: 6-27-85

Sampling Time: 08:30

Sample/Station Location: Sample #2 / G-5

Organic Traffic Number ED803  
Inorganic Traffic Number MEC995  
High Hazard Traffic Number E

Physical Description

At time of collection: clear

Physical Changes (if any)

From time of collection until shipment: NONE

Instrument Readings (i.e. - pH, conductivity...):

pH = 7  
cond = 500  $\mu$ mhos.

Site Name / TDD#: RIPON LANDFILL / R5-8310-01A  
Case Number : 4571  
Sampling Date: 6-27-85  
Sampling Time: 09:25  
Sample/Station Location: Sample #3 / G-6

Organic Traffic Number ED804  
Inorganic Traffic Number MEC996  
High Hazard Traffic Number E

Physical Description  
At time of collection: slightly cloudy

Physical Changes (if any)  
From time of collection until shipment: NONE

Instrument Readings (i.e. - pH, conductivity...):  
pH = 7  
cond = 450  $\mu$ mhos

Sampling Date: 6-27-85  
Sampling Time: 09:35  
Sample/Station Location: Sample #4 / G-8

Organic Traffic Number ED 805  
Inorganic Traffic Number MEC997  
High Hazard Traffic Number E

Physical Description  
At time of collection: Reddish, cloudy

Physical Changes (if any)  
From time of collection until shipment: NONE

Instrument Readings (i.e. - pH, conductivity...):  
pH = 8  
cond = 1500  $\mu$ mhos.

Site Name / TDDI: RIPON LANDFILL / R5-8310-01A  
Case Number: 4571  
Sampling Date: 6-27-85  
Sampling Time: 08:40  
Sample/Station Location: Sample #5 / Dupl-5 (Dup of G-5)

Organic Traffic Number ED806  
Inorganic Traffic Number MEF209  
High Hazard Traffic Number E

Physical Description

At time of collection: clear

Physical Changes (if any)

From time of collection until shipment: NONE

Instrument Readings (i.e. - pH, conductivity...): \_\_\_\_\_

pH = 7  
cond = 500  $\mu$ mhos

Sampling Date: 6-27-85  
Sampling Time: 10:30  
Sample/Station Location: Sample #6 / BLANK

Organic Traffic Number ED807  
Inorganic Traffic Number MEF210  
High Hazard Traffic Number E

Physical Description

At time of collection: clear

Physical Changes (if any)

From time of collection until shipment: NONE

Instrument Readings (i.e. - pH, conductivity...): \_\_\_\_\_

pH = 6  
cond = 0

IIC	MEC 994	MEC 995	MEC 996	MEC 997	MEF 202	MEF 210				
OTC	ED802	ED803	ED804	ED805	ED806	ED807				
SAMPLE	5A	5	6	8	Dup of 5	Blank				

**COMPOUND**

pentachlorophenol										
phenanthrene										
anthracene										
di-n-butylphthalate										
fluoranthene										
benzidine										
pyrene										
butylbenzylphthalate										
3,3'-dichlorobenzidine										
benzo(a)anthracene										
bis(2-ethylhexyl)phthalate										
chrysene										
di-n-octyl phthalate										
benzo(b&k)fluoranthene										
benzo(a)pyrene										
indeno(1,2,3-cd)pyrene										
dibenzo(a,h)anthracene										
benzo(g,h,i)perylene										
alpha-BHC										
beta-BHC										
delta-BHC										
gamma-BHC(lindane)										
heptachlor										
aldrin										
heptachlor epoxide										
endosulfan I										
dieldrin										
4,4'-DDE										
endrin										
endosulfan II										
4,4'-DDD										
endrin aldehyde										
endosulfan sulfate										
4,4'-DDT										
methoxychlor										
endrin ketone										
chlorodane										
toxaphene										
Aroclor-1016										
Aroclor-1221										
Aroclor-1232										
Aroclor-1242										
Aroclor-1248										
Aroclor-1254										
Aroclor-1260										
ELEMENT										
aluminum										
antimony										
arsenic										
barium										
beryllium										
cadmium										
calcium										
chromium										
cobalt										
copper										
iron										
lead		5.8	2.2							
magnesium										
manganese										
mercury										
nickel										
potassium										
selenium										
silver										
sodium										
thallium										
tin										
vanadium										
zinc										
cyanide CHECK IF ANALYZED ( )										
TENTATIVELY IDENTIFIED ORGANICS										

All values ug/l

COMPOUND	ITC	MEC 974	MEC 975	MEC 976	MEC 977	MEF 209	MEF 210	Wisconsin DNR Sampling	
	OTC	ED802	ED803	ED804	ED805	ED806	ED807	11/12/84	5/13/85
	SAMPLE	MW5A	MW5	MW6	MW8	Dup. of MW5	Blank	Falkenberg Well	MW5
chloromethane									
bromomethane						82		45.8; 47.5	120; 210; 200
vinyl chloride			110						
chloroethane									
methylene chloride									
acetone									
carbon disulfide									2.5; 2.1
1,1-dichloroethene									
1,1-dichloroethane						740			460; 340
trans-1,2-dichloroethene			970						
chloroform									
1,2-dichloroethane				20					
2-butanone		14							
1,1,1-trichloroethane									
carbon tetrachloride									
vinyl acetate									
bromodichloromethane									
1,1,2,2-tetrachloroethane									
1,2-dichloropropane									
trans-1,3-dichloropropene									88; 7.1
trichloroethene									
dibromochloromethane									
1,1,2-trichloroethane									3.5; 4.2
benzene									
cis-1,3-dichloropropene									
2-chloroethylvinylether									
bromoform									
2-hexanone									
4-methyl-2-pentanone									30; 2.5
tetrachloroethene									
toluene									
chlorobenzene									
ethylbenzene									
styrene									
total xylenes			37			28			61; 8.6
N-nitrosodimethylamine									
phenol									
aniline									
bis(2-chloroethyl)ether									
2-chlorophenol									
1,3-dichlorobenzene									
1,4-dichlorobenzene									
benzyl alcohol									
1,2-dichlorobenzene									
2-methylphenol									
bis(2-chloroisopropyl)ether									
4-methylphenol									
N-nitroso-di-n-propylamine									
hexachloroethane									
nitrobenzene									
isophrone									
2-nitrophenol									
2,4-dimethylphenol									
benzoic acid									
bis(2-chloroethoxy)methane									
2,4-dichlorophenol									
1,2,4-trichlorobenzene									
naphthalene									
4-chloroaniline									
hexachlorobutadiene									
4-chloro-3-methylphenol									
2-methylnaphthalene									
hexachlorocyclopentadiene									
2,4,6-trichlorophenol									
2,4,5-trichlorophenol									
2-chloronaphthalene									
2-nitroaniline									
dimethyl phthalate									
acenaphthylene									
3-nitroaniline									
acenaphthene									
2,4-dinitrophenol									
4-nitrophenol									
dibenzofuran									
2,4-dinitrotoluene									
2,6-dinitrotoluene									
diethylphthalate									
4-chlorophenyl-phenylether									
fluorene									
4-nitroaniline									
4,6-dinitro-2-methylphenol									
N-nitrosodiphenylamine									
4-bromophenyl-phenylether									
hexachlorobenzene									

57