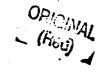
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Site Inspection Report

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Site Investigation of Ohio River Park Neville Island Allegheny County Pennsylvania PA #0170

Prepared by the Commonwealth of Pennsylvania Department of Environmental Resources

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> > April 21, 1989

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1.0 INTRODUCTION

1.1 AUTHORIZATION

As a result of a site preliminary assessment and pursuant to preliminary assessment and site inspection Cooperative Agreement between the Department and the U. S. Environmental Protection Agency, the Department has conducted a Site Inspection at this site. The Department has determined that there is a potential risk of a significant release of hazardous substances from this site. Therefore, a potential risk to human health and the environment exists warranting further study through a Site Inspection.

1.2 SCOPE OF WORK

The Department was tasked to conduct a desktop Site Inspection of the Ohio River Park site. The Site Inspection centers on the possible contamination of surface water, groundwater, soil, and air from chemical substances on the site.

1.3 SUMMARY

Ohio River Park is a 35 acre site located at the western tip of Neville Island in Neville Township, Allegheny County, Pennsylvania (see (Figure 1). In 1976, the Neville Land Company (NLC) donated the land to Allegheny County for development as a park. During construction of the park in 1978, the County came to believe that industrial waste found at the site might represent a health hazard. The County retained Fred C. Hart Associates (FCHA) to conduct an investigation assessing the potential public health hazards at the site. The FCHA report submitted in July 1979 concluded that the site did pose a potential risk to the public health.¹ The County decided at that time to abandon plans to open the park. After discussions between NLC and the County, ownership of the park reverted back to NLC.²

The FCHA studies, and subsequent investigations by Environmental Research and Technology, Inc. (ERT) who were retained by NLC, determined that the site had been used for the disposal of industrial wastes starting in the late 1940's or early 50's and continued until the mid-1960s.³ FCHA estimated that 225,000 cubic yards of waste had been disposed of on the site since 1941.⁴ The effect of this activity on the groundwater and surface water of the site

- ¹ Fred C. Hart Associates. An Investigative Study of Potential Public Health Hazards at Ohio River Park. July 23, 1979, p. 55.
- 2 ERT, Inc. Plan of Study for Neville Island Site Investigation. December, 1980, Section 1-1.
- 3 ERT, Inc. Preliminary Risk Assessment of Neville Island Site. April, 1981, Section 3.3.
- Fred C. Hart Associates. Assessment of Remedial Options at Ohio River Park. January, 1980, p. 40.

is documented by the reports of both FCHA and ERT: elevated levels of organics, $O_{r_{\rm eleval}}$ pesticides, and metals have been detected in both groundwater and surface run-off.1,2 Tables 1 and 2 are summaries of sampling conducted by ERT and FCHA \blacktriangleright (Fers) respectively.

The following is a list of chemicals in excess of detection limits:

- Benzene, toluene, ethylbenzene, xylenes, phenol, 2-chlorophenol, 2,4-dichlorophenol, 2,4,6-trichlorophenol, naphthalene and 2,4-D have been confidently identified with levels greater than 3 times the detection limit and 5 times the background;
- 2) Diethylphthalate and bis(2-ethylhexyl)phthalate have been identified but with levels less than 3 times the detection limits;
- 3) Silvex and 2,4,5-T have been identified using one GC column with levels greater than 3 times the detection limits and 5 times that of background:
- 4) Arsenic, beryllium, cadmium, chromium, copper, iron, lead, mercury, nickel, selenium, thallium, and zinc were reported by ERT to be present on the site with levels greater than 5 times background; and
- 5) Cyanide was detected on the site with levels greater than 5 times background.

Among the chemicals listed above, benzene and arsenic are well known human carcinogens (Group A). 2,4,6-Trichlorophenol, beryllium, and lead are suspected human carcinogens (Group B2).

ERT, Inc. Preliminary Risk Assessment of Neville Island Site, April, 1981, Table 1-2.

Fred C. Hart Associates. An Investigative Study of Potential Public Health Hazardous at Ohio River Park, July 23, 1979, pp. 39-44.

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2.0 SITE DESCRIPTION

2.1 SITE LOCATION

Ohio River Park is located on Neville Island in Neville Township, Allegheny County, Pennsylvania. The site encompasses the western portion of Neville Island along river mile points 9.5-9.9. Coordinates are 40° 31' 06" N latitude and 80° 08' 10" W longitude. The USGS map location is 2.2 inches north and 3.8 inches west of the southeast corner of the USGS Ambridge quadrant.¹ See Figure 2.

2.2 SITE LAYOUT

Ohio River Park consists of a nearly completed park, including an administrative building, two outhouses, bleachers, footers, paved roads and parking areas. The site is relatively flat, with steep banks leading to small or nonexisting beaches. On the back channel, portions of the bank are undercut by the river. Three surface water runoff outfalls, two on the main channel and one on the back channel, are known to exist (see Figure 16). The majority of the facility lies at an elevation of 715' - 720'. The highest point is an observation knoll at 735' near the western tip of the island. The normal pool elevation of the Ohio River at this end of the site is 692'.2

Waste disposal, primarily trenching and end dumping, occurred extensively throughout the southern and western portions of the site (see Figures 9 and 18). Water ponding occurs randomly throughout the site in surface depressions. Areas of hard semi-impermeable layers of solidified waste residues retard drainage and serve as pockets to entrap surface runoff.

The Park is accessed by a gate on Grand Avenue, approximately 300 feet east of the Coraopolis Bridge on its eastern border. See Figure 3.

2.3 OWNERSHIP HISTORY

The Ohio River Park site was acquired by Pittsburgh Coke and Iron Company in the 1920s. On October 19, 1944, Pittsburgh Coke and Iron Company was renamed Pittsburgh Coke and Chemical Company (PC&C).³

PC&C continued to own the property until Aug. 14, 1970 when it conveyed the property to a wholly owned subsidiary, Neville Land Company. At about that time PC&C, a majority owned subsidiary of the Hillman Company, became wholly owned. Pittsburgh Coke and Chemical Company is no longer in existence as a result of a merger.⁴

- United States Geological Survey. Ambridge, Pennsylvania Quadrangle, 7.5 Ninute Series. Topographic Map, 1960, photorevised 1979.
- 2 Ibid.
- 3 ERT, Inc. Detailed Description of Neville Island Site, August, 1981, Section 3.1.1.
- 4 Rittmeyer, Robert W., ERT, Inc. Correspondence to James R. Shack, PADER, May 16, 1988.

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In 1976, NLC donated the site to Allegheny County for the construction of a park. FCHA delivered the report in July, 1979 and concluded that a potential public health threat did exist. In 1980, ownership reverted back to NLC.¹

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2.4 SITE USE HISTORY

The Neville Land Company site was used primarily for agricultural purposes prior to 1947. From the mid-1930s through the mid-1950s a portion of the NLC site was used for the landfilling of municipal refuse.² During 1943 a \times housing project for the Navy was built at the eastern end of the NLC site. These barracks were demolished and removed from the site in 1970.³

In 1947-1948, some of the site's topsoil was removed for use in Forbes Field.⁴ From 1952 until 1965, trenches were dug at the site to dispose of PC&C coking sludges and wastes from the production of cement, and various pesticides. \times In addition, other industrial wastes such as plant demolition materials and slag were disposed of on the site surface.⁵ See Figure 4. A number of trenches have been identified by aerial photographs. However, no records exist of activities that may have occurred between the photograph sessions.

In 1977 the Hillman Company donated the site to Allegheny County⁶ and construction of the park began. During the excavation effort, various wastes, including thirty drums of liquid product, were uncovered.⁷ This alerted the County to the possibility that the site had been used for the disposal of hazardous wastes. After assessments by FCHA in 1979 and 1980, plans to open the virtually completed park were abandoned and ownership of the site reverted back to NLC.⁸ NLC retained ERT to assess the hazards associated with the site and to determine remedial actions.⁹ Table 3 summarizes these activities.

- Rittmeyer, Robert W., ERT, Inc. Correspondence to James R. Shack, PADER, May 16, 1988.
- 2 ERT, Inc. Preliminary Risk Assessment of Neville Island Site, April, 1981, Section 3.3.
- 3 ERT, Inc. Detailed Description of Neville Island Site, August, 1981, Section 3.1.2.
- 4 Ibid., Table 3-6.
- 5 ERT, Inc. Preliminary Risk Assessment of Neville Island Site, April, 1981, Section 1.1.
- ⁶ ERT, Inc. Correspondence to James R. Shack, PADER, May 16, 1988.
- Fred C. Hart Associates. An Investigative Study of Potential Public Health Hazards at Ohio River Park, July 23, 1979, pp. 28-34.
- 8 Rittmeyer, Robert W., ERT, Inc. Correspondence to James R. Shack, PADER, May 16, 1988.

9 ERT, Inc. Detailed Description of Neville Island Site, August, 1981, Section 3.3.4.

2.5 PERMIT AND REGULATORY ACTION HISTORY

To date, no permit or regulatory action has been initiated regarding the Ohio River Park Site.

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2.6 PREVIOUS SITE STUDIES

The Neville Island site has not been used as a waste disposal site since the mid-1960's. Since 1972 four firms have conducted investigations at the site for various purposes. In 1972 and 1973, Pittsburgh Testing Laboratory conducted a subsurface soil investigation for a proposed tank farm under contract to Exxon Company, U.S.A. In 1977, Richardson, Gordon, and Associates, Inc., made a similar study for the construction of a county park under the contract to the Environmental Planning and Design Partnership. FCHA studied the site to determine potential health hazards and to assess remedial options under contract to the Allegheny County Health Department.¹

During the summer of 1980, ERT was engaged by NLC to investigate potential hazards at the site, including air and water monitoring. As a result of ERT's investigation the following reports were prepared:

- 1. Plan of Study for Neville Island Site Investigation. December 1980.
- 2. Concepts for Remedial Action at Neville Island. December 1980.
- 3. Preliminary Risk Assessment of Neville Island Site. April 1980.
- 4. Interim Monitoring Report for Neville Island Site. June 1981.
- 5. Detailed Description of Neville Island Site. August 1981.
- 6. Recommendations for Remedial Action at Neville Island Site. January 1982.
- 7. Security, Monitoring, Inspection and Maintenance Programs for Neville Island Site. March 1983.
- 8. Review of Initial Two Years of Long-Term Groundwater Monitoring Program for the Neville Island Site. April 1985.

In March of 1986, Pa. DER performed a Preliminary Assessment and gave the facility a high priority with regards to the need for a Site Inspection.²

- ERT, Inc. Detailed Description of Maville Island Site, August, 1981., Section 3.3.
- Preliminary Assessment. Pennsylvania Department of Environmental Resources, PA 95, March, 1986.

2.7 REMEDIAL ACTIONS TO DATE

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In September 1981, one barrel of nearly pure 2,4-D and the contaminated surrounding soils were removed from the site and disposed at a landfill in Ohio. In November of that year, a narrow 60 foot long section of the northern shore was stabilized by the construction of a rip-rap buttress in an area of exposed desulfurization waste. In addition, 450 cubic yards of clayey silt was placed \times over a 0.2 acre area of exposed desulfurization waste adjacent to the buttress.¹ \times

Several actions were taken to secure the site from unauthorized and unknowing entry, including the installation of a nine foot high chain link fence with a locked gate on Grand Avenue. The fence runs along Von Stein Lane and Grand Avenue from the Main Channel to the Back Channel. To discourage entry from the northern, western, and southern shorelines, bushes were planted in areas of naturally sparse vegetation or where steep slopes did not exist. ERT decided that areas along the shoreline that had thick existing vegetation and/or steep slopes leading from the shoreline to the park interior were not required to have any modifications to discourage unauthorized entry. The entire perimeter has been posted with warning signs to discourage entry.² The effectiveness of these precautions is uncertain, as access to the site still occurs.

ERT has done monitoring of 10 perimeter wells on a quarterly basis from January 1981 until January 1985. From January 1985 until the present time, monitoring of these wells has continued on a semiannual basis.³ Figure 5 shows the location of the perimeter wells.

ERT, Inc. Recommendations for Remedial Action at Neville Island Site, January, 1982, Section 3.2

2 ERT, Inc. Security, Monitoring, Inspection, and Maintenance Programs for Neville Island Site, March, 1983, Section 3.2.

³ ERT, Inc. Review of Initial Two Years of Long-Term Groundwater Monitoring Program for the Neville Island Site, April, 1985, p.i.

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3.0 ENVIRONMENTAL SETTING

3.1 WATER SUPPLIES

All users of the Ohio River are potentially affected because the surface water and groundwater from the site eventually enters the Ohio River. Of these potential receptors, the largest group affected are the municipal water suppliers. See Figure 6. The water suppliers in the area include:

- A. The Borough of Coraopolis, which draws its water supply from a wellfield on the south side of the back channel. Nine wells exist, but only seven are currently used, the closest of these about 700 feet southwest of the site across the back channel of the Ohio River. The wells range in depth from 60 to 67 feet and tap the Quaternary Sand and Gravel Unit. The Borough distributes an average of 1,000,000 gpd (gallons per day) to approximately 8200 people.¹
- B. The Dixmont State Hospital, which utilized a surface water intake on the north shore of the main channel at mile point 8.0, 1.5 miles upstream from the nearest known outfall on the site. Dixmont is presently closed.
- C. The Moon Township Water Authority, which draws its water from a wellfield on the south side of the Ohio River at river mile point 11.5, 1.8 stream miles downstream of the nearest known outfall at the site. The Authority employs two vertical wells and a Ranney well to tap the Quaternary Sand and Gravel Unit. The Authority distributes an average of 3,000,000 gpd to approximately 30,000 people.²
- D. The Robinson Township Municipal Authority, which draws its water supply from a surface water intake on the south shore of the back channel approximately 0.9 miles upstream of the nearest known outfall of the site. Robinson Township distributes an average of 3,000,000 gpd to approximately 30,000 people.³
- E. The Sewickley Waterworks, which employs a surface water intake located on the north shore of the Ohio River at river mile point 12.8, about 1.6 miles downstream of the nearest known outfall from the site. The Sewickley Waterworks distributes an average of 800,000 gpd to approximately 10,000 people.⁴
- ¹ Barrone, Robert, Coraopolis Department of Public Works, with William Bailey, PADER. Telecon. August 29, 1988.
- ² Mr. Zollinger. Moon Township Municipal Authority, with William Bailey, PADER. Telecon. July 13, 1988.
- ³ Neil, Charles, Robinson Township Municipal Authority, with William Bailey, PADER. Telecon. August 31, 1988.
- 4 Tucci, Ernest. Sewickley Waterworks, with William Bailey, PADER. Telecon. July 11, 1988.

F. West View Municipal Authority, which utilizes a surface water intake and 3 wells at the eastern tip of the Island (river mile point 5.0) and 7 wells on Davis Island (river mile point 4.7). The wells, which are used only in emergency situations such as the Ashland oll spill in January, 1988, tap the Quaternary Sand and Gravel Unit and range in depth from 40 to 74 feet. The surface water intake lies about 4.5 miles upstream from the closest known outfall on the site. West View distributes an average of 18,000,000 gpd to approximately 220,000 people.¹

There are no known private water supplies in the study area.

3.2 SURFACE WATERS

According to the U. S. Army Corps of Engineers, the main channel of the river carries the majority of the flow at an average of 32,500 cfs, while the back channel, under normal flood conditions, carries an average of 2,000 cfs. Approximately 2.8 miles upstream of the site at river mile point 6.4 is the Emsworth Back Channel Dam. This dam is a gated dam that keeps the pool elevation at 710'. Another dam, which is situated on the main channel of the Ohio River, is located at river mile point 13.2, approximately 3.4 miles downstream of the site. Any influence on the groundwater because of the difference in pool elevations would be from northeast to southwest.²

The migration of hazardous wastes from the site into the Ohio River by surface runoff was evidenced by analyses of ponded water and storm run-off at 50,000 ug/l (10-26-79) in ponded water and 24,000 ug/l (0utfall #1 on 10-26-79) at storm sewer outfalls; inorganics, such as sulfide, were also detected in ponded water at 5,500 ug/l (10-26-79) and in outfalls at 335,000 ug/l (0utfall #1, 10-26-79). See Table 4.

Pesticides have also been found at outfall areas. ERT detected 2,4-D levels at 2500 ug/l and 2,4,5-T at 114 ug/l in Outfall $#1^3$.

The Fish and Wildlife Service National Wetlands Inventory Map indicates several small wetland areas are present within the site study area, primarily on the Southern shore of the Ohio River across the back channel from Neville Island. None of the wetland areas within the study area have been field examined.

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3.3 GEOLOGY AND SOILS

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The site is located within the Allegheny Plateau section of the Appalachian Plateaus Physiographic Province. The structure of this area is characterized by gentle, parallel folds trending northeast-southwest.^{3, 4}

- Depp, Donald, Westview Municipal Authority, with William Bailey, PADER. Telecon. November 23, 1988.
- Hein, Paul. U. S. Army Corps of Engineers. Telecon with Deborah McNaughton, PADER, April 20, 1989.
- ³ Wagner, Walter R., et. al. Greater Pittsburgh Region Structure Contour Map, Map 43, 1975. Reprinted 1985.
 - Wagner, Walter R., et al. <u>Geology of the Pittsburgh</u> Area, General Geology 266 Report G 59, 1970, p 3.

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The bedrock in the immediate area of the site is the Conemaugh Group, a heterogeneous unit composed mainly of shales and sandstones of Pennsylvania age (see Figure 17). The Conemaugh Group consists of the Glenshaw (lower) and Casselman (upper) Formations. The lithology of the Glenshaw Formation is characterized by sandstones and shales, and to a lesser extent, limestones and coals. The best water producing unit of this formation is the Saltsburg Sandstone, with an average yield of 55 gpm. The lithology of the Casselman Formation is similarly characterized. The most important water producing unit of the Casselman is the Connellisville sandstone, with an average yield of 25 gpm. 1

Stratigraphically overlying the Conemaugh and outcropping six miles to the south of the site is the Monongahela Group, which is also Pennsylvanian in age. Nearly one half of the Monongahela is comprised of limestones, and the remainder consists of shales, sandstones, and coals. The most economically important unit of the Group is the Pittsburgh coal seam. This Group may not have good aquifers because of dewatering due to coal mining activity.²

Quaternary alluvial deposits which overlie the bedrock are found in major stream valleys. This alluvium, which consists of unconsolidated clay, silt, sand, and gravel, is an important aquifer in the region with yields of 5 to 3000 gpm. The maximum thickness of the valley alluvium is about 60 feet and is derived from two sources: the basal portion of the alluvium is coarsegrained and consists primarily of sand and gravel of glaciofluvial origin (Pleistocene); the upper portion consists of more recently deposited silts and clays eroded from local drainage basins.³

Neville Island is a detached portion of a dissected river terrace deposited by an ancestral Ohio River. The terrace is partly submerged by impoundments on the river, but.remnants of it flank both sides of the river at approximately the same elevation as the island.⁴

The geologic material which comprise Neville Island is a 60 foot sequence of unconsolidated alluvial deposits of clay, sand, and gravel overlying bedrock. See Figure 7 for a generalized stratigraphic column. The alluvium grades fine upward, with the fine-grained fraction ranging from zero thickness to fifteen feet, thickening toward the tip of the site and toward the northeastern corner. Silt and clay are generally found in discrete lenses with the upper part of the aquifer. The underlying coarse sand and gravel unit varies from 25 to 40 feet in thickness and extends laterally beneath the Ohio River, thickening in the central eastern portion of the site.⁵

Gallaher, John T. Summary of Ground Water Resources in Allegheny County, Pennsylvania. Water Resource Report 35, 1973, pp. 53 - 55.

2 Ibid., p. 50 - 53.

3

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Ibid., p. 39 - 47.

Adamson, J. H., et. al. Groundwater Resources of Valley-Fill Deposits of Allegheny County, Pennsylvania. Bulletin W 8, 1949, pp. 17 - 22.

Ibid., pp. 17 - 19.

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The top of the bedrock beneath the site lies at an elevation of approximately 660 feet, rising locally to 670 feet at monitoring well ERT-1. The bedrock is composed of relatively flat-lying, interbedded, micaceous sandstone (argillite). Sandstone is the dominant rock type encountered in borings on the site.

A further investigation would have to be made to determine a hydraulic \times connection between the aquifer and the underlying bedrock at the site.

The Conservation Service of the United States Department of Agriculture has classified the on-site soils as urban type were Figure 8). Urban land is characteristically level land situated on a floca wiain and is comprised of fill material that was placed over natural soils. Urban soils typically have a wide range of pH levels.²

In 1977, RGA prepared a Soil Reconnaissance Report and a subsurface investigation. The results showed that of the 50 test pits dug, only 5 indicated the presence of natural surface topsoil; the other 45 pits contained various amounts of waste material (see Figure 9). In some areas at the site, waste material extended to depths of 10 feet. Half of the test pits appeared to contain less than 5.0 feet of waste material with the rest contained 5.0 feet or more.³

ERT, Inc. Preliminary Risk Assessment of Neville Island Site, April, 1981, Section 4.3.1.4.

2 Coleman, Rita, PADER, with Deborah McNaughton, PADER, December 7, 1988.

³ Fred C. Hart Associates. An Investigative Study of Potential Health Hazards at Ohio River Park, July 23, 1979, pp. 10-11. Х

In 1979 FCHA tested 278 soil samples taken at 19 locations at depths ranging from surface to four feet. Soil was found in only three of the samples.¹ The range of the pH of the soils of the site is 1.5 to $9.1.^2$

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3.4 GROUNDWATERS

The top of the aquifer beneath the site lies about 25 feet below the ground surface.³ The saturated portion extends 35 to 40 feet into fluvial sediments which consist primarily of sand and gravel. This thickness is relatively uniform underneath the island, but decreases abruptly at the Main and Back Channels where erosion of the aquifer by the river current is likely to occur. Drilling records of offshore and onshore borings suggest the river has eroded 10 to 20 feet of coarse-grained part of the aquifer along most of the shoreline of the site. The aquifer, therefore, has a direct hydraulic connection with the river.⁴

Groundwater levels taken from September 1979 to April 1981 (see Figures 10-13) show the aquifer to be elongated in shape, with a slight on-site mounding. Groundwater flow, as derived from water table elevations, is radial from the central part of the mound toward the Back and Main Channels. A groundwater divide parallel to the long axis of the island is apparent: the divide separates north and south flow.⁵

Aquifer recharge is from the Ohio River and from precipitation. The high permeability and ready rechargeability from the Ohio River has made valley deposits a valuable source of groundwater to residents of Neville Island, Coraopolis and vicinity. About 40,000 residents utilize 10 wells for drinking supplies within 1.5 miles downstream of the site. As of 1979, more than 30 wells on the island were used for industrial cooling and municipal supplies, including the reserve wells of West View. All these wells tap the Quaternary Sand and Gravel unit. The induced infiltration is so great that the water pumped from the wells is chemically indistinguishable from water of the Ohio River.⁶

- ¹ Fred C. Hart Associates, An Investigative Study of Potential Health Hazards at Ohio River Park, July 23, 1979, pp. 23-24.
- 2 ERT, Inc. Detailed Description of Neville Island Site, August 1981, Tables 4-1, 4-3, 4-4, 4-6, and 4-7.
- ³ ERT, Inc. Preliminary Risk Assessment of Neville Island Site, April, 1981, Section 4.3.2.
- ⁴ Ibid., Section 4.3.5.1
- ⁵ Ibid., Section 4.3.4.

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⁶ Fred C. Hart Associates. An Investigative Study of Potential Public Health Hazards at Ohio River Park, July 23, 1979, p 8.

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Recharge (infiltration) is calculated using the following equation:

Recharge = (Precipitation - Runoff) - (Potential Evapotranpiration) (Area) (Time)

Typical runoff coefficients for flat sandy soil with grass cover range from 0.05 to 0.10; for flat, clayey soil, typical coefficients range from 0.13 to 0.17 (data selected from Table 3 of EPA Document 530/SW-168, 1975.) Based on observations of site topography, surficial soils, vegetation characteristics, and absence of erosion features related to surface runoff, ERT estimates the recharge to the aquifer to be about 700 gallons per acre per day. The recharge will eventually discharge into the Main or Back Channels.¹

Assuming that input to and discharge from the aquifer are equal over the long term, ERT calculates the discharge from the aquifer of about 10,000 gallons per day along both the north and south shoreline.

Darcy's law states that the rate of flow per unit area of an aquifer is proportional to the gradient of the potential head measured in the direction of flow.² Darcy's Equation describes discharge through a porous medium as a function of hydraulic conductivity, gradient, and area of flow. The following calculations are based on ERT's information:

The general equation is Q = KIA

where: Q is discharge (gal/day) K is hydraulic permeability (gal/day/ft²) i is hydraulic gradient (ft/ft) A is area of cross-section (ft²)

After rearranging terms, the hydraulic conductivity can be determined:

 $K = \frac{0}{1A}$

where: Q = 7000 gal/day calculated on about 10 acres subject to infiltration

i = 0.0025 ft/ft average gradient of water table

- ERT, Inc. Preliminary Risk Assessment of Neville Island Site, April, 1981, p. 4 - 14.
- Wilson, E. M. Engineering Hydrology. A Halsted Press Book, 1974, 2nd ed., p. 74.

$A = 62,700 \text{ ft}^2$

thickness of aquifer unit (38 feet)

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therefore:

K '=

 $\frac{7,000 \text{ gal/day}}{(0.0025 \text{ ft/ft})} (62,700 \text{ ft}^2)$

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or 45 gal/day/ft2

Assuming that Q in is equal to Q out, the hydraulic conductivity of the aquifer is about 6 ft/day.

The linear velocity, v_1 , can be determined by:

 $v_1 = \frac{KI}{n}$

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where: n = porosity estimated to be about 30% therefore:

$$v_1 = \frac{(6 \text{ ft/day}) (0.0025 \text{ ft/ft})}{0.30}$$

or 0.05 ft/day.^1

1 ERT, Inc. Preliminary Risk Assessment of Neville Island Park, April, 1981, Section 4.3.2 and 4.3.5.2.

3.5 CLIMATE AND METEOROLOGY

The average annual temperature for the Pittsburgh area is $61.9^{\circ}F$. The month with the coldest annual temperature is January which has an average temperature of $30.6^{\circ}F$. The warmest month, on average, is July with an average temperature of $74.6^{\circ}F$. The average annual precipitation is 36.30 inches for the area. The month with the lowest precipitation is November, with an average of 2.34 inches; and the highest is July, with an average of 3.83 inches.¹

Although the main concern for monitoring the site has been suspected wastes in the groundwater, surface waters and sediments, the site has also had a history of problems concerning air pollution. In the summer of 1978, the Allegheny County Health Department was called into investigate fumes emanating from areas of the site.² The troubled areas have since been covered with soils. X

3.6 LAND USE

The area of Neville Island immediately east of the site and extending 3/4 of a mile is primarily residential. The remainder of the Island to the east is mostly heavy industry. Coraopolis, on the mainland to the south is mixed use residential and commercial.³ See Figure 14.

3.7 POPULATION DISTRIBUTION

Based on United States Geologic Survey 7.5 minute series topographic maps of the Ambridge, Oakdale, Emsworth, and Pittsburgh West quadrangles and 1980 U.S. Census Bureau data, the population within a 3-mile radius of the site is 27,552.⁴

Eggers, Cindy, National Weather Service, with William Bailey, PADER. Telecon. July 13, 1988.

- Fred C. Hart Associates. An Investigative Study of Health Hazards at Ohio River Park, July 23, 1979, p. 4.
- ³ United States Geological Survey, Ambridge, Pennsylvania Quadrangle 7.5 Minute Series. Topographic map, 1960, photorevised 1979.

⁴ United States Geological Survey, Ambridge, Oakdale, Emsworth, and Pittsburgh West, Pennsylvania 7.5 Minute Series. Topographic maps, 1960, photorevised 1979.

3.8 CRITICAL ENVIRONMENTS

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According to the Pennsylvania Natural Diversity Inventory (PNDI), two species of special concern exist in the Neville Island area. The first is the invertebrate whose common name is Wabash Pigtoe (Fusconia flava). Although the species is not listed by state or federal regulations, it has been identified by the Pennsylvania Biological Survey's Invertebrate Technical Committee as a species of special concern. The evidence for the species location is a specimen collected by A. E. Ortmann in 1919 stored at the Carnegie Museum of Natural History. No modern research on this animal has been done. A second species known from the Island is the plant Vanilla Sweetgrass (<u>Hierochole odorata</u>). The plant is listed in Chapter 82 of the Pennsylvania Code as Pennsylvania Endangered. Evidence for the species exists from a specimen collected by J. A. Shafer in May, 1899 which is stored at the Carnegie Museum.¹

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Two federally listed endangered birds are expected to be found as transient species in the project area. They are the bald eagle (<u>Haliaeetus</u> <u>levcocephalus</u>) and the penegrine falcon (<u>Falco peregrinus</u>). There is no listed critical habitat for these species in the project area.²

McKenna, Kathy, PADER Bureau of Forestry, with Deborah McNaughton, PADER. Telecon. September 1, 1988.

² United States Department of the Interior, Fish and Wildlife Service, November, 1988.

4.0 WASTE TYPES AND QUANTITIES

ERT identified 13 types of waste based on physical description, process from which it was generated, and/or chemical analyses:¹

1. Desulfurization waste consists of iron oxide beds, wood chips and sulfuric acid sludges. Table 5 is a summary of the data collected on this waste by FCHA, RGA, and ERT.

ERT estimated the quantity to be 3,700 cubic yards. Only areas of known contamination were used in their estimates. FCHA estimates the quantity of highly contaminated soil containing this waste to be approximately 10,900 cubic yards.

- 2. Agricultural chemical waste; Pittsburgh Coke and Chemical manufactured, formulated, and resold pesticides. Table 6 summarizes the products involved. Of the 21 products listed, ERT analyzed only for 5: Silvex, 2,4-D, Malathion, 2,4,5-T and parathion. Low concentrations were found in over 50% of the soil samples analyzed. ERT reports that all former PC&C employees interviewed agree agricultural wastes were disposed randomly in the trenches. FCHA reports 4,700 cubic yards of highly contaminated soils are buried on site.
- 3. Coal coking sludges consist of tar decanter sludge and tar acid sludge. Table 7 is a summary of the chemical analysis of these sludges. Figure 15 illustrates the sampling points used by FCHA and ERT.

ERT estimates the quantity of sludges contained in the trenches to be 10,000 - 20,000 cubic yards. FCHA estimates the quantity disposed to be 32,000 cubic yards.

4. Foundry sand was found in large quantities mixed with brick, sawdust, and miscellaneous fill. Table 8 is a summary of data from RGA and ERT samples. Pesticides were found in the ERT sample; RGA sample was not \times analyzed for pesticides. No estimate was made as to the quantity disposed.

ERT, Inc. Detailed Description of Neville Island Site, August, 1981, Section 4.0.

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5. Dry ash is dry, powdery material appearing as a yellow and light gray ash, cinders, and black powdery carbon. ERT analysis demonstrated that it contains volatile organics and pesticides. Table 9 is a summary of ERT samples. No estimate was made as to the quantity disposed.

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- 6. Siag was disposed throughout the site. Analyses performed for ERT show the siag is contaminated with volatile organics and pesticides (see Table 10). ERT did not esimate the quantity of siag disposed at the site, but FCHA calculates the amount to be 90,400 cubic yards.
- 7. Crystalline waste of various types are on site. Only one type was identified. FCHA identified phthalic anhydride. The qualitative description of a white crystalline solid suggested naphthalene. No estimate made as to quantity disposed.
- 8. Leachate samples were collected by FCHA and ERT. See Table 11.
- 9. Calcium carbonate was mixed with the acid sludge in order to "neutralize" the low pH wastes (see Table 12). No estimate was made as to the quantity disposed.
- 10. Miscellaneous wastes have been identified as pitch and epoxy resins. No estimate has been made as to quantity disposed.
- Municipal wastes. Reports that 4 acres of the site was used for municipal waste.
- 12. Demolition waste. No estimate has been made as to quantity disposed.
- 13. Metal barrels reported to be empty or near empty. No estimate has been made as to the number of drums disposed on the site.

FCHA estimates that the total volume of waste disposed since 1941 was approximately 225,000 cubic yards.¹

Fred C. Hart Associates. Assessment of Remedial Options at Ohio River Park, January, 1980, p. 37.

- 17 -

5.0 FIELD TRIP REPORT

5.1 The field trip report is not applicable, as a desk top study was performed. The laboratory analyses submitted by ERT was reviewed according to EPA Functional Guidelines for Evaluating Inorganics and Organic Analyses.

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HAZARDOUS CONDITIONS AND INCO	ENTS				
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DI C 8. SURFACE WATER CONTAMINATION DI POPULATION POTENTIALLY AFFECTED:	10,000	02 COBSERVED IDATE	i _ 2 PG	TENTIAL (T ALLEGE
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1 2 2. FRE/EXPLOSIVE CONOTIONS		02 C OBSERVED IDATE:	, X: PO	TENTIAL	- ALLEGE
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Ohio River				-	02
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a number of suldings with	THO (2) MILES OF SITE		04 DISTANCE TO HE	WEST CFF-SITE BUILDIN	G
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Immediately east township. Acro Residential, Co	st of the site	hannel is the	borough of	Coraopolis.	rea of Neville Mixture of ty of the site.
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		POTENTIAL NAT	ARDOUS WASTES	172	I. DENTIFICATION
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VI. ENVIRONMENT	·				
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01 2 1 ACCESS TO SITE RESTRUCTED CA CESCRIPTION Chain link for * along shorel	ence from main channel to back chaine.	annel Vegetation planted	
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Zollinger. Moon Township Municipal Authority.

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6.0 LABORATORY DATA



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6.1 SAMPLE DATA SUMMARY

The following is the list of chemicals found in the Ohio River Park:

- Benzene, toluene, ethylbenzene, xylenes, phenol, 2-chlorophenol, 2,4-dichlorophenol, 2,4,6-trichlorophenol, naphthalene and 2,4-D have been confidently identified with levels greater than 3 times the detection limits and 5 times the background;
- (2) Diethylphthalate and bis(2-ethylhexyl)phthalate have been identified but with levels less than 3 times the detection limits;
- (3) Silvex and 2,4,5-T have been identified using one GC column with levels greater than 3 times the detection limits and 5 times the background;
- (4) Arsenic, beryllium, cadmium, chromium, copper, iron, lead, mercury, nickel, selenium, thallium and zinc were reported by ERT to be present on the site with levels greater than 5 times the background; and
- (5) Cyanide was detected on the site with levels greater than 5 times the background.

Among the chemicals listed above, benzene and arsenic are well-known human carcinogens (Group A). 2,4,6-Trichlorophenol, beryllium and lead are suspected human carcinogens (Group B2).

The attached data summary reports contain only compounds which were identified in at least one sample.

The following codes are used in the data summary reports to indicate the confidence in the laboratory results:

CODES RELATED TO IDENTIFICATION

(confidence concerning presence or absence of compounds):

U = Not detected. The associated number indicates approximate sample concentration necessary to be detected.

(NO CODE) = Confirmed identification.

- B = Not detected substantially above the level reported in laboratory or field blanks.
- R = Unreliable result. Analyte may or may not be present in sample. Supporting data necessary to confirm result.

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N = Tentative identification. Consider present. More analyses may be needed to confirm its presence or absence.

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D = Data reported by PRP's consultant. Supporting document necessary to confirm result.

CODES RELATED TO QUANTITATING

- J = Analyte present. Reported value may not be accurate or precise.
- K = Analyte present. Reported value may be biased high. Actual value is expected to be lower.
- L = Analyte present. Reported value may be biased low. Actual value is expected to be higher.

UJ = Not detected, quantitation limit may be inaccurate or imprecise.

6.2 QUALITY ASSURANCE REVIEW

6.2.1 ORGANIC DATA:

6.2.1.1 Introduction

Three ground water data packages (January 1981, April 1981, and October 1987), soil and waste data package (November 1981) and Outfall #1 data (included in the January 1981 data package) have been fully reviewed according to the EPA Functional Guidelines for the usability in the Site Inspection report. All data packages are not in the Contract Laboratory Program (CLP) deliverables format. This is understandable because ERT began the investigation of the site in early 1980's, prior to the initiation of CLP.

Since ERT did not have EPA CLP guidelines to follow, some of the data packages provided by ERT might be without some QC information required by present CLP guidelines. If a data package had this kind of problem, the reviewer would use his judgement based on the worst possible cases generally observed in the data package to qualify the data. For example, the January 1981 package had no spike data, the reviewer then assumed the spike recovery was out of QC limit and qualified all positive results with (J) flags and all results below the detection limits with (R) flags.

6.2.1.2 Qualifiers

6.2.1.2.a January 1981 Data Package

The January 1981 data package includes VOA, acid, and B/N results obtained using EPA methods 8240 and 8250 (GC/MS), as well as chlorinated herbicides data obtained using standard method 509B (GC).

Holding times, tuning, calibrations, surrogates, and compound identification have been fully reviewed for VOA, acid and B/N data according to the EPA Functional Guidelines. Problems in VOA's holding times, tuning, calibration, and surrogate recoveries of toluene-D8, nitrobenzene-D5, phenol-D6 and 2-fluorophenol were found.

- -- Some critical BFB and DFTPP ion abundance criteria were not met for some of the samples. The associated data have to be flagged (R) as unreliable.
- -- Since there were no spike, and spike/duplicate in this GC/MS package, and there was no evidence of gross contamination, all data which are above the detection limits are flagged (J) as analyte considered present with the estimated value. All other data below the detection limits are flagged (R) as unreliable results.
- -- Because no raw data come with the ERT-17S data report, all values for this sample are flagged (D) for requiring supporting document.

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ORIGINAL (Post)

Herbicides data have been fully reviewed for holding times, calibrations, and blanks. 2,4-D, silvex and 2,4,5-T were identified in many samples using one GC column and none were detected in the field blanks. However, problems in holding times and calibrations were noted. These might make all values above the detection limits be qualified with (J) flags, but since there was no confirmation experiment, such as second GC column analysis, all data above the detection limits have been flagged (N) instead of (J). All values below the detection limits are unreliable for the holding time violation.

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6.2.1.2.b April 1981 Data Package

The April 1981 data package also includes VOA, acid, B/N, and herbicides data obtained using methods 8240, 8250, and 509B.

For VOA, acid and B/N portion, holding times, tuning, calibrations, blanks, surrogates, and identification have been fully reviewed.

- -- The data which did not pass the tuning criteria are flagged (R) as unreliable.
- -- The benzene, chloroethane and 2,4-dimethylphenol results in ERT-18S, 2,4-dimethylphenol result in MW-4, and phenol result in MW-4 duplicate were flagged (UJ) because spectrum matching criteria were not met. Other positively identified VOCs in ERT-18S, and MW-4/MW-4 duplicate have been flagged (J) for surrogates being out of control limits. Acids in MW-4 duplicate and naphthalene in ERT-18S have been flagged (J) also for the surrogate problems.
- Naphthalene results for ERT-18S and ERT-20 have to be flagged (J) because of poor precision suggested by the ERT-17S duplicate analyses.
- -- The data of phenanthrene, anthracene, butylbenzylphthalate, and 3,3-dichlorobenzidine in ERT-20S have been flagged (UJ) because mass spectroscopic identification criteria were not met.
- -- The data of ERT-17S/ERT-17S duplicate have been flagged (D) for requiring the supporting raw data.

For herbicide analysis, 0.2 ug/L of 2,4-D was detected in the Blank #4. Since no spike/spike duplicate and surrogate were analyzed, all 2,4-D data which are less than the detection limits are unreliable. 2,4-D in ERT-16D has to be flagged (B) for its values being less than 1 ug/L, 5 times the 2,4-D value in the blank. 2,4-D and 2,4,5-T were identified with one GC column. However, no confirmation analyses were performed for herbicides data. Therefore, all other 2,4-D and 2,4,5-T data have been flagged (N).

6.2.1.2.c October 1987 Data Package

The October 1987 data package includes VOA (method 8240) and herbicides analysis (method 8150).

VOA has been fully reviewed for holding times, calibrations, blanks, surrogates, matrix spike/matrix spike duplicate, and compound identification according to the EPA Functional Guidelines. No major problems which might affect the overall usage of the VOA data package were found except the following minor point: the toluene and xylene (o,p) in sample ERT-2D do not meet the standard spectral criteria. Therefore, toluene is flagged (UJ) for not being detected. Total xylene has been identified, but is flagged (K) for its value could be overestimated.

For herbicides analysis, no information about the precision and accuracy was provided. Since there is no evidence of 2,4-D carryover to sample ERT-26, 2,4-D was positively identified in the mass spectrum and its quantification could be unreliable, it is flagged (J).

6.2.1.2.d Soil, Waste, and Leachate Data Package

The soil, waste and leachate data package (November 1980) includes 59 samples analyzed for benzene, toluene, and methylene chloride along with 89 samples analyzed for five different pesticides: 2,4-D, silvex, 2,4,5-T, parathion, malathion.

- -- Many of the volatile and herbicide raw data sheets, chromatograms and log-in record sheets could not be located by ERT. Therefore, all VOA data and large portion of pesticide data are without the supporting document and have been flagged (D).
- For pesticide portion, poor accuracy and poor precision were shown by the results of spike samples. Because the high relative percent difference (>25%) between the initial and continuing calibration response factors was also noted, the following samples have (UJ) flags in their 2,4,5-T and/or silvex data: B-9 (1.3'-2.2'), B-26(1.8'-2.4'), B-28(10.5'-11.5'), B-30(18.0'-19.8'), and B-31(6.0'-9.0').
- -- The parathion and malathion data of the following samples have been flagged (R) for exceeding the holding times: B-5(3.0'-3.4'), B-7(5'-10'), B-8(6.3'-6.7'), B-26(1.8'-2.4'), B-28(10.5'-11.5'), B-30(18.0'-19.8'), and B-31(6.0'-9.0').
- 2,4-D, silvex, 2,4,5-T and parathion have been identified using GC method. The data of the following samples have been flagged (N) because they were above the detection limits but were lack of the confirmation data. These include the 2,4-D data of B-9(1.3'-2.2'), B-26(1.8'-2.4'), B-28(10.5'-11.5'), B-30(18.0'-19.8'), B-31(6.0'-9.0'), TP-200(9.0'), TP-250(1.5'-2.5'), TP-250(9.0'-10.5'), TP-253(8.0'-10'), TP-303(3.5'-6.5'), and TP-303(6.5'-7.0'); the 2,4,5-T data of B-9(1.3'-2.2'), TP-200(9.0'), TP-250(1.5'-2.5'), TP-250(9.0'-10.5'), and TP-303(6.5'-7.0'); the silvex data of TP-200(9.0') and TP-250(9.0'-10.5'); and the parathion data of B-9(1.3'-2.2').
- -- The malathion data of B-9(1.3'-2.2') has a (R) flag for the sample exceeding the holding time.

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6.2.1.3 Summary of Organic Data

6.2.1.3.a Summary of the Groundwater Data Review:

The following compounds have been found and positively identified at least in one of the groundwater samples obtained during the period from January 1981 to October 1987: benzene, toluene, ethylbenzene, xylenes, phenol, 2-chlorophenol, 2,4-dichlorophenol, naphthalene, 2,4,6-trichlorophenol, bis(2-ethylhexyl)phthalate, diethylphthalate, and 2,4-D. Silvex and 2,4,5-T have been identified in 31 groundwater samples using GC method but have never been confirmed by a second GC column or GC/MS.

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6.2.1.3.b Summary of the Surface Water Data Review:

The presence of 2,4-dichlorophenol and 2,4,6-trichlorophenol in Outfall #1 (January 1981) is not confirmed because the DFTPP ion abundance criteria were not met and that made the identification very difficult. 2,4-D and 2,4,5-T in Outfall #1 have been identified using one column in the duplicate GC analyses. Although there was no confirmation column or GC/MS analysis, their presence in the surface water is supported by the existence of 2,4-D in the ground water and the fact that 2,4,5-T is one of the chemicals manufactured or formulated by PC&C.

6.2.1.3.c Summary of the Soil, Waste and Leachate Data Review:

The reviewer cannot validate the volatile data of the soil, waste, and leachate samples because ERT could not locate enough original document other than the raw data sheets. However, the existence of benzene and toluene in the soil and waste samples is supported by earlier results from FCHA report. 2,4-D, 2,4,5-T, and silvex have been identified with one GC column and their quantities reported in many soil and waste samples exceed the detection limits by 3 times. Although no confirmation analyses were performed by ERT, their presence in the soil and waste is also strongly supported by FCHA report.

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6.2.2 INORGANIC DATA:

One usable cyanide analysis data package (August 1980) which has been fully reviewed for the Site Inspection report. Included in this package are eleven samples, one duplicate, two lab blanks, one blind control sample, and one standard control sample. Acceptable precision was demonstrated by the field duplicate results, but poor accuracy is also suggested by the low recovery (76%) of the standard control sample. The lab miscalculated the concentrations of all samples. All reported values in the raw data sheets should be doubled. The corrected values are reported in the data summary. Since the holding times were exceeded for all samples and 0.031 mg/L of cyanide was detected in one of the blanks, all values below 0.155 mg/L are flagged (B) and values above 0.155 mg/L are flagged (J) as estimated. For sample 9210 (MW-3), the lab reported a wrong value (< 0.2 ug/ml) in the raw data sheet. The actual value should be 2.8 mg/L according to the original lab sheet.

Two sets of metal data (January 1981 and April 1981) have been reviewed. The raw data sheets of January 1981 were sent only with the copies of the ERT lab master logbook. No other QC information was provided. All data in this package have been flagged (D) for requiring more supporting documentation.

The April 1981 data was accompanied with the master logbook and QC information which included nineteen lab spikes, along with field duplicates, lab duplicates, blanks, and results of two check standards. All QC data provided by ERT met EPA criteria. According to the correspondence letter sent from Energy Resources Co., Inc. to ERT, samples 656, 974a, 794b, 795, and 834 had especially complex matrix effects and showed multiple peaks for Se. These samples were analyzed by standard addition of a diluted sample. However, no information about the calibrations, serial dilution, interference check sample analysis was provided by the lab for evaluation, all the April 1981 data have been flagged (D).

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6.3 TOXICOLOGICAL EVALUATION

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6.3.1 SUMMARY

Analyses of the groundwater samples collected from the on-site and offshore wells revealed that the groundwater was contaminated with at least twelve target organic compounds and three herbicides. Among these organic contaminants, human carcinogen benzene (up to 49,000 ppb), suspected human carcinogen 2,4,6-trichlorophenol (up to 50,000 ppb), the irritant and organ damaging phenol (up to 50,000 ppb), toxic 2,4-D (up to 25,700 ppb), and organ damaging silvex (up to 143 ppb) are of great concern. In addition, high levels of 2-chlorophenol (up to 5,400 ppb) and 2,4-dichlorophenol (up to 18,000 ppb) are of concern also because these levels are well above EPA water quality criteria.

With regard to inorganic contaminants, notable levels of the following have been found in the groundwater and are of concern because they exceed the enforceable Maximum Contaminant Levels (MCLs), Maximum Contaminant Level Goals (MCLGs) and other water quality criteria (WQC):

	Levels Found	MCLS	MCLGs	WQC
arsenic	up to 5,700 ppb	50 ppb	50 ppb	•
beryllium	up to 33 ppb			0.037 ppb
cadmium	up to 39 ppb	10 ppb	5 ppb	
chromium	up to 387 ppb	50 ppb	120 ppb	
1ron	up to 4,518,000 ppb	•••		300 ppb
nickel	up to 3,600 ppb			13.4 ppb
lead	up to 313 ppb	50 ppb	20 ppb	
selenium	up to 730 ppb	10 ppb	45 ppb	
thallium	up to 178 ppb			13 ppb
cyanide	about 2,800 ppb		· ·	200 ppb

Surface water contamination with 2,4-D (up to 2,500 ppb in Outfall No. 1), 2,4,5-T (up to 114 ppb in Outfall No. 1), bis(2-ethylhexyl)phthalate (93 ppb in Outfall No. 2), beryllium (1 ppb in Outfall No. 1), chromium (31 ppb in Outfall No. 1), nickel (17 ppb in Outfall No. 1), thallium (41 ppb in Outfall No. 1) and cyanide (192 ppb in Outfall No. 2) was reported by ERT. These contaminants have been identified or reported to be on the site. 2,4-D contamination of Outfall No. 1 is of concern because the level exceeds EPA drinking water standard of 100 ppb and NAS SNARL of 90 ppb. The levels of beryllium, chromium, nickel and thallium in Outfall No. 1 are of concern also because they exceed the MCLs, MCLGs, and other water quality criteria just mentioned.

The most significant pathway for potential off-site exposure is a route of leachate contaminating groundwater, groundwater entering the Ohio River, and human consumption of the treated river water via a public water supply system. The pathway of contaminated runoff entering the Ohio River and human ingestion of the treated river water is less significant than the groundwater contamination route.

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6.3.2 SUPPORT DATA

Benzene was identified in samples from four on-site wells (ERT-20, ERT-26, ERT-27, and MW-4) in concentrations that range from 680 ppb to about 49,000 ppb. Benzene was also reported by ERT to be present in other ten on-site wells (ERT-10, ERT-11, ERT-17, ERT-18, ERT-19, MW-1, MW-1A, MW-3, MW-6, and MW-5) in concentrations that range from about 13 ppb to about 67,000 ppb.² Although most of the wells contaminated with benzene are in the south section of the island, three wells (ERT-11, MW-3, and MW-6) in the north sector and two wells (MW-1, and MW-1A) in the east part of the site were also contaminated with benzene. These indicate general groundwater contamination with benzene in most areas of the site.

Since benzene is rather mobile in the environment and wells ERT-26 and ERT-27 are close to the perimeter of Neville Island, they can be assumed to approximate concentrations moving off the site. Benzene is a Group A human carcinogen.^{3,13} Benzene concentrations range at these two wells from 680 ppb to about 73,000 ppb and are from about 100 to about 7,000 times greater than the EPA water quality criterion of 6.6 ppb.⁴ This criterion is based on the estimate of one additional cancer for 100,000 people exposed on a lifetime exposure (1×10^{-5}) to drinking water at this elevated concentration. The potential adverse risks from the benzene-contaminated groundwater moving off the site from the south sector are very significant. One f the potential receptors is the Borough of Coraopolis. Water supply of the Borough of Coraopolis is from the well-field located about 700 feet southwest of the site across the back channel of the Ohio River.

Similar discussions are also applicable to the groundwater moving from the north sector of the island. Since the April 1981 data in the ERT interim monitoring report⁵ indicate the same, if not higher, order of benzene contamination in the ERT-11 well of the north sector as in those wells (ERT-18 and ERT-20) of south sector, and all these wells are close to the perimeter of island, the potential adverse risks from the benzene-contaminated groundwater moving off the site from the north sector should be very significant also.

<u>Toluene</u> was identified in samples from five on-site wells (ERT-18, ERT-20, ERT-26, ERT-27, and MW-4) and one offshore well (ERT-1) in concentrations that range from 10 ppb to about 11,000 ppb. Toluene was also reported by ERT to be present in other five on-site wells (ERT-11, ERT-17, MW-1, MW-1A, and MW-5).⁵ These indicate that toluene contamination is widely distributed at the site although high toluene concentrations (> 1,000 ug/L) are reported in those wells located in the western half of the south sector only.

Since well ERT-26 is close to the island perimeter, it is reasonable to assume that toluene may be moving off the site and entering the river in concentrations of the order of 2,800 to 8,600 ppb. The level is about 8 to 25 times greater than the NAS recommended SNARL for chronic exposure of 340 ppb,⁶ but is less than EPA's water quality criterion of 14,300 ppb.³⁴ Both criteria do include a safety factor of 1,000. In addition, toluene exposure does not appear to be linked with chronic risks such as carcinogenicity and mutagenicity. Off-site risk from toluene should therefore not be significant.

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Ethylbenzene was identified in two on-site wells (ERT-18 and MW-4) and also reported by ERT in two other on-site wells (ERT-11, and ERT-17).7 All these wells are in the west central part of the site. The area of ethylbenzene contamination is likely limited. The concentrations with maximum of about 120 ppb are several orders of magnitude below any expected adverse health effects and an order of magnitude below any recommended criteria. No adverse effects are expected to result from ethylbenzene off-site movement.

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<u>Xylene</u> (ortho, meta, and para) was identified in two on-site wells (ERT-18 and MW-4) and one offshore well (ERT-2). Xylene was also detected in wells ERT-11, ERT-17 and MW-5.⁸ The highest concentration occurred at ERT-18S, where the total xylene concentration was about 2,300 ppb. EPA has suggested a permissible concentration in water of 6,000 ppb based on health effects.⁹ Therefore, no significant effects are expected to result from xylene off-site movement.

<u>Phenol</u> was identified in three on-site wells (ERT-18, ERT-20, and MW-4). Phenol was existing in wells ERT-11, ERT-17 and MW-5 also according to the ERT interim monitoring report.¹⁰ The concentrations ranged from 200 ppb to greater than 50,000 ppb. Concentration of phenol is highest at shallow depth in the south-central area. To protect human health, the EPA water quality criterion for phenol is 300 ppb based on the organoleptic effects of chlorinated phenols inadvertently formed during water purification process.¹¹ Concentrations found at ERT-20 are in the range where some acute effects have been noted. It is of concern for phenol moving off the site at these concentrations.

2,4,6-Trichlorophenol was detected or reported by ERT in six on-site wells (ERT-11, ERT-17, ERT-18, ERT-20, MW-4 and MW-5).¹² The concentrations ranged from 120 ppb to greater than 50,000 ppb. ERT did not analyze the acid portion of the groundwater samples from wells ERT-26 and ERT-27 in the October 1987 monitoring program. Therefore, no data about 2,4,6-trichlorophenol in these two wells are available.

Since wells ERT-18 and ERT-20 are closer to south perimeter of island, these wells can be assumed to approximate closely the movement off site. 2,4,6-Trichlorophenol is suspected of being a human carcinogen (Group B2).13 For a risk corresponding to 1×10^{-5} (one additional lifetime cancer risk for 100,000 people exposed) a level of 12 ppb is considered.¹³ The measured groundwater concentration in ERT-20 exceeds this criterion by more than 4,000 fold. Drinking water contaminated with this level of 2,4,6-trichlorophenol would present a significant adverse health risk.

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<u>2-Chlorophenol</u> was identified in three on-site wells (ERT-18, ERT-20, and MW-4). In addition, 2-chlorophenol was detected by ERT at ERT-11, ERT-17, and MW-5.14 The highest concentration was in ERT-20 where about 5,400 ppb of 2-chlorophenol was detected. EPA water quality criterion of 0.1 ug/L for 2-chlorophenol was set due to its low odor threshold in water and its tainting properties.¹⁵ This criterion is about fifty thousandth of the highest concentration detected on the site. 2-Chlorophenol in the ground water of the site is of concern, although 2-chlorophenol has been reported to be less toxic than the higher chlorophenols.

2.4-Dichlorophenol was identified in three on-site wells (ERT-18, ERT-20, and MW-4). ERT also reported that four other wells (ERT-11, ERT-17, MW-5, and MW-6) had 2.4-dichlorophenol contamination.¹⁴ The highest concentration of 2.4-dichlorophenol was reported at well ERT-18, where 18,000 ppb was detected. This concentration is more than five times the EPA water criterion of 3,090 ppb based on toxicity data¹⁶ and sixty thousand times the EPA criterion of 0.3 ppb based on organoleptic effects.¹⁶ Although 2.4-dichlorophenol is less toxic than the higher chlorinated phenols, its toxicity to certain microorganisms and plant life has been demonstrated and its tumor promoting potential in mice has been reported. Also it can irritate tissue and mucous membranes.¹⁶ 2.4-Dichlorophenol in the ground water of the site is of concern. The 2.4-dichlorophenol is likely to result from 2.4-D decomposition and chlorinated phenols disposed of on the site.

<u>2,4-Dimethylphenol</u> was reported to be present in ERT-17 and M-4.¹⁷ However, its presence in samples taken from ERT-18 and MW-4 in April 1981 could not be confirmed because the GC/MS tuning criteria for the April, 1981 data were not met. The highest concentration of 2,4-dimethylphenol reported for the site is 500 ppb which is a little bit higher than the water quality criterion of 400 ppb based on the organoleptic effects.¹⁸ Little human health data are available and no health-based criteria exist. The potential risks associated with exposure to deimethylphenol are unknown but expected to be small because of its limited on-site quantity.

<u>Naphthalene</u> was identified in three on-site wells (ERT-18, ERT-20, and MW-4). Three other wells (ERT-11, ERT-17, and ERT-19) also were reported by ERT to have naphthalene contamination.¹⁹ The highest concentration of naphthalene, was about 410 ppb, which is less than the short term (1 day and 10 days) Health Advisory of drinking water of 5,300 ppb for naphthalene, as well as the long term, 10-Kg Health Advisory of 5,300 ppb, and the long term, 70-Kg Health Advisory of 18,600 ppb.²⁰ Therefore, the naphthalene in the groundwater of the site is not expected to present a significant risk.

<u>Bis(2-ethylhexyl)phthalate (BEHP)</u> was identified in ERT-20, and was also reported to be present in wells ERT-10, ERT-11, ERT-12, ERT-16, ERT-17, MW-2 and MW-5, along with one Outfall (No. 2).²¹ In all samples, the concentrations were less than 100 ppb. Because the EPA water quality criterion is 15,000 ppb,²² no adverse off-site risk for BEHP is expected.

<u>Diethylphthalate</u> was identified in only one sample from ERT-18 in concentration of 28 ppb. Since the recommended water quality criterion level for protection of human health is 350,000 ppb,¹ the limited quality and low concentration of diethylphthalate on the site should pose no adverse health effects.

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2,4-D was identified with GC/MS in samples from two on-site wells (ERT-26 and ERT-27)³³ and was also identified utilizing one GC column in samples from one offshore well (ERT-3), five on-site wells (ERT-10, ERT-11, ERT-17, ERT-18, and MW-4), and Outfall No. 1 at concentrations greater than 10 ppb. 2,4-D was also identified in samples from eight other wells (ERT-6, ERT-8, ERT-9, ERT-12, ERT-14, ERT-19, ERT-20, and ERT-32) in concentrations that ranged from the detection limit of 1 ppb to 10 ppb. From the data, it appears that 2,4-D is a significant groundwater contaminant widespread on the site.

The maximum concentration detected in the effluent from Outfall No. 1 was about 2,500 ppb. The maximum concentrations detected in ERT-3, ERT-26, and ERT-27 were about 216, 410, and 16 ppb respectively. These values may indicate the concentration level at which 2,4-D moves off-site. The highest concentration detected in the groundwater samples was from ERT-18 and was about 25,700 ppb.

The EPA drinking water standard for 2,4-D is 100 ppb^{23} and the NAS SNARL is 90 $ppb.^6$ Since these criteria are several hundredth the maximum level detected in the groundwater of ERT-18, 2,4-D is of concern for its moderate acute toxicity. No chronic hazards have been shown to be associated with 2,4-D.

2,4,5-T was reported to be present in two offshore wells (ERT-3 and ERT-6), eight on-site wells (ERT-8, ERT-11, ERT-14, ERT-17, ERT-18, ERT-20, MW-4, and MW-5) and Outfall No. 1.24 The reported highest concentration of 2,4,5-T in groundwater was about 266 ppb in sample from MW-4. The effluent from Outfall No. 1, had about 114 ppb of 2,4,5-T. From this information, 2,4,5-T seems to be located over much of the island.

2,4,5-T is an animal suspected carcinogen.²⁵ The EPA Health Advisory for short term exposures (one day and ten days) is 800 ppb. The long term Health Advisories for a 10-Kg child and a 70-Kg adult are 300 ppb and 1,050 ppb respectively.²⁶ In addition, a no-adverse-effect-level in drinking water has been calculated by NAS to be 700 ppb.⁶ Therefore, 2,4,5-T should not present any significant adverse risk off-site.

 $\frac{\text{Silvex (2,4,5-TP)}}{\text{on-site wells (ERT-9, ERT-11, ERT-17, ERT-18, ERT-19, ERT-20, MW-4, and MW-5) with concentrations up to about 143 ppb. The concentration at MW-4 (143 ppb) is about 15 times the EPA drinking water MCL of 10 ppb²⁷ and about 30 times the NAS no-adverse-effect-level of 5.25 ppb.⁶ Samples taken from two other wells (ERT-11 and ERT-18) in April 1981 also exceeded these criteria. Off-site risks due to silvex contaminated groundwater movement are of concern.$

<u>Metal analysis</u> of January and April 1981 water samples revealed notable levels of arsenic (up to 5,700 ppb), beryllium (up to 33 ppb), cadmium (up to 39 ppb), chromium (up to 387 ppb), copper (up to 1,480 ppb), iron (up to 4,518,000 ppb), mercury (up to 2.2 ppb), nickel (up to 3,600 ppb), lead (up to 313 ppb), selenium (up to 730 ppb), thallium (up to 178 ppb), and zinc (up to 19,000 ppb). These levels exceed the enforceable Maximum Contaminant Levels [MCLs (arsenic, 50 ppb; cadmium, 10 ppb; chromium, 50 ppb; lead, 50 ppb; mercury, 2 ppb; and selenium, 10 ppb)],²⁷ Maximum Contaminant Level Goals [MCLGs (arsenic, 50 ppb; cadmium, 5 ppb; chromium, 120 ppb; copper, 1,300 ppb; lead, 20 ppb; and selenium, 45 ppb)], and other water quality criteria (beryllium, 0.0037 ppb; iron, 300 ppb; nickel, 13.4 ppb; thallium, 13 ppb; and zinc, 5,000 ppb).²⁸

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Among these metals, copper, mercury and zinc are of less concern because only one or two samples exceeded criteria mentioned above and their quantities might be limited on the site. Arsenic is of great concern because it is a Group A human carcinogen.^{13,29} The unit risk of arsenic through the oral route is 5 x $10^{-5}/ug/L$,³⁰ i.e. 5 additional lifetime cancer risk for 100,000 people exposed to the potable water with arsenic contamination of 1 ug/L. Approximately 285 additional cancer cases for 1,000 people would occur on exposure to 5,700 ppb arsenic contaminated potable water. Since both beryllium and lead are Group B2 carcinogens, ^{13,31} they are also of great concern.

The area of the most serious metal-contamination of the groundwater seems to be the south-central section of the site because samples from either ERT-18 or MW-4 were consistent in showing the highest levels of all metals just mentioned before except mercury and iron. The highest level or iron was found in sample from MW-6 which is located in the central portion of the site.

The metal analysis of Outfall No. 1 sample taken during the January 1981 sampling round indicated notable levels of beryllium (1 ppb), chromium (31 ppb), nickel (17 ppb), and thallium (41 ppb). All these levels exceed the MCLs, MCLGs or other water quality criteria as mentioned before.

<u>Cyanide</u> was detected in samples taken from MW-3 (about 2,800 ppb) and Outfall No. 2 (about 192 ppb) in August 1980. To protect human health, the EPA has set the water quality criterion at 200 ppb. To protect freshwater aquatic life, maximum of 52 ppb is set.³² Comparing with these criteria, the levels of cyanide in these two water samples are of concern.

Report prepared by

Date:

Sam Fang, Environmental Chemist I

Report reviewed by

Date:

Joe Carpentier, Environmental Chemist II

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Mr. James R. Shack May 16, 1988 Page Nine

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o Herbicide raw data sheets, QC data, chromatograms

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- Volatile raw data sheets and chromatograms. Chromatograms for all analyses could not be located. The GC recorder paper used for this analysis has deteriorated with time, so some of the runs may have been discarded as no longer legible.
- No information other than a summary data sheet could be located for B-13, 13'-13.6' depth.

Sampling Procedures

Sampling procedures followed at the Neville Land Company Site have been consistent with industry standards for this type of work. Beginning in 1983, ERT developed Standard Operating Procedures (SOPs) for soil and water sampling. These SOPs, which are listed in Table 1 and included as Attachment 7, have been followed for all sampling activities since their issuance. Sampling procedures prior to SOP development were not documented in detail, but the SOPs adequately reflect the methods utilized. Some specific details of sampling procedures followed at the site prior to SOP development are presented below.

Ground Water Sampling

Two types of monitoring wells exist at the site: BarCad wells and conventional monitoring wells. BarCads are check valve-type sampling instruments that deliver slugs of water as the system is pressurized, but they cannot deliver a continuous stream of water. Therefore, each time a BarCad is pressurized, one "well volume" is delivered to the surface. Ultra High Purity (UHP) Grade 5 Nitrogen was used to pressurize the BarCads, and samples were collected after the BarCads had been purged three times.

Conventional monitoring wells were purged using bailers, a peristaltic pump; and/or a Johnson-Keck pump. Wells were purged until a minimum of three well volumes were removed. All VOC samples were collected using bailers, while some samples for other parameters were collected directly from the pump discharge. Samples for metals analysis were filtered in the field using 0.45 micron filters. Disposable filters or filters in reusable polycarbonate housing were used.

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Surface Water Sampling

The Ohio River was sampled from an anchored boat using a peristaltic pump and a weighted length of tubing. Samples were collected by compositing water from 1.0 meter off the bottom. the mid-depth point, and 1.0 meter below the surface.

Outfalls were sampled during periods of very low flow in the outfalls. This occurred after a rainfall event, in order to get the "first flush," and also several days after rain, as the discharge slowed to just a trickle. All samples from outfalls were grab samples collected at the end of the pipe. No compositing was done.

Soil and Waste Sampling

Samples were collected from soil borings, as described in ERT's SOP 7115. Split spoons of two- and three-foot lengths were used, as well as a five-foot soil core barrel sampler. Stainless steel utensils were used to handle the soil.

A backhoe was used to excavate numerous test trenches at the site. Many of the trenches provided access for the field team to collect samples directly from the walls of the trench. Other trenches could not be entered, and samples were collected either from the bucket of the backhoe, or by using a long-handled shovel to sample the walls of the trench. Stainless steel utensils and disposable plastic scoops were used to handle the soil samples.

I hope that this information satisfies your needs. Please call me if you have any questions.

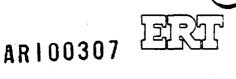
Sincerely,

Bos Kalmeyer

Robert W. Rittmeyer, P.E. Manager Pittsburgh Operations

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cc: M. J. Laskow M. Ferlin



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TABLE 1

SOIL AND WATER SAMPLING SOPS

Number	Title
7110	Surface Soil Sampling
7115	Subsurface Soil Sampling
7120	Surface Water Sampling Techniques
7130	Ground Water Sample Collection from Monitoring Wells
7131	Field Filtration of Water Samples for Inorganics
7220	Monitoring Well Construction and Installation
7230	Test Pits/Trench Subsurface Exploration
7315	Operation/Calibration of HNu Photoionization Analyzer
7320	Calibration and Operation of Hydrolab Water Quality Monitor
7510	Packaging and Shipment of Samples
7600	Decontamination of Equipment



Table 4. continued

Outfall No. 2 (12/13/79) ļ 20 < 200 ۸ 00 00 30 290 37 M M Ę 늘 MT Outfall No. 1 (12/13/79) CONCENTRATION: (ug/1) 2020 3,400 40 18 F ž Ĕ M ħ V • Outfall No. 1 (10/26/79) 20 335,000 90 V 40 180 2 9 220 23 V Ponded Water (12/13/79) 20 V < 200 00 V 100 40 눋 ž NT N Z Ponded Water (10/26/79) < 20 5,500 00 V <u> 006</u> 27 v <u>Metals and Other Organics</u> (continued) Phenolic Cpds (as phenol) Volatile Hydrocarbons 2.4.6-trichlorophenol Parameter 2,4-dichlorophenol 2-chlorophenol (IT) muillent Silver (Ag) 4-nitrophenol Sulfide (S) Zinc (Zn) **Phenols** Benzene Toluene Phenol

AR100309

Сл. П.,

March 15, 1988

4. Operators What are the names, addresses, phone numbers and contact person names of all operators of the site? What is the name of their parent company? What was the time period that they operated the site?

E.

-2-

- 5. What is the relationship between Neville Land Company, Pittsburgh Coke & Chemical, and Hillman Company?
- 6. What happened to MW-2A that made it inoperable?
- 7. Has the site received or applied for any state, county or federal permits? Has the site ever had an effect on any water supply, either private or public?
- 8. Where are the records pertaining to the site, generator, users, transporters, operators and/or owners retained?

As we agreed during our conversation of March 8, 1988, the Department would be willing to accept less than all the QA/QC qualifiers (especially. Organics and Pesticides) as requested in the Department's letter of February 29, 1988.

After discussion with the U. S. EPA, the following is a list of the minimum sample results with QA/QC qualifiers would be needed to conduct the SI:

1. Groundwater results for:

E. R. T.

ERT-1	January	29,	1981	and	April	20,	1981 -
ERT-21-S	January	29,	1981	and	Apr11	20,	1981
MW-4	January	29,	1981.	and	April	20,	1981-
ERT-18-S	January	29,	1981	and	Apr11	20,	1981—
ERT-20-S	January	29,	1981	and	April	20,	1981 -
ERT-17-S	January	29,	1981	and	April	20,	1981

2. Surface Water Samples

Outfall #2	August,	1980
Upstream Sample	August,	1980
Downstream Samples	August,	1980

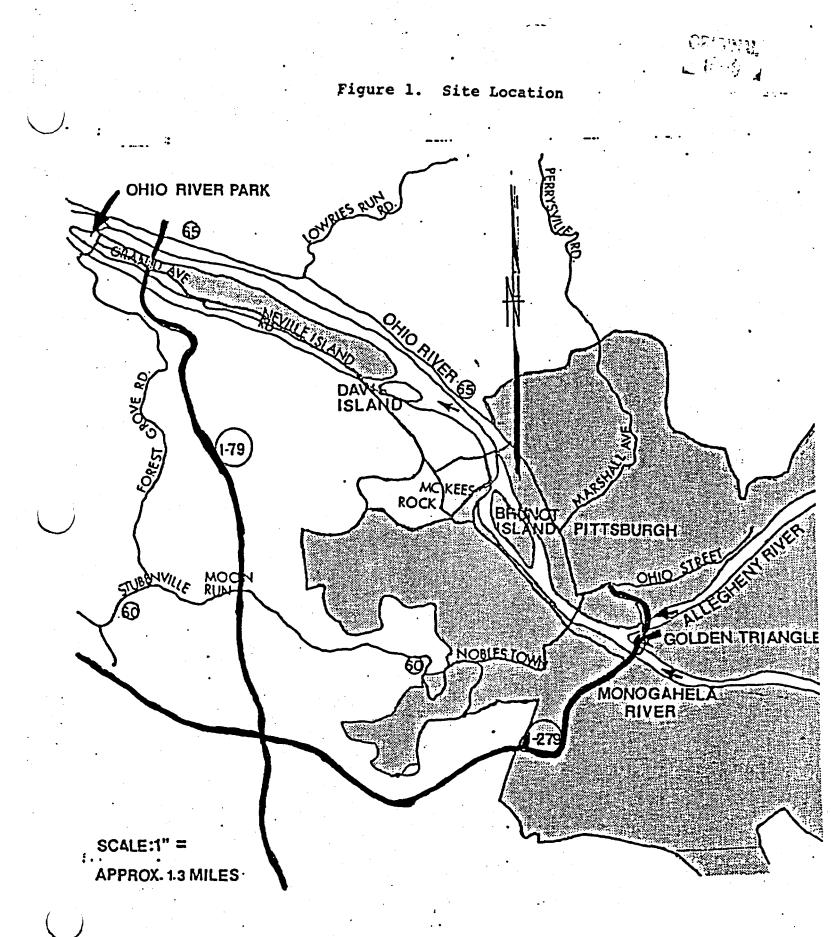
Outfall #1	January,	1981
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3. Soil/Waste:

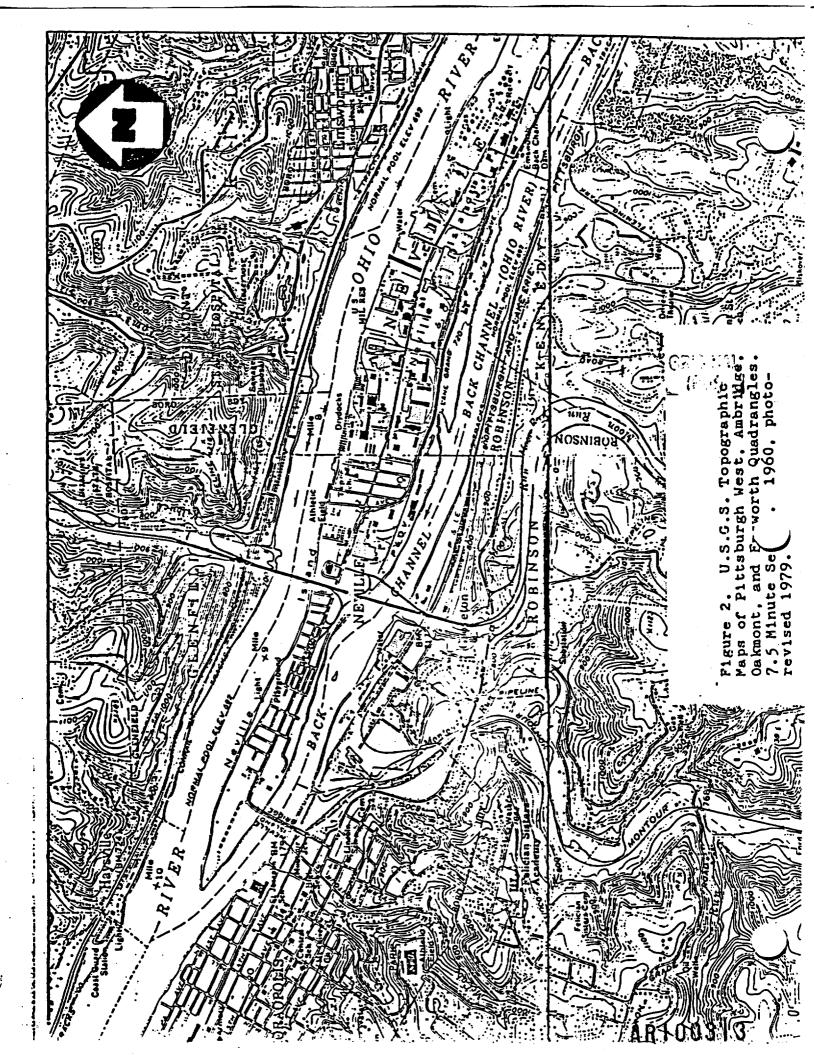
TP275	Drum, 1' depth
B-28	10' - 11.5' depth
B-30	30.0' - 13.6' depth
8-5	3.0' - 3.4' depth
TP-200	Liquid
B-9	1.3' - 2.2' depth

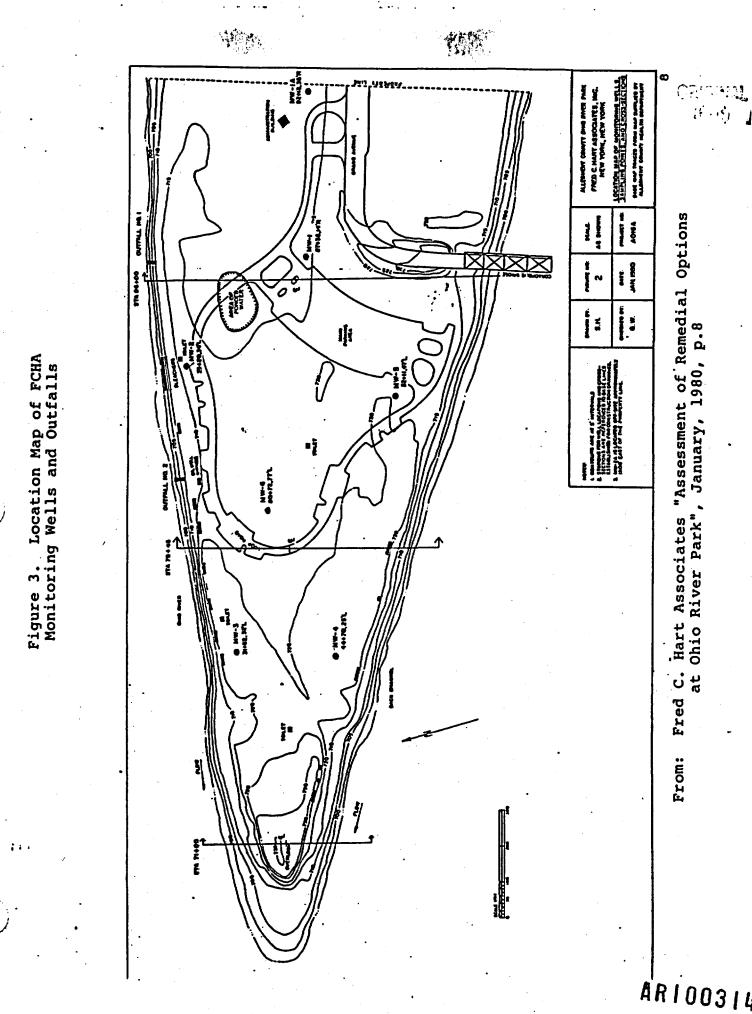
Appendix B

AR100311



From: Fred C. Hart Associates "Assessment of Remedial Options at Ohio River Park", Januarv, 1980, p.2. ARIOO312





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4T-Compucher **16** 1 .1

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ANALYSES PERFORMED :

VCA 2/10/81 ACID 2/15/81 B/N 2/05/81 PEST 2/05/81

Conto HAL

SAMPLE IDENTIFIER: 8074 Compu/Chem Sample Number: 3833

3. PRIORITY POLLUTANT ANALYSIS REPORT

•		COMPOUNDS	CONCENTRATION (UG/L)	DETECTION LIMIT (UG/L)
	1V.	ACROLEIN	BOL	R 100
	2V.	ACRYLONITRILE	BOL	t 100
		BENZENE	BOL	R 10
• •	4V.	BIB (CHLOROMETHYL) ETHER	8DL	R 10
	SV.		BOL	R 10
	6V.		BOL	R 10
	7V.		BDL	R 10
	8V.		BOL	R 10
	9V.		BDL	R 10
• .		2-CHLOROETHYLVINYL ETHER	BDL	R 10
		CHLOROFORM	BDL	2 10
		DICHLOROBROMOMETHANE	BDL	3 10
		DICHLORODIFLUOROMETHANE	BDL	2 10
		1, 1-DICHLOROETHANE	BDL	- 10
	_	1, 2-DICHLOROETHANE	BOL	L 10
		1, 1-DICHLOROETHYLENE	BDL	10
	4 711		BDL	R 10
	18V.			R 10
	19V.	ETHYLBENZENE	BOL	R 10
	20V.	METHYL BROMIDE	BDL	R 10
		METHYL CHLORIDE	BDL	Ř 10
		METHYLENE CHLORIDE	BDL	R 10
		1, 1, 2, 2-TETRACHLOROETHANE	BOL BOL BDL BDL BDL BOL	R 10
		TETRACHLOROETHYLENE	BDL	12 10
				R 10
	267.	TOLUENE 1,2-TRANS-DICHLOROETHYLENE 1,1,1-TRICHLOROETHANE 1,1,2-TRICHLOROETHANE	BDL	R 10
	27V.	1, 1, 1-TRICHLOROETHANE	BDL	R 10
	28V.	1, 1, 2-TRICHLOROETHANE	BDL	r 10
	29V.	TRICHLOROETHYLENE	BDL	R 10
÷	30V.	TRICHLOROFLUOROMETHANE	BOL	R 10
	31V.	VINYL CHLORIDE	BOL	R 10
	1A.	2-CHLOROPHENOL	BDL	°R 25
	2A.	2, 4-DICHLOROPHENOL	BOL	r. 25
	3A.	2, 4-DIMETHYLPHENOL	BDL	25
	4A.	4.6-DINITRO-O-CRESOL	BOL	- 250
	5A.	2, 4-DINITROPHENOL	BOL	r. 250
	6A.	2-NITROPHENOL	BDL	· 25
•	7A.	4-NITROPHENOL	BOL	<u>r.</u> 25
	8A.	P-CHLORO-M-CRESOL	BDL	- 25
	9Å.	PENTACHLOROPHENOL	BOL	25
	10A.	PHENOL	BDL	· 25
•	11A .		BDL	25
		ACENAPHTHENE	BDL	× 10
	2B.		BDL	10
	38.	ANTHRACENE	BDL	10

BOL- BELOW DETECTION LIMIT

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SAMPLE IDENTIFIER: 8074 Compu/Chem Sample Number: 3833

	1	COMPOUNDS BENZIDINE BENZO (A) ANTHRACENE BENZO (A) PYRENE 3,4-BENZOFLUORANTHENE BENZO (GHI) PERYLENE BENZO (K) FLUORANTHENE BIS (2-CHLOROETHOXY) METHANE DIE (2-CHLOROETHOXY) METHANE	CONCENTRATION (UC/L)	DETECTION LIMIT (UG/L)
	4B.	BENZIDINE	BDL	K 10
	58.	BENZO (A) ANTHRACENE	BDL	
	68.	BENZO (A) PYRENE	BOL	R 10
	7B.	3, 4-BENZOFLUORANTHENE	BDL	
	8B.	BENZO (GHI) PERYLENE	BDL	R 25
	98 .	BENZO (K) FLUORANTHENE	BDL	R 10
	10B.	BIS (2-CHLOROETHOXY) METHANE	BDL	K 10
	118.	BIS (2-CHLOROETHYL) ETHER	BDL	r 10
	129.	BIS (2-CHLOROISOPROPYL) ETHER	BDL	R 10
• .	138.	BIS (2-ETHYLHEXYL) PHTHALATE	BDL	R 10
	148.	4-BROMOPHENYL PHENYL ETHER	BDL	r 10
	15B.	BUTYL BENZYL PHTHALATE	BDL	R 10
	16B.	2-CHLORONAPHTHALENE	BOL	R 10
	178.	4-CHLOROPHENYL PHENYL ETHER	BDL	r 10
	188.	CHRYSENE	BDL	弐 10
	19B.	BIS (2-CHLOROETHYL) ETHER BIS (2-CHLOROISOPROPYL) ETHER BIS (2-ETHYLHEXYL) PHTHALATE 4-BROMOPHENYL PHENYL ETHER BUTYL BENZYL PHTHALATE 2-CHLORONAPHTHALENE 4-CHLOROPHENYL PHENYL ETHER CHRYSENE DIBENZO (A, H) ANTHRACENE 1, 2-DICHLOROBENZENE	BDL	r. 25
-	20B.	1, 2-DICHLOROBENZENE	BDL	₹ 10
	218.	1, 3-DICHLOROBENZENE	BDL	R 10
	228.	1. 4-DICHLORBENZENE	BDL	R 10
	23B.	3, 3'-DICHLOROBENZIDINE	BDL	r. 10
	248.	DIETHYL PHTHALATE	BDL	r. 10
	258.	CHRYSENE DIBENZO (A, H) ANTHRACENE 1. 2-DICHLOROBENZENE 1. 3-DICHLOROBENZENE 3. 3'-DICHLOROBENZIDINE DIETHYL PHTHALATE DIMETHYL PHTHALATE DI-N-BUTYL PHTHALATE 2. 4-DINITROTOLUENE 2. 6-DINITROTOLUENE DI-N-OCTYL PHTHALATE	BOL	r. 10
	268.	DI-N-BUTYL PHTHALATE	BDL	10
	278.	2, 4-DINITROTOLUENE	BDL	- 10
	28B.	2. 6-DINITROTOLUENE	BDL	10
	298.	2.6-DINITROTOLUENE DI-N-OCTYL PHTHALATE 1.2-DIPHENYLHYDRAZINE FLUORANTHENE FLUORENE HEXACHLOROBENZENE HEXACHLOROBUTADIENE HEXACHLOROCYCLOPENTADIENE HEXACHLOROETHANE	BDL	r 10
	30B.	1, 2-DIPHENYLHYDRAZINE	BDL	- 10
	318.	FLUORANTHENE	BDL	R 10
	328.	FLUORENE	8DL	r. 10
	33B.	HEXACHLOROBENZENE	BDL	r 10
	348.	HEXACHLOROBUTADIENE	BOL	R 10
	358.	HEXACHLOROCYCLOPENTADIENE	BDL	r 10
	368.	HEXACHLOROCYCLOPENTADIENE HEXACHLOROETHANE INDENO (1.2,3-CD) PYRENE IBOPHORONE NAPHTHALENE NTTROBENZENE	BOL	R 10
	378.	INDENO (1.2.3-CD) PYRENE	BDL	R 25
	389.	IBOPHORONE	BDL	r 10
	378.	NAPHTHALENE	BDL	R 10
	40B.	NITROBENZENE	BDL	R 10
	418.	N-NITROSODIMETHYLAMINE	. BDL	R 10
•	428.	N-NITROSODI-N-PROPYLAMINE	BDL	r 10
	43B.	N-NITROSODIPHENYLAMINE	BOL	R 10
	449.	PHENANTHRENE	BDL	R 10
	458.	PYRENE	BDL	R 10
	468.		BDL	R 10
	1P.		BDL	R 10
	2P.	ALPHA-BHC	BDL	r 10
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EDL= BELOW DETECTION LIMIT

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SAMPLE IDENTIFIER: 8074 COMPU/CHEM SAMPLE NUMBER: 3833

	COMPOUNDS		CONCENTRATION (UG/L)		DETECTION LIMIT (UG/L)
3P.	BETA-BHC		BOL	Я	10
4P.	GAMMA-BHC		801	R	10
5P.	DELTA-BHC		BOL	R	10
6P.	CHLORDANE	•	BOL	8	10
7 P .	4, 4'-DDT		BOL	R	. 10
8 P .	4, 4'-00E		BOL	R	10
9P.	4.4-000		BOL	R	10
10P.	DIELORIN		BOL	R	10
11P.	ALPHA-ENDOSULFAN		BOL	R	10
12P.	BETA-ENDOSULFAN		BOL	R	10
13P.	ENDOSULFAN SULFATE		BDL	R	
149.	ENDRIN	•	BOL	R	10
15P.	ENDRIN ALDEHYDE		BOL	R	
16P.	HEPTACHLOR		BOL	ĸ	
17P.	HEPTACHLOR EPOXIDE		BOL	R	
18P.	PC8-1242		BOL	R	10
19P.	PCB-1254		BOL	R	-
20P.	PCB-1221		BOL	R	
21P.	PCB-1232	•	BDL	R	10
22P.	PCB-1248		BDL	12	10
23P.	PCB-1260		BDL.	K	
24P.	PCB-1016		BOL		
25P.	TOXAPHENE		BDL		

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ERT ID 808 Compuchem 3820

ANALYSES PERFORMED :

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VQA	2/09/81 4
ACID	2/04/81
B/N	205/81
PEST	2/05/81

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SAMPLE IDENTIFIER: 8086 COMPU/CHEM SAMPLE NUMBER: 3826

3. PRIORITY POLLUTANT ANALYSIS REPORT

	COMPOUNDS	CONCENTRATION (UG/L)	DE	TECTION LIMIT (UG/L)	
1V.	ACROLEIN	BDL	R	100	•
27.	ACRYLONITRILE	BDL	R	100	
37.	BENZENE	BDL	R	10	
4V.	BIS (CHLOROMETHYL) ETHER	BOL	R	10	
5V.	BROMOFORM	BOL	R	10	
6V.	CARBON TETRACHLORIDE	BDL	R	10	
7V.	CHLOROBENZENE	BOL	R	10	
8V.	CHLORODIBROMOMETHANE	BOL	R	- 10	
9V.		80L	R	10	
10V.	CHLOROETHANE 2-CHLOROETHYLVINYL ETHER CHLOROEORM	BOL	R	10	
. 117.	CHLOROFORM	BOL	2	10	
127.	DICHLOROBROMOMETHANE	BOL	R	10	
	DICHLORODIFLUOROMETHANE	BOL	r.	10	
	1, 1-DICHLOROETHANE	BOL	R	10	
	1, 2-DICHLOROETHANE	BOL	R	10	
	1, 1-DICHLOROETHYLENE	BDL	R	10	
17V.	1, 2-DICHLOROPROPANE	BOL	R	10	1
18V.	1. 3-DICHLOROPROPYLENE	BOL	R	10	
19V.	ETHYLBENZENE	BOL	R	10	
20V.	METHYL BROMIDE.	BOL	R	10	
21 V.	METHYL CHLORIDE	BDL	R.	10	
22V.		BOL	Ŕ	10	
237.	1, 1, 2, 2-TETRACHLOROETHANE	BOL	12	10	•
244.	TETRACHLOROETHYLENE	BDL	R	10	
25V.	TOLUENE	BDL	2	10	
26V.	1, 2-TRANS-DICHLOROETHYLENE 1, 1, 1-TRICHLOROETHANE 1, 1, 2-TRICHLOROETHANE TRICHLOROETHYLENE TRICHLOROETHYLENE	BDL	-	10	
27V.	1, 1, 1-TRICHLOROETHANE	BDL	ĸ	10	
28V.	1, 1, 2-TRICHLOROETHANE	BOL	R	10	
297.	TRICHLOROETHYLENE	BOL	R	10	
30V.			R	. 19	
31V.	VINYL CHLORIDE	BÓL	R	10	
1A.	2-CHLOROPHENOL	BOL	R	25	
2A.	2. 4-DICHLOROPHENOL	63	2	25	•
3A.	2. 4-DIMETHYLPHENQL	BDL	R	25	
4 A .	4. 6-DINITRO-O-CRESOL	BOL	R	250	
5A.	2, 4-DINITROPHENOL	BDL	R	250	
6A.	2-NITROPHENOL	BDL		25	
7A.	4-NITROPHENOL	BOL	R	25	
BA.	P-CHLORO-M-CRESOL	BDL	R	25	
9A.	PENTACHLOROPHENOL	BDL	R	25	
10A.	PHENOL	BOL	R	25	
11A.	2, 4, 6-TRICHLOROPHENOL	37	しょ	25	١
18.	ACENAPHTHENE	BDL		10	``
28.	ACENAPHTHYLENE	BDL	Ŕ	10	
3B.	ANTHRACENE	BOL	R	10	
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SAMPLE IDENTIFIER: 8086 COMPU/CHEM SAMPLE NUMBER: 3826

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48. BENZIDINE BDL R 10 55. BENZO (A) ANTHRACENE BDL R 10 68. BENZO (A) FURENE BDL R 10 68. BENZO (CH) FURENE BDL R 10 68. BENZO (CH) FLUORANTHENE BDL R 10 68. BENZO (CH) FLUORANTHENE BDL R 10 108. BENZO (CH) FLUORANTHENE BDL R 10 118. BIS (2-CHLOROETHYL) ETHER BDL R 10 128. BIS (2-CHLOROETHYL) ETHER BDL R 10 128. BIS (2-CHLOROETHYL) PHTHALATE BDL R 10 128. BIS (2-CHLOROETHYL) PHTHALATE BDL R 10 128. BIS (2-CHLOROETHYL) PHENYL ETHER BDL R 10 128. BIS (2-CHLOROBENZU PHENYL ETHER BDL R 10 128. DITENZO (A, H) ANTHRACENE BDL R 10 128. 1. 3-DICHLOROBENZENE BDL R 10 128. 1. 3-DICHLOROBENZENE BDL R 10 228. 1. 4-DICHLOROBENZIDINE BDL R 10 238. 3.3'-DICHLOROBENZIDINE BDL R 10 249. DIETHYL PHTHALATE BDL R 10 238. JI-DICHLOROBENZIDINE BDL R 10		COMPOUNDS	CONCENTRATION (UG/L)	DETECTION LIMIT (UG/L)
5B. BENZO (A) ANTHRACENE BDL R 10 6B. BENZO (A) PYRENE BDL R 10 7B. 3. 4-BENZOFLUORANTHENE BDL R 10 8B. BENZO (GHZ) PERYLENE BDL R 10 9B. BENZO (GHZ) PERYLENE BDL R 10 108. BIS (2-CHLOROETHOXY) METHANE BDL R 10 118. BIS (2-CHLOROETHOXY) METHANE BDL R 10 128. BIS (2-CHLOROETHOXY) ETHER BDL R 10 128. BIS (2-CHLOROETHOYL) ETHER BDL R 10 138. BIS (2-CHLOROETHYL) PHTHALATE BDL R 10 148. 4-BROMOPHENYL PHENYL ETHER BDL R 10 158. BUTYL BENZYL PHTHALATE BDL R 10 160. 2-CHLOROBENZENE BDL R 10 1618. CHRYSENE BDL R 10 178. 4-CHLOROBENZENE BDL R 10 189. DIBENZO (A.H) ANTHRACENE BDL R 10 198. DIBENZO (A.H) ANTHRACENE BDL R 10 208. 1.2-DICHLOROBENZENE BDL R 10 210. CHLOROBENZENE BDL R 10 228. J.4-OZCHLOROBENZENE BDL R 10 238. DIZETHYL PHTHA	4B.	BENZIDINE	BDL	R 10
68. BENZO (A) PYRENE BDL R 10 7B. 3, 4-BENZOFLUORANTHENE BDL R 10 B8. BENZO (K1) PERYLENE BDL R 25 98. BENZO (K1) FLUORANTHENE BDL R 10 108. BENZO (K1) FLUORANTHENE BDL R 10 108. BIS (2-CHLOROETHYL) ETHER BDL R 10 128. DIS (2-CHLOROETHYL) PHTHALATE BDL R 10 139. DIS (2-CHLOROETHYL) PHTHALATE BDL R 10 148. 4-BROMOPHENYL PHENYL ETHER BDL R 10 158. DUTYL BENZYL PHTHALATE BDL R 10 168. 2-CHLOROPHENYL PHENYL ETHER BDL R 10 168. 2-CHLOROPAPHINALENE BDL R 10 178. 4-CHLOROPAPHINALENE BDL R 10 181. 3-DICHLOROPENZENE BDL R 10 198. DIBENZO (A. H) ANTHRACENE BDL R 10 218. I. 3-DICHLOROBENZENE BDL R 10 218. I. 4-DICHLORDENZENE BDL R 10 <td< td=""><td></td><td></td><td></td><td></td></td<>				
7B. 3.4-BENZOFLUCRANTHENE BOL R 10 8B. BENZO (GHI) PERYLENE BDL R 10 9B. BENZO (GHI) PERYLENE BDL R 10 10B. BTS (2-CHLOROETHOXY) METHANE BDL R 10 11B. DTS (2-CHLOROETHOXY) ETHER BDL R 10 12B. BTS (2-CHLOROETHOXY) METHALATE BDL R 10 13B. DTS (2-CHLOROETHYL) ETHER BDL R 10 14B. 4-BROMOPHENYL PHENYL ETHER BDL R 10 14B. 4-CHLOROPHYL PHENYL ETHER BDL R 10 16B. CHRYSENE BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 18B. 12-DICHLOROBENZENE BDL R 10 19B. DTENZO (A, H) ANTHRACENE BDL R 10 28D. 12-DICHLOROBENZENE BDL R 10 28D. 12-DICHLOROBENZENE BDL R 10 28D. 3'3'-DICHLOROBENZENE				=-
88. SENZO (CHI) PERVLENE BOL R. 25 98. BENZO (K) FLUORANTHENE BOL R. 10 109. BIS (2-CHLOROETHOXY) METHANE BOL R. 10 118. BIS (2-CHLOROETHVL) ETHER BOL R. 10 128. BIS (2-CHLOROETHVL) ETHER BOL R. 10 138. DIS (2-CHLOROETHVL) ETHER BOL R. 10 138. DIS (2-CHLOROETHVL) ETHER BOL R. 10 148. 4-BROMOPHENYL PHENYL ETHER BOL R. 10 158. BUTYL DENYL PHTHALATE BOL R. 10 168. CREVSENE BOL R. 10 178. 4-CHLOROBENZENE BOL R. 10 189. DIBENZO (A. H) ANTHRACENE BOL R. 10 198. DIECHLOROBENZENE BOL R. 10 218. I. 3-DICHLOROBENZENE BOL R. 10 228. J. 4-DICHLOROBENZENE BOL R. 10 238. J. 3'-DICHLOROBENZENE BOL R. 10 238. J. 3'-DICHLOROBENZENE BOL R. 10 249. DIETHYL PHTHALATE BOL R. 10 258. DIMETOTOLURENE BOL R. 10 258. DIMETOTOLURENE BOL R. 10 258. DIME				
7B. BENZO (K) FLUORANTHÈNE BDL R 10 108. BDB (2-CHLOROETHOXY) METHANE BDL R 10 118. BTB (2-CHLOROETHOXY) METHANE BDL R 10 128. BTB (2-CHLOROETHOXY) PETHER BDL R 10 138. BTS (2-CHLOROETHYL) ETHER BDL R 10 138. BTS (2-CHLOROETHYL) PETHER BDL R 10 148. 4-BROMOPHENYL PHENYL ETHER BDL R 10 148. 4-BROMOPHENYL PHENYL ETHER BDL R 10 168. 2-CHLORONAPHTHALENE BDL R 10 178. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 188. CHRYSENE BDL R 10 198. DIBENZO (A. H) ANTHRACENE BDL R 10 218. 1.3 -DICHLOROBENZENE BDL R 10 228. 1.4 -DICHLOROBENZENE BDL R 10 238. 3.3'-DICHLOROBENZENE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 258. JINTROTOLUENE <td></td> <td></td> <td></td> <td></td>				
10B. BIE (2-CHLOROETHOXY) METHANE BDL R 10 11B. BIE (2-CHLOROETHYL) ETHER BDL R 10 12B. BIE (2-CHLOROEOROPYL) ETHER BDL R 10 13B. BIE (2-CHLOROEOROPYL) PHHALATE BDL R 10 14B. 4-6ROMOPHENYL PHENYL ETHER BDL R 10 15B. BUTYL BENZYL PHTHALATE BDL R 10 15B. BUTYL BENZYL PHTHALATE BDL R 10 16B. 2-CHLOROBENZL PHENYL ETHER BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 19B. CHRYSENE BDL R 10 19B. CHRYSENE BDL R 10 19B. CHRYSENE BDL R 10 20B. 1: 2-DICHLOROBENZENE BDL R 10 21B. 3: 3'-DICHLOROBENZENE BDL R 10 22B. 3: 4-DINTROTOLUERE BDL R 10 23B. 3: -DICHLOROBENZENE BDL R 10 24B. DIETHYL PHTHALATE BDL R 10 25B. DIPHETHYL PHTHALATE BDL R 10 26B. DI -N-BUTYL PHTHALATE BDL R 10 27B. 2: 4-DINITROTOLUENE BDL R 10 26B. DI-N-BUTYL PHTHALAT				
11B. BIS (2-CHLOROETHYL): ETHER BDL R. 10 12B. BIS (2-CHLOROISOPROPYL) ETHER BDL R. 10 13B. BIS (2-CHLOROISOPROPYL) ETHER BDL R. 10 14B. 4-BROMOPHENYL PHENYL ETHER BDL R. 10 15B. BUTYL BENZYL PHTHALATE BDL R. 10 16B. 2-CHLOROMAPHTHALATE BDL R. 10 16B. 2-CHLOROMAPHTHALENE BDL R. 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R. 10 17B. 4-CHLOROBENZENE BDL R. 10 19B. DIDENZO (A, H) ANTHRACENE BDL R. 10 20B. 1, 2-DICHLOROBENZENE BDL R. 10 21B. 1, 3-DICHLOROBENZENE BDL R. 10 22B. 1, 4-DICHLOROBENZENE BDL R. 10 23B. 3, 3'-DICHLOROBENZENE BDL R. 10 24B. DIETHYL PHTHALATE BDL R. 10 25B. DIMETHYL PHTHALATE BDL R. 10 25B. DIMETHYL PHTHALATE BDL R. 10 26B. 0I-N-ROTYL PHTHALATE BDL R. 10 26B. 1, 2-DIPHENYLHYDRAZINE BDL R. 10 27B. 2, 4-DINITROTOLUENE BDL R. 10				
12B. BIE (2-CHLOROISOPROPYL) ETHER BDL R 10 13B. BIS (2-ETHYLHEXYL) PHTHALATE BDL R 10 14B. 4-BROMOPHENYL PHENYL ETHER BDL R 10 15B. BUTYL BENZYL PHTHALATE BDL R 10 16B. 2-CHLORONAPHTHALENE BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 19B. CHRYSENE BDL R 10 19B. CHRYSENE BDL R 10 20B. I. 2-DICHLOROBENZENE BDL R 10 21B. 1. 3-DICHLOROBENZENE BDL R 10 22B. 3. 3'-DICHLOROBENZENE BDL R 10 22B. 3. 3'-DICHLOROBENZIENE BDL R 10 23B. 3. 3'-DICHLOROBENZIENE BDL R 10 24B. DIETHYL PHTHALATE BDL R 10 25B. DIPHENYL PHTHALATE BDL R 10 26B. DIP-N-BUTYL PHTHALATE BDL R 10 27B. 2. 4-DINITROTOLUENE BDL R 10 27B. 2. 4-DINITROTOLUENE BDL R 10 308. 1. 2-DIPHENYL HYTHALATE BDL R 10 328. 1. 2-DIPHENYL	-			
13B. BIS (2-ETHYLHEXYL) PHTHALATE BDL R 10 14B. 4-BROMOPHENYL PHENYL ETHER BDL R 10 15B. BUTYL BENYL PHTHALATE BDL R 10 16B. 2-CHLORONAPHTHALENE BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 17B. 01BENZO (A.H) ANTHRACENE BDL R 10 18B. CHRYSENE BDL R 10 18B. 1.2-DICHLOROBENZENE BDL R 10 21B. 1.3-DICHLOROBENZENE BDL R 10 22B. 1.4-DICHLOROBENZENE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 258. J.4-DICHLOROBENZIDINE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 268. DI-N-BUTYL PHTHALATE BDL R 10 276. Z.4-DINITROTOLUENE BDL R 10 288. J.2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORENE BDL R 10 338. HEXACHLOROCYCLOPENTADIENE <td< td=""><td></td><td></td><td></td><td></td></td<>				
148. 4-BROMOPHENYL PHENYL ETHER BDL R 10 15B. BUTYL BENZYL PHTHALATE BDL R 10 16B. 2-CHLORONAPHTHALENE BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 17B. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 18B. CHRYSENE BDL R 10 19B. DIBENZO (A. H) ANTHRACENE BDL R 10 22B. 1. 2-DICHLOROBENZENE BDL R 10 23B. 3. 3'-DICHLOROBENZENE BDL R 10 23B. 3. 3'-DICHLOROBENZENE BDL R 10 24B. DIETHYL PHTHALATE BDL R 10 25B. DIMETHYL PHTHALATE BDL R 10 26B. DI-N-BUTYL PHTHALATE BDL R 10 27B. 2.4-DINITROTOLUENE BDL R 10 30B. J.2-OZPHENYLHYDRAZINE BDL R 10 32B. FLUOR	_			
15B. BUTYL BENZYL PHTHALATE BDL R 10 16B. 2-CHLORONAPHTHALENE BDL R 10 17B. 4-CHLORONAPHTHALENE BDL R 10 17B. 4-CHLORONAPHTHALENE BDL R 10 19B. CHRYSENE BDL R 10 19B. DIBENZO (A. H) ANTHRACENE BDL R 10 20B. 1. 2-DICHLOROBENZENE BDL R 10 21B. 1. 3-DICHLOROBENZENE BDL R 10 22B. 1. 4-DICHLOROBENZENE BDL R 10 23B. 3. 3'-DICHLOROBENZENE BDL R 10 24B. DIETHYL PHTHALATE BDL R 10 25B. DIMETHYL PHTHALATE BDL R 10 26B. J. 4-DINITROTOLUENE BDL R 10 26B. J. 4-DINITROTOLUENE BDL R 10 27B. 2. 4-DINITROTOLUENE BDL R 10 28B. 1. 2-DIPHENYLHYDRAZINE BDL R 10 30B. 1. 2-DIPHENYLHYDRAZINE BDL R 10 30B. 1. 2-DIPHENYLHYDRAZINE BDL R 10 31B. FLUORANTHENE BDL R 10 328. HEXACHLOROBUTADIENE BDL R 10 358. HEXACHLOROBUTADIENE BDL R				•
168. 2-CHLORONAPHTHALENE BDL R 10 178. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 188. CHRYSENE BDL R 10 198. DIBENZO (A, H) ANTHRACENE BDL R 10 208. 1. 2-DICHLOROBENZENE BDL R 10 218. 1. 3-DICHLOROBENZENE BDL R 10 228. 1. 4-DICHLOROBENZENE BDL R 10 239. 3. 3'-DICHLOROBENZENE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 268. 2. 4-DINITROTOLUENE BDL R 10 278. 0I-N-OCTYL PHTHALATE BDL R 10 289. 0I-N-OCTYL PHTHALATE BDL R 10 308. FLUORANTHENE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORAN				
178. 4-CHLOROPHENYL PHENYL ETHER BDL R 10 188. CHRYSENE BDL R 10 198. DIBENZO (A, H) ANTHRACENE BDL R 10 208. 1. 2-DICHLOROBENZENE BDL R 10 218. 1. 3-DICHLOROBENZENE BDL R 10 228. 1. 4-DICHLOROBENZENE BDL R 10 238. 3. 3'-DICHLOROBENZENE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. OIMETHYL PHTHALATE BDL R 10 268. DI-N-BUTYL PHTHALATE BDL R 10 276. Z. 4-DINITROTOLUENE BDL R 10 289. J. 2-DIPHENYLHYDRAZINE BDL R 10 308. J. 2-DIPHENYLHYRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. HEXACHLOROBUTADIENE BDL R 10 339. HEXACHLOROBENZENE BDL R 10				
188. CHRYSENE BDL R 10 198. DIBENZO (A, H) ANTHRACENE BDL R 25 208. 1. 2-DICHLOROBENZENE BDL R 10 218. 1. 3-DICHLOROBENZENE BDL R 10 228. 1. 4-DICHLOROBENZENE BDL R 10 238. 3. 3'-DICHLOROBENZENE BDL R 10 238. 0.3'-DICHLOROBENZIDINE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 268. 2. 4-DINITROTOLUENE BDL R 10 278. 2. 4-DINITROTOLUENE BDL R 10 278. 2. 4-DINITROTOLUENE BDL R 10 278. 2. 4-DINITROTOLUENE BDL R 10 308. 1. 2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORANTHENE BDL R 10 339. MEXACHLOROBUTADIENE BDL R 10 348. MEXACHLOROBUTADIENE BDL R 10 358. HEXACHLOROBUTADIENE BDL R 10 368. IESOPHOROME BDL R 10 </td <td></td> <td></td> <td></td> <td></td>				
198. DIBENZO (A, H) ANTHRACENE BDL R 25 208. I. 2-DICHLOROBENZENE BDL R 10 218. I. 3-DICHLOROBENZENE BDL R 10 228. I. 4-DICHLOROBENZENE BDL R 10 238. 3. 3'-DICHLOROBENZENE BDL R 10 238. 3. 3'-DICHLOROBENZIDINE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 268. DI-N-BUTYL PHTHALATE BDL R 10 278. 2.4-DINITROTOLUENE BDL R 10 288. 2.6-DINITROTOLUENE BDL R 10 278. 2.4-DINITROTOLUENE BDL R 10 308. 1.2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORENE BDL R 10 338. HEXACHLOROBENZENE BDL R 10 349. MEXACHLOROBENZENE BDL R 10 358. HEXACHLOROBENZENE BDL R 10 358. HEXACHLOROBENZENE BDL R 10 368. NEXACHLOROCYCLOPENTADIENE BDL R 10 378. INDENO (1, 2, 3-CD) PYRENE BDL R				
208. 1. 2-DICHLOROBENZENE BDL R 10 218. 1. 3-DICHLOROBENZENE BDL R 10 228. 1. 4-DICHLOROBENZENE BDL R 10 238. 3. 3'-DICHLOROBENZENE BDL R 10 238. 3. 3'-DICHLOROBENZENE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 258. 2. 4-DINITROTOLUENE BDL R 10 276. 2. 4-DINITROTOLUENE BDL R 10 278. 0OCTYL PHTHALATE BDL R 10 278. 0OCTYL PHTHALATE BDL R 10 308. 1. 2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORENE BDL R 10 339. MEXACHLOROBENZENE BDL R 10 358. HEXACHLOROBENZENE BDL R 10 358. HEXACHLOROBENZENE BDL R 10 358. INSOPHORONE BDL R 10 358. INSOPHORONE BDL R 10 378. INDENO (1.2.3-CD) PYRENE BDL R 10 378. INPETHALÈNE BDL R 10				
218. 1. 3-DICHLOROBENZENE BDL R 10 228. 1. 4-DICHLOROBENZENE BDL R 10 238. 3. 3'-DICHLOROBENZIDINE BDL R 10 248. DIETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 258. DIMETHYL PHTHALATE BDL R 10 268. DI-N-BUTYL PHTHALATE BDL R 10 268. DI-N-OCTYL PHTHALATE BDL R 10 278. 2. 4-DINITROTOLUENE BDL R 10 288. 2. 4-DINITROTOLUENE BDL R 10 308. 1. 2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORENE BDL R 10 338. MEXACHLOROBUTADIENE BDL R 10 348. MEXACHLOROBUTADIENE BDL R 10 358. MEXACHLOROBENZENE BDL R 10 364. MEXACHLOROBUTADIENE BDL R 10 378. INDENO (1, 2, 3-CD) PYRENE BDL R 10 378. NAPHTHALÈNE BDL R 10 378. NAPHTHALÈNE BDL R 10			BDL	r 25
22B. 1. 4-DICHLORBENZENE BDL R 10 23B. 3. 3'-DICHLOROBENZIDINE BDL R 10 24B. DIETHYL PHTHALATE BDL R 10 25B. DIMETHYL PHTHALATE BDL R 10 26B. DI-N-BUTYL PHTHALATE BDL R 10 26B. DI-N-BUTYL PHTHALATE BDL R 10 27B. 2. 4-DINITROTOLUENE BDL R 10 28B. 01-N-BUTYL PHTHALATE BDL R 10 29B. 01-N-CTYL PHTHALATE BDL R 10 29B. 1. 2-DIPHENYLHYDRAZINE BDL R 10 30B. FLUORANTHENE BDL R 10 31B. FLUORANTHENE BDL R 10 32B. FLUORENE BDL R 10 33B. HEXACHLOROBENZENE BDL R 10 34B. HEXACHLOROBENZENE BDL R 10 35B. HEXACHLOROBENZENE BDL R 10 36B. NEXACHLOROBENZENE BDL R 10 37B. INDENO (1, 2, 3-CD) PYRENE BDL R 10 37B. NAPHTHALÈNE BDL R 10 37B. NAPHTHALÈNE BDL R 10 37B. NAPHTHALÈNE BDL R 10	208.	1, 2-DICHLOROBENZENE	BDL	R . 10
23B. 3, 3'-DICHLOROBENZIDINE BDL R. 10 249. DIETHYL PHTHALATE BDL R. 10 25B. DIMETHYL PHTHALATE BDL R. 10 25B. DIMETHYL PHTHALATE BDL R. 10 25B. DIMETHYL PHTHALATE BDL R. 10 26B. DI-N-BUTYL PHTHALATE BDL R. 10 27B. 2, 4-DINITROTOLUENE BDL R. 10 27B. 01-N-OCTYL PHTHALATE BDL R. 10 27B. 01-N-OCTYL PHTHALATE BDL R. 10 32B. FLUORANTHENE BDL R. 10 32B. FLUORANTHENE BDL R. 10 32B. FLUORANTHENE BDL R. 10 32B. HEXACHLOROBUTADIENE BDL R. 10 33B. MEXACHLOROBUTADIENE BDL R. 10 34B. HEXACHLOROCYCLOPENTADIENE BDL R. 10 35B. INDENO (1,2,3-CD) PYRENE BDL R. 10 37B. NAPHTHALENE BDL R. 10 448. PHENANTHRENE BDL R. 10	218.	1, 3-DICHLOROBENZENE	BDL	r 10
248. DIETHYL PHTHALATE BDL 2 10 258. DIMETHYL PHTHALATE BDL 2 10 268. DI-N-BUTYL PHTHALATE BDL 2 10 278. 2. 4-DINITROTOLUENE BDL 2 10 268. 2. 4-DINITROTOLUENE BDL 2 10 278. 2. 4-DINITROTOLUENE BDL 2 10 278. 01-N-OCTYL PHTHALATE BDL 2 10 308. 1. 2-DIPHENYLHYDRAZINE BDL 2 10 328. FLUORENE BDL 2 10 328. FLUORENE BDL 2 10 339. HEXACHLOROBENZENE BDL 2 10 348. HEXACHLOROBUTADIENE BDL 2 10 358. HEXACHLOROCYCLOPENTADIENE BDL 2 10 368. IEOPHORONE BDL 2 25 388. IEOPHORONE BDL 2 25 388. IEOPHORONE BDL 2 10 378. NAPHTHALÈNE BDL 2 10 378. NAPHTHALÈNE BDL 2 10 378. N-NITROSODIMETHYLAMINE BDL 2 10 448. PHENANTHRENE BDL 2 10 458. N-NITROSODIPHENYLAMINE BDL 2 10 448. PHENANTHRENE	228.	1. 4-DICHLORBENZENE	' BDL	R 10
248. DIETHYL PHTHALATE BDL 2 10 258. DIMETHYL PHTHALATE BDL 2 10 268. DI-N-BUTYL PHTHALATE BDL 2 10 278. 2. 4-DINITROTOLUENE BDL 2 10 268. 2. 4-DINITROTOLUENE BDL 2 10 278. 2. 4-DINITROTOLUENE BDL 2 10 278. 01-N-OCTYL PHTHALATE BDL 2 10 308. 1. 2-DIPHENYLHYDRAZINE BDL 2 10 328. FLUORENE BDL 2 10 328. FLUORENE BDL 2 10 339. HEXACHLOROBENZENE BDL 2 10 348. HEXACHLOROBUTADIENE BDL 2 10 358. HEXACHLOROCYCLOPENTADIENE BDL 2 10 368. IEOPHORONE BDL 2 25 388. IEOPHORONE BDL 2 25 388. IEOPHORONE BDL 2 10 378. NAPHTHALÈNE BDL 2 10 378. NAPHTHALÈNE BDL 2 10 378. N-NITROSODIMETHYLAMINE BDL 2 10 448. PHENANTHRENE BDL 2 10 458. N-NITROSODIPHENYLAMINE BDL 2 10 448. PHENANTHRENE	23B.	3, 3'-DICHLOROBENZIDINE	BDL	R 10
25B. DIMETHYL PHTHALATE BDL R 10 26B. DI-N-BUTYL PHTHALATE BDL R 10 27B. 2. 4-DINITROTOLUENE BDL R 10 28B. 2. 4-DINITROTOLUENE BDL R 10 28B. 2. 4-DINITROTOLUENE BDL R 10 27B. 2. 4-DINITROTOLUENE BDL R 10 28B. 2. 4-DINITROTOLUENE BDL R 10 30B. 1. 2-DIPHENYL PHTHALATE BDL R 10 30B. 1. 2-DIPHENYL HYDRAZINE BDL R 10 32B. FLUORANTHENE BDL R 10 32B. FLUORANTHENE BDL R 10 32B. HEXACHLOROBENZENE BDL R 10 34B. MEXACHLOROCYCLOPENTADIENE BDL R 10 35B. HEXACHLOROCYCLOPENTADIENE BDL R 10 36B. IEOPHORONE BDL R 10 37B. INDENO (1, 2, 3-CD) PYRENE BDL R 10 37B. INDENO (1, 2, 3-CD) PYRENE BDL R 10 37B. NAPHTHALENE BDL R 10 408. NTROSODIMETHYLAMINE BDL R 10 418. N-NITROSODIMETHYLAMINE BDL R 10 428. N-NITROSODIPHENYLAMINE BDL R <td></td> <td></td> <td></td> <td></td>				
268. DI-N-BUTYL PHTHALATE BDL R 10 278. 2. 4-DINITROTOLUENE BDL R 10 288. 2. 6-DINITROTOLUENE BDL R 10 288. 2. 6-DINITROTOLUENE BDL R 10 298. 0I-N-OCTYL PHTHALATE BDL R 10 308. 1. 2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORENE BDL R 10 339. HEXACHLOROBENZENE BDL R 10 348. HEXACHLOROBUTADIENE BDL R 10 358. HEXACHLOROCYCLOPENTADIENE BDL R 10 368. IEOPHORONE BDL R 10 378. INDENO (1.2.3-CD) PYRENE BDL R 10 378. NAPHTHALENE BDL R 10 378. NAPHTHALENE BDL R 10 408. NITROBENZENE BDL R 10 418. N-NITROSODIMETHYLAMINE BDL R 10 428. N-NITROSODIMETHYLAMINE BDL R 10 428. N-NITROSODIMETHYLAMINE BDL R 10 438. N-NITROSODIPHENYLAMINE BDL R 10 448. PHENANTHRENE BDL R 10 <tr< td=""><td>258.</td><td>DIMETHYL PHTHALATE</td><td></td><td>R 10</td></tr<>	258.	DIMETHYL PHTHALATE		R 10
27B. 2. 4-DINITROTOLUENE BDL 4 10 28B. 2. 6-DINITROTOLUENE BDL R 10 27B. DI-N-OCTYL PHTHALATE BDL R 10 30B. 1. 2-DIPHENYLHYDRAZINE BDL R 10 31B. FLUORANTHENE BDL R 10 32B. FLUORANTHENE BDL R 10 32B. HEXACHLOROBENZENE BDL R 10 33B. HEXACHLOROBUTADIENE BDL R 10 34B. HEXACHLOROCYCLOPENTADIENE BDL R 10 35B. HEXACHLOROETHANE BDL R 10 36B. INDENO (1, 2, 3-CD) PYRENE BDL R 10 37B. INDENO (1, 2, 3-CD) PYRENE BDL R 10 37B. INDENO (1, 2, 3-CD) PYRENE BDL R 10 37B. INDENO (1, 2, 3-CD) PYRENE BDL R 10 37B. INDENONE BDL R 10 37B. NAPHTHALENE BDL R 10 </td <td></td> <td></td> <td></td> <td></td>				
28B. 2, 6-DINITROTOLUENE BDL R 10 27B. DI-N-OCTYL PHTHALATE BDL R 10 30B. 1, 2-DIPHENYLHYDRAZINE BDL R 10 31B. FLUORANTHENE BDL R 10 32B. FLUORANTHENE BDL R 10 32B. FLUORANTHENE BDL R 10 32B. FLUORENE BDL R 10 33B. HEXACHLOROBENZENE BDL R 10 34B. HEXACHLOROBUTADIENE BDL R 10 35B. HEXACHLOROCYCLOPENTADIENE BDL R 10 36B. HEXACHLOROCTCLOPENTADIENE BDL R 10 36B. HEXACHLOROCTRANE BDL R 10 37B. INDENO (1.2, 3-CD) PYRENE BDL R 10 37B. INDENO (1.2, 3-CD) PYRENE BDL R 10 37B. INDENO (1.2, 3-CD) PYRENE BDL R 10 37B. NAPHTHALÈNE BDL R 10 408. NITROSODIMETHYLAMINE BDL R 10 408. N-NITROSODIMETHYLAMINE BDL R 10 428. N-NITROSODIPHENYLAMINE BDL R 10 438. N-NITROSODIPHENYLAMINE BDL R 10 448. PHENANTHRENE BDL R 10	_			
278. DI-N-OCTYL PHTHALATE BDL R 10 308. 1.2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORENE BDL R 10 329. HEXACHLOROBENZENE BDL R 10 339. HEXACHLOROBUTADIENE BDL R 10 348. HEXACHLOROCYCLOPENTADIENE BDL R 10 358. HEXACHLOROCYCLOPENTADIENE BDL R 10 368. HEXACHLOROCTHANE BDL R 10 368. HEXACHLOROETHANE BDL R 10 378. INDENO (1, 2, 3-CD) PYRENE BDL R 10 378. ISOPHORONE BDL R 10 378. ISOPHORONE BDL R 10 378. NAPHTHALENE BDL R 10 408. NITROBODIMETHYLAMINE BDL R 10 408. N-NITROSODIPHENYLAMINE BDL R 10 438. N-NITROSODIPHENYLAMINE BDL R 10 438. PHENANTHRENE BDL R 10 458.				
308. 1.2-DIPHENYLHYDRAZINE BDL R 10 318. FLUORANTHENE BDL R 10 328. FLUORENE BDL R 10 338. HEXACHLOROBENZENE BDL R 10 348. HEXACHLOROBUTADIENE BDL R 10 348. HEXACHLOROCYCLOPENTADIENE BDL R 10 358. HEXACHLOROCYCLOPENTADIENE BDL R 10 368. HEXACHLOROCYCLOPENTADIENE BDL R 10 368. HEXACHLOROCYCLOPENTADIENE BDL R 10 378. INDENO (1, 2, 3-CD) PYRENE BDL R 10 378. ISOPHORONE BDL R 10 378. ISOPHORONE BDL R 10 378. NAPHTHALENE BDL R 10 408. NITROBENZENE BDL R 10 418. N-NITROSODIMETHYLAMINE BDL R 10 428. N-NITROSODIPHENYLAMINE BDL R 10 438. N-NITROSODIPHENYLAMINE BDL R 10 448. PHENANTHRENE BDL R 10 <td< td=""><td></td><td>•</td><td></td><td></td></td<>		•		
31B.FLUORANTHENEBDLR1032B.FLUORENEBDLR1033B.HEXACHLOROBENZENEBDLR1034B.HEXACHLOROBUTADIENEBDLR1035B.HEXACHLOROCYCLOPENTADIENEBDLR1036B.HEXACHLOROETHANEBDLR1037B.INDENO (1, 2, 3-CD) PYRENEBDLR1037B.INDENO (1, 2, 3-CD) PYRENEBDLR1037B.NAPHTHALÉNEBDLR1039B.NAPHTHALÉNEBDLR1040B.NITROBENZENEBDLR1040B.NITROSODIMETHYLAMINEBDLR1042B.N-NITROSODIMETHYLAMINEBDLR1043B.N-NITROSODIPHENYLAMINEBDLR10448.PHENANTHRENEBDLR10458.PYRENEBDLR10468.1. 2, 4-TRICHLOROBENZENEBDLR101P.ALDRINBDLR102P.ALPHA-EHCBDLR10				
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AR100321	2P.	ALPHA-BHC	BDL	R 10
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BDL= BELOW DETECTION LIMIT

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SAMPLE IDENTIFIER: 8086 COMPU/CHEM SAMPLE NUMBER: 3826

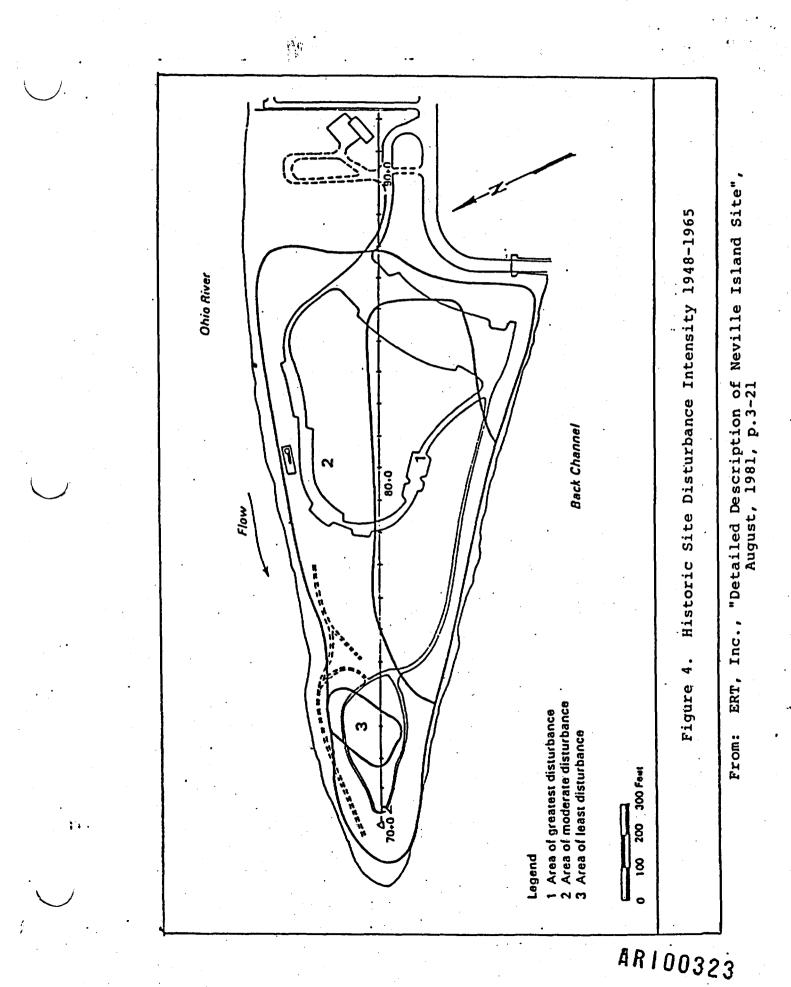
	COMPOUNDS	CONCENTRATION (UG/L)	DETECTION LIMIT (UG/L)	
3 P .	BETA-BHC	BOL	R	10
4P.		BDL	R	10
SP.	DELTA-BHC	BOL	R	10
6P.		BOL	R	10
7P.		BOL	R	10
8P.		BOL	R	10
9P.		BOL	8	10
10P.	DESLORIN	BOL	R	10
11P.		BDL	R	10
12P.	-	BOL	R	10
13P.		801	R	10
14P.	ENDRIN	BOL	R	10
15P.		BOL	R	10
16P.		BDL	ñ	10
17P.		BDL	R	10
18P.		BOL	R	10
19P.		BDL	રે	10
20P.	PC8-1221	BDL	રે	10
21P.	PC8-1232	BOL		10
		•	R	
22P.	PC8-1248	BDL	2	10
23P.	PCB-1260	BOL	R	10
24P.	PCB-1016	BOL	R	10
25P.	TOXAPHENE	BOL	R	10

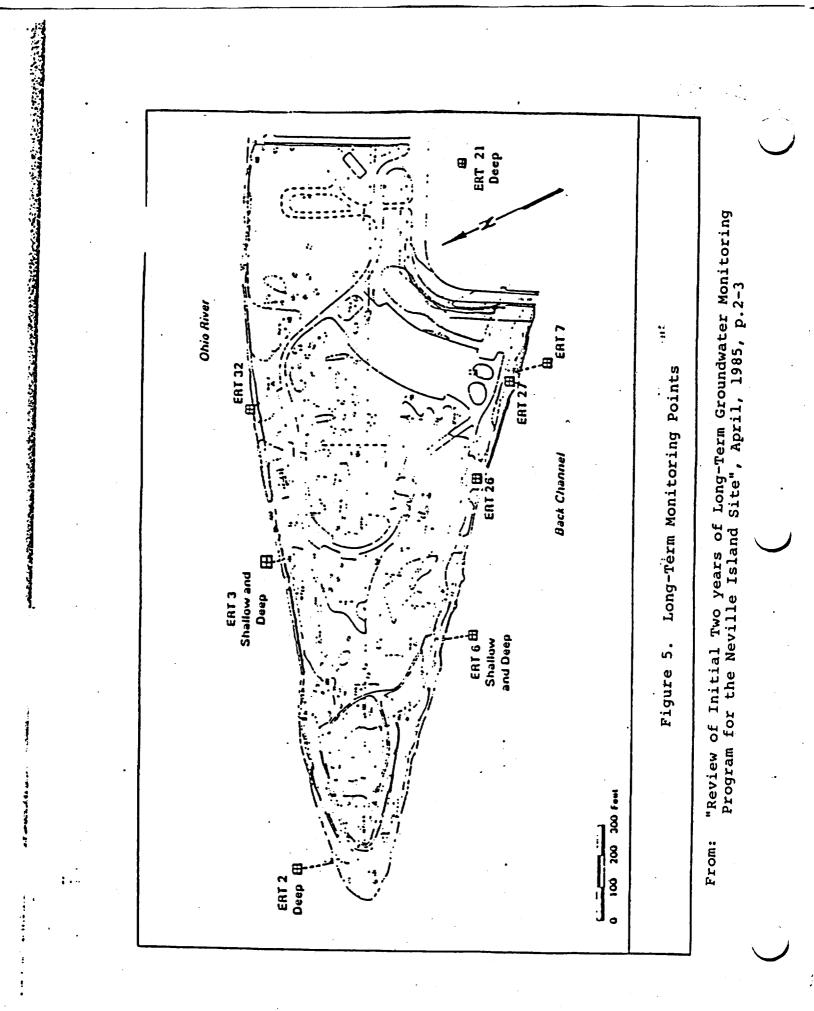
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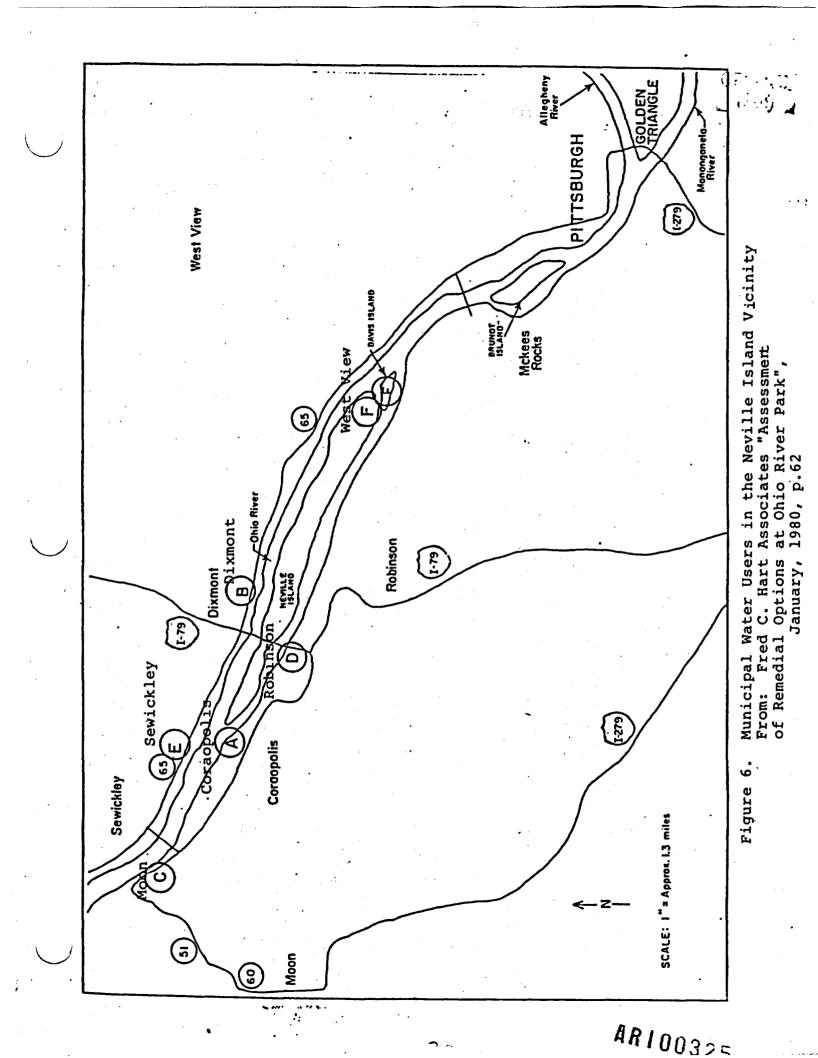
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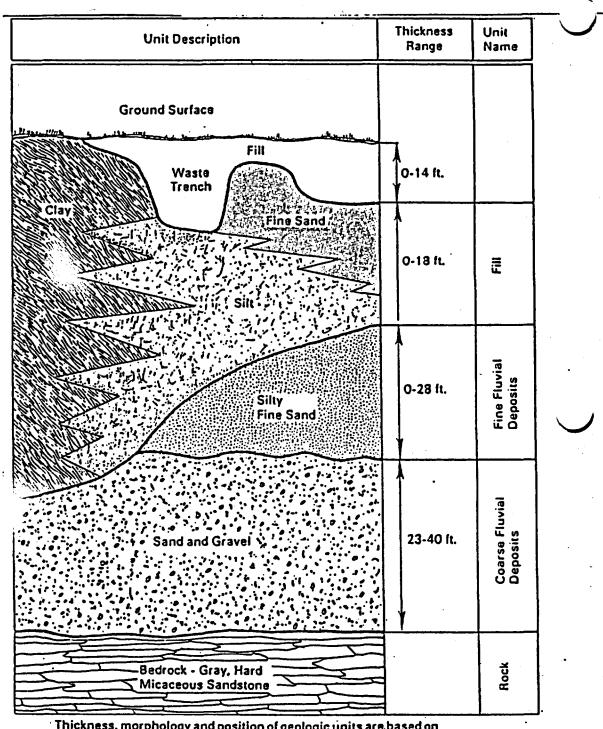
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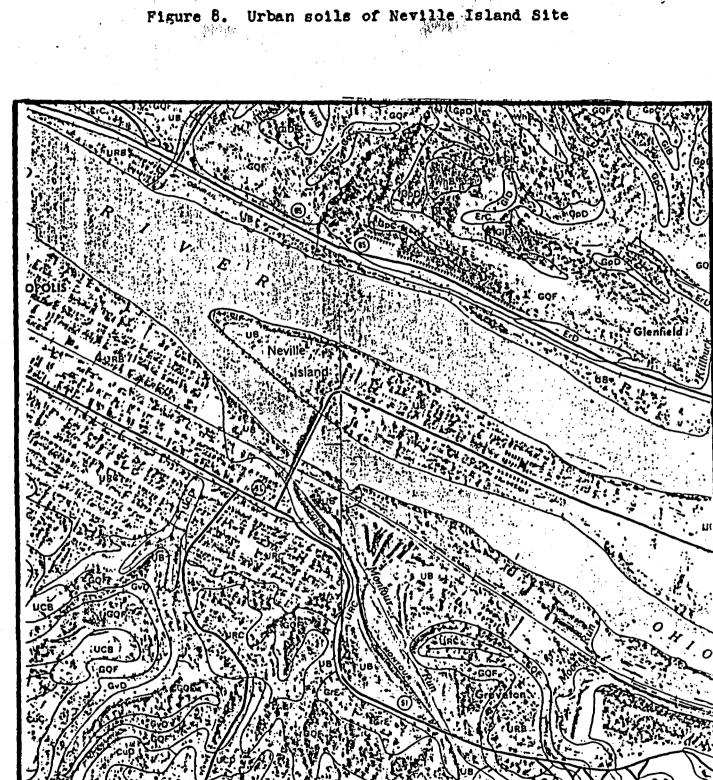
Figure 7. Schematic Stratigraphic Column



Thickness, morphology and position of geologic units are based on boring information. This column represents a variety of conditions that may be encountered beneath the site; it may not depict all possible conditions.

From: ERT, Inc. "Preliminary Risk Assessment of Neville Island Site", April, 1981, p.4-7

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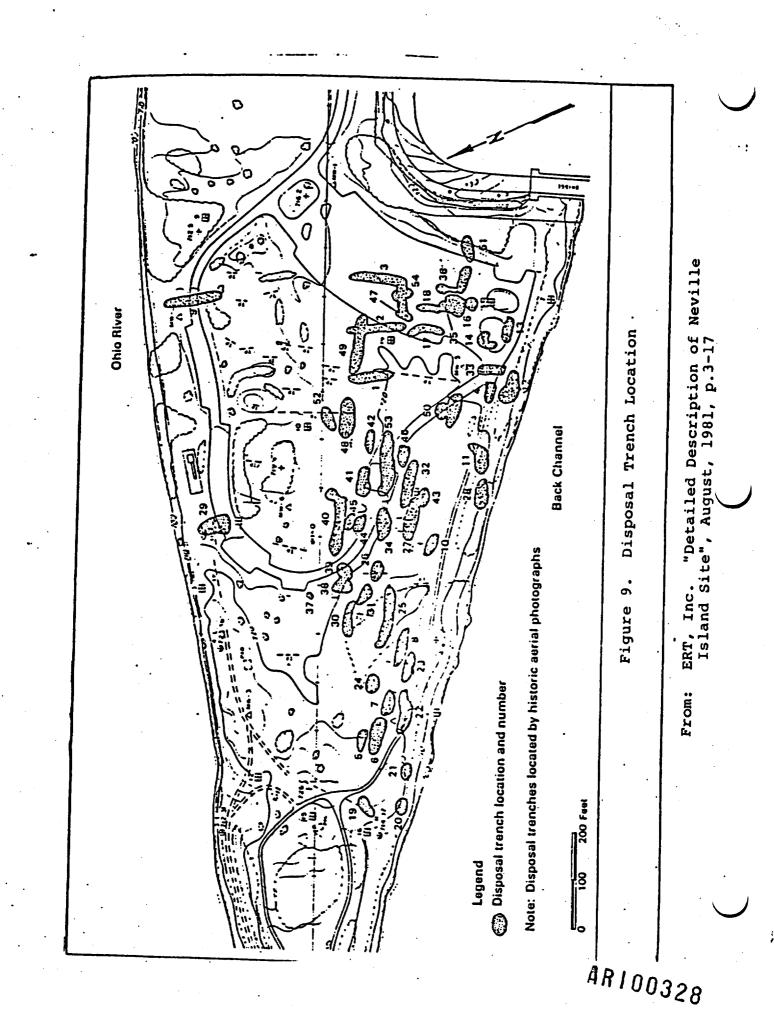


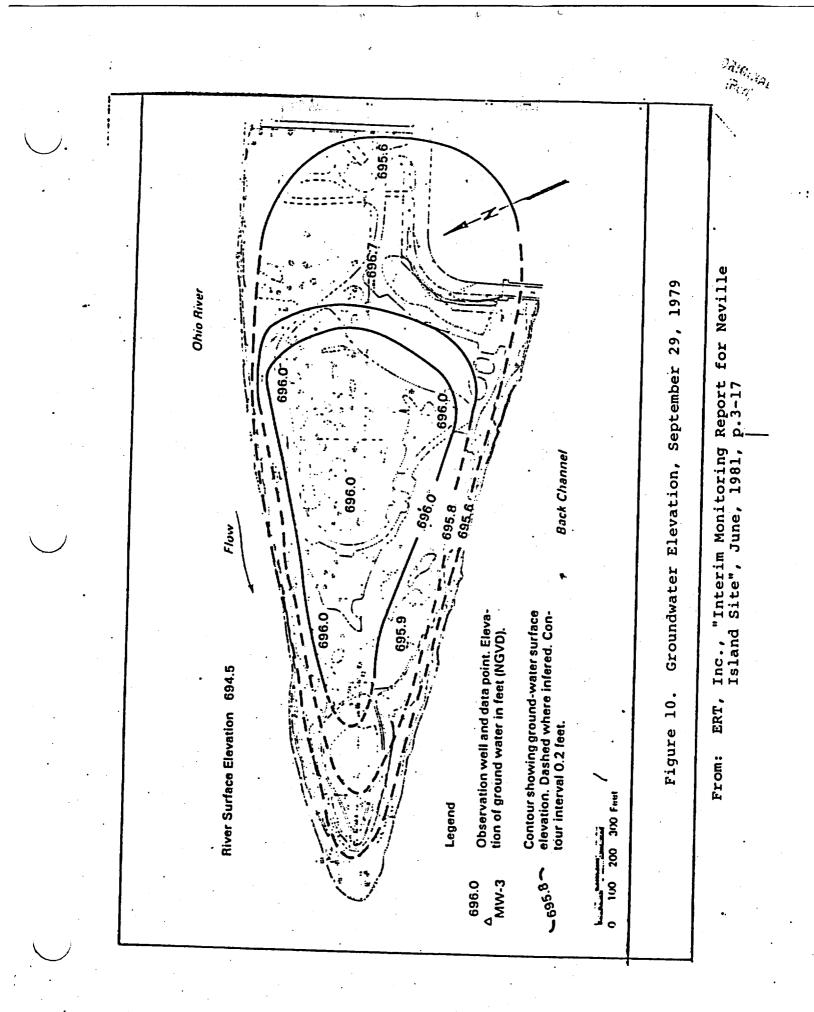
From: U.S.D.A. Soil Conservation Service Soil Survey of Allegheny County, Pa. PADER State Conservation Commission, March, 1988. RICINAL IP.ca;

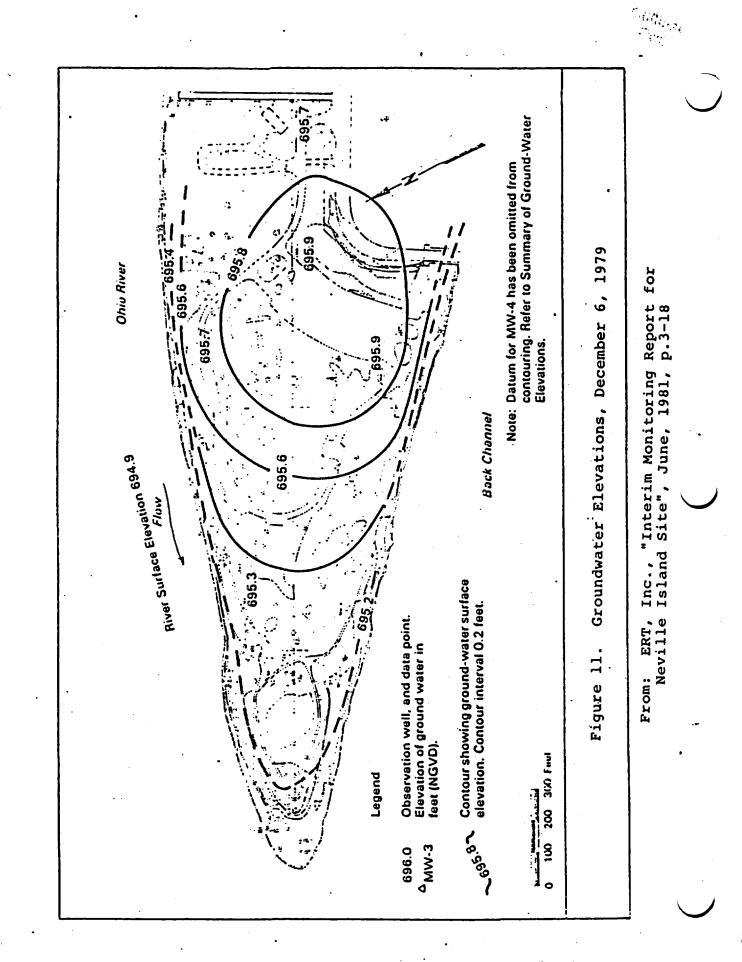
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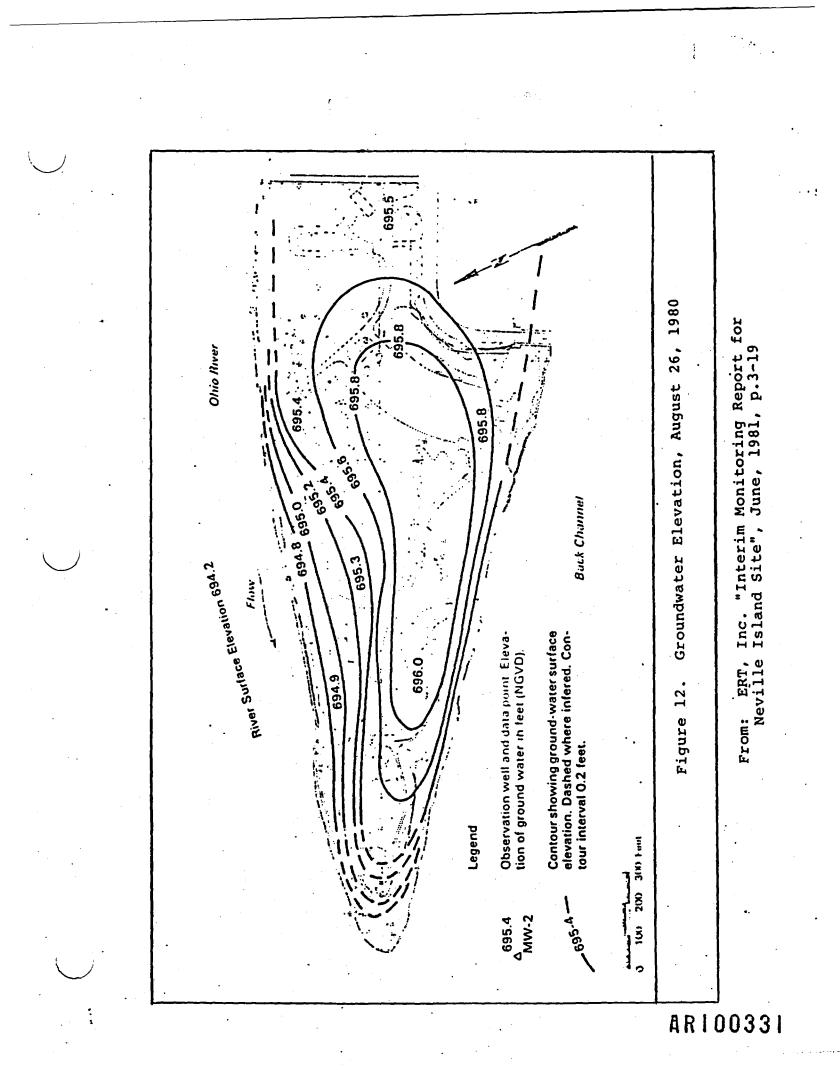


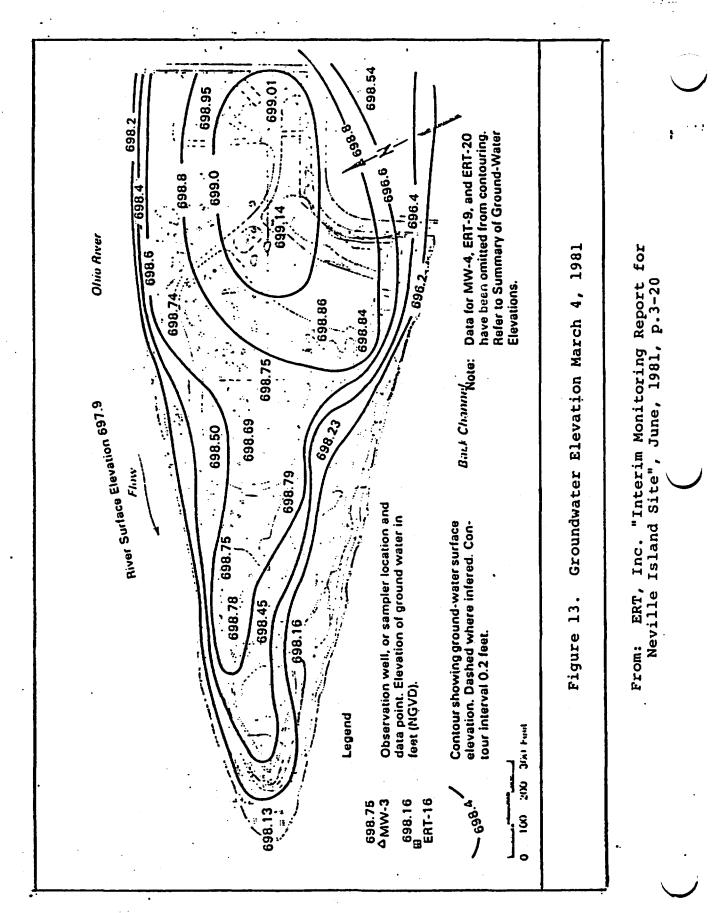




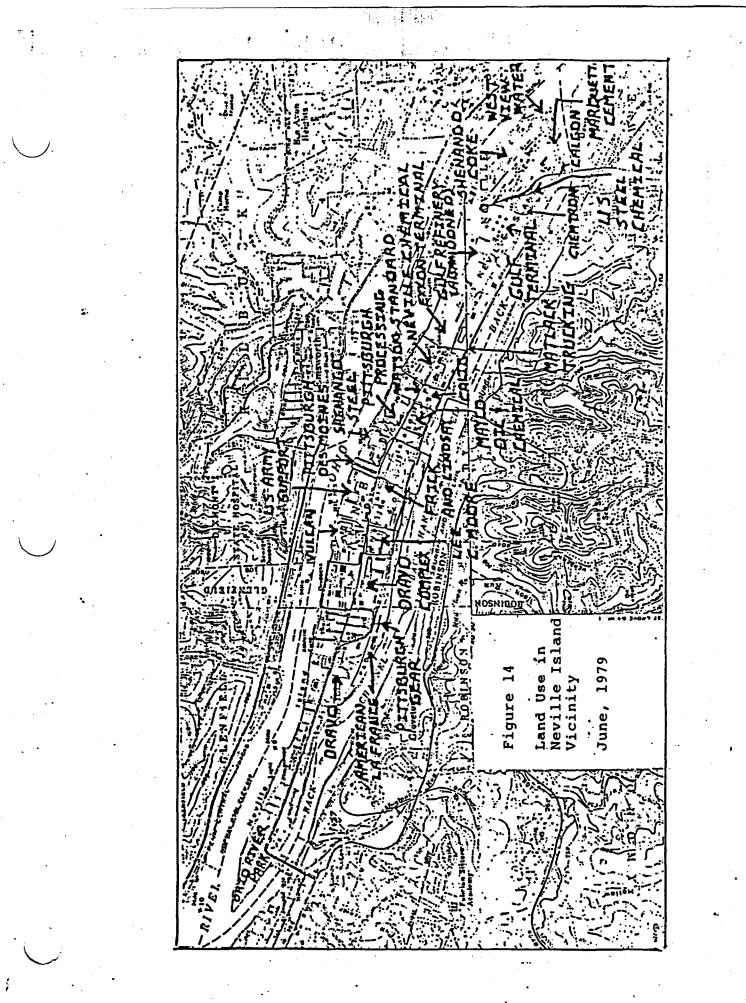
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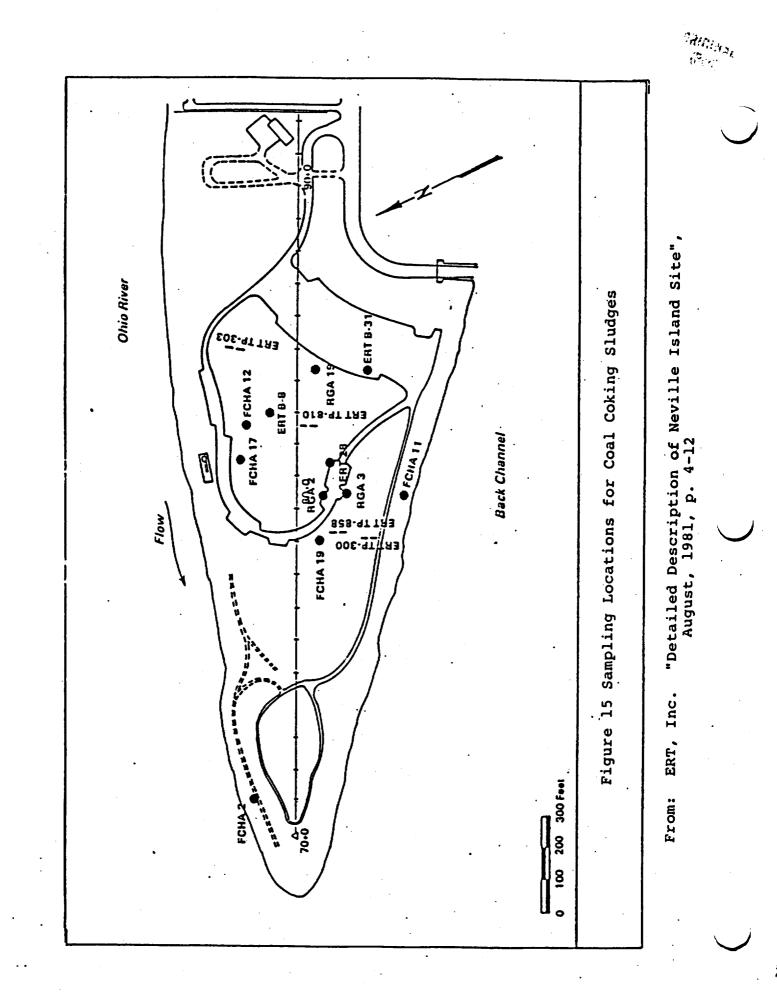
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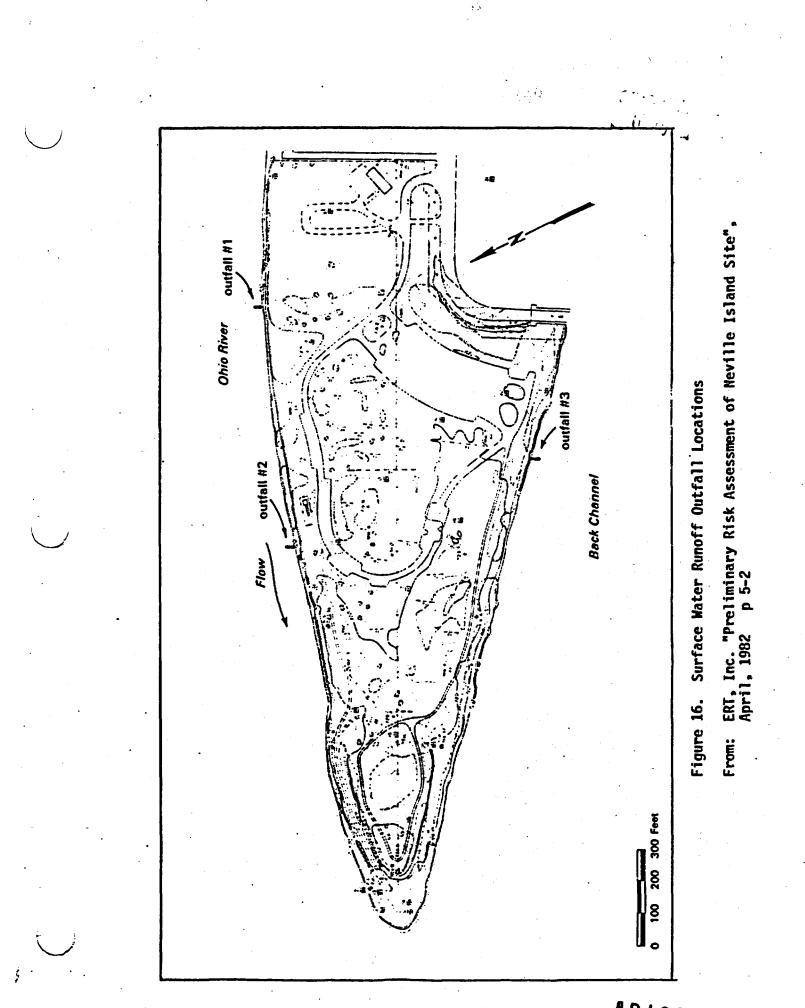




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- 8. Ibid., p. 3 27.
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- 11. Sittig, M., Handbook of Toxic and Hazardous Chemicals and Carcinogens, second edition, Noyes Publications, Park Ridge, N.J., 1985, p. 704.
- 12. Environmental Research & Technology, Inc., Interim Monitoring Report for Neville Island Site, June 1981, p. 3 - 30.
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- 15. Sittig, N., Handbook of Toxic and Hazardous Chemicals and Carcinogens, second edition, Noyes Publications, Park Ridge, N.J., 1985, p. 237.

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- 16. Ibid., page 327.
- 17. Environmental Research & Technology, Inc., Interim Monitoring Report for Neville Island Site, June 1981, page 3-29.
- 18. U. S. Environmental Protection Agency, 2,4-Dimethylphenol: Ambient Water Quality Criteria, Washington, D.C., 1980.
- 19. Environmental Research & Technology, Inc., Interim Monitoring Report for Neville Island Site, June 1981, page 3-31.
- 20. U. S. Environmental Protection Agency, Naphthalene: Health Advisory, Office of Drinking Water, Wawshington, D.C., 1988.
- 21. Environmental Research & Technology, Inc., Interim Monitoring Report for Neville Island Site, June 1981, page 3-30.
- 22. U. S. Environmental Protection Agency, Phthalate Esters: Ambient Water Quality Criteria, Washington, D.C., 1980.
- 23. Sittig, M., Handbook of Toxic and Hazardous Chemicals and Carcinogens, second edition, Noyes Publications, Park Ridge, N.J., 1985, page 288.
- 24. Environmental Research & Technology, Inc., Interim Monitoring Report for Neville Island Site, June 1981, page 3-33.
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- 27. Federal Register. 1985. National Primary Drinking Water Regulations; Synthetic Organic Chemicals, Inorganic Chemicals, and Microorganisms; Final Rule and Proposed Rule. Vol. 50, No. 219. Wednesday, November 13, 1985.
- 28. (a) U. S. Environmental Protection Agency, Quality Criteria for Water, Office of Water and Hazardous Materials, Washington, D.C., July 1976.
 (b) U. S. Environmental Protection Agency, Ambient Water Quality Criteria, Washington, D.C., 1980.
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- 31. (a) International Agency for Research on Cancer, IARC Monographs on the Carcinogenic Risks of Chemicals to Humans, Lyon, France <u>1</u>, 17 (1972). (b) 1bid., <u>23</u>, 325 (1980).
- 32. U. S. Environmental Protection Agency, Cyanides: Ambient Water Quality Criteria, Washington, D.C., 1980.
- 33. Environmental Research & Technology, Inc., Review of Initial Two Years of Long-Term Groundwater Monitoring Program for the Neville Island Site, April 1985.
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Appendix A

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An ENSR Company 601 GRANT STREET, PORTER BUILDING, 10th FLOOR, PITTSBURGH, PA 15219, (412) 261-2910



environmental and engineering excellence

May 16, 1988

Mr. James R. Shack Project Officer Division of Remedial Response Pennsylvania DER Bureau of Waste Management Highland Building 121 South Highland Avenue Pittsburgh, PA 15206-3988

SUBJECT: NEVILLE ISLAND SITE

Dear Jim:

The enclosed report entitled "Security, Monitoring, Inspection and Maintenance Programs for Neville Island Site" was inadvertently left out of the information sent to you earlier today. My apologies for any inconvenience that this oversight may have caused.

Sincerely yours, itmo

Robert W. Rittheyer, P.E. Manager Pittsburgh Operations

RWR/kml

Enclosure

cc: M.J. Laskow M.A. Ferlin



An ENSR Company 601 GRANT STREET, PORTER BUILDING, 10th FLOOR, PITTSBURGH, PA 15219, (412) 261-2910

environmental and engineering excellence

May 16, 1988 ERT Project No.: 4920-001-400

Mr. James R. Shack Project Officer Division of Remedial Response Pennsylvania DER Bureau of Waste Management Highland Building 121 South Highland Avenue Pittsburgh, PA 15206-3988

SUBJECT: NEVILLE ISLAND SITE

Dear Jim:

This letter and its attachments are in response to your 15 March and 29 February 1988 letters requesting additional information regarding the Neville Land Company (NLC) site on Neville Island. The items listed in your 15 March 1988 letter are addressed separately below. A copy of the report entitled "Security, Monitoring, Inspection, and Maintenance Programs for Neville Island Site," as requested in your 29 February letter, is enclosed.

<u>Ownership History</u>

The site was farmland until it was acquired by Pittsburgh Coke & Iron Company in the 1920's. On October 19, 1944, Pittsburgh Coke & Iron Company was renamed to Pittsburgh Coke & Chemical Company.

Title to the site was probably held at one time or another by subsidiaries, Green <u>Bag</u> Cement Company and later Neville Island Land Company, which merged into Pittsburgh Coke & Chemical Company on December 10, 1964.

Pittsburgh Coke & Chemical Company continued to own the property until August 14, 1970, when it conveyed the property to a wholly-owned subsidiary, Neville Land Company. At about that time Pittsburgh Coke & Chemical Company, a majority-owned subsidiary of The Hillman Company, became wholly-owned. Pittsburgh Coke & Chemical Company is no longer in existence as a result of a merger. (See Company Relationships, below).

ALASKA · CALIFORNIA · COLORADO · ILLINOIS · MASSACHUSETTS · MINNESOTA · NEW JERSEY · PENNSYLVANIA · AP-10034

CRICINAL .

Mr. James R. Shack May 16, 1988 Page Two

The site was donated to Allegheny County on March 4, 1977. In June, 1980 Allegheny County reconveyed the site to Neville Land Company.

Owner contacts are as follows:

o Neville Land Company 1900 Grant Building Pittsburgh, Pennsylvania 15219 Contact: Mark J. Laskow (412) 281-2620

> Parent Corporation: Wilmington Securities, Inc. 1006 Wilmington Trust Center Wilmington, Delaware 19801 Contact: Mark J. Laskow (412) 281-2620

Allegheny County
 Department of Planning and Development
 429 Forbes Avenue
 Pittsburgh, Pennsylvania 15219
 Contact: Director of Department of Planning and
 Development (412) 355-5960

<u>Generators/Users</u>

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Pittsburgh Coke & Iron Company This corporation merged into Pittsburgh Coke & Chemical Company which went through a series of mergers and as a result merged into Wilmington Securities, Inc. whose principle office is located at:

1006 Wilmington Trust Center Rodney Square North Wilmington, Delaware 19801 Contact: Mark J. Laskow (412) 281-2620

- o Neville Township 3rd Street and Grand Avenue Pittsburgh, Pennsylvania 15225 Contact: Supervisor (412) 264-8235
- Pittsburgh Coke & Chemical Company
 This corporation went through a series of mergers and as a result merged into Wilmington Securities, Inc.
 whose principle office if located at:

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Mr. James R. Shack May 16, 1988 Page Three

> 1006 Wilmington Trust Center Rodney Square North Wilmington, Delaware 19801 Contact: Mark J. Laskow (412) 281-2620 Parent Corporation: The Hillman Company

No formal records exist regarding the types and quantities of wastes disposed at the site. The report entitled "Detailed Description of Neville Island Site" (ERT Document No. P-A616-435, August 1981) presents the best available information regarding this issue. Section 5.1 of this report summarizes the types and estimated quantities of wastes presently found at the site.

Transporters

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Philips Contracting Company 88 Beaver Grade Road Pittsburgh, Pennsylvania 15108 Contact: Jake Philips (412) 923-1717

Philips Contracting Company hauled only cinders, foundry sand, brick, etc. from 1936 - 1945. He also hauled and buried domestic garbage for Neville Township.

Operators

The site was operated by Pittsburgh Coke & Iron Company from the 1920's to 1944. Pittsburgh Coke & Iron Company merged into Pittsburgh Coke & Chemical Company who continued to operate the site until the 1950's. During the years that material was disposed of at the site, Pittsburgh Coke & Chemical Company was a publicly owned corporation whose largest stockholder was a corporate predecessor of The Hillman Company. This corporation's successor is now a wholly-owned subsidiary of The Hillman Company, Wilmington Securities, Inc. Its office is located at:

1006 Wilmington Trust Center Rodney Square North Wilmington, Delaware 19801 Contact: Mark J. Laskow (412) 281-2620 Parent Corporation: The Hillman Company Mr. James R. Shack May 16, 1988 Page Four

NLC. PC&C and Hillman Company Relationship

The Hillman Company owns all of the capital stock of Wilmington Securities, Inc. which is the successor by merger of Pittsburgh Coke & Chemical Company. Neville Land Company is a wholly-owned subsidiary of Wilmington Securities, Inc.

MW-2A Status

Monitoring well MW-2A was installed on property east of the site by Fred C. Hart during their investigations conducted for Allegheny County. This well was found to be damaged beyond repair when ERT conducted its first field sampling effort in August 1980. It was apparently run over by a grass cutter.

Site Permits - Water Supply Effects

No permits have been applied for or received for the site. A Notification of Hazardous Waste Activity was filed in June 1981 to secure an EPA Identification Number for use in manifesting off-site shipments of materials generated during TP-275 excavation (PAT 44-001-3688). NLC subsequently "denotified" in October 1981.

No evidence has been found suggesting that the site has had any effect on any private or public water supply. As discussed in the report entitled "Preliminary Risk Assessment of Neville Island Site" (ERT Document No. P-4616-721, April 1981), contaminated ground water is the most likely pathway for engendering adverse health effects off site. Hydrogeologic evidence, as presented in the report entitled "Interim Monitoring Report for Neville Island Site" (ERT Document No. P-4616-336), suggests that contaminated ground water beneath the site flows radially from the center of the site and enters the Ohio River. Once in the river, volatilization, degradation, and dilution reduce the concentrations of potentially harmful materials to levels below those at which adverse health effects have been demonstrated.

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Mr. James R. Shack May 16, 1988 Page Five

Records Retention

No operational records have been retained. Financial and corporate documents have been retained at the following locations:

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o Generator:

Pittsburgh Coke & Chemical Company This company has merged out of existence

- o Owner/Operator: Neville Land Company 1900 Grant Building Pittsburgh, Pennsylvania 15219
- o Generator: Neville Township 3rd Street and Grand Avenue Pittsburgh, Pennsylvania 15225

o Transporter:

Philips Contracting Company 88 Beaver Grade Road Pittsburgh, Pennsylvania 15108

Analytical OA/OC Information

Information is enclosed regarding the following samples:

- Perimeter ground water monitoring wells, April and October, 1987 (Attachments 1 and 2);
- o Selected ground water monitoring wells, April 1981
 (Attachment 3);
- o Selected ground water monitoring wells and Outfall 1, January 1980 (Attachment 4);
- Outfall 2, upstream and downstream samples, August 1980 (Attachment 5); and
- Selected soil and waste samples, November 1980 (Attachment 6).

Mr. James R. Shack May 16, 1988 Page Six

The content of each attachment is presented below:

Attachment 1: Perimeter Monitoring Wells, October 1987

- o ERT Report submitted to Program Manager
- Data report for volatiles and herbicides from ENSECO (subcontract laboratory)
- o ENSECO full data packages, including chromatograms, mass spectra, quality control data, instrument tune and calibration
- Metals data, including laboratory notebook pages, computer printouts, calibration data, quality control data

ERT laboratory and custody records .

Attachment 2: Perimeter Monitoring Wells, May 1987

- o ERT Report as submitted to Program Manager
- Herbicide report and raw data from ENSECO (subcontract laboratory) for GC analyses, chromatograms, computer printouts, QC data
- Herbicide report and raw data for GC/MS/SIM analyses from ENSECO, chromatograms, QC data
- o Metals analyses raw data, computer printouts, QC data
- Volatile organics analyses, chromatograms, QC data
- o ERT laboratory and custody records

Attachment 3: Selected Monitoring Wells, April 1981

o GC/MS analyses for volatiles, acids, base/neutrals, pesticides, and PCB. Full data packages, including data report sheets, chromatograms, computer printouts, mass spectra of all detected priority pollutants, instrument tune and calibration data are provided for a

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Mr. James R. Shack May 16, 1988 Page Seven

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total of 9 samples. These include two field blanks and a submitted duplicate of MW-4. Data reports for two duplicate analyses of ERT-17S include only the final report sheets; COMFU/CHEM (subcontract laboratory) could not provide the full package within the time frame of our request.

Copies of Master Logbook pages for samples submitted to the ERT Laboratory

- o Summary data sheets for metals, herbicides, water quality parameters
- o Raw data and laboratory notebook pages for water quality measurements
- Metals data report with QC data from ERCO (subcontract laboratory)
- o Herbicide data sheets

o Herbicide chromatograms

Attachment 4: Selected Monitoring Wells and Outfall 1, January 1981

> GC/MS analyses for volatiles, acids, base/neutrals, pesticides, and PCB. Full copies of data reports with chromatograms, mass spectra of all identified priority pollutants, computer printouts, instrument tune, and calibration data are provided for a total of 6 samples. COMPU/CHEM (subcontract laboratory) could not provide the full data package for 17S within the time frame of our request. The data report included for this sample does contain copies of the reconstructed total ion chromatograms and mass spectra of identified priority pollutants. Sample ERT-21S was not submitted for analysis.

> Copies of the ERT Master Log for samples submitted to the ERT Laboratory

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o Data sheets for metals, TOC, cyanide analyses

o Data sheets for herbicide analyses, chromatograms for herbicides

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Attachment 5: Outfall 2, Upstream and Downstream Samples, August 1980

 GC/MS analyses for volatile organics, acids, base/neutrals, pesticides, PCB, acrolein, and acrylonitrile. Copies of data report sheets only are included. COMPU/CHEM (subcontract laboratory) could not provide backup data within the required time frame of our request. Analyses were performed following the protocol of EPA Methods 624 and 625. Quality assurance requirements were met or exceeded according to COMPU/CHEM documentation.

o TCDD analyses. Analyses were performed by Monsanto. The laboratory listed on the report no longer provides this service, and its equipment and records were relocated several years ago. Attempts to track through Monsanto were unsuccessful. The report, as submitted to ERT, contains quality control results and method followed.

- o ERT Master Log pages for samples submitted to ERT Laboratory
- o Results for water quality analyses performed at ERT

Attachment 6: Selected Soil and Waste Samples, November 1980

- o ERT Log Book for samples submitted to the ERT Laboratory
- TCDD analytical results and report. Analyses were performed at Battelle. The report includes a full discussion of the method, quality control measures, and results obtained for the samples.
- o Herbicide and volatiles summary data sheets

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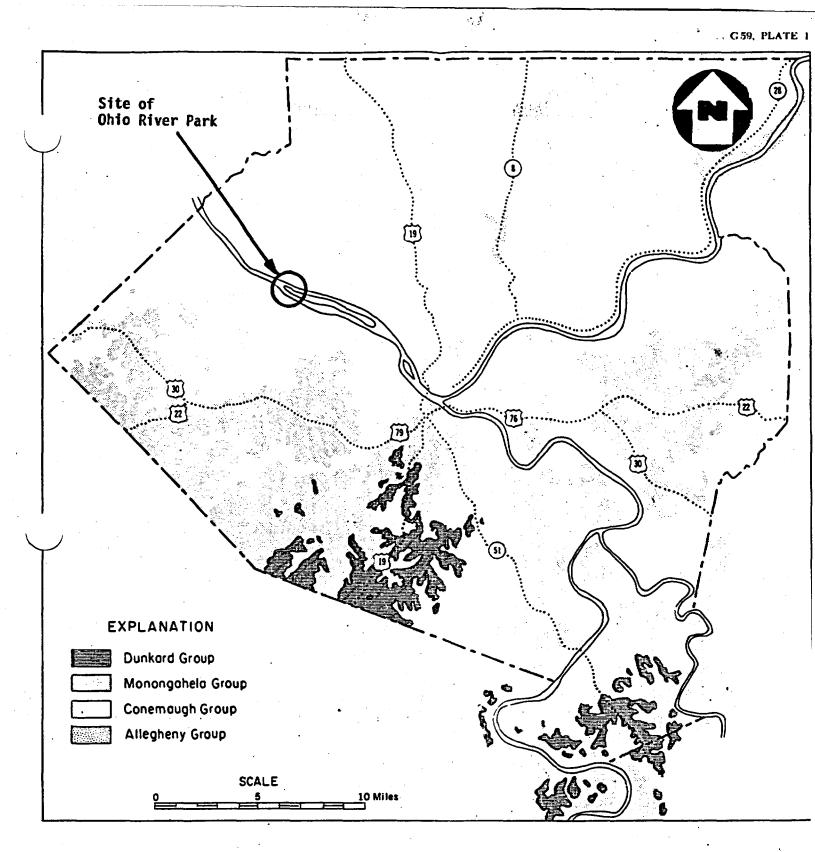
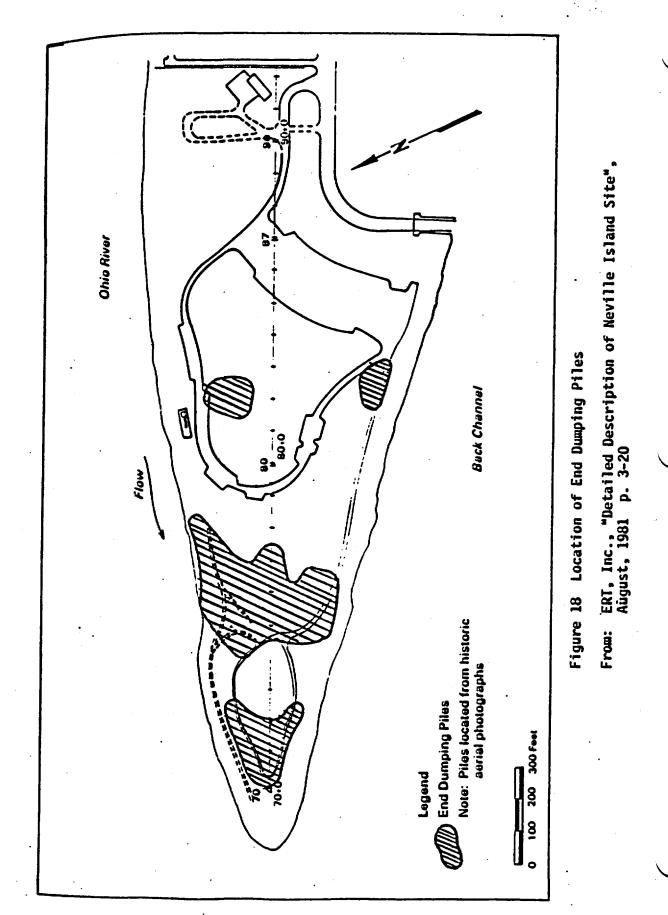


Figure 17

From: Walter R. Wagner <u>Geology of the Pittsburgh Area</u>, General Geology Report G 59, 1970. Plate No. 1.



Appendix C

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Table 1

OHIO RIVER PARK ANALYSIS OF GROUNDWATER SAMPLES FROM TEST WELLS:

6.9 **×0.2** 90,000 24,000 180 200 300 × 20 430 × 20 12 **V**10 s V 1× 20 20 49 14 **MH-6** From: Fred C. Hart Associates, "Assessment of Remedial Options at Ohio River Park" 6.3 <0.2 80,000 410 260 100,000 958 **Z** 200 480 V 25 20 66 80 V 20 33 300 NW-5 Concentration (ug/1) in Test Well × 0.2 3.5 75,000 25,600 4,300 210 120 A 60 150 2,900 605,000 20 **A** 40 60 <u> 00</u> 24 MM-4 6.8 16.5 ×0.2 78,000 13,000 410 ۷10 ۱0 140 100 36 **v** 20 160 < 5 < A 20 300 ۸ 20 MW-3 7.2 20,000 **v** 5 < 50 100,000 2021< < 20 < <20 < 680 s V 200 2020 27 30 MW-2 4 <a>0.2 6.3 43,000 11,000 80 150 A 20 A 10 202 .40 160 s V 20 V 9 9 1-1-1-1 0(0.02)^(c) 0(0.087) Water Qual Criteria 50^(d) 0.2 2 1,000 50 133 2 145 2 Phenolic Cpds. (Phenol) Parameter Carbon Inorganic (C) Chromium Total (Cr) Carbon Organic (C) Cyanide Total (CN) Beryllium. (Be) Antimony (Sb) Selenium (Se) Arsenic (As) Mercury (Hg) Cadmium (Cd) Silver (Ag) Copper (Cu) Nickel (Ni) <u>Meta</u>ls^(b) Lead (Pb) 퓜

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January, 1980, pp.42-46

		MM-6	•	100	410		ന	4	-8	•		'n	2	F 1		•		ŝ	•		
\bigcirc		MM-5 MM		100	910 4		,530	2	•		•	•	25 .	l		Ę		ſ	•	LI L	
·	n Test Well	MW-4 M		300	9,400		.	71	•	₹2	170		300	r	, NI	17 8		00	37	1,060	
	i (l/pu) noi	M E-M		<100	470 9,4		38 48,000	m	r	t	2	I	9	6	41	•		8 6,200	v	7 1,0	•
	continued continued contration (ug/1) in Test Well	MW-2 M		×100 ×	250		. 61	6			-	~ ~	, S	۲.		ŧ	V	S	۲. ۲	9	
	Table 1, conti	4	•	< 100 ×	190		2	4	6	യ	2	•	2	17 *		D	7	13	59	0	•
: : : :		Water Quality . Criteria (a)		4	5,000		0(15)	20 ^(e)	15,700	0(1.8)	0(2.1)	0.63	1,100	2			0(2.0)	17,400	0(21)		
		Parameter	Metals ^(b) (Con't)	Thallium (T1)	Zinc (Zn)	Volatile Organics	Benzene	Ch1orobenzene	l,l,l-trichloroethane	l,l,2,2-tetrachloroethane	Chloroform	l,3-dichloropropylene-cis	E thy 1 benzene	Methylene chloride	Dichlorobromomethane	Chlorodibromomethane	Tetrachloroethylene	Toluene	Trichloroethylene	Xylene	

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Table 1, continued

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			Conc	Concentration (ug/l) in Test Well	/1) in Test !	del 1	
Parameter	Water Quality Criteria (a)	I-WM	MW-2	MM-3	HH-4	MIV-S	MW-6
Phthalate Esters (cont'd)							
Diethyl phthalate	. 60,000	NT	2	e N	230	360	NT
Bis (2-ethylhexyl) phthalate		NT	6	25	18	28	NT
Butyl benzyl phthalate		NT .	ī	~` `	0	v	NT
Polychlorinated Biphenyls (PCB)				•			
PCB-1254 (Aroclor 1254)	0(0.2 ng/1)	NT	ı	0.20	•	0.64	 - 1
PCB-1248 (Aroclor 1248)	0(0.2 ng/1)	NT	ı	0.19	8.	0.54	ИТ
Other Organics							
Bis(2-chloroethyl) ether	0(0.42)	NT	1	•	ŀ	22	NT
Bis(2-chloroethoxy) methane		NT	ı	Ì	50 ::	9	NT
Isophorone	460	NT	1		011	7	NT
l,2-diphenyl hydrazine	0(0.4)	NT	v	t .		ı	NT
(a) Criteria for protection of human health. the Maximum Contaminant Level (MCL) spec Safe Drinking Water Act is given. All l given in ug/l unless otherwise noted.	th. pecif] lev	For 2,4-D ied in the els are	(P)	Water Quality Criteria hexavalent. The Water hexavalent chromium is target risk level of 1		for chromium other than Quality Criteria for 0(8 ng/l for an interim in 100,000).	nium other than Critería for for an interím 200).

Analysis performed on filtered samples. (q)

interim target risk level: concentration estimated to result in additional lifetime cancer risk of l in 100,000. Potential carcinogen. Numbers in parentheses give <u>(</u>)

- (e) For tainting -- 450 ug/l, toxic.
- Not detected. 1

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NT: Not tested for in sample.

	1 1							i					·:	
	MW-6	NT	NT	NT .	NT	NT	N		۲ ۱	N . TN	NT	NT	L. L.	
	4e11 	0.05	- 0.0	€0.03	41	8.3	1.5	•	ſ	' VI	1	- IV 58	~	
~ .	Concentration (ug/l) in Test Well 2 MW-3 MW-4 I	1.8	0.2	≤0.08	1,140	<0.05	<0.50		۱ (۲) g	v	-		m
	tration (ug MM-3	0.03	≰0.02 0.04	€0.01		<0.05	< 0.50	- -	- 7	الا الا	v	ন ।	~ ~	· · · ·
l, continued	Concen Mil-2	0.04	0.92	•	< 0.50	1.2		•	•	8	. 1	• •	. 2	v
Table l, co	L-WM	. IN	TN TN	TN :	TN TN	IN TN		NT	NT	NT	NT	TN TN	NT	NT
Ĥ	Water Quality 	:		100 / 421 /	100 (MCL)	•	•	20	•		-	143	•	
•	Parameter Pesticides (cont'd)	ot- BHC B- BHC	۲- BHC (lindane)	е- вис 2,4-D	Silvex	2,4,5-T	<u>Polycyclic Aromatic Hydro- carbons (PAH) and Other Coal</u> Tar Hydrocarbons	Acenaphthene	Acenaphthalene	Anthracene	Phenanthrene	Naphthalene Phthalato Ectore	Di-n-buty] phthalate	Ul-n-octyl phthalate

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Table 1, continued

Water Quality WM-1 MM-1 Criteria (a) NT 3.400 NT 30 NT 30 NT 30 NT 0.5 NT 30 NT NT NT 31 0.5 NT NT 32 NT NT NT 33 0.5 NT NT 30 NT NT NT 31 100 NT NT 33 0(0.046 ng/1) NT NT 34 0(0.098 ng/1) NT NT 0(0.098 ng/1) NT 0(0.23 ng/1) NT 0(0.23 ng/1) NT 0(0.23 ng/1) NT	Water (uality Criteria (a) Mu-1 Mu-2 Mu-3 Mu-4 Mu-4 <t< th=""><th>•</th><th></th><th></th><th>Concer</th><th>tration (ug</th><th>Concentration (ug/l) in Test Well</th><th>Well</th><th></th></t<>	•			Concer	tration (ug	Concentration (ug/l) in Test Well	Well	
3,400 MT - 2,300 9 0.3 NT - - 41 9 30 NT - - 91 9 30 NT - - 91 9 0.5 NT - - 91 9 0.5 NT - - 91 9 0.5 NT - - - 94 0.5 NT - - - 94 NT - - - - 12 NT - - - - - 12 000 NT - - - - 27 010 NT - - - - 27 $0(0.046 ng/1)$ NT - - - - 27 $0(0.098 ng/1)$ NT - - - - - - - $0(0.098 ng/1)$ NT - - - - <td< th=""><th>3,400 NT - 2,300 blenol 0.3 NT - - 41 blenol 30 NT - - 41 blenol 30 NT - - 41 blenol 30 NT - - 41 blenol 0.5 NT - - - - bytphenol 0.5 NT -</th><th>Parameter</th><th>Water Quality Criteria (a)</th><th>I-WM</th><th>MW-2</th><th>MW-3</th><th>MM-4</th><th>MW-5</th><th>9-WM</th></td<>	3,400 NT - 2,300 blenol 0.3 NT - - 41 blenol 30 NT - - 41 blenol 30 NT - - 41 blenol 30 NT - - 41 blenol 0.5 NT - - - - bytphenol 0.5 NT -	Parameter	Water Quality Criteria (a)	I-WM	MW-2	MW-3	MM-4	MW-5	9-WM
3,400 NT - 2,300 9 0.3 NT - - 41 9 30 NT - - 41 9 30 NT - - 94 9 30 NT - - 230 9 NT - - - 94 9 NT - - - 230 29 NT - - - 270 4 NT - - - 12 NT - - - 27 27 $0.0 NT - - 27 27 0.10 NT - - 27 27 0.0 NT - - 16 15 0.0 NT - - 63 1 0.0 0.0 - - 53 1 0.0 0.0 - - 50 - - 0$	3,400 NT - 2,300 0.3 NT - - 41 30 NT - - 41 30 NT - - 41 30 NT - - 41 0.5 NT - - 41 0.5 NT - - 41 0.5 NT - - - - NT N - - - - - 88.6 NT -	<u>Phenols</u>							
0.3 NT - - 41 9 30 NT - - 9 9 30 NT - - - 94 0.5 NT - - 20 29 0.5 NT - - 210 4 0.5 NT - - 212 27 86.6 NT - - 27 27 27 88.6 NT - - 27 27 27 88.6 NT - - 27 27 27 88.6 NT - - 27 27 27 100 NT 1 - 26 15 16 $0.0.044 ng/1) NT - - 60.01 - - 26 - - 0(0.048 ng/1) NT - - 20.01 - - - - - - - - - - - - -<$	0.3 NT $ 30$ NT $ 0.5$ NT $ 0.5$ NT $ NT$ $ NT$ $ 68.6$ NT $ NT$ $ 0.0006 ng/1)$ NT $ 0(0.044 ng/1)$ NT $ 0(0.098 ng/1)$ NT $ 0(0.098 ng/1)$ NT $ -$ <t< td=""><td>Phenal</td><td>3,400</td><td>NT</td><td>I</td><td>t</td><td>2,300</td><td>6</td><td>NT</td></t<>	Phenal	3,400	NT	I	t	2,300	6	NT
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NT - - 270 4 NT - - - 12 NT - - - 27 NT - - - 27 68.6 NT - - 27 68.6 NT - - 27 68.6 NT - - 27 61.00 NT - - 27 00 NT 14 - 63 1 00 0.044 ng/1) NT - 63 1 00 0.044 ng/1) NT - 63 1 00 0.098 ng/1) NT - $< < 0.01$ - - 01 01 0.42 0.02 - - - - - 01 NT - $< < < 0.01$ - - - - - - - - - - - - - - - - - - - - <td< td=""><td>$\begin{array}{llllllllllllllllllllllllllllllllllll$</td><td>2,4-dichlorophenol</td><td>0.5</td><td>NT</td><td>G</td><td>S</td><td>220</td><td>29</td><td>NT</td></td<>	$ \begin{array}{llllllllllllllllllllllllllllllllllll$	2,4-dichlorophenol	0.5	NT	G	S	220	29	NT
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0(0.23 ng/l) NT - ≤0.02	0(0.23 ng/l) NT - 0(0.23 ng/l) NT -	Endrin aldehyde		NT	0.10	0.07	I	3	NT
0(0.23 ng/l) NT - <0.01	0(0.23 ng/l) NT -	Heptachlor	0(0.23 ng/l)	NT	,	≤0.02	ı	,	NT
		Heptachlor epoxide	0(0.23 ng/1)	NT	·	<0.01	,	•.	NT

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Major Organic Compounds - South-Central Tract

• •	Approximate	
•	Maximum	
Organic Constituent	Concentration	General Range
	(ppb) ;	(ppb)
•	•	
Benzene	100,000	1,000 - 50,000
Toluene	30,000	100 - 6,000
Phenol	10,000	1,500 - 4,000
Chlorophenols		
2,4,6-trichloro -	10,000	50 - 150
2,4-dichloro -	8,000	500 - 3,000
2,chloro -	1,000	100 - 400
Napthalene	400	50 - 100
Ethyl benzene	100	20 - 30
Pesticides	•	
2,4-D	9,000	1 - 1,000
2,4,D-T	250	0.1 - 10
2,4,5-TP (Silvex)	140 .	0.1 - 10
•		

From:

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a: ERT, Inc.,

"Preliminary Risk Assessment of Neville Island Site", April, 1981, pp 4-40 and 5-6

Table 3

CHRONOLOGICAL SUMMARY OF NLC SITE HISTORY

	1930's	1940's	1950's	1960's	1970's	<u>1.980's</u>
Farmland ^(a)	×	x	۴		•	
Navy barracks		. x		•	x	
Topsoil removed for Forbes field		x	•		- 	
Municipal waste disposal	x		x	-		
Industrial waste disposal			x	x		
Ohio River Park construction					x x	
ERT, Inc. field s	tudies					×
		7				-

'a) Neville Island was reported as a fertile farmland in 1880 and was free from any industrial development until after World War I.

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From: ERT, Inc., "Detailed Description of Neville Island Site", August, 1981, p.3-26

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Table

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Parameter	Ponded Water (10/26/79)	Ponded Water (12/13/79)	Outfall No. 1 (10/26/79)	Outfall No. 1 (12/13/79)	Outfall No. (12/13/79
H	6.8	7.5	11.5	6.6	7.8
Total Cyanide	m	S	*_ V	- v	207
Carbon Organic	50,000	24,000	24,000	7,000	17,000
Metals and Other Inorganics	•	•		•	
Antimony (Sb)	<100 ∧	- 200	100	001 >	300
Beryllium (Be)	20	< 20		۰ . د	< 20
Arsenic (As)	Λ5	5	2	ы V	5 V
Cadmium (Cd)	۸ 10 ا	A 10	 10 10 	, 10 10	01
Chromium Total (Cr)	< 20	30 V	40	< 30	20 V
Copper (Cu)	50	< 20	40	< 20	< 20
Lead (Pb)	< 50	< 50	< 50		50
Mercury (Hg)	< 0.2	< 0.2	0.7	۸ 0.2	< 0.2
Nickel (Ni)	20	20	80	< 20	
Selenium (Se)	۲ 5	ы У	9	ы V	so V

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Table 4, continued

		C 0 N	C O N C E N T R A T I O N: (U9/1)	(l/6n) :N 0	
Parameter	Ponded Water (10/26/79)	Ponded Water (12/13/79)	Outfall No. 1 (10/26/79)	Outfall No. 1 (12/13/79)	Outfall No. 2 (12/13/79)
Phenols (continued)	•	·			
2,4-dinitrophenol	•	NT	•	ŃT	NT
2.4-dimethylphenol	t	NT	- v	NT	NT
Chlorinated Hydrocarbons			•		
Chloroform	8	7	v	4	2
Methylene chloride	12	8	NA	7	C.
Polycyclic Aromatic and Other Coal Tar Hydrocarbons		•			
Anthracene	e	NT	آلا	·· NT	NT
Fluorene	I	NT	۲	NT	NT
Phenanthrene	•	NT	1	NT	NT
Naphthalene	-	NT	61	NT	. TN
Phthalate Esters					
Bis (2-ethylhexyl) phthalate	73	NT	29	NT	NT
Di-n-octyl phthalate	8	NT	ł	NT	IN I

Table 4, continued

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and the second sec

Outfall No. 2 (12/13/79) M F Outfall No. 1 (12/13/79) CONCENTRATION: (ug/1) ž M Outfall No. 1 (10/26/79) V Ponded Water (12/13/79) .₽ ž Ponded Water (10/26/79) 9 Bis (2-chloroethoxy) methane Parameter 2.6-dinitrotoluene Other Organics

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Not available due to methylene chloride contamination of sample bottle. Interference (sulfide). Not detected. NT: *

"Not detected".

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NA:

Villa Samples		NiA Samples		÷		EAT Sound ca	•	
1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	Scampel or Low at from Despects		11.1	43574	Sumple Lucat ten Urpth	1.0-24	aurface	
Nincellanerus Analynin	Alac's	1 Lawrence	Harrillananan Analysis	1	Hurell	Hlucellancous Analymia	nu lynta	
2 2 25 250 255 10 - 21 270 - 51 270 - 51	ett Claritate Extension Loss on Lass on	3 1.41 1.42	2.8 212 192	4.4 U.ST 151		1.5 	1.5 260 260	200
Hend Space(n) Analysis	<u>15 A cranera</u>	tal And	Klemental Analynia (pph)	3	helninge Aub	• • •	• •	1.4
ı	2002	=	2	2	U 31 -	• •	• •	1.1
i .		~ 4	-2	• ±	3 72	. 1	• •	0.4 0.4
Trace - Truce Trace - Saul Trace - Tearr	Ngi Pras Sang Caul	2-4-	:-	2+4-	5 ktw/16(wet) ktw/16(Jry)			141
	14.14	-	•		Pent fell	Fentielde Analynia (ppm)	(mid) =1	
Vulutile Urganica Analynin			•		2.4-4	•	•	.
•					Silver I	lateriurene	, , z	••
••					Rolafklow	-	• •	• •
					Volattle Organica	Organtra	Analynia (pim)	(
					Benzerne Ted incom	12	• •	•
					Not hy Lone Chilor Ide	-	•	•
	• •			•	Set ann I n	Seraming Electrum Hirrowenpy (SEM) Analywiw	m Hirran Lyulu	Å L 11 31
				•				unud filter vich high s vich bigh s Hysol, er Hysol, er (Hu,)sol,
Beflact as the "elf" ubuve the a. Trimethylbensene Mumhthalean		oluer.						

	Table 12		•.			CKC to mp less	and the second					
Sample 172127 Lacartion 172127 Depth 30-40*	8-8 1.7-2.0'		TP-200 1.2-1.4		TF-2.0 1.5-5.5'	i	11-220 7.0'		TP-250 1-3.0°		TP-250 1.5-2.5'	
Mucellamous Analysta	Pesticide Analysis (pim)	Analysi.	(SEM Analywin			est le ld	Peuticide Analysis (ppu)	(ndd)		
ut 9.9 Chlurufern 9.34 Extructablua Less of 412 Iguleiun	2.4-h 0.64 2.4.5-T 0.16 511ver 0.94 Parathion 0.13 Mulathion 0.26		2,4-19 1.0 2,4,5-T 1.0 511ven 1.0 Parathim 0.5 Muluthim 1.0		Fine Caleton carbonate part letus containing sume gypuum	I C C	2.4-7 2.4.3-7 511vex Parechion Malechion	8 0 0 0 0	2.4-6 2.4.5-T 511vex Parethion Malathion		2,4-D 2,4,5-T 511vex Purathion Halathion	4.4 1.2 1.0 1.0
Clumical Analyuim (2)		Velay	Valut Ile Vrganic Analyuis	,			Vol :	1110 00	Volatile Organic Analysis (ppm)	yata (p	Ĩ	
5102 9.0 Fr ₂ 03 1.0 Cao 45.0		J I I I I I I I I I I I I I I I I I I I	Benzenu 6 Taluenu 10 Methylenu l Chlaride				Benzune 200 Tuluune 10.000 HeCl ₂ 15	000 000 1 5	henzene Talvene McCl ₂	- 11 -	Renzene Tolvene NaCl ₂	
•				•			SEN Analyala	_			. '	
P ₂ 0 ₅ 1.6 50 ₃ 1.0 Cu0 1.0					•		Fine calcium carbinate perticles containing some Bypsum	urticle. Jone	•			•
			TAT	Sample	FRT Samples (Cuntinued)							
Lucation 17-25) Depth B-10.0'	TP-275 5.0'		77-275 2.5-4.0°		TF-290 8-8.5°		17-290 5.5-7.0°		TP-500 2-3.0°		TF-500 2.5-3.0*	00
	Pesticide Analysi	(mng) ala	î			1	SEM Analysis	••	Pesticide Analysis (ppm)	Analys	is (ppm)	
2.4-0 5.9 2.4.5-T 1.0 Stitus 1.0	2,4-0 2,4,5-T 51)vex		2,4-D 2,4,5-T 5,11vea	12.0	2.4-0 4 2.4.5-7 1 Silver 1	-	Flae calcium cariumate particles containing some		2.4-D 2.4.5-T 511vex	40.0 2.1 10.1	2.4-0 2.4.5-T 511ve#	1.0
55	Parathlun Nalathlun	0.5	Pacat blow Halachton	0.15	<u> </u>	0.047	the set		Parachion Nalachion		Parathion Nalathion	
Volat []e	Volatile Organic Analysiu (ppm)		•						•		SEM Analysis	Į
Benzene 1 Toluene 2	Benzei Tolvene	-?					•				Mixture of Rypuum quarks, and some	Ryruu Bone
Mechylene 2	Nethylene	n									fron exide	

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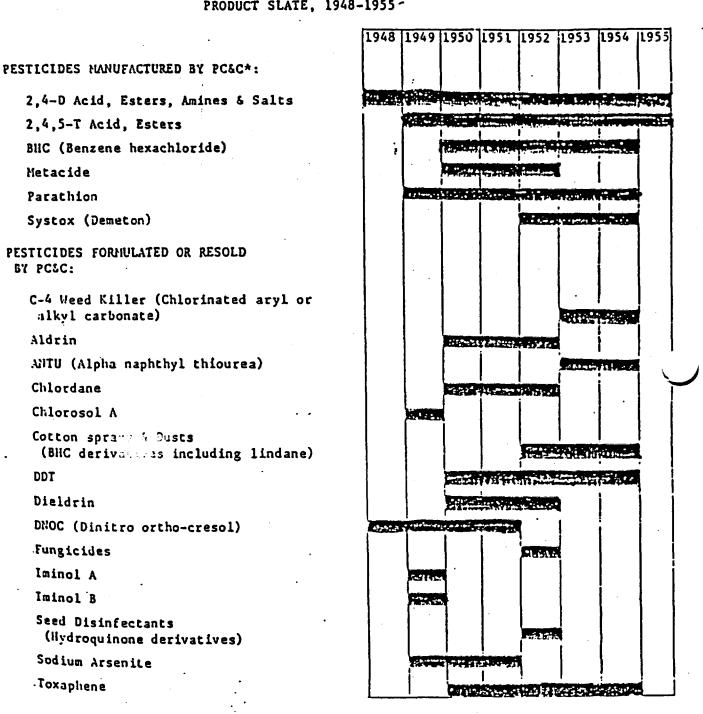
From: ERT, Inc., "Detailed Description of Neville Island Site", August, 1981, p.4-29

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AR100363

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Table 6



SUMMARY OF AGRICULTURAL CHEMICALS LISTED ON PC&C PRODUCT SLATE, 1948-1955-

Available evidence suggests that only these pesticides were actually produced by PCLC. Other chamicals were formulated or resold.

From: ERT, Inc. Detailed Description of the Neville Island Site, AR100364

21.0 A cerburu. cuntaining ceal. Loom dust and fuldepar.broken and burned plastic 2.2 2 TP-858 Muluture 8.0. Ulcimate Analysin (I) pleces. SEM Analysia 0.1 6.2 0.02 32.6 Saturated Saturated 5.5-6.5' 400 ppb 17-010 Hedica "Detailed Description of Neville Island Site" August, 1981, p.4-9 Halature Summary of Analytical Data for Coal Coking Sludges Bunzene Tuluene Xylenes TMB į Parathiun <0.18 Malmthiun <0.06 . 00C-4T 0.76 60.05 Hend Space Analysis 2 aturated Saturated acuraced Jaturated Methylene Chloride Extraction in Soil led I un 1.11 22 61 .0.1 19150 Smull' TH Nap (c) Smill Phenanthrene Anthracene 51 Lvex 2-7.2 Pusticide Analysis (pps) Fluntine Bunzene Kylenes **faluene** Indane Parathium 1.6 Muluchium «0.097 DBC Nup. 2,4-B 0.48 2,4,5-T <0.07 511vex <0.085 Ē TP-300 Saturated Mechylene Chloride Extraction in Soll Photo Com Larga Small Trace No trganles detected Talura. Xylenes THB Nap. Benzene Volatile Arganic 0.01 6.0 farathion of 5 Walathim et.d 12.0 Methylene chlorlde el.0 EKT Sumples Andlysis 8-31 6.0-0.0* ···· M9-94 H-1 [41] I.M. 1944 Suturated SILVIE 2-4.0 Large Large Small 22 Analysia (ppm) Toluune Xylunes THB 1000 2400 ç Bunsene Vulatile Organic 1-28 9.0-9.3 Methylene Chinride ī 20-28-11 in him me Saturated Saturated 160 ppm Nudium ERT, Inc., 2.e, Nued Spece^(a) Anulysis Nincelluncous Analysis Parachiun <0,84 Benzene Toluene Xylenes ThB 5.0 Malachiun <2.6 10.5-11.5 Table 7. 튚 1-2-4 8--28 I I vez 2.4-0 Sacuraced Sucuraced 25 PT (Traditional data and the second seco Large Nedium Smail ==;; From: Large Methylene Chloridu Extraction 65.0 Salathion <1.0 Pesticide Analysis Parachion <0. 6.3-6.7 Toluene Xylenes THB pli Phenul Jenzene Tulwine Nap. 4. S-T 8-8 -8 1 Lvin Q-3. 5-6.31 ۱. Paruthion 17.0 Muluchion 1(6) 0.00 Trace Trace Smull Smull frace 0.0 TACE 1 Plummathrene Naphahahau Acouphthene // I wor much hone Fluerine Chrysene Sumple Execution In-puts yrene 2.4.5-T Locution \$11vex Xylenea Thib(b) 2.4-0 Benzend ol wend Nap. (c) Indune Bepch E

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Table 7, continued

BTU/1b(wet) 10800 BTU/1b(dry) (f)

BTU/lb(wet) 2200 BTU/lb(dry) 3300

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EP Toxic Extraction Test (ppm)

ERT Sample TP-858 Ultimate Analysis (Z) (CONT) 9 % \$

Solut Ion: Benzene Toluene Xylenes

8

Solution: Benzene Toluene Xylenes

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9 9 9 9 7 9 8

Solid: Benzene Toluene Aylenes

> 2000 420

Solid: Benzene Toluene Xylenes

	•			$ \begin{array}{cccccccccccccccccccccccccccccccccccc$
	77125 TP-19 4.0-9.0'	2.8 202 372		23 23 21 16 16 1.5 1.5 1.5 7 equil
		pil Chloreforn Extractables Lous on Junifion	(=dd)	2 39 5102 14 5102 23 3 4 $Fe203$ 3 $Fe203$ 1 4 $Fe203$ 3 $Fe203$ 1 4 $Fe203$ 4 $Fe203$ 1 4 $Fe203$ 4 $Fe203$ 1 4 $Fe203$ 4 $Fe203$ 1 6 -1 $Fe203$ 4 -10 4 6 -1 $Fe203$ 4 6 -1 $Fe203$ 4 6 -1 $Fe203$ 4 6 -1 -103 4 6 -1033 4 7 -10333 4 7 -10333 4
ples	77124 77-3 0.0-4.0'		alysis	14 1 14 1 15 1 15 1 15 1 15 1 14 1 14 1
KCA Sumples	77124 77-3 1 72-3 1 0.0-4.0' Hiscelleneous Analysis	pH Chloroform Extractables Luss on Luss on Luss los	Cliumical Analysis (ppm)	 39 \$102 34 Fr203 4 Fr203 4 Fr203 4 Hr0 4 Hr0 4 Pr203 6 Cu0 9.0 S03 9.
	77124 77-3 3.0-7.5'	5.4 II II		$\begin{array}{cccccccccccccccccccccccccccccccccccc$
	Sumple Location Depth	pH Chloroform Extructables Lous un Ignicion		1102 39 12203 4 20 3.4 160 3.4 1.60 41 2.05 41 2.0 41 1.20 41 1.20 41 1.20 1.1 1.20 1.20 1.20 1.20 1.20 1.20 1.20 1.20

From: ERT, Inc., "Detailed Description of Neville Island Site", August, 1981, p.4-10 Charles and the

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TABLE 8

1. 4

SUMMARY OF ANALYTICAL DATA FOR FOUNDRY SAND

RGA	Sample		ERT Sa	mple	
Location (a) Depth	TP-32 0-4.0'		ocation (b) Pepth ;	B-5 15-16'	
Miscella	neoue Analysis		Pesticide Ana	lysis (ppm)	
pH Chloroform Extractabl Loss on Ignition	9.1 es 0.46% 51%		2,4-D 2,4,5-T Silvex Parathion Malathion	12 [.] 0.12 <0.08 <0.12 <0.71	
Elementa	1 Analysis (%)	_	•		•
S102	55		·		·
Fe203	• 6				
CaO	<1				•
MgO	<1	•	· · · · · · · · · · · · · · · · · · ·	• •	
P205	<1		•		

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From ERT, Inc., "Detailed Description of Neville Island Site", August, 1981, 4-13

AR100367

A of Dry Ash

Data Summar

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Table

34/680 4/80 13.5 <0.07/1 4.7 Volatile Organic Analysis (b) ----6300 10700 0 <u>Ultimute Analysis (2)</u> SEM Analysis 3.9-5.61 **B-30** Molsture Content (net) DTU/1b (dry) Toluene Benzene Xy lenes BTU/1b Ash 5 t n ERT Samples (a) <0.5 ī 47 ī \$ \$ 7 7 5-10.0' 11-5 **Hutly Jone** Parathion **Nul athlon** Volatile Organic Analysis (ppm) Chlor ide 2,4-D 2,4,5-T Tolucine Bunzene SILVEX Pesticide Analysis (ppm) 3000 6000 5.2 Ś <0.5 3.5-3.4 7 Ļ 2-1 ocation Methy]ene Parathion Malathlon Depth Chloridu 2,4-1) 2,4,5-T Benzene Toluene Silvex

clinker maturial, coal impurities, sulfur crystals, aluminum and silica present

> Only ERT sampled dry ash. ઉઉ

The first number is the result from UPLC analysis of EP extract (µg/ml) The second number is a calculated concentration for the solid $(\mu g/g)$. "Detailed Description of ^august, 1981, p.4-16

Neville Island Site",

From: ERT, Inc.,

Table 10. Analytic Data Summary of Industrial Slag

222 Volatile Organic Analynis (ppm) TP-800 1.5-2.5 Benzenu Tuluene Methylene Chloride iigh temperature real taining irms, sulfur, and aluminum <u> Ittimite Analynin (2)</u> 0.01 1650 2510 0.16 we rich in enicium and milica, forming srittle composite co e 26.1 0.31 0.099 0.099 0.13 24.6 SEM Analyula TP-800 2.5-3.5' BTU/II (uci) BTU/Ib (dry) EKT S Imples 2,4,5-T 511vex Parachion Malachion Hulsture 9-3.5 <u>Peritetue Andruin (pp</u> 8--24 4 **- 5-**-6 - 5 -2,4 · 10 2, 4, 5 · 1 SI Iven Parathim Mahathim M - Nut detectable Scomple – Lauration 18-25 Inspile – 18,5-18,6* 0.13 0.13 0.15 2,5-D 2,5,5-F Stivex Parathion Molathion Emineion Spectography (2) 3 ŝ Chemberth Analynta (pem) X-RAY Speet Concention (2) 9.6 9.6 1-1 1 ŝ <u>-</u> E-1 01-1 9**-**-= <u>=</u> ₹. Ŧ τ Ê Ē êêê ê £ pli Louis on ignition 1.51 F. 22 F⁴20 r, 05 510, Miscellaurous Analysin ຣິວ Fill'A Sampelers Falution Spectography (2) 1n7 X-Kay Spectrography (1) 9.0 0.0 2 Lunn on Ignitiun 01-1 <u>- 10</u> <u>ei-i</u> 01-1 01-1 9, 510 7 Ţ Ŧ Ţ Ţ £ 2 2 2 ĝ Sample Location Depth M10 Zn0 Cr2⁰3 Flu0 20°2 Fe203 N320 ر⁰2 ۱۸ 510] 5002

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From: ERT, Inc., "Detailed Description of Neville Island Site", August, 1981, p.4-18

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TABLE 11

SUMMARY OF ANALYTICAL DATA FOR LEACHATE SAMPLES

FCHA SAMPLES

	Concentrat	ion (ppm)
<u>Metals and Pesticides</u>	Sample 20 Location 13 Depth 4'	25 7 3.5'
Arsenic (As) Cadmium (Cd) Chromium, total (Cr) Copper (Cu) Cyanide, free (CN) Cyanide, total (CN) Lead (Pb) Manganese (Mn) Mercury (Mg) Nickei (Ni)1 pH Phenolic Compounds Selenium (Se) Silver (Ag) Zinc (Zn) Antimony (Sb) Beryllium (Be) Thallium (T1) Parathion 2,4-D 2,4,5-T Silvex	$\begin{array}{c} 0.075 \\ < 0.01 \\ < 0.02 \\ 3.2 \\ \hline \\ 0.015 \\ 0.19 \\ 4.8 \\ 1200 \\ 0.05 \\ 7 \\ 2.6 \\ < 0.02 \\ < 0.02 \\ < 0.02 \\ < 0.02 \\ < 0.1 \\ < 0.02 \\ < 0.1 \\ 4.5 \\ \hline \\ - \\ - \\ - \\ - \end{array}$	<pre><0.1 <0.01 0.04 0.23 Present 1.32 0.23 16 2.6 0.16 7 36.5 <0.1 <0.02 0.35 0.4 <0.02 0.19 0.039 48 1.2 0.05</pre>
Organic Parameters	Concentrat	<u>ton (ppb)</u>
Acenaphthene Benzene Carbon tetrachloride Chlorobenzene 1,1,2,2-tetrachloroethane 2,4,6-trichlorophenol Parachlorometa cresol Chloroform 2-chlorophenol 2,4-dichlorophenol 1,2-dichloropropylene, cis Ethylbenzene Fluoranthene Methylene chloride	420 4 8 6 - 22 200 - 23 240 2 115 50000	54 5100 2 8 5 2400 - 90 800 15000 - 43 150 210 A D

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SUPPORT DOCUMENTATION FOR THE REVIEW OF ORGANIC ANALYSIS LAB DATA PACKAGE

CASE:	Ohio River Park (January 1981)
TYPE OF	ANALYSIS: Volatile, BNA, Herbicide
CONTRAC	T LABORATORY: ERT (Subcontractor)
REVIEWE	R: Sam Fring
REVIEW	DATE:

APPLICABLE SAMPLE NO'S .:	614, 8040, 8054, 80t
fost9, for6, 10237, 10	240, 10239, 1-268;
1-769, 1-295, 1-296, 1	
10/79, 10181, 10182,10	184,1-185,10186, 10206
1-2=7, 10205, 10209, 10234	1-235 1-226 102271-22
10229, 1-187, and 10223.	

THE FOLLOWING TABLE INDICATES AREAS WHICH WERE EXAMINED IN DETAIL, THE IDENTIFIED PROBLEM AREAS, AND SUPPORT DOCUMENTATION ATTACHMENTS:	Cho	EAS EXA IN DETA eck 1f otnote Commer	VIL yes or letter	-	, Ci	ROBLEM IDENTI heck 11 footnot	FIED yes o te numt	ber	·	Check	NTATION HMENTS if yes ntify	or
	volatile	Acid	BIN	Herbicide	Volntile	Acid	N/51	Herbitide	Volatile	Acid	B/N	Herbicide
HOLDING TIMES	V	V	~	V	V	· · · · ·		V	V	V	~	$\overline{1}$
BLANK ANALYSIS RESULTS: TARGET COMPOUNDS	IV.	V	~	Y								1
BLANK ANALYSIS RESULTS: TENTATIVE IDS.		·										
SURROGATE SPIKE RESULTS	V	V	V	\mathcal{D}	V	V	V	\bigcirc		V		
ATRIX SPIKE RESULTS	\square	\bigcirc	\mathcal{O}	V	\bigcirc	\mathcal{O}	\mathcal{P}	V.				1
DUPLICATE ANLAYSIS RESULTS	\bigcirc	\odot	\bigcirc	V	\square	\mathcal{O}	\bigcirc	V.				
TARGET COMPOUND MATCHING QUALITY	V	2	V	V					V		\checkmark	
TENTATIVELY IDENTIFIED COMPOUNDS		V	~									
DFTPP 8 BFB SPECTRUM TUNE RESULTS			V	(¢)	<u>.</u>						~	
GC INSTRUMENT PERFORMANCE	$\downarrow \downarrow$									<u> </u>		1_
INITIAL CALIBRATIONS	$\mathbf{\hat{e}}$	2	\bigcirc		<u> </u>				ļ	I	<u> </u>	
CONTINUING CALIBRATIONS	<u> </u>	V	V	V		<u> </u>		<u> </u>		ļ		Ŀ
QUANTITATION OF RESULTS		$\perp \checkmark$					·		ļ	<u> </u>	ļ	
OTHERS Congregated Confirmentions	1 .		i	(3)		l	L	(3)				I

(Not applicable .

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•/ please see the Analytical Quality Assurance section of this report. delected Nore cletrited REMARKS 1981 SITE NAME Ohio River Park Nac January DATE OF SAMPLE 1-2/L 2000 1300 32 D 1900 D 400 100 D 160 D 160 D 10 D his(2. Ellistheryl) philalate > 000 20J DETECTED 140 R Bue R 410 J P3 R 37 R 718 928 90 J 478 24,6-Trichlaropher COMPOLINDS Naphthalese INORGANIC Youch Youch Ilook 63R 24-Dichlorophenol 2-Chlorophenol of this data and non-target, tentatively identified compour Π 93 R 3920 R 32 8 2508 Phenol Ethylbenzene ORGANIC 10000 SPUK 6700 R 900 R 141 เริ่ Toluene [ano/<] 7/84 Benzene D 7/51 7/8-1 Units P3/L どう 7/6N JIEN 처 wite ייזיאט White Lutu 111 Phase Later **Vatu** Blank (field) + 1 Block (field) #2 Sample Description ER7-205 ERT-185 CUTFALL # and Location 4-M~ ERT-175 ロバーー يەتەر 12 Number ARIOO Sample 6503 et.o 45.28 8-24 6903 325

SAMPLE DATA SUMMARY TARGET COMPOUNDS

				14. az																					-
	981)	•	· .		-				• :					•							(_	39 (†	*) *)*	Nal 1) 1.:	3
	SITE NAME <u>Ohio River Park (Jannary</u> 1981) OF SAMPLE 1/13/81 - 1/28/81		1 4 4 5 4 5 10 10 10 10 10 10 10 10 10 10 10 10 10	REMARKS																				ed compounds, please see the Analytical Quality Assurance section of this report.	
•	<u>Ohio River</u> 1/13/81 -					÷	+						-					-				-	-	urance sec	
	NAME <u>Oh</u> MPLE <u>1/13</u>	-				╎╾┥							┟╴											ality Assi	<u> </u>
•	SITE NAME Date of Sample			<u> </u>				. 		-	-	+				-		<u> </u>	-		-		_	tical Qu	
	D	DETECTED	· ·			 -	+						<u> </u> 										-	he Analy	
8	AHIC	COMPOUNDS										ŀ												se see t	
Harbicides] INORGANIC	3					-			•		+	 -					$\frac{1}{1}$	+					lds, plea	1 1
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•	lic sides		2,4,5-T		4 1,3 N	V 0.16 N			NC1.0 N CC0	N #11	N 23	2	NZO		95N	I agen					1 266 N	N920	N 64		
-) ORGANIC Herbicides		Silvex (2,4,5-1 2,4-D	7P) 3	2	2	1.5 N		N 7.50 N 77 - C	2000N	NorSz	NAK	1,6 N	N CO NOI			29N ISN	<1.6 R	-	V4N C/1N	N CHI IN DEI		92co N	ively 10	
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•	· · · ·	•			44-5 ERT-20D	ERT-20C	ERT-12 D		ERT-6D			ERT-3S	ERT - PS	ERT -95	EKT-90	ERT-10M	ERT- 10D	ERT- 160	ERT-160 0	ERT- 19M	ERT-19 D	MW-4	ERT - 18 S		Nova: For a review of this data and houstanget, tentering
	•		•	Sample Number	10251	142 01	10207	10269	10 295	4/20/	10308	10309	10179	19101	C8)0	10185	10186	9211	10207	10208	0°.7			03	

SAMPLE DATA SUMMARY TARGET-COMPOUNDS

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·		SITE NAVE Ohio River Park (January 1981) OF SAMPLE 1/13181 - 1/28/81			REMARKS																50 20		tion of thy report.	
	·	<u>Ohio River</u> 1/13/81 - 1			-																		please see the Analytical Quality Assurance section of th	
	. •	site Nave _	-	•																			cal Quality /	
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hple data summary target-compounds-	Herbicides (continued)	INORGANIC	N				╈		╋	┢		╈						╋		+	╋		olease	
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	•	ORGANIC Herbicides		Silvex (1,4,5-TH		0.SeN 0	C42N	+		+				┼╢	╉	+-	┠╌┨	╉	╏		+	┼┤	1 dent	•
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	•				Phase	Water	MIT																dog pu-	
. :					Sample Description and Location	ERT-18M FRT-18D	ERT- 19 S			-									•					
			•	• • •	Sample Number	1420	9441		ļ							A	R	0	0	37	4			Note: For a

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PESTICIDE-	BRAUTT-	a f							:			T	T										-				-				
	2.4.5 TRIBROWO	(14-123)																										c limits C limits			
	1	(801-12)	598	1354	754	*~5	41.5																				- 5 outside of QC limits	; outside of QC limits : outside of QC limits			
	91- 10H3H4	(¥++)	29.9	2.44	36.6	291	223																•			-	10	30			
SEM-VOLATHE														T													La Ma		1.		
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	1.111)	(811-62)	88	46.2	43. P	49.9	41.3																				ITS		•		
	-041M	(11-114)	54	39.2	<u>ہ</u> *	• *	* 0																				RED OC LIMITS				
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	TOLUENE-DA	(811-90)	11.*	119 \$	100	106/102	99													-							RE OUTSIDI	LIMITS ON	•		
ت ب	are Are Are		22	8-74 / 383		Prs7/1865 106/102	S. 86/256										-										* VALUES ARE OUTSIDE OF CONTRACT REQUIRED QC	** ADVISORY LIMITS ONLY	1	Comments:	
				EK7-1 2		h-NW		<u> </u>		_1_		<u> </u>	ما در ب	1	3-1	.1			<u>e</u>	•	• <u> </u>	4	<u> </u>	<u>ل</u> ے۔	#	`		#			

WATER SUBROGATE PERCENT RECOVERY SUMMARY

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1/21/81 +1/22/81 HERBILIDE RECOURRY HERBICION SPIKE LEVIELS ENPRET 1.0 ug/2 level soont Alignot = 0.504grecovery + Sml finatuol = 0.10 ug/ml 1.0 yg/ml _ 5.0 uz 10.041Z 25.00 5.043/ml ____ 50.049 1R ... 0.0 BLONK soon water COMPARE RECOVERY SULA to STIMAMO SOLN MADE to EXPERTIS LEVIL .0.3391 Suver 1.4816 1.5621 1.5219 2,4,5-T 1:3979 1.4604 1.4242 -----<u>_____RPD</u> 157% * 300 0.4037 119% × 27 1.043 1 50 A D 9.5342 98.8%. 1.2296 B0.8%. 20076 t 5 1.5020 T. 1.4219 100 % # 0.9818 68.7% 37.1% # sta-1.2.7.673. 2.7389 2.753 2.7404 2.9500 5 13,041 12.9043 11.5776 12.8974 14.1036 12.3859 10.5318 12.1062 T 12.1971 13.1500 ____ RPD 56.0% * RECA 2.2525 ____ 82.2% ___ 37.9%. _____ 1.4-D___ 10412 226A 1.5335 13:5347 104:9.64 0.8% 5:1vex 13:4547 104 13 *-72,80%. 2.45-7 7.5728 62,2013 * 11.2953 39.5% * STD_ 5.042/ml _____ 11.4054 ____ 11.0176 11.2115 50.2451 48.2176 49.2313 49,7472 50.8364 48.6571 RPD 2.4-12 504318 A 8.5036 15 521" * 36.0% * 42.000 3 17.3% Silvex 71.17. * 22.2% * 35.0348 2,4,5°T 38.6432 77.18 28.6796 57.5% × 29.5% × QC_Limits_____Recovery____ RPD 2,4-0 63-87 47. - * Values are cutsiAP=110097815-Silvex 73 -103. 5%

	10. Dhis River Pack Contractor ERT (Subcen	A Chever Contract NO	بني من
Instru	ment ID Date/23/8/	Time 8:09	
Lab II	D Data Release Authorized By	/:	
m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE	•
51	30.0 - 60.0% of mass 198	55,13	
68	less than 2.0% of mass 69	0.67	(136)
69	mass 69 relative abundance	49,10	
70	less than 2.0% of mass 69	0	(0)
127	40.0 - 60.0% of mass 198	41.45	1
19 7	less than 1.0% of mass 198	0	
198	base peak, 100% rélative abundance	/00	
199	5.0 - 9.0% of mass 198	5.68	
275	10.0 - 30.0% of mass 198	20.83	
365	greater than 1.00% of mass 198	1.97	
441	present, but less than mass 443	8.67	
442	greater than 40.0% of mass 198	66.61	•
443	17.0 - 23.0% of mass 442	1.2.24	(18,38)

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

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¹Value in parenthesis is % mass 69. ²Value in parenthesis is % mass 442.

SAMPLE ID	LAB ID	DATE OF ANALYSIS	TIME OF ANALYSIS
RT-185	Fo54/3654	1/22/81	12:48
MW-4	8059 13655	1/23/51	14:03
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asa N	to Unit FATE Contractor ERTISCO	Contract	No
nstrui	GC/MS TUNING AND MA Decafluorotriphenylph No. Ohio River Park Contractor ERT(Side ment ID Date/27/81	Time	9:05
	D Data Release Authorized E		· _ · · · · · · · · · · · · · · · · · ·
m/e		%RELATIVE ABUNDANC	E
51	30.0 - 60.0% of mass 198	61.61	meets entended cr
68	less than 2.0% of mass 69	0,19	(0.3 <u>9</u> 1
69	mass 69 relative abundance	56.58	
70	less than 2.0% of mass 69	0	(0)'
127	40.0 - 60.0% of mass 198	44.19	
197	less than 1.0% of mass 198	0	•
198	base peak, 100% relative abundance	100	
199	5.0 - 9.0% of mass 198	6.69	
275	· 10.0 - 30.0% of mass 198	20.38	·
365	greater than 1.00% of mass 198	1.35	
441	present, but less than mass 443	6.44	
442	greater than 40.0% of mass 198	57.71	
443	17.0 - 23.0% of mass 442	10.35	(2,y) ²



SAMPLES, BLANKS AND STANDARDS.

²Value in parenthesis is % mass 442.

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	TIME OF ANALYSIS	DATE OF ANALYSIS	LAB ID	SAMPLE ID
-	2:54	1/28/21	8060 / 3662	EX7 -205
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Instru	ment ID Date	ERT (Subcontinutor) Contract No
Lad II	D Data Relea	ase Authorized By:
m/e	ION ABUNDANCE CRITERIA	*RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	57.89
68	less than 2.0% of mass 69	0 (0)
69	mass 69 relative abundance	67.64
70	less than 2.0% of mass 69	0 (0)
127	40.0 - 60.0% of mass 198	43.24
197	less than 1.0% of mass 198	0
198	base peak, 100% relative abundance	/00
199	5.0 - 9.0% of mass 198	5.89
275	10.0 - 30.0% of mass 198	16.36
365	greater than 1.00% of mass 198	0.94 *
441	present, but less than mass 443	P.53
442	greater than 40.0% of mass 198	63.73
443	17.0 - 23.0% of mass 442	11.66 (18:39)

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SAMPLE ID	LAB ID	DATE OF ANALYSIS	TIME OF ANALYSIS
RT -185	8-54/2654	1/27181	٤٦:21
MIX-4	8-59/3655	1/27/8/	16.14
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Instru	ment ID	Decailuorotriphen Buer Park Contractor ERT 5 Date 2/4/0	81	Time	6-19			
		Data Release Author						
m/e			•	RELATIVE ABUNDANC	E E			
51	30.0 - 60.0	0% of mass 198		75.05	*			
68	less than 2	2.0% of mast 69		1.67	(2,•8) ¹			
69	mass 69 re	elative abundance		80.25				
70	less than 2	2.0% of mass 69		υ	(o) ¹			
127	40.0 - 60.0	0% of mass 198		50.88				
197	less than 1	1.0% of mass 198		0				
198	base peak,	, 100% relative abundance		100				
199	5.0 - 9.0%	i of mass 198	·	5.61				
275	10.0 - 30.0	0% of mass 198		15.6	2			
365	greater that	an 1.00% of mass 198		1.32				
441	present, b	out less than mass 443		5.95				
442	greater that	an 40.0% of mass 198		49.69				
443	17.0 - 23.0	0% of mass 442		8.37				
		TUNE APPLIES TO THE FOLLOWING ID STANDARDS. $\# f_A$		¹ Value in parenthesis is % mass 69 ² Value in parenthesis is % mass 44				
SAMP	LEID	LAB ID		DATE OF ANALYSIS	TIME OF ANALYSIS			
UTFA	121	8086 13826		2/4/81	2/:40			
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		River Park Contractor ER						
Instru	ment ID _	<u>5</u> Date <u>2/11/2</u>	8.1	Time	18:23			
Lab II	D	Data Release Autho	prized By: _	<u></u>				
<u>m/e</u>				*RELATIVE ABUNDANC				
51		.0% of mass 198		74,04	*			
68	less than :	2.0% of mass 69		0	(o) ¹			
69	mass 69 r	elative abundance		78.01				
70	less thari	2.0% of mass 69	·	0	ا(ق)ا			
127	40.0 - 60.	.0% of mass 198		5043				
197	less than	1.0% of mass 198		3.14	*			
198	base peak	, 100% relative abundance		100				
199	5.0 - 9.09	6 of mass 198		5.62				
275	10.0 - 30	.0% of mass 198		14.49				
365	greater th	an 1.00% of mass 198		0*				
441	present, b	out less than mass 443		6.04				
442	greater th	an 40.0% of mass 198		47.27				
443	17.0 - 23	.0% of mass 442	<u> </u>	7.61 * (16.1.5)2				
		TUNE APPLIES TO THE FOLLOWIN ND STANDARDS. + Fail	G	1 Value	in parenthesis is % mass 69. In parenthesis is % mass 442.			
SAMP	LE ID	LAB ID		DATE OF ANALYSIS	TIME OF ANALYSIS			
EK1 -	205	8060 / 3662		2/11/81	20:50			

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7/85 Form V

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•		Bromofluoroben)			
lase N	No. Ohio Ria	erPark Contractor ERT (Subo	mps Chem)	Contract No.			
nstru	ment 10 2	Date Date		Time 16:30			
<u>ا مد</u>)	Data Release Authorized I	Βγ:				
n/e			XRELATIV				
	[2640			
50	├ ─────	if the base peak					
	<u> </u>	f the base peak	_	50.49			
5		3% relative abundance		100			
8	5.0 - 9.0% of 1			5.91			
73	}	i of the base peak		0			
74		0.0% of the base peak		81.97			
75	5.0 - 9.0% of a			436 (532)1			
78	 	5.0%, but less than 101.0% of mass 174		P1.57 (99.5)			
77	5.0 - 9.0% of	masa 178		4.07 (4.99) 2			
	FORMANCE TU B, BLANKS AND	INE APPLIES TO THE FOLLOWING STANDARDS. + Fail		¹ Value in parenti ² Value in parenti	hesis is % mass 174. hesis is % mass 176.		
SAN	APLE ID	LAB ID	DATE OF AN	ALYSIS TIME	OF ANALYSIS		
EK1	-185	8054 / 3654	1/22/81		21:05 V		
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		Bromofluorobenz Park Contractor ERT(Sub ERT(Ce Date 1/3-/8-1	contractor)			
	-	Data Release Authorized By	•	<u>-</u> -		
m/s		CRITERIA	RELATIVE ABUNDA	INCE		
50	15.0 - 40.0% of the	base peak	29,35			
75	30.0 - 60.0% of the	base peak	51.90			
9 5	Base peak, 100% rel	ative abundance	100			
96	5.0 - 9.0% of the bi	se peak	7.70			
173	Less than 1.0% of t	ne base peak	0			
174	Greater than 50.0%	of the base peak	74.67			
175	5.0 - 9.0% of mass	74	. 4.58	(6,13)		
176	Greater than 95.0%	but less than 101.0% of mass 174	73,33	(98,21)]		
177	5.0 - 9.0% of mass	176	4.24	(<i>5.78</i>) ²		
IS PEF	FORMANCE TUNE	APPLIES TO THE FOLLOWING	¹ Value in	parenthesis is % mass 174.		

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	TIME OF ANALYSIS	DATE OF ANALYSIS	LAB ID	SAMPLE ID
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	7/85 Form V	AR100383	B-16	

Instrum	ient ID	<u>Parke</u> Contractor <u>Subcontruiter</u> <u>Date</u> <u>1/22/8/</u> Data Release Authorized B	Time2					
		Data Release Authorized E						
m/s			By:					
-	ION ABUNE	ANCE CRITERIA	SRELATIVE ABUND	ANCE				
50	15.0 - 40.0%	of the base peak	2/.6	5				
75	30.0 - 60.0%	of the base peak	39.8	9				
95	Base peak, 1	00% relative abundance	100					
96	5.0 - 9.0% of	the base peak	50	4				
173	Less than 1.	% of the base peak	υ	υ				
174	Greater than	50.0% of the base peak	65.19					
175	5.0 - 9.0% o	mass 174	3.42 (5.251					
176	Greater than	95.0%, but less than 101.0% of mass 174	61.72 (9462)					
177	5.0 - 9.0% a	f mass 178	3,	2	(5,06) 2			
		TUNE APPLIES TO THE FOLLOWING D STANDARDS. $# f_a; I$	¹ Value i 2 Value i	n parenthe: n parenthe:	sis is % mass 174. sis is % mass 176.			
SAMP	LEID	LAB ID	DATE OF ANALYSIS	TIME C	F ANALYSIS			
MN	-4	8-59/3655	1/23/81	2:08	/ 3:58			

B-16

AR100384

7/85 Form V ţ

Case No Dete Date Date				trade (<u>willham</u>) Contrac	OFIGINAL (P.05) 1 1:13		
	· · · · ·	Data Release	•				
m/s	ION ABUND	ANCE CRITERIA	·	RELATIVE ABUND	ANCE		
50	15.0 - 40.0%	of the base peak		19,5	٤		
75	30.0 - 60.0%	of the base peak		41.4	6		
9 5	Base peak, 16	00% relative abundance		100			
96	5.0 · 9.0% of	the base peak		5,59			
173	Less than 1.0	% of the base peak	· · ·	0			
174	Greater than	50.0% of the base peak		90,19			
175	5.0 - 9.0% of	mass 174	•••	46	o (5./•)}		
178	Greater than	95.0%, but less than 101.0	% of mass 174	91.1	ويعور الشروي فسيبد والتسبيب الشفاع فتعف المتبه والشفاع المتبه		
177	5.0 • 9.0% of	f mass 176		5.0	×3 (5.92) ²		
		TUNE APPLIES TO THE FO	ollowing *fail		in parenthesis is % mass 17 in parenthesis is % mass 17		
- SAI	MPLE ID	LAB ID		DATE OF ANALYSIS	TIME OF ANALYSIS		
ERT -1 5074		£074 / 3833		2/10/81	4:45		

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7/85 Form V

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AR100385

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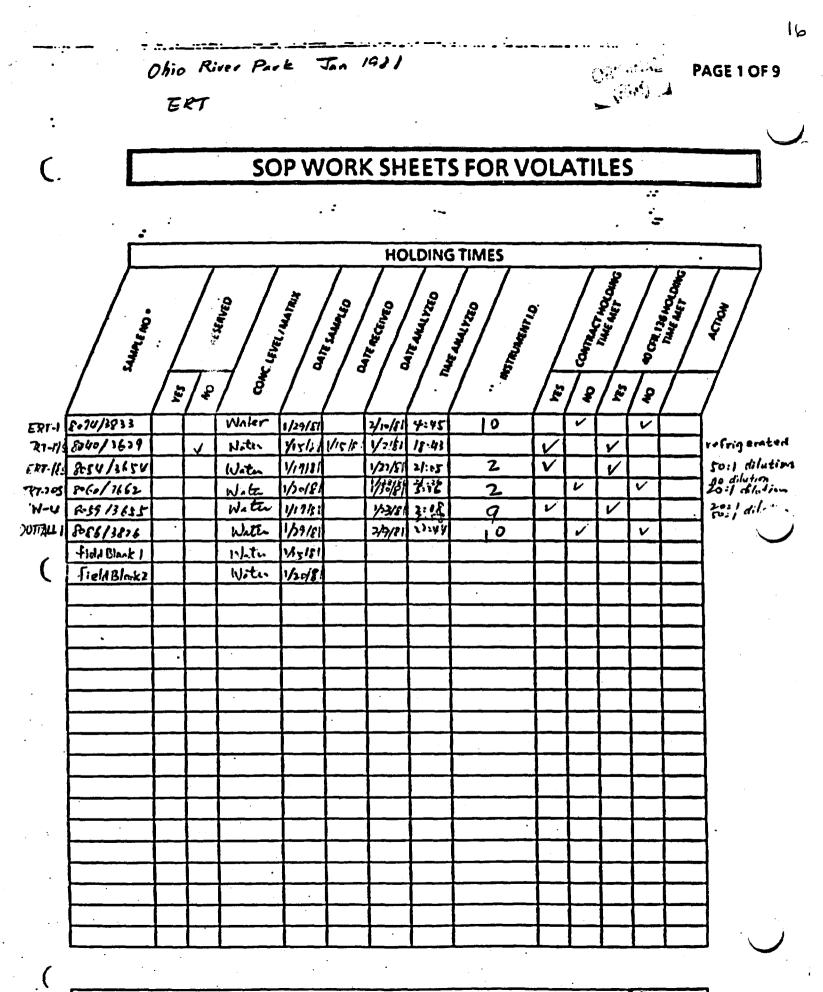
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* INCLUDE MATRIX SPIKES, BLANKS AND RE-RUNS HERE AR 100386

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	C. A. 1. 022	Water	1/24/80			2/13/8	15:46		+		-	[Acid
-77-1 -RT-1	Fe14/257.3 E074/3823		1/29/81			3/5/81		<u> </u>					- BIN
· L	5-40/3623			1/15/8	Viela				17				Aeid
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	2054/215V	Nati.	1/19/81	<u> </u>		1/2 7/51	15:53	5			1		Acid
7	205U/1ASU		1/17/8			1/23/81		4	V		1	· ·	B/N
	Folo/3602		1/20/81			2/1/11	70:5+	ζ			1		Acid
	8-6- 13662		1/20/51			178/81	2:5¥	ú					B/N
/	8059/3655	Water	VigR			1/27/81	16-34	5					Acid
/[8059/3155	NAteu	1/19/21			1/22/81	14:•3	4	\vee	V			R/N
1152121	8+86/3176	Neta	1/29/5			2/4/8/	21:4+	5	V	V	1		Acid
1FALL /	8-86/31-6	Water	1/29/81			45/81	15.42	3	V	<u> </u>	′·		B/N
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* INCLUCE MATRIX SPIKES, BLANKS AND RE-RUNS HERE

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	LA	UNIS KINGI FARE BORATORY <u>FRT</u> Jan	1981
		SOP WORK SHEETS FOR PESTICIDES /	PCBs
,		Co _{llan}	. 1
		HOLDING TIMES	··.
	Salange No.		⁴⁰ CM 136 MOLONIC
			8
·1W-5 [10237	Water 1/191 1/20/2 2/20/21 1+34	
RT-2.0	10240	- 1/19/8/ 1/2-/8/ 2/5/2	
ERTZOD DUP	10241	··· 1/14/51 /20/81 2/5/81 2/19/81	
RTJAN	1-239	··· 1/19151 1/2018 2/5181 2/19181	
ERT-120 -4W - 2	10269	·· /2-/21 //2/21 2/5/2 1/2//24 1640	
40	10295	·· 1/21/11 1/25/61 2/21/51 11:12	<u>+</u> +
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UTIALLA	10309	· 4/28/68 4/29/28 2/21/68 9:59	
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OUP TRT-35	10309	· v> \$/\$ 1/31/2 1/21 1/21/2 -	
RT-85	10179	·· VIS181 VIS181 2/1/81	
ER7-95	10181	· VIBISI VISTER 2/1/EI : 1	
R1-90	1-182	·· 1/13181 4/15/81 2/1/81 2/14/81	
ERT-195	10184	· VIVI81 VISI81 2/181 2/181 10:29	!
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EXT-110		· VI4181 VI5181 2/1/81 2/3/161 1+:26	
ERT-160		· · YISISH VISIS 3/1/81 2/18/81	
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Ext 494		* 1/15/81 1/15/81 2/1/81 2/18/81 10:18	
EGT-190 MN-4	10209	·· 1/1518 1/15181 7/1181 2/2181 10.56 ·· 1/16181 1/181 2/1/81 2/156 16:08	<u>├</u>
ER1-145		·· VI6/81 VI9/81 2/1/81 2/18/8 15:00	
E		· VI6121 /11/21 2/1/21 2/0-121 15:44	
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AT-19		~ 1/16181 1/19/21 2/13/21 12:04	
,	· · · · · · · · · · · · · · · · · · ·	* INCLUDE MATRIX SPIKES, BLANKS AND RE-RUNS HERE	
Field Blo	.k#1 10187	18/11/2 13/21/2 13/11/2 13/11/2 13/11/2	
	k#2 1013	" YIE/SI 1/19/81 -415/81	
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SAMPLE IDENTIFIER: 8040 COMPU/CHEM SAMPLE NUMBER: 3629

1. ANALYTICAL METHODOLOGY

THE SAMPLES WERE PREPARED AND ANALYZED ACCORDING TO TWO (2) CENERAL PROCEDURES: (1) "SAMPLING AND ANALYSIS PROCEDURES FOR SCREENING OF INDUSTRIAL EFFLUENTS FOR PRIORITY POLLUTANTS," REVISED APRIL 1977, US-EPA, AND (2) EPA METHOD 624, "ORGANICS BY PURGE AND TRAP," AND METHOD 625, "BASE/NEUTRALS, ACIDS, AND PESTICIDES," US-EPA, REVISED DECEMBER 3, 1979, FEDERAL REGISTER (GUIDELINES ESTABLISHING TEST PROCEDURES FOR THE ANALYSIS OF POLLUTANTS). THE LABORATORY PROCEDURES USED FOLLOW THOSE IN METHODS 608, 624, OR 625. QUALITY ASSURANCE, SAMPLE CUSTODY, AND DOCUMENT CONTROL PROCEDURES WERE FOLLOWED WHICH MEET OR EXCEED EPA REQUIREMENTS.

2. SAMPLE RECORD DATE

- A. RECEIVED/REFRIGERATED
- B. ORGANICS
 - 1. EXTRACTED 2. ANALYZED

VOLATILES BASE/NEUTRALS ACIDS PESTICIDES/PCBS

C. METALS

ANALYZED

NOT REQUESTED

01/15/81

01/16/81

01/21/81

01/22/81

02/04/81

02/04/81

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SAMPLE IDENTIFIER: 8040 COMPU/CHEM SAMPLE NUMBER: 3629

	COMPOUNDS	· · ·	CONCENTRATION (UG/L)	DETE	CTION LIMIT (UG/L)
2P.	BETA-BHC		SOL	D	10
4F.	CAMMA-BHC		80L	Þ	10
3P.	DELTA-BHC		EDL	D	10
6P.	CHLORDANE		BDL	D	10
7P.	4, 4'-00T		BDL	D	10
8P.	4,4'-DDE		BOL	5 D	10
9P.	4.4'-000		80L	5	10
10P.	DIELDRIN		SDL .	5	10
11P.	ALPHA-ENDOSULFAN		BDL	5	10
12P.	BETA-ENDOSULFAN		BDL	3	10
13P.	ENDOSULFAN SULFATE		EDL	2	10
14F.	ENDRIN		BOL	D	10
13P.	ENDRIN ALDEHYDE		EDL		10
16P.	HEPTACHLOR		EDL	5	10
17F.	HEPTACHLOR EPOXIDE	•	EDL	5	10
-18P.	PCB-1242		BDL		10
19P.	PCB-1254	•	BDL	•	10
20P.	PC8-1221		BDL		10
21P.	FCB-1232		BDL	•	10
22P.	PCB-1248		BDL	· ·	10
23P.	PCB-1260		BDL	þ	10
24P.	PCB-1016	:	BOL	•	10
25P.	TOXAPHENE		BDL		10
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SAMPLE IDENTIFIER: 8040 COMPU/CHEM SAMPLE NUMBER: 3629

3. PRIORITY POLLUTANT ANALYSIS REPORT

CET

	COMPOUNDS	CONCENTRATION (UG/L)		DETECTION LIMIT (UG/L)
10	ACROLEIN	BOL	P	199
ZV.	ACRYLONITRILE	BDL	0	100
JV.		>10000*	2	10
4V.		BDL	0	10
57.	BROMOFORM	BOL	2	10
6V.	CARBON TETRACHLORIDE	BOL		10
77.		BOL	0	10
87.		BOL	2	
9V.		BOL	5	• .
10V.		BOL	D	
	CHLOROFORM	BOL	Þ	
12V.		BOL	Þ	10
137.		BOL	D	10
	1, 1-DICHLOROETHANE	BOL)	10
	1, 2-DICHLOROETHANE	BOL	9	19
	1. 1-DICHLOROETHYLENE	BOL	Þ	10
177.		BOL	D	10
18V.		BOL	Ð	10
19V.		32		10
204.		BOL	D	10
217	METHYL CHLORIDE	BOL	Ż	10
227.	METHYLENE CHLORIDE	BOL	7	10
23V.	1, 1, 2, 2-TETRACHLOROETHANE	BOL	"	
	TETRACHLOROETHYLENE	BOL	ż	10
25V.	TOLUENE	1300*	ي بر	10
26V.	1. 2-TRANS-DICHLOROETHYLENE	BOL		10
27V.	1, 1, 1-TRICHLOROETHANE	8 0L	مذُ	10
28V.	1, 1, 2-TRICHLOROETHANE	BOL	5	10
297.	TRICHLOROETHYLENE	BOL	٠,	10
30V.	TRICHLOROFLUOROMETHANE	80L	5	10
31V.	VINYL CHLORIDE	BOL	•	10
1A.	2-CHLOROPHENOL	440	ワ	25
2A.	2, 4-DICHLOROPHENOL	500	1	25
·3A.	2, 4-DIMETHYLPHENOL	BOL	3	25
4 A .	4. 5-DINITRO-O-CRESOL	BOL	ر	250
5A.	2, 4-DINITROPHENOL	BOL	5	250
6A.	2-NITROPHENOL	BOL	5	23
7A.	4-NITROPHENOL	BOL	•	25
8 A .	P-CHLORO-M-CRESOL	BOL	1.2	25
9A.	PENTACHLOROPHENOL	BOL	• •	25
10A.	PHENOL	1700+		25
11A.		160		25
	ACENAPHTHENE	BOL		10
29.		BOL		10
38.	ANTHRACENE	BOL		10
			•	

BDL= BELOW DETECTION LIMIT * Saturated Ions

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SAMPLE IDENTIFIER: 8040 COMPU/CHEM SAMPLE NUMBER: 3629

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	COMPOUNDS	CONCENTRATION (UG/L)	
4E.	PENZIDINE	EDL	> 10
56.		BDL	
	EENZO (A) PYRENE	BDL	
	3, 4-BENZOFLUORANTHENE	BDL	
85.		BDL	
96.		BDL	
	SIS (2-CHLOROETHOXY) METHANE	SDL	
110	BIS (2-CHLOROETHYL) ETHER	BDL	
	SIS (2-CHLOROISOPROPYL) ETHER		
	BIS (2-ETHYLHEXYL) PHTHALATE	10	
	4-BROMOPHENYL PHENYL ETHER		D 10
145. 160	ATORUMUTICATE THENTL EINER	BOL	
135.	EUTYL EENZYL PHTHALATE 2-CHLORONAPHTHALENE	BDL	
168.	2-CHLORONAPHTHALENE	BOL	
175.	4-CHLOROPHENYL FHENYL ETHER	BOL	
188.	CHRYSENE	BOL	
175.	DIBENZO (A, H) ANTHRACENE	BOL	
208.	1.2-DICHLOROBENZENE	BOL	
216.	1, 3-DICHLOROBENZENE	BOL	9 10
22B.	1,4-DICHLORBENZENE	BDL	D 10
238.	3,3'-DICHLOROBENZÍDINE	BOL	
248.	DIETHYL PHTHALATE	BDL	
256.	1, 2-DICHLOROBENZENE 1, 3-DICHLOROBENZENE 1, 4-DICHLOROBENZENE 3, 3'-DICHLOROBENZIDINE DIETHYL PHTHALATE DIMETHYL PHTHALATE	BOL	-
26B.	DI-N-BUTYL PHTHALATE	BDL	
276.	2, 4-DINITROTOLUENE	BDL	
	2, 6-DINITROTOLUENE	BDL	
	DI-N-OCTYL PHTHALATE	BDL	
	1, 2-DIPHENYLHYDRAZINE	BDL	
	FLUORANTHENE	SOL	
	FLUORENE	80L	
	HEXACHLOROBENZENE	BOL	· > 10
	HEXACHLOROBUTADIENE	BOL BOL BOL	D 10
	HEXACHLOROCYCLOPENTADIENE		D 10
	HEXACHLOROETHANE		
308. 378.	INDENO (1,2,3-CD) PYRENE	BOL	
		BDL	> 25
388.	ISOPHORONE	BOL	-, 10
398.	NAFHTHALENE	180	> 10
40B.		BDL	> 10
41B.		BDL	
428.	N-NITROSODI-N-PROPYLAMINE	80L	> 10
438.	N-NITROSODIPHENYLAMINE	BDL	> 10
448.		BDL	5 10
45B.		BDL) 10
46B.	1, 2, 4-TRICHLOROBENZENE	80L	. 10
19.		BDL	- 10
2P.	ALPHA-BHC	BOL	~ 10
· •	<i>.</i>		·- · ·

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ERT ID # ERT-185 8054 Compuchen # ______

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ANALYSES PERFORMED

VOA	1 22 81; 5031 1	DILUTION - 1/23/81		
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B/N	123/61	180 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190 - 190		
PEST	1/23/81			

CPROVAL SAMPLE IDENTIFIER: 8054 COMPU/CHEM SAMPLE NUMBER: 3654

3. PRIORITY POLLUTANT ANALYSIS REPORT

	COMPOUNDS	CONCENTRATION (UG/L)		DETECTION LIMIT (UG/L)
1V.	ACROLEIN	• BDL	R	100
-	ACRYLONITRILE	BDL	R	100
	BENZENE	67000* **	R	10
-	BIS (CHLOROMETHYL) ETHER	BDL	ĸ	10
	BROMOFORM	BDL	R	10
_	CARBON TETRACHLORIDE	BDL	R	10
	CHLOROBENZENE	BDL	R	10
	CHLORODIBROMOMETHANE	BDL	R	10
	CHLORDETHANE	BDL	R	10
	2-CHLORDETHYLVINYL ETHER	BDL	R	10
	CHLOROFORM	BDL	R	10
	DICHLOROBROMOMETHANE	BDL	R	10
	DICHLORODIFLUOROMETHANE	BDL		10
	1, 1-DICHLORDETHANE	BDL	R	
	1.2-DICHLOROETHANE		R	10
	1, 1-DICHLOROETHYLENE	BDL	R	10
	1, 2-DICHLOROPROPANE	BDL	-	10
		BDL		10
	1, 3-DICHLOROPROPYLENE	BDL	R	10
	ETHYLBENZENE	73	R	10
	METHYL BROMIDE	BDL	R	10 III
	METHYL CHLORIDE	BDL	R	
227.		BDL	R	
	1, 1, 2, 2-TETRACHLORDETHANE	BDL	R	
	TETRACHLOROETHYLENE	6DL	R	
	TOLUENE	9000* **	R	10
	1, 2-TRANS-DICHLOROETHYLENE	BDL	R	10
	1, 1, 1-TRICHLORDETHANE	BDL	R	10
26V.	1, 1, 2-TRICHLORDETHANE	BDL	ĸ	10
	TRICHLORDETHYLENE	BDL	R	10
30V.	TRICHLOROFLUOROMETHANC	BDL	R	10
31V.	VINYL CHLORIDE	BDL	R	10
1A.	2-CHLOROPHENOL	140	R	. 25
2A.	2, 4-DICHLOROPHENOL	8000 **	R	25
JA.	2, 4-DIMETHYLPHENOL	BDL	R	25
4A.	4,6-DINITRO-O-CRESOL	·		
5A.	2.4-DINITROPHENOL	BDL	R	
	2-NITROPHENOL	BDL		
	4-NITROPHENOL	BDL		25
	P-CHLORO-M-CRESOL	BDL	A.	25
	PENTACHLOROPHENOL	BDL	ĸ	25
	PHENOL	3900 - **	R	25
	2, 4, 6-TRICHLOROPHENOL	83	R	25
	ACENAPHTHENE	BDL	R	10
	ACENAPHTHYLENE	BDL	R	
	ANTHRACENE	BDL		. 10
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BDL= BELOW DETECTION LIMIT * Value determined from 50:1 dilution ** Quantitated from Secondary Ion

SAMPLE IDENTIFIER: 8054 COMPU/CHEM SAMPLE NUMBER: 3654

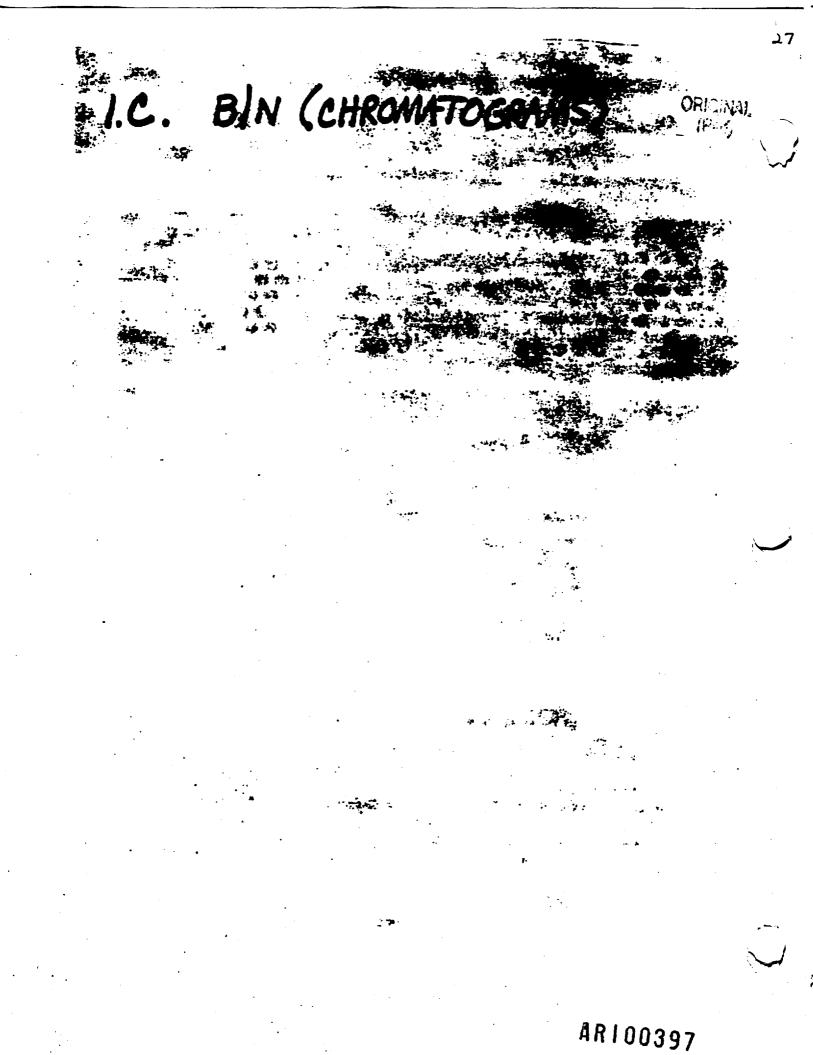
				Carl State
	COMPOUNDS	CONCENTRATION (UG/L)		DETECTION LIMIT (UG/L)
	BENZIDINE	BDL		
	BENZO (A) ANTHRACENE	BDL	R	
	BENZO (A) PYRENE	BDL	R	
	3. 4-BENZOFLUORANTHENE	BDL	R	
	BENZO (GHI) PERYLENE	BDL	R	
	BENZO (K) FLUORANTHENE	BDL	R	
	BIS (2-CHLOROETHOXY) METHANE BIS (2-CHLOROETHYL) ETHER	BDL	- r¢i	
	BIS (2-CHLORDISOPROPYL) ETHER	BDL	R	
	BIS (2-ETHYLHEXYL) PHTHALATE	BDL BDL	RR	
	4-BROMOPHENYL PHENYL ETHER	BDL	~ ~	
	ENTYL BENZYL PHTHALATE	BDL	R	
	2-CHLORONAPHTHALENE	BDL	R	
	4-CHLOROPHENYL PHENYL ETHER	BDL	R.	
	CHRYSENE	BDL	R.	
	DIBENZO (A, H) ANTHRACENE	BDL	ĸ	
20E.	1, 2-DICHLOROBENZENE	BDL	R	
21 B.	1, 3-DICHLOROBENZENE	BDL	.	
22B.	1,4-DICHLORBENZENE	BDL	Ř	
233.	3, 3'-DICHLORGBENZIDINE	BDL	R	
	DIETHYL PHTHALATE	BDL	R	10
	DIMETHYL PHTHALATE	BDL	R	10
	DI-N-BUTYL PHTHALATE	BDL	R	10
	2.4-DINITROTOLUENE	BDL	R	10
	2, 6-DINITROTOLUENE	BDL	R	10
•	DI-N-OCTYL PHTHALATE	BDL	え	10
	1, 2-DIPHENYLHYDRAZINE	BDL	R	. 10
	FLUCRANTHENE	BDL	R	10
	FLUGRENE	BDL	હ	
339.		BDL	R	10
343.	HEXACHLOROBUTADIENE	BDL	R	
	HEXACHLOROCYCLOPENTADIENE	BDL	R	
	HEXACHLOROETHANE	BDL	R	
	INDENO (1,2,3-CD) PYRENE Isophorone	BDL	R	
	NAPHTHALENE	BDL 410	RJ	
-	NITROBENZENE	BDL	- A	
	N-NITROSODIMETHYLAMINE	BDL	H	· • •
	N-NITROSODI-N-PROPYLAMINE	BDL		
	N-NITROSODIPHENYLAMINE	BDL	R R	
	PHENANTHRENE	BDL	к. 12	
	PYRENE	BDL	R	
	1, 2, 4-TRICHLOROBENZENE	BDL	R	
	ALDRIN	BDL	R	
22.		BDL	R	
			~	

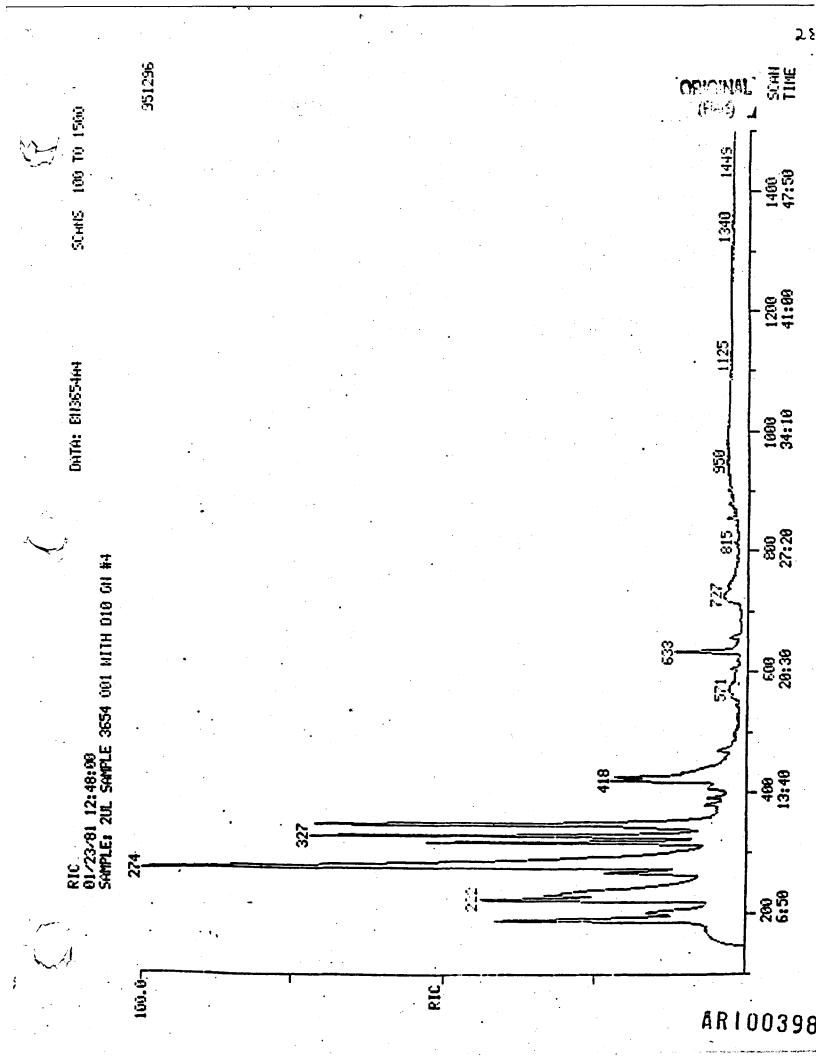
BDL= BELOW DETECTION LIMIT

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SAMPLE IDENTIFIER: 8054 COMPU/CHEM SAMPLE NUMBER: 3654

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	•		annal .		L (Faire Sa
	COMPOUNDS	•	CONCENTRATION (UG/L)		DETECTION LIMIT
3P.	BETA-BHC		BDL	R	10
4P.	GAMMA-BHC		BDL	R	10
5P.	DELTA-BHC	•	BDL	IC.	10
6P.	CHLORDANE	1	BDL	Ř	10
7P.	4,4'-DDT		BDL	R	10
8P.	4,4'-DDE		BDL	R	10
9P.	4,4'-DDD		BDL	R	10
10P.	DIELDRIN		BDL	R	10
11P.	ALPHA-ENDOSULFAN	•	BDL	R.	10
12P.	BETA-ENDOSULFAN		BDL	R	10
13P.	ENDOSULFAN SULFATE		BDL	R	10
14P.	ENDRIN		BDL	R	10
15P.	ENDRIN ALDEHYDE		BDL	8	10
16P.	HEFTACHLOR		BDL	R	10
17P.	HEPTACHLOR EPOXIDE		BDL	R	10
18P.	PCB-1242		BDL	R	10
19P.	PCB-1254	•.	BDL	R	10
20P.	PC8-1221	•	BDL	R	10
21P.	PCB-1232	. •	BDL	A	10
22P.	PCB-124E		BDL	R	10
23P.	FCB-1260		BDL	Ŕ.	10
24P.	PCB-1016		BDL	ĸ	10
25P.	TOXAFHENE		BDL	R	10





FINNIGAN TARGET COMPOUND ANALYEIS QUANTITATION REPORT FILE: BN365444

DATA: BN3654A4 TI 01/23/81 12:48:00 SAMPLE: 201 SAMPLE 3654 001 WITH D10 ON #4 SUBMITTED BY: #4 ANALYST: AC

AMOUNT=AREA(HCHT) * REF AMNT/(REF. AREA(HGHT)* RESP. FACT) RESP. FAC. FROM LIBRARY ENTRY

NO NAME D10-ANTHRACENE (INTERNAL STANDARD) 1 2 D-5 NITROBENZENE (SURROGATE STANDARD) 2-FLUOROBIPHENYL (SURROGATE STANDARD) З D8-NAPHTHALENE (SUFROGATE STANDARD) 4 5 D-10 ANTHRACENE (INTERNAL STANDARD) 6 1, 3-DICHLOROBENZENE 7 HEXACHLOROETHANE 8 NITROBENZENE 9 NAPHTHALENE 10 2-CHLORONAPHTHALENE 2, 6-DINITROTOLUENE 11 4-CHLOROPHENYL PHENYL ETHER 12 13 1, 2-DIPHENYLHYDRAZINE 14 N-NITROSODIFHENYL MINE 15 HEXACHLOROSENZEN 16 ANTHRACENE 17 PHENANTHRENE 18 4-BROMOPHENYL PHENIL ETHER 19 FLUORANTHENE 20 CHRYSENE D-10 ANTHRACENE (INTERNAL STANDARD) 21 22 1, 4-DICHLORGEENTENE 23 BIS (2-CHLORGETHYL' ETHER 24 HEXACHLOROBUTADIENE 25 BIS (2-CHLORCETHOXY)METHANT 26 ACENAPHTHYLENE 27 FLUGRENE 28 2, 4-DINITROTCLUENE 29 PYRENE 30 BENZO(E)FLUORANTHEME 31 BENZO(A)PYREME 32 BENZIDINE 33 1, 2, 4-TRICHLOROBENZENE 34 DIBENZO (A, H) ANTHRACENE 35 BENZO (A) ANTHRACENE 34 INDENO (1, 2, 3-CD) FYRENE 37 D-10 ANTHRACENE (INTERNAL STANDARD) 32 1, 2-DICHLORGEENZENE 39-ISOPHORONE 40 ACENAPHTHENE 41 DIMETHYLFHTHALATE 42 DIETHYLPHTHALATE 43. DI-N-BUTYLPHTHALATE 44 BUTYLBENEYE PHTHA: ATE 45 BIS (2-ETHYLHEAND OPHTHALSOF 4ć -DI-CCTYLEHTHALAT:

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-	NO	NAME	, [.]					$T_{1}^{2} \in \mathcal{F}_{1}^{2} \oplus \mathcal{F}_{2}^{2}$	•					
	47	EIS(2-CHLC						•	c.		<u>.</u>			
	48	N-NITROSOD			INE				N.	ن د با توقع من				
1	49	BENZO (K)								_ (his	9 .			
•	50	3, 3'-DICHL								-				
	51	BENZO (G. H												
	52	BIS (2-ETH	YLHEXYL) FHTH	IALATE	(SEC	CONDARY ION)						
			****		007	54								
	NO	M/E SCAN	TIME 21:38	REF	. 006		TH	AREA(HGHT)					TO	
	1	188 633	21:30	T . T	. 000	A	BB	169577.	100.	000	UG/L	10.	83 .	
	. 5	NOT FOUND	44.54				-	10000				1		
	3	172 425	14:31			A .	68	100032.	43.	781	UG/L*	∠ 4 .	74	
	4		11:00			A		136612.	50.	526	UG/L-	5.	47	
	5	188 633	21:38	2 1	999	A	BE	169577.	100.	000	UG/L	10.	83	
	6	NOT FOUND												
	7	NOT FOUND					· ·	•						
	8	NOT FOUND			_				•					
	9		11:10	5 0). 517	A	68	1095490. 413,	,6 368.	064	UG/L	39.	87 -	, y e
	10	NOT FOUND	•		•			•	* x	6-	1 pm			
	11	NOT FOUND	•							ſ	- A Barrer			
	12	NOT FOUND							-			_		
		77 556					BB	1491.			UG/L			
			19:28	5 0	. 900	A	88	1082.	Ø.	694	UG/L	Ø.	68	
	15	NOT FOUND	64.64	E 0		•		07/0	-			•		
	16	178 632	21:36	5 6	000		EB				UG/L			
	17	178 632	21:36	5 0	. 778	A .	BB	2763.	0.	472	UG/L	Ø.	65	
	18	NOT FOUND												
	19	NOT FOUND				•						•		
1	20 21	NOT FOUND 188 633	21:38	D 1 1	0.070		-	140677	100		115 21		~~	
1	55	188 633 Not Found	E1:30	<u> </u>	000	F.	55	169577.	100.	666	UG/L	10	83	
	23	NOT FOUND	•											
	24	NOT FOUND												
		•	11-27	7 0		t.	ບຣ	2533.	2	541	UG/L	<u>م</u>	37	
		152 484	16:32		. 765						UG/L		02	
	27	166 537	18:21		. 848		BE	96.			UG/L	-	00	
	28	NOT FOUND				-		70.	, v .	V+1		C.	v.e	
	29	202 769	26:16	21 1	215	Α.	86	2156.	6	549	UG/L	0	06	
	30	NOT FOUND		4					Ψ.	u -77				
	31	NOT FOUND			•									
	32	NOT FOUND	•								•	-		
	33	NOT FOUND					·			•	•			
	34	NOT FOUND												
	35	NOT FOUND						•			•			
	36	NOT FOUND							ė.					
	37	188 633	21:38	37 1	. 000	A	68	169577.	100.	000	UG/L	10.		
	38	146 216			. 341		BB	669.			UG/L			
	39	NOT FOUND	_	-										
	40	NOT FOUND				-								
-	41	NOT FOUND							•					
•	42	NOT FOUND			•									
·	43	149 686	23: 26	37 1	. 084	Α	BB	3314.	0.	573	UG/L	e	05	
~ <u>.</u> .	44	149 835	28:32		. 31		BF:	1841			UGZE	Ċ.	65	
-	45	145 855	27:15		. 33		65	14363.			UG.AL		<u> </u>	
~	46	145 513	31.12		4-		EF	617.			UG/L		01	
• •	47		6 C-		273		VE:	13426.			UGZO	C,	ė 1	
•	· 46	130 304	16 22		450		E:E	3892	41	06-	06.4	4	45 .	א.
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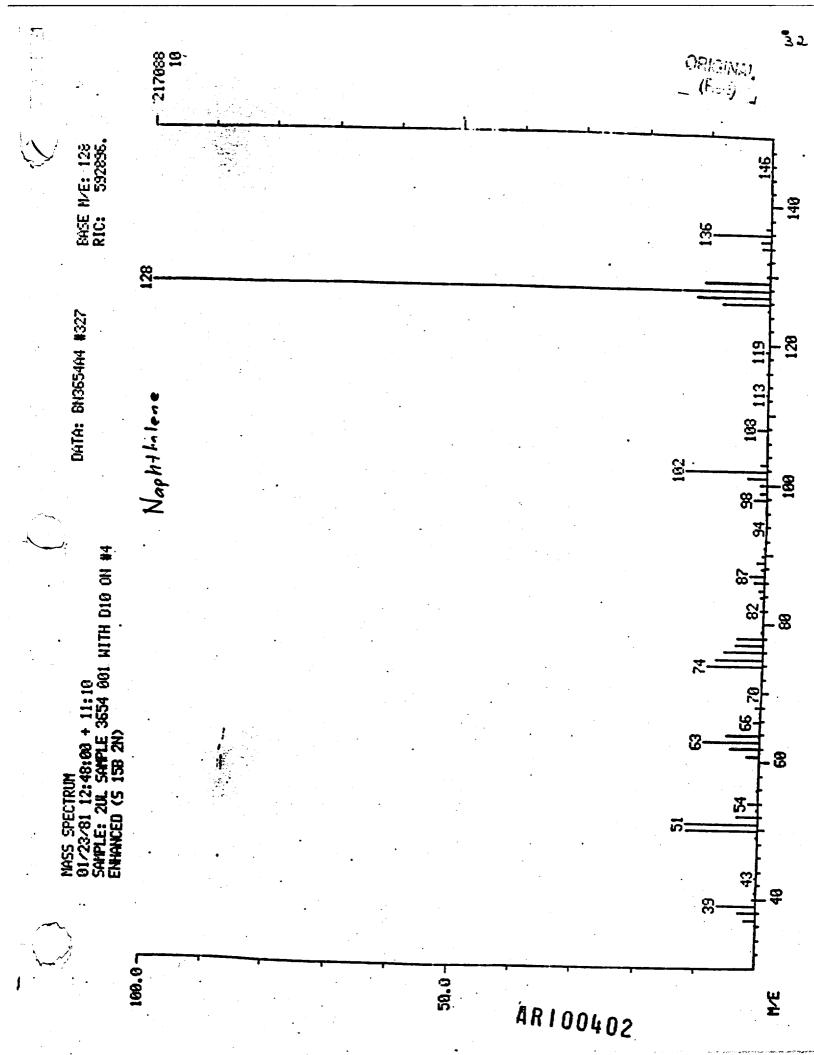
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							OF! MADE	:
NO 49	M/E SCAI Not foun		REF	RRT	METH	AREA (HGHT)	AMOUNT	XTOT
50 51 52	NOT FOUN NOT FOUN 167 85	D	37	1. 352	A BB	4132.	3. 449 UG/L	0 3
	NO Pest.	or Tadd	form	. اص	,	•		- •

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33 E CMPLICH 3655 ANALYSE'S PERFORM VOA 1 23 81; 2011 DILUTION + 123 81 5011 DILUTION + 123 81 ACID III B/N 1/23/81 PEST 123/81

SAMPLE IDENTIFIER: 8059 COMPU/CHEM SAMPLE NUMBER: 3655

3. PRIORITY POLLUTANT ANALYSIS REPORT

	COMPOUNDS	CONCENTRATION (UG/L)	DETECTION LIMIT (UG/L)
1V.	ACROLEIN	BOL	R 100
27.	ACRYLONITRILE	BOL	R 199
3V.	BENZENE	4000001	r 10
4V.	BIS (CHLOROMETHYL) ETHER	BOL	« <u>10</u>
5V.	BROMOFORM	BOL	
	CARDON TETRACHLORIDE	BOL	R 10
	CHLOROBENZENE	ODL	R 19
	CHLORODIBROMOMETHANE	BOL	ñ 10
	CHLOROETHANE	BOL	R 10
10V.		BOL	R 10
	CHLOROFORM	BOL	R 10
	DICHLOROBROMOMETHANE	BOL	R 10
	DICHLORODIFLUOROMETHANE	BOL	`R 10 3
	1. 1-DICHLOROETHANE	SOL	R 10
	1, 2-DICHLOROETHANE	BOL	A 10
444	1, 1-DICHLOROETHYLENE	BOL	2 10 2 10
	1, 2-DICHLOROPROPANE	801	R 10
187.	1, 3-DICHLOROPROPYLENE	BOL	R 10
19V.	ETHYLBENZENE	32	
20V.	METHYL BROMIDE	JE	
	METHYL CHLORIDE	32 32 80L 80L 80L 80L	R 10 R 10
		BUL	
22V.	METHYLENE CHLORIDE	BDL	r 10 r 10
_	TETRACHLOROETHYLENE		K 10
	TOLUENE	5800**	R 10
26V.			R 10
277.		BDL	R 10
	1, 1, 2-TRICHLOROETHANE	BDL	R 10
297.	TRICHLOROETHYLENE	BDL	R 10
30V.	TRICHLOROFLUOROMETHANE	BDL	2 10
	VINYL CHLORIDE	BOL	Q. 10
	2-CHLOROPHENOL	71	R 25
	2.4-DICHLOROPHENOL	720	× 25
	2.4-DIMETHYLPHENOL	BDL	R 25
44.	4. 6-DINITRO-O-CRESOL	BDL	x 250°
5A.	2.4-DINITROPHENOL	BOL	r 250
6A.	2-NITROPHENOL	BOL	r 25
7A.	4-NITROPHENOL	BOL	r 25
BA.	P-CHLORO-M-CRESOL	BOL	r 25
9A.	PENTACHLOROPHENOL	BOL	r 25
10A.		2500	r. 25
11A.		47	ir 25
1B.		BOL	8 10
28.	ACENAPHTHYLENE	BOL	- 10
3B.	ANTHRACENE	BDL	÷ 10

BDL= BELOW DETECTION LIMIT * Value determined from 50:1 dilution * Value determined from 20:1 dilution

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ORIGINAL (Fig.)

SAMPLE IDENTIFIER: 8059 COMPU/CHEM SAMPLE NUMBER: 3655 35

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	COMPOUNDS BENZIDINE BENZO (A) ANTHRACENE BENZO (A) PYRENE 3. 4-BENZOFLUORANTHENE BENZO (CHI) PERYLENE BENZO (K) FLUORANTHENE BIS (2-CHLOROETHOXY) METHANE BIS (2-CHLOROETHOXY) METHANE BIS (2-CHLOROISOPROPYL) ETHER BIS (2-CHLOROISOPROPYL) ETHER BIS (2-CHLOROISOPROPYL) ETHER BIS (2-CHLOROISOPROPYL) ETHER BUTYL BENZYL PHTHALATE 2-CHLORONAPHTHALENE 4-GROMOPHENYL PHENYL ETHER CHRYSENE DIBENZO (A H) ANTHRACENE 1. 2-DICHLOROBENZENE 1. 3'-DICHLOROBENZENE 3. 3'-DICHLOROBENZIDINE DIETHYL PHTHALATE DIMETHYL PHTHALATE 2. 4-DINITROTOLUENE 2. 4-DINITROTOLUENE 2. 4-DINITROTOLUENE DI-N-OCTYL PHTHALATE 1. 2-DIPHENYLHYDRAZINE FLUORANTHENE FLUORENE HEXACHLOROBENZENE HEXACHLOROBENZENE HEXACHLOROBURADIEN	CONCENTRATION (UG/L)		DETECTION LIMIT (UG/L)
48.	BENZIDINE	BOL	R	10
58.	BENZO (A) ANTHRACENE	BOL	R	10
68.	BENZO (A) PYRENE	BOL	R	10
78.	3. 4-BENZOFLUORANTHENE	BOL	R	10
8 8 .	BENZO (GHI) PERYLENE	BOL	R	25
98.	BENZO (X) FLUORANTHENE	BOL	R	10
198.	BIS (2-CHLOROETHOXY) METHANE	BOL	ぇ	10
118.	BIS (2-CHLOROETHYL) ETHER	BOL	R	19
128.	BIS (2-CHLOROISOPROPYL) ETHER	80L	×	10
138.	BIS (2-ETHYLHEXYL) PHTHALATE	BOL	.•	10
148.	4-BROMOPHENYL PHENYL ETHER	BOL	Ŕ	• 10
158.	BUTYL BENZYL PHTHALATE	BOL	R	10
16B.	2-CHLORONAPHTHALENE	801	R	10
178.	4-CHLOROPHENYL PHENYL ETHER	BDL	R	10
188.	CHRYSENE	BOL	ぇ	10
198.	DIBENZO (A.H) ANTHRACENE	BOL	R.	25
20B.	1, 2-DICHLOROBENZENE	BOL	R	10
219.	1. 3-DICHLOROBENZENE	BOL	R	10
229.	1 DICHLORBENZENE	BOL	R	10 -
238.	3. J'-DICHLOROBENZIDINE	BOL	R	10
248.	DIETHYL PHTHALATE	BOL	R	. 10
258.	DIMETHYL PHTHALATE	BOL	R	10
26B.	DI-N-BUTYL PHTHALATE	BOL	, in the second s	10
27B.	2. 4-DINITROTOLUENE	BOL	R	10
288.	2. 6-DINITROTOLUENE	BOL	2	10
278.	DI-N-OCTYL PHTHALATE	BOL	R	10,
308.	1, 2-DIPHENYLHYDRAZINE	BOL	Ŕ	10
318.	FLUORANTHENE	BDL		10
32B.	FLUORENE	BDL	2	10
33B.	HEXACHLOROBENZENE	BDL	3	10
348.	HEXACHLOROBUTADIENE	BOL	a	10
35B.	HEXACHLOROCYCLOPENTADIENE	BDL	3	10
368.	HEXACHLOROETHANE	BOL	ିତ	10
37B.	INDENO (1, 2, 3-CD) PYRENE	BOL	Â	25
388	ISOPHORONE	BDL	R	10
39B.	NAPHTHALENE	90	1	10 .
40B.	NITROBENZENE	BOL	્રે	10
418.	N-NITROSODIMETHYLAMINE	BOL	R	10
428	N-NITROSODI-N-PROPYLAMINE	BOL	R	10
438.		BOL	2	10
448.		BOL		
458.		BOL	~	10
46B.		BOL	g	10
	ALDRIN	BOL	2	10
2P.	ALPHA-BHC	80L		,
6F .		BUL	Ś	70

BDL= BELOW DETECTION LIMIT

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SAMPLE IDENTIFIER: 8059 COMPU/CHEM BAMPLE NUMBER: 3655

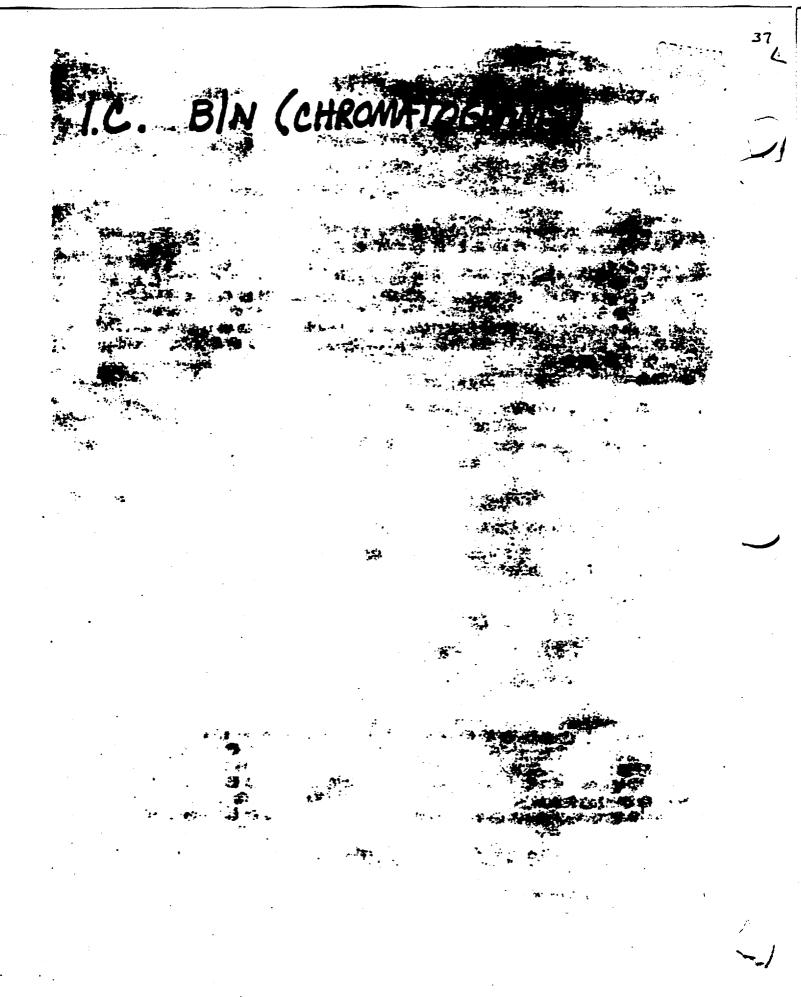
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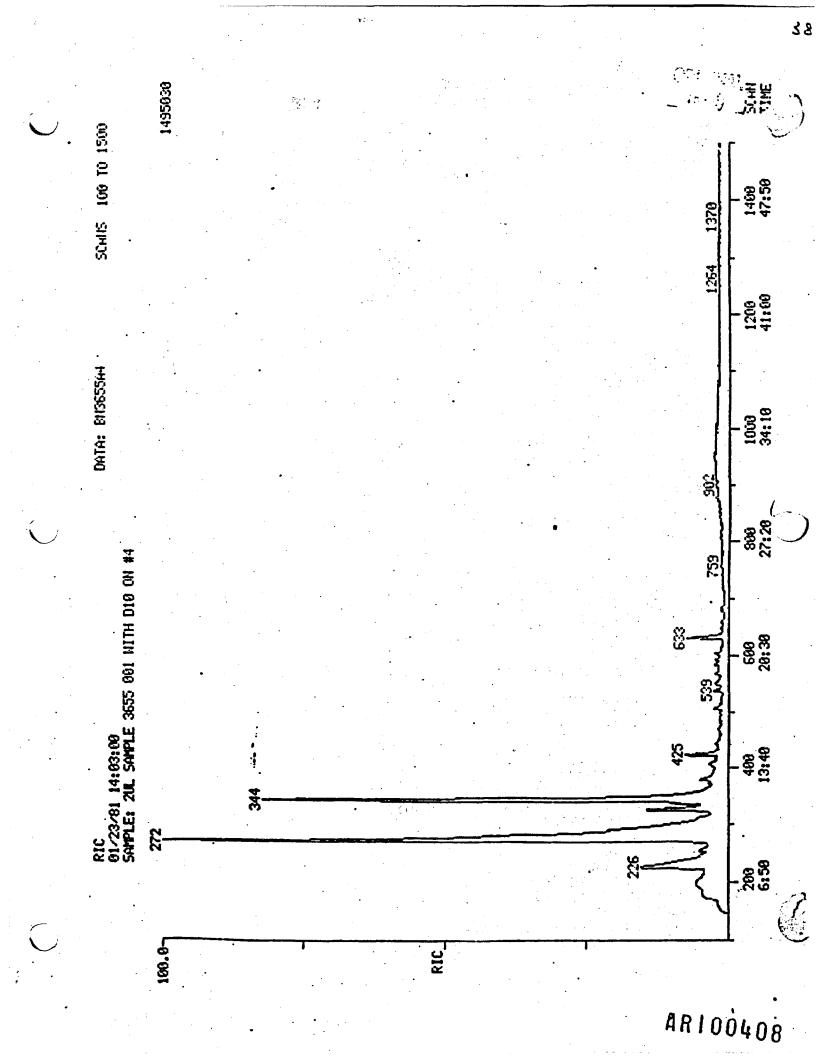
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	COMPOUNDS	CONCENTRATION (UG/L)	DETECTION LIMIT (UG/L)
3P.	BETA-BHC	80L	R 10
4P.	CAMMA-BHC	BOL.	r 10
5P.	DELTA-BHC	BOL	r 10
6P.	CHLORDANE	BOL	R 10
· 7P.	4, 4'-DDT	BOL	R 10
8P.	4, 4'-0DE	BDL	R 19
9P.	4, 4'-DDD	BOL	R 10
10P.	DIELDRIN	BOL	R 10
11P.	ALPHA-ENDOBULFAN	BOL	R 10
12P.	BETA-ENDOSULFAN	BOL	r 10
13P.	ENDOSULFAN SULFATE	BOL	R 10
14P.	ENDRIN	BOL	R 10
1SP.	ENDRIN ALDEHYDE	BÖL	R 10 ±
16P.	HEPTACHLOR	BOL	
17P.	HEPTACHLOR EPOXIDE	BOL	
18P.	PCB-1242		r 10 5
		BOL	
19P.	PCB-1254	BOL	R 10
20P.	PCB-1221	BOL	R 10
21P.	PCB-1232	BDL	r 10
22P.	PCB-1248	BDL	R 10
23P.	PCB-1260	BDL.	#. 10
24P.	PCB-1016	BDL	a 10
25P.	TOXAPHENE	BOL	10

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1. n g	FINNIGAN TARGET COMPOUND ANALYSIS QUANTITATION REPORT FILE BN3655A4	•
1		•
8 J.	DATA: BN3655A4. TI	-
	01/23/81 14:03:00	
	SAMPLE: 2UL SAMPLE 3655 001 WITH D10 ON #4 SUBMITTED BY: #4 ANALYST: AC	
	SARUTIED RI: 44 HAVELOI VC	
	AMOUNT #AREA(HCHT) # REF. AMNT/(REF. AREA(HCHT) # F	ESP. FACT)
	RESP. FAC. FROM LIBRARY ENTRY	
	NO NAME 1 D19-ANTHRACENS (INTERNAL STANDARD)	
	2 D-5 NITROBENZENE (SURROGATE STANDARD)	
	3 2-FLUOROBIPHENYL (SURROGATE STANDARD)	
	4 DB-NAPHTHALENE (SURROGATE STANDARD)	
	5. D-10 ANTHRACENE (INTERNAL STANDARD)	· ·
	6 1,3-DICHLOROBENZENE 7 HEXACHLOROETHANE	
	7 HEXACHLOROETHANE 8 NITROBENZENE	
	9 NAPHTHALENE	
	10 2-CHLORONAPHTHALENE	
	11 2.6-DINITROTOLUENE	
	12 4-CHLOROPHENYL PHENYL ETHER	•
· .	13 1.2-DIPHENYLHYDRAZINE	•
	14 N-NITROSODIPHENYLAMINE 15 Hexachlorobenzene	
	15 ANTHRACENE	
	17 PHENANTHRENE	
	18 4-BROMOPHENYL PHENYL ETHER	
	17 FLUORANTHENE	
	20 CHRYSENE	
	21 D-10 ANTHRACENE (INTERNAL STANDARD) 22 1.4-DICHLOROBENZENE	
	23 BIS (2-CHLOROETHYL) ETHER	· ·
	24 HEXACHLOROBUTADIENE	1. A.
	25 BIS (2-CHLOROETHOXY)METHANE	
	26 ACENAPHTHYLENE 27 FLUORENE	
	27 FLUORENE 28 2,4-DINITROTOLUENE	
	29 PYRENE	
•	30 BENZO(B)FLUORANTHENE	1
	31 BENZO(A)PYRENE	·
•	32 BENZIDINE	
•	33 1,2,4-TRICHLOROBENZENE 34 DIBENZO (A,H) ANTHRACENE	
	35 BENZO (A) ANTHRACENE	
	36 INDENO (1, 2, 3-CD) PYRENE	
	37 D-10 ANTHRACENE (INTERNAL STANDARD)	
	38 1, 2-DICHLOROBENZENE	· .
	39 ISOPHORONE	•
	40 ACENAPHTHENE 41 DIMETHYLPHTHALATE	
-	42 DIETHYLPHTHALATE	
ř.	43 DI-N-BUTYLPHTHALATE	
(A)	44 BUTYLEENZYLFHTHALATE	
S)	45 BIS (2-ETHYLHEXYL) FHTHALATH	
	46 DI-OCTYLFHTHALATE	•

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18 - 5 NAME NO BIS(2-CHLOROISOPROFYL)ETHER 47 N-NITROSODI-N-PROPYLAMINE 48 BENZO (K) FLUORANTHENE 49 3. 3'-DICHLOROBENZIDINE 50 BENZO (G, H, I) PERYLENE 51 BIS (2-ETHYLHEXYL) PHTHALATE (SECONDARY ION) 52 NO M/E SCAN TIME REF RRT METH AREA(HGHT) AMOUNT %TOT 633 21:38 1.000 A BB 1 - 188 1 156356. . 100.000 UC/L 16. 57 NOT FOUND 21 2 1220 3 172 425 14:31 1 0.671 A BB 105120. 49.898 UG/L 8. 27 326 11:08 A BV 4 136 1 0. 515 144745. 58.061 UG/L 9.62 168 633 21:38 5 1.000 A BB 5 156356. 100.000 UG/L 16. 57 NOT FOUND 6 NOT FOUND NOT FOUND 8 128 328 Q 11:12 5 0.518 A BB 223414. 90 81. 409 UC/L 13.49 -705. NOT FOUND 10 NOT FOUND 11 12 NOT FOUND 13 A VV 77 551 19:50 0. 670 13861. 3. 534 UC/L 0.59 NOT FOUND 14 15 NOT FOUND 21:34 5 0. 997 16 178 631 A BE 1094. • ' 0.211 UG/L 0.04 17 631 178 21:34 5 6. 997 A BB 1094. 0. 211 UC/L 0.04 NOT FOUND 18 19 NOT FOUND NOT FOUND 20 21 188 633 21:38 21 1. 600 A EB 156356. 100 000 UC/L 16 57 22 NOT FOUND 23 NOT FOUND 24 NOT FOUND 25 NOT FOUND 26 152 484 487. 0.179 UG/L 16:32 21 0.765 A BB 0.03 27 528 166 18:02 21 0.834 A EB 1366. 0. 633 UG/L 0.10 28 NOT FOUND 29 0.070 UG/L 202 770 26:18 21 1. 216 A BB 325. 0.01 NOT FOUND 30 31 NOT FOUND 32 NOT FOUND 33 NOT FOUND NOT FOUND 34 35 NOT FOUND NOT FOUND 36 37 188 633 21:38 37 1.000 A BB 156356. 100.000 UC/L 16. 57 38 NOT FOUND 39 NOT FOUND 40 NOT FOUND 3. 815 UG/L 41 163 508 17:21 0.803 A EB 8509. 37 0.63 42 149 554 18:56 37 0.875 A BV - 782. 0. 267 UG/L 0.04 43 149 23:26 1. 084 A BB 1. 542 UG/L 0.26 686 37 8229. 835 37 44 149 28, 33 1. 319 A BE 1019. 0. 426 UG/L 0 07 45 149 655 29.13 1. 351 4 EE 2324. 0. 595 UG/L 0.10 37 145 37 305. 46 913 1 442 0.047 UC/L 31:12 A 68 C CL 27. 45 229 3927. 1. 798 UG/L 0 50 47 7.49 0.342 A EE 48 NOT FOUND + 9 fr 1 m - 1/1.

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AR100410

NO	M/E 50		REF	RRT	METH	AREA (HGHT)	AMOUNT	•	%TOT	
49 50 . 51	NOT FOU NOT FOU NOT FOU	JND			•		•	· · ·		
52	167 E	356 29:15	37	1. 352	A BB	875.	0. 792	UG/L	0 13	

42 1 46090 10 \rangle 845E N E: 128 RIC: 155904. <u>z</u> 160 140 DHTH: E113655HH #323 ਨ 8 120 <u>8</u> 102 Nophiliclene 8 റ്റ. 5 001 WITH DIO GN #4 8 11:12 ដ 8 MHSS SPECTRUM SAMPLE: **31/23/8**] ន្លៈ **6** 23 106.07 50.6ž ۶ AR100412

43 5 ERT- 205 8060 ERT ID COMPUCHEM # 3662 ANALYSES VOA 113081; 20:1 DILUTION + 1/308 ACID 21116 BIN 120/81 PEST 128181 AR100413

SAMPLE IDENTIFIER: 8060 COMPU/CHEM SAMPLE NUMBER: 3662

3. PRIORITY POLLUTANT ANALYSIS REPORT

•	COMPOUNDS	CONCENTRATION (UG/L)	DETECTION LIMIT (UG/L)
1V.	ACROLEIN	BOL	r 160
2V.	ACRYLONITRILE	BOL	2 100
JV.	BENZENE	>10000 **	J 10
4V.	BIS (CHLOROMETHYL) ETHER	BOL	R . 10
-5V.	BROMOFORM	É BOL	2 10
	CARBON TETRACHLORIDE	BDL	R 10
7V.	CHLOROBENZENE	BOL	R 10
EV.	CHLORODIBROMOMETHANE	BOL	R 10
9V.	CHLOROETHANE	· BDL	R 10
	2-CHLOROETHYLVINYL ETHER	BOL	R 10
	CHLOROFORM	BOL	R 10
	DICHLOROBROMOMETHANE	BDL	3 10
	DICHLORODIFLUOROMETHANE	. BDL	R 10
	1, 1-DICHLOROETHANE	BDL	R 10
	1, 2-DICHLOROETHANE	BDL	R 10.
	1, 1-DICHLOROETHYLENE	BOL	R 10
	1. 2-DICHLOROPROPANE	BOL	R 10
18V.	1, 3-DICHLOROPROPYLENE	BDL	x 10
19V.	ETHYLBENZENE	BDL	~ 10
204.	METHYL BROMIDE	BOL	·~ 10
21V.	METHYL CHLORIDE	BDL	R . 10
22V.	METHYLENE CHLORIDE	BOL	8 10
	1, 1, 2, 2-TETRACHLOROETHANE	BOL	- 10
	TETRACHLOROETHYLENE	BDL	~ 10
25V.	TOLUENE	14	J 10
26V.	1, 2-TRANS-DICHLOROETHYLENE	BDL	R 10
27V.	1, 1, 1-TRICHLOROETHANE	BDL	R 10
28V.	1, 1, 2-TRICHLOROETHANE	BDL	R 10
29V.	TRICHLOROETHYLENE	BOL	R 10
30V.	TRICHLOROFLUOROMETHANE	BDL	e 10
31V.	VINYL CHLORIDE	BDL	·~. 10
1A.	2-CHLOROPHENOL	>1000+	R 25
2A.	2, 4-DICHLOROPHENOL	1100	r 25
ЗA.	2, 4-DIMETHYLPHENOL	- BDL	r 25
4A.	4, 6-DINITRO-O-CRESOL	BDL	r 250
5 A.	2, 4-DINITROPHENOL	BOL	r 250
6A.	2-NITROPHENOL	BDL	* 25
7A .	4-NITROPHENOL	BOL	
BA.	P-CHLORO-M-CRESOL	BDL	v 25
9A.	PENTACHLOROPHENOL	BDL	、 25
10A.	PHENOL	>10000 *	r 25
11A .	2, 4, 6-TRICHLOROPHENOL	>10000 *	r 25
18.	ACENAPHTHENE	BDL	<u>e 10</u>
28.	ACENAPHTHYLENE	BOL	÷ 10
38.	ANTHRACENE	BDL	~ 10

BDL= BELOW DETECTION LIMIT * Saturated Ion ** Value determined by Secondary Ion in a 20:1 dilution

AR100414

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SAMPLE IDENTIFIER: 8060 COMPU/CHEM SAMPLE NUMBER: 3662

		. •	•
• •	COMPOUNDS	CONCENTRATION (UG/L)	DETECTION LIMIT (UG/L)
78. 99. 109. 129. 139. 148. 159. 148. 159. 148. 159. 148. 159. 148. 159. 148. 159. 148. 159. 149. 159. 159. 159. 159. 159. 159. 159. 15	BENZIDINE BENZO (A) ANTHRACENE BENZO (A) PYRENE 3.4-BENZOFLUORANTHENE BENZO (GHI) PERYLENE BENZO (K) FLUORANTHENE BIS (2-CHLOROETHOXY) METHANE	(UC/L) BOL BOL BOL BOL BOL BOL BOL 20 EOL BOL BOL BOL BOL BOL BOL	(UC/L) R 10 R 10
428. 438. 448. 458. 468. 1P.	N-NITROSODIHETHYLAMINE N-NITROSODI-N-PROPYLAMINE N-NITROSODIPHENYLAMINE PHENANTHRENE PYRENE 1, 2, 4-TRICHLOROBENZENE ALDRIN ALPHA-EHC	80L 80L 80L 80L 80L 80L 80L	•

BDL= BELOW DETECTION LIMIT

AR100415

46.

SAMPLE IDENTIFIER: 8060 Compu/Chem Sample Number: 3662

•	COMPOUNDS	CONCENTRATION (UG/L)		DETECTION LIMIT (UG/L)
3P.	BETA-BHC	BOL	R	10
4P.	GAMMA-BHC	BOL	R	10
5P.	DELTA-BHC	BOL	R	10
6P.	CHLORDANE	BOL	R	10
7P.	4, 4'-DDT	BOL	R	10
8P.	4, 4'-DDE	BOL	R	10
9P.	4.4'-000	BOL	R	10
10P.	DIELDRIN	BDL	R	10
11P.	ALPHA-ENDOSULFAN	BOL	Ŕ	10
12P.	BETA-ENDOSULFAN	BOL	12	10
13P.	ENDOSULFAN SULFATE	BOL	R	10
14P.	ENDRIN	BDL	R	10
15P.	ENDRIN ALDEHYDE	BOL	R	10
16P.	HEPTACHLOR	BOL	R	10
17P.	HEPTACHLOR EPOXIDE	BOL	R	10
18P.	PCB-1242	BDL	15	10
17P.	PCB-1254	BOL		10
20P.	PCB-1221	BOL	•	10
21P.	PCB-1232	801	. • .	10
22P.	PCB-1248	BDL	×	10
23P.	PCB-1260	BDL	÷	10
24P.	PC8-1016	BOL	•	10
25P.	TOXAPHENE	BOL		10
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